Classification

To attempt classification, one method is to use linear regression and map all predictions greater than 0.5 as a 1 and all less than 0.5 as a 0. However, this method doesn't work well because classification is not actually a linear function.

The classification problem is just like the regression problem, except that the values we now want to predict take on only a small number of discrete values. For now, we will focus on the **binary classification problem** in which y can take on only two values, 0 and 1. (Most of what we say here will also generalize to the multiple-class case.) For instance, if we are trying to build a spam classifier for email, then X(i) may be some features of a piece of email, and y may be 1 if it is a piece of spam mail, and 0 otherwise. Hence, $y \in \{0,1\}$. 0 is also called the negative class, and 1 the positive class, and they are sometimes also denoted by the symbols "-" and "+." Given X(i), the corresponding Y(i) is also called the label for the training example.

Hypothesis Representation

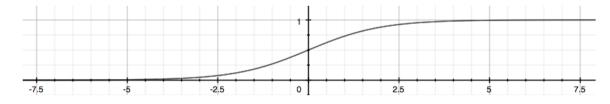
We could approach the classification problem ignoring the fact that y is discrete-valued, and use our old linear regression algorithm to try to predict y given x. However, it is easy to construct examples where this method performs very poorly. Intuitively, it also doesn't make sense for he(x) to take values larger than 1 or smaller than 0 when we know that $y \in \{0, 1\}$. To fix this, let's change the form for our hypotheses he(x) to satisfy $0 \le he(x) \le 1$. This is accomplished by plugging $\theta \tau x$ into the Logistic Function.

Our new form uses the "Sigmoid Function," also called the "Logistic Function":

$$h_{ heta}(x) = g(heta^T x)$$

$$z = heta^T x \ g(z) = rac{1}{1 + e^{-z}}$$

The following image shows us what the sigmoid function looks like:



The function g(z), shown here, maps any real number to the (0, 1) interval, making it useful for transforming an arbitrary-valued function into a function better suited for classification.

 $h\theta(x)$ will give us the **probability** that our output is 1. For example, $h\theta(x)=0.7$ gives us a probability of 70% that our output is 1. Our probability that our prediction is 0 is just the complement of our probability that it is 1 (e.g. if probability that it is 1 is 70%, then the probability that it is 0 is 30%).

$$h_{\theta}(x) = P(y = 1|x; \theta) = 1 - P(y = 0|x; \theta)$$

 $P(y = 0|x; \theta) + P(y = 1|x; \theta) = 1$

Decision Boundary

In order to get our discrete 0 or 1 classification, we can translate the output of the hypothesis function as follows:

$$h_{ heta}(x) \geq 0.5
ightarrow y = 1 \ h_{ heta}(x) < 0.5
ightarrow y = 0$$

The way our logistic function g behaves is that when its input is greater than or equal to zero, its output is greater than or equal to 0.5:

$$g(z) \ge 0.5$$

when $z > 0$

Remember.

$$egin{aligned} z=0, e^0=1 &\Rightarrow g(z)=1/2 \ z & o \infty, e^{-\infty} & o 0 \Rightarrow g(z)=1 \ z & o -\infty, e^\infty & o \infty \Rightarrow g(z)=0 \end{aligned}$$

So if our input to g is $\theta^{\Lambda}TX$, then that means:

$$h_{ heta}(x) = g(heta^T x) \geq 0.5$$
 when $heta^T x \geq 0$

From these statements we can now say:

$$heta^T x \geq 0 \Rightarrow y = 1 \ heta^T x < 0 \Rightarrow y = 0$$

The **decision boundary** is the line that separates the area where y = 0 and where y = 1. It is created by our hypothesis function.

Example

$$egin{aligned} heta = egin{bmatrix} 5 \ -1 \ 0 \end{bmatrix} \ y = 1 \ if \ 5 + (-1)x_1 + 0x_2 \geq 0 \ 5 - x_1 \geq 0 \ -x_1 \geq -5 \ x_1 \leq 5 \end{aligned}$$

In this case, our decision boundary is a straight vertical line placed on the graph where $x_1=5$, and everything to the left of that denotes y=1, while everything to the right denotes y=0.

Again, the input to the sigmoid function g(z) (e.g. $\theta^{\Lambda}TX$) doesn't need to be linear, and could be a function that describes a circle (e.g. $z=\theta_0+\theta_1x_{12}+\theta_2x_{22}$) or any shape to fit our data.