

# Lecture 2: Supervised learning continued

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Send questions to [6.867-staff@lists.csail.mit.edu](mailto:6.867-staff@lists.csail.mit.edu) for faster responses!

All homework-related questions should be posted by Friday!

**Supervised learning:** Given a few examples find a predictor function that works for new examples.

We are given a training set. We'll primarily talk about batch / offline supervised classification where training data is given upfront.

**Training set:**  $S_n = \{(x^i, y^i), i = 1, \dots, n\}, (x^i, y^i) \sim p^*$

We select  $\hat{h}: X = R^d \rightarrow Y = \{-1, 1\}$

Goal is to minimize **generalization error** (risk):

$$R(\hat{h}) = E_{(x, y) \sim p^*} \{ \text{Loss}(y, \hat{h}(x)) \}$$

(expected value of loss on pairs sampled over  $p^*$ )

$$\text{Loss}(y, \hat{h}(x)) = \begin{cases} 1, & \text{if } y \neq \hat{h}(x) \\ 0, & \text{o. w.} \end{cases}$$

## Generative approach to solve the supervised learning problem

If I knew  $p^*$  I could perform optimal. I could evaluate what the generalization error is and find an  $\hat{h}$  that would minimize it.

Estimate  $\hat{p}(x, y)$  based on the training set  $S_n$ . We need constraints for estimating.

Once we have a guess on  $p^*$  we'll use a predictor  $\hat{h}(\cdot) = \underset{h}{\operatorname{argmin}} E_{(x, y) \sim \hat{p}} \{ \text{Loss}(y, h(x)) \}$

## Discriminative approach

$$\hat{R}_n(h) = \frac{1}{n} \sum_{i=1}^n \text{Loss}(y^i, h(x^i))$$

$h \in H$  – set of classifiers

We want to find an  $h$  minimize  $\hat{R}_n(h)$

Use  $\hat{h}(\cdot) = \underset{h \in H}{\operatorname{argmin}} \hat{R}_n(h)$

## Linear classifiers

$$X = R^d$$

$$Y = \{-1, 1\}$$

$$h(x; \theta, \theta_0) = \text{sign}(\theta x + \theta_0) = \begin{cases} +1, & \text{if } \theta x + \theta_0 > 0 \\ -1, & \text{otherwise} \end{cases}$$

You can get arbitrarily complex classifiers if you know how to handle linear classifiers.

$$\theta \vec{x} + \theta_0 = 0 \Leftrightarrow \theta(\vec{x} - \vec{x}_0) = 0 \Rightarrow \theta \vec{x}_0 = -\theta_0 \Rightarrow \theta_0 = -\theta \vec{x}_0$$

Distance to boundary  $\frac{y(\theta x)}{\|\theta\|}$ , where  $\|\theta\|$  = norm of theta

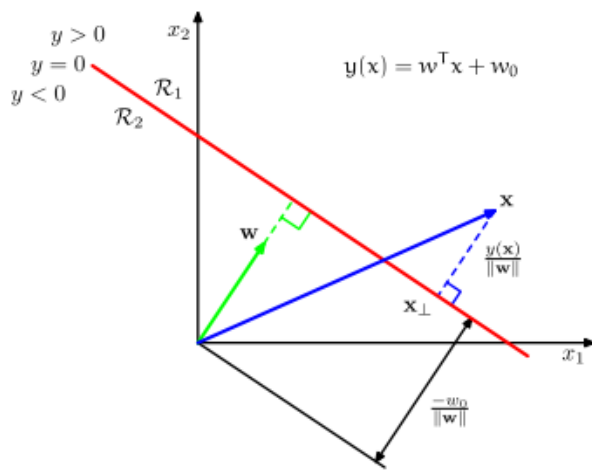


Figure 1: Geometry of linear discriminant functions (Bishop Figure 4.1).

TODO: Put graph in

## Training without errors

**Assumption 1:** Training examples are linearly separable with margin  $\gamma$ :

$$\exists \theta^* \text{ s.t. } \forall i, \frac{y^i \theta^* x^i}{\|\theta^*\|} > \gamma, \gamma > 0, i = 1, \dots, n$$

This says that all training examples are at least distance  $\gamma$  from the boundary.

Put another way, this means the examples are **linearly separable**.

**Assumption 2:** Training examples are bounded by a sphere/circle of radius  $r$ :  $\|x^{(i)}\| \leq r, i = 1, \dots, n$

## Perceptron algorithm

- Start at step 0:  $\theta^{(0)} = 0$  (vector)
- Cycle through training samples correcting errors

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- If  $y^{(i)}(\theta^{(k)} x^{(i)}) \leq 0$  (mistake), then  $\theta^{(k+1)} = \theta^{(k)} + y^{(i)} x^{(i)}$

Thus, if our two assumptions hold, then our perceptron algorithm makes at most  $\frac{r^2}{\gamma^2}$  mistakes.

The number of mistakes does not depend on the number of training examples or on the dimension of  $X$ ,  $\dim(X)$

## Training without errors online

Assumption 1: Training examples are linearly separable with margin  $\gamma$ :

$$\exists \theta^* s. t. \forall \frac{y^i \theta^* x^i}{\|\theta^*\|} > \gamma, \gamma > 0, i = 1, \dots, \infty$$

Assumption 2: Training examples are bounded by a sphere/circle or radius  $r$ :  $\|x^{(i)}\| \leq r, i = 1, \dots, \infty$

## Perceptron algorithm online

- Start at step 0:  $\theta^{(0)} = 0$  (vector)
- Cycle through training samples correcting errors
- If  $y^{(i)}(\theta^{(k)} x^{(i)}) \leq 0$  (mistake), then  $\theta^{(k+1)} = \theta^{(k)} + y^{(i)} x^{(i)}$

Once again, if our two assumptions hold, then our perceptron algorithm makes at most  $\frac{r^2}{\gamma^2}$  mistakes.

Why the  $r^2/\gamma^2$  bound?

$\cos(\theta^k, \theta^*) = \frac{\theta^k \theta^*}{\|\theta^k\| \times \|\theta^*\|}$ , where  $\theta^k$  is theta after  $k$  updates and theta star is the theta we assume exists

**Step 1:** Show that as we keep updating  $\frac{\theta^k \theta^*}{\|\theta^*\|} \geq k\gamma$

$$\frac{\theta^k \theta^*}{\|\theta^*\|} = \frac{(\theta^{k-1} + y^i x^i) \times \theta^*}{\|\theta^*\|} = \frac{\theta^{k-1} \times \theta^*}{\|\theta^*\|} + \frac{y^i x^i \times \theta^*}{\|\theta^*\|}$$

$$\frac{y^i x^i \times \theta^*}{\|\theta^*\|} \geq \gamma$$

$$\Rightarrow \frac{\theta^k \theta^*}{\|\theta^*\|} \geq k\gamma$$

**Step 2:** Norm of our parameter vector does not increase too high:  $\|\theta^k\|^2 \leq kr^2$

- We only update based on a mistake. We are correcting mistakes, which keeps the norm in check.

$$\cos(\theta^k, \theta^*) = \frac{\theta^k \theta^*}{\|\theta^k\| \times \|\theta^*\|} \geq \frac{k\gamma}{\|\theta^k\|} \geq \frac{k\gamma}{\sqrt{k} \times r} = \sqrt{k} \frac{\gamma}{r}$$

Why use the perceptron algorithm when we can find the maximum margin linear separator directly? We know  $\gamma$ .

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## **Maximum margin linear separator**

$$\text{minimizing } \frac{1}{2} \|\theta\|^2$$

$$y^i \theta x^i \geq 1, \forall i$$

Unique answer