Memory-Based Learning Instance-Based Learning K-Nearest Neighbor

Motivating Problem

Inductive Assumption

Similar inputs map to similar outputs

- If not true => learning is impossible
- If true => learning reduces to defining "similar"

Not all similarities created equal

 predicting a person's weight may depend on different attributes than predicting their IQ

1-Nearest Neighbor

$$Dist(c_{1}, c_{2}) = \sqrt{\sum_{i=1}^{N} \left(attr_{i}(c_{1}) - attr_{i}(c_{2})\right)^{2}}$$

$$NearestNeighbor = MIN_{j}(Dist(c_{j}, c_{test}))$$

$$prediction_{test} = class_{j} (or value_{j})$$

works well if no attribute or class noise as number of training cases grows large, error rate of 1-NN is at most 2 times the Bayes optimal rate

k-Nearest Neighbor

$$Dist(c_1, c_2) = \sqrt{\sum_{i=1}^{N} \left(attr_i(c_1) - attr_i(c_2)\right)^2}$$

$$k - NearestNeighbors = \left\{k - MIN(Dist(c_i, c_{test}))\right\}$$

$$prediction_{test} = \frac{1}{k} \sum_{i=1}^{k} class_i \ (or \frac{1}{k} \sum_{i=1}^{k} value_i)$$

Average of k points more reliable when:

- noise in attributes
- noise in class labels
- classes partially overlap

How to choose "k"

Large k:

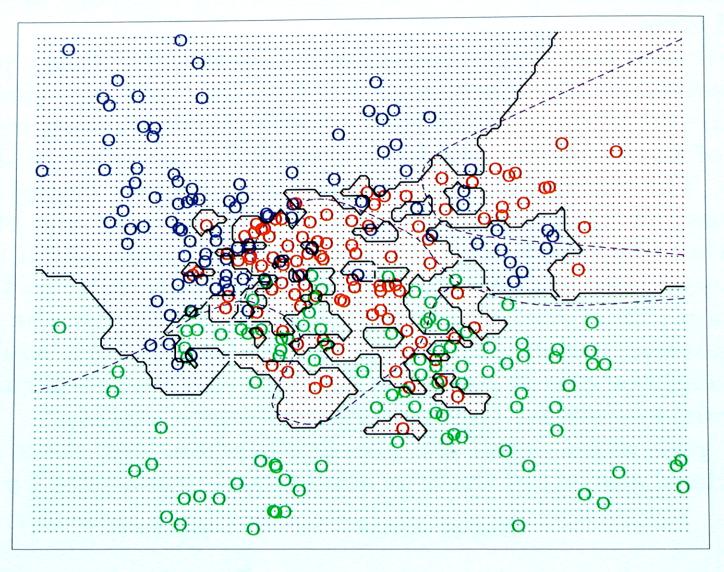
- less sensitive to noise (particularly class noise)
- better probability estimates for discrete classes
- larger training sets allow larger values of k

Small k:

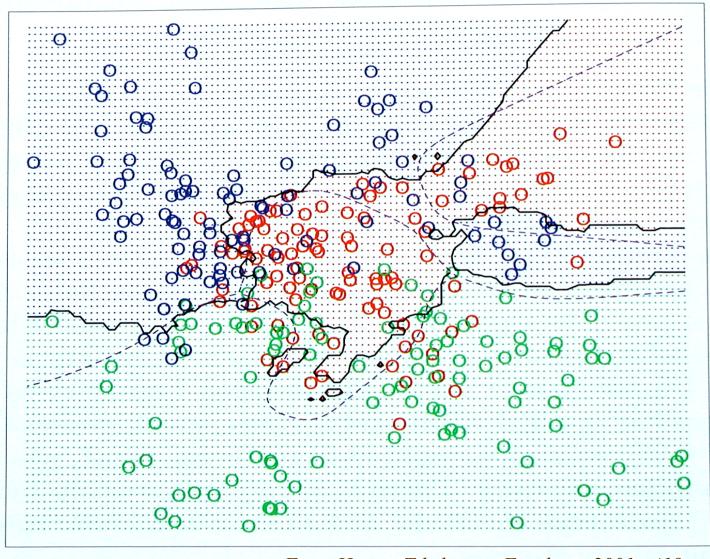
- captures fine structure of space better
- may be necessary with small training sets

Balance must be struck between large and small k As training set approaches infinity, and k grows large, kNN becomes Bayes optimal

1-Nearest Neighbor



15-Nearest Neighbors



From Hastie, Tibshirani, Friedman 2001 p418

Cross-Validation

Models usually perform better on training data than on future test cases

1-NN is 100% accurate on training data!

Leave-one-out-cross validation:

- "remove" each case one-at-a-time
- use as test case with remaining cases as train set
- average performance over all test cases

LOOCV is impractical with most learning methods, but extremely efficient with MBL!

Distance-Weighted kNN

tradeoff between small and large k can be difficult

– use large k, but more emphasis on nearer neighbors?

$$prediction_{test} = \frac{\sum_{i=1}^{k} w_i * class_i}{\sum_{i=1}^{k} w_i} (or \frac{\sum_{i=1}^{k} w_i * value_i}{\sum_{i=1}^{k} w_i})$$

$$w_k = \frac{1}{Dist(c_k, c_{test})}$$

Locally Weighted Averaging

Let k = number of training points Let weight fall-off rapidly with distance

$$prediction_{test} = \frac{\sum_{i=1}^{k} w_i * class_i}{\sum_{i=1}^{k} w_i} (or \frac{\sum_{i=1}^{k} w_i * value_i}{\sum_{i=1}^{k} w_i})$$

$$w_k = \frac{1}{e^{KernelWidth \cdot Dist(c_k, c_{test})}}$$

KernelWidth controls size of neighborhood that has large effect on value (analogous to k)

Locally Weighted Regression

All algs so far are strict averagers: interpolate, but can't extrapolate

Do weighted regression, centered at test point, weight controlled by distance and KernelWidth

Local regressor can be linear, quadratic, n-th degree polynomial, neural net, ...

Yields piecewise approximation to surface that typically is more complex than local regressor

Euclidean Distance

$$D(c1, c2) = \sqrt{\sum_{i=1}^{N} (attr_{i}(c1) - attr_{i}(c2))^{2}}$$

gives all attributes equal weight?

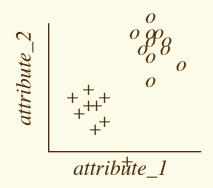
- only if scale of attributes and differences are similar
- scale attributes to equal range or equal variance

assumes spherical classes

$$attribute_{-1}$$

$$attribute_{-1}$$

Euclidean Distance?



if classes are not spherical?

if some attributes are more/less important than other attributes?

if some attributes have more/less noise in them than other attributes?

Weighted Euclidean Distance

$$D(c1,c2) = \sqrt{\sum_{i=1}^{N} w_i \cdot \left(attr_i(c1) - attr_i(c2)\right)^2}$$

large weights => attribute is more important small weights => attribute is less important zero weights => attribute doesn't matter

Weights allow kNN to be effective with elliptical classes Where do weights come from?

Learning Attribute Weights

Scale attribute ranges or attribute variances to make them uniform (fast and easy)

Prior knowledge

Numerical optimization:

- gradient descent, simplex methods, genetic algorithm
- criterion is cross-validation performance

Information Gain of single attributes

Information Gain

Information Gain = reduction in entropy due to splitting on an attribute

Entropy = expected number of bits needed to encode the class of a randomly drawn + or – example using the optimal info-theory coding

$$Entropy = -p_{+} \log_2 p_{+} - p_{-} \log_2 p_{-}$$

$$Gain(S, A) = Entropy(S) - \sum_{v \in Values(A)} \frac{\left|S_{v}\right|}{\left|S\right|} Entropy(S_{v})$$

Booleans, Nominals, Ordinals, and Reals

Consider attribute value differences:

 $(attr_i(c1) - attr_i(c2))$

Reals: easy! full continuum of differences

Integers: not bad: discrete set of differences

Ordinals: not bad: discrete set of differences

Booleans: awkward: hamming distances 0 or 1

Nominals? not good! recode as Booleans?

Curse of Dimensionality

as number of dimensions increases, distance between points becomes larger and more uniform if number of relevant attributes is fixed, increasing the number of less relevant attributes may swamp distance

$$D(c1,c2) = \sqrt{\sum_{i=1}^{relevant} \left(attr_i(c1) - attr_i(c2)\right)^2 + \sum_{j=1}^{irrelevant} \left(attr_j(c1) - attr_j(c2)\right)^2}$$

when more irrelevant than relevant dimensions, distance becomes less reliable

solutions: larger k or KernelWidth, feature selection, feature weights, more complex distance functions

Advantages of Memory-Based Methods

Lazy learning: don't do any work until you know what you want to predict (and from what variables!)

- never need to learn a global model
- many simple local models taken together can represent a more complex global model
- better focussed learning
- handles missing values, time varying distributions, ...
- Very efficient cross-validation
- Intelligible learning method to many users
- Nearest neighbors support explanation and training
- Can use any distance metric: string-edit distance, ...

Weaknesses of Memory-Based Methods

Curse of Dimensionality:

often works best with 25 or fewer dimensions

Run-time cost scales with training set size

Large training sets will not fit in memory

Many MBL methods are strict averagers

Sometimes doesn't seem to perform as well as other methods such as neural nets

Predicted values for regression not continuous

Combine KNN with ANN

Train neural net on problem

Use outputs of neural net or hidden unit activations as new feature vectors for each point

Use KNN on new feature vectors for prediction

Does feature selection and feature creation

Sometimes works better than KNN or ANN

Current Research in MBL

Condensed representations to reduce memory requirements and speed-up neighbor finding to scale to 10⁶–10¹² cases

Learn better distance metrics

Feature selection

Overfitting, VC-dimension, ...

MBL in higher dimensions

MBL in non-numeric domains:

- Case-Based Reasoning
- Reasoning by Analogy

References

Locally Weighted Learning by Atkeson, Moore, Schaal Tuning Locally Weighted Learning by Schaal, Atkeson, Moore

Closing Thought

In many supervised learning problems, all the information you ever have about the problem is in the training set.

Why do most learning methods discard the training data after doing learning?

Do neural nets, decision trees, and Bayes nets capture *all* the information in the training set when they are trained?

In the future, we'll see more methods that combine MBL with these other learning methods.

- to improve accuracy
- for better explanation
- for increased flexibility