The Bayes classifier

Theorem

The classifier
$$h^*(\mathbf{x}) := \arg \max_k \eta_k(\mathbf{x})$$
 satisfies
 $R(h^*) = \min R(h)$

where the min is over all possible classifiers.

To calculate the Bayes classifier/Bayes risk, we need to know $\eta_k(\mathbf{x}) = \mathbb{P}[Y = k | X = \mathbf{x}]$

Alternatively, since $\pi_k f_{X|Y}(\mathbf{x}|k) = \eta_k(\mathbf{x}) f_X(\mathbf{x})$, to find the maximum $\eta_k(\mathbf{x})$ it is sufficient to know $\pi_k f_{X|Y}(\mathbf{x}|k)$

Linear discriminant analysis (LDA)

In linear discriminant analysis (LDA), we make the (strong) assumption that

$$X|Y = k \sim \mathcal{N}(\boldsymbol{\mu}_k, \boldsymbol{\Sigma})$$

for k = 0, ..., K - 1

Here $\mathcal{N}(\mu,\Sigma)$ is the multivariate Gaussian/normal distribution with mean μ and covariance matrix Σ

$$\phi(\mathbf{x}; \boldsymbol{\mu}, \boldsymbol{\Sigma}) = \frac{1}{(2\pi)^{d/2} |\boldsymbol{\Sigma}|^{1/2}} \exp\left\{-\frac{1}{2}(\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1}(\mathbf{x} - \boldsymbol{\mu})\right\}$$

Note: Each class has the same covariance matrix Σ

Example

Suppose that K = 2

$$d_M^2(\mathbf{x}; \widehat{\boldsymbol{\mu}}_0, \widehat{\boldsymbol{\Sigma}}) - 2\log \widehat{\pi}_0 \underset{1}{\overset{0}{\underset{1}{\leq}}} d_M^2(\mathbf{x}; \widehat{\boldsymbol{\mu}}_1, \widehat{\boldsymbol{\Sigma}}) - 2\log \widehat{\pi}_1$$

It turns out that by setting

$$\mathbf{w} = \widehat{\boldsymbol{\Sigma}}^{-1} (\widehat{\boldsymbol{\mu}}_1 - \widehat{\boldsymbol{\mu}}_0)$$
$$b = \frac{1}{2} \widehat{\boldsymbol{\mu}}_0^T \widehat{\boldsymbol{\Sigma}}^{-1} \widehat{\boldsymbol{\mu}}_0 - \frac{1}{2} \widehat{\boldsymbol{\mu}}_1^T \widehat{\boldsymbol{\Sigma}}^{-1} \widehat{\boldsymbol{\mu}}_1 + \log \widehat{\frac{\pi_1}{\pi_0}}$$

we can re-write this as

$$\mathbf{w}^T \mathbf{x} + b \underset{1}{\stackrel{0}{\leq}} \mathbf{0}$$
 linear classifier

Challenges for LDA

The generative model is rarely valid

Moreover, the number of parameters to be estimated is

- class prior probabilities: K 1
- means: Kd
- covariance matrix: $\frac{1}{2}d(d+1)$

If d is small and n is large, then we can accurately estimate these parameters (provably, using Hoeffding)

If n is small and d is large, then we have more parameters than observations, and will likely obtain very poor estimates

- first apply a dimensionality reduction technique to reduce $d \ \$ (more on this later)
- assume a more structured covariance matrix

Another possible escape

Recall from the very beginning of the lecture that the Bayes classifier can be stated either in terms of maximizing $\pi_k f_{X|Y}(\mathbf{x}|k)$ or $\eta_k(\mathbf{x})$

In LDA, we are estimating $\pi_k f_{X|Y}(\mathbf{x}|k)$, which is equivalent to the full joint distribution of (X, Y)

All we *really* need is to be able to estimate $\eta_k(\mathbf{x})$

- we don't need to know $f_X(\mathbf{x})$

LDA commits one of the cardinal sins of machine learning:

Never solve a more difficult problem as an intermediate step

Is there a better approach?

Another look at plugin methods

Suppose
$$K = 2$$

Define $\eta(\mathbf{x}) = \eta_1(\mathbf{x})$
 $= 1 - \eta_0(\mathbf{x})$

In this case, another way to express the Bayes classifier is as $h^*(\mathbf{x}) = \begin{cases} 1 & \text{if } \eta(\mathbf{x}) \ge 1/2 \\ 0 & \text{if } \eta(\mathbf{x}) < 1/2 \end{cases}$

Note that we do not actually need to know the full distribution of (X, Y) to express the Bayes classifier

All we really need is to decide if $\eta(\mathbf{x}) \geq 1/2$

Gaussian case

Suppose that K = 2 and that $X|Y = k \sim \mathcal{N}(\mu_k, \Sigma)$ $\eta(\mathbf{x}) = \frac{\pi_1 \phi(\mathbf{x}; \boldsymbol{\mu}_1, \boldsymbol{\Sigma})}{\pi_1 \phi(\mathbf{x}; \boldsymbol{\mu}_1, \boldsymbol{\Sigma}) + \pi_0 \phi(\mathbf{x}; \boldsymbol{\mu}_0, \boldsymbol{\Sigma})}$ $\pi_1 e^{-\frac{1}{2}(\mathbf{x}-\boldsymbol{\mu}_1)^T \Sigma^{-1}(\mathbf{x}-\boldsymbol{\mu}_1)}$ $\frac{1}{\pi_1 e^{-\frac{1}{2}(\mathbf{x}-\boldsymbol{\mu}_1)^T \boldsymbol{\Sigma}^{-1}(\mathbf{x}-\boldsymbol{\mu}_1)} + \pi_0 e^{-\frac{1}{2}(\mathbf{x}-\boldsymbol{\mu}_0)^T \boldsymbol{\Sigma}^{-1}(\mathbf{x}-\boldsymbol{\mu}_0)}}$ 1 $1 + \frac{\pi_0}{\pi_1} e^{\frac{1}{2} (\mathbf{x} - \boldsymbol{\mu}_1)^T \boldsymbol{\Sigma}^{-1} (\mathbf{x} - \boldsymbol{\mu}_1) - \frac{1}{2} (\mathbf{x} - \boldsymbol{\mu}_0)^T \boldsymbol{\Sigma}^{-1} (\mathbf{x} - \boldsymbol{\mu}_0)}$ $=\frac{1}{1+e^{-(\mathbf{w}^T\mathbf{x}+b)}}$

Logistic regression

This observation gives rise to another class of plugin methods, the most important of which is logisitic regression, which implements the following strategy

1. Assume
$$\eta(\mathbf{x}) = \frac{1}{1 + e^{-(\mathbf{w}^T \mathbf{x} + b)}}$$
 ($\mathbf{w} \in \mathbb{R}^d$, $b \in \mathbb{R}$)

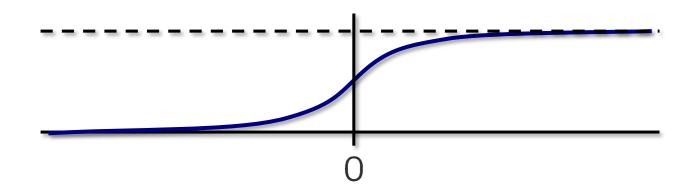
- 2. Directly estimate \mathbf{w}, b (somehow) from the data
- 3. Plug the estimate

$$\widehat{\eta}(\mathbf{x}) = \frac{1}{1 + e^{-(\widehat{\mathbf{w}}^T \mathbf{x} + \widehat{b})}}$$

into the formula for the Bayes classifier

The logistic function

The function $\frac{1}{1+e^{-t}}$ is called a *logistic* function (or a *sigmoid* function in other contexts)



The logistic regression classifier

Denote the logistic regression classifier by

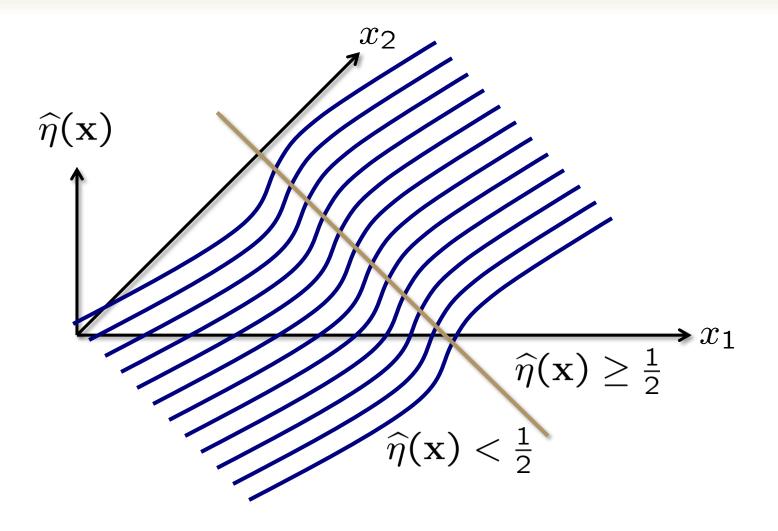
$$\widehat{h}(\mathbf{x}) = \mathbf{1}_{\left\{\widehat{\eta}(\mathbf{x}) \ge 1/2\right\}}(\mathbf{x})$$
Note that $\widehat{h}(\mathbf{x}) = \mathbf{1} \iff \widehat{\eta}(\mathbf{x}) \ge \frac{1}{2}$

$$\iff \frac{1}{1 + \exp\left(-(\widehat{\mathbf{w}}^T \mathbf{x} + \widehat{b})\right)} \ge \frac{1}{2}$$

$$\iff \exp\left(-(\widehat{\mathbf{w}}^T \mathbf{x} + \widehat{b})\right) \le \mathbf{1}$$

$$\iff (\widehat{\mathbf{w}}^T \mathbf{x} + \widehat{b}) \ge \mathbf{0}$$
So $\widehat{h}(\mathbf{x}) = \begin{cases} \mathbf{1} & \text{if } \widehat{\mathbf{w}}^T \mathbf{x} + b \ge \mathbf{0} \\ \mathbf{0} & \text{otherwise} \end{cases}$ linear classifier

Example



Estimating the parameters

Challenge: How to estimate the parameters for

$$\eta(\mathbf{x}) = \frac{1}{1 + e^{-(\mathbf{w}^T \mathbf{x} + b)}}$$

One possibility: $\mathbf{w} = \widehat{\boldsymbol{\Sigma}}^{-1} (\widehat{\boldsymbol{\mu}}_1 - \widehat{\boldsymbol{\mu}}_0)$
 $b = \frac{1}{2} \widehat{\boldsymbol{\mu}}_0^T \widehat{\boldsymbol{\Sigma}}^{-1} \widehat{\boldsymbol{\mu}}_0 - \frac{1}{2} \widehat{\boldsymbol{\mu}}_1^T \widehat{\boldsymbol{\Sigma}}^{-1} \widehat{\boldsymbol{\mu}}_1 + \log \frac{\widehat{\pi}_1}{\widehat{\pi}_0}$

Alternative: Maximum likelihood estimation

For convenience, let's let $\theta = (b, \mathbf{w})$

Note that $\eta(\mathbf{x})$ is really a function of both \mathbf{x} and $\boldsymbol{\theta}$, so we will use the notation $\eta(\mathbf{x}; \boldsymbol{\theta})$ to highlight this dependence

The *a posteriori* probability of our data

Suppose that we knew θ . Then we could compute

$$\mathbb{P}[y_i | \mathbf{x}_i; \boldsymbol{\theta}] = \mathbb{P}[Y_i = y_i | X_i = \mathbf{x}_i; \boldsymbol{\theta}]$$
$$= \begin{cases} \eta(\mathbf{x}_i; \boldsymbol{\theta}) & \text{if } y_i = 1\\ 1 - \eta(\mathbf{x}_i; \boldsymbol{\theta}) & \text{if } y_i = 0 \end{cases}$$

$$=\eta(\mathbf{x}_i; \boldsymbol{ heta})^{y_i}(1-\eta(\mathbf{x}_i; \boldsymbol{ heta}))^{1-y_i}$$

n

Because of independence, we also have that

$$\mathbb{P}[y_1,\ldots,y_n|\mathbf{x}_1,\ldots\mathbf{x}_n;oldsymbol{ heta}] = \prod_{i=1}^n \mathbb{P}[y_i|\mathbf{x}_i;oldsymbol{ heta}] \ = \prod_{i=1}^n \eta(\mathbf{x}_i;oldsymbol{ heta})^{y_i}(1-\eta(\mathbf{x}_i;oldsymbol{ heta}))^{1-y_i}$$

Maximum likelihood estimation

We don't actually know θ , but we do know y_1, \ldots, y_n

Suppose we view y_1, \ldots, y_n to be fixed, and view $\mathbb{P}[y_1, \ldots, y_n | \mathbf{x}_1, \ldots, \mathbf{x}_n; \boldsymbol{\theta}]$ as just a function of $\boldsymbol{\theta}$

When we do this, $\mathcal{L}(\theta) = \mathbb{P}[y_1, \dots, y_n | \mathbf{x}_1, \dots, \mathbf{x}_n; \theta]$ is called the *likelihood* (or likelihood function)

The method of *maximum likelihood* aims to estimate θ by finding the θ that *maximizes* the *likelihood* $\mathcal{L}(\theta)$

In practice, it is often more convenient to focus on maximizing the *log-likelihood*, i.e., $\log \mathcal{L}(\theta)$

The log-likelihood

To see why, note that the likelihood in our case is given by

$$\mathcal{L}(\boldsymbol{ heta}) = \prod_{i=1}^n \eta(\mathbf{x}_i; \boldsymbol{ heta})^{y_i} (1 - \eta(\mathbf{x}_i; \boldsymbol{ heta}))^{1-y_i}$$

Thus, the log-likelihood is given by

$$\ell(\boldsymbol{\theta}) = \log \mathcal{L}(\boldsymbol{\theta})$$

= $\sum_{i=1}^{n} y_i \log \eta(\mathbf{x}_i; \boldsymbol{\theta}) + (1 - y_i) \log(1 - \eta(\mathbf{x}_i; \boldsymbol{\theta}))$

It is often easier to work with summations instead of products, and since the log is a monotonic transformation, maximizing $\ell(\theta)$ is equivalent to maximizing $\mathcal{L}(\theta)$

Simplifying the log-likelihood

Notation

- $\widetilde{\mathbf{x}} = [1, x(1), \dots, x(d)]^T$
- $\boldsymbol{\theta} = [b, w(1), \dots, w(d)]^T$
- This means that $\mathbf{w}^T \mathbf{x} + b = \boldsymbol{\theta}^T \widetilde{\mathbf{x}}$, which lets us write

$$\eta(\mathbf{x}_i; oldsymbol{ heta}) = rac{1}{1 + e^{-oldsymbol{ heta}^T ilde{\mathbf{x}}_i}}$$

Thus, if we let $g(t) = \frac{1}{1+e^{-t}}$, then we can write

$$\ell(\boldsymbol{\theta}) = \sum_{i=1}^{n} y_i \log g(\boldsymbol{\theta}^T \widetilde{\mathbf{x}}_i) + (1 - y_i) \log(1 - g(\boldsymbol{\theta}^T \widetilde{\mathbf{x}}_i))$$

Simplifying the log-likelihood

Facts:

$$\log g(t) = \log \left(\frac{1}{1+e^{-t}}\right) = -\log(1+e^{-t})$$

$$\log(1 - g(t)) = \log\left(1 - \frac{1}{1 + e^{-t}}\right)$$
$$= \log\left(\frac{e^{-t}}{1 + e^{-t}}\right) = -t - \log(1 + e^{-t})$$
$$= \log\left(\frac{1}{1 + e^{t}}\right) = -\log(1 + e^{t})$$

Simplifying the log-likelihood

Facts:

$$\log g(t) = -\log(1 + e^{-t})$$

$$\log(1 - g(t)) = -t - \log(1 + e^{-t}) = -\log(1 + e^{t})$$
Thus
$$\ell(\theta) = \sum_{i=1}^{n} y_i \log g(\theta^T \tilde{\mathbf{x}}_i) + (1 - y_i) \log(1 - g(\theta^T \tilde{\mathbf{x}}_i))$$

$$= \sum_{i=1}^{n} -y_i \log(1 + e^{-\theta^T \tilde{\mathbf{x}}_i}) - \log(1 + e^{\theta^T \tilde{\mathbf{x}}_i})$$

$$-y_i(-\theta^T \tilde{\mathbf{x}}_i - \log(1 + e^{-\theta^T \tilde{\mathbf{x}}_i}))$$

$$= \sum_{i=1}^{n} y_i \theta^T \tilde{\mathbf{x}}_i - \log(1 + e^{\theta^T \tilde{\mathbf{x}}_i})$$

Maximizing the log-likelihood

How can we maximize

$$\ell(\boldsymbol{\theta}) = \sum_{i=1}^{n} y_i \boldsymbol{\theta}^T \widetilde{\mathbf{x}}_i - \log(1 + e^{\boldsymbol{\theta}^T \widetilde{\mathbf{x}}_i})$$

with respect to
$$\theta$$
?

Find a
$$\theta$$
 such that $\nabla \ell(\theta) = \begin{bmatrix} \frac{\partial \ell(\theta)}{\partial \theta_1} \\ \vdots \\ \frac{\partial \ell(\theta)}{\partial \theta_{d+1}} \end{bmatrix} = 0$

(i.e., compute the partial derivatives and set them to zero)

Computing the gradient

It is not too hard to show that

$$egin{split}
abla \ell(oldsymbol{ heta}) &= \sum_{i=1}^n
abla \left(y_i oldsymbol{ heta}^T \widetilde{\mathbf{x}}_i - \log(1+e^{oldsymbol{ heta}^T \widetilde{\mathbf{x}}_i})
ight) \ &= \sum_{i=1}^n y_i \widetilde{\mathbf{x}}_i - \widetilde{\mathbf{x}}_i e^{oldsymbol{ heta}^T \widetilde{\mathbf{x}}_i} (1+e^{oldsymbol{ heta}^T \widetilde{\mathbf{x}}_i})^{-1} \ &= \sum_{i=1}^n \widetilde{\mathbf{x}}_i (y_i - g(oldsymbol{ heta}^T \widetilde{\mathbf{x}}_i)) = 0 \end{split}$$

This gives us d + 1 equations, but they are **nonlinear** and have no closed-form solution

Optimization

Throughout signal processing and machine learning, we will very often encounter problems of the form

 $\underset{\mathbf{x} \in \mathbb{R}^{d}}{\mathsf{minimize}} f(\mathbf{x})$

(or minimize $-\ell(\boldsymbol{\theta})$ for today) $_{\boldsymbol{\theta}\in\mathbb{R}^{d+1}}$

In many (most?) cases, we cannot compute the solution simply by setting $\nabla f(\mathbf{x}) = 0$ and solving for \mathbf{x}

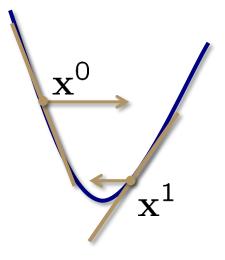
However, there are many powerful algorithms for finding \mathbf{x} using a computer

Gradient descent

A simple way to try to find the minimum of our objective function is to iteratively *"roll downhill"*

From \mathbf{x}^{0} , take a step in the direction of the negative gradient

$$\begin{aligned} \mathbf{x}^{1} &= \mathbf{x}^{0} - \alpha_{0} \nabla f(\mathbf{x})|_{\mathbf{x} = \mathbf{x}^{0}} & \alpha_{0} : \text{``step size''} \\ \mathbf{x}^{2} &= \mathbf{x}^{1} - \alpha_{1} \nabla f(\mathbf{x})|_{\mathbf{x} = \mathbf{x}^{1}} \end{aligned}$$



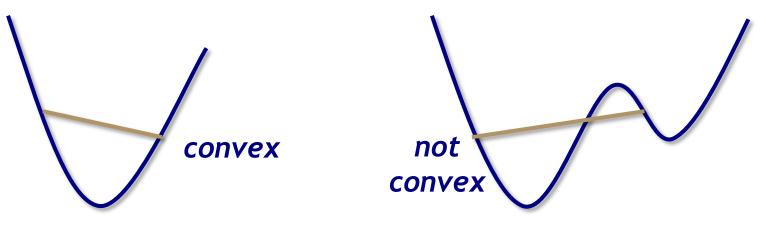
Convergence of gradient descent

The core iteration of gradient descent is to compute

$$\mathbf{x}^{j+1} = \mathbf{x}^j - \alpha_j \nabla f(\mathbf{x})|_{\mathbf{x} = \mathbf{x}^j}$$

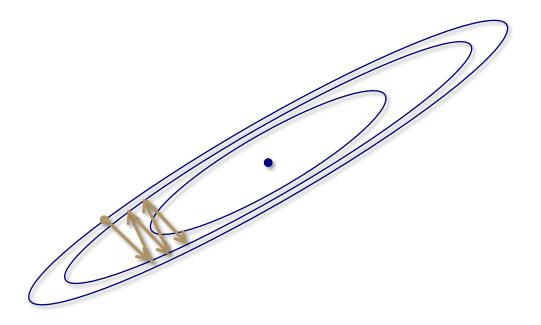
Note that if $\nabla f(\mathbf{x})|_{\mathbf{x}=\mathbf{x}^{j}} = 0$, then we have found the minimum and $\mathbf{x}^{j+1} = \mathbf{x}^{j}$, so the algorithm will terminate

If f is convex and sufficiently smooth, then gradient descent (with a fixed step size α) is guaranteed to converge to the global minimum of f



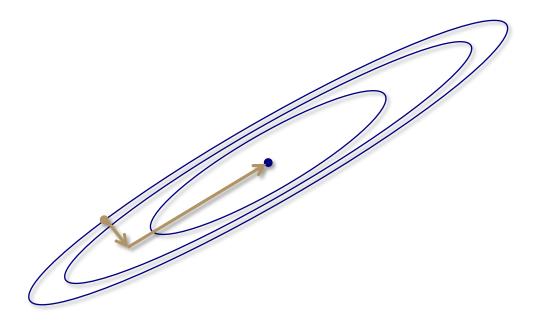
Step size matters!

Even though gradient descent provably converges, it can potentially take a while



Step size matters!

Even though gradient descent provably converges, it can potentially take a while



Newton's method

Also known as the Newton-Raphson method, this approach can be viewed as simply using the second derivative to automatically select an appropriate step size

$$\mathbf{x}^{j+1} = \mathbf{x}^{j} - \left(\nabla^{2}f(\mathbf{x})\right)^{-1}\nabla f(\mathbf{x})|_{\mathbf{x}=\mathbf{x}^{j}}$$

Hessian
matrix
$$\nabla^{2}f(\mathbf{x}) = \begin{bmatrix} \frac{\partial^{2}f}{\partial x_{1}^{2}} & \frac{\partial^{2}f}{\partial x_{1}\partial x_{2}} & \cdots & \frac{\partial^{2}f}{\partial x_{1}\partial x_{d}}\\ \frac{\partial^{2}f}{\partial x_{2}\partial x_{1}} & \frac{\partial^{2}f}{\partial x_{2}^{2}} & \cdots & \frac{\partial^{2}f}{\partial x_{2}\partial x_{d}}\\ \vdots & \vdots & \ddots & \vdots\\ \frac{\partial^{2}f}{\partial x_{d}\partial x_{1}} & \frac{\partial^{2}f}{\partial x_{d}\partial x_{2}} & \cdots & \frac{\partial^{2}f}{\partial x_{d}^{2}} \end{bmatrix}$$

Optimization for logistic regression

The negative log-likelihood in logistic regression is a convex function

Both gradient descent and Newton's method are common strategies for setting the parameters in logistic regression

Newton's method is much faster when the dimension d is small, but is impractical when d is large

Why?

More on this on next week's homework

Comparison of plugin methods

Naïve Bayes, LDA, and logistic regression are all *plugin methods* that result in *linear* classifiers

Naïve Bayes

- plugin method based on density estimation
- scales well to high-dimensions and naturally handles mixture of discrete and continuous features

Linear discriminant analysis

- better if Gaussianity assumptions are valid

Logistic regression

- models only the distribution of Y|X, not(X, Y)
- valid for a larger class of distributions
- fewer parameters to estimate

Beyond plugin methods

Plugin methods can be useful in practice, but ultimately they are very limited

- There are always distributions where our assumptions are violated
- If our assumptions are wrong, the output is totally unpredictable
- Can be hard to verify whether our assumptions are right
- Require solving a more difficult problem as an intermediate step

For most of the remainder of this course will focus on *nonparametric* methods that avoid making such strong assumptions about the (unknown) process generating the data