# **CS4220: Machine Learning**

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#### 1. Bayesian Classifier Theorem

#### 1.1. Decision Boundary

$$p(x|w_1)p(w_1) \ge p(x|w_2)p(x|w_2)$$
 (1)

#### 1.2. Minimizing the Classification Error Probability (Bayes Error)

$$P_e = p(w_2) \int_{-\infty}^{x_0} p(x|w_2) dx + p(w_1) \int_{x_0}^{\infty} p(x|w_1) dx$$
 (2)

where  $p(x_0|w_1)p(w_1) = p(x_0|w_2)p(x|w_2)$ 

#### 1.3. Minimizing the Average Risk

Loss Matrix

$$L = \begin{bmatrix} 0 & \lambda_{12} \\ \lambda_{21} & 0 \end{bmatrix} \tag{3}$$

Risk Function

$$r = \lambda_{21} p(w_2) \int_{-\infty}^{x_0} p(x|w_2) dx + \lambda_{12} p(w_1) \int_{x_0}^{\infty} p(x|w_1) dx$$
 (4)

#### 1.4. Gaussian pdf in the l-dimensional space

$$p(x) = \frac{1}{(2\pi)^{l/2} |\Sigma|^{l/2}} exp(-\frac{1}{2}(x-\mu)^T \Sigma^{-1}(x-\mu))$$
 (5)

where  $\Sigma = E[(x - \mu)(x - \mu)^T]$ 

#### 1.5. Bayesian Classifier

discriminant functions

$$g_i(x) = ln(p(x|w_i)p(w_i)) = lnp(x|w_i) + lnp(w_i)$$
(6)

Normally Distributed Classifier

$$g_i(x) = -\frac{1}{2}x^T \Sigma_i^{-1} x + \frac{1}{2}x^T \Sigma_i^{-1} \mu_i - \frac{1}{2}\mu_i^T \Sigma_i^{-1} \mu_i + \frac{1}{2}\mu_i^T \Sigma_i^{-1} x + lnp(w_i) + c_i$$
 (7)

where  $c_i = -(l/2)ln2\pi - (1/2)ln(det(\Sigma_i))$ 

- if l=2, corelation=0

$$g_i(x) = -\frac{1}{2\sigma_i^2}(x_1^2 + x_2^2) + \frac{1}{\sigma_i^2}(\mu_i 1x_1 + \mu_i 2x_2) - \frac{1}{2\sigma_i^2}(\mu_{i1}^2 + \mu_{i2}^2) + lnp(w_i) + c_i$$
 (8)

 $g_i(x) - g_j(x) = 0$  are quadratics (i.e., ellipsoids, parabolas, hyperbolas, pairs of lines)

- if covariance matrix is the same in all classes

$$g_i(x) = w_i^T x + b (9)$$

where  $w_i = \Sigma^{-1}\mu_i$  and  $b = lnp(w_i) - \frac{1}{2}\mu_i^T\Sigma^{-1}\mu_i$ 

- if Diagonal covariance matrix with equal elements ( $\Sigma = \sigma^2 I$ )

$$g_i(x) = \frac{1}{\sigma^2} \mu_i^T x + b \tag{10}$$

#### 2. ESTIMATION OF UNKNOWN PROBABILITY DENSITY FUNCTIONS

#### 2.1. ML

we considered  $\theta$  as an unknown parameter.

$$\hat{\theta}_{ML} = \arg \max_{\theta} \prod_{k=1}^{N} p(x_k; \theta)$$
(11)

ML estimate of  $\sigma^2$ 

$$\hat{\sigma}_{ML}^2 = \frac{1}{N} \sum_{k=1}^{N} (x_k - \mu)^2 \tag{12}$$

ML estimate of  $\mu$ 

$$\hat{\mu}_{ML} = \frac{1}{N} \sum_{k=1}^{N} x_k \tag{13}$$

# 2.2. MAP

we considered  $\theta$  as a random vector.

$$p(\theta|X) = \frac{p(\theta)p(X|\theta)}{P(X)} \tag{14}$$

then,

$$\hat{\theta}_{MAP}: \frac{\partial}{\partial \theta} p(\theta|X) = 0 \text{ or } \frac{\partial}{\partial \theta} p(X|\theta) p(\theta) = 0$$
 (15)

# 2.3. Bayesian Inference

Given the set X of the N training vectors and the *a priori information* about the pdf  $p(\theta)$ , the goal is to compute the conditional pdf p(x|X).

$$p(x|X) = \int p(x|\theta)p(\theta|X)d\theta \tag{16}$$

with

$$p(\theta|X) = \frac{p(X|\theta)p(\theta)}{p(X)} \tag{17}$$

$$p(X|\theta) = \prod_{k=1}^{N} p(x_k|\theta)$$
(18)

# 3. Normal-based Classifier: Quadratic Discriminant, Linear Discriminant and Nearest Mean

Let's assume that we have two classes:

# 3.1. Quadratic Discriminant

by eq.7, the quadratic classifier,

$$f(x) = x^T W x + w^T + w_0 (19)$$

with

$$W = \frac{1}{2}(\Sigma_2^{-1} - \Sigma_1^{-1}) \tag{20}$$

$$w = \mu_1^T \Sigma_1^{-1} - \mu_2^T \Sigma_2^{-1} \tag{21}$$

$$w_0 = -\frac{1}{2}ln(det(\Sigma_1)) - \frac{1}{2}\mu_1^T \Sigma_1^{-1} \mu_1 + lnp(y_1) + \frac{1}{2}ln(det(\Sigma_2)) + \frac{1}{2}\mu_2^T \Sigma_2^{-1} \mu_2 - lnp(y_2)$$
(22)

(i.e., ellipsoids, parabolas, hyperbolas, pairs of lines)

#### 3.2. Linear Discriminant

by eq.9, the linear classifier,

$$f(x) = w^T x + w_0 (23)$$

with

$$w = \Sigma^{-1}(\mu_2 - \mu_1) \tag{24}$$

$$w_0 = \frac{1}{2}\mu_2^T \Sigma^{-1} \mu_2 - \frac{1}{2}\mu_1^T \Sigma^{-1} \mu_1 + \ln \frac{p(y_1)}{p(y_2)}$$
 (25)

Therefore, the Linear Discriminant has a strong assumption on that  $\Sigma_1 = \Sigma_2$ 

#### 3.3. Nearest Mean Classifier

by eq.10, the nearest mean classifier,

$$f(x) = w^{T}x + w_0 = \sigma^2(g_1(x) - g_2(x))$$
(26)

with

$$w = \mu_2 - \mu_1 \tag{27}$$

$$w_0 = \frac{1}{2}\mu_2^T \mu_2 - \frac{1}{2}\mu_1^T \mu_1 + \sigma^2 \ln \frac{p(y_1)}{p(y_2)}$$
(28)

Therefore, the Nearest Mean has a **strong assumption** on mutually uncorrelated and of the same variance  $(\Sigma_1 = \Sigma_2 = \sigma^2 I)$ 

#### 4. More Parametric Classifiers

#### 4.1. Logistic Classifier

$$\begin{cases} p(y_1|x) = \frac{1}{e^{-(w^T x + w_0)} + 1} \\ p(y_2|x) = \frac{1}{e^{(w^T x + w_0)} + 1} \end{cases}$$
 (29)

Maximize Log Likelihood

$$lnp(y|x) = \sum_{i=1}^{N} ln(\frac{1}{e^{-y_i(w^T x_i + w_0)} + 1})$$
(30)

#### 4.2. Fisher Classifier

$$y_i = w^T x_i \begin{cases} \ge 0 & \text{if class 1} \\ < 0 & \text{if class 2} \end{cases}$$
 (31)

Minimize Square Loss

$$L(w) = \sum_{i=1}^{N} (w^{T} x_{i} - y_{i})^{2}$$
(32)

#### 4.3. The Perception

Forward&backward propagation and update w

$$w \leftarrow w + \eta y x \tag{33}$$

# 5. Non-parametric Classification

In the above sections, all the classifiers are based on **Parametric Classification** method, more precisely, based on normal distribution and Bayes Theorem. In this section, it will mainly demonstrate two non-parametric classifiers - Parzen Classifier and Nearest Neighbour Classifier (Both methods are sensitive to the scaling of the features).

#### 5.1. Parametric vs. Non-parametric

- Parametric: Assumptions can greatly simplify the learning process, but can also limit what can be learned. Algorithms that simplify the function to a known form are called parametric machine learning algorithms.
- Non-parametric: Algorithms that do not make strong assumptions about the form
  of the mapping function are called non-parametric machine learning algorithms.
  By not making assumptions, they are free to learn any functional form from the
  training data.

#### 5.2. Histogram method

it's easy, so I just skip it.

#### 5.3. Parzen Density Estimation

- Kernel

$$f(x) = \begin{cases} 0 & \text{if } |\mathbf{r}| > \mathbf{h} \\ \frac{1}{V} & \text{if } |\mathbf{r}| \le \mathbf{h} \end{cases}$$
 (34)

- Parzen Classifier

$$p(z|h) = \frac{1}{n} \sum_{i=1}^{n} K(||z - x_i||, h)$$
(35)

- Parzen plugs in the Gaussian density:

$$p(x|w_i) = \frac{1}{n_i} \sum_{i=1}^{n_i} N(x|x_j^{(i)}, hI)$$
(36)

#### 5.4. Nearest Neighbour Classification

$$p(x) = p(x|w_m)p(w_m) = \frac{k_m}{n_m V_k} \cdot \frac{n_m}{n}$$
(37)

where  $V_k$  is the volume of the sphere centered at x with radius r (the distance to the k-th nearest neighbor)

# 6. More Non-parametric Classifiers

GitHub Markdown

#### 6.1. SVM

By putting some constraints on the linear classifier, the VC dimension can be reduced. Why do that? Ans: When h is small, the true error is close to the apparent error

$$\begin{cases} w^{T} x_{i} + b \ge +1 & \text{for } y_{i} = +1 \\ w^{T} x_{i} + b \le -1 & \text{for } y_{i} = -1 \end{cases}$$
 (38)

Core Idea of SVM: Find the decision boundary, while maximize the margin

1. The above two equation can be merged into one

$$y_i(w^T x_i + b) - 1 \ge 0 (39)$$

2. The distance between the two boundaries

$$maximize \frac{2}{||w||} \to minimize \frac{1}{2}||w||^2 \tag{40}$$

3. by Lagrange Multiplier

$$L = \frac{1}{2}||w||^2 - \sum \alpha_i [y_i(wx_i + b) - 1]$$
(41)

4. by 
$$\frac{\partial L}{\partial w} = 0$$

$$w = \sum \alpha_i y_i x_i \tag{42}$$

5. by 
$$\frac{\partial L}{\partial b} = 0$$
 
$$\sum \alpha_i y_i = 0 \tag{43}$$

6. put eq.42 back to eq.41

$$L = \sum_{i} \alpha_{i} - \frac{1}{2} \sum_{i} \sum_{j} \alpha_{i} \alpha_{j} y_{i} y_{j} x_{i} \cdot x_{j}$$

$$\tag{44}$$

7. put eq.42 back to decision rule

$$\begin{cases} \sum \alpha_i y_i x_i \cdot \mathbf{u_i} + b - 1 \ge 0 & \text{Then, +} \\ \sum \alpha_i y_i x_i \cdot \mathbf{u_i} + b - 1 \le 0 & \text{Then, -} \end{cases}$$
(45)

8. kernelize

$$K(x_i, x_j) = \phi(x_i) \cdot \phi(x_j) \tag{46}$$

## 6.2. Agglomerative Hierarchical Clustering

- 1. Determine distances between all clusters
  - Two nearest objects in the clusters: single linkage
  - Two most remote objects in the clusters: complete linkage
  - Cluster centers: average linkage
- 2. Merge clusters that are closes
- 3. IF #clusters>1 THEN GOTO 1
- Dendrogram: Cut at "largest jump"→ Clustering
- Fusion Graph: Cut at "largest drop" → Clustering

## 7. Regression

#### 7.1. Intuitively Understanding

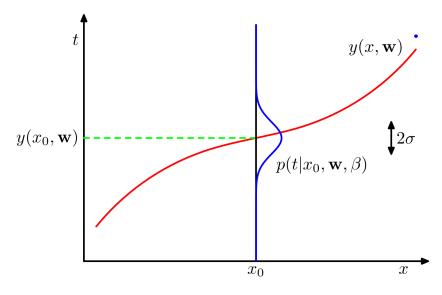


Figura 1. Regression Curve

Suppose that there is a fixed  $x_0$  and a bunch of choices of  $\theta$  (here are  $\mu$  and  $\sigma^2(\beta^{-1})$ ). If we expect the predicted value  $\hat{y}(x_0, w)$  to locate on the true value  $t_0$ , the probability of this occurrence  $p(t_0|x_0, w, \beta)$  should be maximized.

# 7.2. Maximum Likelihood Regression

$$p(t|x, w, \beta) = \prod_{n=1}^{N} N(t_n|y(x_n, w), \beta^{-1})$$
(47)

Log Likelihood function

$$lnp(t|x, w, \beta) = -\frac{\beta}{2} \sum_{n=1}^{N} y(x_n, w) - t_n^2 + \frac{N}{2} ln\beta - \frac{N}{2} ln(2\pi)$$
 (48)

To solve

$$\frac{\partial}{\partial w} lnp(t|x_0, w, \beta) = 0 \tag{49}$$

and

$$\frac{\partial}{\partial \beta^{-1}} lnp(t|x_0, w, \beta) = 0 \tag{50}$$

If we use linear regression, the solutions of eq.49 and eq.50

$$w_{ML} = (X^T X)^{-1} X^T Y \beta_{ML}^{-1} = \frac{1}{N} (w_{ML}^T X - Y)$$
 (51)

## 7.3. Max a Posterior Regression

Suppose that we have some knowledge about  $w N(0, \alpha I)$ 

$$w_{MAP}: (\prod_{i=1}^{N} p(y_i|w^T x_i, \sigma^2)) p(w|0, \alpha I)$$
 (52)

$$\frac{\partial}{\partial w} \left( \prod_{i=1}^{N} p(y_i | w^T x_i, \sigma^2) \right) p(w|0, \alpha I) = 0 \to w_{MAP} = \left( X^T X + \frac{\sigma^2}{\alpha} I \right)^{-1} X^T Y \tag{53}$$

## 8. Regularization

# 8.1. Keep Eigenvalues Away From Zero

Add identity to  $XX^T$ 

$$\hat{w} = (X^T X + \lambda I)^{-1} X^T Y \tag{54}$$

#### 8.2. LASSO, L1 Norm

$$min_{\beta} \frac{1}{N} \sum_{i=1}^{n} (y_i - w^T x_i)^2 + \lambda ||w||$$
 (55)

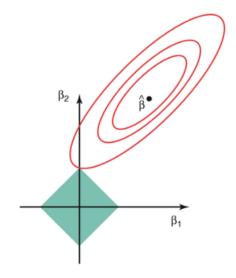


Figura 2. LASSO/L1 Regularization

$$\lambda \propto \frac{1}{ au}$$

# 8.3. Ridge, L2 Norm

$$min_{\beta} \frac{1}{N} \sum_{i=1}^{n} (y_i - w^T x_i)^2 + \lambda ||w||^2$$
 (56)

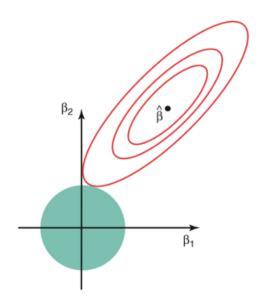


Figura 3. Ridge/L2 Regularization

$$\lambda \propto \frac{1}{ au}$$

# 8.4. L1 vs. L2

• L1 is for feature selection

• L2 is for avoiding overfitting

## Read More

# 9. Curves

# 9.1. Bias-Variance Decomposition

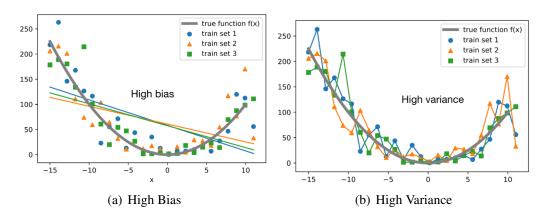


Figura 4. High Bias vs. High Variance

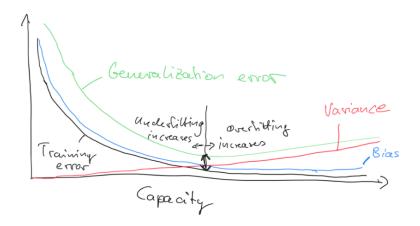


Figura 5. Decomposition Of Loss

### 9.2. Cross Validation Curve

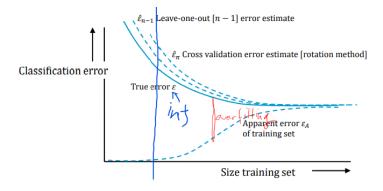


Figura 6. Cross Validation Curve

# 9.3. Learning Curve

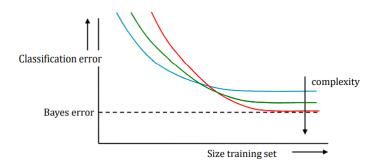


Figura 7. Learning Curve

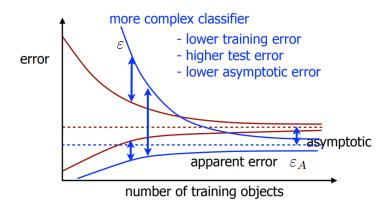


Figura 8. Learning Curve 2

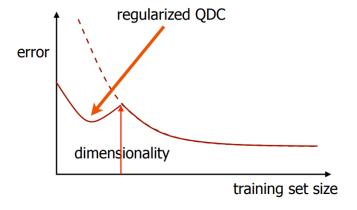


Figura 9. Regularized QDC

# 9.4. Feature Curve

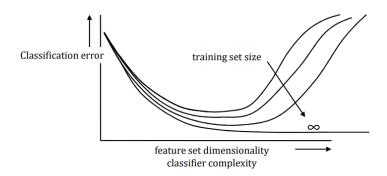


Figura 10. Feature Curve

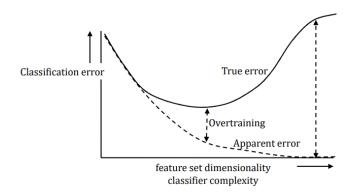


Figura 11. Curse of Dimensionality