

COSC6328.3
Speech & Language Processing



No.3

Pattern Classification (I)

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Outline

- **Pattern Classification Problems**
- **Bayesian Decision Theory**
- **How to build an optimal classifier?**
 - **Minimum-Error-Rate Classification**
 - **Maximum *a posterior* (MAP) decision rule**
 - **Model estimation: maximum likelihood, Bayesian learning, discriminative training, etc.**
- **Design classifier based on Discriminant**
 - **Linear-Discriminant-Functions-based classifier**
 - **Neural Networks**
 - **Support vector machines (SVM), large-margin classifiers**

Pattern Classification Problem

- Given a fixed set of finite number of classes: $\omega_1, \omega_2, \dots, \omega_N$.
- We have an unknown pattern/object P without its class identity.
- BUT we can measure (observe) some feature(s) X about P :
 - X is called feature or observation of the pattern P .
 - X can be scalar or vector or vector sequence
 - X can be continuous or discrete
- Pattern classification problem: $X \rightarrow \omega_i$
 - Determine the class identity for any a pattern based on its observation or feature.
- Fundamental issues in pattern classification
 - How to make an optimal classification?
 - In what sense is it optimal?

Examples of pattern classification(I)

- Speech recognition:
 - Pattern: voice spoken by a human being
 - Classes: language words/sentences used by the speaker
 - Features: speech signal characteristics measured by a microphone \rightarrow a sequence of feature vectors
 - Each vector: continuous, high-dimensional, real-valued numbers
- Natural language understanding:
 - Pattern: written or spoken languages of human
 - Classes: all possible semantic meanings or intentions
 - Features: the used words or word-sequences (sentences)
 - Discrete, scalars or vector

Examples of pattern classification(II)

- Image understanding:
 - Pattern: given images
 - Classes: all known object categories
 - Features: color or gray scales in all pixels
 - Continuous, multiple vectors/matrix
 - Examples: face recognition, OCR (optical character recognition).
- Gene finding in bioinformatics:
 - Pattern: a newly sequenced DNA sequence
 - Classes: all known genes
 - Features: all nucleotides in the sequence
 - Discrete; 4 types (adenine, guanine, cytosine, thymine)
- Protein classification in bioinformatics:
 - Pattern: protein primary 1-D sequence
 - Classes: all known protein families or domains
 - Features: all amino acids in the sequence: discrete; 20 types

Bayesian Decision Theory(I)

- Bayesian decision theory is a fundamental statistical approach to all pattern classification problems.
- Pattern classification problem is posed in probabilistic terms.
 - Observation X is viewed as random variables (vectors,...)
 - Class id ω is treated as a discrete random variable, which could take values $\omega_1, \omega_2, \dots, \omega_N$.
 - Therefore, we are interested in the joint probability distribution of X and ω which contains all info about X and ω .

$$p(X, \omega) = p(\omega) \cdot p(X | \omega)$$

- If all the relevant probability values and densities are known in the problem (we have complete knowledge of the problem), Bayesian decision theory leads to the optimal classification
 - Optimal \rightarrow Guarantee minimum average classification error
 - The minimum classification error is called the Bayes error.

Bayesian Decision Theory(II)

- In pattern classification, all relevant probabilities mean:
 - Prior probabilities of each class $P(\omega_i)$ ($i=1,2,\dots,N$): how likely any a pattern comes from each class before observing any features \rightarrow prior knowledge from previous experience
 - All priors sum to one: $\sum_{i=1}^N P(\omega_i) = 1$
 - Class-conditional probability density functions of the observed feature, X , $p(X | \omega_i)$ ($i=1,2,\dots,N$): how the feature distributes for all patterns belonging to one class ω_i .
 - If X is continuous, $p(X | \omega_i)$ is a continuous p.d.f. distribution
For every class ω_i : $\int_X p(X | \omega_i) \cdot dX = 1$
 - If X is discrete, $p(X | \omega_i)$ is discrete probability mass function (p.m.f.) distribution. For every class ω_i : $\sum_X p(X | \omega_i) = 1$

Example of class-conditional p.d.f.

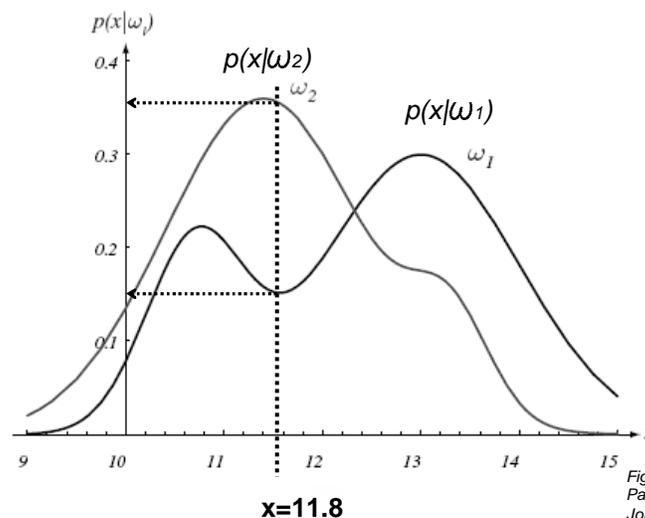


Figure from Duda et. al.,
Pattern classification
John Wiley & Sons®, Inc.

Bayes Decision Rule: Maximum *a posterior* (MAP) (I)

- If not observe any feature of an incoming unknown pattern P , classify it based on prior knowledge only

$$\omega_p = \arg \max_{\omega_i} P(\omega_i)$$

- Roughly guess it as the class with largest prior probability

- If observe some features X of the unknown patten P , we can convert the prior probability $P(\omega_i)$ into a posterior probability based on the Bayes theorem:

$$\text{posterior} = \frac{\text{prior} \times \text{likelihood}}{\text{evidence}}$$

Bayes decision rule: Maximum *a posterior* (MAP) (II)

- Where
 - Prior $P(\omega_i)$: probability of receiving a pattern from class ω_i before observing anything. (prior knowledge)
 - Likelihood $p(X | \omega_i)$: probability of observing feature X if assume X comes from a pattern in class ω_i . (if assume X is given, treat it as a function of ω_i , it is called likelihood function)
 - Posterior $p(\omega_i | X)$: probability of getting a pattern from class ω_i after observing its features as X .
 - Evidence $p(x)$: a scalar factor to guarantee posterior probabilities sum to one regarding ω_i .

$$p(\omega_i | X) = \frac{P(\omega_i) \cdot p(X | \omega_i)}{p(X)} = \frac{P(\omega_i) \cdot p(X | \omega_i)}{\sum_i P(\omega_i) \cdot p(X | \omega_i)}$$

Bayes decision rule: Maximum *a posterior* (MAP) (II)

- If we observe some features X of an unknown pattern, the observation can convert the prior into posterior. Intuitively, we can class the pattern based on the posterior probabilities, resulting in *the maximum a posterior (MAP) decision rule*, also called *Bayes decision rule*.
- For an unknown pattern P , after observing some features X , we classify it into the class with the largest posterior probability:

$$\begin{aligned}\omega_p &= \arg \max_{\omega_i} p(\omega_i | X) = \arg \max_{\omega_i} \frac{P(\omega_i) \cdot p(X | \omega_i)}{p(X)} \\ &= \arg \max_{\omega_i} P(\omega_i) \cdot p(X | \omega_i)\end{aligned}$$

The MAP decision rule is optimal (I)

- How well the MAP decision rule behaves??
- **Optimality:** assume we have complete knowledge, including $P(\omega_i)$ and $p(X | \omega_i)$ ($i=1,2,\dots,N$), the MAP decision rule is optimal to classify patterns, which means it will achieve the lowest average classification error rate.
- **Proof of optimality of the MAP rule:**
 - Given a pattern P , if its true class id is ω_i , but we classify it as ω_p , then the classification error is counted as $l(\omega_p | \omega_i)$:

$$l(\omega_p | \omega_i) = \begin{cases} 0 & (\omega_p = \omega_i) \\ 1 & (\omega_p \neq \omega_i) \end{cases}$$

which is also known as 0-1 loss function.

The MAP decision rule is optimal (II)

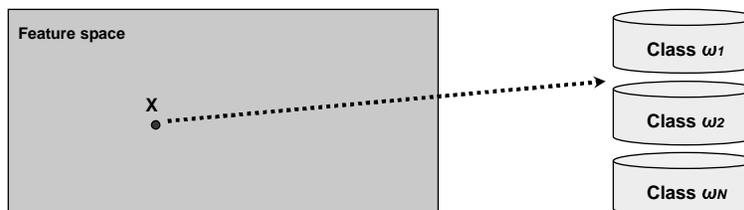
- Proof of optimality of the MAP rule (cont')
 - For a pattern P , after observing X , the posterior $p(\omega_i | X)$ is the probability that the true class id of P is ω_i . Thus the expected (average) classification error associated with classifying P as ω_P is calculated:

$$\begin{aligned} R(\omega_P | X) &= \sum_{i=1}^N l(\omega_P | \omega_i) \cdot p(\omega_i | X) \\ &= \sum_{\omega_i \neq \omega_P} p(\omega_i | X) \\ &= 1 - p(\omega_P | X) \end{aligned}$$

- The optimal classification is to minimize the above average classification error, i.e., if observing X , we classify P as ω_P to minimize $R(\omega_P | X) \rightarrow$ maximize $p(\omega_P | X)$
 - \rightarrow the MAP decision rule is optimal, which achieves the minimum average error rate. The minimum error is called Bayes error.

The MAP decision rule

- A general decision rule is a mapping function, given any an observation X , output a class id ω_P : $X \rightarrow \omega_P$
- If we totally have N classes, a decision rule will partition the entire feature space of X into N different regions, O_1, O_2, \dots, O_N . If X is located in the region O_i , we classify it as class ω_i .
- Each region O_i could consist of many contiguous areas.
- The MAP decision rule (the Bayes decision rule) is optimal among all possible decision rules in terms of minimizing average classification errors conditional on that we have complete and precise knowledge about the underlying problem.



The MAP decision rule example

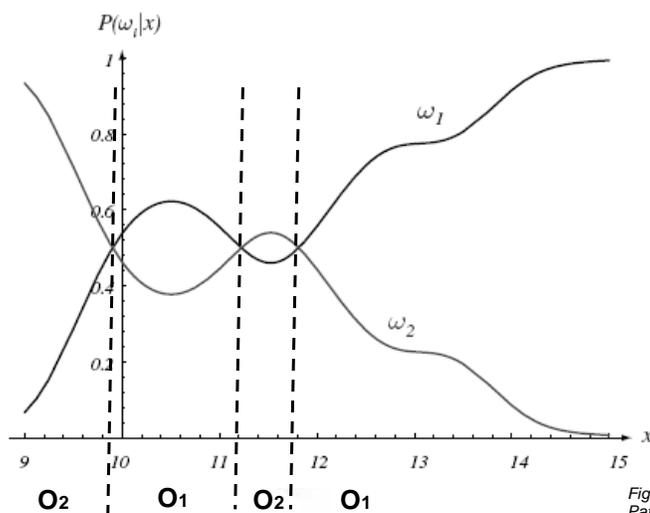


Figure from Duda et. al.,
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Classification Error Probability of a decision rule

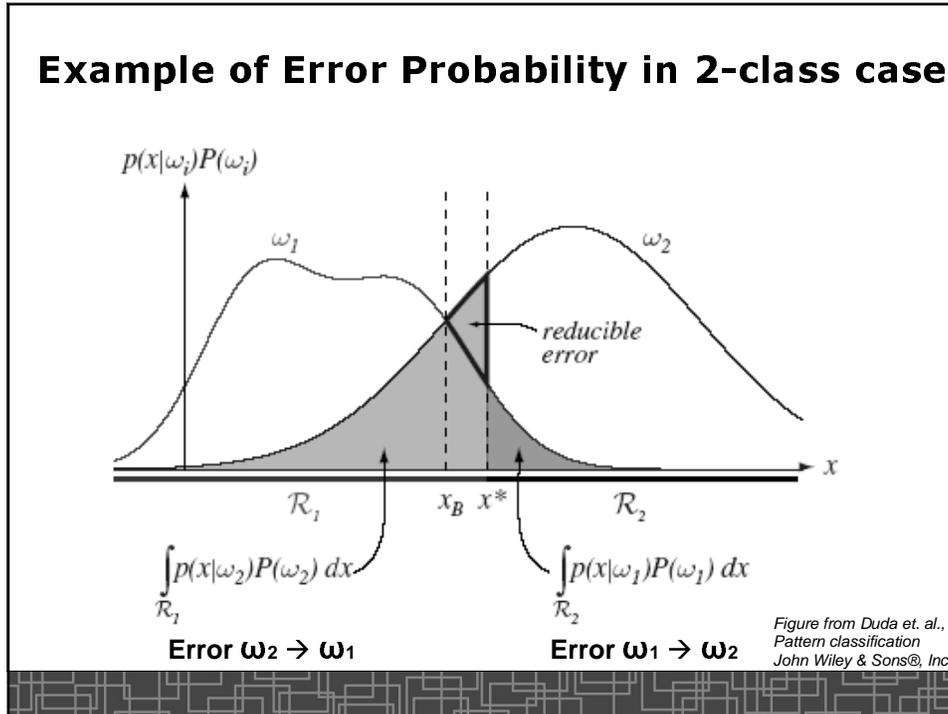
- Assume N -class problem, any a decision rule partition the feature space into N regions, O_1, O_2, \dots, O_N .
- $\Pr(X \in O_i, \omega_j)$ denotes the probability of the observation X of a pattern (its true class id is ω_j) falls in the region O_i .
- The overall classification error probability of the decision rule is:

$$\Pr(\text{error}) = 1 - \Pr(\text{correct})$$

$$= 1 - \sum_{i=1}^N \Pr(X \in O_i, \omega_i)$$

$$= 1 - \sum_{i=1}^N \Pr(X \in O_i | \omega_i) \cdot P(\omega_i)$$

$$= 1 - \sum_{i=1}^N \int_{O_i} p(X | \omega_i) \cdot P(\omega_i) dX$$



Bayes Error

- Bayes error: error probability of the Bayes (MAP) decision rule.
- Since Bayes decision rule guarantees the minimum error, the Bayes error is the lower bound of all possible error probabilities.
- It is difficult to calculate what the probability actually is, even for the very simple cases because of discontinuous nature of the decision regions in the integral, especially in high dimensions.
- Some approximation methods to estimate an upper bound.
 - Chernoff bound
 - Bhattacharyya bound
- Evaluate on an independent test set.

Example: Bayes decision for independent binary features

- Bayes decision rule (the MAP rule) is also applicable when feature X is discrete.
- A simple case: 2-class (ω_1, ω_2), feature vector is d -dimensional vector, whose components are binary-valued and conditionally independent.

$$X = (x_1, x_2, \dots, x_d)^t \quad x_i = 0, 1 \quad (1 \leq i \leq d)$$

$$p_i = \Pr(x_i = 1 | \omega_1) \quad \text{and} \quad q_i = \Pr(x_i = 1 | \omega_2)$$

$$p(X | \omega_1) = \prod_{i=1}^d p_i^{x_i} (1 - p_i)^{1-x_i}$$

$$p(X | \omega_2) = \prod_{i=1}^d q_i^{x_i} (1 - q_i)^{1-x_i}$$

Example: Bayes decision for independent binary features

- The MAP decision rule:

classify to ω_1 if $P(\omega_1) \cdot p(X | \omega_1) \geq P(\omega_2) \cdot p(X | \omega_2)$, otherwise ω_2

Equivalently, we have the decision function :

$$g(X) = \sum_{i=1}^d \left[x_i \ln \frac{p_i}{q_i} + (1 - x_i) \ln \frac{1 - p_i}{1 - q_i} \right] + \ln \frac{P(\omega_1)}{P(\omega_2)} = \sum_{i=1}^d \lambda_i x_i + \lambda_0$$

$$\lambda_i = \ln \frac{p_i(1 - q_i)}{q_i(1 - p_i)} \quad \lambda_0 = \sum_{i=1}^d \ln \frac{1 - p_i}{1 - q_i} + \ln \frac{P(\omega_1)}{P(\omega_2)}$$

If $g(X) \geq 0$, classify to ω_1 , otherwise ω_2 .

Missing features/data (I)

- If we know the full probability structure of a problem, we can construct the optimal Bayes decision rule.
- In some practical situations, for some patterns, we can't observe the full feature vector described in the probability structure. Only partial information of the feature vector is observed, but some components are missing.
- How to classify such corrupted inputs to obtain minimum average error?
- Let the full feature vector $X=[X_g, X_b]$, X_g represents the observed or good features, X_b represents the missing or bad ones.
- In this case, the optimal decision rule is constructed based on the posterior, $p(\omega_i | X_g)$, as follows:

$$\omega_p = \arg \max_{\omega_i} p(\omega_i | X_g)$$

Missing features/data(II)

• Where

$$p(\omega_i | X_g) = \frac{p(\omega_i, X_g)}{p(X_g)} = \frac{\int p(\omega_i, X_g, X_b) dX_b}{p(X_g)}$$

$$(1) = \frac{\int p(\omega_i | X_g, X_b) \cdot p(X_g, X_b) dX_b}{p(X_g)}$$

$$= \frac{\int p(\omega_i | X) \cdot p(X) dX_b}{\int p(X) dX_b}$$

$$(2) = \frac{\int P(\omega_i) \cdot p(X_g, X_b | \omega_i) dX_b}{\sum_{\omega_i} \int P(\omega_i) \cdot p(X_g, X_b | \omega_i) dX_b}$$

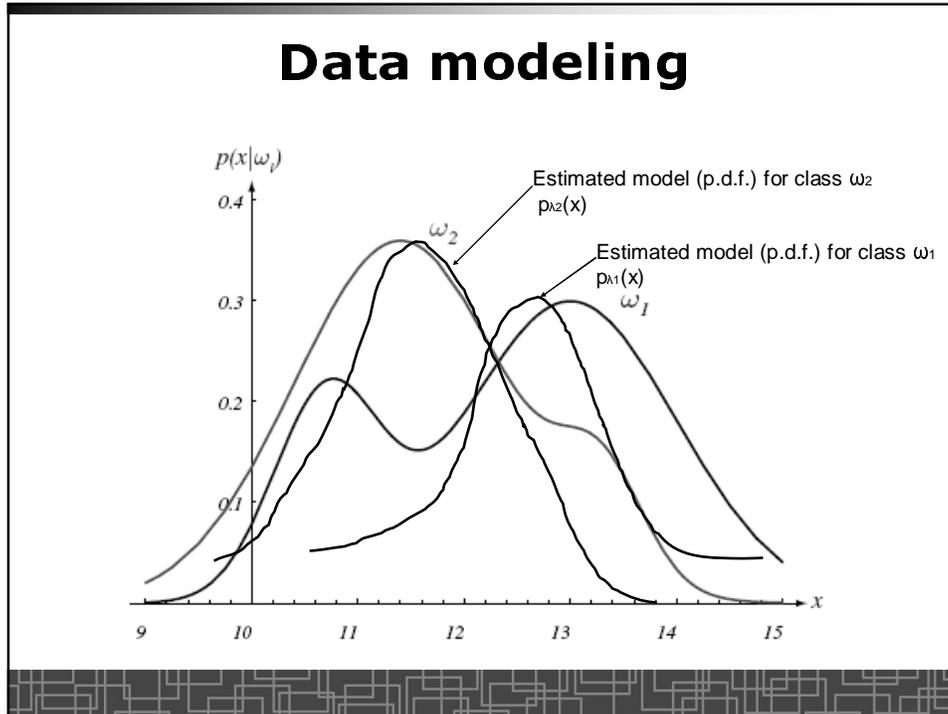
$$= \frac{\int P(\omega_i) \cdot p(X | \omega_i) dX_b}{\sum_{\omega_i} \int P(\omega_i) \cdot p(X | \omega_i) dX_b}$$

Optimal Bayes decision rule is not achievable in practice

- The optimal Bayes decision rule is not feasible in practice.
 - In any practical problem, we can not have a complete knowledge about the problem.
 - e.g., the class-conditional probability density functions (p.d.f.), $p(X | \omega_i)$, are always unavailable and extremely hard to estimate.
- However, possible to collect a set of sample data (a set of feature observations) for each class in question.
 - The sample data are always far from enough to estimate a reliable p.d.f. by using sample data themselves ONLY, e.g., some nonparametric methods → sampling density / histogram.
- Question: How to build a reasonable classifier based on a limited set of sample data, instead of the true p.d.f.'s?

Statistical Data Modeling

- For any real problem, the true p.d.f.'s are always unknown, neither the form of the functions nor the parameters.
- Our approach – **statistical data modeling** : based on the available sample data set, choose a proper statistical model to fit into the available data set.
 - **Data Modeling stage**: once the statistical model is selected, its function form becomes known except a set of model parameters associated with the model are unknown to us.
 - **Learning (training) stage**: the unknown parameters can be estimated by fitting the model into the data set based on certain estimation criterion.
 - the estimated statistical model (assumed model format + estimated parameters) will give a parametric p.d.f. to approximate the real but unknown p.d.f. of each class.
 - **Decision (test) stage**: the estimated p.d.f.'s are plugged into the optimal Bayes decision rule in place of the real p.d.f.'s
 - plug-in MAP decision rule
 - Not optimal any more but performs reasonably well in practice



Plug-in MAP decision rule

- Once the statistical models are estimated, they are treated as if they were true distributions of the data, and plug into the form of the optimal Bayes (MAP) decision rule in place of the unknown true p.d.f.'s.
- The plug-in MAP decision rule:

$$\begin{aligned}
 \omega_p &= \arg \max_{\omega_i} p(\omega_i | X) = \arg \max_{\omega_i} \frac{P(\omega_i) \cdot p(X | \omega_i)}{p(X)} \\
 &= \arg \max_{\omega_i} P(\omega_i) \cdot p(X | \omega_i) \\
 &\approx \arg \max_{\omega_i} \bar{P}_{\Gamma_i}(\omega_i) \cdot \bar{p}_{\Lambda_i}(X | \omega_i)
 \end{aligned}$$

Some useful models(I)

- A proper model must be chosen based on the nature of observation data.
- Some useful statistical models for a variety of data
 - Normal (Gaussian) distribution
 - uni-modal continuous feature scalars
 - Multivariate normal (Gaussian) distribution
 - uni-modal continuous feature vectors
 - Gaussian Mixture models (GMM)
 - continuous feature scalars/vectors with multi-modal distribution nature
 - For speaker recognition/verification distribution of speech features over a large population



Some useful models (II)

- Some useful models (cont'd)
 - Markov chain model: discrete sequential data
 - N-gram model in language modeling
 - Hidden Markov Models (HMM's): ideal for various kinds of sequential observation data; provides better modeling capability than simple Markov chain model.
 - Model speech signals for recognition (one of the most successful story of data modeling)
 - Model language/text data for part-of-speech tagging, shallow language understanding, etc.
 - Model biological data (DNA & protein sequence): profile HMM.
 - Lots of other application domains.

Some useful models (III)

- Some useful models (*not covered in this course*)
 - Random Markov Field: multi-dimensional spatial data
 - Model image data: e.g., used for OCR, etc.
 - HMM is a special case of random Markov field
 - Graphical models (a.k.a., Bayesian networks, Belief networks)
 - High-dimensional data (discrete or continuous)
 - To model a very complex stochastic process
 - Automatically learn dependency from data
 - Used widely in machine learning, data mining
 - HMM is also a special case of graphical model
- Neural networks, support vector machine (SVM) DON'T fit here.
 - Not to model the distribution (p.d.f.) of data directly.
 - Discriminative method: model the boundaries of data sets

Generative vs. discriminative models

- Posterior probability $p(\omega_i|X)$ plays the key role in pattern classification.
 - Generative Models: focus on probability distribution of data
$$p(\omega_i|X) \sim p(\omega_i) \cdot p(X|\omega_i)$$
$$\approx p'(\omega_i) \cdot p'(X|\omega_i) \quad (\text{the plug-in MAP rule})$$
 - Discriminative Models: directly model discriminant function:
$$p(\omega_i|X) \sim g_i(X)$$

Pattern classification based on Discriminant Functions (I)

- Instead of designing a classifier based on probability distribution of data, we can build an ad-hoc classifier based on some discriminant functions to model class boundary info directly.
- Classifier based on discriminant functions:
 - For N classes, we define a set of discriminant functions $g_i(X)$ ($i=1,2,\dots,N$), one for each class.
 - For an unknown pattern with feature vector Y , the classifier makes the decision as

$$\omega_Y = \arg \max_i g_i(Y)$$

- Each discriminant function $g_i(X)$ has a pre-defined function form and a set of unknown parameters θ_i , rewrite it as $g_i(X; \theta_i)$.
- Similarly θ_i ($i=1,2,\dots,N$) need to be estimated from some training data.

Pattern classification based on Discriminant Functions (II)

- Some common forms for discriminant functions:
 - Linear discriminant function:

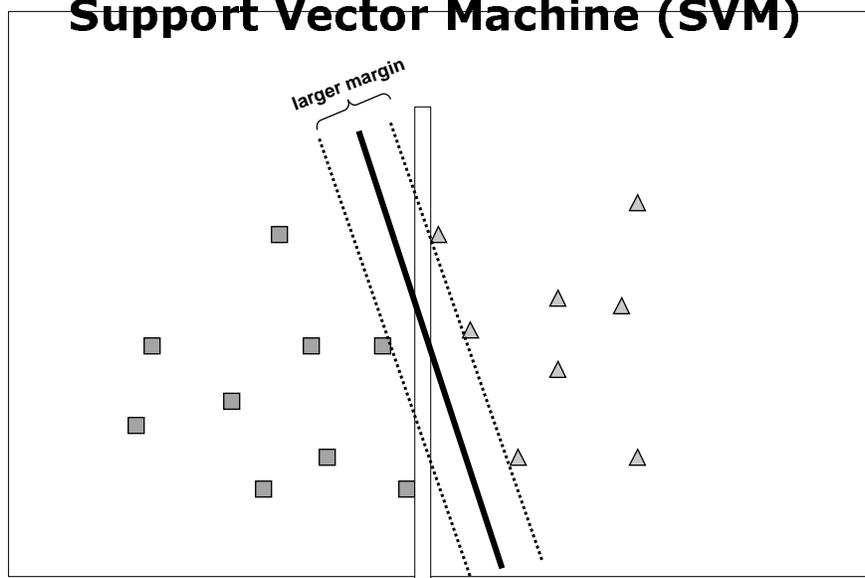
$$g(X) = w^t \cdot X + w_0$$

- Quadratic discriminant function: (2nd order)
- Polynomial discriminant function: (N-th order)
- *Neural network*: (arbitrary nonlinear functions)
- Optimal discriminant functions: optimal MAP classifier is a special case when choosing discriminant functions as class posterior probabilities.

Pattern classification based on Discriminant Functions (III)

- Unknown parameters of discriminant functions are estimated by some gradient descent method to minimize an objective function, such as empirical classification errors in training set, etc.
 - minimizing empirical classification errors
 - maximize separation margin: support vector machine

Large-Margin Classifier: Support Vector Machine (SVM)

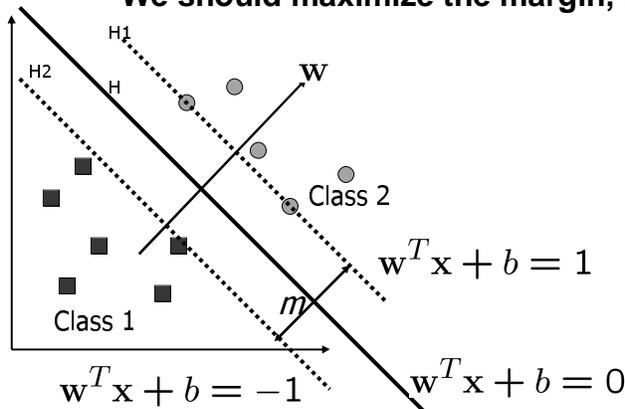


Support Vector Machine (I)

- The decision boundary H should be as far away from the data of both classes as possible

– We should maximize the margin, m

$$m = \frac{2}{\|w\|}$$



Because the perpendicular distance from H_1, H_2 to the origin is $|1-b|/\|w\|$ and $|-1-b|/\|w\|$

Support Vector Machine (II)

- The decision boundary can be found by solving the following constrained optimization problem:

$$\text{Minimize } \frac{1}{2} \|w\|^2 \quad \|w\|^2 = w^T w$$

$$\text{subject to } y_i(w^T x_i + b) \geq 1 \quad \forall i$$

- Convert to its dual problem:

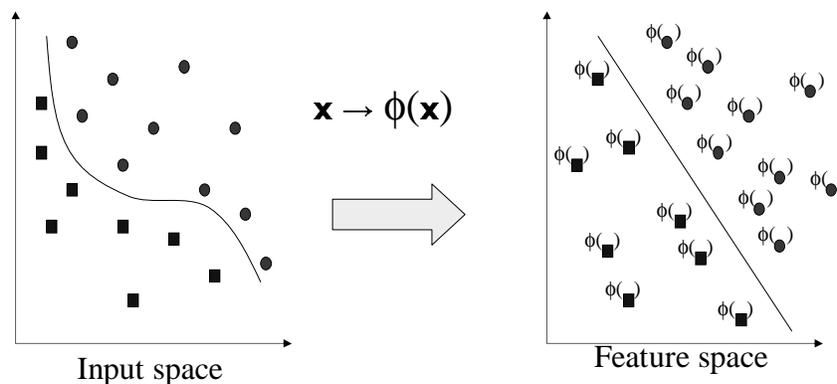
$$\text{max. } W(\alpha) = \sum_{i=1}^n \alpha_i - \frac{1}{2} \sum_{i=1, j=1}^n \alpha_i \alpha_j y_i y_j x_i^T x_j$$

$$\text{subject to } \alpha_i \geq 0, \sum_{i=1}^n \alpha_i y_i = 0$$

– This is a standard quadratic programming (QP) problem.

Support Vector Machine (III)

- For nonlinear separation boundary:
 - use a Kernel function



Linearly Non-Separable cases

- We allow “error” ξ_i in classification

