

Machine learning: lecture 11

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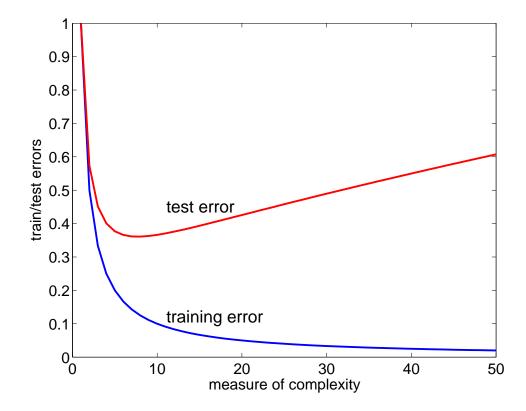


Topics

- Complexity and generalization
 - finite set of classifiers
 - VC-dimension, learning



Why care about "complexity"?



• We need a quantitative measure of complexity in order to be able to relate the training error (which we can observe) and the test error (that we'd like to optimize)



Finite case

- We'll start by considering only a finite number of possible classifiers, $h_1(\mathbf{x}), \ldots, h_M(\mathbf{x})$ (e.g., randomly chosen linear classifiers)
- Key questions:
 - 1. Given n training examples and M possible classifiers how far can the training and test errors be?
 - 2. How many training examples do we need so that the errors are close?
 - The answers will depend on M.



Finite case: definitions

$$\hat{\mathcal{E}}_{n}(i) = \frac{1}{n} \sum_{t=1}^{n} \underbrace{\mathsf{Loss}(y_{t}, h_{i}(\mathbf{x}_{t}))}_{t=1} = \text{empirical error of } h_{i}(\mathbf{x})$$
$$\mathcal{E}(i) = E_{(\mathbf{x}, y) \sim P} \{ \mathsf{Loss}(y, h_{i}(\mathbf{x})) \} = \text{expected error of } h_{i}(\mathbf{x})$$



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• Suppose we choose the classifier that minimizes the training error, $\hat{i}_n = \arg\min_{i=1,...,M} \hat{\mathcal{E}}_n(i)$, then

Training error
$$= \hat{\mathcal{E}}_n(\hat{i}_n)$$

Test error $= \mathcal{E}(\hat{i}_n)$



Finite case: errors

• The training and test errors,

Fraining error
$$= \hat{\mathcal{E}}_n(\hat{i}_n)$$

Test error $= \mathcal{E}(\hat{i}_n)$

are necessarily close if we can show that the errors are close for all the classifiers in our set:

$$|\hat{\mathcal{E}}_n(i) - \mathcal{E}(i)| \le \epsilon$$
, for all $i = 1, \dots, M$

• We can now express our key questions more formally in terms of $n,\ M,$ and ϵ



Finite case: key questions revisited

- Key questions (rewritten):
 - 1. Given n training examples and M possible classifiers, what is the smallest ϵ such that

$$\max_{i=1,\dots,M} |\hat{\mathcal{E}}_n(i) - \mathcal{E}(i)| \le \epsilon$$

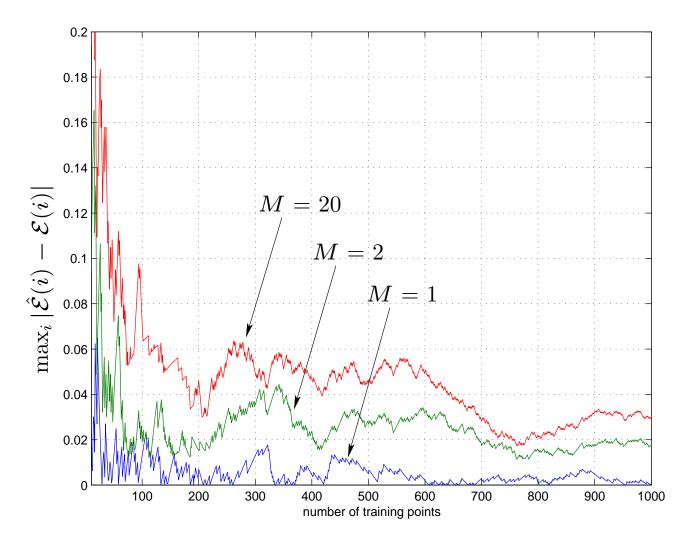
2. For a given ϵ how many training examples do we need so that

$$\max_{i=1,\dots,M} |\hat{\mathcal{E}}_n(i) - \mathcal{E}(i)| \le \epsilon$$

Since training examples are sampled at random from some underlying distribution, we can only answer these questions probabilistically.



Finite case: errors





Finite case: probabilistic statement

• We can relate n, M, and ϵ by requiring that with high probability, the empirical errors of all the classifiers in our set are ϵ -close to their expected errors:

$$P\left(\max_{i=1,\dots,M} |\hat{\mathcal{E}}_n(i) - \mathcal{E}(i)| \le \epsilon\right) \ge 1 - \delta$$

The probability is taken over the choice of the training set and $1-\delta$ specifies our confidence in the probabilistic statement.



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• Equivalently, we can bound the probability that the empirical error of some classifier in our set deviates more than ϵ from the expected error:

$$P\left(\max_{i=1,\dots,M} |\hat{\mathcal{E}}_n(i) - \mathcal{E}(i)| > \epsilon\right) \le \delta$$



 \bullet Let's fix $n,~M,~{\rm and}~\epsilon$ and try to find δ so that

$$P\left(\max_{i=1,\dots,M} |\hat{\mathcal{E}}_n(i) - \mathcal{E}(i)| > \epsilon\right) \le \delta$$

still holds. The probability is take over the choice of the training set.



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still holds. The probability is take over the choice of the training set.

By using the fact that $P(A \, {\rm or} \, B) \leq P(A) + P(B)$ we get

$$P\left(\max_{i} |\hat{\mathcal{E}}_{n}(i) - \mathcal{E}(i)| > \epsilon\right) \leq \sum_{i=1}^{M} P\left(|\hat{\mathcal{E}}_{n}(i) - \mathcal{E}(i)| > \epsilon \right)$$



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$$\leq \sum_{i=1}^{M} 2\exp(-2n\epsilon^{2}) \quad \text{(Chernoff)}$$



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$$\leq \sum_{i=1}^{M} 2\exp(-2n\epsilon^{2}) \quad \text{(Chernoff)}$$
$$= M \cdot 2\exp(-2n\epsilon^{2}) = \delta$$



• We are now able to relate n, M, ϵ , and δ :

$$M \cdot 2 \exp(-2n\epsilon^2) = \delta$$
, or $\epsilon = \sqrt{\frac{\log(M) + \log(2/\delta)}{2n}}$

• We can restate our result in terms of a bound on the expected error of any classifier in our set.

Theorem: With probability at least $1 - \delta$ over the choice of the training set, for all $i = 1, \ldots, M$

$$\mathcal{E}(i) \le \hat{\mathcal{E}}_n(i) + \epsilon(n, M, \delta)$$

where $\epsilon = \epsilon(n,M,\delta)$ is a "complexity penalty".



Measures of complexity

- Typically the set of classifiers is not a finite nor a countable set (e.g., the set of linear classifiers)
- There are still many ways of trying to capture the "effective" number of classifiers in such a set:
 - degrees of freedom (number of parameters)
 - Vapnik-Chervonenkis (VC) dimension
 - description length

etc.



VC-dimension: preliminaries

• A set of classifiers F: For example, this could be the set of all possible linear classifiers, where $h \in F$ means that

$$h(\mathbf{x}) = \operatorname{sign}\left(w_0 + \mathbf{w}_1^T \mathbf{x}\right)$$

for some values of the parameters w_0, \mathbf{w}_1 .



VC-dimension: preliminaries

• Complexity: how many different ways can we label n training points $\{\mathbf{x}_1, \ldots, \mathbf{x}_n\}$ with classifiers $h \in F$?

In other words, how many distinct binary vectors

$$[h(\mathbf{x}_1) h(\mathbf{x}_2) \dots h(\mathbf{x}_n)]$$

do we get by trying out each $h \in F$ in turn?

$$\begin{bmatrix} -1 & 1 & \dots & 1 &] & h_1 \\ [& 1 & -1 & \dots & 1 &] & h_2 \end{bmatrix}$$



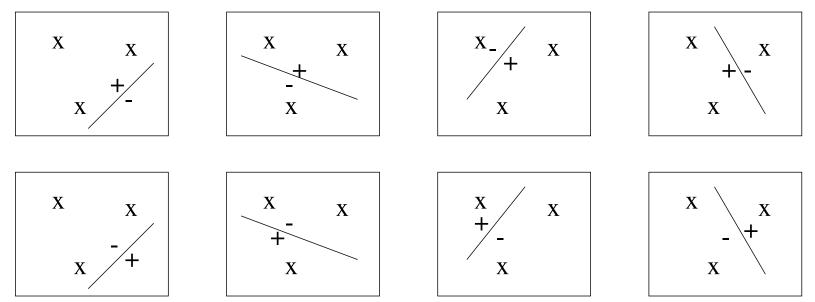
VC-dimension: shattering

• A set of classifiers F shatters n points $\{\mathbf{x}_1, \ldots, \mathbf{x}_n\}$ if

 $[h(\mathbf{x}_1) h(\mathbf{x}_2) \dots h(\mathbf{x}_n)], h \in F$

generates all 2^n distinct labelings.

 Example: linear decision boundaries shatter (any) 3 points in 2D

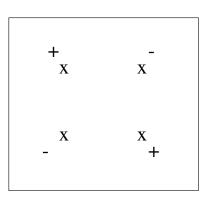


but not any 4 points...



VC-dimension: shattering cont'd

 We cannot shatter any set of 4 points in 2D with linear classifiers. For example, we cannot generate the following XOR-labeling:

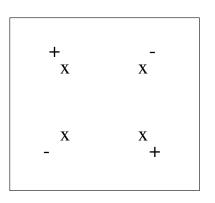


• More generally: the set of all d-dimensional linear classifiers can shatter exactly d + 1 points



VC-dimension: shattering cont'd

 We cannot shatter any set of 4 points in 2D with linear classifiers. For example, we cannot generate the following XOR-labeling:



- More generally: the set of all d-dimensional linear classifiers can shatter exactly d + 1 points
- **Definition:** The VC-dimension d_{VC} of a set of classifiers F is the number of points F can shatter



Learning and VC-dimension

• We learn something only after we no longer can shatter the training points (have more than d_{VC} training examples)

Rationale: suppose we have *n* training examples and labels $(\mathbf{x}_1, y_1), \ldots, (\mathbf{x}_n, y_n)$ and $n < d_{VC}$. Does the training set constrain our prediction for \mathbf{x}_{n+1} ?

Because we expect to be able to shatter n+1 points ($\leq d_{VC}$) it follows that we can find $h_1, h_2 \in F$, both consistent with training labels, but

$$h_1(\mathbf{x}_{n+1}) = 1, \quad h_2(\mathbf{x}_{n+1}) = -1$$

We therefore cannot determine which label to predict for \mathbf{x}_{n+1} .

Learning and VC-dimension

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