

CSC 411: Lecture 16: Kernels

Class based on Raquel Urtasun & Rich Zemel's lectures

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- Kernel trick

Summary of Linear SVM

- Binary and linear separable classification
- Linear classifier with maximal margin
- Training SVM by maximizing

$$\max_{\alpha_i \geq 0} \left\{ \sum_{i=1}^N \alpha_i - \frac{1}{2} \sum_{i,j=1}^N t^{(i)} t^{(j)} \alpha_i \alpha_j (\mathbf{x}^{(i)T} \cdot \mathbf{x}^{(j)}) \right\}$$

$$\text{subject to } \alpha_i \geq 0; \quad \sum_{i=1}^N \alpha_i t^{(i)} = 0$$

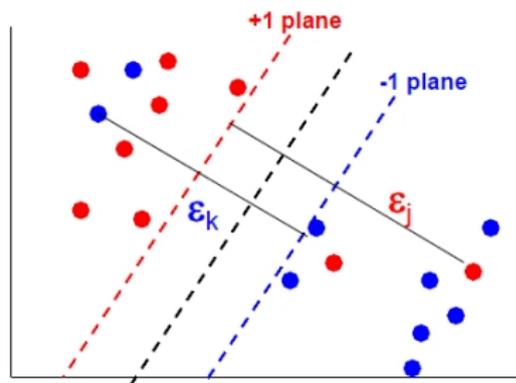
- The weights are

$$\mathbf{w} = \sum_{i=1}^N \alpha_i t^{(i)} \mathbf{x}^{(i)}$$

- Only a small subset of α_i 's will be nonzero, and the corresponding $\mathbf{x}^{(i)}$'s are the **support vectors** \mathbf{S}
- Prediction on a new example:

$$y = \text{sign} \left[b + \mathbf{x} \cdot \left(\sum_{i=1}^N \alpha_i t^{(i)} \mathbf{x}^{(i)} \right) \right] = \text{sign} \left[b + \mathbf{x} \cdot \left(\sum_{i \in \mathbf{S}} \alpha_i t^{(i)} \mathbf{x}^{(i)} \right) \right]$$

What if data is not linearly separable?



- Introduce **slack variables** ξ_i

$$\min \frac{1}{2} \|\mathbf{w}\|^2 + \lambda \sum_{i=1}^N \xi_i$$

$$\text{s.t. } \xi_i \geq 0; \quad \forall i \quad t^{(i)}(\mathbf{w}^T \mathbf{x}^{(i)}) \geq 1 - \xi_i = 0$$

- Example lies on wrong side of hyperplane $\xi_i > 1$
- Therefore $\sum_i \xi_i$ upper bounds the number of training errors
- λ trades off training error vs model complexity
- This is known as the **soft-margin** extension

Non-linear Decision Boundaries

- Note that both the learning objective and the decision function depend only on dot products between patterns

$$\ell = \sum_{i=1}^N \alpha_i - \frac{1}{2} \sum_{i,j=1}^N t^{(i)} t^{(j)} \alpha_i \alpha_j (\mathbf{x}^{(i)T} \cdot \mathbf{x}^{(j)})$$

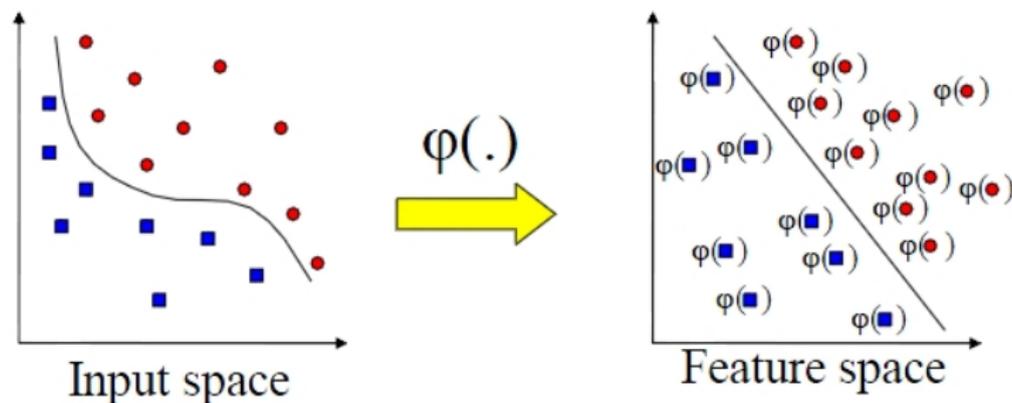
$$y = \text{sign}[b + \mathbf{x} \cdot (\sum_{i=1}^N \alpha_i t^{(i)} \mathbf{x}^{(i)})]$$

- How to form non-linear decision boundaries in input space?
 - Map data into feature space $\mathbf{x} \rightarrow \phi(\mathbf{x})$
 - Replace dot products between inputs with feature points

$$\mathbf{x}^{(i)T} \mathbf{x}^{(j)} \rightarrow \phi(\mathbf{x}^{(i)})^T \phi(\mathbf{x}^{(j)})$$

- Find linear decision boundary in feature space
- Problem: what is a good feature function $\phi(\mathbf{x})$?

Input Transformation



- Mapping to a feature space can produce problems:
 - ▶ High computational burden due to high dimensionality
 - ▶ Many more parameters
- SVM solves these two issues simultaneously
 - ▶ “Kernel trick” produces efficient classification
 - ▶ Dual formulation only assigns parameters to samples, not features

- **Kernel trick:** dot-products in feature space can be computed as a kernel function

$$K(\mathbf{x}^{(i)}, \mathbf{x}^{(j)}) = \phi(\mathbf{x}^{(i)})^T \phi(\mathbf{x}^{(j)})$$

- Idea: work directly on \mathbf{x} , avoid having to compute $\phi(\mathbf{x})$
- Example:

$$\begin{aligned} K(\mathbf{a}, \mathbf{b}) &= (\mathbf{a}^T \mathbf{b})^3 = ((a_1, a_2)^T (b_1, b_2))^3 \\ &= (a_1 b_1 + a_2 b_2)^3 \\ &= a_1^3 b_1^3 + 3a_1^2 b_1^2 a_2 b_2 + 3a_1 b_1 a_2^2 b_2^2 + a_2^3 b_2^3 \\ &= (a_1^3, \sqrt{3}a_1^2 a_2, \sqrt{3}a_1 a_2^2, a_2^3)^T (b_1^3, \sqrt{3}b_1^2 b_2, \sqrt{3}b_1 b_2^2, b_2^3) \\ &= \phi(\mathbf{a}) \cdot \phi(\mathbf{b}) \end{aligned}$$

- Examples of kernels: **kernels measure similarity**

1. Polynomial

$$K(\mathbf{x}^{(i)}, \mathbf{x}^{(j)}) = (\mathbf{x}^{(i)T} \mathbf{x}^{(j)} + 1)^d$$

where d is the degree of the polynomial, eg $d = 2$ for quadratic

2. Gaussian

$$K(\mathbf{x}^{(i)}, \mathbf{x}^{(j)}) = \exp\left(-\frac{\|\mathbf{x}^{(i)} - \mathbf{x}^{(j)}\|^2}{2\sigma^2}\right)$$

3. Sigmoid

$$K(\mathbf{x}^{(i)}, \mathbf{x}^{(j)}) = \tanh(\beta(\mathbf{x}^{(i)T} \mathbf{x}^{(j)} + a))$$

- Each kernel computation corresponds to dot product
 - ▶ calculation for particular mapping $\phi(\mathbf{x})$ implicitly maps to high-dimensional space
- Why is this useful?
 1. Rewrite training examples using more complex features
 2. Dataset not linearly separable in original space may be linearly separable in higher dimensional space

Kernel Functions

- Mercer's Theorem (1909): any reasonable kernel corresponds to some feature space
- Reasonable means that the Gram matrix is positive definite

$$K_{ij} = K(\mathbf{x}^{(i)}, \mathbf{x}^{(j)})$$

- Feature space can be very large
 - ▶ polynomial kernel $(1 + (\mathbf{x}^{(i)})^T \mathbf{x}^{(j)})^d$ corresponds to feature space exponential in d
 - ▶ Gaussian kernel has infinitely dimensional features
- Linear separators in these super high-dim spaces correspond to highly nonlinear decision boundaries in input space

Classification with Non-linear SVMs

- Non-linear SVM using kernel function $K()$:

$$\ell = \sum_{i=1}^N \alpha_i - \frac{1}{2} \sum_{i,j=1}^N t^{(i)} t^{(j)} \alpha_i \alpha_j K(\mathbf{x}^{(i)}, \mathbf{x}^{(j)})$$

- Maximize ℓ w.r.t. $\{\alpha\}$ under constraints $\forall i, \alpha_i \geq 0$
- Unlike linear SVM, cannot express \mathbf{w} as linear combination of support vectors
 - ▶ now must retain the support vectors to classify new examples
- Final decision function:

$$y = \text{sign}\left[b + \sum_{i=1}^N t^{(i)} \alpha_i K(\mathbf{x}, \mathbf{x}^{(i)})\right]$$

- Advantages:
 - ▶ Kernels allow very flexible hypotheses
 - ▶ Poly-time exact optimization methods rather than approximate methods
 - ▶ Soft-margin extension permits mis-classified examples
 - ▶ Variable-sized hypothesis space
 - ▶ Excellent results (1.1% error rate on handwritten digits vs. LeNet's 0.9%)
- Disadvantages:
 - ▶ Must choose kernel parameters
 - ▶ Very large problems computationally intractable
 - ▶ Batch algorithm

- Software:
 - ▶ A list of SVM implementations can be found at <http://www.kernel-machines.org/software.html>
 - ▶ Some implementations (such as LIBSVM) can handle multi-class classification
 - ▶ SVMLight is among the earliest implementations
 - ▶ Several Matlab toolboxes for SVM are also available
- Key points:
 - ▶ Difference between logistic regression and SVMs
 - ▶ Maximum margin principle
 - ▶ Target function for SVMs
 - ▶ Slack variables for mis-classified points
 - ▶ Kernel trick allows non-linear generalizations