

CS 446/ECE 449: Machine Learning

Lecture 9: PAC Learning Theory (I)

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Recap: Supervised Learning Algorithms

Models we have learned so far:

Model	Linear?	Parametric?	Loss	Generative/ Discriminative
K-nearest neighbor	N	N	N/A	Discriminative
Naive Bayes	Y	Y	NLL	Generative
Logistic regression	Y	Y	Logistic/NLL	Discriminative
Linear SVM	Y	Y	Hinge	Discriminative
Kernelized SVM	N	N	Hinge	Discriminative
Decision Tree	N	N	N/A	Discriminative
AdaBoost	N	N	Exp	Discriminative

Note:

- NLL = negative log-likelihood
- Generative = modeling $\Pr(X, Y)$
- Discriminative = modeling $\Pr(Y|X)$

Lecture Today

- Bayes Error, Bayes Predictor
- Error Decomposition

Bayes Error Rate

So far we have learned many different classification algorithms. Beyond their different design choices, how should we compare their performance **theoretically**?

- For a given prediction problem, what is the optimal error that we can hope to achieve? Which predictor will achieve the optimal error?
- Given a problem and a model, how far is our model from the optimal predictor?

Bayes Error Rate

The learning process:

- We can choose a predictor f from some pre-defined class of functions \mathcal{F} , e.g., the class of linear predictors, decision trees, kernel machines, neural networks, etc.

We also have our training data $\mathcal{D} := \{(x^{(i)}, y^{(i)})\}_{i=1}^n \sim \mu$ sampled independently and identically (iid) from the underlying distribution μ over $\mathcal{X} \times \mathcal{Y}$

We can then talk about two error measures (classification):

$$\text{Training error: } \hat{\varepsilon}_{\mathcal{D}}(f) := \frac{1}{n} \sum_{i=1}^n \mathbb{I}(f(x^{(i)}) \neq y^{(i)})$$

$$\text{Test error: } \varepsilon_{\mu}(f) := \mathbb{E}_{\mu} [\mathbb{I}(f(X) \neq Y)] = \Pr_{\mu}(f(X) \neq Y)$$

We are interested in finding f that minimizes the test error but we can only observe the training error

Bayes Error Rate

Bayes error rate: the theoretically minimum test error that can be achieved:

$$\text{Bayes error: } \varepsilon_{\mu}^* := \inf_{f: \mathcal{X} \rightarrow \mathcal{Y}} \varepsilon_{\mu}(f)$$

Assuming X is a continuous RV and let $p(x)$ be the probability density of X . Then for any classifier $f: \mathcal{X} \rightarrow \{0,1\}$, we have:

$$\begin{aligned} \varepsilon_{\mu}(f) &= \Pr_{\mu}(f(X) \neq Y) \\ &= \int_{\mathcal{X}} \left(\Pr_{\mu}(Y = 1 | X = x) \cdot \mathbb{I}(f(x) = 0) + \Pr_{\mu}(Y = 0 | X = x) \cdot \mathbb{I}(f(x) = 1) \right) p(x) \, dx \\ &\geq \int_{\mathcal{X}} \min \left\{ \Pr_{\mu}(Y = 1 | X = x), \Pr_{\mu}(Y = 0 | X = x) \right\} p(x) \, dx \\ &= \mathbb{E}_{\mu} \left[\min \left\{ \Pr_{\mu}(Y = 1 | X), \Pr_{\mu}(Y = 0 | X) \right\} \right] \\ &= \frac{1}{2} - \frac{1}{2} \mathbb{E}_{\mu} [|2\eta(X) - 1|] \end{aligned}$$

where $\eta(X) := \Pr_{\mu}(Y = 1 | X)$ is the conditional probability

Bayes Error Rate

Bayes error depends on the distribution μ :

$$\begin{aligned}\varepsilon_\mu(f) &= \Pr_\mu(f(X) \neq Y) \\&= \int_{\mathcal{X}} \left(\Pr_\mu(Y = 1 | X = x) \cdot \mathbb{I}(f(x) = 0) + \Pr_\mu(Y = 0 | X = x) \cdot \mathbb{I}(f(x) = 1) \right) p(x) \, dx \\&\geq \int_{\mathcal{X}} \min \left\{ \Pr_\mu(Y = 1 | X = x), \Pr_\mu(Y = 0 | X = x) \right\} p(x) \, dx \\&= \mathbb{E}_\mu \left[\min \left\{ \Pr_\mu(Y = 1 | X), \Pr_\mu(Y = 0 | X) \right\} \right] \\&= \boxed{\frac{1}{2} - \frac{1}{2} \mathbb{E}_\mu [|2\eta(X) - 1|]} \quad \text{Bayes error rate: } \varepsilon_\mu^*\end{aligned}$$

- The Bayes error only depends on the distribution μ
- It's unknown since we don't know μ in practice
- It's always ≤ 0.5

Bayes Error Rate

The classifier that achieves the Bayes error is called the **Bayes classifier**, and it has the following form:

$$\eta(X) := \Pr(Y = 1 \mid X)$$

$$f_{\text{Bayes}}(X) := \begin{cases} 1 & \text{if } \eta(X) \geq \frac{1}{2} \\ 0 & \text{otherwise} \end{cases}$$

- Again, this is unknown since we don't know μ
- Recall the proof:

$$\begin{aligned} \varepsilon_{\mu}(f) &= \int_{\mathcal{X}} \left(\Pr(Y = 1 \mid X = x) \cdot \mathbb{I}(f(x) = 0) + \Pr(Y = 0 \mid X = x) \cdot \mathbb{I}(f(x) = 1) \right) p(x) \, dx \\ &\geq \int_{\mathcal{X}} \min \left\{ \Pr(Y = 1 \mid X = x), \Pr(Y = 0 \mid X = x) \right\} p(x) \, dx \\ &= \frac{1}{2} - \frac{1}{2} \mathbb{E}_{\mu} [|2\eta(X) - 1|] \end{aligned}$$

Think: when will $\varepsilon_{\mu}^* = 0$? when will $\varepsilon_{\mu}^* = 0.5$?

Bayes Error Rate

Intuitively, the Bayes error is a measure of the “noise” in the underlying distribution:

Bayes error rate: $\varepsilon_{\mu}^* = \frac{1}{2} - \frac{1}{2} \mathbb{E} [|2\eta(X) - 1|]$

$$\eta(X) := \Pr(Y = 1 \mid X)$$

- If $\forall x, \eta(x) = 1$, or $\eta(x) = 0$, then $\varepsilon_{\mu}^* = 0$
- If $\forall x, \eta(x) = \frac{1}{2}$, then $\varepsilon_{\mu}^* = \frac{1}{2}$

Bayes Error Rate

Intuitively, the Bayes error is a measure of the “noise” in the underlying distribution:

Bayes error rate: $\epsilon_{\mu}^* = \frac{1}{2} - \frac{1}{2} \mathbb{E} [|2\eta(X) - 1|]$

$$\eta(X) := \Pr(Y = 1 \mid X)$$

Example: Suppose we have the following data generative process. There exists a vector w^* , such that for each x , the labels are generated in the following process:

- First, compute the label $y = \text{sgn}(w^{*\top} x)$
- Then, with probability $0 < p < 0.5$, flip the label y

Question: What's the Bayes error rate for this example?

Bayes Error Rate

The concept is not unique to classification problems. For regression problems, under the squared loss:

$$\begin{aligned}\forall f, \varepsilon_\mu(f) &= \mathbb{E}_\mu [(f(X) - Y)^2] \\ &= \mathbb{E}_X \mathbb{E}_Y [(f(X) - Y)^2 | X] \\ &= \mathbb{E}_X \mathbb{E}_Y [(f(X) - \mathbb{E}[Y|X] + \mathbb{E}[Y|X] - Y)^2 | X] \\ &= \mathbb{E}_X \mathbb{E}_Y [(f(X) - \mathbb{E}[Y|X])^2 + (\mathbb{E}[Y|X] - Y)^2 \\ &\quad + \cancel{2(f(X) - \mathbb{E}[Y|X])(\mathbb{E}[Y|X] - Y)} | X] \\ &\geq \mathbb{E}_X \mathbb{E}_Y [(\mathbb{E}[Y|X] - Y)^2 | X]\end{aligned}$$

Bayes error rate: $= \mathbb{E}_X \text{Var}[Y | X]$

Bayes optimal regressor: $f_{\text{Bayes}}(X) = \mathbb{E}[Y | X]$

Bayes Error Rate

Again, the Bayes error in regression can also be understood as a measure of the “noise” in the underlying distribution:

Bayes error rate: $\epsilon_{\mu}^* = \mathbb{E} \text{Var}[Y|X]$

Example: Suppose we have the following data generative process. There exists a vector w^* , such that for each x , the labels are generated in the following process:

- First, compute the label $y = w^{*\top} x$
- Then, inject a white noise $\epsilon \sim \mathcal{N}(0, \delta^2)$ into the label so that $y \leftarrow y + \epsilon$

Question: What's the Bayes error under the squared loss for this example?

Bayes Error Rate

The optimal error a learner can hope to achieve also depends on the class of functions \mathcal{F} it can choose from, called **hypothesis class**

For binary classification problems:

- If \mathcal{F} contains all the binary functions, then $\inf_{f \in \mathcal{F}} \varepsilon_{\mu}(f) = \varepsilon_{\mu}^*$
- If \mathcal{F} is very restricted, e.g., only contains **constant functions**, then $\inf_{f \in \mathcal{F}} \varepsilon_{\mu}(f) = \min\{\Pr(Y = 0), \Pr(Y = 1)\}$

Clearly, in the second case, the error is larger than the Bayes error.

Bayes Error Rate

The optimal error a learner can hope to achieve also depends on the class of functions \mathcal{F} it can choose from, called **hypothesis class**

For **regression problems** under mean-squared error:

- If \mathcal{F} contains all the real-valued functions, then $\inf_{f \in \mathcal{F}} \varepsilon_{\mu}(f) = \varepsilon_{\mu}^*$
- If \mathcal{F} is very restricted, e.g., only contains constant functions, then $\inf_{f \in \mathcal{F}} \varepsilon_{\mu}(f) = \text{Var}[Y]$

In the second case, the error is larger than the Bayes error by the law of total variance: $\text{Var}[Y] = \mathbb{E}\text{Var}[Y|X] + \text{Var}\mathbb{E}[Y|X] \geq \mathbb{E}\text{Var}[Y|X]$

Bayes Error Rate

Summary:

Binary classification:

Bayes error rate: $\varepsilon_{\mu}^* = \mathbb{E} \min \{ \Pr(Y = 1 | X), \Pr(Y = 0 | X) \}$

Bayes optimal classifier: $f_{\text{Bayes}}(X) := \begin{cases} 1 & \text{if } \Pr(Y = 1 | X) \geq \frac{1}{2} \\ 0 & \text{otherwise} \end{cases}$

Regression with squared loss:

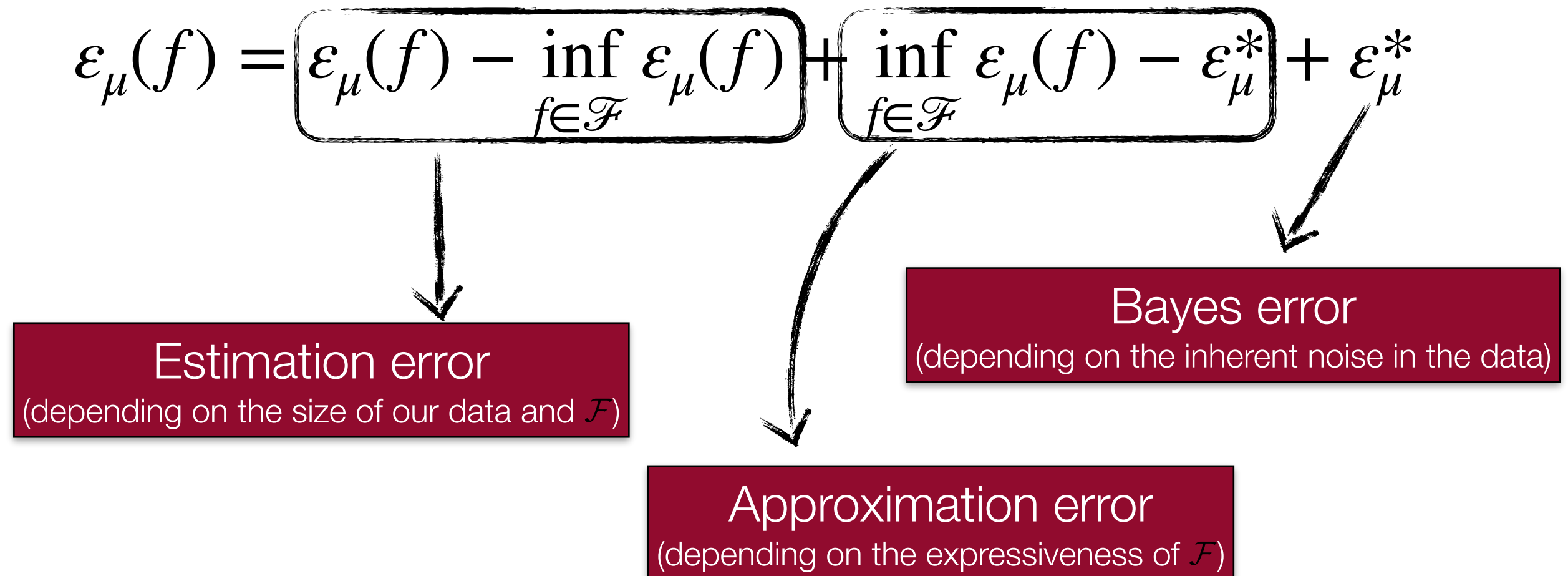
Bayes error rate: $\varepsilon_{\mu}^* = \mathbb{E} \text{Var}[Y | X]$

Bayes optimal regressor: $f_{\text{Bayes}}(X) = \mathbb{E}[Y | X]$

Error Decomposition

For a given hypothesis class \mathcal{F} , we may have $f_{\text{Bayes}} \notin \mathcal{F}$. In this case we cannot hope to achieve ε_{μ}^* , but instead $\inf_{f \in \mathcal{F}} \varepsilon_{\mu}(f)$.

Error decomposition: $\forall f \in \mathcal{F}$:



Error Decomposition

Error decomposition: $\forall f \in \mathcal{F}$:

$$\varepsilon_{\mu}(f) = \underbrace{\varepsilon_{\mu}(f) - \inf_{f \in \mathcal{F}} \varepsilon_{\mu}(f)}_{\text{Estimation error}} + \underbrace{\inf_{f \in \mathcal{F}} \varepsilon_{\mu}(f) - \varepsilon_{\mu}^*}_{\text{Approximation error}} + \underbrace{\varepsilon_{\mu}^*}_{\text{Bayes error}}$$

Estimation error
(depending on the size of our data and \mathcal{F})

Approximation error
(depending on the expressiveness of \mathcal{F})

Bayes error
(depending on the inherent noise in the data)

- Often the case, there is a trade-off between the estimation error and the approximation error
- If \mathcal{F} is more expressive, then the approximation error gets smaller but the estimation error gets larger
- If \mathcal{F} is more restricted, then the approximation error gets larger but the estimation error gets smaller (assume the size of training data is fixed)

Error Decomposition

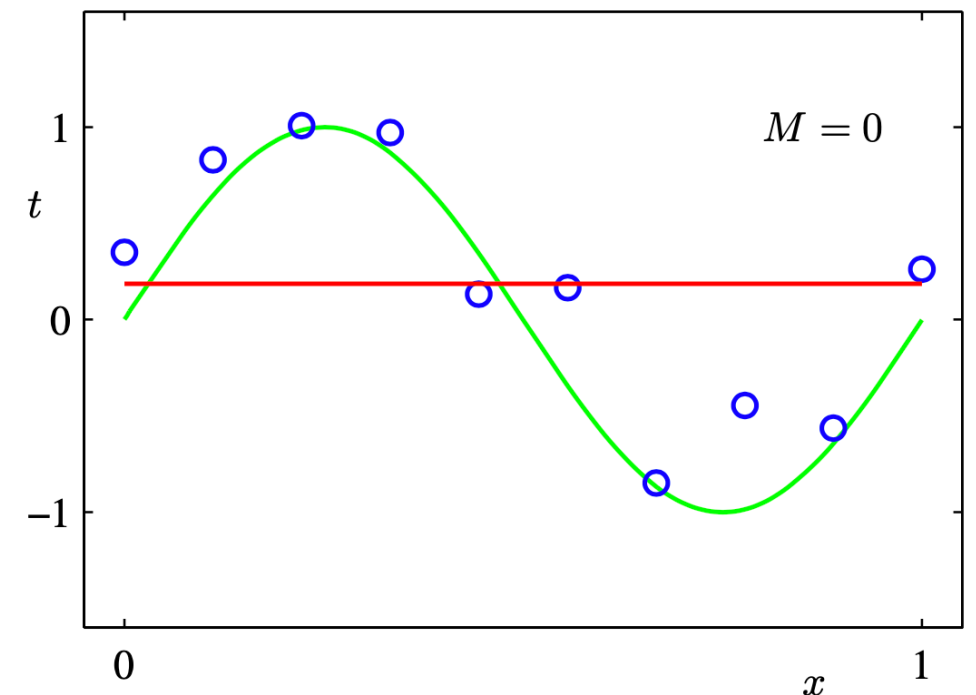
Example: fitting a trigonometric function with polynomials
(degree = M)

Degree 0

Bayes classifier



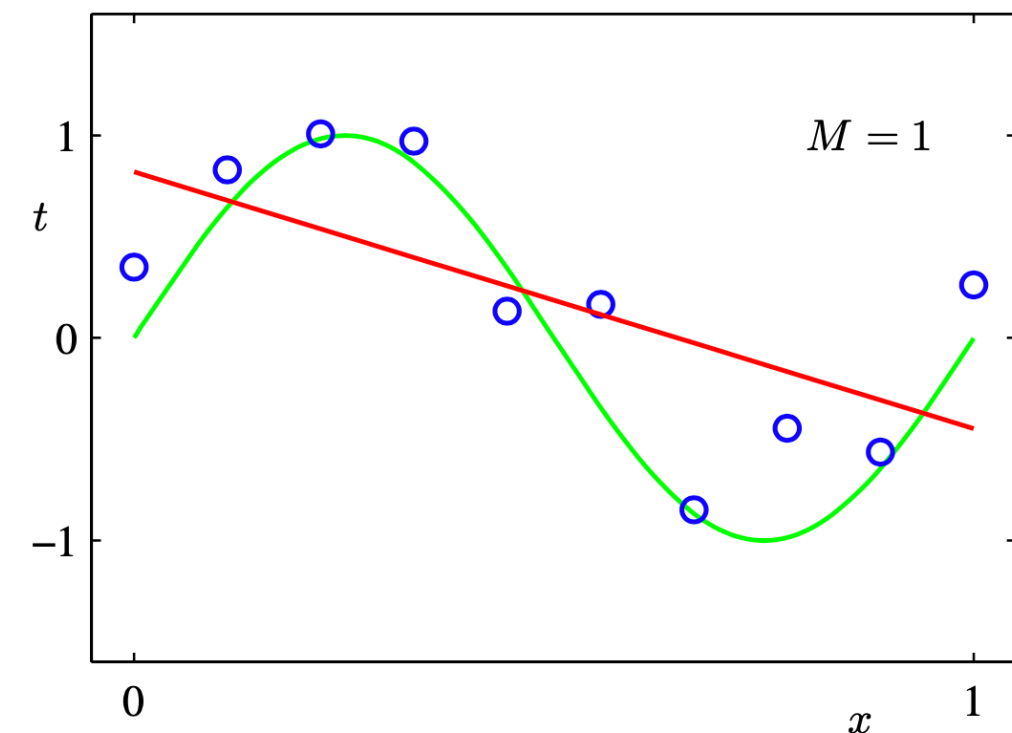
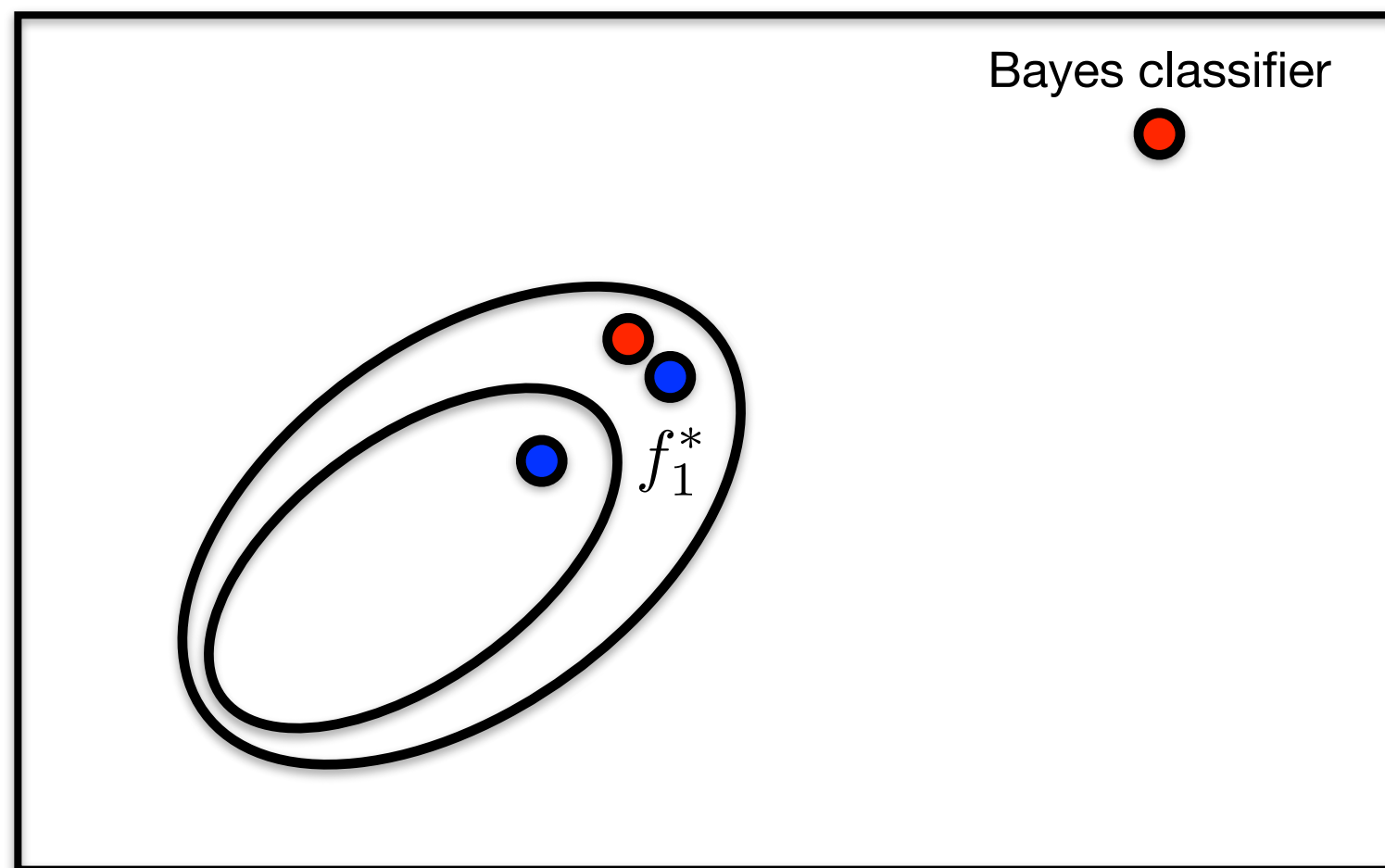
f_0^*



Error Decomposition

Example: fitting a trigonometric function with polynomials
(degree = M)

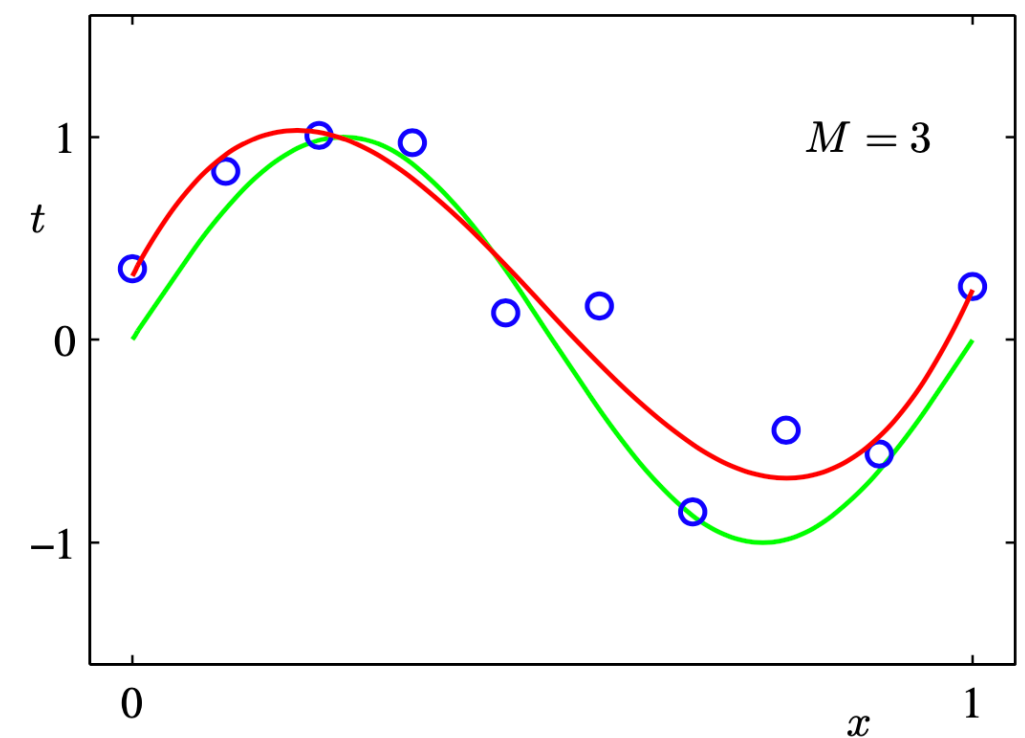
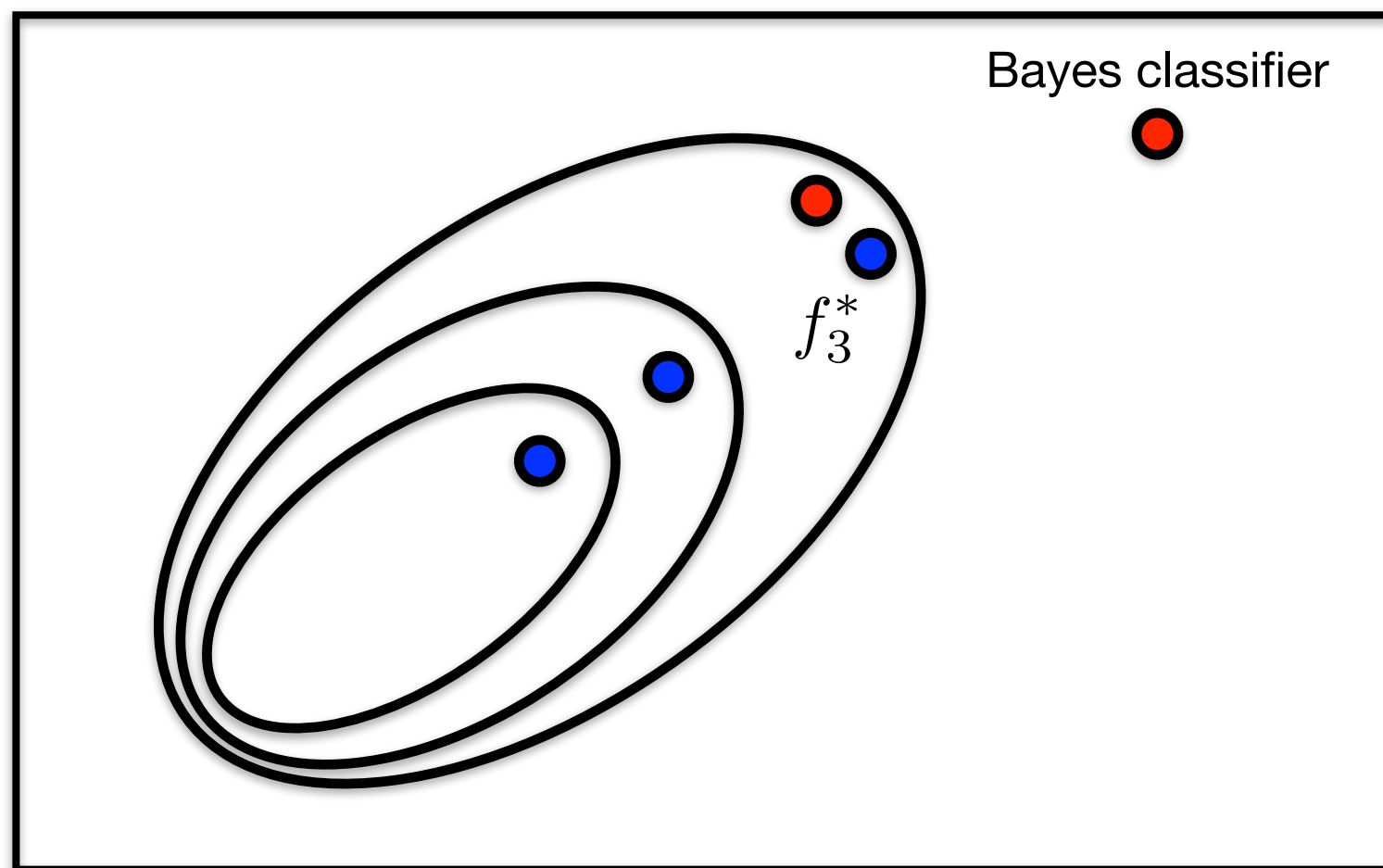
Degree 1



Error Decomposition

Example: fitting a trigonometric function with polynomials
(degree = M)

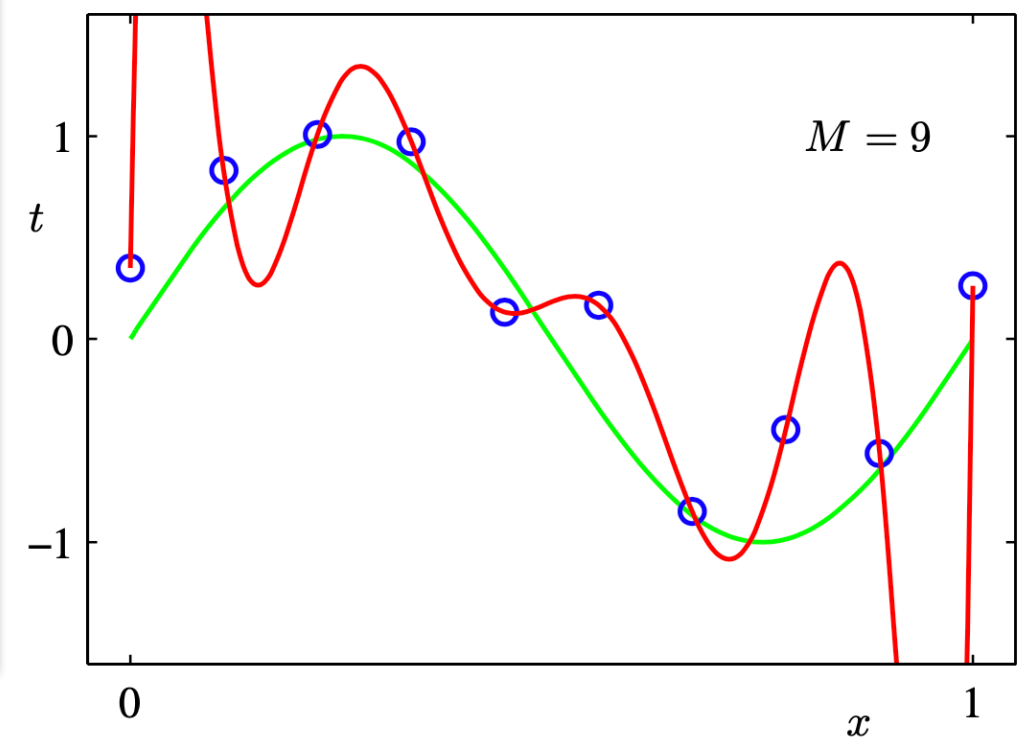
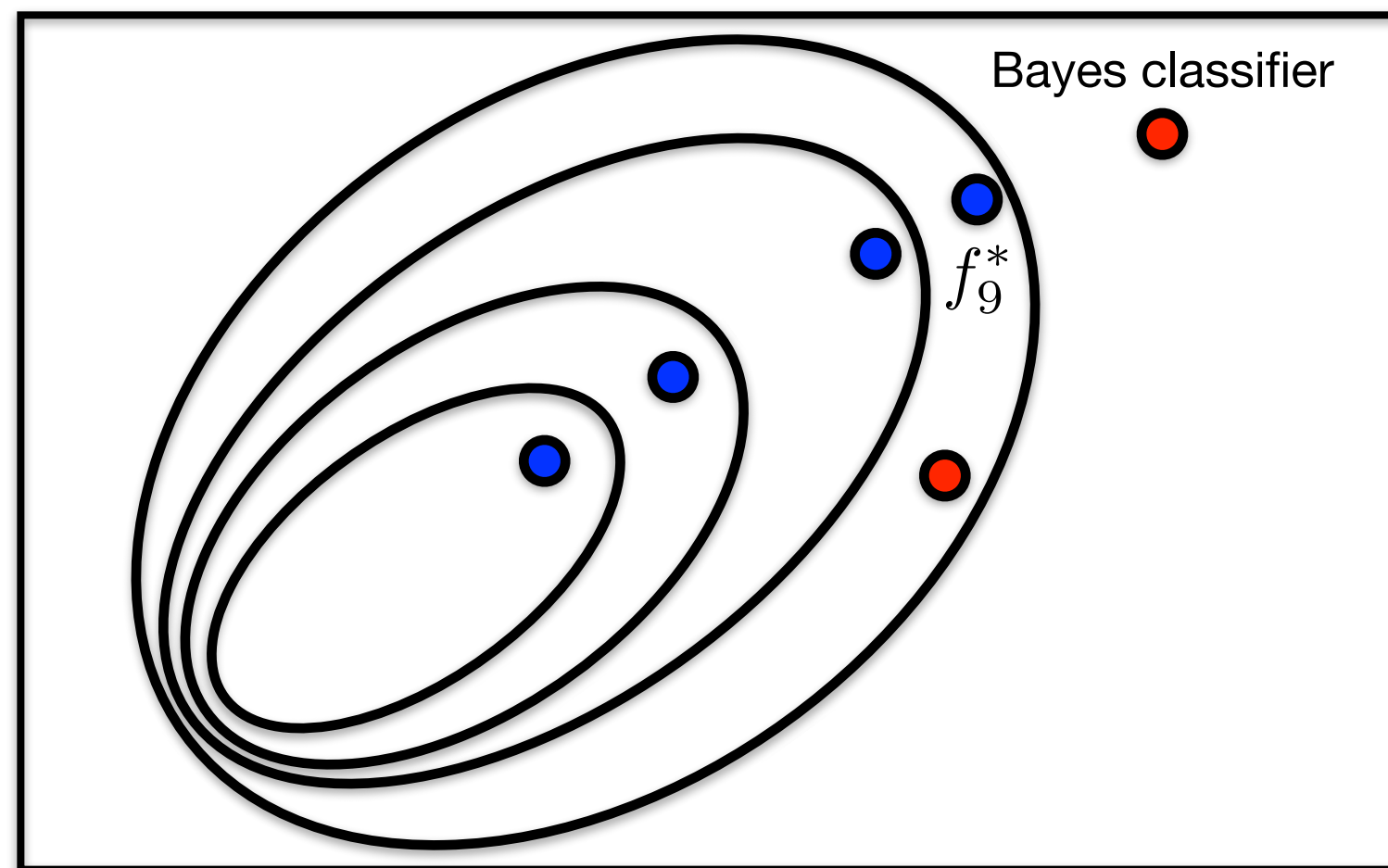
Degree 3



Error Decomposition

Example: fitting a trigonometric function with polynomials
(degree = M)

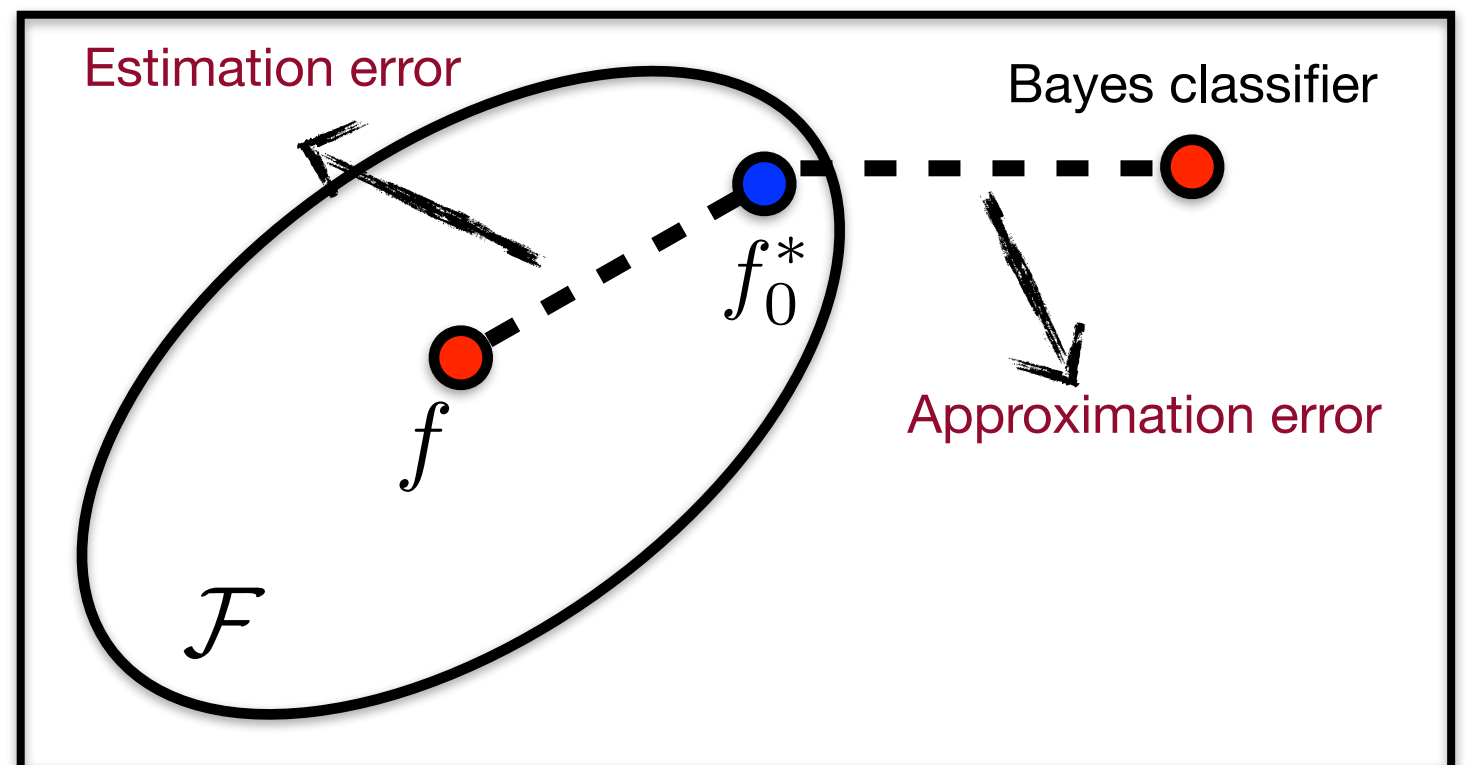
Degree 9



Error Decomposition

Reminder: the approximation error only depends on \mathcal{F} while the estimation error depends on both \mathcal{F} and data

- We should aim to minimize the estimation error
- How does the estimation error depend on the sample size, the expressiveness/richness of \mathcal{F} , or the distribution μ ?
- Ideally, for a fixed hypothesis class \mathcal{F} , could we ensure that the estimation error goes to 0 as the sample size n increases?



Next Time

- Probably Approximately Correct (PAC) framework
- High-probability generalization bound
- Vapnik–Chervonenkis dimension (VC dim)