## Classification: basics

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### **OVERVIEW**

Binary Classification

Multi-class Classification

LDA and QDA

Logistic Regression: binary case

Logistic Regression: multi-class case, sparsity

Others: nonparametric classifier, robust loss, perceptrons

## Binary Classification

## Sensitivity and Specificity

Imagine a scenario where people are tested for a disease:

- ► The test outcome: positive (sick) or negative (healthy)
- ► The actual status: positive (sick) or negative (healthy)

There are four possible scenarios:

- ► True positive (TP): sick people correctly identified as sick
- ▶ False positive (FP) : healthy people incorrectly identified as sick
- ▶ True negative (TN): healthy people correctly identified as healthy
- ▶ False negative (FN): sick people incorrectly identified as healthy

	Test Outcome		
True outcome	Positive	Negative	Total
Positive	True Pos. (TP)	False Neg. (FN)	P
Negative	False Pos. (FP)	True Neg. (TN)	N
	$P^*$	$N^*$	

	Test Outcome		
True outcome	Positive	Negative	Total
Positive	True Pos. (TP)	False Neg. (FN)	P
Negative	False Pos. (FP)	True Neg. (TN)	N
	$P^*$	$N^*$	

Accuracy (1-Error): 
$$(TP + TN)/(TP + TN + FP + FN)$$
.

Sensitivity (true positive rate/power/recall): the proportions of positives that are correctly identified

Sensitivity = 
$$TP/P = TP/(TP + FN)$$

Specificity (true negative rate): the proportions of negative that are correctly identified

Specificity = 
$$TN/N = TN/(FP + TN)$$

- ▶ False positive rate (Type I error): = 1 -Specificity.
- ▶ False negative rate (Type II error): = 1 Sensitivity
- ▶ False discovery rate (FDR/precision): the proportion of predicted positives that are in fact false positives  $FDR = FP/P^*$

## example 1

	Predicted	
True	email	spam
email	573	40
spam	53	334

spam =presence of disease, email=absence of disease

$$\begin{array}{ll} \text{specificity} &= 100 \times \frac{573}{573+40} = 93.4\% \\ \text{sensitivity} &= 100 \times \frac{334}{334+53} = 86.3\% \end{array}$$

### example 2

Threshold A is chosen to balance sensitivity and specificity without leaning too heavily towards either.

True Positive (TP)	False Negative (FN)	
40	5	
False Positive (FP)	True Negative (TN)	
10	45	

With Threshold A, we have: - Sensitivity: 88.9% - Specificity: 81.8%

Threshold B (it has a lower criterion for a positive test result)—adjusted to make the test more sensitive to detecting the disease.

True Positive (TP)	False Negative (FN)	
45	0	
False Positive (FP)	True Negative (TN)	
20	35	

With Threshold B, we have: - Sensitivity: 100% - Specificity: 63.6%

## Binary classification: problem

- ▶ input vector  $X \in \mathcal{X} \subset \mathbb{R}^d$
- ightharpoonup output  $Y \in \{0,1\}$
- ▶ "hard classifier"  $h: \mathcal{X} \longrightarrow \{0, 1\}$

The rule is characterized as

$$h(X) = 1(b(X) > 0)$$

where b is the boundary function (or discriminant function) that gives the decision boundary  $\{x:b(x)=0\}$ .

- ▶ If b(X) is a linear in X, then the classifier has a linear boundary (in X-space).
  - ▶ With transformed X included, the classifier can have a nonlinear boundary (in X-space).

## Examples of linear boundary

Linear logit model:

Assume that the **logit function** is linear in x, i.e.,

$$b(x) = \log \frac{\Pr(Y = 1 \mid X = x)}{\Pr(Y = 0 \mid X = x)} = \beta_0 + \beta_1^{\top} x$$

Thus the classification boundary is given by  $\{x: \beta_0 + \beta_1^\top x = 0\}$ 

Examples: LDA, Logistic regression (see later)

The classification error rate, of h is defined as

$$R(h) = E_{X,Y}(1(Y \neq h(X))) = P(Y \neq h(X))$$

The rule h that minimizes R(h) is

$$h^*(x) = \begin{cases} 1 & \text{if } m(x) > \frac{1}{2} \\ 0 & \text{otherwise} \end{cases}$$

where  $m(x) = P(Y = 1 \mid X = x) = E(Y \mid X = x)$ .

- ► This optimal rule is called the **Bayes rule (classifier)** (under equal costs).
- ▶ The risk  $R(h^*)$  is called the **Bayes risk**.
- ▶ The set  $\{x: m(x) \frac{1}{2} = 0\}$  is called the **Bayes decision** boundary.

Alternatively, the Bayes rule is  $h^*$ , is given by

$$h^*(x) = \begin{cases} 1 & \text{if} \quad P(Y=1 \mid X=x) > P(Y=0 \mid X=x) \\ 0 & \text{if} \quad P(Y=1 \mid X=x) < P(Y=0 \mid X=x) \end{cases}$$

The classification boundary of the Bayes rule is

$$\{x: P(Y = 1 \mid X = x) = P(Y = 0 \mid X = x)\}\$$
$$= \{x: P(Y = 1 \mid X = x) - 0.5 = 0\}\$$

From Bayes' theorem

$$p(Y = 1 \mid X = x) = \frac{\pi_1 p_1(x)}{\pi_1 p_1(x) + (1 - \pi_1) p_0(x)}$$

- $\pi_1 = p(Y = 1), \pi_0 = p(Y = 0)$ : the marginal distribution of Y (prior class probabilities)
- ▶  $p_j(x) = p(x \mid Y = j)$ : the conditional density of X given that Y = j.

$$h^*(x) = \begin{cases} 1 & \text{if } \frac{p_1(x)}{p_0(x)} > \frac{\pi_0}{\pi_1} \\ 0 & \text{otherwise.} \end{cases}$$

This decision-making process balances two types of information:

- ▶ Likelihood ratio (evidence)  $\frac{p_1(x)}{p_0(x)}$ : compares how probable it is that the observed data x comes from class 1 as opposed to class 0.
- Prior ratio  $\frac{\pi_0}{\pi_1}$ : our initial belief about the relative frequency of class 0 to class 1 before seeing any data (say x).

## Unequal Losses

For any decision function, there are two possible errors:

- ▶ misclassifying a sample in class 0 to 1 (false positive)
- ▶ misclassifying a sample in class 1 to 0 (false negative)

Each type of error is associated with a cost (the price to pay for the consequence):

- $\blacktriangleright$  L(1,0) is the cost of misclassifying a sample in class 1 to 0
- ightharpoonup L(0,1) is the cost of misclassifying a sample in class 0 to 1.

We assume L(j,j) = 0 for j = 0, 1; but it may not be L(0,1) = L(1,0).

The loss becomes

$$L(Y, h(X)) = L(1, 0)1(Y = 1, h(X) = 0) + L(0, 1)1(Y = 0, h(X) = 1)$$

For fixed x, the Bayes rule is given as

$$h^*(x) = \begin{cases} 1 & \text{if } L(1,0)P(Y=1 \mid X=x) > L(0,1)P(Y=0 \mid X=x) \\ 0 & \text{if } L(1,0)P(Y=1 \mid X=x) < L(0,1)P(Y=0 \mid X=x) \end{cases}$$

Equivalently,

$$h^*(x) = \begin{cases} 1 & \text{if } & \frac{P(Y=1|X=x)}{P(Y=0|X=x)} > \frac{L(0,1)}{L(1,0)} \\ 0 & \text{if } & \frac{P(Y=1|X=x)}{P(Y=0|X=x)} < \frac{L(0,1)}{L(1,0)} \end{cases}$$

the Bayes rule

$$h^*(x) = 1 \left\{ x : P(Y = 1 \mid X = x) > \frac{L(0,1)}{L(0,1) + L(1,0)} \right\}.$$

In light of the Bayes' theorem,

- $\pi_1 = p(Y = 1), \pi_0 = p(Y = 0)$ : the marginal distribution of Y (prior class probabilities)
- ▶  $p_j(x) = p(x \mid Y = j)$ : the conditional density of X given that Y = j.

$$h^*(x) = \begin{cases} 1 & \text{if } & \frac{p_1(x)}{p_0(x)} > \frac{\pi_0 L(0,1)}{\pi_1 L(1,0)} \\ 0 & \text{if } & \frac{p_1(x)}{p_0(x)} < \frac{\pi_0 L(0,1)}{\pi_1 L(1,0)} \end{cases}$$

By changing the weights for L(0,1) and L(1,0), we can effectively change the classification threshold.

- ightharpoonup L(0,1) = the cost of predicting a "non-disease" example to "disease"
- ightharpoonup L(1,0) = the cost of predicting a "disease" example to "non-disease"

$$h^*(x) = 1 \left\{ x : P(Y = 1 \mid X = x) > \frac{L(0, 1)}{L(0, 1) + L(1, 0)} \right\}.$$

How to increase the sensitivity and decrease the specificity of the rule?

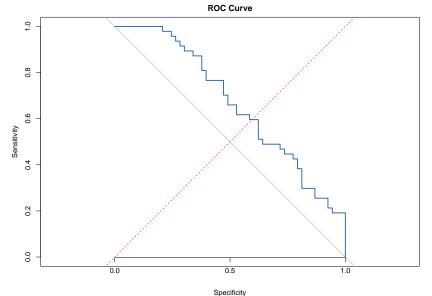
▶ Increase L(1,0) and decrease L(0,1).

How to increase the specificity and decrease the sensitivity of the rule?

▶ Increase L(0,1) and decrease L(1,0).

## Receiver Operating Characteristic (ROC) curve

A ROC curve is a plot of sensitivity v.s. specificity:



- ▶ An ideal ROC curve will hug the top right corner.
- ▶ An alternative ROC curve will be a curve plotting the sensitivity (true positive rate or 1-Type II error) against the false positive rate (Type I error).
- ► The area under curve (AUC) is a commonly used quantitative measure of overal predictive performance.
  - ▶ A value of 0.5 means the predictions were no better than random guessing.

### Multi-class Classification

### Multi-class Classification

- ightharpoonup Class label  $Y \in \{1, \dots, K\}, K \geq 3$ .
- ▶ The classifier  $h : \mathbb{R}^d \longrightarrow \{1, \dots, K\}$ .

The loss function  $L(Y, h(X)) = \sum_{k=1}^{L} \sum_{l=1}^{K} C(l, k) I(Y = l, h(X) = k)$  where C(l, k) = cost of classifying a sample in class l to class k.

The classification risk, or error rate, of h is defined as

$$R(h) = E_{X,Y}(L(Y, h(X)))$$

Using the 0-1 loss, C(k, k) = 0 for any  $k = 1, \dots, K$ , but equal to 1 otherwise, the rule h that minimizes R(h) is

$$h^*(x) = \arg\max_{k=1,\dots,K} P(Y = k \mid x)$$

i.e., assign x to the most probable class using  $P(Y \mid x)$ .

We generally need to estimate multiple discriminant functions  $\delta_k(x), k = 1, \dots, K$ 

- ▶ Each  $\delta_k(x)$  is associated with class k.
- ▶  $\delta_k(x)$  represents the evidence strength of a sample (x, y) belonging to class k.

The decision rule constructed using  $\delta_k$  's is

$$\hat{h}(x) = k^*, \quad \text{ where } \quad k^* = \arg\max_{k=1,\dots,K} \delta_k(x)$$

The decision boundary of the classification rule  $\hat{h}$  between class k and class l is defined as

$$\{x:\delta_k(x)=\delta_l(x)\}$$

Note:  $\delta_k(x)$  is related but need not be exact  $P(Y = k \mid x)$ .

# LDA and QDA

# Gassuain discriminant analysis: Binary classification

If 
$$X \mid Y = 0 \sim N(\mu_0, \Sigma_0)$$
 and  $X \mid Y = 1 \sim N(\mu_1, \Sigma_1)$ ,

Using the Bayes' Theorem,

$$h^*(x) = \begin{cases} 1 & \text{if } r_1^2 < r_0^2 + 2\log\left(\frac{\pi_1}{1-\pi_1}\right) + \log\left(\frac{|\Sigma_0|}{|\Sigma_1|}\right) \\ 0 & \text{otherwise} \end{cases}$$

 $r_i = \sqrt{(x - \mu_i)^\top \Sigma_i^{-1} (x - \mu_i)}$  for i = 0, 1 is the Mahalanobis distance between x and  $\mu_i$ .

Note: LDA is a special case where  $\Sigma_1 = \Sigma_0$ .

# Quadratic discriminant analysis (QDA)

$$\log \frac{\Pr(Y=1 \mid X=x)}{\Pr(Y=0 \mid X=x)} = \log \frac{\pi_1 \phi(x; \mu_1, \Sigma_1)}{\pi_0 \phi(x; \mu_0, \Sigma_0)} = \delta_1(x) - \delta_0(x).$$

The Bayes rule is

$$h^*(x) = \operatorname{argmax}_{k \in \{0,1\}} \delta_k(x)$$

where

$$\delta_k(x) = -\frac{1}{2}\log|\Sigma_k| - \frac{1}{2}(x - \mu_k)^{\top} \Sigma_k^{-1}(x - \mu_k) + \log \pi_k$$

This is called the Gaussian discriminant function

- ▶ The decision boundary:  $\{x \in \mathcal{X} : \delta_1(x) = \delta_0(x)\}$ 
  - quadratic discriminant analysis (QDA): boundary is quadratic

To estimate  $\pi_0, \pi_1, \mu_0, \mu_1, \Sigma_0, \Sigma_1$ :

$$\widehat{\pi}_{0} = \frac{1}{n} \sum_{i=1}^{n} (1 - Y_{i}), \quad \widehat{\pi}_{1} = \frac{1}{n} \sum_{i=1}^{n} Y_{i}$$

$$\widehat{\mu}_{0} = \frac{1}{n_{0}} \sum_{i:Y_{i}=0} X_{i}, \quad \widehat{\mu}_{1} = \frac{1}{n_{1}} \sum_{i:Y_{i}=1} X_{i}$$

$$\widehat{\Sigma}_{0} = \frac{1}{n_{0} - 1} \sum_{i:Y_{i}=0} (X_{i} - \widehat{\mu}_{0}) (X_{i} - \widehat{\mu}_{0})^{\top}$$

$$\widehat{\Sigma}_{1} = \frac{1}{n_{1} - 1} \sum_{i:Y_{i}=1} (X_{i} - \widehat{\mu}_{1}) (X_{i} - \widehat{\mu}_{1})^{\top}$$

## Linear discriminant analysis (LDA)

**LDA** assumes both classes are from Gaussian and they have the same covariance matrix

$$\Sigma_k = \Sigma, \quad k = 0, 1$$

Note that

$$\log \Pr(Y = k \mid X = x) = -\frac{1}{2} (x - \boldsymbol{\mu}_k)^{\top} \Sigma^{-1} (x - \boldsymbol{\mu}_k) + \log \pi_k + \text{ const.}$$

If prior probabilities are same, the LDA classifies x to the class with centroid closest to x, using the squared Mahalanobis distance, based on the common covariance matrix.

Alternatively,

$$h^*(x) = \begin{cases} 1 & \text{if } \delta_1(x) > \delta_0(x) \\ 0 & \text{otherwise} \end{cases}$$

where the Gaussian discriminant function can be simplified

$$\delta_k(x) = x^{\top} \Sigma^{-1} \mu_k - \frac{1}{2} \mu_k^{\top} \Sigma^{-1} \mu_k + \log \pi_k.$$

- ▶ The decision boundary:  $\{x \in \mathcal{X} : \delta_1(x) = \delta_0(x)\}$ 
  - ▶ linear discriminant analysis (LDA): boundary is linear

Pooled estimate of the  $\Sigma$ :

$$\widehat{\Sigma} = \frac{(n_0 - 1)\,\widehat{\Sigma}_0 + (n_1 - 1)\,\widehat{\Sigma}_1}{n_0 + n_1 - 2}$$

## Multi-class classification (trivial extension)

QDA assume that  $X \mid Y = k \sim N(\mu_k, \Sigma_k)$ .

$$h^*(x) = \operatorname{argmax}_k \delta_k(x)$$

where

$$\delta_k(x) = -\frac{1}{2}\log|\Sigma_k| - \frac{1}{2}(x - \mu_k)^{\top} \Sigma_k^{-1}(x - \mu_k) + \log \pi_k.$$

If all Gaussians assumed to have equal variance  $\Sigma$ ,

$$\delta_k(x) = x^{\top} \Sigma^{-1} \mu_k - \frac{1}{2} \mu_k^{\top} \Sigma^{-1} \mu_k + \log \pi_k.$$

The corresponding estimates are given by

$$\widehat{\pi}_{k} = \frac{1}{n} \sum_{i=1}^{n} 1 (y_{i} = k), \quad \widehat{\mu}_{k} = \frac{1}{n_{k}} \sum_{i:Y_{i} = k} X_{i}$$

$$\widehat{\Sigma}_{k} = \frac{1}{n_{k} - 1} \sum_{i:Y_{i} = k} (X_{i} - \widehat{\mu}_{k}) (X_{i} - \widehat{\mu}_{k})^{\top}$$

$$\widehat{\Sigma} = \frac{\sum_{k=0}^{K-1} (n_{k} - 1)\widehat{\Sigma}_{k}}{n - K}.$$

## Logistic Regression: binary case

## Binary case

The logistic regression assumes that

$$p_1(x; \beta_0, \boldsymbol{\beta}_1) := P(Y = 1 \mid X = x) = \frac{\exp(\beta_0 + x^{\top} \boldsymbol{\beta}_1)}{1 + \exp(\beta_0 + x^{\top} \boldsymbol{\beta}_1)}$$

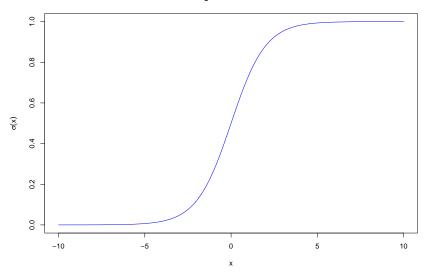
The model can be written as

$$logit(x) := logit(Pr(Y = 1 \mid X = x)) = log \frac{Pr(Y = 1 \mid X = x)}{Pr(Y = 0 \mid X = x)} = \beta_0 + \beta_1^{\top} x$$

- ▶ logit function: logit(a) = log(a/(1-a)) : (0,1)  $\mapsto \mathbb{R}$ 
  - $\triangleright$   $\beta_0 + \beta_1^\top x$ : logits, net input or pre-activation value
- ► The inverse of logit function (**logistic function** or **sigmoid function**):

$$\sigma(a) = \exp(a)/(1 + \exp(a)) : \mathbb{R} \mapsto (0, 1)$$

#### Sigmoid Function



## MLE for logistic models

Notations: assuming  $x_i$  contains the constant term 1 (thus a p+1 vector).

$$\boldsymbol{\beta} := \{\beta_0, \boldsymbol{\beta}_1^\top\}^\top$$

$$\mathbf{y} := [y_1, \cdots, y_n]^\top$$

$$\mathbf{p} := \mathbf{p}(\boldsymbol{\beta}) = [p(x_1; \boldsymbol{\beta}), \cdots, p(x_n; \boldsymbol{\beta})]^\top$$

$$\mathbf{W} := \mathbf{W}(\boldsymbol{\beta}) = \operatorname{diag} \{p(x_i; \boldsymbol{\beta}) (1 - p(x_i; \boldsymbol{\beta}))\} : n \times n$$

The log (conditional) likelihood function is

$$\ell(\beta) = \sum_{i=1}^{n} \{ y_i \log p(x_i; \beta) + (1 - y_i) \log [1 - p(x_i, \beta)] \}$$

The loss function, as a negative loglikelihood function, is called **binomial deviance** (loss):

$$L\left(Y, p_{Y}(X)\right) = -\{1(Y=0)\log(\Pr(Y=0\mid X; \pmb{\theta})) + 1(Y=1)\log(\Pr(Y=1\mid X; \pmb{\theta}))\}$$

$$-\ell(\boldsymbol{\beta}) = -\sum_{i=1}^{n} \left\{ y_i \log p\left(x_i; \boldsymbol{\beta}\right) + (1 - y_i) \log \left[1 - p\left(x_i, \boldsymbol{\beta}\right)\right] \right\}$$
$$= -\sum_{i=1}^{n} \left\{ y_i \boldsymbol{\beta}^{\top} x_i - \log \left[1 + \exp\left(\boldsymbol{\beta}^{\top} x_i\right)\right] \right\}.$$

The score and Hessian are given by

$$\frac{\partial(-\ell(\boldsymbol{\beta}))}{\partial \boldsymbol{\beta}} = -\sum_{i=1}^{n} x_i \left[ y_i - p\left(x_i; \boldsymbol{\beta}\right) \right] = -\mathbf{X}^{\top}(\mathbf{y} - \mathbf{p})$$

$$\frac{\partial^2(-\ell(\boldsymbol{\beta}))}{\partial \boldsymbol{\beta} \partial \boldsymbol{\beta}^{\top}} = \sum_{i=1}^{n} x_i x_i^{\top} p\left(x_i; \boldsymbol{\beta}\right) \left[ 1 - p\left(x_i; \boldsymbol{\beta}\right) \right] = \mathbf{X}^{\top} \mathbf{W} \mathbf{X} \quad (p.s.d)$$

Gradient descent step: In the k-th sep,

$$\boldsymbol{\beta}^{(k+1)} = \boldsymbol{\beta}^{(k)} + \alpha \times \mathbf{X}^{\top} (\mathbf{y} - \mathbf{p}^{(k)})$$

Newton-Raphson step: In the k-th sep,

$$\boldsymbol{\beta}^{(k+1)} = \boldsymbol{\beta}^{(k)} + \left(\mathbf{X}^{\top}\mathbf{W}^{(k)}\mathbf{X}\right)^{-1}\mathbf{X}^{\top}(\mathbf{y} - \mathbf{p}^{(k)})$$

$$= \left(\mathbf{X}^{\top}\mathbf{W}^{(k)}\mathbf{X}\right)^{-1}\mathbf{X}^{\top}\mathbf{W}^{(k)}\left(\mathbf{X}\boldsymbol{\beta}^{(k)} + \mathbf{W}^{(k)^{-1}}(\mathbf{y} - \mathbf{p}^{(k)})\right)$$

$$= \left(\mathbf{X}^{\top}\mathbf{W}^{(k)}\mathbf{X}\right)^{-1}\mathbf{X}^{\top}\mathbf{W}^{(k)}\mathbf{z}^{(k)}$$

where we defined the adjusted response

$$\mathbf{z}^{(k)} = \mathbf{X}\boldsymbol{\beta}^{(k)} + \mathbf{W}^{(k)^{-1}}(\mathbf{y} - \mathbf{p}^{(k)})$$

The Newton-Raphson's approach is equivalent to **Iteratively** Reweighted Least Squares Algorithm:

$$\boldsymbol{\beta}^{(k+1)} = \arg\min_{\boldsymbol{\beta}} (\mathbf{z}^{(k)} - \mathbf{X}\boldsymbol{\beta})^{\top} \mathbf{W}^{(k)} (\mathbf{z}^{(k)} - \mathbf{X}\boldsymbol{\beta}).$$

# Interpretation of $\beta_j$

$$e^{\beta_j} = \underbrace{\frac{P\left(Y=1|\ldots,X_j=x+1,\ldots\right)/P\left(Y=0|\ldots,X_j=x+1,\ldots\right)}{P\left(Y=1|\ldots,X_j=x,\ldots\right)/P\left(Y=0|\ldots,X_j=x,\ldots\right)}_{\text{odds ratio}}}$$

When an increase of  $X_j$  by one unit from x to x + 1, while keeping all other predictors fixed, it multiplies the odds by  $e^{\beta_j}$  (relative change from the odds when  $X_j = x$ );

e.g.

- ▶ If  $X_j = 0$  or 1, then for the group with  $X_j = 1$ , the odds of the event are  $e^{\beta_j}$  times that of the group with  $X_j = 0$ , with other values of  $X_{-j}$  fixed.
  - When  $\beta_j > 0$ , the group with  $X_j = 1$  has  $100(e^{\beta_j} 1)\%$  more odds than the group with  $X_j = 0$ , with other values of  $X_{-j}$  fixed.
  - ▶ When  $\beta_j < 0$ , then it is a decrease in the odds by  $100(1 e^{\beta_j})\%$ .

For intercept,  $e^{\beta_0} \div (1 + e^{\beta_0})$  is the probability of the event for the base group when all X's = 0.

## Inferences: logistic regression

Assuming correct model specification, by central limit theorem, the MLE estimator

$$\hat{\boldsymbol{\beta}} \to N\left(\boldsymbol{\beta}^*, \left(\mathbf{X}'\mathbf{W}(\boldsymbol{\beta}^*)\mathbf{X}\right)^{-1}\right).$$

Here,  $\boldsymbol{\beta}^*$  is the truth. The estimator for the variance of  $\hat{\boldsymbol{\beta}}$  is given by

$$v\hat{a}r(\hat{\boldsymbol{\beta}}) = \left(\mathbf{X}'\mathbf{W}(\hat{\boldsymbol{\beta}})\mathbf{X}\right)^{-1}.$$

So as  $n \to \infty$ ,

$$\frac{\hat{\beta}_j - \beta_j}{se(\hat{\beta}_j)} \sim N(0, 1),$$

where

$$se(\hat{\beta}_j) = \left[ \left( \mathbf{X}' \mathbf{W} \left( \hat{\boldsymbol{\beta}} \right) \mathbf{X} \right)^{-1} \right]_{j,j}^{1/2}$$

# Alternative formulation: logistic regression

Suppose that

$$Y_i = 1\{\beta_0 + X_i^{\top} \boldsymbol{\beta}_1 + \varepsilon_i\}$$

where  $\varepsilon_i$  is independent of  $X_i$ , following a logistic distribution (mean 0 and standard deviation  $\pi/\sqrt{3}$ ), aka, the c.d.f. of  $\varepsilon_i$  given by

$$F(\epsilon) = \frac{exp(\epsilon)}{1 + exp(\epsilon)}.$$

Then

$$P(Y_i = 1 | X_i = x) = \frac{exp(\beta_0 + x^\top \boldsymbol{\beta}_1)}{1 + exp(\beta_0 + x^\top \boldsymbol{\beta}_1)}.$$

If important predictors are omitted from the model (model misspecified), the usual interpretation linking the coefficients to (true) log-odds will break down.

Logistic Regression: multi-class case, sparsity

# Logistic Regression: multi-class case, sparsity

Suppose there are K groups. Let the K-th group be the base group. One may model  $Pr(Y = k|x; \beta_0, \beta)$  as

$$Pr(Y = k|x; \boldsymbol{\beta}_0, \boldsymbol{\beta}) = \frac{\exp\left(x^{\top} \boldsymbol{\beta}_k + \beta_{k0}\right)}{\sum_{k'=1}^{K} \exp\left(x^{\top} \boldsymbol{\beta}_{k'} + \beta_{k'0}\right)}$$

- $\triangleright \beta_0 := (\beta_{10}, \dots, \beta_{K0})$  the vector of K intercepts.
- $\beta := (\beta_1, \dots, \beta_K)^{\top}$ , a K by p matrix.

The **S** is the **softmax function**  $\mathbb{R}^K \mapsto (0,1)^K$ , defined as

$$\mathbf{S}(\boldsymbol{\eta})_k = \frac{e^{\eta_k}}{\sum_{k'}^K e^{\eta_{k'}}}, \quad k = 1, \dots, K; \quad \boldsymbol{\eta} = (\eta_1, \dots, \eta_K)^{\top}$$

$$Pr(Y = k|x; \boldsymbol{\beta}_0, \boldsymbol{\beta}) := Pr(Y = k|\boldsymbol{\eta}) = \mathbf{S}(\boldsymbol{\eta})_k$$

The multi-class logistic regression (or multinomial logistic regression) models K-1 logits:

- $\triangleright$  set  $\beta_{K0} = 0, \beta_K = 0$  to avoid overparametrization.
- $\triangleright$  treat K as the base group

$$\log \frac{\Pr(Y = 1 \mid X = x)}{\Pr(Y = K \mid X = x)} = \beta_{10} + \beta_1^{\top} x$$

$$\log \frac{\Pr(Y = 2 \mid X = x)}{\Pr(Y = K \mid X = x)} = \beta_{20} + \beta_2^{\top} x$$

$$\log \frac{\Pr(Y = K - 1 \mid X = x)}{\Pr(Y = K \mid X = x)} = \beta_{(K-1)0} + \beta_{K-1}^{\top} x$$

Note: if prediction is the only concern, we may fit a overparametrized model. This is the approach taken in the perceptron-based model.

Equivalently,

$$p_k(x) \equiv \Pr(Y = k \mid x) = \frac{\exp(\beta_{k0} + \beta_k^{\top} x)}{1 + \sum_{l=1}^{K-1} \exp(\beta_{l0} + \beta_l^{\top} x)} \text{ for } k = 1, \dots, K-1$$

$$p_K(x) \equiv \Pr(Y = K \mid x) = \frac{1}{1 + \sum_{l=1}^{K-1} \exp(\beta_{l0} + \beta_l^{\top} x)}$$

Clearly  $\sum_{k=1}^{K} p_k(x) = 1$ . The parameter vector

$$\boldsymbol{\theta} = \left\{ \beta_{10}, \boldsymbol{\beta}_1^{\top}, \dots, \beta_{(K-1)0}, \boldsymbol{\beta}_{K-1}^{\top} \right\}^{\top}$$

Let 
$$p_{k,i} := Pr(Y_i = k | X = x_i, \boldsymbol{\theta})$$
 and  $p_{y_i}(x_i; \boldsymbol{\theta}) = Pr(Y_i = y_i | X = x_i, \boldsymbol{\theta}).$ 

$$\ell(\theta) = \sum_{i=1}^{n} \log p_{y_i}(x_i; \theta) = \log \left( \prod_{i=1}^{n} \prod_{k=1}^{K} p_{k,i}^{1(y_i = k)} \right)$$
$$= \sum_{i=1}^{n} \sum_{k=1}^{K} 1(y_i = k) \log(p_{k,i})$$

Since  $\beta_{K0} := 0$  and  $\beta_K := 0$ , the log-likelihood:

$$\textstyle \sum_{i=1}^{n} \left\{ \beta_{y_i0} + \beta_{y_i}^\top x_i - \log \left[ 1 + \sum_{k=1}^{K-1} \exp \left( \beta_{k0} + \beta_k^\top x_i \right) \right] \right\}$$

The loss function, as a negative loglikelihood function, is called **multinomial deviance** (loss):

$$L(Y, p_Y(X)) = -\log p_Y(X; \theta) = -\sum_{k=1}^{K} 1(Y = k) \log(Pr(Y = k|X; \theta)).$$

## Compare logistic regression with LDA

### Logistic regression:

- Maximizing the conditional likelihood, the multinomial likelihood with probabilities  $Pr(Y = k \mid \mathbf{X})$
- ightharpoonup The marginal density Pr(X) is ignored (fully nonparametric)
  - **discriminative approach**: only modelling  $Pr(Y = k \mid x)$

#### LDA:

Maximizing the full log-likelihood based on the joint density

$$\Pr(X, Y = k) = \phi(X; \boldsymbol{\mu}_k, \Sigma) \, \pi_k$$

- ▶ Marginal density does play a role  $\Pr(\mathbf{X}) = \sum_{k} \pi_{k} \phi(X; \boldsymbol{\mu}_{k}, \Sigma)$ 
  - **generative approach**: modelling  $Pr(x \mid Y = k)$  (usually joint modelling of Pr(X, Y = k))

# Regularized logistic regression

The idea is to minimize the penalized negative likelihood function (binary response):

$$\min_{\beta_0, \boldsymbol{\beta}_1} \sum_{i=1}^n \left( -y_i \left( \beta_0 + \mathbf{x}_i^\top \boldsymbol{\beta}_1 \right) + \log \left( 1 + e^{\beta_0 + \mathbf{x}_i^\top \boldsymbol{\beta}_1} \right) \right) + \lambda J(\boldsymbol{\beta}_1)$$

The update is equivalent to solving the weighted LS till convergence (Iteratively Reweighted Least Squares Algorithm):

$$(\boldsymbol{\beta}_0^{(k+1)}, \boldsymbol{\beta_1}^{(k+1)}) = \arg\min_{\boldsymbol{\beta}} \left\{ (\mathbf{z}^{(k)} - \mathbf{X}\boldsymbol{\beta})^\top \mathbf{W}^{(k)} (\mathbf{z}^{(k)} - \mathbf{X}\boldsymbol{\beta}) + \lambda J(\boldsymbol{\beta}_1) \right\}.$$

For the Lasso penalty, it can be solved using coodinate descent.

# Sparse logistic regression

## **Algorithm** (Coordinate descent for sparse logistic regression):

Let 
$$\hat{\boldsymbol{\beta}}^{(0)} = (\hat{\beta}_0^{(0)}, \hat{\beta}_1^{(0)}, \dots, \hat{\beta}_p^{(0)})^T$$
  
For  $k = 0, 1, 2, \dots$ ,

P Compute  $p_1(x_i; \hat{\boldsymbol{\beta}}^{(k)}), z_i^{(k)}, w_{ii}^{(k)}, i = 1, \dots, n$ .

Let  $\beta_0 = \sum_i [w_{ii}^{(k)}(z_i^{(k)} - \sum_{l=1}^p \hat{\beta}_l^{(k)}x_{il})]/\sum_i w_{ii}^{(k)}, \quad \beta_l = \hat{\beta}_l^{(k)}, l = 1, \dots, p$ .

P for  $j = 1, \dots, p$  do

Compute  $r_{ij} = z_i^{(k)} - \beta_0 - \sum_{l \neq j} \beta_l x_{il}, i = 1, \dots, n$ 

Compute  $u_j^{(k)} = \sum_{i=1}^n w_{ii}^{(k)} r_{ij} x_{ij},$ 

Compute  $v_j^{(k)} = \sum_{i=1}^n w_{ii}^{(k)} x_{ij}^2,$ 

Compute  $\beta_j = \operatorname{soft}(u_j^{(k)}/v_j^{(k)}, \lambda/v_j^{(k)})$ 

P  $\hat{\beta}_0^{(k+1)} = \beta_0, \hat{\beta}_j^{(k+1)} = \beta_j, j = 1, \dots, p$ 

#### until convergence

Others: nonparametric classifier, robust loss, perceptrons

## **KNN**

For any given  $X = x_0$ , we find the K closest neighbors to  $X = x_0$  in the training data, and examine their corresponding Y:

$$P(Y = j \mid X = x_0) = \frac{1}{K} \sum_{i \in N_K(x_0)} 1(y_i = j)$$

Estimate the conditional probability for group j by the proportion out of the k neighbors that are in group j.

The smaller that K is the more flexible the method will be.

Note: more on nonparametric method (e.g., nonparametric logistic regression) in future lessons.

## Others: Alternative loss functions

For binary classification,

$$\min_{f \in \mathcal{F}_{\pm 1}} \mathcal{E}_{X,Y}(1(Y \neq f(X)))$$

where  $\mathcal{F}_{\pm 1}$  consists of function that maps to  $\{-1,1\}$ .

The ERM solution is

$$\hat{f}_n = \arg\min_{f \in \mathcal{F}_{\pm 1}} \frac{1}{n} \sum_{i=1}^n 1(y_i \neq f(x_i))$$

• the choice of  $\mathcal{F}_{\pm 1}$ : hard-classifier  $f \in \mathcal{F}_{\pm 1}$ 

 $\triangleright$  perceptron:  $\mathcal{F}_{\pm 1} = \{ \operatorname{sign}(\beta_0 + \boldsymbol{\beta}^\top x) : \beta_0, \boldsymbol{\beta} \}$ 

Equivalent formulation, but using  $y_i \in \{-1, 1\}$ ,

$$\hat{f}_n = \arg\min_{f \in \mathcal{F}_{\pm 1}} \frac{1}{n} \sum_{i=1}^n 1(-y_i f(x_i) > 0)$$

Using some smooth and convex surrogate function  $\psi(z)$  for 1(z > 0) and relaxing the class of functions  $\mathcal{F}$ , solve

$$\hat{f}_n = \arg\min_{f \in \mathcal{F}} \frac{1}{n} \sum_{i=1}^n L(y_i, f(x_i)) = \frac{1}{n} \sum_{i=1}^n \psi(-y_i f(x_i))$$

- ightharpoonup Choice of L(y, f):
  - Squared error:  $(y-f)^2 = (1-yf)^2$
  - ▶ Binomial deviance (logistic): log(1 + exp(-2yf))
  - ightharpoonup Exponential loss:  $\exp(-yf)$
  - $\triangleright$  SVM loss:  $(1 yf)_+$
- ▶ Choice of  $\mathcal{F}$ : soft-classifier  $f \in [-1, 1]$  or  $\mathbb{R}$ 
  - decision:  $\hat{Y}_i = \operatorname{sign}(\hat{f}_n(X_i))$

Let z = -yf > 0: indicate misclassification.

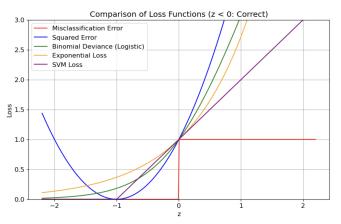
ightharpoonup Misclassification error: 1(z > 0)

▶ Squared error:  $(1+z)^2$ 

▶ Binomial deviance (scaled):  $\log(1 + \exp(2z))/\log(2)$ 

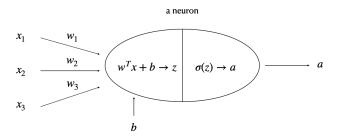
ightharpoonup Exponential loss:  $\exp(z)$ 

▶ SVM loss:  $(1+z)_+$ 



More in future lessons...

## Perceptron



input layer

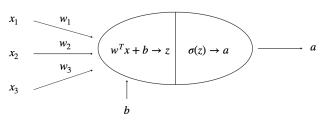
first hidden layer

Here,  $\sigma(\cdot)$  is an activation function: three simplest cases

- i. identity  $\sigma(v) = v$
- ii. indicator (Heaviside step)  $\sigma(v) = 1(v > 0)$
- iii. sigmoid  $\sigma(v) = 1/(1 + e^{-v})$

The output  $\hat{a}$  will be measured against the actual outcome value under some loss function.

#### a neuron



input layer

first hidden layer

## Formally,

$$\sigma\left(\sum_{j=1}^p x_j w_j + b\right) = \sigma\left(\boldsymbol{x}^{\top} \mathbf{w} + b\right), \qquad \boldsymbol{\beta} := (\boldsymbol{w}, b)$$

Note: OLS in regression is a special case with  $\sigma(\cdot)$  being identity function and the squared error loss (see Adaline).

## Rosenblatt's Perceptron

Rosenblatt's Perceptron:  $\sigma(v) = 1(v > 0)$ 

Let  $\beta$  include the b, x includes the intercept,  $\alpha>0$  small constant (learning rate). Start with random weight, then iteratively update the weight

- ▶ For k = 1, 2, ..., K:
  - ightharpoonup for each  $i = 1, \ldots, n$ :
    - ightharpoonup compute  $\hat{y}_i^{(k)} = 1 \left( \boldsymbol{x}_i^{\top} \boldsymbol{\beta}^{(k)} > 0 \right)$  (Heaviside)
    - ightharpoonup compute  $\mathbf{e}_i^{(k)} = y_i \hat{y}_i^{(k)}$
    - update  $\boldsymbol{\beta}^{(k+1)} = \boldsymbol{\beta}^{(k)} + \alpha \times \mathbf{e}_i^{(k)} \times \boldsymbol{x}_i$
- ▶ a **training epoch**: *one* loop over the whole training data
- $\triangleright$  the whole process repeats for K times or **epochs**

A training algorithm that updates the parameter after seeing one example is called **on-line training** (stochastic gradient descent).

## Compared with logistic regression

Recall logistic regression, the update at the k-th iteration

$$\boldsymbol{\beta}^{(k+1)} = \boldsymbol{\beta}^{(k)} + \alpha \times \mathbf{X}^{\top} (\mathbf{y} - \mathbf{p}^{(k)}) = \boldsymbol{\beta}^{(k)} + \alpha \times \sum_{i}^{n} (y_i - p(x_i; \boldsymbol{\beta}^{(k)})) \boldsymbol{x}_i$$

## The **on-line training** process is

- ▶ For k = 1, 2, ..., K:
  - - compute  $\hat{y}_i^{(k)} = p(x_i; \boldsymbol{\beta}^{(k)})$  (sigmoid)
    - $compute e_i^{(k)} = y_i \hat{y}_i^{(k)}$
    - $\beta^{(k+1)} = \beta^{(k)} + \alpha \times e_i^{(k)} \times x_i$

#### Note

- ▶ Perceptron uses hard-classifier, while logistic uses soft-classifier
- ▶ Perceptron algorithm does not converge if the data is not linearly separable
- ▶ SGD (WNLS) for logistic regression converges to global minimum of the binary entropy loss function (even when the data is not linearly separable)

# Adaline (ADAptive LInear NEuron)/OLS

Adaline for classification is a special case when  $\sigma(\cdot)$  is identity function and the loss is squared error loss.

$$MSE = \frac{1}{n} \sum_{i=1}^{n} (\hat{a}_i - y_i)^2, \qquad \hat{a}_i = \boldsymbol{x}_i^{\top} \hat{\beta}$$

It can be implemented using gradient descent for OLS.

$$\triangleright \boldsymbol{\beta}^{(k+1)} = \boldsymbol{\beta}^{(k)} + \alpha \times 2(y_i - \boldsymbol{x}_i^{\top} \hat{\boldsymbol{\beta}}^{(k)}) \times \boldsymbol{x}_i$$

The Adaline classifies based on the hard rule

$$\hat{y}_i = 1 \left( \boldsymbol{x}_i^{\top} \hat{\boldsymbol{\beta}} > 0.5 \right)$$

A major difference from Rosenblatt's Perceptron is that

- $\hat{a}_i$  is the continuous output (as opposed to  $\hat{y}_i$ ) that is measured against the true  $y_i$  in the loss calculation
  - ▶ That is, the hard-classification step is not backpropagated.