

## Instance Based Learning

- $k$ -Nearest Neighbor
- Locally weighted regression
- Radial basis functions
- Case-based reasoning
- Lazy and eager learning

## Instance-Based Learning

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

Nearest neighbor (1- Nearest neighbor) :

- Given query instance  $x_q$ , locate nearest example  $x_n$ , estimate

$$\hat{f}(x_q) \leftarrow f(x_n)$$

$k$  - Nearest neighbor :

- Given  $x_q$ , take vote among its  $k$  nearest neighbors (if discrete - valued target function)
- Take mean of  $f$  values of  $k$  nearest neighbors (if real - valued)

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

## When to Consider Nearest Neighbor

- Instance map to points in  $R^n$
- Less than 20 attributes per instance
- Lots of training data

Advantages

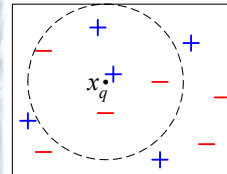
- **Training is very fast**
- Learn complex target functions
- Do not lose information

Disadvantages

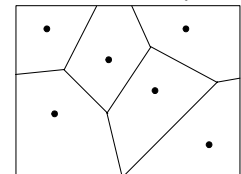
- **Slow at query time**
- Easily fooled by irrelevant attributes

## $k$ -NN Classification

5-Nearest Neighbor



1-NN Decision Surface



## Behavior in the Limit

Define  $p(x)$  as probability that instance  $x$  will be labeled 1 (positive) versus 0 (negative)

Nearest Neighbor

- As number of training examples approaches infinity, approaches Gibbs Algorithm
- Gibbs: with probability  $p(x)$  predict 1, else 0

$k$ -Nearest Neighbor:

- As number of training examples approaches infinity and  $k$  gets large, approaches Bayes optimal
- Bayes optimal: if  $p(x) > 0.5$  then predict 1, else 0
- Note Gibbs has at most twice the expected error of Bayes optimal

## Distance-Weighted $k$ -NN

Might want to weight nearer neighbors more heavily ...

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

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

## Curse of Dimensionality

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

**Curse of dimensionality:** nearest neighbor is easily misled when high-dimensional  $X$

One approach:

- Stretch  $j$ th axis by weight  $z_j$ , where  $z_1, z_2, \dots, z_n$  chosen to minimize prediction error
- Use cross-validation to automatically choose weights  $z_1, z_2, \dots, z_n$
- Note setting  $z_j$  to zero eliminates dimension  $j$  altogether see (Moore and Lee, 1994)

## Locally Weighted Regression

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

Why not form explicit approximation  $\hat{f}(x)$  for region around  $x_q$ ?

- Fit linear function to  $k$  nearest neighbors
- Or fit quadratic, etc.
- Produces "piecewise approximation" to  $f$

Several choices of error to minimize:

- Squared error over  $k$  nearest neighbors

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

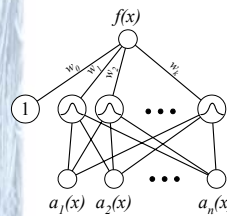
- Distance - weighted squared error over all neighbors

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

## Radial Basis Function Networks

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

## Radial Basis Function Networks



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

$$f(x) = w_0 + \sum_{u=1}^k w_u K_u(d(x_u, x))$$

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

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

## Training RBF Networks

Q1: What  $x_u$  to use for kernel function  $K_u(d(x_u, x))$ ?

- Scatter uniformly through instance space
- Or use training instances (reflects instance distribution)

Q2: How to train weights (assume here Gaussian  $K_u$ )?

- First choose variance (and perhaps mean) for each  $K_u$ 
  - e.g., use EM
- Then hold  $K_u$  fixed, and train linear output layer
  - efficient methods to fit linear function

## Case-Based Reasoning

Can apply instance-based learning even when  $X \neq R^n$

→ need different "distance" metric

Case-Based Reasoning is instance-based learning applied to instances with symbolic logic descriptions:

```
((user-complaint error53-on-shutdown)
```

```
(cpu-model PowerPC)
```

```
(operating-system Windows)
```

```
(network-connection PCIA)
```

```
(memory 48meg)
```

```
(installed-applications Excel Netscape
```

```
VirusScan)
```

```
(disk 1Gig)
```

```
(likely-cause ???)
```

## Case-Based Reasoning in CADET

CADET: 75 stored examples of mechanical devices

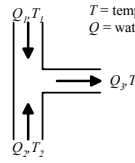
- each training example:
  - <qualitative function, mechanical structure>
- new query: desired function
- target value: mechanical structure for this function

Distance metric: match qualitative function descriptions

## Case-Based Reasoning in CADET

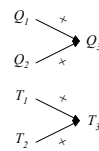
A stored case: T-junction pipe

Structure:



$T$  = temperature  
 $Q$  = waterflow

Function:

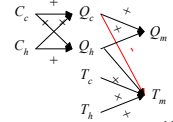


A problem specification: Water faucet

Structure:

?

Function:



## Case-Based Reasoning in CADET

- Instances represented by rich structural descriptions
- Multiple cases retrieved (and combined) to form solution to new problem
- Tight coupling between case retrieval and problem solving

Bottom line:

- Simple matching of cases useful for tasks such as answering help-desk queries
- Area of ongoing research

## Lazy and Eager Learning

**Lazy:** wait for query before generalizing

- k-Nearest Neighbor, Case-Based Reasoning

**Eager:** generalize before seeing query

- Radial basis function networks, ID3, Backpropagation, etc.

Does it matter?

- 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)

## kd-trees (Moore)

- *Eager* version of  $k$ -Nearest Neighbor
- Idea: decrease time to find neighbors
  - train by constructing a lookup ( $kd$ ) tree
  - recursively subdivide space
    - ignore class of points
    - lots of possible mechanisms: grid, maximum variance, etc.
  - when looking for nearest neighbor search tree
  - nearest neighbor can be found in  $\log(n)$  steps
  - $k$  nearest neighbors can be found by generalizing process (still in  $\log(n)$  steps if  $k$  is constant)
- Slower training but faster classification

## kd Tree

