Lecture 20

## Statistical Learning for Neural Networks

### Guaranteed Success for ERM with Finite Hypotheses

#### \* Accuracy $\epsilon$ and Tolerance $\delta$

- $\epsilon$  is the desired maximum error rate of the hypothesis found by ERM.
- $\delta$  is the tolerance for the probability of failure, i.e., the probability that the ERM hypothesis will exceed the error rate  $\epsilon$ .

#### The Bound

With  $m > \frac{\log(|H|/\delta)}{\epsilon}$ , ERM will find a hypothesis with error less than  $\epsilon$  with probability greater than  $1 - \delta$ .

For a large enough m, ERM produces a hypothesis with error under  $\epsilon$  at probability  $1-\delta.$ 

### VC Dimension

- Definition:
  - VC Dimension is the maximum number of data points that a model class can shatter.
- Shattering:
  - A model class can shatter a set of data points if it can perfectly classify every possible arrangement of labels for that set.

### VC Dimension of a Line in 2D

• In a two-dimensional space, the VC dimension of a line is 3.



# PAC (Probably Approximately Correct) learnability

- Being able to learn a good-enough hypothesis with high probability given enough examples.
- If with enough data, a model from H can be learned that is probably correct (within  $\epsilon$  error) with high confidence (probability >  $1 \delta$ ), then H is PAC Learnable.

H is PAC Learnable if:

- there is a function  $m_H: (0,1)^2 o \mathbb{N}$
- ${}^{\bullet}$  and a learning algorithm A,
- ${}^{\bullet}$  such that for every distribution D over X,
- every  $\epsilon,\delta>0$  and every f in H,
- for samples S of size m ( $m=m_{H}(\epsilon,\delta)$ )
- ${}^{ullet}$  generated by D and labeled by f,
- $\Pr[L_D(A(S)) > \epsilon] < \delta.$

### **Relaxing the Realizability Assumption**

#### • Realizability Assumption:

- Assumes True Function in Hypothesis Class H
- Hypotheses Contain Exact Solution
- Realistic Setup:
  - Lack of A Priori Knowledge:
    - Learner Doesn't Know if True Classifier in  ${\cal H}$

#### **Unpredictability of Labels**:

- Another aspect of more realistic scenarios: labels (the outputs we're trying to predict) might not be fully determined by the instance attributes (the input data).
  - Various reasons: noise in the data, unobserved variables, or inherently stochastic processes.
  - In such cases, even the best possible model in *H* might not perfectly predict the labels for all instances.

#### **Implications for Learning**:

- When these more realistic conditions are assumed, learning becomes more challenging.
- The learner must now find the best possible hypothesis within the class, even if none of the hypotheses can perfectly predict all instances.

### **General Loss Functions**

- Beyond Classification Errors:
  - Our learning approach goes beyond counting classification errors.
- Domain Set Z:
  - Z represents the set of all possible instances or data points.
- Loss Function *l*:
  - $l:H imes Z o \mathbb{R}$
  - Quantifies model h's loss on instance z.

### **General Loss Functions**

#### • Probability Distribution P:

- P is a probability distribution over Z.
- Defines the likelihood of each instance.
- Expected Loss  $L_P(h)$ :
  - $L_P(h) = E_{z \sim P}[l(h,z)]$
  - \* Average loss of model h under P.
- Flexible Learning Formalism:
  - Allows us to assess model performance with various loss functions and distributions.

### **Agnostic PAC Learnability**

- H is Agnostic PAC Learnable if:
  - There exists a function  $m_H: (0,1)^2 o \mathbb{N}$  and a learning algorithm A.
  - For every distribution P over X imes Y and every  $\epsilon,\delta>0.$
  - For samples S of size  $m>m_H(\epsilon,\delta)$  generated by P.

$$\Pr[L_P(A(S)) \leq \inf_{h \in H} L_P(h) + \epsilon] \geq 1 - \delta$$

### **Guaranteeing Learnability**

- **Question**: Can such learnability be guaranteed?
- **Crucial Factor**: The VC (Vapnik-Chervonenkis) dimension of the class *H*.
- Fundamental Theorem of Statistical Learning:
  - "A class *H* is PAC (Probably Approximately Correct) learnable if and only if its VC dimension is finite."

### **Fundamental Theorem - Quantitative Version**

• The number of random labeled samples needed for learning a class of predictors *H* is given by:  $\int VCdim(H) + \log(\frac{1}{2})$ 

$$O\left(\frac{VCdim(H) + \log(\frac{1}{\delta})}{\epsilon^2}\right)$$

### **Complexity-Dependent Error Bound**

- ${}^{\bullet} \ L_P(h) \leq \inf_{h \in H} L_P(h) + f(c(h);m)$
- The true error of a hypothesis h is bounded by a function of its complexity and training sample size.

#### $\operatorname{Function} f(c(h);m)$

- Dependent on the complexity measure c(h) of hypothesis h and sample size m.
- Indicates a relationship between hypothesis complexity, sample size, and the achievable error.

Integrates the concept of hypothesis complexity into the error bound.

### **Expressive Power of Neural Networks**

**Theorem**: Fix some  $\epsilon$  in the range (0, 1), and let s(n) denote the minimal number of nodes.

- There exists a neural network with s(n) nodes that can approximate up to  $\epsilon$  every function from  $[0,1]^n$  to [0,1].
- s(n) is exponential in n.

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- s(n) is exponential in n.
- Key Insight:
- Neural networks possess remarkable expressive power.
- They can approximate a wide range of functions with high accuracy.
- However, achieving this expressive power may require a large number of nodes, which can grow exponentially with the input dimension *n*.

### Measuring Error Guarantees For Neural Networks

$$L_P(h) \leq ext{inf}_{h \in H} \, L_S(h) + \sqrt{rac{|E| + \log(1/\delta)}{m}}$$

Where:

- |E| is the number of edges (parameters) in the hypothesis h.
- $\bullet m$  is the sample size.

### **Rethinking Generalization in Deep Learning**

- Authors: Chiyuan Zhang, Samy Bengio, Moritz Hardt, Benjamin Recht, Oriol Vinyals
- Key Insights:
- A paradox in deep learning: large neural networks manage to generalize well and maintain a minimal gap between training and testing performance, contrary to what might be expected given their size.
- Small generalization errors are commonly linked to model traits or training regularization.
- The authors' experiments indicate these reasons don't fully account for the networks' generalization success.

### Experiment on Generalization in Deep Networks

- State-of-the-art convolutional neural networks for image classification, when trained with stochastic gradient methods, can easily fit to randomly labeled training data.
- This ability to fit random labels is not significantly impacted by the use of explicit regularization techniques.
- The phenomenon persists even when true image data is replaced with unstructured random noise.
- Theoretical work supports these findings by showing that neural networks with a sufficient number of parameters relative to data points can perfectly express any finite sample set.

### **Experimental Details**

#### • Data Modification:

- True labels in standard datasets (CIFAR10, ImageNet) were replaced with random labels.
- In an extension of the experiment, actual image pixels were replaced with completely random noise.

#### • Training Results:

- Neural networks reached zero training error on randomly labeled data, suggesting they can memorize the dataset.
- Test error was equivalent to random guessing due to lack of correlation between training and test labels.
- Training times were only slightly longer than with true labels, indicating ease of optimization even with random data.

### **Experimental Details**

#### • Impact of Noise:

- The introduction of noise to the images did not prevent neural networks from fitting the data.
- A progressive increase in noise led to a corresponding increase in generalization error, yet networks could still capture any signal left in the labels.

#### • Implications:

- These results challenge the role of VC-dimension, Rademacher complexity, and uniform stability in explaining generalization in neural networks.
- The networks' ability to fit random labels and noise points to a high capacity for memorization, which is not accounted for by traditional learning complexity measures.

Uniform convergence may be unable to explain generalization in deep learning

#### **Authors:**

Vaishnavh Nagarajan, Zico Kolter

#### Main Claim:

- The paper challenges the adequacy of uniform convergence as a tool for explaining the generalization behavior in overparameterized deep neural networks.
- It highlights a key finding: generalization bounds based on uniform convergence can paradoxically increase with the size of the training dataset.

### **Understanding Uniform Convergence**

- Uniform convergence is a concept in statistical learning theory. It describes how closely the empirical loss (loss on training data) of a learning algorithm converges to the expected loss (loss on the entire data distribution) uniformly over all hypotheses in a hypothesis class.
- This concept is crucial for establishing generalization bounds, which predict how well a model trained on a finite dataset will perform on unseen data.
- The idea is that if the empirical loss converges uniformly to the true loss across all hypotheses, one can be confident that a hypothesis with low empirical loss will also have low true loss, hence will generalize well.

### **Implications of the Findings**

- The paper demonstrates scenarios with overparameterized models (like deep neural networks with more parameters than training data points) where uniform convergence fails to explain generalization.
- It shows that even when considering only hypotheses output by gradient descent with low test errors, uniform convergence provides vacuous (ineffective) generalization guarantees.
- These findings cast doubt on the ability of uniform convergencebased bounds to fully explain why large neural networks generalize effectively, suggesting the need for alternative or additional theoretical frameworks.