Linear Regression

our dataset format is $(x^{(i)}, y^{(i)})$, and we have dataset size m, so where i is ith data, and for $x^{(i)}$, it may have n features, that to say, for any x,

$$x = \{x_0, x_1, ..., x_n\}$$

Note that dimension of x is n+1, that is to simplify the calculation, x_0 often equal to 1, to make θ_0 a constant term.

To fit LR on our dataset, we assume:

$$h_{ heta} = heta_0 x_0 + heta_1 x_1 + .., + heta_n x_n = heta^T x_n$$

and $h_{ heta}^{(i)}$ is LR predicted output for $x^{(i)}$, there $x_0=1.$

So how do we know our predition is reliable? we introduce cost function here.

$$J(heta) = rac{1}{m} \sum_{i=0}^m (h_ heta(x)^{(i)} - y^{(i)})^2$$

the larger the $J(\theta)$, the more inaccurate the prediction result is. so we must choose a $\theta=[\theta_0,\theta_1,...,\theta_n]$ to minimise $J(\theta)$

Gradient Descent

we update θ to minimise $J(\theta)$ by

$$\theta_j = \theta_j - \alpha \cdot \frac{1}{m} \frac{\partial J(\theta)}{\partial \theta_j}$$

and when m=1

$$egin{aligned} rac{\partial J(heta)}{\partial heta_i} &= 2 \cdot rac{1}{2} (h_ heta - y) \cdot rac{\partial}{\partial heta j} (h_ heta x - y) \ &= 2 \cdot rac{1}{2} (h_ heta - y) rac{\partial}{\partial heta j} (\sum_{i=0}^m heta_i x_i - y) \ &= (h_ heta - y) x_j \end{aligned}$$

$$heta_j = heta_j - lpha \cdot rac{1}{m} (h_ heta - y) x_j$$

it is called gradient descent.

Normal equation

it is other way to get appropriate θ .

$$\theta = (X^T X)^{-1} \cdot (X^T y)$$

if X^TX is noninvertible, it may caused by:

- Redundant features, where two features are very closely related (i.e. they are linearly dependent)
- Too many features (e.g. m ≤ n). In this case, delete some features or use "regularization" (to be explained in a later lesson).

Reference

Mechine Learning