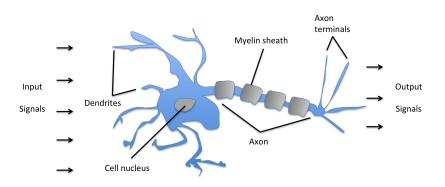
Chapter 2

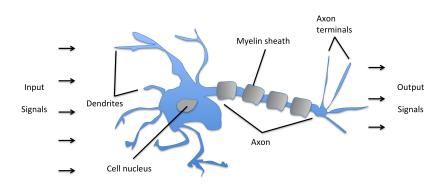
Training Machine Learning Algorithms for Classification

August 14, 2017

Biology



Logic Gate



- Simple logic gate with binary outputs
- Signals arrive at dendrites
- Integrated into cell body
- If signal exceeds threshold, generate output, and pass to axon

Rosenblatt Perceptron

- Binary classification task
- Positive class (1) vs. negative class (-1)
- Define activation function $\phi(z)$
- Takes as input a dot product of input and weights
- Net input: $z = w_1x_1 + \cdots + w_mx_m$

$$\mathbf{w} = \begin{bmatrix} w^{(1)} \\ w^{(2)} \\ \vdots \\ w^{(m)} \end{bmatrix}, \mathbf{x} = \begin{bmatrix} x^{(1)} \\ x^{(2)} \\ \vdots \\ x^{(m)} \end{bmatrix}$$

Heaviside step function

- $\phi(z)$ known as activation
- if activation above some threshold, predict class 1
- predict class -1 otherwise

Heaviside Step Function

$$\phi(z) = egin{cases} 1 & ext{if } z \geq heta \ -1 & ext{otherwise} \ . \end{cases}$$

Step function simplified

Bring the threshold θ to the left side of the equation and define a weight-zero as $w_0=-\theta$ and $x_0=1$, so that we write ${\bf z}$ in a more compact form

$$z = w_0 x_0 + w_1 x_1 + \cdots + w_m x_m = \mathbf{w}^\mathsf{T} \mathbf{x}$$

and

$$\phi(z) = egin{cases} 1 & ext{if } z \geq 0 \ -1 & ext{otherwise} \ . \end{cases}$$

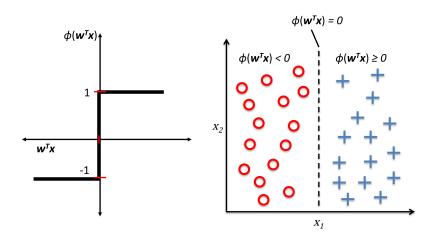
Basic Linear Algebra

Vector dot product

$$z = \mathbf{w}^\mathsf{T} \mathbf{x} = \sum_{j=0}^m \mathbf{w_j} \mathbf{x_j}$$

$$\begin{bmatrix} 1 & 2 & 3 \end{bmatrix} \times \begin{bmatrix} 4 \\ 5 \\ 6 \end{bmatrix} = 1 \times 4 + 2 \times 5 + 3 \times 6 = 32.$$

Input squashed into a binary output



Rosenblatt perceptron algorithm

- Initialize the weights to 0 or small random numbers.
- ② For each training sample $\mathbf{x}^{(i)}$, perform the following steps:
 - **1** Compute the output value \hat{y} .
 - Opdate the weights.

Weight update

Weight update rule:

$$w_j := w_j + \Delta w_j$$

Perceptron learning rule:

$$\Delta w_j = \eta \left(y^{(i)} - \hat{y}^{(i)} \right) x_j^{(i)}$$

Where η is the learning rate (a constant between 0.0 and 1.0), $y^{(i)}$ is the true class label of the *i*th training sample, and $\hat{y}^{(i)}$ is the predicted class label.

Update rule examples

Correct prediction, weights unchanged:

$$\Delta w_j = \eta \bigg(-1 - 1 \bigg) x_j^{(i)} = 0$$

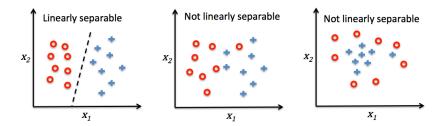
$$\Delta w_j = \eta \bigg(1 - 1 \bigg) x_j^{(i)} = 0$$

Wrong prediction, weights pushed towards the positive or negative class:

$$\Delta w_j = \eta \left(1 - 1\right) x_j^{(i)} = \eta(2) x_j^{(i)}$$

$$\Delta w_j = \eta \left(-1 - 1 \right) x_j^{(i)} = \eta(-2) x_j^{(i)}$$

Linear separability



Convergence

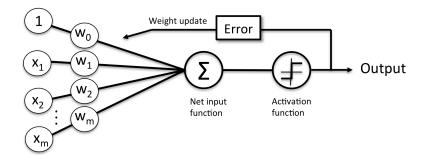
Convergence guaranteed if

- The two classess linearly separable
- Learning rate is sufficiently small

If classes cannot be seprated:

- Set a maximum number of passes over the training dataset (epochs)
- Set a threshold for the number of tolerated misclassifications
- Otherwise, it will never stop updating weights (converge)

Linear separability



Perceptron implementation

→ iPython notebook on github

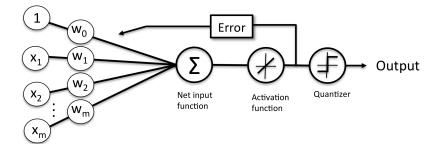
ADAPtive LInear NEuron (Adaline)

- Weights updated based on a linear activation function
- Remember that perceptron used a unit step function
- \bullet $\phi(z)$ is simply the identity function of the net input

$$\phi(\mathbf{w}^T\mathbf{x}) = \mathbf{w}^T\mathbf{x}$$

A quantizer is then used to predict class label

Adaline: notice the difference with perceptron



Cost functions

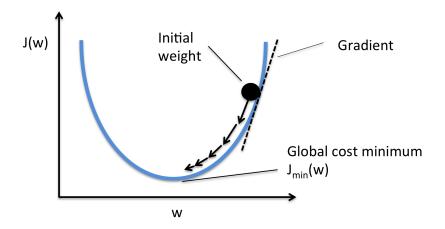
- ML algorithms often define an objective function
- This function is optimized during learning
- It is often a cost function we want to minimize
- Adaline uses a cost function $J(\cdot)$
- Learns weights as the sum of squared errors (SSE)

$$J(\mathbf{w}) = \frac{1}{2} \sum_{i} \left(y^{(i)} - \phi(z^{(i)}) \right)^{2}$$

Advantages of Adaline cost function

- The linear activation function is differentiable
- Unlike the unit step function
- It is convex
- Can use gradient descent to learn the weights

Gradient Descent



Gradient Descent

- Weights updated by taking small steps
- Step size determined by learning rate
- Take a step away from the gradient $\nabla J(\mathbf{w})$ of the cost function

$$\mathbf{w} := \mathbf{w} + \Delta \mathbf{w}$$
.

• The weight change is defined as follows:

$$\Delta \mathbf{w} = -\eta \nabla J(\mathbf{w})$$

Gradient computation

To compute the gradient of the cost function, we need to compute the partial derivative of the cost function with respect to each weight w_j ,

$$\frac{\partial J}{\partial w_j} = -\sum_i \left(y^{(i)} - \phi(z^{(i)}) \right) x_j^{(i)},$$

Weight update of weight w_j

$$\Delta w_j = -\eta \frac{\partial J}{\partial w_j} = \eta \sum_i \left(y^{(i)} - \phi(z^{(i)}) \right) x_j^{(i)}$$

We update all weights simultaneously, so Adaline learning rule becomes

$$\mathbf{w} := \mathbf{w} + \Delta \mathbf{w}$$
.



Partial derivatives

$$\frac{\partial J}{\partial w_j} = \frac{\partial}{\partial w_j} \frac{1}{2} \sum_i \left(y^{(i)} - \phi(z^{(i)}) \right)^2
= \frac{1}{2} \frac{\partial}{\partial w_j} \sum_i \left(y^{(i)} - \phi(z^{(i)}) \right)^2
= \frac{1}{2} \sum_i 2(y^{(i)} - \phi(z^{(i)})) \frac{\partial}{\partial w_j} \left(y^{(i)} - \phi(z^{(i)}) \right)
= \sum_i \left(y^{(i)} - \phi(z^{(i)}) \right) \frac{\partial}{\partial w_j} \left(y^{(i)} - \sum_i \left(w_j^{(i)} x_j^{(i)} \right) \right)
= \sum_i \left(y^{(i)} - \phi(z^{(i)}) \right) \left(- x_j^{(i)} \right)
= -\sum_i \left(y^{(i)} - \phi(z^{(i)}) \right) x_j^{(i)}$$

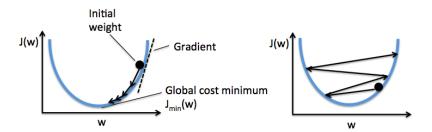
Adaline learning rule vs. Perceptron rule

- Looks identical
- $\phi(z^{(i)})$ with $z^{(i)} = \mathbf{w}^T \mathbf{x}^{(i)}$ is a real number
- And not an integer class label as in Perceptron
- The weight update is done based on all samples in training set
- Perceptron updates weights incrementally after each sample
- This approach is known as "batch" gradient descent

Perceptron implementation

→ iPython notebook on github

Lessons learned



- Learning rate too high: error becomes larger (overshoots global min)
- Learning rate too low: takes many epochs to converge
- Feature normalization

Stochastic gradient descent (SGD)

- Large dataset with millions of data points ("big data")
- Batch gradient descent costly
- Need to compute the error for the entire dataset ...
- ... to take one step towards the global minimum!

$$\Delta \mathbf{w} = \eta \sum_{i} \left(y^{(i)} - \phi(z^{(i)}) \right) \mathbf{x}^{(i)}.$$

SGD updates the weights incrementally for each training sample

$$\Delta \mathbf{w} = \eta \left(y^{(i)} - \phi(z^{(i)}) \right) \mathbf{x}^{(i)}.$$

SGD details

- Approximation of gradient descent
- Reaches convergence faster because of frequent weight updates
- Important to present data in random order
- Learning rate often gradually decreased (adaptive learning rate)
- Can be used for online learning
- Middle ground between SGD and batch GD is known as mini-batch learning
 - E.g. 50 examples at a time
 - Can use vector/matrix operations rather than loops as in SGD
 - Vectorized operations highly efficient