

The perceptron algorithm consists of one of the simplest types of neural network architectures. It is a single-layer neural network with a single neuron. The neuron is a linear combination of the input variables, with a bias term, and then passed through an activation function. The activation function is used to determine the output of the neuron. The algorithm itself is a supervised learning algorithm, meaning that it requires labeled training data to train the model. The algorithm is iterative, hence it'll continue to update the weights until the model converges. The learning rate is a hyperparameter that controls the step size of the weight updates. The activation function is a hyperparameter that determines the output of the neuron. The learning rule is as follows:

$$w^{new} \leftarrow w^{old} + \eta \cdot (z - \hat{z}(x)) \cdot x, \quad \hat{z}(x) = f(net(x)), \quad net(x) = \sum_{i=1}^n w_i \cdot x^{(i)} + w_0$$

where w_i is the weight for the i th input variable, η is the learning rate, z is the true label, \hat{z} is the predicted label, and x_i is the i th input variable. The learning rule is applied for each training example, and an epoch is a single pass through the entire training set.

1. Considering the following linearly separable training data:

	y_1	y_2	y_3	z
x_1	0	0	0	-1
x_2	0	2	1	1
x_3	1	1	1	1
x_4	1	-1	0	-1

Given the perceptron learning algorithm with a learning rate $\eta = 1$, sign activation and all weights initialized to one (including the bias):

- Considering y_1 and y_2 , apply the algorithm until convergence. Draw the separation hyper-plane.
- Considering all input variables, apply one epoch of the algorithm. Do weights change for an additional epoch?
- Identify the perceptron output for $x_{new} = [0 \ 0 \ 1]^T$.
- What happens if we replace the sign function with the step function? Specifically, how would you change η to ensure the same results?

(a) As per the question statement, we're working with $\eta = 1$ and sign activation:

$$\hat{z} = f(net) = \begin{cases} 1 & net \geq 0 \\ -1 & net < 0 \end{cases}$$

Moreover, we're considering weights (and the bias, w_0) initialized at one. Performing epochs following the $\{x_1, \dots, x_4\}$ order, we get the following weight updates:

$$\begin{aligned} \hat{z}(x_1) &= f(net(x_1)) = f(1 + 1 \cdot 0 + 1 \cdot 0) = f(1) = 1 \\ \hat{z}(x_2) &= f(net(x_2)) = f(1 + 1 \cdot 0 + 1 \cdot 2) = f(3) = 1 \\ \hat{z}(x_3) &= f(net(x_3)) = f(1 + 1 \cdot 1 + 1 \cdot 1) = f(3) = 1 \\ \hat{z}(x_4) &= f(net(x_4)) = f(1 + 1 \cdot 1 + 1 \cdot -1) = f(1) = 1 \end{aligned}$$

$$\begin{aligned}
x_1 : \quad w &\leftarrow \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} + 1 \cdot (-1 - 1) \cdot [1 \quad 0 \quad 0] = \begin{bmatrix} -1 \\ 1 \\ 1 \end{bmatrix} \\
x_2 : \quad w &\leftarrow \begin{bmatrix} -1 \\ 1 \\ 1 \end{bmatrix} + 1 \cdot (1 - 1) \cdot [1 \quad 0 \quad 2] = \begin{bmatrix} -1 \\ 1 \\ 1 \end{bmatrix} \\
x_3 : \quad w &\leftarrow \begin{bmatrix} -1 \\ 1 \\ 1 \end{bmatrix} + 1 \cdot (1 - 1) \cdot [1 \quad 1 \quad 1] = \begin{bmatrix} -1 \\ 1 \\ 1 \end{bmatrix} \\
x_4 : \quad w &\leftarrow \begin{bmatrix} -1 \\ 1 \\ 1 \end{bmatrix} + 1 \cdot (-1 - 1) \cdot [1 \quad 1 \quad -1] = \begin{bmatrix} -1 \\ 1 \\ 1 \end{bmatrix}
\end{aligned}$$

After entering a new epoch, we'd update the weights again, this time for x_1 ; since such an update wouldn't lead to an actual update on the weight matrix, that'd make a full pass on the training set without changes, and, as such, the algorithm would converge. Therefore, the regression hyperplane is given by:

$$-1 + x_1 + x_2 = 0 \leftrightarrow x_2 = 1 - x_1$$

(b) Here, we apply the same algorithm as before, but now considering all input variables.

$$\begin{aligned}
\hat{z}(x_1) &= f(\text{net}(x_1)) = f(1 + 1 \cdot 0 + 1 \cdot 0 + 1 \cdot 0) = f(1) = 1 \\
\hat{z}(x_2) &= f(\text{net}(x_2)) = f(1 + 1 \cdot 0 + 1 \cdot 2 + 1 \cdot 1) = f(4) = 1 \\
\hat{z}(x_3) &= f(\text{net}(x_3)) = f(1 + 1 \cdot 1 + 1 \cdot 1 + 1 \cdot 1) = f(4) = 1 \\
\hat{z}(x_4) &= f(\text{net}(x_4)) = f(1 + 1 \cdot 1 + 1 \cdot -1 + 1 \cdot 0) = f(2) = 1
\end{aligned}$$

$$\begin{aligned}
x_1 : \quad w &\leftarrow \begin{bmatrix} 1 \\ 1 \\ 1 \\ 1 \end{bmatrix} + 1 \cdot (-1 - 1) \cdot [1 \quad 0 \quad 0 \quad 0] = \begin{bmatrix} -1 \\ 1 \\ 1 \\ 1 \end{bmatrix} \\
x_2 : \quad w &\leftarrow \begin{bmatrix} -1 \\ 1 \\ 1 \\ 1 \end{bmatrix} + 1 \cdot (1 - 1) \cdot [1 \quad 0 \quad 2 \quad 1] = \begin{bmatrix} -1 \\ 1 \\ 1 \\ 1 \end{bmatrix} \\
x_3 : \quad w &\leftarrow \begin{bmatrix} -1 \\ 1 \\ 1 \\ 1 \end{bmatrix} + 1 \cdot (1 - 1) \cdot [1 \quad 1 \quad 1 \quad 1] = \begin{bmatrix} -1 \\ 1 \\ 1 \\ 1 \end{bmatrix} \\
x_4 : \quad w &\leftarrow \begin{bmatrix} -1 \\ 1 \\ 1 \\ 1 \end{bmatrix} + 1 \cdot (-1 - 1) \cdot [1 \quad 1 \quad -1 \quad 0] = \begin{bmatrix} -1 \\ 1 \\ 1 \\ 1 \end{bmatrix}
\end{aligned}$$

Once again, we'd enter a new epoch and update the weights considering the first sample, but since such an update wouldn't lead to an actual update on the weight matrix, that'd make a full pass on the training set without changes, and, as such, the algorithm would converge. **An additional epoch wouldn't, therefore, change the weights.**

- (c) As we've mentioned before, the perceptron's output is given by the activation function - here, the sign of the net input. Therefore, the perceptron's output is given by:

$$\hat{z} = f(net) = \begin{cases} 1 & net \geq 0 \\ -1 & net < 0 \end{cases}$$

Here, considering the weights computed in the previous question, we'll have the following:

$$net(x_{new}) = -1 + 1 \cdot 0 + 1 \cdot 0 + 1 \cdot 1 = 0$$

As we know, $f(0) = 1$, and, as such, the perceptron would classify the new sample as belonging to the binary class 1.

- (d) As we know, the step function is rather similar to the sign function, defined as follows:

$$f(net) = \begin{cases} 1 & net \geq 0 \\ 0 & net < 0 \end{cases}$$

Let's consider the following learning rules (the first one considers the activation function as the sign function, and the second one considers the activation function as the step function):

$$\text{Sign function: } w \leftarrow w + \eta \cdot (z - \text{sign}(net)) \cdot x$$

$$\text{Step function: } w \leftarrow w + \eta \cdot (z - \text{step}(net)) \cdot x$$

We can note, of course, that the only differing term between both rules is $(z - \hat{z})$; therefore, to make it so that both rules are equivalent, we'd have to make sure that the step function's output is correctly balanced with the sign function's output, utilizing a scalar factor (be it τ) for such purpose, effectively altering the learning rate to be $\eta_{sign} = \tau \cdot \eta_{step}$. Let's try to find τ :

$$z - \text{sign}(net) = \tau(z - \text{step}(net))$$

We know that $z \in \{-1, 1\}$, and, as such, we can write each side of the equation's interval as follows:

$$[-1, 1] - [-1, 1] = \tau([-1, 1] - [0, 1]) \leftrightarrow [-2, 2] = \tau[-1, 1]$$

As such, if we want an equal learning rule utilizing both step and sign functions, we'd have to set $\tau = 2$, and, as such, get the following learning rule:

$$w \leftarrow w + 2\eta_{sign} \cdot (z - \text{step}(net)) \cdot x \quad \leftrightarrow \quad w \leftarrow w + \eta_{sign} \cdot (z - \text{sign}(net)) \cdot x$$

2. Let us consider the following activation function and training set:

$$\hat{z}(x, w) = \frac{1}{1 + e^{-2wx}}$$

	y_1	y_2	z
x_1	1	1	1
x_2	2	1	1
x_3	1	3	0
x_4	3	3	0

Consider also the half sum of squared errors as the loss function:

$$E(w) = 1/2 \sum_{i=1}^N (z_i - \hat{z}(x_i, w))^2$$

- Determine the gradient descent learning rule for this unit.
- Compute the first gradient descent update, assuming an initialization of all ones.
- Compute the first stochastic gradient descent update assuming an initialization of all ones.

As we should know by now, (activation) functions of the type $\frac{1}{1+e^{-k}}$ are called sigmoid functions - functions that are bounded by 0 and 1, with $\sigma(0) = 0.5$. $\sigma(x)$ is plotted below:

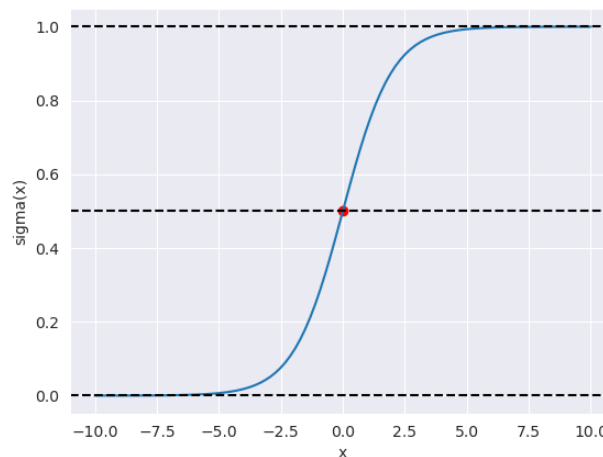


Figure 1: Sigmoid function

Sigmoid functions are widely used in Machine Learning (and, in particular, in perceptrons/neural network nodes) since they're able to map any real number to the interval $[0, 1]$, and, as such, they're able to represent probabilities fairly well. The sigmoid's derivative is rather famous (and must be known for the exam), given by:

$$\sigma'(x) = \sigma(x) \cdot (1 - \sigma(x)) \cdot x'$$

Gradient descent is essentially a linear regression model where, once again, we try to minimize a given error function - if in the case of the previous perceptron learning rule this error function was simply the difference between the desired output and the perceptron's output, in the case of gradient descent we have a more complex error function, such as the half sum of squared errors. We utilize this more complex learning rule when, for example, we're in the presence of non-linearly separable data, and, as such, we can't use the perceptron learning rule to find a separating hyperplane. Here, the learning rule is written as follows (considering E as the error function):

$$w \leftarrow w - \eta \cdot \frac{\partial E}{\partial w}$$

Note that here, since we're working with gradient **descent**, we're now subtracting the gradient of the error function from the weights, instead of adding it - we're trying to minimize the error function, after all.

(a) Note that the activation function, here, is a sigmoid that not only depends on x but also w :

$$\hat{z}(x, w) = \sigma(2wx) = \frac{1}{1 + e^{-2wx}}$$

Let's try to derivate the error function with respect to w :

$$\begin{aligned} \frac{\partial E}{\partial w} &= \frac{\partial}{\partial w} \frac{1}{2} \sum_{i=1}^N (z_i - \hat{z}(x_i, w))^2 \\ &= \frac{2}{2} \sum_{i=1}^N (z_i - \hat{z}(x_i, w)) \frac{\partial}{\partial w} (z_i - \hat{z}(x_i, w)) \\ &= - \sum_{i=1}^N (z_i - \sigma(2wx_i)) \sigma(2wx_i) (1 - \sigma(2wx_i)) 2x_i \\ &= -2 \sum_{i=1}^N x_i (z_i - \sigma(2wx_i)) \sigma(2wx_i) (1 - \sigma(2wx_i)) \end{aligned}$$

Therefore, our learning rule will be written as:

$$w \leftarrow w + 2\eta \sum_{i=1}^N x_i (z_i - \sigma(2wx_i)) \sigma(2wx_i) (1 - \sigma(2wx_i))$$

(b) **Note:** On all matrix dot products between x_i 's and w , in this and the following questions, we're considering here for it to be the dot product between x_i and the **transposed** of w , since otherwise the dimensions wouldn't match (probably an issue regarding the original question's statement).

Considering an initialization of all ones, we have the following:

$$\eta = 1, \quad w = \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix}, \quad X = \begin{bmatrix} 1 & 1 & 1 & 1 \\ 1 & 2 & 1 & 3 \\ 1 & 1 & 3 & 3 \end{bmatrix}$$

Here, we're asked to compute the first gradient descent update - i.e, the first update of the weights after the first epoch. As such, we'll have the following (computations in this sheet's notebook):

$$w \leftarrow w + 2\eta \sum_{i=1}^N x_i (z_i - \sigma(2wx_i)) \sigma(2wx_i) (1 - \sigma(2wx_i))$$

$$\leftarrow \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} + 2 \cdot 1 \cdot \left(\begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \cdot \left(1 - \sigma\left(2 \cdot \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \cdot \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix}\right)\right) \cdot \sigma\left(2 \cdot \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \cdot \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix}\right) \cdot \left(1 - \sigma\left(2 \cdot \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \cdot \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix}\right)\right) \right) + \dots = \begin{bmatrix} 0.99992 \\ 0.999917 \\ 0.999735 \end{bmatrix}$$

- (c) If in the previous question, with a batch gradient descent update, we computed the weight update for the whole first epoch at once, with stochastic gradient descent we'll compute the weight update for each sample individually - the "sum" in the learning rule, here, will be replaced by a single sample, as can be seen below:

$$w \leftarrow w + 2\eta x_1 (z_1 - \sigma(2wx_1)) \sigma(2wx_1) (1 - \sigma(2wx_1))$$

This will be essentially calculating the expression written in the question above (before the \dots), which ends up updating the weight matrix to the following:

$$w_{new} = \begin{bmatrix} 1.00001 \\ 1.00001 \\ 1.00001 \end{bmatrix}$$

3. Let us consider the following activation function, with the training data set from the previous exercise:

$$\hat{z}(x, w) = \frac{1}{1 + e^{-wx}}$$

Here, we'll be using the cross-entropy loss function:

$$E(w) = - \sum_{i=1}^N z_i \log \hat{z}(x_i, w) + (1 - z_i) \log(1 - \hat{z}(x_i, w))$$

- (a) Determine the gradient descent learning rule for this unit.
- (b) Compute the first gradient descent update, assuming an initialization of all ones.
- (c) Compute the first stochastic gradient descent update assuming an initialization of all ones.

- (a) Here, we'll compute the learning rule for the gradient descent the same way, with the exception of the gradient of the error function, of course: here, we're working with the cross-entropy loss function, instead of the squared error function. Thankfully, and even though the deduction for such derivative is not at all trivial, it is also possible to bring it in our note sheet to the exam. As such, we'll have the following:

$$E(w) = - \log(P(z|x, w)) = - \sum_{i=1}^N (z_i \log \hat{z}(x_i, w) + (1 - z_i) \log(1 - \hat{z}(x_i, w)))$$

$$\frac{\partial E}{\partial w} = - \sum_{i=1}^N (x_i, w)' (z_i - \hat{z}(x_i, w))$$

In our case, considering that the activation function is a sigmoid once again (this time $\sigma(wx)$), we'll have:

$$\frac{\partial E}{\partial w} = - \sum_{i=1}^N (w \cdot x_i)' (z_i - \sigma(wx_i)) = - \sum_{i=1}^N x_i (z_i - \sigma(wx_i))$$

As such, the learning rule will be:

$$w \leftarrow w + \eta \sum_{i=1}^N x_i (z_i - \sigma(wx_i))$$

(b) Once again, regarding an initialization of all ones, we'll have the following:

$$\eta = 1, \quad w = \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix}$$

Now, with the new learning rule, we'll be able to write the first batch gradient descent update as:

$$w \leftarrow w + \eta \sum_{i=1}^N x_i (z_i - \sigma(wx_i)) = \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} + 1 \cdot \left(\begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \cdot \left(1 - \sigma \left(\begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \cdot \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \right) \right) \right) + \dots = \begin{bmatrix} -0.926984 \\ -2.90718 \\ -4.91178 \end{bmatrix}$$

(c) Once again, with a stochastic gradient descent update, we'll compute the weight update for each sample individually - here, considering the **first** update, we're only going to consider the first sample. As such, we'll have:

$$w \leftarrow w + \eta x_1 (z_1 - \sigma(wx_1)) = \begin{bmatrix} 1.04743 \\ 1.04743 \\ 1.04743 \end{bmatrix}$$

4. Consider now the half sum of squared errors loss function, paired with the following activation function:

$$\hat{z}(x, w) = e^{(wx)^2}$$

(a) Determine the gradient descent learning rule for this unit.

(b) Compute the stochastic gradient descent update for input $x_{new} = [1 \ 1]^T$, $z_{new} = 0$, with initial weights $w = [0 \ 1 \ 0]^T$ and learning rate $\eta = 2$.

(a) Once again, we'll have to compute the derivative of the loss function with respect to the weights. As such, we'll have:

$$E(w) = \frac{1}{2} \sum_{i=1}^N (z_i - \hat{z}(x_i, w))^2$$

$$\frac{\partial E}{\partial w} = \frac{\partial}{\partial w} \frac{1}{2} \sum_{i=1}^N (z_i - \hat{z}(x_i, w))^2 = \sum_{i=1}^N (z_i - \hat{z}(x_i, w)) \frac{\partial}{\partial w} (z_i - \hat{z}(x_i, w)) = -2 \sum_{i=1}^N x_i (wx_i) (z_i - e^{(wx_i)^2}) e^{(wx_i)^2}$$

Our learning will be, then:

$$w \leftarrow w + 2\eta \sum_{i=1}^N x_i (wx_i) (z_i - e^{(wx_i)^2}) e^{(wx_i)^2}$$

- (b) We're asked to perform a stochastic gradient descent update for a new query vector $x_{new} = [1 \ 1]^T$, with label 0. Moreover, the weights are initialized as $w = [0 \ 1 \ 0]^T$ and the learning rate is $\eta = 2$. As such, and plugging these values onto our learning rule, we'll have the following:

$$\begin{aligned} w &\leftarrow w + 2\eta x_{new} (wx_{new}) (z_{new} - e^{(wx_{new})^2}) e^{(wx_{new})^2} \\ &\leftarrow \begin{bmatrix} 0 \\ 1 \\ 0 \end{bmatrix} + 2 \cdot 2 \cdot \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \cdot \left(\begin{bmatrix} 0 \\ 1 \\ 1 \end{bmatrix} \cdot \begin{bmatrix} 1 \\ 1 \\ 1 \end{bmatrix} \right) \dots \\ &\leftarrow \begin{bmatrix} -29.5562 \\ -28.5562 \\ -29.5562 \end{bmatrix} \end{aligned}$$

5. Consider the sum squared and cross-entropy loss functions. Any stands out? What changes when one changes the loss function?

Note: this answer was mostly written by [João Rocha](#).

The cross-entropy loss function stands out as it corresponds to making a MLE of the parameter w under the assumption that the data is generated by a **Bernoulli** distribution, such that the outputs z are sampled from a random variable Z with probability $p(Z = 1) = \sigma(wx)$ and $p(Z = 0) = 1 - \sigma(wx)$. As such, the cross-entropy loss function is the negative log-likelihood of the data under such distribution.

The sum of squares loss function would be more fitted to a regression scenario, since it matches a MLE estimation of w under the assumption that the outputs z are sampled from a random variable $Z \sim \mathcal{N}(wx, \sigma^2)$. Nevertheless, one can use the sum of squares loss function on classification problems as well - as a matter of fact, this function is also convex, thus the use of gradient descent will also find a global minimum. In this case, the difference between the functions will be the value of the minimal parameter: since the functions are different, most likely their minimal values will be different as well, leading to different estimations of the parameter w ; regarding classification problems, like the one in hands, the cross-entropy loss function should generally lead to a better estimation of the parameter w .