Nearest Neighbor

Key idea: just store all training examples $\langle x_i, f(x_i) \rangle$

Need some <u>distance measure</u> between instances (e.g. Euclidean distance, Hamming distance)

Nearest neighbor:

• Given query instance x_q , first locate nearest training example x_n , then estimate $\hat{f}(x_q) = f(x_n)$

k-Nearest neighbor:

- Given xq, take vote among its k nearest neighbors (if discrete-valued target function)
 - Let \boldsymbol{k} not be divisible by number of possible labels
- Take mean of f values of k nearest neighbors if f real-valued

$$\hat{f}(x_q) = \frac{\sum_{i=1}^k f(x_i)}{k}$$

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Outline

• k-Nearest Neighbor

· Locally weighted regression

· Radial basis functions

· Case-based reasoning

· Lazy and eager learning

CSCE 478/878 Lecture 7: Instance-Based Learning

Stephen D. Scott (Adapted from Tom Mitchell's slides)

When To Consider Nearest Neighbor

- Instances are in \mathbb{R}^n (or, one can define some distance measure between instances; can use kernels)
- Less than 20 attributes per instance
 - To avoid <u>curse of dimensionality</u>, where many irrelevant attributes causes distance to be large, but distance is small if only relevant attributes used
 - Also, large number of attributes increases classification complexity
- Lots of training data

Advantages:

- Robust to noise
- Stable
- Training is very fast
- Learn complex target functions
- Don't lose information

Disadvantages:

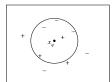
Slow at query time (active research area: fast indexing and accessing algorithms)

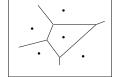
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· Easily fooled by irrelevant attributes

Voronoi Diagram

Decision surface for 1-NN





Nearest Neighbor's Behavior in the Limit

Consider p(x) defines probability that instance x will be labeled positive (versus negative)

Nearest neighbor (k = 1):

 \bullet As number of training examples $\to \infty,$ approaches Gibbs Algorithm

Recall Gibbs has at most twice the expected error of Bayes optimal

k-Nearest neighbor:

 As number of training examples → ∞ and k gets large, approaches Bayes optimal (best possible with given hyp. space and prior information)

Bayes optimal: if p(x) > 0.5 then predict +, else -

Distance-Weighted k-NN

Might want weight nearer neighbors more heavily:

$$\hat{f}(x_q) = \operatorname*{argmax}_{v \in V} \sum_{i=1}^k w_i \, \delta(v, f(x_i))$$

for discrete-valued $(\delta(v,f(x_i))=1$ if $v=f(x_i)$ and 0 otherwise), and

$$\hat{f}(x_q) = \frac{\sum_{i=1}^{k} w_i f(x_i)}{\sum_{i=1}^{k} w_i}$$

for continuous

where

$$w_i \equiv \frac{1}{d(x_q, x_i)^2}$$

and $d(x_q, x_i)$ is distance between x_q and x_i

Note now it makes sense to use all training examples instead of just k (Shepard's method), but then get increased time to classify instances

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Radial Basis Function (RBF) Networks

- Global approximation to target function, in terms of linear combination of local approximations
- Used, e.g., for image classification
- A different kind of neural network
- Closely related to distance-weighted regression, but "eager" instead of "lazy"

Curse of Dimensionality

Imagine instances described by 20 attributes, but only 2 are relevant to target function

<u>Curse of dimensionality</u>: nearest neighbor is easily misled by high-dimensional X

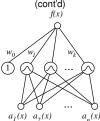
One approach:

- Stretch jth axis by weight z_j , where z_1, \ldots, z_n chosen to minimize prediction error
- Use cross-validation to automatically choose weights z_1, \ldots, z_n
- ullet Setting z_j to zero eliminates this dimension altogether see [Moore and Lee, 1994]

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RBF Networks (cont'd)



where $a_i(x)$ are the attributes describing instance x, and

$$\hat{f}(x) = w_0 + \sum_{i=1}^{k} w_i K_u(d(x_i, x))$$

(Note no weights from input to hidden layer)

One common choice for $K_u(d(x_u, x))$ is

$$K_u(d(x_u, x)) = \exp\left(-\frac{1}{2\sigma_u^2}d^2(x_u, x)\right),$$

i.e. Gaussian with mean at \boldsymbol{x}_u and variance $\sigma_u^2,$ all features independent

[note bug on p. 239]

Locally Weighted Regression

Note k-NN forms local approximation to f for each query point x_q

Why not form an explicit approximation $\hat{f}(x)$ for region surrounding x_a ?

- Fit linear, quadratic, etc. function to k nearest neighbors
- Produces "piecewise approximation" to f
- Do this for each new query point x_a

Several choices of error to minimize:

• Squared error over k nearest neighbors

$$E_1(x_q) \equiv \frac{1}{2} \sum_{x \in k \text{ nearest nbrs of } x_q} \left(f(x) - \hat{f}(x) \right)^2$$

• Distance-weighted squared error over all nbrs

$$E_2(x_q) \equiv \frac{1}{2} \sum_{x \in D} \left(f(x) - \hat{f}(x) \right)^2 K(d(x_q, x))$$

(*K* is decreasing in its argument)

• Combine E_1 and E_2

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Training Radial Basis Function Networks

- 1. Choose number of kernel functions (hidden units)
 - If = number training instances, can fit training data exactly by placing one center per instance
 - Using fewer ⇒ more efficient, less chance of overfitting
- 2. Choose center (= mean for Gaussian) x_u of kernel function $K_u(d(x_u,x))$
 - Use all training instances if enough kernels avail.
 - Use subset of training instances
 - Scatter uniformly throughout instance space
 - Can cluster data and assign one per cluster (helps answer step 1 also)
 - Can use EM to find means of mixture of Gaussians
 - Can also use e.g. EM to find σ_u 's (for Gaussian)
- Hold kernels fixed and train weights to fit linear function (output layer), e.g. GD or EG

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Case-Based Reasoning and CADET

Can apply instance-based learning even when \boldsymbol{X} much more complex

Need different "distance" metric

Case-Based Reasoning is instance-based learning where instances have symbolic logic descriptions

```
((user-complaint error53-on-shutdown)
  (cpu-model PowerPC) (operating-system Windows
  (memory 48meg)
  (installed-apps Excel Netscape VirusScan)
  (disk lgig)
  (likely-cause ???))
```

<u>CADET</u>: 75 stored examples of mechanical devices, e.g. water faucets

- Training instance: \(\)qualitative function, mech. structure \(\)
- · New query: desired function
- Target value: mechanical structure for this function

Distance metric: match qualitative function descriptions

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Lazy and Eager Learning

Lazy: Wait for query before generalizing

 k-NN, locally weighted regression, case-based reasoning

Eager: Generalize before seeing query

 Radial basis function networks, ID3, backpropagation, naïve Bayes

Does it matter?

- Computation time for training and generalization
- Eager learner must create global approximation, lazy learner can create many local approximations
- If they use same H, lazy can represent more complex functions (e.g. consider H = linear functions) since it considers the query instance x_q before generalizing, i.e. lazy produces a new hypothesis for each new x_q

Case-Based Reasoning in CADET

Example

A stored case: T-junction pipe

Structure: Function: $Q_1.T_1 \qquad T = \text{temperature} \qquad Q_1 \sim Q_1 \sim Q_2 \sim Q_3.T_3 \qquad Q_2 \sim Q_3.T_3 \sim Q_2.T_1 \sim Q_2.T_2 \sim Q_3.T_2 \sim Q_2.T_2 \sim Q_3.T_2 \sim Q_3.T_3 \sim Q_3.T_$

A problem specification: Water faucet

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E.g. distance measure = size of largest isomorphic subgraph

Case-Based Reasoning in CADET

(cont'd)

- Instances represented by rich structural (symbolic) descriptions, vs. e.g. points in \mathbb{R}^n for k-NN
- Multiple cases retrieved (and combined) to form solution to new problem: Similar to k-NN, except combination procedure can rely on knowledge-based reasoning (e.g. can two components be fit together?)
- Tight coupling between case retrieval, knowledge-based reasoning, and problem solving, e.g. application of rewrite rules in function graphs and backtracking in search space

Bottom line:

- Simple matching of cases useful for tasks such as answering help-desk queries
- Area of ongoing research, including improving indexing and search methods

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