# CS4442/9542b Artificial Intelligence II prof. Olga Veksler

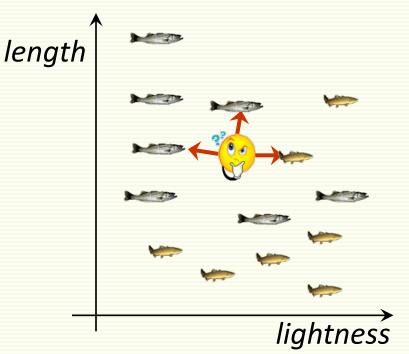
# Lecture 3 Machine Learning K Nearest Neighbor Classifier

# Today

- kNN classifier the simplest classifier on earth
- matlab implementation of kNN

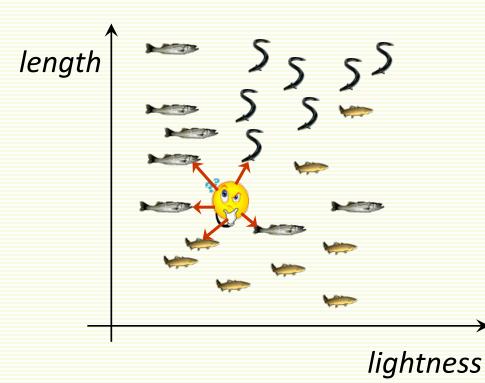
## k-Nearest Neighbors

- classify an unknown example with the most common class among k closest examples
  - "tell me who your neighbors are, and I'll tell you who you are"
- Example:
  - *k* = 3
  - 2 sea bass, 1 salmon
  - Classify as sea bass



# kNN: Multiple Classes

- Easy to implement for multiple classes
- Example for *k* = 5
  - 3 fish species: salmon, sea bass, eel
  - 3 sea bass, 1 eel, 1 salmon  $\Rightarrow$  classify as sea bass

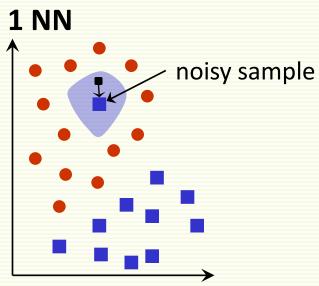


# kNN: How to Choose k?

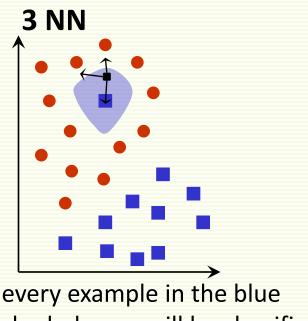
- In theory, if infinite number of samples available, the larger is *k*, the better is classification
- But the caveat is that all k neighbors have to be close
  - Possible when infinite # samples available
  - Impossible in practice since # samples is finite

# kNN: How to Choose k?

- Rule of thumb is *k* = sqrt(*n*), *n* is number of examples
  - interesting theoretical properties
- In practice, *k* = 1 is often used for efficiency, but can be sensitive to "noise"

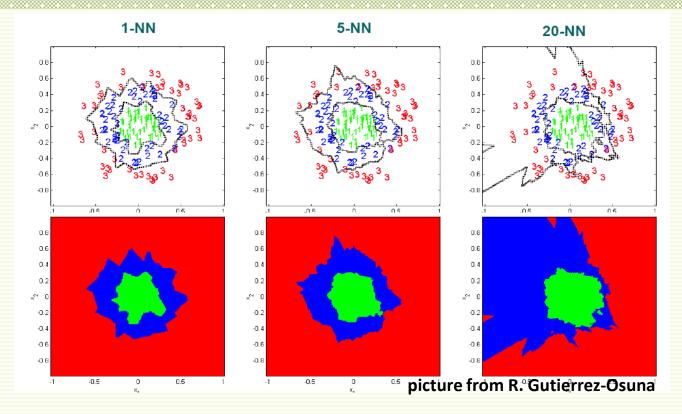


every example in the blue shaded area will be misclassified as the blue class



shaded area will be classified correctly as the red class

## kNN: How to Choose k?



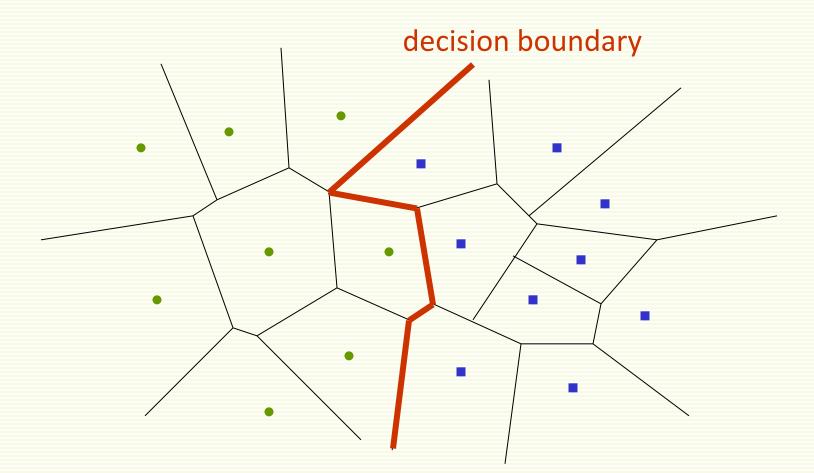
- Larger **k** gives smoother boundaries, better for generalization
  - But only if *locality* is preserved. Locality is not preserved if end up looking at samples too far away, not from the same class.
- Interesting theoretical properties if k < sqrt(n), n is # of examples</li>
- Can choose k through cross-validation (study soon)

# kNN: How Well does it Work?

- kNN is simple and intuitive, but does it work?
- Theoretically, the best error rate is the Bayes rate **E**\*
  - Bayes error rate is the best (smallest) error rate a classifier can have, for a given problem, but we do not study it in this course
- Assume we have an unlimited number of samples
- kNN leads to an error rate greater than E\*
- But even for k =1, as n → ∞, it can be shown that kNN error rate is smaller than 2E\*
- As we increase k, the upper bound on the error gets better, that is the error rate (as n→∞) for the kNN rule is smaller than cE\*, with smaller c for larger k
- If we have lots of samples, kNN works well

#### **1NN Visualization**

#### Voronoi tesselation is useful for visualization



### kNN Selection of Distance

• So far we assumed we use Euclidian Distance to find the nearest neighbor:

$$D(a,b) = \sqrt{\sum_{k} (a_k - b_k)^2}$$

- Euclidean distance treats each feature as equally important
- However some features (dimensions) may be much more discriminative than other features

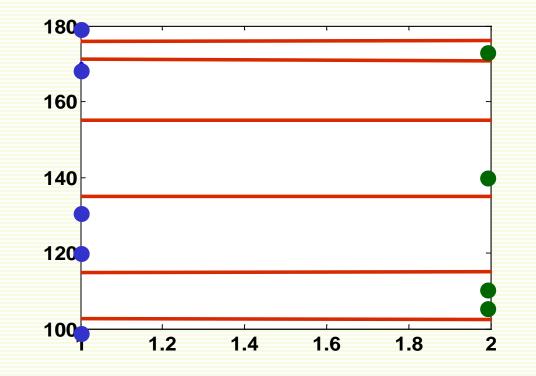
#### kNN Distance Selection: Extreme Example

- feature 1 gives the correct class: 1 or 2
- feature 2 gives irrelevant number from 100 to 200
- dataset: [1 150]
  - [2 110]
- classify **[1 100]**

$$D(\begin{bmatrix} 1\\100 \end{bmatrix}, \begin{bmatrix} 1\\150 \end{bmatrix}) = \sqrt{(1-1)^2 + (100-150)^2} = 50$$
$$D(\begin{bmatrix} 1\\100 \end{bmatrix}, \begin{bmatrix} 2\\110 \end{bmatrix}) = \sqrt{(1-2)^2 + (100-110)^2} = 10.5$$

- [1 100] is misclassified!
- The denser the samples, the less of this problem
- But we rarely have samples dense enough

#### kNN Distance Selection: Extreme Example



• Decision boundary is in red, and is really wrong because

- feature 1 is discriminative, but it's scale is small
- feature 2 gives no class information but its scale is large, it dominates distance calculation

- Notice that 2 features are on different scales:
- First feature takes values between 1 or 2
- Second feature takes values between 100 to 200
- Idea: normalize features to be on the same scale
- Different normalization approaches
- Linearly scale the range of each feature to be, say, in range [0,1]

$$f_{new} = \frac{f_{old} - f_{old}^{\min}}{f_{old}^{\max} - f_{old}^{\min}}$$

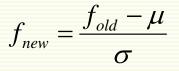
- Linearly scale to **0** mean variance **1**:
- If Z is a random variable of mean μ and variance σ<sup>2</sup>, then (Z μ)/σ has mean 0 and variance 1
- For each feature **f** let the new rescaled feature be

$$f_{new} = \frac{f_{old} - \mu}{\sigma}$$

• C is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously

**C**<sub>new</sub>=(**C**-repmat(mean(**C**), size(**C**,1), 1))\*diag(1./std(**C**))

 C is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously
 C<sub>new</sub>=(C-repmat(mean(C), size(C,1), 1))\*diag(1./std(C))



 $C = \begin{bmatrix} 1 & 180 \\ 1 & 100 \\ 1 & 160 \\ 2 & 120 \\ 2 & 150 \\ 2 & 170 \end{bmatrix}$  size(C,1) = 6

• **C** is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously

**C**<sub>new</sub>=(**C**-repmat(mean(**C**), **6**, 1))\*diag(1./std(**C**))

$$C = \begin{bmatrix} 1 & 180 \\ 1 & 100 \\ 1 & 160 \\ 2 & 120 \\ 2 & 150 \\ 2 & 170 \end{bmatrix}$$

$$f_{new} = \frac{f_{old} - \mu}{\sigma}$$

• **C** is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously

 $\boldsymbol{C}_{new} = (\boldsymbol{C} - repmat(mean(\boldsymbol{C}), 6, 1)) * diag(1./std(\boldsymbol{C}))$ 

$$C = \begin{bmatrix} 1.5 & 146.7 \end{bmatrix}$$

$$C = \begin{bmatrix} 1 & 180 \\ 1 & 100 \\ 1 & 160 \\ 2 & 120 \\ 2 & 150 \\ 2 & 170 \end{bmatrix}$$

$$f_{new} = \frac{f_{old} - \mu}{\sigma}$$

### repmat Function

$$repmat \left( \begin{bmatrix} 1 & 2 \\ 3 & 4 \end{bmatrix}, 2, 3 \right) = \begin{bmatrix} 1 & 2 \\ 3 & 4 \end{bmatrix} \begin{bmatrix} 1 & 2$$

• **C** is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously

 $f_{new} = \frac{f_{old} - \mu}{\sigma}$ 

C\_new=(C-repmat(mean(C), 6, 1))\*diag(1./std(C))

1.5	146.7
1.5	146.7
1.5	146.7
1.5	146.7
1.5	146.7
1.5	146.7

mean(*C*) = [1.5 146.7]

 $f_{new} = \frac{f_{old} - \mu}{\sigma}$ 

• **C** is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously

 $C_{new} = (C - repmat(mean(C), 6, 1))*diag(1./std(C))$ 

<b>[</b> 1	180	1.5	146.7	-0.5	33.3
1	100	1.5	146.7	-0.5	-46.7
1	160	1.5	146.7	-0.5	13.3
2	120	 1.5	146.7	-0.5	-26.7
2	150	1.5	146.7	-0.5	3.3
2	170	1.5	146.7	1.5	23.3

• C is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously

Cnew = (C-repmat(mean(C), 6, 1))\*diag(1./std(C))

 $f_{new} = \frac{f_{old} - \mu}{\sigma}$ 

std(*C*) = [0.55 30.8]

$$C = \begin{bmatrix} 1 & 180 \\ 1 & 100 \\ 1 & 160 \\ 2 & 120 \\ 2 & 150 \\ 2 & 170 \end{bmatrix}$$

• C is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously

Cnew = (C-repmat(mean(C), 6, 1))\*diag(1./std(C))

1./std(*C*) = [ 1.83 0.03 ]

 $f_{new} = \frac{f_{old} - \mu}{\sigma}$ 

std(*C*) = [0.55 30.8]

• **C** is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously

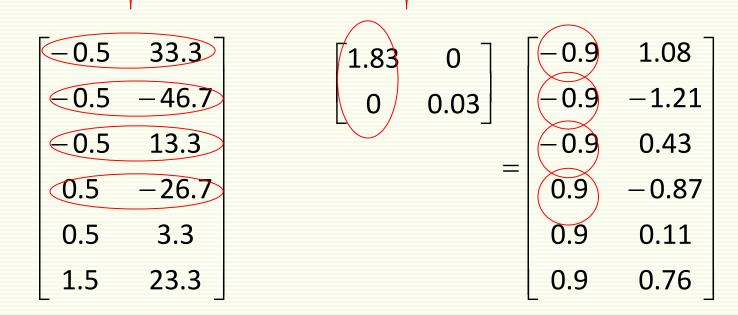
Cnew = (C-repmat(mean(C), 6, 1))\*diag(1./std(C))

 $f_{new} = \frac{f_{old} - \mu}{\tau}$ 

1./std(*C*) = [ 1.83 0.03 ]

• **C** is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously

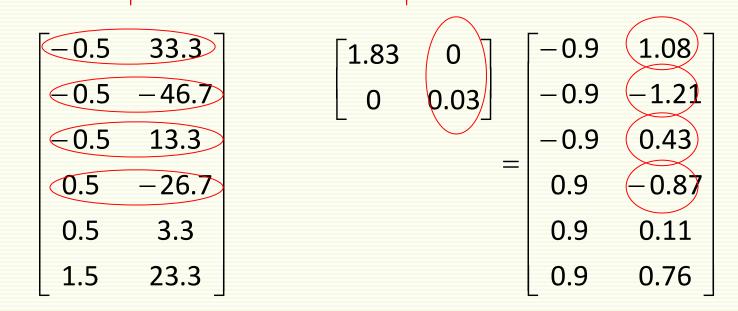
Cnew = (C-repmat(mean(C), 6, 1))\*diag(1./std(C))



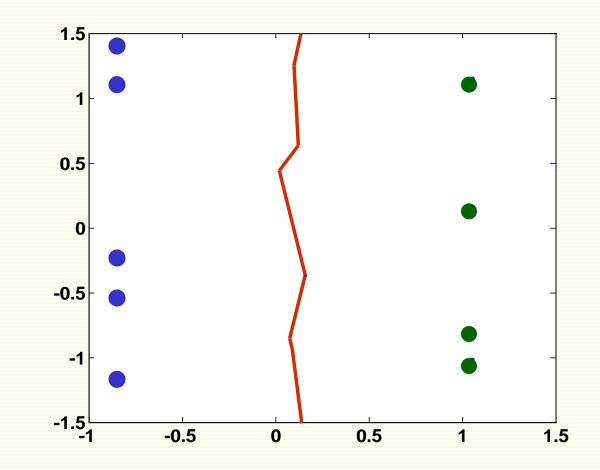
 $f_{new} = \frac{J_{old} - \mu}{2}$ 

• **C** is a matrix with all samples stored as rows, in Matlab can normalize all features simultaneously

Cnew = (C-repmat(mean(C), 6, 1))\*diag(1./std(C))



 $f_{new} = \frac{J_{old} - \mu}{2}$ 



## kNN: Selection of Distance

 Feature normalization does not help in high dimensional spaces if most features are irrelevant

$$D(a,b) = \sqrt{\sum_{k} (a_{k} - b_{k})^{2}} = \sqrt{\sum_{i} (a_{i} - b_{i})^{2} + \sum_{j} (a_{j} - b_{j})^{2}}$$
  
discriminative noisy  
features features

 If the number of useful features is smaller than the number of noisy features, Euclidean distance is dominated by noise

## kNN: Feature Weighting

• Scale each feature by its importance for classification

$$D(a,b) = \sqrt{\sum_{k} w_{k} (a_{k} - b_{k})^{2}}$$

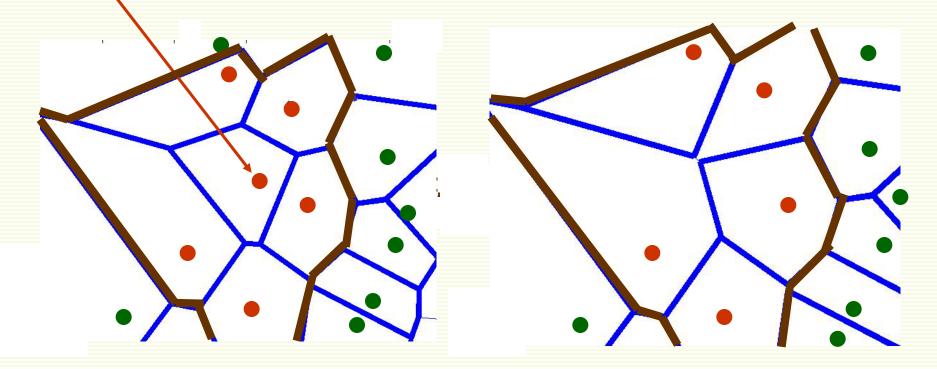
- Can use our prior knowledge about which features are more important
- Can learn the weights w<sub>k</sub> using cross-validation (to be covered later)

# kNN: Computational Complexity

- Basic kNN algorithm stores all examples
- Suppose we have **n** examples each of dimension **d**
- **O(d)** to compute distance to one example
- O(nd) to find distances to all examples
- **O(knd)** to find **k** closest examples examples
  - O(nd) + O(kn) if careful
- Thus total complexity is **O(knd)**
- Very expensive for a large number of samples
- But we need a large number of samples for kNN to work well!

# Reducing Complexity: editing 1NN

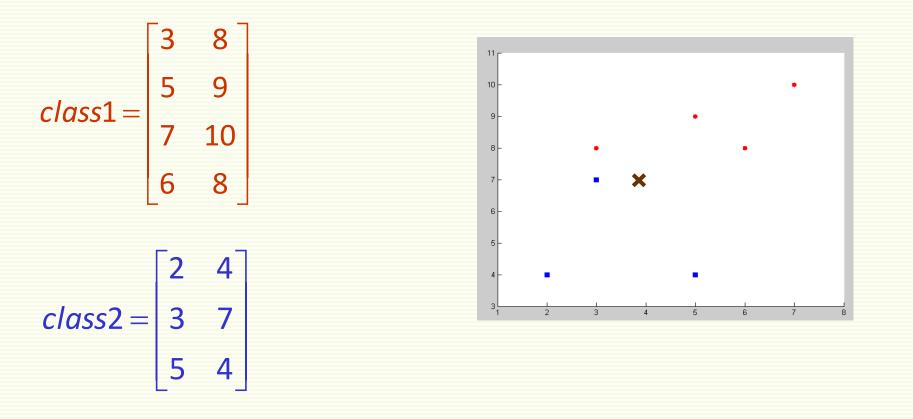
 If all Voronoi neighbors have the same class, a sample is useless, remove it



- Number of samples decreases
- Decision boundary does not change

# **Reducing Complexity: Partial Distance**

- Have current *k* closes samples
- Abort distance computation if partial distance is already greater than the full distance to the current k closest samples
- Advantages:
  - complexity decreases
  - we are guaranteed to find closes neighbor(s)
- Disadvantages:
  - how much complexity decreases depends on our luck and data layout



• Want to classify **newSample** = [4 7]

### kNN in Matlab without Loops

```
numClass1 = size(Class1,1);
numClass2
             = size(Class2,1);
totalSamples = numClass1+numClass2;
combinedSamples = [Class1;Class2];
trueClass = [zeros(numClass1,1)+1;zeros(numClass2,1)+2];
testMatrix = repmat(newSample,totalSamples,1);
absDiff = abs(combinedSamples-testMatrix);
absDiff = absDiff.^2;
dist = sum(absDiff,2);
[Y,I] = sort(dist);
neighborsInd = I(1:k);
neighbors = trueClass(neighborsInd);
class1 = find(neighbors == 1);
class2 = find(neighbors == 2);
      = [size(class1,1);size(class2,1)];
ioint
[value class] = max(joint);
```

 $class1 = \begin{bmatrix} 3 & 8 \\ 5 & 9 \\ 7 & 10 \\ 6 & 8 \end{bmatrix}$  $class2 = \begin{bmatrix} 2 & 4 \\ 3 & 7 \\ 5 & 4 \end{bmatrix}$ 

```
newSamle = \begin{bmatrix} 4 & 7 \end{bmatrix}
```

$$k = 3$$

<pre>numClass1 = size(Class1,1); numClass2 = size(Class2,1); totalSamples = numClass1+numClass2; combinedSamples = [Class1;Class2]; trueClass = [zeros(numClass1,1)+1;zeros(numClass2,1)+2];</pre>	class1 =	3 5 7 6	8 9 10 8	
		2	4	
numClass1 = 4	class2 =	3	7	
numClass2 = 3		_5	4_	
totalSamples = 7	newSamp	ole=	4	7]
$combinedSamples = \begin{bmatrix} 3 & 8 \\ 5 & 9 \\ 7 & 10 \\ 6 & 8 \\ 2 & 4 \\ 3 & 7 \\ 5 & 4 \end{bmatrix} trueClass = \begin{bmatrix} 1 \\ 1 \\ 1 \\ 2 \\ 2 \\ 2 \end{bmatrix}$				

<pre>testMatrix = repmat(newSample,totalSamples,1); absDiff = abs(combinedSamples-testMatrix); absDiff = absDiff.^2; dist = sum(absDiff,2);</pre>	newSample=[4 7]
$testMatrix = \begin{bmatrix} 4 & 7 \\ 4 & 7 \\ 4 & 7 \\ 4 & 7 \\ 4 & 7 \\ 4 & 7 \\ 4 & 7 \\ 4 & 7 \\ 4 & 7 \end{bmatrix}  absDiff = \begin{bmatrix} 1 & 1 \\ 1 & 2 \\ 3 & 3 \\ 2 & 1 \\ 2 & 3 \\ 1 & 0 \\ 1 & 3 \end{bmatrix}$	$combinedSamples = \begin{bmatrix} 3 & 8 \\ 5 & 9 \\ 7 & 10 \\ 6 & 8 \\ 2 & 4 \\ 3 & 7 \\ 5 & 4 \end{bmatrix}$
$absDiff = \begin{bmatrix} 1 & 1 \\ 1 & 4 \\ 9 & 9 \\ 4 & 1 \\ 4 & 9 \\ 1 & 0 \\ 1 & 9 \end{bmatrix} \qquad dist = \begin{bmatrix} 2 \\ 5 \\ 18 \\ 5 \\ 13 \\ 1 \\ 10 \end{bmatrix}$	$trueClass = \begin{bmatrix} 1\\1\\1\\1\\2\\2\\2\\2\end{bmatrix}$

[Y,I] = sort(dist neighborsInd = I(1:k); neighbors = trueClas	t); ss(neighborsInd);	dist = $\begin{bmatrix} 2 \\ 5 \\ 18 \\ 5 \\ 13 \\ 1 \end{bmatrix}$
$\mathbf{Y} = \begin{bmatrix} 1\\2\\5\\5\\10\\13\\18 \end{bmatrix}$	$I = \begin{bmatrix} 6\\1\\2\\4\\7\\5\\3 \end{bmatrix}$	$\begin{bmatrix} 1\\ 1\\ 1\\ 1\\ 1\\ 2\\ 2\\ 2 \end{bmatrix}$
$neighborsInd = \begin{bmatrix} 6\\1\\2\end{bmatrix}$	$neighbors = \begin{bmatrix} 2\\ 1\\ 1 \end{bmatrix}$	$\begin{bmatrix} 2\\2 \end{bmatrix}$ k = 3

class1 = find(neighbors == 1); class2 = find(neighbors == 2); joint = [size(class1,1);size(class2,1)];

[value class] = max(joint);

$$class1 = \begin{bmatrix} 2 \\ 3 \end{bmatrix}$$
$$class2 = \begin{bmatrix} 1 \end{bmatrix}$$
$$joint = \begin{bmatrix} 2 \\ 1 \end{bmatrix}$$

class = 1

 $neighbors = \begin{bmatrix} 2 \\ 1 \\ 1 \end{bmatrix}$ 

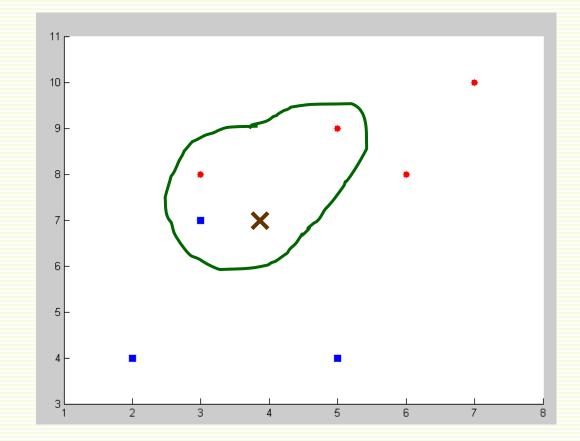
class1 = find(neighbors == 1); class2 = find(neighbors == 2); joint = [size(class1,1);size(class2,1)];

[value class] = max(joint);

Also can use class = mode(neighbors) instead

$$class1 = \begin{bmatrix} 2\\3 \end{bmatrix}$$
$$class2 = \begin{bmatrix} 1\\1 \end{bmatrix}$$
$$joint = \begin{bmatrix} 2\\1 \end{bmatrix}$$
$$class = 1$$

neighbors =  $\begin{bmatrix} 2 \\ 1 \\ 1 \end{bmatrix}$ 



# Video

#### <u>http://videolectures.net/aaai07\_bosch\_knnc/</u>

# kNN Summary

- Advantages
  - Can be applied to the data from any distribution
    - for example, data does not have to be separable with a linear boundary
  - Very simple and intuitive
  - Good classification if the number of samples is large enough
- Disadvantages
  - Choosing k may be tricky
  - Test stage is computationally expensive
    - No training stage, all the work is done during the test stage
    - This is actually the opposite of what we want. Usually we can afford training step to take a long time, but we want fast test step
  - Need large number of samples for accuracy