CS840a Fall 2006 Learning and Computer Vision Prof. Olga Veksler

Lecture 3

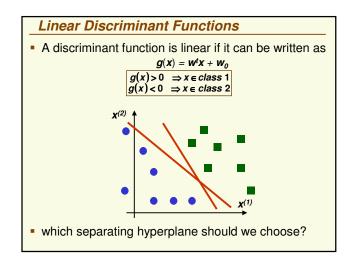
SVM Information Theory (a little BIT) Some pictures from C. Burges

SVM

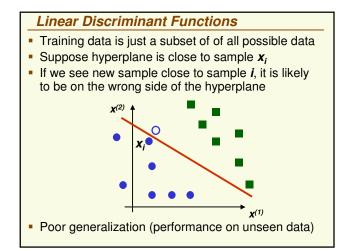
- Said to start in 1979 with Vladimir Vapnik's paper
- Major developments throughout 1990's
- Elegant theory
 - Has good generalization properties
- Have been applied to diverse problems very successfully in the last 10-15 years
- One of the most important developments in pattern recognition in the last 10 years

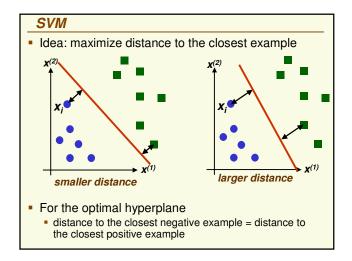
Today

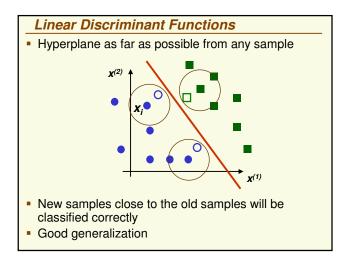
- Support Vector Machines
- Mutual Information
- Preparation for the next time:
 - papers: "Object Recognition with Informative Features and Linear Classification" by M. Naquet and S. Ullman
 - Ignore section of tree-augmented network
 - "Face Recognition with Support Vector Machines: Global vs. Component-based Approach"

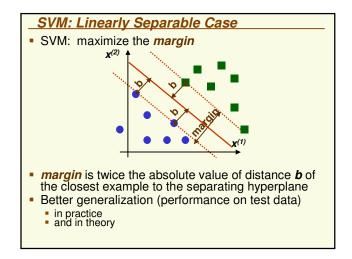


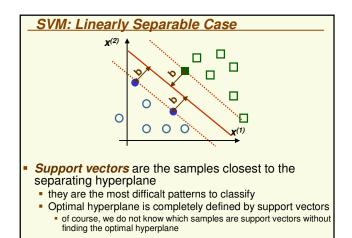


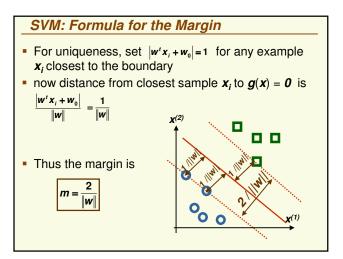


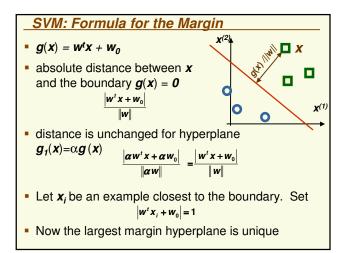


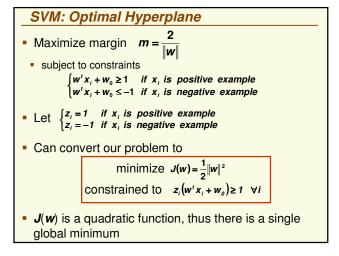


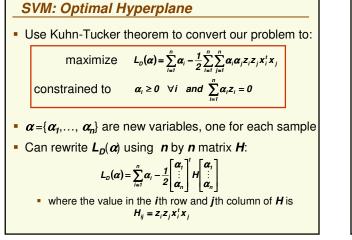


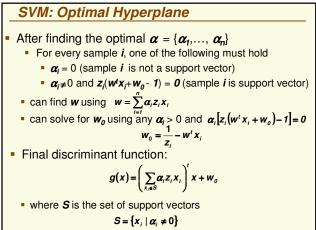












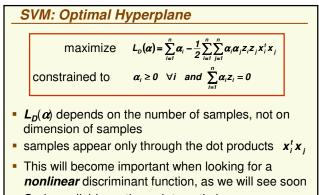
SVM: Optimal Hyperplane

Use Kuhn-Tucker theorem to convert our problem to:

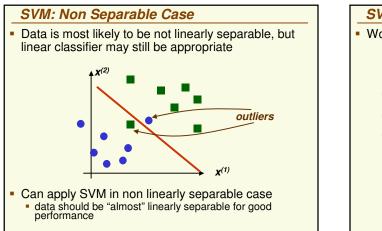
maximize
$$L_{D}(\alpha) = \sum_{i=1}^{m} \alpha_{i} - \frac{1}{2} \sum_{i=1}^{m} \sum_{j=1}^{n} \alpha_{i} \alpha_{j} z_{i} z_{j} x_{i}^{t} x_{j}$$

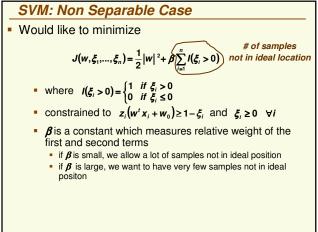
constrained to $\alpha_{i} \ge 0 \quad \forall i \text{ and } \sum_{i=1}^{n} \alpha_{i} z_{i} = 0$

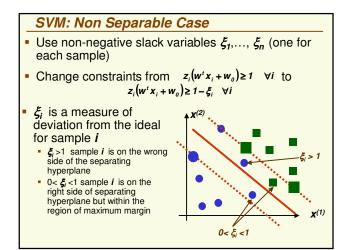
- $\alpha = \{\alpha_1, ..., \alpha_n\}$ are new variables, one for each sample
- $L_D(a)$ can be optimized by quadratic programming
- $L_{D}(\alpha)$ formulated in terms of α
 - it depends on w and wo indirectly

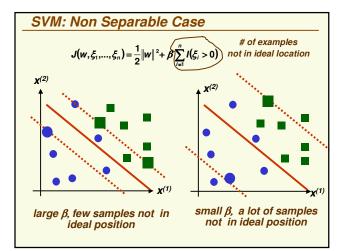


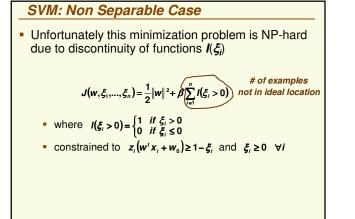
Code available on the web to optimize

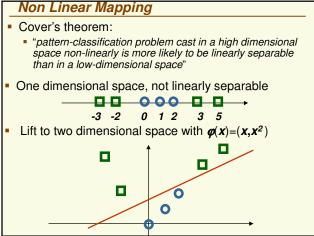


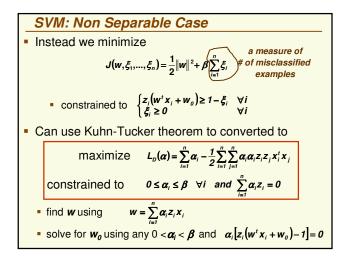


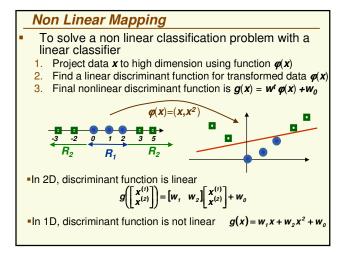


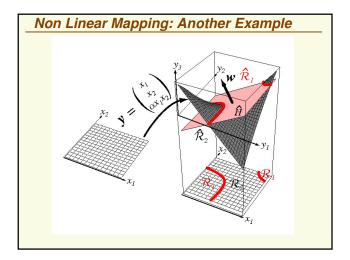












	Non Linear SVM: Kernels
	Recall SVM optimization
	maximize $L_D(\alpha) = \sum_{i=1}^n \alpha_i - \frac{1}{2} \sum_{i=1}^n \sum_{j=1}^n \alpha_i \alpha_i z_j x_j^* x_j$
•	Note this optimization depends on samples x_i only through the dot product $x_i^t x_j$

If we lift x_i to high dimension using $\varphi(x)$, need to compute high dimensional product $\varphi(\mathbf{x}_i)^t \varphi(\mathbf{x}_i)$

maximize
$$L_D(\alpha) = \sum_{i=1}^n \alpha_i - \frac{1}{2} \sum_{i=1}^n \sum_{j=1}^n \alpha_j \alpha_i z_i z_j \varphi(x_i)^{*} \varphi(x_j)$$

 $K(x_n, x_j)$

Idea: find *kernel* function $K(x_i, x_j)$ s.t. $K(\mathbf{x}_i, \mathbf{x}_i) = \boldsymbol{\varphi}(\mathbf{x}_i)^{t} \boldsymbol{\varphi}(\mathbf{x}_i)$

New Lineau OV/M. Kernels

Non Linear SVM

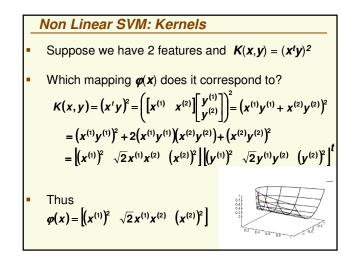
- Can use any linear classifier after lifting data into a higher dimensional space. However we will have to deal with the "curse of dimensionality"
 - 1. poor generalization to test data
 - 2. computationally expensive
- SVM avoids the "curse of dimensionality" problems by
 - enforcing largest margin permits good generalization

 It can be shown that generalization in SVM is a function of the margin, independent of the dimensionality
 - computation in the higher dimensional case is performed only implicitly through the use of *kernel* functions

Non Linear SVM: Kernels

maximize
$$L_{D}(\alpha) = \sum_{i=1}^{n} \alpha_{i} - \frac{1}{2} \sum_{i=1}^{n} \sum_{j=1}^{n} \alpha_{i} \alpha_{i} z_{i} z_{j} \varphi(x_{i})^{*} \varphi(x_{j}) \frac{1}{K(x_{i}, x_{i})}$$

- Then we only need to compute $K(x_i, x_i)$ instead of
 - $\varphi(x_i)^t \varphi(x_j)$ "kernel trick": do not need to perform operations in high dimensional space explicitly



Non Linear SVM

- search for separating hyperplane in high dimension $w\varphi(x) + w_o = 0$
- Choose φ(x) so that the first ("0"th) dimension is the augmented dimension with feature value fixed to 1

$$\varphi(x) = \begin{bmatrix} 1 & x^{(1)} & x^{(2)} & x^{(1)}x^{(2)} \end{bmatrix}^{1}$$

• Threshold parameter w_0 gets folded into the weight vector w

$$\begin{bmatrix} \mathbf{w}_0 & \mathbf{w} \end{bmatrix} \begin{bmatrix} \mathbf{x} \\ \mathbf{\varphi}(\mathbf{x}) \end{bmatrix} = \mathbf{0}$$

Non Linear SVM: Kernels

- How to choose kernel function $K(x_i, x_i)$?
- *K*(*x_i*,*x_j*) should correspond to product φ(*x_i*)^tφ(*x_j*) in a higher dimensional space
- Mercer's condition tells us which kernel function can be expressed as dot product of two vectors
- Kernel's not satisfying Mercer's condition can be sometimes used, but no geometrical interpretation
- Some common choices (satisfying Mercer's condition):
- Polynomial kernel $K(x_i, x_j) = (x_i^t x_j + 1)^p$
- Gaussian radial Basis kernel (data is lifted in infinite dimension)

$$\mathcal{K}(\boldsymbol{x}_i, \boldsymbol{x}_j) = \exp\left(-\frac{1}{2\sigma^2} \|\boldsymbol{x}_i - \boldsymbol{x}_j\|^2\right)$$

Non Linear SVM

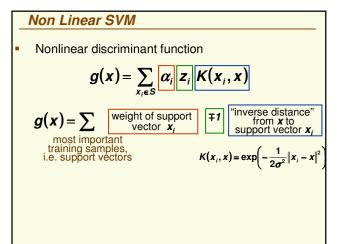
- Will not use notation a = [w₀ w], we'll use old notation w and seek hyperplane through the origin wp(x) = 0
- If the first component of $\varphi(x)$ is not **1**, the above is equivalent to saying that the hyperplane has to go through the origin in high dimension
 - removes only one degree of freedom
 - But we have introduced many new degrees when we lifted the data in high dimension



- Start with data x₁,...,x_n which lives in feature space of dimension d
- Choose kernel K(x_i,x_j) or function φ(x_i) which takes sample x_i to a higher dimensional space
- Find the largest margin linear discriminant function in the higher dimensional space by using quadratic programming package to solve:

maximize $L_D(\alpha) = \sum_{i=1}^n \alpha_i - \frac{1}{2} \sum_{i=1}^n \sum_{j=1}^n \alpha_j \alpha_j z_j z_j K(x_i, x_j)$

constrained to
$$0 \le \alpha_i \le \beta \quad \forall i \text{ and } \sum_{i=1}^{n} \alpha_i z_i = 0$$



Non Linear SVM Recipe

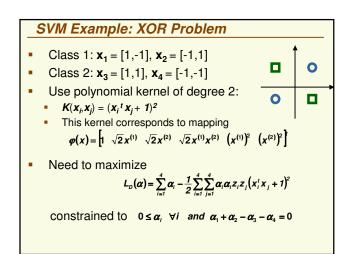
- Weight vector **w** in the high dimensional space: $w = \sum_{x_i \in S} \alpha_i z_i \varphi(x_i)$
 - where **S** is the set of support vectors $S = \{x_i \mid \alpha_i \neq 0\}$
- Linear discriminant function of largest margin in the high dimensional space:

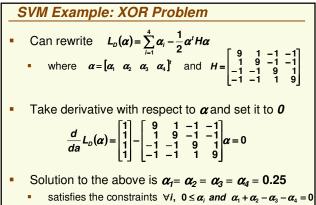
$$g(\varphi(x)) = w^t \varphi(x) = \left(\sum_{x_i \in S} \alpha_i z_i \varphi(x_i)\right)^t \varphi(x)$$

Non linear discriminant function in the original space $a(x) = \left(\sum \alpha_{i,z,m}(x_{i})\right)^{t} a(x) = \sum \alpha_{i,z,m}(x_{i}) a(x) = \sum \alpha_{i,z,k}(x_{i}, x)$

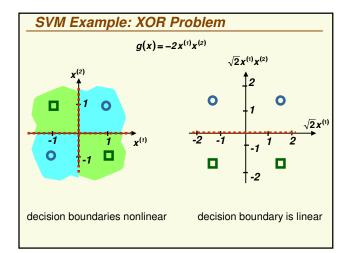
$$g(x) = \left(\sum_{x_i \in S} \alpha_i z_i \varphi(x_i)\right) \varphi(x) = \sum_{x_i \in S} \alpha_i z_i \varphi^i(x_i) \varphi(x) = \sum_{x_i \in S} \alpha_i z_i \kappa(x_i, x)$$

• decide class 1 if g(x) > 0, otherwise decide class 2





all samples are support vectors



SVM Example: XOR Problem

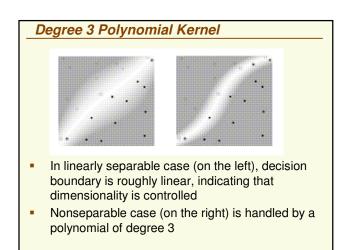
 $\varphi(x) = \begin{bmatrix} 1 & \sqrt{2} x^{(1)} & \sqrt{2} x^{(2)} & \sqrt{2} x^{(1)} x^{(2)} & (x^{(1)})^2 & (x^{(2)})^2 \end{bmatrix}$

Weight vector w is:

$$w = \sum_{i=1}^{4} \alpha_i z_i \varphi(x_i) = 0.25(\varphi(x_1) + \varphi(x_2) - \varphi(x_3) - \varphi(x_4))$$
$$= \begin{bmatrix} 0 & 0 & 0 & -\sqrt{2} & 0 & 0 \end{bmatrix}$$

• Thus the nonlinear discriminant function is:

$$g(x) = w\varphi(x) = \sum_{i=1}^{\circ} W_i \varphi_i(x) = -\sqrt{2} \left(\sqrt{2} x^{(1)} x^{(2)} \right) = -2 x^{(1)} x^{(2)}$$



SVM Summary

Advantages:

- Based on nice theory
- excellent generalization properties
- objective function has no local minima
- can be used to find non linear discriminant functions
- Complexity of the classifier is characterized by the number of support vectors rather than the dimensionality of the transformed space
- Disadvantages:
- tends to be slower than other methods
- quadratic programming is computationally expensive
- Not clear how to choose the Kernel

Information theory

• Suppose we toss a **fair** die with 8 sides

- need 3 bits to transmit the results of each toss
- 1000 throws will need 3000 bits to transmit
- Suppose the die is biased
 - side A occurs with probability 1/2, chances of throwing B are 1/4, C are 1/8, D are 1/16, E are 1/32, F 1/64, G and H are 1/128
 - Encode A= 0, B = 10, C = 110, D = 1110,..., so on until G =
 - 1111110, H = 1111111
 - We need, on average, 1/2+2/4+3/8+4/16+5/32+6/64+7/128+7/128
 - = 1.984 bits to encode results of a toss1000 throws require 1984 bits to transmit
 - Less bits to send = less "information"
 - Biased die tosses contain less "information" than unbiased die tosses (know in advance biased sequence will have a lot of A's)
- What's the number of bits in the best encoding?
- Extreme case: if a die always shows side A, a sequence of 1,000 tosses has no information, 0 bits to encode

Information theory

- Information Theory regards information as only those symbols that are uncertain to the receiver
- only infrmatn esentil to understnd mst b tranmitd
 Shannon made clear that uncertainty is the very commodity of communication
- The amount of information, or uncertainty, output by an information source is a measure of its entropy
- In turn, a source's entropy determines the amount of bits per symbol required to encode the source's information
- Messages are encoded with strings of 0 and 1 (bits)

Information theory

- if a die is fair (any side is equally likely, or uniform distribution), for any toss we need log(8) = 3 bits
- Suppose any of n events is equally likely (uniform distribution)
 P(x) = 1/n, therefore -log P = -log(1/n) = log n
- In the "good" encoding strategy for our biased die example, every side x has -log p(x) bits in its code
- Expected number of bits is



Shannon's Entropy

$H[p(x)] = -\sum_{x} p(x) \log p(x) = \sum_{x} p(x) \log \frac{1}{p(x)}$

- How much randomness (or uncertainty) is there in the value of signal x if it has distribution p(x)
 - For uniform distribution (every event is equally likely), H[x] is maximum
 - If p(x) = 1 for some event x, then H[x] = 0
 - Systems with one very common event have less entropy than systems with many equally probable events
- Gives the expected length of optimal encoding (in binary bits) of a message following distribution p(x)
 - doesn't actually give this optimal encoding

Mutual Information of X and Y

$I[x, y] = H(x) - H(x \mid y)$

- Measures the average reduction in uncertainty about x after y is known
- or, equivalently, it measures the amount of information that y conveys about x
- Properties
 - I(x,y) = I(y,x)
 - I(x,y) ≥ 0
 - If x and y are independent, then I(x,y) = 0
 - I(x,x) = H(x)

Conditional Entropy of X given Y

$$H[x | y] = \sum_{x,y} p(x,y) \log \frac{1}{p(x | y)} = -\sum_{x,y} p(x,y) \log p(x | y)$$

- Measures average uncertainty about x when y is known
- Property:
 - H[x] ≥ H[x|y], which means after seeing new data (y), the uncertainty about x is not increased, on average

MI for Feature Selection

$$I[x,c] = H(c) - H(c | x)$$

- Let x be a proposed feature and c be the class
- If I[x,c] is high, we can expect feature x be good at predicting class c