# Learning Curve Theory

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#### Abstract

Recently a number of empirical "universal" scaling law papers have been published, most notably by OpenAI. 'Scaling laws' refers to power-law decreases of training or test error w.r.t. more data, larger neural networks, and/or more compute. In this work we focus on scaling w.r.t. data size n. Theoretical understanding of this phenomenon is limited, except in finite-dimensional models for which error typically decreases with  $n^{-1/2}$  or  $n^{-1}$ , where n is the sample size. We develop and theoretically analyse the simplest possible (toy) model that can exhibit  $n^{-\beta}$  learning curves for arbitrary power  $\beta > 0$ , and determine to which extent power laws are universal or depend on the data distribution or loss function: Roughly, learning curves exhibit a power law with  $\beta = \frac{\alpha}{1+\alpha}$  for Zipf-distributed data with exponent  $1 + \alpha$ , independent of the choice of loss. Furthermore, noise rapidly deteriorates/improves in instantaneous/time-averaged learning curves for increasing n, suggesting that model selection should be based on cumulative (AUC) or time-averaged error, not final test error.

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## Power Laws in Large-Scale Machine Learning

- 'Mantra' of modern machine learning: 'bigger is better'.
- The larger and deeper *Neural Networks (NNs)* are, the more data they are fed, the longer they are trained, the better they perform.
- *Quantification:* Test error decreases as a *power law*, with the *data size*, with the *model size* (number of NN parameters), as well as with the *compute budget* used for training ...
- assuming one factor is not "*bottlenecked*" by the other two factors, -or- all three factors are increased appropriately in tandem.
- *Note: Subtract irreducible error* due to intrinsic noise in the data and/or non-vanishing model mis-specification.

#### Power Laws in Deep Learning

**DeepLearning Scaling [KMH<sup>+</sup>20] – Log-log Plots** Test loss of a Transformer trained to autoregressively model language



### Ubiquity/Universality of Power Laws

#### Power laws have been observed for many

- problem types (supervised, unsupervised, transfer learning)
- data types (images, video, text, even math)
- many NN architectures (Transformers, ConvNets, ...)
- *different loss functions* (cross-entropy, log, logistic, 0-1)

#### [HNA<sup>+</sup>17, RRBS19, HGLS20, HKK<sup>+</sup>20, KMH<sup>+</sup>20]

- This has *led some to the belief that power laws might be universal*: Whatever the problem, data, model, learning algorithm, or loss, learning curves follow power laws.
- To which extent this conjecture is true, we do not know, since *theoretical understanding* of this phenomenon is *limited*.

### This Talk

- Scaling with data size n.
- *Problem*: Classical learning theory leads to scaling laws  $n^{-\beta}$  with  $\beta = \frac{1}{2}$  or  $\beta = 1$ , not the observed  $\beta \approx 0.05...0.35 < \frac{1}{2}$ .
- Conjecture: Any theoretical explanation of  $\beta < \frac{1}{2}$  requires real-world data and models of unbounded complexity.

#### Possible suitable model choices:

- (a) *scaling up the model* (e.g. NN) with data, as done in the experiments [intertwines scaling with data and scaling with model size]
- (b) *non-parametric models* (e.g. kNN [SB14], Kernel regression [BCP20]) [more sophisticated analysis, manifold explanation]
- (c) a model with (countably-)infinitely-many parameters (this talk)
  [Hut21] [more accurate analysis. Zipf explanation]

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### General Findings within our Toy Model

- For domains of unbounded complexity, a *variety of learning curves* are possible, not only power-laws.
- Real *data* is often *Zipf distributed* (e.g. the frequency of words in text), which is itself a power law. This *implies power law learning* curves with "interesting"  $\beta < \frac{1}{2}$ ,
- Though many (even *non-Zipf*) distributions *also* lead to *power laws* but with "uninteresting"  $\beta = 1$ .

It is plausible that these findings remain true for most infinite models.

### Key Findings within our Toy Model

#### In general, learning curves consist of 3 terms

- 1. a *data-independent loss-dependent* power law (usually  $n^{-1/2}$  or  $n^{-1}$ ),
- 2. a *data-dependent loss-independent* power law  $n^{-\beta}$  for  $0 < \beta \le 1$ , with (typically small)  $\beta = \frac{\alpha}{1+\alpha}$  for  $(\alpha + 1)$ -Zip-distributed data,
- 3. an *irreducible term* due to noise and/or model approximation error.

#### The signal-to-noise ratio

- rapidly *deteriorates* with *n* in instantaneous learning curves.
- rapidly *improves* with *n* in time-averaged learning curves.
- Consistent with arguments by [Hut06] for log-loss, Model selection should be based on cumulative (AUC) or time-averaged error, rather than final test error.

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### Scaling with Model Size

- Consider a function  $f : [0; 1]^d \to \mathbb{R}$  which we wish to approximate.
- A naive approximation is to discretize the hyper-cube to an  $\varepsilon$ -grid. This constitutes a model with  $m = (1/\varepsilon)^d$  parameters.
- If f is 1-Lipschitz, it can approximate f to accuracy  $\varepsilon = m^{-1/d}$ , i.e. the (absolute) error scales with model size m as a power law with exponent -1/d.
- More generally, if first k derivatives of f are bounded, m parameters suffice and are necessary for Θ(m<sup>-k/d</sup>) approximation accuracy [Mha96, DHM89]
- Adapted to NNs by [Pin99] and empirically verified and extended by [SK20] to using the dimension of the data distribution in the penultimate layer of the NN.

### $Data \ Size {\leftrightarrow} Iterations {\leftrightarrow} Compute$

- (i) Usually in deep learning, *compute is proportional to the number of learning iterations*, since/provided batch and model size are kept fixed.
- (ii) in *online learning*, every data item is used only once, hence the size of data used up to iteration *n* is proportional to *n*.
- (iii) This is also true for *stochastic learning* algorithms for some recent networks, such as GPT-3, trained on massive data sets, where every data item is used at most once (with high probability).
- (iv) When generating *artificial data*, it is natural to generate a new data item for each iteration.

Hence in these 4 settings, the *learning curves*, *error-with-data-size*, *error-with-iterations*, *and error-with-compute*, *are scaled versions of each other*. For this reason, *scaling of error with iterations also tells us how error scales with data size and even with compute*, but scaling with model size is different.

### Scaling with Data Size

- This is the traditional domain of *Statistical Learning Theory (SLT)* [SB14], *online learning* [GPS18], and *online convex optimization* [Haz16].
- The *fundamental (PAC) theorem of SLT* states that the empirical error converges to the generalization error at a rate of  $n^{-1/2}$  for models of finite VC-dimension, and n i.i.d. samples.
- Applies to many models (SVMs, regression, NNs, finite decision trees, ...), many algorithms (Empirical Risk Minimization (ERM), (stochastic) gradient descent approximations, ...) many losses (convex-Lipschitz-bounded, convex-smooth-bounded, ...).

### Scaling with Data Size (ctd)

- n<sup>-1/2</sup> scaling also trivially follows from the central limit theorem for virtually any finitely-parameterized model in the under-parameterized regime of more-data-than-parameters:
  Parameters can be estimated to accuracy n<sup>-1/2</sup> hence absolute (locally quadratic loss) decays with n<sup>-1/2</sup> (n<sup>-1</sup>).
- We could easily create power laws with any  $\beta$  by choosing *exotic loss*  $|\hat{y} y_t|^{\beta/2}$ , but this would *not explain* the observed  $\beta$  for the used standard losses.
- The average *regret* considered *in online learning* theory and online convex optimization has similar requirements on the model (e.g. finite-dimensional) and *exhibits the same rates*  $n^{-1/2}$  or  $n^{-1}$  (or  $\frac{1}{n} \ln n$  due to the time-average), *under similar conditions*.

## Interesting Scaling with Data Beyond $n^{-1/2}$

- An example of a non-parametric model whose sample complexity has been analysed with "interesting" rate, is k-nearest neighbors (kNN).
- For *d*-dimensional Lipschitz functions, the *error of kNN* is bounded by  $n^{-1/(d+1)}$  [SB14, Thm.19.3&19.5].
- Power  $-1/(d+1) \approx -1/d$  is *due to density of data points* being  $n^{-1/d}$  similar to discretization discussed before in terms of model size.
- Learning curves  $n^{-\alpha/d}$  for kernel regression [BCP20, SGW20, BDK<sup>+</sup>21].
- Also hold for *infinitely wide NNs*, since equivalent to kernel regression with a Neural Tangent Kernel (NTK)
- $\alpha$  depends on target smoothness and choice of loss function.
- The underlying mechanism of  $\varepsilon$ -covering a *d*-dimensional *data* manifold with  $n \stackrel{\times}{\approx} (1/\varepsilon)^{d/\alpha}$  data points is the same.
- The origin of the power law in our toy model is very different.

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### The Goal of this Work

Identify and study the simplest model that is able to exhibit power-law learning curves as empirically observed in Deep Learning.

- Toy model: i.i.d. classification problems with countable feature space.
- A natural practical *example* application would be *classifying words* w.r.t. some criterion.
- Slides: deterministic labels and 0-1 loss
- *Toy algorithm* predicts/recalls the *class* for a new *feature* from a previously observed (*feature*, *class*) pair, or acts randomly on a novel *feature*.
- Paper: Extension to noisy labels and general loss.

### The Toy Model

- Classification:  $h \in \mathcal{H} := \mathcal{X} \to \mathcal{Y}$ , e.g.  $\mathcal{Y} = \{0, 1\}$  for binary.
- Classifier h learnt from data  $\mathcal{D}_n := \{(i_1, y_1), ..., (i_n, y_n)\} \in (\mathcal{X} \times \mathcal{Y})^n$ .
- We need *infinite*  $\mathcal{X}$  for interesting learning curves.
- Smallest suitable  $\mathcal{X} \simeq \mathbb{N}$ , which we henceforth assume.
- Model class H := N → Y is uncountable and has ∞ VC-dim., hence is not PAC learnable, but still can be learnt consistently.
- *Features*  $i_t \in \mathbb{N}$  are drawn i.i.d. with  $\mathbb{IP}[i_t = i] =: \theta_i \ge 0$  $(\sum_{i=1}^{\infty} \theta_i = 1).$
- $\infty$  vector  $\boldsymbol{\theta} \equiv (\theta_1, \theta_2, ...)$  characterizes the *feature distribution*.
- Noise-free: Label y<sub>t</sub> = h<sub>0</sub>(i<sub>t</sub>), where h<sub>0</sub> ∈ H is unknown true deterministic labelling function.
- Results change little for noisy labels.

### The Toy Algorithm

#### *Toy Algorithm A* : $\mathbb{N} \times (\mathbb{N} \times \mathcal{Y})^* \to \mathcal{Y}$

- *memorizes* all past labelled features  $\mathcal{D}_n$ .
- on next feature  $i_{n+1} = i$  recalls  $y_t$  if  $i_t = i$  for some  $i \le n$ ,
- or outputs undefined if  $i \notin i_{1:n}$  i.e. if i is new.

Formally:

$$A(i, \mathcal{D}_n) := \begin{cases} y_t & \text{if } i = i_t \text{ for some } t \leq n \\ \bot & \text{else } \text{i.e. if } i \notin i_{1:n} \end{cases}$$

#### Error

- Algorithm A only makes an *error* predicting label  $y_{n+1}$  if  $i_{1:n} \notin i_{1:n}$ .
- Formally, the *(instantaneous) error*  $E_n$  of A when predicting label  $y_{n+1}$  for feature  $i_{n+1}$  from  $\mathcal{D}_n$  is  $E_n := [[i_{n+1} \notin i_{1:n}]]$ .
- Expected (instantaneous) error (w.r.t.  $\mathcal{D}_n$  and  $i_{n+1}$ ):  $\mathbb{E}_n := \mathbb{E}[\mathsf{E}_n] = \mathbb{P}[i_{n+1} \notin i_{1:n}] = \sum_{i=1}^{\infty} \theta_i (1-\theta_i)^n$
- *Intuition:* If feature *i* has not been observed so far (happens with prob.  $(1 \theta_i)^n$ ), then feature *i* is observed (happens with prob.  $\theta_i$ ), the algorithm makes an error.
- $\mathbb{E}_n$  as a function of *n* constitutes an *(expected) learning curve*.
- Cf. probability of discovering a new species from data [Cha81], but usage&analyses of model & resulting expressions are totally different.
- Results change little for most other loss functions.

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### **Exponential Decay**

- Very simple case: *m* of the  $\theta_i$  are equal, the rest are 0.
- Error  $\mathbb{E}_n = (1 \frac{1}{m})^n \le e^{-n/m}$  decays exponentially with *n*.

This case is *not too interesting* to us, since

- (a) this case corresponds to a *finite model*
- (b) exponential decay is an "*artifact*" of the deterministic label and discontinuous 0-1 error.
- (c) becomes a power law 1/n after time-averaging (see later).
- (d) does not explain the Deep Learning power law learning curves.



### **Superposition of Exponentials**

- Expected Error  $\mathbb{E}_n$  is invariant under *bijective renumbering* of features  $i \in \mathbb{N}$
- Hence we can *w.l.g.* assume  $\theta_1 \ge \theta_2 \ge \theta_3 \ge \dots$
- Some  $\theta$ s may be equal. *Group equal*  $\theta$ s together into  $\overline{\overline{\theta}}_j$  with multiplicity  $m_j > 0$

• 
$$\mathbb{E}_n = \sum_{j=1}^M m_j \bar{ar{ heta}}_j e^{-n \bar{ar{ heta}}_j}$$
, where  $\bar{ar{ heta}}_j := -\ln(1 - ar{ar{ heta}}_j) pprox ar{ar{ heta}}_j$ 

•  $M \in \mathbb{N} \cup \{\infty\}$  is the number of *different*  $\theta_i > 0$ .

### **Superposition of Exponentials**

- $\mathbb{E}_n = \sum_{j=1}^M m_j \bar{\theta}_j e^{-n\bar{\vartheta}_j}$ is a superposition of exponentials in *n* with different decay rates  $\bar{\vartheta}_j$
- Sum will be *dominated* by different terms at different "times" *n*.
- Different *phases* of exponential decay



- For  $M < \infty$ , eventually exponential decay  $e^{-m\bar{\vartheta}_M}$  will dominate  $\mathbb{E}_n$ .
- The same "caveats" (a)-(d) apply as for M = 1 two slides ago.

### Approximation

 Let f : ℝ → ℝ be a smooth and monotone decreasing *interpolation* of θ : ℕ → ℝ, i.e. f(i) := θ<sub>i</sub> and f'(x) < 0:</li>

$$\mathbb{E}_{n} = \sum_{i=1}^{\infty} f(i)(1-f(i))^{n} \approx \int_{1}^{\infty} f(x)e^{-nf(x)}dx$$
  
$$\stackrel{(a)}{=} \int_{0}^{\theta_{1}} \frac{ue^{-nu}du}{|f'(f^{-1}(u))|} \approx \frac{1}{n^{2}|f'(f^{-1}(\frac{1}{n}))|} = \frac{d}{dn}f^{-1}(\frac{1}{n})$$

(a) Reparametrization u = f(x) and  $f(1) = \theta_1$  and  $f(\infty) = 0$  and dx = du/f'(x) and f' < 0.

- (×) Numerator  $ue^{-nu}$  concentrated around u = 1/n, hence can replace u by 1/n in denominator.
  - Intuition:  $\mathbb{E}_n$  is dominated by samples  $i_0$  for which  $\theta_{i_0} \approx \frac{1}{n}$ .

• Accuracy of the integral representation is 1/en + o(1/n).

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### Zipf-distributed data

- Empirically many *data* follow a power-law distribution called *Zipf distr.* in this context:
- The *frequency* of the *i*th most frequent item is approximately  $\theta_i \propto i^{-(\alpha+1)}$  for some  $\alpha > 0$ .

• 
$$\mathbb{E}_n \stackrel{\times}{=} n^{-\beta}$$
 where  $\beta := \frac{\alpha}{1+\alpha}$ 



• That is, Zipf-distributed data (with power  $\alpha + 1$ ) lead to a *power-law learning curve* (with power  $\beta = \frac{\alpha}{1+\alpha} < 1$ ).

### (Super)Exponentially-Distributed Data

- *Exponential data distr.*  $\theta_i \propto e^{-\gamma i}$  is more skewed than any power law.
- Still  $\mathbb{E}_n \approx 1/\gamma n$ , i.e. still leads to a power law learning curve.
- But exponent  $\beta = 1$  is "uninteresting" (much larger than observed)
- Surprise: Any super-exponential data (e.g. θ<sub>i</sub> ∝ e<sup>-γi<sup>2</sup></sup>, but quite unrealistic) always leads to a (sort of) power law as long as θ<sub>i</sub> > 0 for infinitely many i, unlike finite model which gives exponential decay:
- $\mathbb{E}_n \stackrel{\times}{\leq} n^{-1}$  for all *n* and  $\mathbb{E}_n \stackrel{\times}{\geq} n^{-1}$  for infinitely many *n*.

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#### Instantaneous Variance

- Variance  $\mathbb{V}_n$  of  $\mathsf{E}_n := \llbracket i_{n+1} \notin i_{1:n} \rrbracket$  as a function of n is important.
- Useful learning curve requires *Standard Error (STE)*  $\sqrt{\mathbb{V}_n/k} < \mathbb{E}[\mathbb{E}_n] \equiv \mathbb{E}_n =: \mu_n$  when averaging over k runs.
- $\mathsf{E}_n \in \{0,1\}$  hence  $\mathsf{E}_n^2 = \mathsf{E}_n$  hence  $\mathbb{V}[\mathsf{E}_n] = \operatorname{I\!E}[\mathsf{E}_n^2] \operatorname{I\!E}[\mathsf{E}_n]^2 = \mu_n(1 \mu_n)$

• Since  $\mu_n \to 0$  for  $n \to \infty$ , the Standard Deviation (STD)  $\sigma_n := \sqrt{\mathbb{V}[\mathbb{E}_n]} = \sqrt{\mu_n(1 - \mu_n)}$  $\approx \sqrt{\mu_n} \gg \mu_n = \mathbb{E}_n$ 

 For good signal-to-noise ratio we need k ≫ μ<sub>n</sub><sup>-1/2</sup> runs (increasing with n!)



### **Time-Averaged Mean and Variance**

- Alternative: Report the time-averaged error  $\overline{\mathsf{E}} := \frac{1}{N} \sum_{n=0}^{N-1} \mathsf{E}_n$ , rather than the instantaneous error  $\mathsf{E}_n$ .
- Expectation:  $\mathbb{E}[\overline{\mathsf{E}}_N] = \frac{1}{N} \sum_{i=1}^{\infty} [1 (1 \theta_i)^N]$

• Variance: 
$$\mathbb{V}[\overline{\mathsf{E}}_{N}] = \frac{1}{N^{2}} \sum_{i=1}^{\infty} (1-\theta_{i})^{N} [1-(1-\theta_{i})^{N}] - \frac{1}{N^{2}} \sum_{i\neq j} [(1-\theta_{i})^{N} (1-\theta_{j})^{N} - (1-\theta_{i}-\theta_{j})^{N}]$$

## Uniform Case $\theta_i = \frac{1}{m} \llbracket i \leq m \rrbracket$



 I.e. Standard Deviation is (much) smaller than the mean for N ≫ m, so the time-averaged learning curves have a much better signal-to-noise ratio.

## Zipf Case $\theta_i \propto i^{-(\alpha+1)}$

- Recall expected error:  $\mathbb{E}_n \approx c_\alpha n^{-\beta}$ , where  $0 < \beta = \frac{\alpha}{1+\alpha} < 1$ .
- Time-averaged expected error:  $\mathbb{E}[\overline{\mathsf{E}}_N] \approx \frac{c_{\alpha}}{N} \int_0^N n^{-\beta} dn = \frac{c_{\alpha}}{1-\beta} N^{-\beta}$
- Same power law with the same exponent  $\beta$  (generic property)
- STD  $\sigma[\overline{\mathsf{E}}_N] \stackrel{\times}{\approx} N^{-\frac{1/2+\alpha}{1+\alpha}} \ll N^{-\frac{\alpha}{1+\alpha}} \stackrel{\times}{\approx} \operatorname{I\!E}[\overline{\mathsf{E}}_N]$
- Signal-to-noise ratio is  $\sigma[\overline{E}_N]/\mathbb{E}[\overline{E}_N] \stackrel{\times}{\approx} N^{-1/(2+2\alpha)}$ . STD much smaller than Mean.
- Single run suffices to get a good (and excellent for n ≥ 500) signal-to-noise ratio for ave. and cum. error



#### General $\theta$ Case

- Signal-to-noise ratio:  $\frac{\sigma[\overline{E}_N]}{\operatorname{I\!E}[\overline{E}_N]} \leq \frac{\sqrt{\frac{1}{N} \operatorname{I\!E}_N}}{\operatorname{I\!E}[\overline{E}_N]} = \frac{\sqrt{N \operatorname{I\!E}_N}}{\sum_{n=0}^{N-1} \operatorname{I\!E}_n} \xrightarrow{N \to \infty} 0$
- *Proof* requires to distinguish two cases:
- 1)  $\sum_{n=0}^{\infty} \mathbb{E}_n \leq c$  (e.g. exponential error decay in finite models),
- 2)  $\sum_{n=0}^{N-1} \mathbb{E}_n \to \infty$  (most  $\infty$  models, e.g. Zipf, even exponential  $\theta_i$ )

### Instantaneous vs. Time-Averaged Error

- Trivial observation: For θ<sub>0</sub> = 1, we have i<sub>n</sub> = 1 ∀n, hence E<sub>0</sub> = 1 and E<sub>n</sub> = 0 ∀n ≥ 1 and V[E<sub>n</sub>] = 0 ∀n.
- This is the fastest any error can decay, 0 after 1 observation, hence always  $\overline{E}_n = \Omega(1/n)$ . Fazit:

If 
$$\mathbb{E}_n = o(1/n)$$
, report  $\mathbb{E}_n$ , since  $\ll \overline{\mathbb{E}}_n$ .

If  $\mathbb{E}_n = \tilde{\Omega}(1/n)$ , report  $\overline{\mathbb{E}}_n$ , since  $\stackrel{\times}{\approx} \mathbb{E}_n$  but variance is smaller.

- Esp. in Deep Learning with small  $\beta$ , we have  $\overline{E}_n \approx E_n$ .
- Low variance does not follow directly from law of large numbers, since E<sub>1</sub>, E<sub>2</sub>, E<sub>3</sub>, ... are not independent.

## Zipf-Distributed Words in Typical Texts



Relative (left scale) and absolute (right *Power law* fit to learning curve for this scale) word frequency, and fitted Zipf law.

data set for a word classification task.

- The *power-law fit is good* if *n* is not too large.
- For large *n*, the error decays exponentially as  $\exp(-\theta_{min}n)$ , since word frequency is quantized  $(\in \mathbb{N})$ .

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### Noisy Labels or Targets – Implications

- (a) Need "smarter" "learning" algorithm, e.g. predicting the average.
- (b) Subtract irreducible error due to label noise before studying scaling.
- (c) Extra  $n^{-1/2}$   $(n^{-1})$  additive error term for absolute (square loss) due to parameter estimation error, hence
- (d) Inst. loss will not decay expon. anymore even if model is finite.
- (e) Otherwise the scaling laws for Zipf data are unchanged.

In summary, conceptually error/loss is a sum of 3 terms:

- (1) The parameter learning rate  $n^{-1/2}$  (squared for locally quadratic loss)
- (2) the same power law  $n^{-\beta}$  as in the deterministic case,
- (3) the inherent "entropy" in the data.
  - *Remarkably:* Instantaneous square  $\text{Loss}_n^{noisy}(A) \stackrel{\times}{=} \mathbb{E}_n^{det.} + \mathbb{E}[\overline{\mathbb{E}}_n^{det.}].$
  - This "magically" ensures (c,d,e), since  $\mathbb{E}[\overline{\mathsf{E}}_n] \stackrel{\times}{\approx} \max\{\mathbb{E}_n, \frac{1}{n}\}$ .
  - For instance, for a finite model,  $\operatorname{Loss}_n(A) \stackrel{\times}{\approx} \operatorname{I\!\!E}[\overline{\mathsf{E}}_n] \stackrel{\times}{\approx} \frac{1}{n}$ .

### **Other Loss Functions**

- Deterministic toy model:  $\mathbb{E}[\text{Loss}_n] \stackrel{\times}{=} \mathbb{E}_n$  for most loss functions
- Noisy labels: Same, but extra n<sup>-1</sup> or n<sup>-1/2</sup> term (now fastest possible decay)
- Universality at least within toy model: For large models, scaling laws are indep. of loss function and not affected by noise.

#### **Continuous Features**

- Feature spaces are most often vector spaces  $\mathbb{R}^d$ .
- No feature ever repeats exactly  $(x_n \neq x_m \text{ for } n \neq m)$ .
- *Simple processes*: Dirichlet = Chinese Restaurant = Stick-Breaking.
- Leads to power law learning curves  $n^{-1}$ , but  $\beta = 1$  is uninteresting.
- Generalized 2-parameter *Poison Dirichlet Process* [BH10] also only leads to β = 1.
- *Open problem:* Finding analytically tractable models with continuous features that exhibit interesting learning curves.

### **Generalizing Algorithms**

- Proper models/algorithms for continuous features *need to generalize* from observed inputs to similar future not-yet-observed inputs.
- Simple model: Partition domain into countably many cells
- If done a-priori and independent  $\mathcal{D}_n$  reduces back to toy model
- More realistically, if partitioning, e.g. clustering of data, is data (size) dependent, it will affect the scaling.
- 'perfect prediction for exact repetition' *abstracts* 'classify features in the same cell alike' *abstracts* 'classify similar observations alike or similarly'.
- So maybe some of our findings or analysis tools approximately *transfer*.

### Deep learning

- (Deep) neural networks are a particularly powerful class of models/algorithms that *can generalize*,
- But they are notoriously *difficult to theoretically analyse*.
- It may be a long way from our toy model to a similar analysis of NNs.
- Furthermore we have not at all considered the equally interesting questions of *scaling with model size*.

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- We introduced and analyzed the simplest model that can exhibit *power laws (decrease of error with data size)* consistent with recent findings in deep learning.
- Many but not all *data distributions* lead to power laws.
- Zipf data with exponent  $\alpha + 1$  lead to power law with exponent  $\beta = \alpha/(1 + \alpha)$ . Artifact of the model or wider validity?
- The *signal-to-noise ratio* for the time-averaged error tends to zero, which implies that a single experimental run suffices for stable results.
- *Model selection* should be based on cumulative (AUC) error, rather than final test error [Hut06].

### Limitations

- The toy model is totally unrealistic as a Deep Learning model,
- but we believe it captures the (or at least a) true reason for the observed scaling laws w.r.t. data.
- Hopefully can be generalized to NNs
- We have not addressed *scaling laws w.r.t. model size*.

### **Applications**

- May help making better or more principled choices for network architecture (depth, with, and beyond), hyper-parameters, fine-tuning, data augmentation, pre-training, etc. [CJS<sup>+</sup>93, HGLS20].
- Being able to extrapolate the consequences of such choices from *cheap training on a small subset of the data* to the whole corpus by simply fitting power laws can save significant compute.
- The *cost of training recent models has reached millions of dollars* and can exhaust and exceed even FAANGs computational resources.

### **List of Notation**

#### Symbol Explanation

- [Bool] 1 if Bool=True, 0 if Bool=False
- $\mathbb{E}, \mathbb{V}$  Expectation, Variance
- $\stackrel{\times}{=}$  Equal within a multiplicative constant
- $\theta_i$  probability of feature *i*
- $\mathcal{D}_n$  Data consisting of *n* (feature *i*, label *y*) pairs
- $E_n$  Instantaneous Error of A on  $i_{n+1}$  predicting  $y_{n+1}$  from  $\mathcal{D}_n$
- **E**<sub>n</sub> Expectation of Instantaneous Error  $E_n$  w.r.t.  $\mathcal{D}_{n+1}$
- $\overline{\mathsf{E}}_N$  Time-Averaged Error  $\mathsf{E}_n$  from n = 0, ..., N 1
- $\alpha + 1$  Exponent of Zipf distributed data frequency  $i^{-(\alpha+1)}$
- β Exponent of power law  $n^{-β}$  for error as a function of data size nγ Decay rate for exponential data distribution  $e^{-γi}$

#### References

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