

# CSCE 978 Lecture 3: Risk and Loss Functions\*

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## Introduction

- In Lecture 1 we mentioned our desire to infer a “good” classifier
- What does this mean?!?!?
- There are many ways to define “goodness”, even for binary classification

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## Outline

- Loss functions
  - Binary classification
  - Regression
- Expected risk
- Sections 1.3, 3.1–3.2 (also read Section 3.5)

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## Loss Functions

- D3.1** Let  $(x, y, f(x)) \in \mathcal{X} \times \mathcal{Y} \times \mathcal{Y}$  be the pattern  $x$ , its true label  $y$  and a prediction  $f(x)$  of  $y$ . A [loss function](#) is a mapping  $c : \mathcal{X} \times \mathcal{Y} \times \mathcal{Y} \rightarrow [0, \infty)$  with the property  $c(x, y, y) = 0$  for all  $x \in \mathcal{X}$  and  $y \in \mathcal{Y}$
- $c$  is always  $\geq 0$  so we can't use good predictions to “undo” bad ones
  - It is always possible to get 0 loss on pattern  $x$  by predicting correctly
  - Our choice of loss function will depend on considerations of computational complexity and statistical properties

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## Loss Functions

### Binary Classification

- Count number of misclassifications:

$$c(x, y, f(x)) = \begin{cases} 0 & \text{if } y = f(x) \\ 1 & \text{otherwise} \end{cases}$$

- Same as above, but penalty is input-dependent:

$$c(x, y, f(x)) = \begin{cases} 0 & \text{if } y = f(x) \\ \tilde{c}(x) & \text{otherwise} \end{cases}$$

- E.g. if  $y \in \{\text{rocks, diamonds}\}$  then penalty for “false diamond” classification depends on  $x$ ’s weight

- Can also have different values for false positive ( $y = -1$ ) and false negative ( $y = +1$ ) errors

- If  $y \in \{\text{cancer}, \neg\text{cancer}\}$  then FP results in unnecessary treatment, but FN can be fatal

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## Loss Functions

### Binary Classification

(cont’d)

- If  $f(x)$  is real-valued and  $y \in \{-1, +1\}$ , can think of  $\text{sign}(f(x))$  as prediction and  $|f(x)|$  as a confidence. Then a highly confident incorrect prediction can be penalized more, as can low-confidence correct predictions:

- Soft margin loss:

$$\begin{aligned} c(x, y, f(x)) &= \max(0, 1 - yf(x)) \\ &= \begin{cases} 0 & \text{if } yf(x) \geq 1 \\ 1 - yf(x) & \text{otherwise} \end{cases} \end{aligned}$$

- Logistic loss:

$$c(x, y, f(x)) = \ln(1 + \exp(-yf(x)))$$

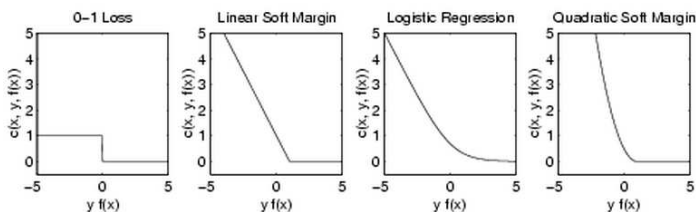
- Both penalize a lot for confident, incorrect predictions, penalize a little for low confidence, and don’t penalize much or at all for confident, correct predictions

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## Loss Functions

### Binary Classification

(cont’d)



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## Loss Functions

### Regression

- In regression,  $\mathcal{Y} \subseteq \mathbb{R}$  rather than  $\mathcal{Y} = \{-1, +1\}$

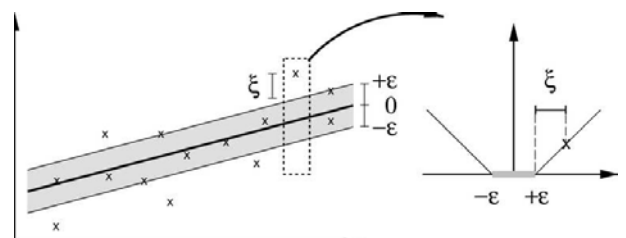
- Thus we’re interested in how far off our prediction  $f(x)$  is

- Squared loss (very popular):

$$c(x, y, f(x)) = (f(x) - y)^2$$

- Can extend soft margin loss to  $\epsilon$ -insensitive loss, which doesn’t penalize for close predictions:

$$c(x, y, f(x)) = |f(x) - y|_\epsilon = \max(|f(x) - y| - \epsilon, 0)$$



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## Loss Functions

### Practical Considerations

- Want loss function to be:
  - Cheap to compute
  - Have few discontinuities in first derivative
  - Convex (to ensure unique global optimum)
  - Yield computationally efficient solutions for learning
  - Resistant to outliers/noise

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## Risk

- A loss function measures error on individual examples
- Our ultimate goal is to minimize loss on new (yet unseen) examples
- How do we measure this?
  - Without making certain assumptions, this is very difficult or even impossible
  - Assume that there is a probability distribution  $P(x, y)$  on  $\mathcal{X} \times \mathcal{Y}$  that governs generation of patterns and labels
    - \* Assume the pairs  $(x, y)$  are drawn iid (independent and identically distributed) according to  $P(x, y)$
    - \* Generally, we won't make specific assumptions about the nature of  $P(x, y)$
  - $P(y | x)$  = conditional probability of getting label  $y$  given that  $x$  is the pattern (so  $x$  could have a different label on each draw)

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## Risk Definitions

- For now, assume we know all the new patterns we'll ever classify; call these the test patterns  $x'_1, \dots, x'_{m'}$  (note we do not know the labels until after we make predictions)

**D3.2** When test set  $x'_1, \dots, x'_{m'}$  already known, goal is to minimize the expected error on the test set:

$$R_{\text{test}}[f] := \frac{1}{m'} \sum_{i=1}^{m'} \int_{\mathcal{Y}} c(x'_i, y, f(x'_i)) dP(y | x'_i)$$

- Often, minimizing  $R_{\text{test}}[f]$  not realistic since typically don't know test set a priori
  - One exception: querying fixed collection of images, biological sequences, etc.

**D3.3** The expected risk (expected loss) wrt  $P$  &  $c$ :

$$\begin{aligned} R[f] &:= \mathbf{E}[R_{\text{test}}[f]] = \mathbf{E}[c(x, y, f(x))] \\ &= \int_{\mathcal{X} \times \mathcal{Y}} c(x, y, f(x)) dP(x, y) \end{aligned}$$

- Not realistic since we don't know  $P(x, y)$

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## Risk Definitions (cont'd)

- To get a handle on  $P(x, y)$ , assume it's the same one that generated the training set
- Now use the training patterns to estimate  $P(x, y)$

**D3.4** The empirical risk is

$$\begin{aligned} R_{\text{emp}}[f] &:= \int_{\mathcal{X} \times \mathcal{Y}} c(x, y, f(x)) p_{\text{emp}}(x, y) dx dy \\ &= \frac{1}{m} \sum_{i=1}^m c(x_i, y_i, f(x_i)) \end{aligned}$$

- Easy to compute and generally straightforward to minimize (depending on  $c$ )
- So now all we have to do is find an  $f$  that minimizes  $R_{\text{emp}}[f]$ , use that as our predictor, and we're done, right?

(Can we go home now?)

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**NO!**

- We have to appropriately restrict the set of functions  $\mathcal{F}$  from which we choose  $f$ 
  - Otherwise,  $R_{\text{emp}}[f]$  won't approximate  $R[f]$ , which is what we want to minimize
- E.g. what if  $\mathcal{F}$  is the set of all functions from  $\mathcal{X}$  to  $\mathcal{Y}$ ?
  - Then our learning algorithm could get  $R_{\text{emp}}[f] = 0$  by simply storing the  $(x, y)$  pairs in a table (i.e. memorization)
  - Is this learning? Will it generalize well?
- Restricting  $\mathcal{F}$  has been looked from many perspectives: e.g. VC dimension, bias, structural risk minimization
- Our approach (called regularization) will quantify the “power” (“expressiveness”) of each  $f$  and minimize a sum of this and  $R_{\text{emp}}[f]$ 
  - Special case: minimum description length principle

**Topic summary (over Lectures 2 and 3) due in 1 week!**