Projects **Machine Learning Course**Fall 2018

**EPFL** 

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Mock Midterm Exam - Nov 19, 2018

# 1 Subgradient Descent

Derive the (sub)gradient descent update rule for a one-parameter linear model using the Mean Absolute Error,

$$\mathcal{L}_{\mathsf{MAE}}(\mathbf{X}, \mathbf{y}, w) = \frac{1}{N} \sum_{n=1}^{N} |wx_n - y_n|.$$

Hint: The function f(x) = |ax| is a composition of two simpler function. Use the chain rule!

### 2 Multiple-Output Regression

Let  $S = \{(\mathbf{y}_n, \mathbf{x}_n)\}_{n=1}^N$  be our training set for a regression problem with  $\mathbf{x}_n \in \mathbb{R}^D$  as usual. But now  $\mathbf{y}_n \in \mathbb{R}^K$ , i.e., we have K outputs for each input. We want to fit a linear model for each of the K outputs, i.e., we now have K regressors  $f_k(\cdot)$  of the form

$$f_k(\mathbf{x}) = \mathbf{x}^\top \mathbf{w}_k,$$

where each  $\mathbf{w}_k^{\top} = (w_{k1}, \cdots, w_{kD})$  is the weight vector corresponding to the k-th regressor. Let  $\mathbf{W}$  be the  $D \times K$  matrix whose columns are the vectors  $\mathbf{w}_k$ .

Our goal is to minimize the following cost function  $\mathcal{L}$ :

$$\mathcal{L}(\mathbf{W}) = \sum_{k=1}^{K} \sum_{n=1}^{N} \frac{1}{2\sigma_{k}^{2}} (y_{nk} - \mathbf{x}_{n}^{\mathsf{T}} \mathbf{w}_{k})^{2} + \frac{1}{2} \sum_{k=1}^{K} \|\mathbf{w}_{k}\|_{2}^{2},$$

where the  $\sigma_k$  are known real-valued scalars. Let  $\boldsymbol{\sigma}=(\sigma_1,\cdots,\sigma_K)$ .

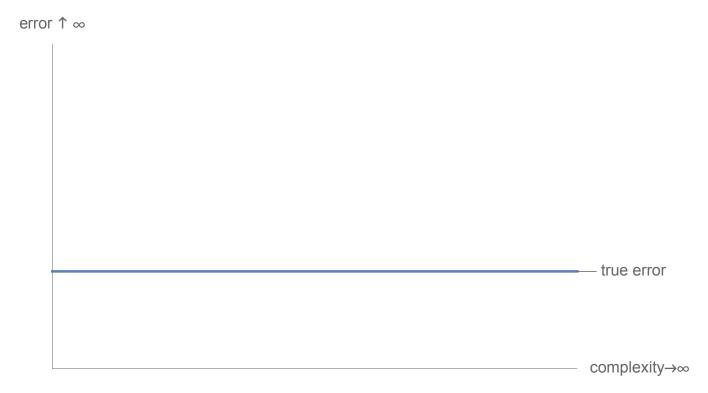
For the solution, let X be the  $N \times D$  matrix whose rows are the feature vectors  $\mathbf{x}_n$ .

- 1. Write down the normal equations for  $\mathbf{W}^*$ , the minimizer of the cost function. I.e., what is the first-order condition that  $\mathbf{W}^*$  has to fulfill in order to minimize  $\mathcal{L}(\mathbf{W})$ .
- 2. Is the minimum  $\mathbf{W}^{\star}$  unique? Assuming it is, write down an expression for this unique solution.
- 3. Write down a probabilistic model, so that the MAP solution for this model coincides with minimizing the above cost function. Note that this will involve specifying the likelihoods as well as a suitable prior (which will give you the regression term).

### 3 Bias Variance Trade-off (Due to Alex Smola)

Assume that you have two data sets that contain iid samples from the same distribution, call them  $S_1$  and  $S_2$ .  $S_1$  contains 5000 samples, whereas  $S_2$  contains 100000 samples. You randomly split each of the data sets into a training and a testing set, where eighty percent of the data is assigned to the training set. You then train and test on a family of increasing complexity.

In the figure below draw four curves, two that show the *training error* as a function of the model complexity (for  $S_1$  and  $S_2$ ) and two that show the *testing error* as a function of the model complexity (for  $S_1$  and  $S_2$ ). Label each of the 4 curves clearly. The constant curve labeled "true error" corresponds to the error due to the inherent noise in the samples and is drawn as a reference curve.



### 4 Exponential Families

Consider the Poisson distribution with parameter  $\lambda$ . It has a probability mass function given by  $p(i) = \frac{\lambda^i e^{-\lambda}}{i!}$ ,  $i = 0, 1, \cdots$ .

- (i) Write p(i) in the form of an exponential distribution  $p(i) = h(i)e^{\eta\phi(i)-A(\eta)}$ . Explicitly specify  $h, \eta, \phi$ , and  $A(\eta)$ .
- (ii) Compute  $\frac{dA(\eta)}{d\eta}$  and  $\frac{d^2A(\eta)}{d\eta^2}$ ? Is this the result you expected?

# 5 Multiple Choice Questions and Simple Problems

Mark the correct answer(s). More than one answer can be correct!

- In regression, "complex" models tend to
  - 1. overfit
  - 2. have large bias
  - 3. have large variance
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- We are given a data set  $S = \{(\mathbf{x}_n, y_n)\}$  for a binary classification task where  $\mathbf{x}_n$  in  $\mathbb{R}^D$ . We want to use a nearest-neighbor classifier. In which of the following situations do we have a reasonable chance of success with this approach? [Ignore the issue of complexity.]
  - 1. n is fixed,  $D \to \infty$
  - 2.  $n=D^2$ ,  $D\to\infty$
  - 3.  $n \to \infty$ ,  $D \ll \ln(n)$
  - 4.  $n \to \infty$ , D is fixed
- We add a regularization term because
  - 1. this sometimes renders the minimization problem of the cost function into a strictly convex/concave problem
  - 2. this tends to avoid overfitting
  - 3. this converts a regression problem into a classification problem
- The k-nearest neighbor classifier
  - 1. typically works the better the larger the dimension of the feature space
  - 2. can classify up to k classes
  - 3. typically works the worse the larger the dimension of the feature space
  - 4. can only be applied if the data can be linearly separated
  - 5. has a misclassification rate of at most two times the one of the Bayes classifier if we have lots of data
  - 6. has a misclassification rate that is two times better than the one of the Bayes classifier
- A real-valued scalar Gaussian distribution
  - 1. is a member of the exponential family with one scalar parameter
  - 2. is a member of the exponential family with two scalar parameters
  - 3. is not a member of the exponential family
- Which of the following statements is correct, where we assume that all the stated minima and maxima are in fact taken on in the domain of relevance.
  - 1.  $\max\{0, x\} = \max_{\alpha \in [0, 1]} \alpha x$
  - 2.  $\min\{0, x\} = \min_{\alpha \in [0, 1]} \alpha x$
  - 3. Let  $g(x) := \min_{y} f(x, y)$ . Then  $g(x) \le f(x, y)$
  - 4.  $\max_{x} g(x) \leq \max_{x} f(x, y)$
  - 5.  $\max_{x} \min_{y} f(x, y) \leq \min_{y} \max_{x} f(x, y)$
- Which of the following statements are correct?
  - 1. The training error is typically smaller than the test error.
  - 2. The SVM (support vector machine) formulation we discussed can be optimized using SGD.
  - 3. One iteration of SGD for ridge regression costs roughly  $\Theta(ND)$ , where N is the number of samples and D is the dimension.
  - 4. Logistic regression as formulated in class can be optimized using SGD.
- You have given the 2D data shown in Figure 1. You are allowed to add one component to your data (in addition to a constant component) and then must use a linear classifier. What component should you pick?
  - 1.  $x_1 + x_2$
  - 2.  $1/|x_1+x_2|$
  - 3.  $x_1 + 4x_2$

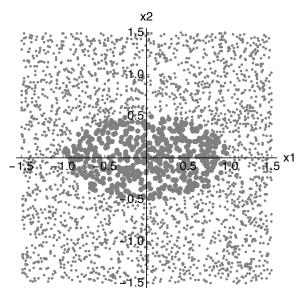


Figure 1: Some 2D data for classification. The two classes are indicated by different point sizes.

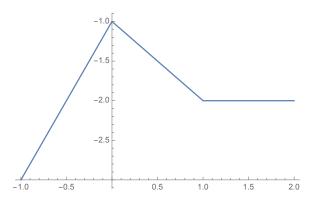


Figure 2: What is the subgradient of this function at x = 1?

- 4.  $4x_1 + x_2$
- 5.  $4x_1^2 + x_2^2$
- 6.  $x_1^2 x_2^2$
- 7.  $x_1^2 + 4x_2^2$
- The following functions are convex:
  - 1.  $f(x) := x^2, x \in \mathbb{R}$
  - 2.  $f(x) := x^3$ ,  $x \in [-1, 1]$
  - 3.  $f(x) := -x^3, x \in [-1, 0]$
  - 4.  $f(x) := e^{-x}, x \in \mathbb{R}$
  - 5.  $f(x) := e^{-x^2/2}, x \in \mathbb{R}$
  - 6.  $f(x) := \ln(1/x), x \in (0, \infty)$
  - 7.  $f(x) := g(h(x)), x \in \mathbb{R}$ , where g, h are convex and increasing over  $\mathbb{R}$
- Let  $f: \mathbb{R}^D \to \mathbb{R}$  be the function  $f(\mathbf{w}) := \exp(\mathbf{x}^\top \mathbf{w})$ , where  $\mathbf{x} \in \mathbb{R}^D$ . What is  $\nabla_{\mathbf{w}} f$ ?
- Which of the following scalars g is a subgradient for the function shown in Figure 2 at the point x = 1?
  - 1. g = -1
  - 2.  $g = -\frac{1}{2}$
  - 3. none exists
  - 4. g = 0