Problem Set #3: Deep Learning & Unsupervised Learning

Problem 1 A simple neural network

Let $X = \{x^{(1)}, x^{(2)}, \dots, x^{(m)}\}$ be dataset of m examples with 2 features. That is, $x^{(i)} \in \mathbb{R}^2$. Samples are classified into 2 categorie with labels $y \in \{0, 1\}$, as shown in Figure 1. Want to perform binary classification using a simple neural networks with the architecture shown in Figure 2.

Two features x_1 and x_2 , the three neurons in the hidden layer h_1 , h_2 , h_3 , and the output neuron as o. Weight from x_i to h_j be $w_{i,j}^{[1]}$ for i = 1, 2 and j = 1, 2, 3, and weight from h_j to o be $w_j^{[2]}$. Finally, denote intercept weight for h_j as $w_{0,j}^{[1]}$ and the intercept weight for o as $w_0^{[2]}$. Use average squared loss instead of the usual negative log-likelihood:

$$l = \frac{1}{m} \sum_{i=1}^{m} (o^{(i)} - y^{(i)})^{2}.$$

(a) Suppose we use sigmoid function as activation function for h_1 , h_2 , h_3 , and o. We have

$$h_1 = g(w_1^{[1]}x), \quad h_2 = g(w_2^{[1]}x), \quad h_3 = g(w_3^{[1]}x), \quad o = g(w_3^{[2]}h).$$

Hence,

$$\frac{\partial l}{\partial w_{1,2}^{[1]}} = \frac{1}{m} \sum_{i=1}^{m} 2(o^{(i)} - y^{(i)}) o^{(i)} (1 - o^{(i)}) w_2^{[2]} h_2^{(i)} (1 - h_2^{(i)}) x_1^{(i)},$$

where $h_2^{(i)} = g(w_{0,2}^{[1]} + w_{1,2}^{[1]} x_1^{(i)} + w_{2,2}^{[1]} x_2^{(i)})$ and g is the sigmoid function. Therefore, the gradient descent update to $w_{1,2}^{[1]}$, assuming learning rate α is

$$w_{1,2}^{[1]} := w_{1,2}^{[1]} - \frac{2\alpha}{m} \sum_{i=1}^{m} (o^{(i)} - y^{(i)}) o^{(i)} (1 - o^{(i)}) w_2^{[2]} h_2^{(i)} (1 - h_2^{(i)}) x_1^{(i)}$$

where $h_2^{(i)} = g(w_{0,2}^{[1]} + w_{1,2}^{[1]}x_1^{(i)} + w_{2,2}^{[1]}x_2^{(i)}).$

(b) Now, suppose the activation function for h_1 , h_2 , h_3 , and o is the step function f(x), defined as

$$f(x) = \begin{cases} 1, & (x \ge 0), \\ 0, & (x < 0). \end{cases}$$

Is it possible to have a set of weights that allow the neural network to classify this dataset with 100% accuracy? If so, provide a set of weights by completing optimal_step_weights wihin src/p01_nn.py and explain your reasoning for those weights. If not, please explain the reasoning.

There is a set of weights that allow the neural network to classify this dataset with 100% accuracy. For the step function activation, we have

$$h_{1} = f(w_{1}^{[1]}x) = f(w_{0,1}^{[1]} + w_{1,1}^{[1]}x_{1} + w_{2,1}^{[1]}x_{2})$$

$$h_{2} = f(w_{2}^{[1]}x) = f(w_{0,2}^{[1]} + w_{1,2}^{[1]}x_{1} + w_{2,3}^{[1]}x_{2})$$

$$h_{3} = f(w_{3}^{[1]}x) = f(w_{0,3}^{[1]} + w_{1,3}^{[1]}x_{1} + w_{2,3}^{[1]}x_{2})$$

$$o = f(w^{[2]}h) = f(w_{0}^{[2]} + w_{1}^{[2]}h_{1} + w_{2}^{[2]}h_{2} + w_{3}^{[2]}h_{3}).$$

Notice from Figure 1 that the label $y^{(i)} = 0$ if and only if $x^{(i)}$ satisfies

$$\begin{cases} x_2^{(i)} > 0.5, \\ x_1^{(i)} > 0.5, \\ x_1^{(i)} + x_2^{(i)} < 4. \end{cases}$$

Now, let

$$w_1^{[1]} = \begin{bmatrix} 0.5 \\ 0 \\ -1 \end{bmatrix}, \quad w_2^{[1]} = \begin{bmatrix} 0.5 \\ -1 \\ 0 \end{bmatrix}, \quad w_3^{[1]} = \begin{bmatrix} -4 \\ 1 \\ 1 \end{bmatrix}, \quad w_1^{[1]} = \begin{bmatrix} -0.5 \\ 1 \\ 1 \\ 1 \end{bmatrix}.$$

Under this set of weights, if all inequalities are satisfied, then $h_1 = h_2 = h_3 = 0$ and $w^{[2]}h = -0.5$. Otherwise, $h_1 + h_2 + h_3 \ge 1$ and $w^{[2]}h \ge -0.5$. Hence, This set of weights will capture all the conditions and allow the nerual network to classify this dataset with 100% accuracy.

(c) Let the activation function for h_1 , h_2 , h_3 , and o is the linear function f(x) = x, and the activation function for o be the same step function as before. Is it possible to have a set of weights that allow the neural network to classify this dataset with 100% accuracy? If so, provide a set of weights by completing optimal_linear_weights wihin src/p01_nn.py and explain your reasoning for those weights. If not, please explain the reasoning.

Problem 2 KL divergence and maximum likelihood

Kullback-Leibler (KL) divergence is a measure of how much one probability distribution is different from a second one. The KL divergence between two discrete-valued distribution P(X), Q(X) over the outcome space \mathcal{X} is defined as follows:

$$D_{\mathrm{KL}}(P \parallel Q) = \sum_{x \in \mathcal{X}} P(x) \log \frac{P(x)}{Q(x)}.$$

Assuem P(x) > 0 for all x. (One other standard thing to do is adopt the convention that $0 \log 0 = 0$.) Sometimes, we also write the KL divergence more explicitly as $D_{\text{KL}}(P \parallel Q) = D_{\text{KL}}(P(X) \parallel Q(X))$.

Background on Information Theory

The *entropy* of a probability distribution P(X), defined as

$$H(P) = -\sum_{x \in \mathcal{X}} P(x) \log P(x).$$

measures how dispersed a probability distribution is. Notably, $\mathcal{N}(\mu, \sigma^2)$ has the highest entropy among all possible continuous distribution that has mean μ and variance σ^2 . The entropy H(P) is the best possible long term average bits per message (optimal) that can be achieved under probability distribution P(X).

The *cross entropy* is defined as

$$H(P,Q) = -\sum_{x \in \mathcal{X}} P(x) \log Q(x).$$

The cross entropy H(P,Q) is the long term average bits per message (suboptimal) that results under a distribution P(X), by reusing an encoding scheme designed to be optimal for a scenario with probability distribution Q(X).

Notice that

$$D_{\mathrm{KL}}(P \parallel Q) = \sum_{x \in \mathcal{X}} P(x) \log P(x) - \sum_{x \in \mathcal{X}} P(x) \log Q(x) = H(P, Q) - H(P).$$

If H(P,Q)=0, then it necessarily means P=Q. In ML, it is common task to find distribution Q that is close to another distribution P. To achieve this, we optimize $D_{\mathrm{KL}}(P\parallel Q)$. Later we will see that Maximum Likelihood Estimation turns out to be equivalent minimizing KL divergence between the training data and the model.

(a) **Nonnegativity.** Prove that

$$D_{\mathrm{KL}}(P \parallel Q) \ge 0$$

and $D_{KL}(P \parallel Q) = 0$ if an only if P = Q.

Hint: Use Jensen's inequality.

Proof. By definition,

$$D_{\mathrm{KL}}(P \parallel Q) = \sum_{x \in \mathcal{X}} P(x) \log \frac{P(x)}{Q(x)} = -\sum_{x \in \mathcal{X}} P(x) \log \frac{Q(x)}{P(x)}.$$

Since $-\log x$ is strictly convex, by Jensen's inequality, we have

$$D_{\mathrm{KL}}(P \parallel Q) = -\sum_{x \in \mathcal{X}} P(x) \log \frac{Q(x)}{P(x)} \ge -\log \sum_{x \in \mathcal{X}} P(x) \frac{Q(x)}{P(x)} = 0.$$

When the equality holds,

$$\log \frac{Q(x)}{P(x)} = 0$$

with probability 1. That is, Q = P with probability 1. This completes the proof.

(b) Chain rule for KL divergence. The KL divergence between 2 conditional distributions $P(X \mid Y)$, $Q(X \mid Y)$ is defined as follows:

$$D_{\mathrm{KL}}(P(X\mid Y)\parallel Q(X\mid Y)) = \sum_{y} P(y) \left(\sum_{x} P(x\mid y) \log \frac{P(x\mid y)}{Q(x\mid y)}\right).$$

This can be thought of as the expected KL divergence between the corresponding conditional distributions on x. That is, between $P(X \mid Y = y)$ and $Q(X \mid Y = y)$, where the expectation is taken over the random y.

Prove the following chain rule for KL divergence:

$$D_{\mathrm{KL}}(P(X,Y) \parallel Q(X,Y)) = D_{\mathrm{KL}}(P(X) \parallel Q(X)) + D_{\mathrm{KL}}(P(Y \mid X) \parallel Q(Y \mid X)).$$

Proof.

LHS =
$$\sum_{x} \sum_{y} P(x, y) \log \frac{P(x, y)}{Q(x, y)}$$

= $\sum_{x} \sum_{y} P(y \mid x) P(x) \left[\log \frac{P(y \mid x)}{Q(y \mid x)} + \log \frac{P(x)}{Q(x)} \right]$
= $\sum_{x} \sum_{y} P(y \mid x) P(x) \log \frac{P(y \mid x)}{Q(y \mid x)} + \sum_{x} P(x) \log \frac{P(x)}{Q(x)} \sum_{y} P(y \mid x)$
= $\sum_{x} \sum_{y} P(y \mid x) P(x) \log \frac{P(y \mid x)}{Q(y \mid x)} + \sum_{x} P(x) \log \frac{P(x)}{Q(x)}$
= $D_{KL}(P(X) \parallel Q(X)) + D_{KL}(P(Y \mid X) \parallel Q(Y \mid X))$
= RHS.

(c) **KL** and maximum likelihood. Consider density estimation problem and suppose we are given training set $\{x^{(i)}\}_{i=1}^m$. Let the empirical distribution be $\hat{P}(x) = \frac{1}{m} \sum_{i=1}^m \mathbb{I}\{x^{(i)} = x\}$. (\hat{P} is just the uniform distribution over the training set; i.e., sampling from the empirical distribution is the same as picking a random example from the training set.)

Suppose we have a family of distributions P_{θ} parametrized by θ . Prove that finding the maximum likelihood estimates for the parameter θ is equivalent to finding P_{θ} with minimal KL divergence from \hat{P} . That is, prove that

$$\underset{\theta}{\operatorname{argmin}} D_{\mathrm{KL}}(\hat{P} \parallel P_{\theta}) = \underset{\theta}{\operatorname{argmax}} \sum_{i=1}^{m} \log P_{\theta}(x^{(i)}).$$

Proof. Notice that \hat{P} is the uniform distribution over the training set, thus $\hat{P}(x^{(i)}) = \frac{1}{m}$ for i = 1, ..., m. It follows that

$$D_{\mathrm{KL}}(\hat{P} \parallel P_{\theta}) = \sum_{x} \hat{P}(x) \log \frac{\hat{P}(x)}{P_{\theta}(x)} = -\log m - \frac{1}{m} \sum_{i=1}^{m} \log P_{\theta}(x^{(i)}).$$

Hence,

$$\underset{\theta}{\operatorname{argmin}} D_{\mathrm{KL}}(\hat{P} \parallel P_{\theta}) = \underset{\theta}{\operatorname{argmax}} \sum_{i=1}^{m} \log P_{\theta}(x^{(i)}),$$

as desired. \Box

Remark: Consider the relationship between parts (b-c) and multi-variate Bernoulli Naive bayes parameter estimation. In Naive Bayes model we assumed P_{θ} is the following form: $P_{\theta}(x,y) = p(y) \prod_{i=1}^{n} p(x_i \mid y)$. By the chain rule for KL divergence, we therefore have

$$D_{\mathrm{KL}}(\hat{P} \parallel P_{\theta}) = D_{\mathrm{KL}}(\hat{P}(y) \parallel p(y)) + \sum_{i=1}^{n} D_{\mathrm{KL}}(\hat{P}(x_i \mid y) \parallel p(x_i \mid y)).$$

This shows that finding the maximum likelihood/minimum KL divergence estimates of the parameters decomposes into 2n + 1 independent optimization problems: One for the class priors p(y), and one for each conditional distributions $p(x_i \mid y)$ for each feature x_i given each of the two possible labels for y. Specifically, finding the maximum likelihood estimates for each of these problems individually results in also maximizing the likelihood of the joint distribution. This similarly applies to bayesian networks.

Problem 3 KL divergence, Fisher Information, and the Natural Gradient

KL divergence between the two distributions is an asymetric measure of how different two distributions are. Consider two distributions over the same space given by densities p(x), q(x). The KL divergence between two continuous distributions is defined as

$$D_{\mathrm{KL}}(p \parallel q) = \int p(x) \log \frac{p(x)}{q(x)} dx$$
$$= \mathbb{E}_{x \sim p(x)}[\log p(x)] - \mathbb{E}_{x \sim p(x)}[\log q(x)].$$

A nice property of KL divergence is that it is invariant to parametrization. This means, KL divergence evaluates to the same value no matter how we parametrize the distribution P and Q. For example, if P and Q are in exponential family, the KL divergence between them is the same whether we are using natural parameters, natural parameters, or canonical parameters, or any arbitrary parametrization.

Now consider the problem of fitting model parameters using gradient descent. While KL divergence is invariant to parametrization, the gradient w.r.t the model parameters gradient is *invariant to parametrization*. We need to use *naturla gradient*. This will make the optimization process invariant to the parametrization.

We will construct and derive the natural gradient update rule. Along the way, we will introduce *score function* and *Fisher Information*. Finally, we will see how this new natural gradient based optimization is actually equivalent to Newton's method for Generalized Linear Models.

Let the distribution of a random variable Y parametrized by $\theta \in \mathbb{R}^n$ be $p(y;\theta)$.

(a) **Score function.** The score function with $p(y;\theta)$ is defined as $\nabla_{\theta} \log p(y;\theta)$, which signifies the sensitivity of the likelihood function with respect to the parameters. Show that the expected value of the score is 0.

Proof. The expected value of the score

$$\mathbb{E}_{y \sim p(y;\theta)}[[\nabla_{\theta'} \log p(y;\theta)]_{\theta'=\theta}] = \int p(y;\theta)[\nabla_{\theta'} \log p(y;\theta)]_{\theta'=\theta} dy$$

$$= \int p(y;\theta) \frac{1}{p(y;\theta)} [\nabla_{\theta'} p(y;\theta)]_{\theta'=\theta} dy$$

$$= \left[\nabla_{\theta'} \int p(y;\theta) dy\right]_{\theta'=\theta}$$

$$= 0$$

(b) **Fisher information.** Fisher information is defined as the covariance matrix of the score function,

$$\mathcal{I}(\theta) = \operatorname{Cov}_{y \sim p(y;\theta)} \left[\nabla_{\theta'} \log p(y;\theta') \right]_{\theta' = \theta}.$$

Intuitively, the Fisher information represents the amount of information that a random variable Y carries about a parameter θ of interest. Show that the Fisher information can be equivalently given by

$$\mathcal{I}(\theta) = \mathbb{E}_{y \sim p(y;\theta)} \left[\nabla_{\theta'} \log p(y;\theta') \nabla_{\theta'} \log p(y;\theta')^T \right]_{\theta' = \theta}.$$

Note that the fisher information is a function of the parameter. The parameter is both a) the parameter value at which the score function is evaluated, and b) the parameter of the distribution with respect to which the expectation and variance is calculated.

Proof.