

CS 188: Artificial Intelligence

Spring 2010

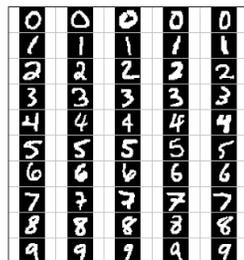
Lecture 24: Perceptrons and More!

4/22/2010

Pieter Abbeel – UC Berkeley
Slides adapted from Dan Klein

Announcements

- W7 due tonight [this is your last written for the semester!]
- Project 5 out tonight --- Classification!



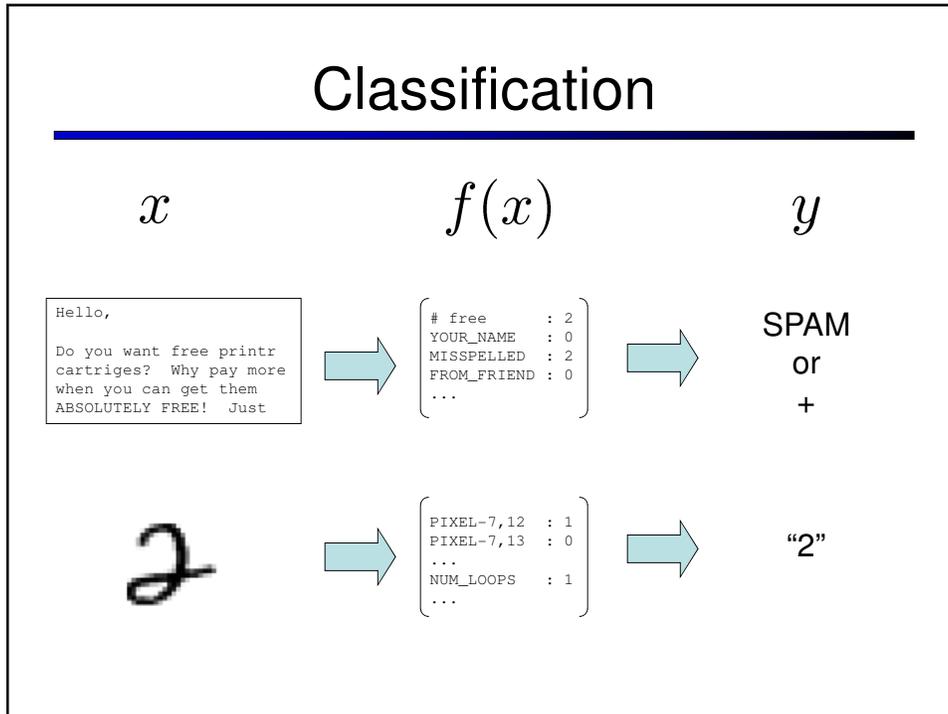
Announcements (2)

- Contest logistics
 - Up and running!
 - Tournaments every night
 - Final tournament: We will use submissions received by Thursday May 6, 11pm.
- Contest extra credit through bonus points on final exam [all based on final ranking]
 - 0.5pt for beating Staff
 - 0.5pt for beating Fa09-TeamA (top 5), Fa09-TeamB (top 10), and Fa09-TeamC (top 20) from last semester [total of 1.5pts to be earned]
 - 1pt for being 3rd
 - 2pts for being 2nd
 - 3pts for being 1st

Where are we and what's left?

- So far:
 - Search
 - CSPs
 - Adversarial search
 - MDPs and RL
 - Bayes nets, probabilistic inference
 - Machine learning
- Today: Machine Learning part III:
 - kNN and kernels
- Tuesday: Applications in Robotics
- Thursday: Applications in Vision and Language
+ Conclusion + Where to learn more

Classification



Classification overview

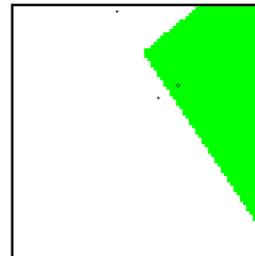
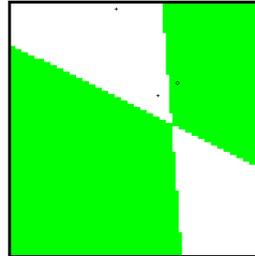
- **Naïve Bayes:**
 - Builds a model training data
 - Gives prediction probabilities
 - Strong assumptions about feature independence
 - One pass through data (counting)
- **Perceptron:**
 - Makes less assumptions about data
 - Mistake-driven learning
 - Multiple passes through data (prediction)
 - Often more accurate
- **SVM:**
 - Properties similar to perceptron
 - Convex optimization formulation
- **Nearest-Neighbor:**
 - Non-parametric: more expressive with more training data
- **Kernels**
 - Efficient way to make linear learning architectures into nonlinear ones

Case-Based Reasoning

- **Similarity for classification**
 - Case-based reasoning
 - Predict an instance's label using similar instances

- **Nearest-neighbor classification**
 - 1-NN: copy the label of the most similar data point
 - K-NN: let the k nearest neighbors vote (have to devise a weighting scheme)
 - Key issue: how to define similarity
 - Trade-off:
 - Small k gives relevant neighbors
 - Large k gives smoother functions
 - Sound familiar?

- [Demo]

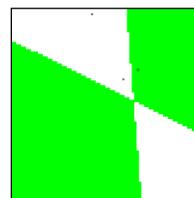


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<http://www.cs.cmu.edu/~zhuxj/courseproject/knndemo/KNN.html>

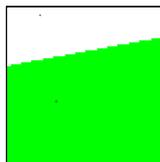
Parametric / Non-parametric

- **Parametric models:**
 - Fixed set of parameters
 - More data means better settings
- **Non-parametric models:**
 - Complexity of the classifier increases with data
 - Better in the limit, often worse in the non-limit
- (K)NN is **non-parametric**

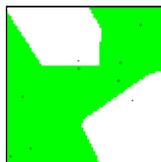


Truth

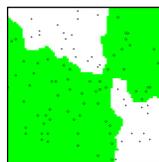
2 Examples



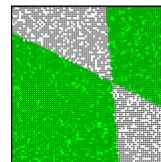
10 Examples



100 Examples



10000 Examples



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Nearest-Neighbor Classification

- Nearest neighbor for digits:
 - Take new image
 - Compare to all training images
 - Assign based on closest example

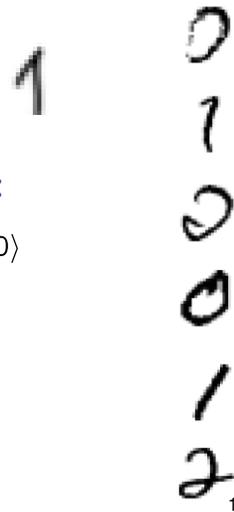
- Encoding: image is vector of intensities:

$$\mathbf{1} = \langle 0.0 \ 0.0 \ 0.3 \ 0.8 \ 0.7 \ 0.1 \dots 0.0 \rangle$$

- What's the similarity function?
 - Dot product of two images vectors?

$$\text{sim}(x, x') = x \cdot x' = \sum_i x_i x'_i$$

- Usually normalize vectors so $\|x\| = 1$
- min = 0 (when?), max = 1 (when?)



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Basic Similarity

- Many similarities based on **feature dot products**:

$$\text{sim}(x, x') = f(x) \cdot f(x') = \sum_i f_i(x) f_i(x')$$

- If features are just the pixels:

$$\text{sim}(x, x') = x \cdot x' = \sum_i x_i x'_i$$

- Note: not all similarities are of this form

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Invariant Metrics

- Better distances use knowledge about vision
- Invariant metrics:
 - Similarities are invariant under certain transformations
 - Rotation, scaling, translation, stroke-thickness...
 - E.g: 
 - $16 \times 16 = 256$ pixels; a point in 256-dim space
 - Small similarity in \mathbb{R}^{256} (why?)
 - Variety of invariant metrics in literature
- Viable alternative: transform training examples such that training set includes all variations

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Classification overview

- Naïve Bayes
- Perceptron
- SVM
- Nearest-Neighbor
- **Kernels**

A Tale of Two Approaches ...

- Nearest neighbor-like approaches
 - Can use fancy similarity functions
 - Don't actually get to do explicit learning
- Perceptron-like approaches
 - Explicit training to reduce empirical error
 - Can't use fancy similarity, only linear
 - Or can they? Let's find out!

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Perceptron Weights

- What is the final value of a weight w_y of a perceptron?
 - Can it be any real vector?
 - No! It's built by adding up inputs.

$$w_y = 0 + f(x_1) - f(x_5) + \dots$$

$$w_y = \sum_i \alpha_{i,y} f(x_i)$$

- Can reconstruct weight vectors (the **primal representation**) from update counts (the **dual representation**)

$$\alpha_y = \langle \alpha_{1,y} \ \alpha_{2,y} \ \dots \ \alpha_{n,y} \rangle$$

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Dual Perceptron

- How to classify a new example x ?

$$\begin{aligned}\text{score}(y, x) &= w_y \cdot f(x) \\ &= \left(\sum_i \alpha_{i,y} f(x_i) \right) \cdot f(x) \\ &= \sum_i \alpha_{i,y} (f(x_i) \cdot f(x)) \\ &= \sum_i \alpha_{i,y} K(x_i, x)\end{aligned}$$

- If someone tells us the value of K for each pair of examples, never need to build the weight vectors!

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Dual Perceptron

- Start with zero counts (alpha)
- Pick up training instances one by one
- Try to classify x_n ,

$$y = \arg \max_y \sum_i \alpha_{i,y} K(x_i, x_n)$$

- If correct, no change!
- If wrong: lower count of wrong class (for this instance), raise score of right class (for this instance)

$$\alpha_{y,n} = \alpha_{y,n} - 1$$

$$w_y = w_y - f(x_n)$$

$$\alpha_{y^*,n} = \alpha_{y^*,n} + 1$$

$$w_{y^*} = w_{y^*} + f(x_n)$$

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Kernelized Perceptron

- If we had a black box (**kernel**) which told us the dot product of two examples x and y :
 - Could work entirely with the dual representation
 - No need to ever take dot products (“kernel trick”)

$$\begin{aligned}\text{score}(y, x) &= w_y \cdot f(x) \\ &= \sum_i \alpha_{i,y} K(x_i, x)\end{aligned}$$

- Like nearest neighbor – work with black-box similarities
- Downside: slow if many examples get nonzero alpha

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Kernels: Who Cares?

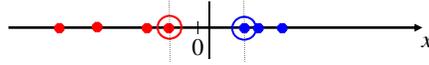
- So far: a very strange way of doing a very simple calculation
- “Kernel trick”: we can substitute any* similarity function in place of the dot product
- Lets us learn new kinds of hypothesis

* Fine print: if your kernel doesn't satisfy certain technical requirements, lots of proofs break. E.g. convergence, mistake bounds. In practice, illegal kernels *sometimes* work (but not always).

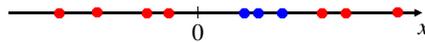
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Non-Linear Separators

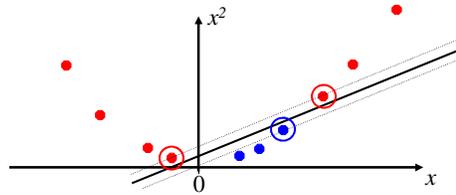
- Data that is linearly separable (with some noise) works out great:



- But what are we going to do if the dataset is just too hard?



- How about... mapping data to a higher-dimensional space:

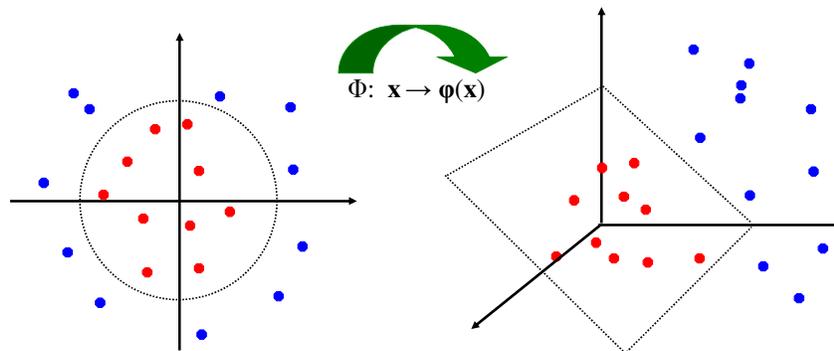


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This and next few slides adapted from Ray Mooney, UT

Non-Linear Separators

- General idea: the original feature space can always be mapped to some higher-dimensional feature space where the training set is separable:



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Some Kernels

- Kernels **implicitly** map original vectors to higher dimensional spaces, take the dot product there, and hand the result back

- Linear kernel:
$$K(x, x') = x \cdot x' = \sum_i x_i x'_i$$

$$\phi(x) = x$$

- Quadratic kernel:
$$K(x, x') = (x \cdot x' + 1)^2$$

$$= \sum_{i,j} x_i x_j x'_i x'_j + 2 \sum_i x_i x'_i + 1$$

For $x \in \mathbb{R}^3$:

$$\phi(x) = [x_1 x_1 \ x_1 x_2 \ x_1 x_3 \ x_2 x_1 \ x_2 x_2 \ x_2 x_3 \ x_3 x_1 \ x_3 x_2 \ x_3 x_3 \ \sqrt{2}x_1 \ \sqrt{2}x_2 \ \sqrt{2}x_3 \ 1]$$

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Some Kernels (2)

- Polynomial kernel:
$$K(x, x') = (x \cdot x' + 1)^d$$

For $x \in \mathbb{R}^3$:

$$\phi(x) = [x_1^d \ x_2^d \ x_3^d \ \sqrt{d}x_1^{d-1}x_2 \ \sqrt{d}x_1^{d-1}x_3 \ \dots \ \sqrt{d}x_1 \ \sqrt{d}x_2 \ \sqrt{d}x_3 \ 1]$$

For $x \in \mathbb{R}^n$ the d -order polynomial kernel's implicit feature space is $\binom{n+d}{d}$ dimensional.

By contrast, computing the kernel directly only requires $O(n)$ time.

Some Kernels (3)

- Kernels **implicitly** map original vectors to higher dimensional spaces, take the dot product there, and hand the result back
- Radial Basis Function (or Gaussian) Kernel: infinite dimensional representation

$$K(x, x') = \exp(-\|x - x'\|^2)$$

- Discrete kernels: e.g. string kernels
 - Features: all possible strings up to some length
 - To compute kernel: don't need to enumerate all substrings for each word, but only need to find strings appearing in both x and x'

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Why Kernels?

- Can't you just add these features on your own (e.g. add all pairs of features instead of using the quadratic kernel)?
 - Yes, in principle, just compute them
 - No need to modify any algorithms
 - But, number of features can get large (or infinite)
- Kernels let us compute with these features implicitly
 - Example: implicit dot product in polynomial, Gaussian and string kernel takes much less space and time per dot product
 - Of course, there's the cost for using the pure dual algorithms: you need to compute the similarity to every training datum

Recap: Classification

- Classification systems:
 - Supervised learning
 - Make a prediction given evidence
 - We've seen several methods for this
 - Useful when you have labeled data



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Where are we and what's left?

- So far foundations: Search, CSPs, Adversarial search, MDPs and RL, Bayes nets and probabilistic inference, Machine learning
- Tuesday: Applications in Robotics



- Thursday: Applications in Vision and Language + Conclusion + Where/How to learn more