

Lecture - 10 (Logistic Regression)

probabilistic :-

$$P(Y=y | X=x) = \frac{1}{1 + \exp(y\langle \theta, x \rangle)}$$

$$y = (+1) \text{ or } (-1) \Rightarrow \frac{1}{1 + \exp(\langle \theta, x \rangle)} \quad \frac{1}{1 + \exp(-\langle \theta, x \rangle)}$$

$$\rightarrow P(Y=1/x) + P(Y=-1/x) = 1$$

$$\rightarrow \text{Decision boundary} \quad P(Y=1/x) = P(Y=-1/x) = 0.5$$

$$\rightarrow P(Y=1/x) > 0.5 \Rightarrow \text{greater confidence class 1 than -1}$$

Prediction Confidence

Task	Or
\rightarrow predict the label associated with a feature vector x	\rightarrow predict value $\text{Sign}(\langle \theta, x \rangle)$
\rightarrow Measure of confidence of that prediction	$\rightarrow P(Y=y X=x) = \frac{1}{1 + \exp(y\langle \theta, x \rangle)}$

properties of Linear classifiers

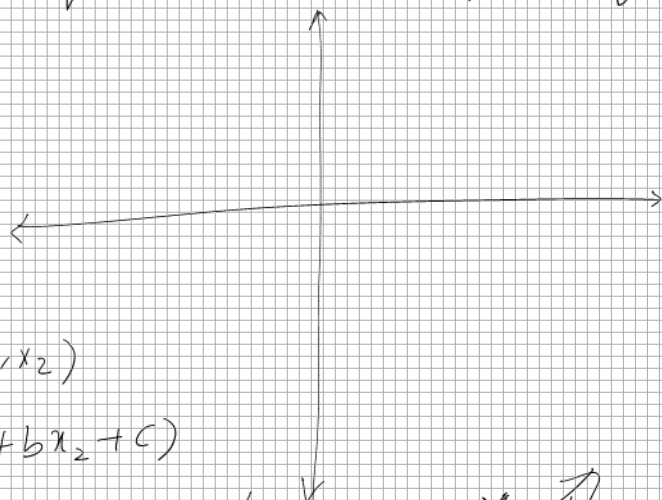
- ↳ ① Easy to train
- ↳ ② Can predict labels very fast
- ↳ ③ Statistical theory of linear classifiers leading to effective modeling strategies.
- ↳ ④ Linear classifiers - Extend to higher dimensions
↳ ~~extra~~ computational load
- ↳ (If x is sparse, the inner product $\langle \theta, x \rangle$ is extremely fast)

$$\text{Bias Term} = \langle \theta, x \rangle + b = x_1 \theta_1 + x_2 \theta_2 + \dots + x_d \theta_d + b$$

Bias term makes our model

→ linear classifiers have the form $y = \text{sign}(\langle \theta, x \rangle)$

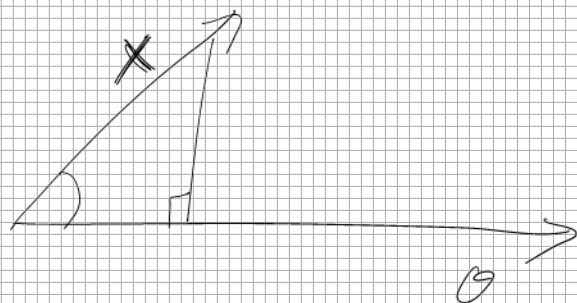
→



→ Given $x = (x_1, x_2)$

→ $\text{Sign}(ax_1 + bx_2 + c)$

→ classification problem /
decision boundary: don't
depend whether θ is normalized
or not.



→ Should the decision boundary pass
through origin?

↳ x a parameter
vector is zero.

Answer: Not necessarily.

If we don't require that (by requiring
one features to value 1) the classifier
becomes considerably more powerful.
i.e. we can learn much more general
decision boundaries

$$x_1\theta_1 + x_2\theta_2 + \dots + x_d\theta_d + c$$

without effort, the decision boundary has to to
pass through origin. i.e. $\langle \theta, x \rangle = 0$
at origin $x=0, \theta=0 \Rightarrow \langle \theta, x \rangle = 0$

→ Decision boundary passing through origin restricts the model in a significant way.

$$n_1 \theta_1 + n_2 \theta_2 + \dots + n_d \theta_d + c = \langle x, \theta \rangle + c$$

$$(x_1, x_2, \dots, x_d, 1)$$

$\langle x, \theta \rangle$ \leftarrow θ is $d+1$ dimensional.

→ MLE

Frequentists: MLE uses pairs of feature vectors & labels, usually from historic data to estimate a model classifier or feature.

Bayesian: A single classifier cannot represent the "truth". Estimate the mixed probability that each classifier is correct & use them all.

Justification for using MLE

→ It converges to the optimal solution in the limit of large data (consistency).

→ Convergence occurs at the fastest possible rate of convergence (statistical efficiency).

Gradient Descent: is a suitable to large data as long as it is sparse, not so much dependent on the dimensionality.
Independent (θ, x) = easy to calculate

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$$\theta_j \leftarrow \theta_j - \alpha \frac{\sum_{i=1}^n \log(1 + \exp(y^{(i)} \langle \theta, x^{(i)} \rangle))}{2 \theta_j}$$

when $x^{(i)}$ are sparse, partial derivatives can be made particularly fast.

→ Stochastic Gradient Converges faster than Gradient descent.
