

Mathematics of Deep Learning

Stability and robustness

Lessons: Kevin Scaman
TDs: Mathieu Even



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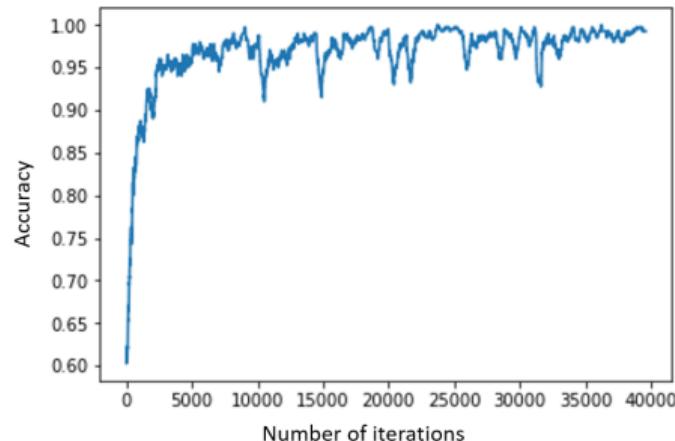
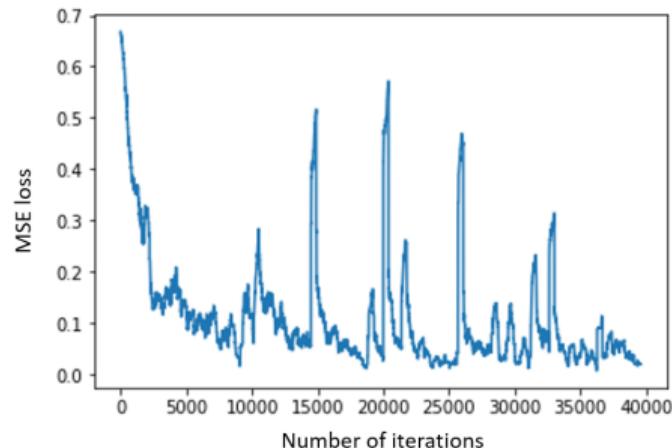
Stability during training

Weights initialization, gradient vanishing and explosion

Stability during training

Example with simple RNNs (Elman networks, no gating mechanisms)

- ▶ The gradients are sometimes **very large**.
- ▶ This leads to a large **drop in accuracy**.
- ▶ Results are **quite random**, final performance depends on initialization.



Gradient vanishing and explosion

Breaking gradient descent

- ▶ If θ_t are the iterates of the parameters learned using stochastic gradient descent on minibatches $(x_{t,i}, y_{t,i})_{i \in [1, K]}$ at time t , then we have

$$\theta_{t+1} = \theta_t - \frac{\eta}{K} \sum_i \nabla \mathcal{L}_{x_{t,i}, y_{t,i}}(\theta),$$

where $\mathcal{L}_{x,y}(\theta) = \ell(g_\theta(x), y)$.

- ▶ **Gradient vanishing:** When the gradients $\nabla \mathcal{L}_{x_{t,i}, y_{t,i}}(\theta)$ are very small compared to θ_t , the iteration does not modify the parameters.
- ▶ **Gradient explosion:** When the gradients $\nabla \mathcal{L}_{x_{t,i}, y_{t,i}}(\theta)$ are very large compared to θ_t , the iteration will push the parameters to extreme values.

Gradient vanishing and explosion

Why is it a problem for deep learning?

- ▶ By chain rule, the gradient tends to multiply along the layers.
- ▶ Example: If $g^{(L)}(x) = f^{(L)} \circ f^{(L-1)} \circ \dots \circ f^{(1)}(x)$ where $f^{(L)} : \mathbb{R} \rightarrow \mathbb{R}$, then

$$g^{(L)'}(x) = \prod_{l=1}^L f^{(l)'}(g^{(l-1)}(x))$$

- ▶ If $f^{(l)'}(g^{(l-1)}(x)) \approx c$, then $g^{(L)'}(x) \approx c^L$.
- ▶ **Exponentially small** w.r.t. L if $c < 1$ (gradient vanishing).
- ▶ **Exponentially large** w.r.t. L if $c > 1$ (gradient explosion).

Mitigation techniques: how to avoid this?

Gradient clipping

- ▶ `torch.nn.utils.clip_grad_norm_(model.parameters(), threshold)`
- ▶ **Pros:** Easiest method, just limits the gradient norm to a fixed value.
- ▶ **Cons:** Only for gradient explosion, adds an extra hyper-parameter.

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Architecture changes

- ▶ Gates in RNNs, residuals in CNNs, dropout, batch normalization, ...
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- ▶ **Cons:** Requires to change the network architecture, application dependent.

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Weight initialization

- ▶ Automatically implemented, but can have an **large impact** on performance

Weights initialization

Ideal initialization scheme

- ▶ The better the model is at initialization, the more changes we have of find good weights.
- ▶ We would like to have values that are reasonable, $\forall i \in [\![1, d^{(L)}]\!], |g_{\theta}(x)_i| \approx 1$.
- ▶ We would like to have gradients that are neither too large nor too small

$$\forall i \in [\![1, p]\!], \quad |\nabla \mathcal{L}_{x,y}(\theta)_i| \approx 1$$

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Simple solution

- ▶ Set $b^{(l)} = 0$ and sample the weights $W_{ij}^{(l)} \sim \mathcal{P}$ i.i.d. with expectation 0 and variance $V^{(l)}$.
- ▶ Choose $V^{(l)}$ so that the variance is constant across layers.

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- ▶ Technical assumptions:
 - ▶ The probability distribution is symmetric w.r.t. 0 and $\mathcal{P}(\{0\}) = 0$.
 - ▶ The activation function is ReLU $\sigma(x) = \max\{0, x\}$.

Derivation of optimal weights

Preliminary results

- ▶ Let $x \in \mathbb{R}^{d^{(0)}}$ a fixed input and, $\forall l \in \llbracket 1, L \rrbracket$, $X^{(l)} = g_\theta^{(2l-1)}(x)$.
- ▶ For any $l \in \llbracket 1, L \rrbracket$, the variables $(X_i^{(l)})_{i \in \llbracket 1, d^{(2l-1)} \rrbracket}$ are identically distributed.
- ▶ The distribution of $X_i^{(l)}$ is symmetric w.r.t. 0 (and thus $\mathbb{E}(X_j^{(l)}) = 0$).

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- ▶ If the properties are verified for l , then $X_i^{(l)} = \sum_j W_{ij}^{(l)} \sigma(X_j^{(l-1)})$, which is identically distributed and symmetric.



Derivation of optimal weight variance

Variance of the intermediate outputs

- ▶ For any $l \in \llbracket 2, L \rrbracket$ and $i \in \llbracket 1, d^{(l)} \rrbracket$, we have

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 \end{aligned}$$

- Hence, the **variance is constant across layers** if $V^{(l)} = 2/d^{(l-1)}$, and

$$\text{var}(g_\theta(x)_i) = 2\|x\|_2^2/d^{(0)}$$

Kaiming initialization (Kaiming He et.al., 2015)

Gaussian weights

Our assumptions are satisfied if we use Gaussian weights $W_{ij}^{(l)} \sim \mathcal{N}\left(0, \frac{2}{d^{(l-1)}}\right)$.

Uniform weights

If we take uniform weights $W_{ij}^{(l)} \sim \mathcal{U}([-r^{(l)}, r^{(l)}])$, then $V^{(l)} = r^{(l)}/3$ and

$$r^{(l)} = \sqrt{\frac{6}{d^{(l-1)}}}$$

Variance of the gradient

Variance propagation during backprop

- ▶ Same analysis for backprop, but in **reverse**.
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Xavier initialization (Xavier Glorot & Yoshua Bengio, 2010)

Let $c > 0$ be a hyper-parameter. The weights are initialized using the heuristic

$$W_{ij}^{(l)} \sim \mathcal{U}([-r^{(l)}, r^{(l)}]) \quad \text{and} \quad r^{(l)} = \sqrt{\frac{6c^2}{d^{(l)} + d^{(l-1)}}}$$

Batch normalization

Idea

- ▶ Normalize the input of each layer by **removing mean and dividing by std.**
- ▶ Also uses a **learnable affine map**.

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Definition

- ▶ If $(x_i)_i$ is a batch of b inputs (to the layer), then the output is:

$$y_i = \frac{x_i - E}{\sqrt{V + \varepsilon}} \cdot \gamma + \beta$$

where $E = \frac{1}{b} \sum_i x_i$ and $V = \frac{1}{b} \sum_i (x_i - E)^2$ (coord.-wise), γ and β are learnable vectors.

Batch normalization



The output depends on the whole batch, not just single inputs!

Train and eval

- ▶ The behavior of batch norm is different between training and evaluation (e.g. `model.train()` and `model.eval()` in Pytorch).
- ▶ At evaluation, the model uses a (moving) average of **all training batches**.
- ▶ Stores E and V for each training batch, and then computes

$$(1 - \rho) \sum_t \rho^t E_t \quad \text{and} \quad (1 - \rho) \sum_t \rho^t V_t$$

where (typically) $\rho = 0.9$.

Recap

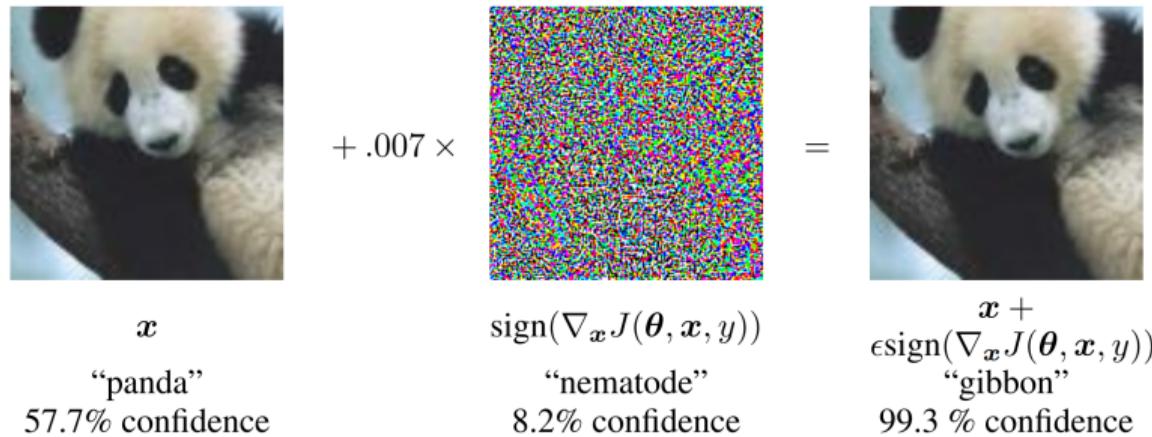
- ▶ Gradient **vanishing** and **explosion** can happen during training of **deep** NNs.
- ▶ **Gradient clipping, batch normalization, regularisation** and proper **weight initialization** can help stabilize training.
- ▶ The variance of the weights at initialization should be **inversely proportional to the layer width**.

Robustness and adversarial attacks

Confusing a neural network with noise

Adversarial attacks

- ▶ Can a small (invisible) noise change the prediction of a vision model?
- ▶ Vision models are robust to random input noise.
- ▶ Vision models are **extremely fragile** to **well-crafted** input noise.



source: Explaining and Harnessing Adversarial Examples, Goodfellow et al, ICLR 2015.

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source: Robust Physical-World Attacks on Deep Learning Visual Classification, Eykholt et al, CVPR 2018.

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source: Accessorize to a crime: Real and stealthy attacks on state-of-the-art face recognition, Sharif et.al., CCS 2016.

Adversarial attacks: examples

Fast gradient sign method (Goodfellow et.al., 2014)

- ▶ **Idea:** Take one gradient step in the direction that **maximizes the loss**.
- ▶ To control the maximum pixel noise, use the coordinates' sign instead of value.
- ▶ **Limitations:** Destroys performance, but cannot target a specific class.

$$x^{\text{att}} = x^{\text{true}} + \varepsilon \operatorname{sign}(\nabla_x \mathcal{L}(\theta, x^{\text{true}}, y^{\text{true}}))$$

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Iterative Target Class Method (Kurakin et.al., 2016)

- ▶ **Idea:** Perform gradient descent on the loss with **labels swaped**.
- ▶ To control the maximum pixel noise, project on a ball of radius ε around x .
- ▶ **Limitations:** Requires to know the model weights (white box setting).

$$x_{k+1}^{\text{att}} = \operatorname{Clamp}_{x^{\text{true}}, \varepsilon} (x_k^{\text{att}} + \varepsilon \operatorname{sign}(\nabla_x \mathcal{L}(\theta, x_k^{\text{att}}, y^{\text{att}})))$$

Beyond the white box setting

White-box attacks

- ▶ Use the knowledge of the model to create the perturbation.
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Defenses

- ▶ Augment the dataset with adversarial attacks (brute-force).
- ▶ Control the smoothness of the model (see next).

Robustness of neural networks

What makes a model robust?

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- ▶ For piece-wise linear interpolation, Lipschitz constant is **smaller than target function**.
- ▶ For neural networks: $L_{g_\theta} \leq \prod_l L_{f^{(l)}} \dots$ can be exponential in number of layers!

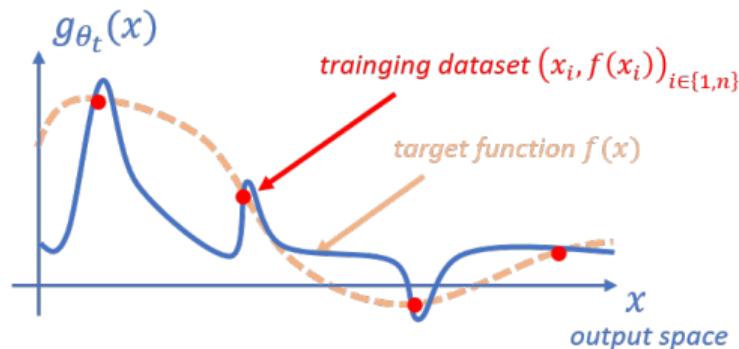
Generalization beyond the training samples

From train accuracy to test accuracy

Beyond the training samples



Good generalization



Poor generalization

- ▶ **Left model:** More regular, worst on the training set, better on the whole space.
- ▶ **Right model:** Less regular, better on the training set, worst on the whole space.
- ▶ How does the model behaves when the **test samples are different from the training samples?**

Beyond the training samples

Training objective and risk minimization

- Let $g_\theta : \mathcal{X} \rightarrow \mathcal{Y}$ be a model and \mathcal{D} be a distribution of data points in $\mathcal{X} \times \mathcal{Y}$.

$$\min_{\theta \in \mathbb{R}^d} \mathcal{L}_{\mathcal{D}}(\theta) \triangleq \mathbb{E}_{(X,Y) \sim \mathcal{D}}(\ell(g_\theta(X), Y))$$

- During training we minimize $\mathcal{L}_{\hat{\mathcal{D}}_n}(\theta)$ where $\hat{\mathcal{D}}_n = \frac{1}{n} \sum_i \delta_{(x_i, y_i)}$ is the empirical distribution over the training dataset $(x_i, y_i)_{i \in [\![1,n]\!]}$.

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Statistical error

- If $\theta \in \mathbb{R}^d$ is independent of the training samples, then, with probability $1 - \delta$,

$$\left| \mathcal{L}_{\hat{\mathcal{D}}_n}(\theta) - \mathcal{L}_{\mathcal{D}}(\theta) \right| \leq \|\ell\|_\infty \sqrt{\frac{2 \ln(2/\delta)}{n}}$$

- Unfortunately, the SGD iterates $\hat{\theta}_{n,t}$ depend on the training dataset $\hat{\mathcal{D}}_n$...

Decomposition of the error

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- Let $\hat{\theta}_{n,t}$ be the parameters after t training steps and $\theta^* \in \operatorname{argmin}_\theta \mathcal{L}_{\mathcal{D}}(\theta)$. Then,

$$\mathcal{L}_{\mathcal{D}}(\hat{\theta}_{n,t}) = \mathcal{L}_{\mathcal{D}}(\hat{\theta}_{n,t}) - \mathcal{L}_{\hat{\mathcal{D}}_n}(\hat{\theta}_{n,t}) + \mathcal{L}_{\hat{\mathcal{D}}_n}(\hat{\theta}_{n,t}) - \mathcal{L}_{\hat{\mathcal{D}}_n}(\theta^*) + \mathcal{L}_{\hat{\mathcal{D}}_n}(\theta^*) - \mathcal{L}_{\mathcal{D}}(\theta^*) + \mathcal{L}_{\mathcal{D}}(\theta^*)$$

Generalization error **Optimization error** **Statistical error** **Approx.**

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Generalization error
 Optimization error
 Statistical error
 Approx.

- Approximation error:** by the universality of MLPs, is arbitrarily small.

$d \searrow$

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- Statistical error:** Convergence in $O\left(\frac{1}{\sqrt{n}}\right)$ by Tchebyshev concentration. $n \searrow$
- Generalization error:** Difficult part. Depends on the model and opt. $d \nearrow, t \nearrow, n \searrow$

Overfitting in ML

Usual analysis

- ▶ Optimization error decreases
- ▶ Generalization error increases
- ▶ There is a trade-off

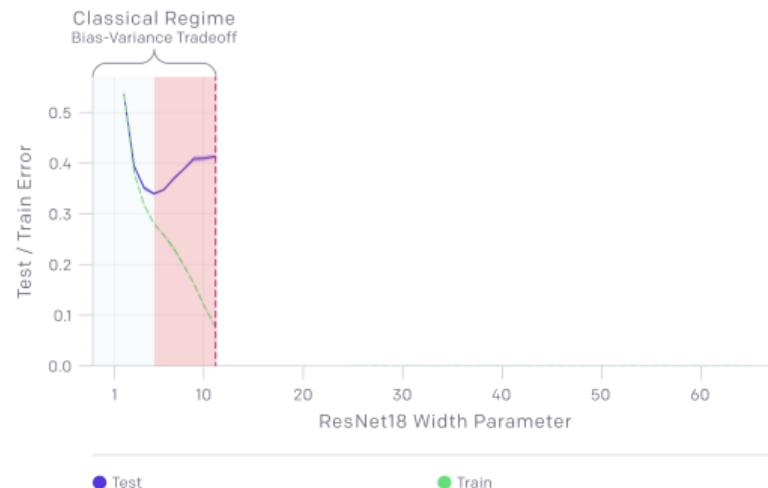
Usual mitigation strategies

- ▶ Early stopping
- ▶ Hyper-parameter selection via cross-validation
- ▶ Regularization: $\min_{\theta} \mathcal{L}_{\hat{\mathcal{D}}_n}(\theta) + g(\theta)$ (usually $g(\theta) = \gamma \|\theta\|_2^2$).

But...Double descent!

Overfitting mitigated by over-parameterization

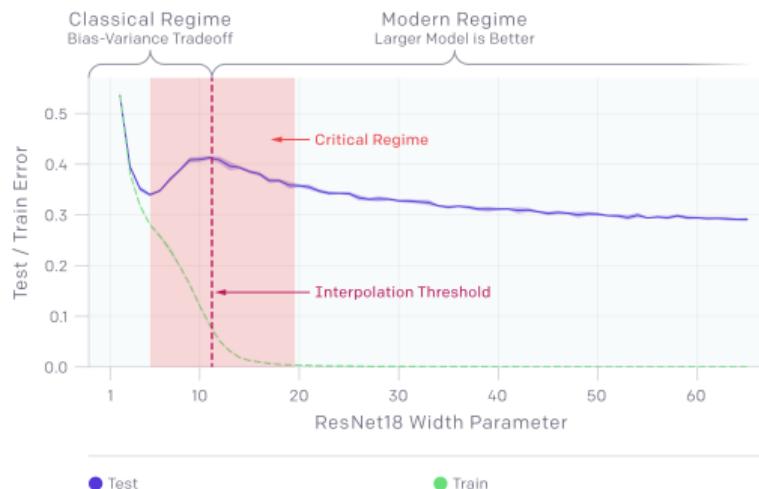
- ▶ After a certain model size, test error starts decreasing again.
- ▶ Over-parameterizing tends to create **implicit regularization**.



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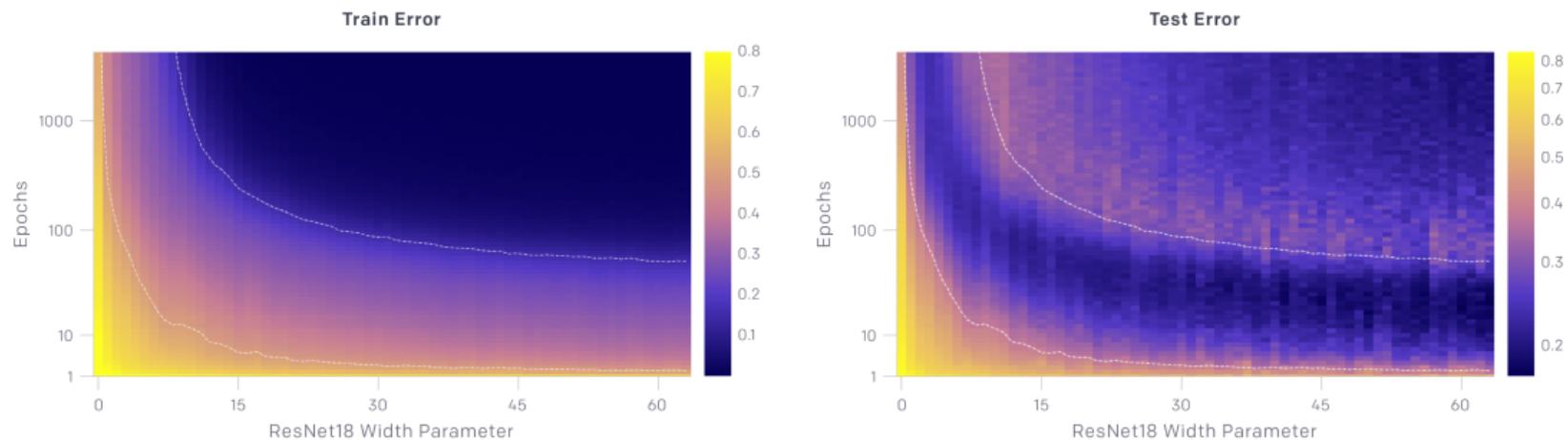
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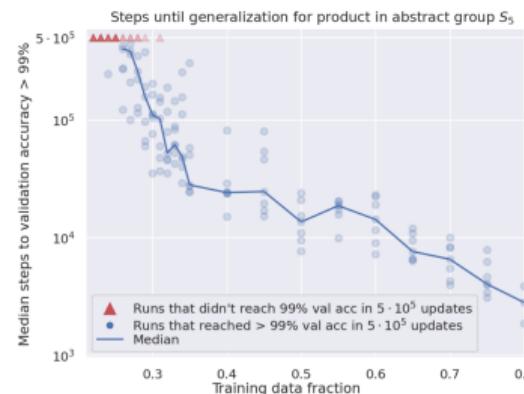
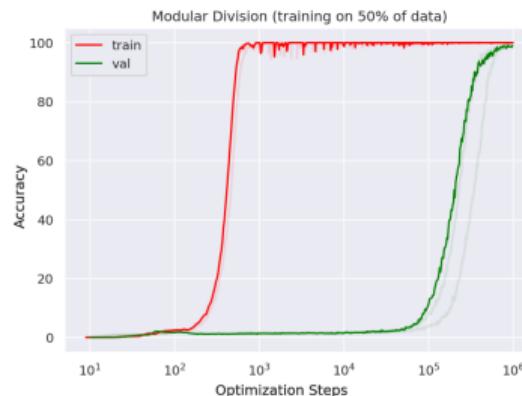


source: <https://openai.com/blog/deep-double-descent/>

But... Grokking!?

Generalization beyond overfitting

- ▶ All hope is lost... until you forgot to turn your computer off during the holidays.
- ▶ Very (very) large plateaux during training.
- ▶ Still not a satisfactory explanation (don't do this at home. ;-)).



★	a	b	c	d	e
a	a	d	?	c	d
b	c	d	d	a	c
c	?	e	d	b	d
d	a	?	?	b	c
e	b	b	c	?	a

source: Grokking: Generalization Beyond Overfitting on Small Algorithmic Datasets, Power et.al., 2022.

Recap

- ▶ **Overfitting** to the training dataset can be an issue when the number of parameters is larger than the number of samples.
- ▶ In practice, **overparameterization can help generalization** (often called implicit regularization).
- ▶ The training curves can exhibit a **double descent behavior**.
- ▶ Long plateaux can appear on the test loss/accuracy.