

Homework 2 - Introduction to Machine Learning for Engineers

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February 12, 2025

1 Naive Bayes Parameters

Problem Statement

The naïve Bayes approach assumes that the feature vectors are independent given the label, that is, for any given data point $\mathbf{x}_i = [x_{i1}, x_{i2}, \dots, x_{id}]^\top$ and its label y_i , we have

$$P(x_{i1}, x_{i2}, \dots, x_{id}, y_i) = P(y_i) \prod_{j=1}^d P(x_{ij} | y_i)$$

Suppose that we are given the data set $\mathbf{X} = [\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_N]^\top$, where $\mathbf{X} \in \mathbb{R}^{N \times 4}$, consisting of N data points and 4 features, and its label $\mathbf{y} = [y_1, y_2, \dots, y_N]^\top$. Assume that the first feature has 2, the second has 3, the third has 4, and the fourth has 5 possible values: Let $X_1 \in \{1, 2\}$, $X_2 \in \{1, 2, 3\}$, $X_3 \in \{1, 2, 3, 4\}$, and $X_4 \in \{1, 2, 3, 4, 5\}$. Also, there are 4 possible labels, i.e., $Y \in \{1, 2, 3, 4\} \forall i$.

This is what I have:

Feature 1: $X_1 \in \{1, 2\}$

Feature 2: $X_2 \in \{1, 2, 3\}$

Feature 3: $X_3 \in \{1, 2, 3, 4\}$

Feature 4: $X_4 \in \{1, 2, 3, 4, 5\}$

Label: $Y \in \{1, 2, 3, 4\}$

Naive's bayes classifier follow Bayesian rule:

$$P(Y|X) = \frac{P(X|Y)P(Y)}{P(X)}$$

Every data point has 4 features and 1 label:

$$x_i = [x_{i1}, x_{i2}, x_{i3}, x_{i4}] \quad \text{and} \quad y_i$$

- a. [5 points] Determine the number of free parameters θ and π that one has to estimate using the naïve Bayes framework. (We use θ for the probabilities of a feature given a label (e.g., $\theta_{3,2,4} = P(X_3 = 2 | Y = 4)$) and π for the class priors (e.g., $\pi_3 = P(Y = 3)$) similar to the lecture slides.)

Free parameters means those that we need to estimate.

class priors:

$$\pi_1 = P(Y = 1), \quad \pi_2 = P(Y = 2), \quad \pi_3 = P(Y = 3), \quad \pi_4 = P(Y = 4)$$

We need to estimate only three of these probabilities, since probability always sum up to one. So for class priors we have:

3 parameters to estimate

We need to estimate how likely each feature is given a label

Feature 1: it has 2 possible values, so we need to estimate 2 probabilities for each label

$$\theta_{1,1} = P(X_1 = 1 | Y = 1), \quad \theta_{1,2} = P(X_1 = 2 | Y = 1)$$

probability should sum up to one, so we only need to estimate one of the probabilities. So for feature 1 given 4 labels, we have:

4 parameters to estimate

Feature 2: it has 3 possible values, so we need to estimate 3 probabilities for each label

$$\theta_{2,1} = P(X_2 = 1 | Y = 1), \quad \theta_{2,2} = P(X_2 = 2 | Y = 1), \quad \theta_{2,3} = P(X_2 = 3 | Y = 1)$$

probability should sum up to one, so we only need to estimate two of the probabilities. So for feature 2 given 4 labels, we have:

$4 \times 2 = \mathbf{8}$ parameters to estimate

Feature 3: it has 4 possible values, so we need to estimate 4 probabilities for each label:

$$\theta_{3,1} = P(X_3 = 1 | Y = 1), \quad \theta_{3,2} = P(X_3 = 2 | Y = 1), \quad \theta_{3,3} = P(X_3 = 3 | Y = 1), \quad \theta_{3,4} = P(X_3 = 4 | Y = 1)$$

probability should sum up to one, so we only need to estimate three of the probabilities. So for feature 3 given 4 labels, we have:

$4 \times 3 = \mathbf{12}$ parameters to estimate

Feature 4: it has 5 possible values, so we need to estimate 5 probabilities for each label:

$$\theta_{4,1} = P(X_4 = 1 | Y = 1), \quad \theta_{4,2} = P(X_4 = 2 | Y = 1), \quad \theta_{4,3} = P(X_4 = 3 | Y = 1), \quad \theta_{4,4} = P(X_4 = 4 | Y = 1), \quad \theta_{4,5} = P(X_4 = 5 | Y = 1)$$

probability should sum up to one, so we only need to estimate four of the probabilities. So for feature 4 given 4 labels, we have:

$$4 \times 4 = \mathbf{16} \text{ parameters to estimate}$$

So the total number of parameters to estimate is:

$$3 + 4 + 8 + 12 + 16 = \mathbf{43} \text{ parameters to estimate}$$

- b. [4 points] If the features are not independent conditioned on the label, one has to estimate the entire joint distribution $P(X_1, X_2, X_3, X_4 | Y = i)$, for all $i \in \{1, 2, 3, 4\}$. Determine the number of free parameters that one has to estimate in such a scenario.
for each label, we have 4 features, so we need to estimate the entire joint distribution for each label. For each label, we have:
For each class $Y = i$:

Feature 1: $X_1 \in \{1, 2\}$

Feature 2: $X_2 \in \{1, 2, 3\}$

Feature 3: $X_3 \in \{1, 2, 3, 4\}$

Feature 4: $X_4 \in \{1, 2, 3, 4, 5\}$

So the total number of possible combinations for each label is:

$$2 \times 3 \times 4 \times 5 = 120$$

When we add the joint distribution for every label, it should sum up to one. So the total number of parameters to estimate is:

$$120 - 1 = \mathbf{119} \text{ parameters to estimate}$$

We have 4 labels, so the total number of parameters to estimate is:

$$4 \times 119 = \mathbf{476} \text{ parameters to estimate}$$

- c. [1 point] Based on the numbers of parameters you found in parts (a) and (b), explain one advantage of assuming conditional independence.

2 Naive Bayes in Practice

In this problem we will use the naïve Bayes algorithm to classify movie reviews as positive, neutral or negative. A simple approach involves maintaining a vocabulary of words that commonly occur in movie reviews and using the frequency of their occurrence in the three classes to classify movie reviews.

We are given the vocabulary $V = \{1 : \text{“incredible”}, 2 : \text{“plot”}, 3 : \text{“great”}, 4 : \text{“amazing”}, 5 : \text{“okay”}, 6 : \text{“decent”}, 7 : \text{“movie”}, 8 : \text{“no”}, 9 : \text{“acting”}, 10 : \text{“waste”}\}$. We will use V_i to represent the i th word in V . The training set provided includes four positive reviews:

- “great movie amazing”
- “incredible movie”
- “great acting amazing plot”
- “amazing acting amazing plot”

two neutral reviews:

- “okay movie”
- “decent no amazing acting”

and two negative reviews:

- “amazing waste”
- “no movie plot”

Recall that the naïve Bayes classifier is a generative classifier, where the probability of an input $\mathbf{x} = [x_1, x_2, \dots, x_n]^\top$ depends on its class y . In our case the input vector \mathbf{x} corresponding to each movie review has $n = 10$ equal to the number of words in the vocabulary V , where each entry x_i is equal to the number of times word V_i occurs in \mathbf{x} .

- a. [2 points] Calculate the naïve Bayes estimates of $\Pr(y = 1)$, $\Pr(y = 2)$ and $\Pr(y = 3)$ from the training data, where $y = 1$ corresponds to positive reviews, $y = 2$ to neutral reviews and $y = 3$ to negative reviews.
- b. [1 point] List the feature vector \mathbf{x} for each positive review in the training set.
- c. [2 points] In the naïve Bayes model, the likelihood of a sentence with feature vector \mathbf{x} given a class c is

$$\Pr(\mathbf{x} \mid y = c) = \prod_{k=1}^n (\theta_{c,k})^{x_k}$$

where $\theta_{c,k}$ is the weight of word k among all words of class c . Calculate the maximum likelihood estimate of $\theta_{1,4}$, $\theta_{1,7}$, $\theta_{2,4}$, $\theta_{2,7}$, $\theta_{3,4}$ and $\theta_{3,7}$.

- d. [3 points] Given a new review “amazing movie”, decide whether it is positive, neutral or negative, based on the naïve Bayes classifier, learned from the above data.
- e. [4 points] Use Laplacian smoothing with $\alpha = 1$ to decide whether the review “decent movie” is positive, neutral or negative. In one sentence, describe the problem we would encounter if we had not used Laplacian smoothing.

3 Logistic Regression in Practice

Given a training set $D = \{(\mathbf{x}^{(1)}, y^{(1)}), \dots, (\mathbf{x}^{(N)}, y^{(N)})\}$ where $(\mathbf{x}^{(i)} \in \mathbb{R}^d, y^{(i)} \in \{0, 1\})$ is the feature vector and the binary label for data point i , we want to find the parameters $\hat{\mathbf{w}}$ that maximize the likelihood for the training set, assuming a parametric model of the form

$$p(y = 1 \mid \mathbf{x}, \mathbf{w}) = \frac{1}{1 + \exp(-\mathbf{w}^T \mathbf{x})}.$$

The conditional log likelihood of the training set is

$$L(\mathbf{w}) = \sum_{i=1}^N [y^{(i)} \log p(y^{(i)} \mid \mathbf{x}^{(i)}, \mathbf{w}) + (1 - y^{(i)}) \log (1 - p(y^{(i)} \mid \mathbf{x}^{(i)}, \mathbf{w}))],$$

and the gradient is

$$\nabla L(\mathbf{w}) = \sum_{i=1}^N (y^{(i)} - p(y^{(i)} \mid \mathbf{x}^{(i)}, \mathbf{w})) \mathbf{x}^{(i)}. \quad (2)$$

a. [2 points] Is it possible to get a closed form for the parameters $\hat{\mathbf{w}}$ that maximize the conditional log likelihood? If it is possible, give the closed-form solution. If not, how would you compute $\hat{\mathbf{w}}$ in practice? Explain your method to find $\hat{\mathbf{w}}$ in a few sentences or provide a short pseudo-code.

For a binary logistic regression model, we predict $y = 1$ when $p(y = 1 \mid \mathbf{x}) \geq 0.5$. Assume that the decision boundary occurs when $P(y = 1 \mid \mathbf{x}, \mathbf{w}) = P(y = 0 \mid \mathbf{x}, \mathbf{w})$.

b. [4 points] Find the decision boundary, which is the set of \mathbf{x} satisfying $P(y = 1 \mid \mathbf{x}, \mathbf{w}) = P(y = 0 \mid \mathbf{x}, \mathbf{w})$.

c. [1 point] Is this model a linear classifier?

d. [2 points] Now, let us assume that in our application of this model, our tolerance of mis-classifying samples with the true label $y = 0$ is much lower than our tolerance to mis-classifying samples with true label $y = 1$ (i.e., we do not want to estimate a data point as belonging to class 0 if its true class is 1. However, we can tolerate the error of estimating a data point belong to class 1 if its true class is 0.) For example, in predicting the presence of a disease, we may be much more tolerant of mis-classifying a healthy person (true label $y = 0$) as having the disease ($y = 1$) compared to mis-classifying a person with the disease (true label $y = 1$) as not having it ($y = 0$). Now suppose you have trained a logistic regression model to obtain the model weights \mathbf{w} . How would you adjust the prediction with weights \mathbf{w} to decrease the mis-classification of samples with true label $y = 1$?

4 Solving Logistic Regression

The cross-entropy loss function for a logistic regression task on a dataset $\{(\mathbf{x}_1, y_1), (\mathbf{x}_2, y_2), \dots, (\mathbf{x}_n, y_n)\}$ can be written as

$$L(\mathbf{w}) = - \sum_{i=1}^n [y_i \log p(y_i | \mathbf{x}_i, \mathbf{w}) + (1 - y_i) \log(1 - p(y_i | \mathbf{x}_i, \mathbf{w}))],$$

In this question, we will show that finding the MLE (maximum likelihood estimator) is equivalent to minimizing the cross-entropy loss in Eq. (3). However, unlike linear regression, it is difficult to derive a closed-form solution for logistic regression.

- a. [3points] Starting from the definition of negative log-likelihood (Eq. (4) below), show that it is equal to the cross-entropy loss (Eq. (3)).

The negative log-likelihood is given by:

$$-\ell(\mathbf{w}) = -\log \left(\prod_{i=1}^n \left(\frac{1}{1 + \exp(-\mathbf{w}^\top \mathbf{x}_i)} \right)^{y_i} \left(1 - \frac{1}{1 + \exp(-\mathbf{w}^\top \mathbf{x}_i)} \right)^{1-y_i} \right). \quad (4)$$

Show that this is equivalent to the cross-entropy loss:

$$L(\mathbf{w}) = - \sum_{i=1}^n \left[y_i \log \left(\frac{1}{1 + \exp(-\mathbf{w}^\top \mathbf{x}_i)} \right) + (1 - y_i) \log \left(1 - \frac{1}{1 + \exp(-\mathbf{w}^\top \mathbf{x}_i)} \right) \right]. \quad (3)$$

- b. [4 points] Since it is difficult to directly optimize Eq. (3), gradient descent algorithms are usually used to find the optimum, as is discussed in the lecture. Show that the negative log-likelihood is a convex function. You may use the facts that the sum of convex functions is also convex, and that if f and g are both convex, twice differentiable and g is non-decreasing, then $g(f)$ is convex.

Hint: A function f is convex if for any $x_1, x_2 \in \text{Domain}(f)$,

$$f(tx_1 + (1-t)x_2) \leq tf(x_1) + (1-t)f(x_2), \quad \forall t \in [0, 1] \quad (5)$$

Also, a twice-differentiable function is convex if and only if the Hessian is positive semi-definite.

We now turn our attention to another algorithm known as iterative weighted least squares, which is based on the Newton-Raphson algorithm. Let $\mathbf{w}^{(k)}$ denote the parameter vector \mathbf{w} at the k th iteration. Then, the update rule of iterative weighted least-squares is as follows:

$$\mathbf{w}_{(k+1)} = \mathbf{w}_{(k)} - (\nabla^2 \mathcal{L}(\mathbf{w}_{(k)}))^{-1} \nabla L(\mathbf{w}_{(k)}),$$

where $\nabla L(\mathbf{w})$ and $\nabla^2 L(\mathbf{w})$ are the gradient and Hessian of the negative log-likelihood in Eq. (3), respectively. Consider the following notations:

- \mathbf{y} is an $N \times 1$ column vector with the i -th element being the label y_i ,
- \mathbf{X} is an $N \times d$ matrix with \mathbf{x}_i^\top being the i -th row,
- \mathbf{W}_k is an $N \times N$ diagonal matrix with the i -th diagonal element being $\frac{\exp(-\mathbf{w}_k^\top \mathbf{x}_i)}{(1 + \exp(-\mathbf{w}_k^\top \mathbf{x}_i))^2}$,
- \mathbf{p}_k is an $N \times 1$ column vector with the i -th element being $\frac{1}{1 + \exp(-\mathbf{w}_k^\top \mathbf{x}_i)}$,
- $\mathbf{z}_k = \mathbf{X}\mathbf{w}_k + \mathbf{W}_k^{-1}(\mathbf{y} - \mathbf{p}_k)$.

It can be shown that:

- $\nabla L(\mathbf{w}_k) = -\mathbf{X}^\top(\mathbf{y} - \mathbf{p}_k)$,
- $\nabla^2 L(\mathbf{w}_k) = \mathbf{X}^\top \mathbf{W}_k \mathbf{X}$.

- c [5 points] Using the above, prove that

$$\mathbf{w}_{k+1} = (\mathbf{X}^\top \mathbf{W}_k \mathbf{X})^{-1} \mathbf{X}^\top \mathbf{W}_k \mathbf{z}_k.$$

Note: The name of the algorithm comes from the observation that the algorithm at each step is solving the weighted least squares problem,

$$\mathbf{w}_{k+1} = \arg \min_{\mathbf{w}_k} (\mathbf{z}_k - \mathbf{X}\mathbf{w}_k)^\top \mathbf{W}_k (\mathbf{z}_k - \mathbf{X}\mathbf{w}_k).$$

5 SVMs: Hinge loss and mistake bounds

Suppose we build a predictive model $\hat{y} = \text{sgn}(\mathbf{w}^\top \mathbf{x})$, where $\text{sgn}(z) = 1$ if z is positive and -1 otherwise. Here we are dealing with binary predictions where each label is in $\{-1, 1\}$. The classification loss counts the number of mistakes (i.e., points where $y \neq \text{sgn}(\mathbf{w}^\top \mathbf{x})$) that we make with \mathbf{w} on a dataset. When we train the SVM model on N data points $\{(\mathbf{x}_1, y_1), (\mathbf{x}_2, y_2), \dots, (\mathbf{x}_N, y_N)\}$, the SVM loss function can be viewed as a relaxation to the classification loss.

The hinge loss on a data point (\mathbf{x}, y) is defined as:

$$\ell((\mathbf{x}, y), \mathbf{w}) = \max(0, 1 - y\mathbf{w}^\top \mathbf{x}),$$

where $\mathbf{x} \in \mathbb{R}^D$ and $y \in \{-1, 1\}$. The SVM attempts to minimize:

$$L(\mathbf{w}) = \frac{1}{N} \sum_{i=1}^N \ell((\mathbf{x}_i, y_i), \mathbf{w}) + \lambda \|\mathbf{w}\|_2^2,$$

where $\lambda > 0$ is the regularization parameter.

a. [4 points] Suppose that for some \mathbf{w} we have a correct prediction of y_i with \mathbf{x}_i , i.e., $y_i = \text{sgn}(\mathbf{w}^\top \mathbf{x}_i)$. What range of values can the hinge loss, $\ell((\mathbf{x}, y), \mathbf{w})$, take on this correctly classified example? Points that are classified correctly and which have non-zero hinge loss are referred to as margin mistakes. Also, what is the possible range of the hinge loss for an incorrectly classified example?

b. [5 points] Let $M(\mathbf{w})$ be the number of mistakes made when we use the weight vector \mathbf{w} to classify our dataset (i.e., the number of training data points for which $y_i \neq \text{sgn}(\mathbf{w}^\top \mathbf{x}_i)$). Note that \mathbf{w} can be an arbitrary weight vector (not necessarily the one that solves the SVM optimization problem). Show that:

$$\frac{1}{N} M(\mathbf{w}) \leq \frac{1}{N} \sum_{i=1}^N \max(0, 1 - y_i \mathbf{w}^\top \mathbf{x}_i). \quad (11)$$

In other words, the average hinge loss on our dataset is an upper bound on the average number of mistakes we make on our dataset. By minimizing the average hinge loss with \mathbf{w} , we can ensure that the SVM model makes few mistakes.

c. [3 points] Assume that the training data is separable. However, it turns out that the minimizer \mathbf{w}^* of the objective in Eq. (10) mis-classifies some training samples. How should we adjust λ (i.e., make it larger or smaller) to make the SVM work properly on training data? Why? Explain in 1-2 sentences.

6 Support Vector Machine - Slack Variables & Duality Intuition

In hard SVM, our aim is to maximize the margin when samples are linearly separable. The margin is defined as the distance of the closest point to the classification boundary. In lecture, we showed that this is equivalent to solving the following hard-SVM optimization problem:

$$\min_{w,b} \frac{1}{2} \|w\|_2^2 \quad \text{s.t.} \quad y_i(w^\top x_i + b) \geq 1, \quad \forall i \in [n] \quad (12)$$

where $w, x_i \in \mathbb{R}^d$, $y_i \in \{-1, +1\}$, and n denotes the number of data points. Here, each (x_i, y_i) pair denotes a training data point.

- a.** [1 point] Explain in one or two sentences why the hard-SVM formulation given above may fail when applied to real-world datasets. To address this issue, we use the soft-SVM formulation by introducing slack variables ξ_i for each data point $i = 1, 2, \dots, n$:

$$\min_{w,b,\xi_1,\dots,\xi_n} \left(\frac{1}{2} \|w\|_2^2 + C \sum_{i=1}^n \xi_i \right)$$

$$\text{s.t.} \quad \forall i, \quad y_i(w^\top x_i + b) \geq 1 - \xi_i, \quad \xi_i \geq 0.$$

- b.** [4 points] Discuss how the introduction of slack variables, ξ_i , resolves the impracticalities of the hard-SVM formulation that you identified in part (a). Why do we use “ $\geq 1 - \xi_i$ ” instead of “ ≥ 1 ” in the constraints and why is it necessary to include “ $C \sum_i \xi_i$ ” in the minimization objective? Your answer should be no more than three or four sentences long.
- c.** [4 points] Further, as we see in the class, the dual formulation of the soft-SVM problem can be preferred.
- d.** [3 points] Find the number of variables to be optimized in the primal and dual forms of SVM. Then, suppose that we decide to use some other kernels (e.g., radial basis kernel, polynomial kernel) instead of the linear kernel $k(\mathbf{x}_i, \mathbf{x}_j) = \mathbf{x}_i^\top \mathbf{x}_j$. Does this change the number of dual variables we need to use?
- e.** [1 point] Why might you prefer to solve the SVM in the dual formulation as opposed to the primal formulation?

7 Implementing SVMs

In this problem, you will experiment with SVMs on a real-world dataset. You will implement a linear SVM (i.e., an SVM using the original features). You should implement it from scratch without using any libraries like scikit-learn, although you can use existing packages to solve quadratic programs. If you are not sure if the library you want to use is allowed, please ask the TAs or post a question on Piazza. We provide you with a guided Python Notebook (hw2_svm_imp.ipynb) on which you will find parts you need to fill in. You can use this notebook on JupyterLab / JupyterNotebook / GoogleColab.

Please append your notebook file as a PDF at the end of your submission so that your code and cell outputs can be seen on the same submission PDF. If you use JupyterLab / JupyterNotebook / GoogleColab, you can save the notebook as PDF from File→Print. You can also use online ‘ipynb to pdf’ file converters. Please do not forget to append your notebook as a PDF to the end of your submission PDF for us to grade this question.

Dataset: We have provided the Heart Disease dataset from UCI’s machine learning data repository. The provided binary classification dataset has 13 input features, and 303 samples.

- By running the first cell of the given notebook, you download the ucimlrepo package. It will help us download the dataset. For this step, you should have an internet connection.
- In the second cell, we provide all required libraries to complete the homework.
- In the third cell, we download the dataset (again, you need an internet connection), one-hot encode some categorical features, and split it into train/test parts using a set seed.

You should run those three cells without any modification. After running the first three cells, you should have `X_train` (shape of (216, 22)), `X_test` (shape of (83, 22)), `y_train` (shape of (216,)), and `y_test` (shape of (83,)) variables.

1. Let $x_k^{(1)}, \dots, x_k^{(N)}$ be the values of feature k for all training set (X_{train}) points where N is the number of samples in the training set. Preprocess the training (X_{train}) and test data (X_{test}) by:

- (a) Computing the mean $\bar{x}_k = \frac{1}{N} \sum_{i=1}^N x_k^{(i)}$ of each feature and subtracting it from all values of this feature.
- (b) Dividing each feature by its standard deviation, defined as

$$s_k = \sqrt{\frac{1}{N} \sum_{i=1}^N (x_k^{(i)} - \bar{x}_k)^2}$$

for feature k . Here, \bar{x}_k is the sample mean of this feature.

Save the normalized data samples as variables named `X_train_normalized` for the training set and `X_test_normalized` for the test set. Note that the shape of these variables should be the same as the shapes of `X_train` and `X_test`, respectively.

This type of preprocessing is useful for SVMs, as SVMs attempt to maximize the distance between the separating hyperplane and the support vectors. If one feature (i.e., one dimension in this space) has very large values, it will dominate the other features when calculating this distance. Rescaling the features will ensure that they all have the same influence on the distance metric.

Note that the mean and standard deviation should be estimated from the training data (`X_train`) and then applied to both datasets. This is because using the statistics from the training data ensures that the model is not biased by the test data, which should remain unseen during training.

Report the mean and the standard deviation of the first and the last features computed on the training data (`X_train`):

- Mean of the first feature: \bar{x}_1
- Standard deviation of the first feature: s_1
- Mean of the last feature: \bar{x}_{22}
- Standard deviation of the last feature: s_{22}