Learning From Data
Lecture 17
Memory and Efficiency in Nearest Neighbor

Memory
Efficiency

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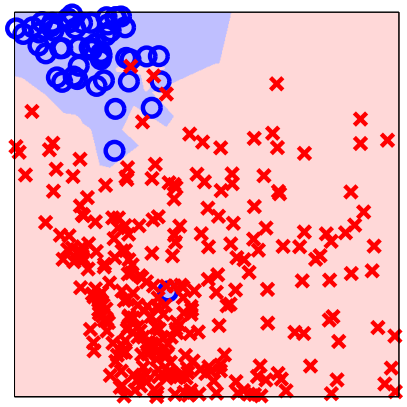
CSCI 4100/6100

RECAP: Similarity and Nearest Neighbor

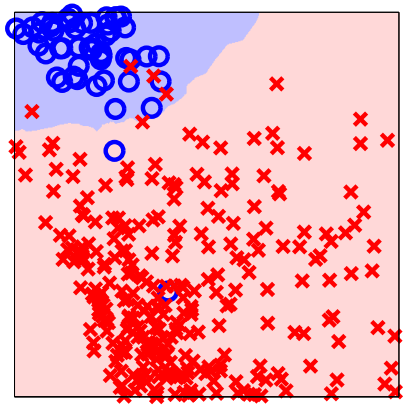
Similarity

$$d(\mathbf{x}, \mathbf{x}') = \| \mathbf{x} - \mathbf{x}' \|$$

1-NN rule



21-NN rule



1. Simple.
2. No training.
3. Near optimal E_{out}:
$$k \to \infty,\ k/N \to 0 \implies E_{\text{out}} \to E^*_{\text{out}}.$$
4. Good ways to choose k:
$$k = 3;\ k = \left\lceil \sqrt{N} \right\rceil;\ \text{validation/cross validation.}$$
5. Easy to justify classification to customer.
6. Can easily do multi-class.
7. Can easily adapt to regression or logistic regression
$$g(\mathbf{x}) = \frac{1}{k} \sum_{i=1}^{k} y_{[i]}(\mathbf{x}) \qquad g(\mathbf{x}) = \frac{1}{k} \sum_{i=1}^{k} [\![y_{[i]}(\mathbf{x}) = +1]\!]$$
8. **Computationally demanding**.

Computational Demands of Nearest Neighbor

Memory.

Need to store all the data, $O(Nd)$ memory.

$N = 10^6$, $d = 100$, double precision$\approx$ 1GB

Finding the nearest neighbor of a test point.

Need to compute distance to every data point, $O(Nd)$.

$N = 10^6$, $d = 100$, 3GHz processor

$\approx$ 3ms (compute $g(\mathbf{x})$)

$\approx$ 1hr (compute CV error)

$>$ 1month (choose best k from among 1000 using CV)

Two Basic Approaches

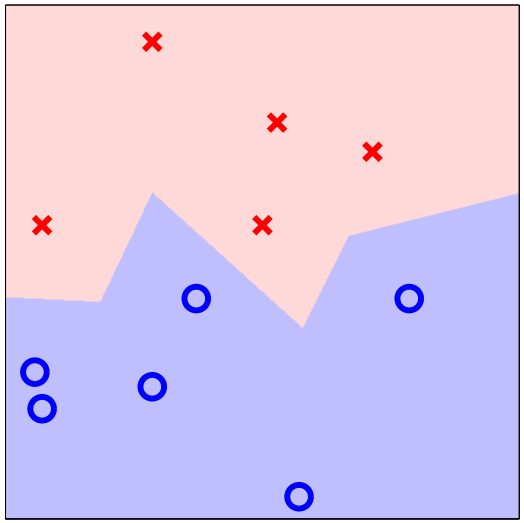
Reduce the amount of data.

The 5-year old does not remember every horse he has seen, only a few *representative* horses.

Store the data in a specialized data structure.

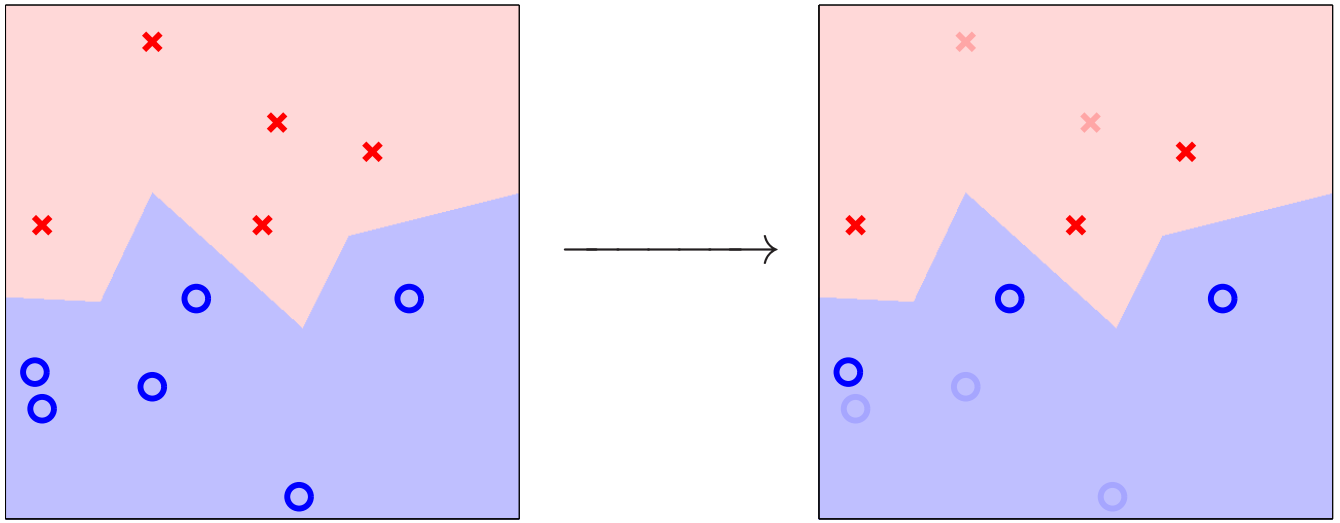
Ongoing research field to develop geometric data structures to make finding nearest neighbors fast.

Throw Away Irrelevant Data



$$k = 1$$

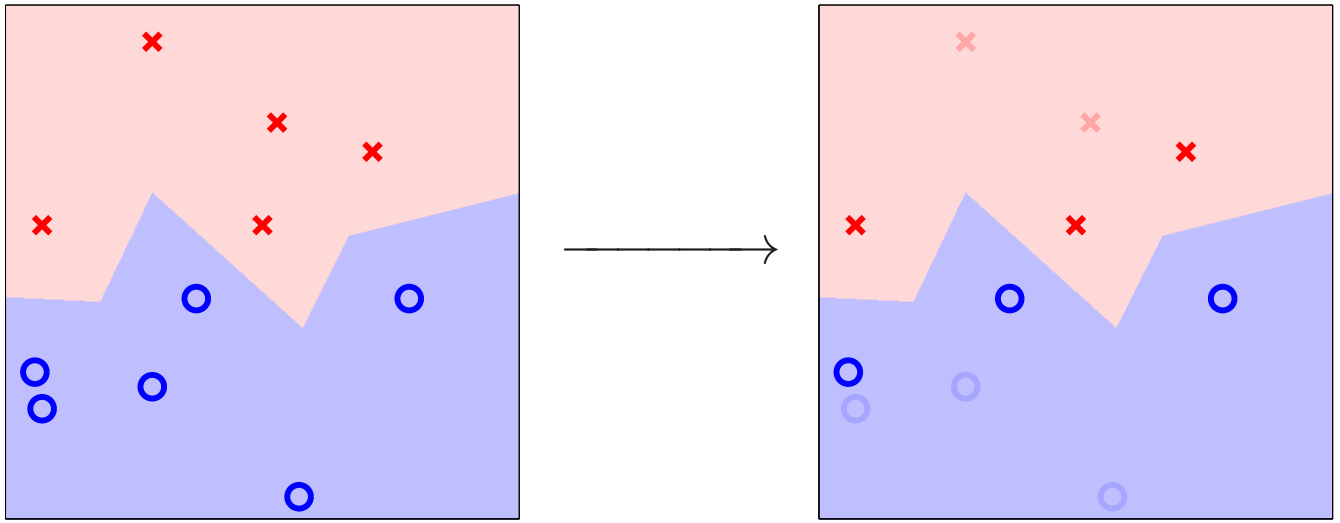
Decision Boundary Consistent



$g(\mathbf{x})$ unchanged

Does E_{in} change?
Does E_{out} change?

Decision Boundary Consistent

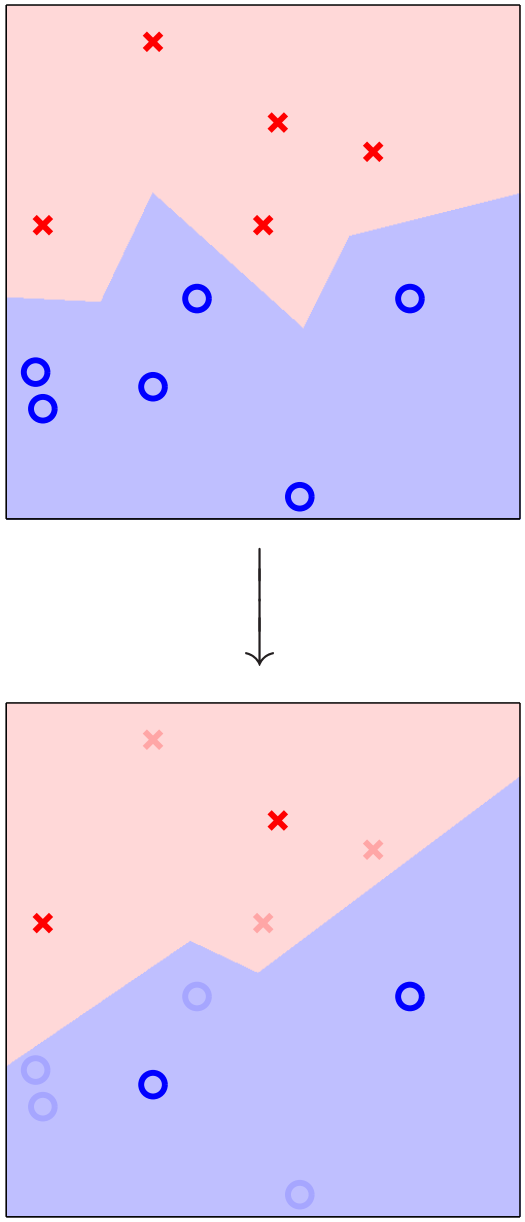


$g(\mathbf{x})$ unchanged

Does E_{in} change? No

Does E_{out} change? No

Training Set Consistent

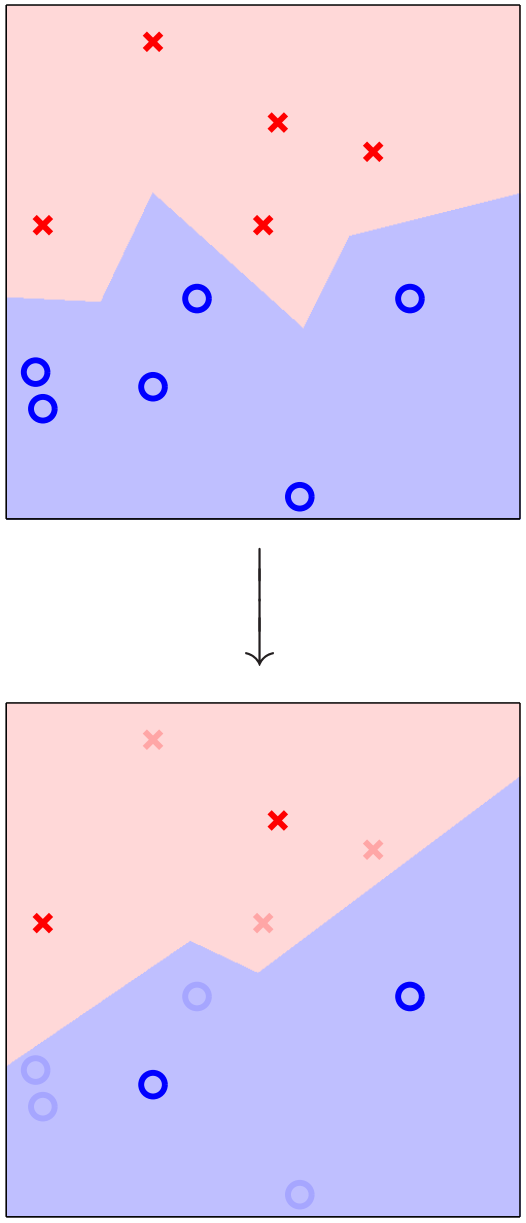


$g(\mathbf{x}_n)$ unchanged

Does E_{in} change?

Does E_{out} change?

Training Set Consistent

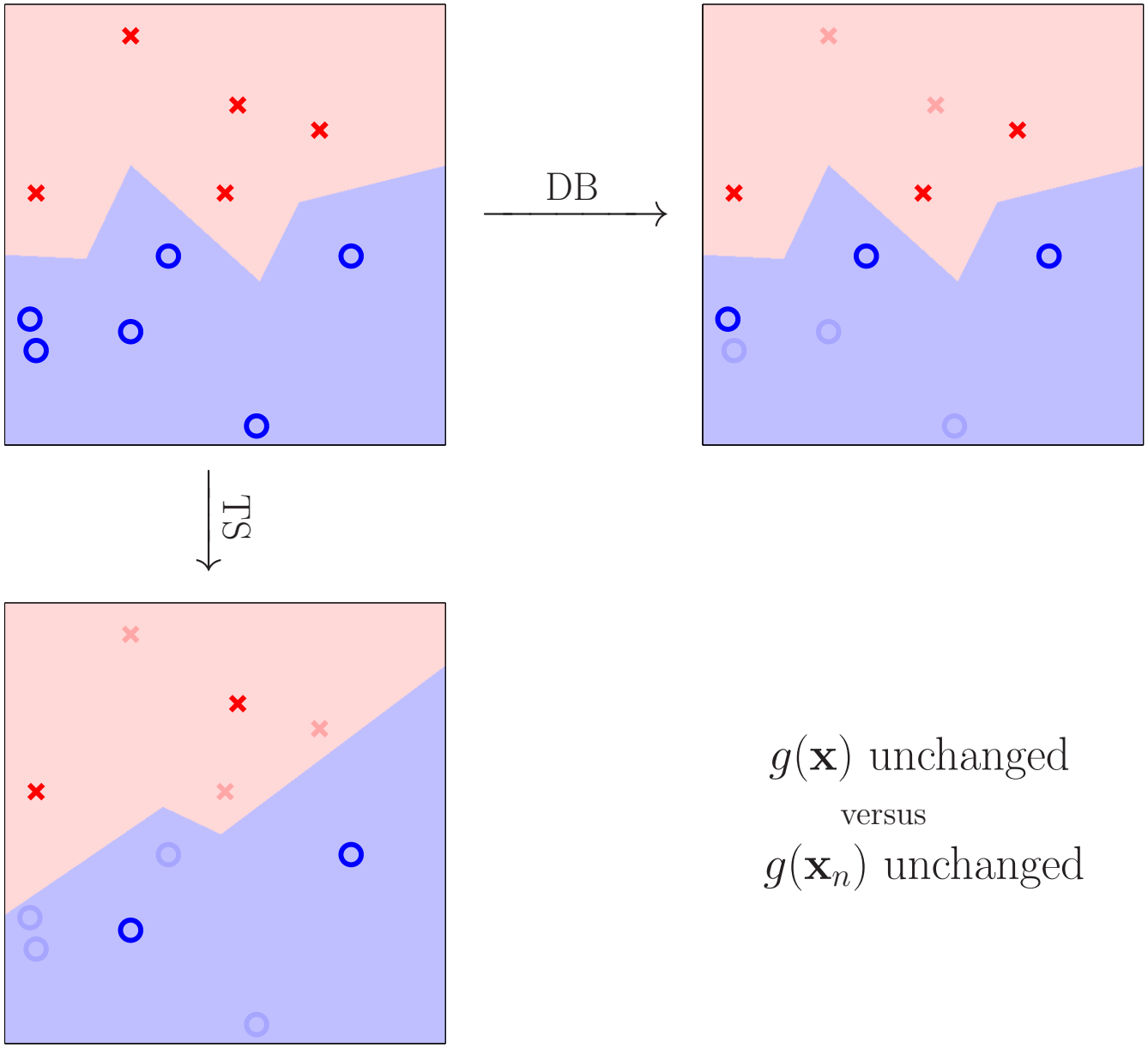


$g(\mathbf{x}_n)$ unchanged

Does E_{in} change? No

Does E_{out} change? Usually

Decision Boundary Vs. Training Set Consistent

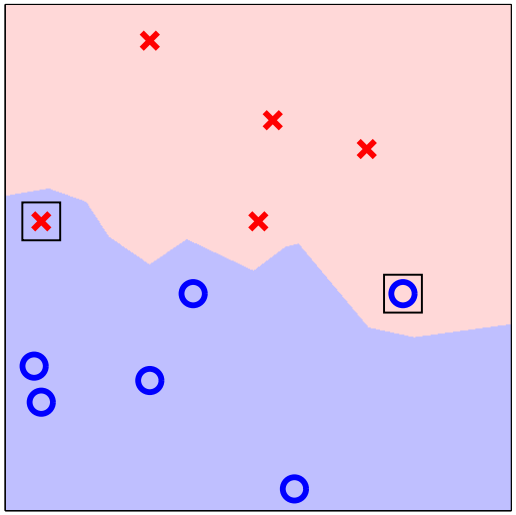


$g(\mathbf{x})$ unchanged

versus

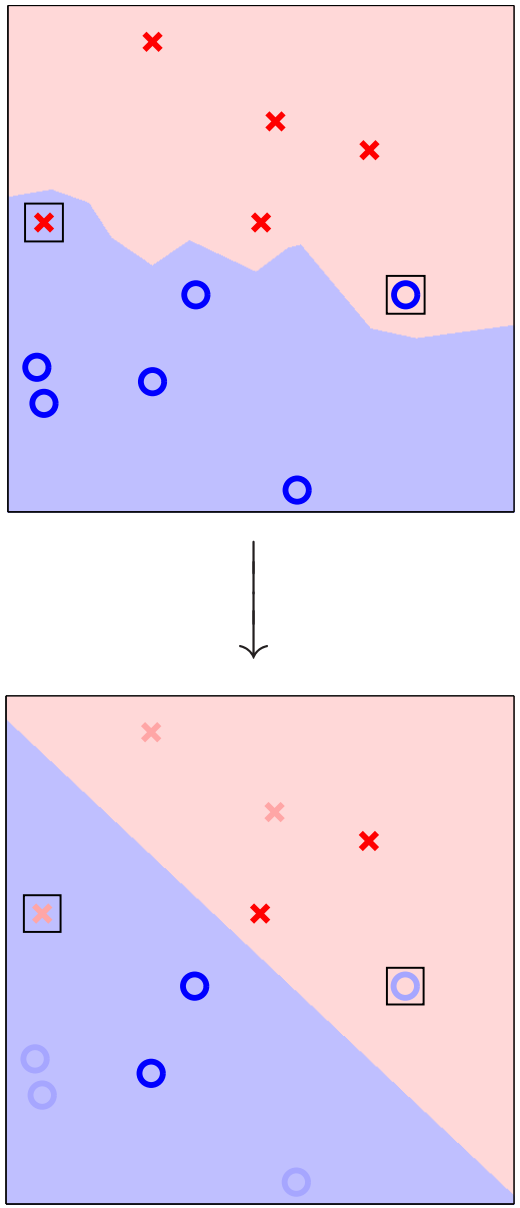
$g(\mathbf{x}_n)$ unchanged

Consistent Does Not Mean $g(\mathbf{x}_n) = y_n$



$$k = 3$$

Training Set Consistent ($k = 3$)

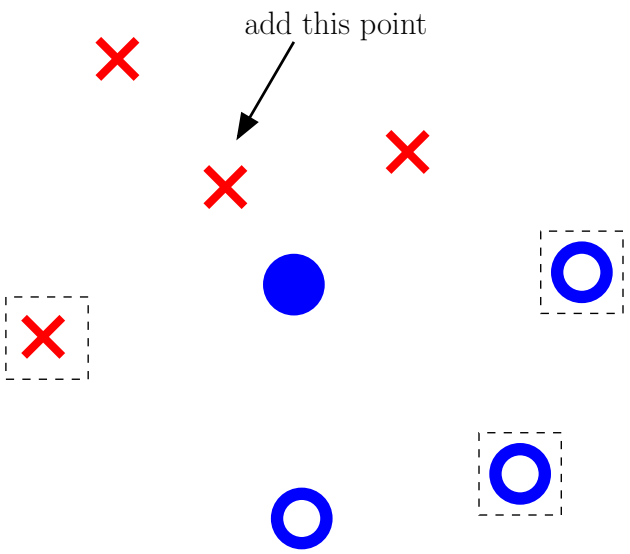


$g(\mathbf{x}_n)$ unchanged

Does E_{in} change? No

Does E_{out} change? Usually

CNN: Condensed Nearest Neighbor ($k = 3$)

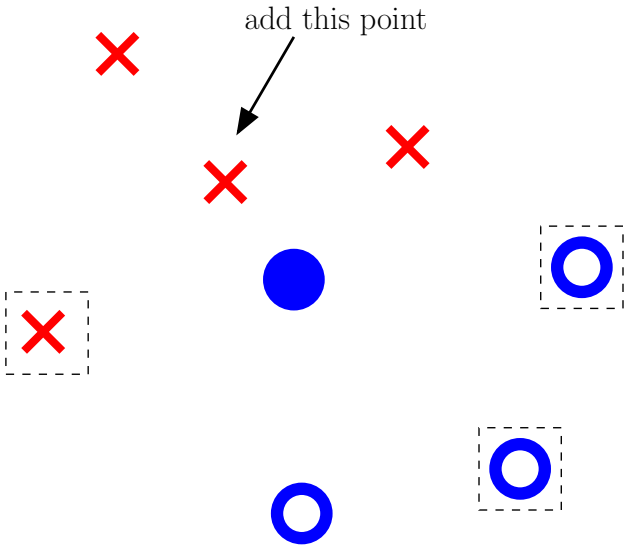


Consider the solid blue point:

i. blue w.r.t. selected points

ii. red w.r.t. $\mathcal{D}$

CNN: Condensed Nearest Neighbor



Consider the solid blue point:

i. blue w.r.t. selected points

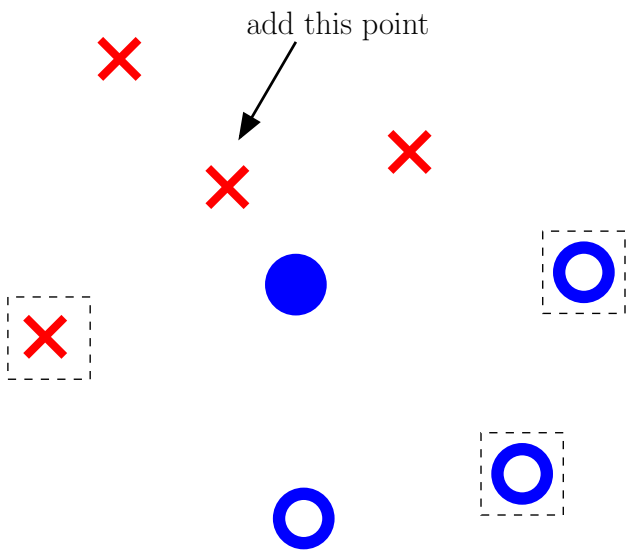
ii. red w.r.t. $\mathcal{D}$

Add a red point:

i. not already selected

ii. closest to the inconsistent point

CNN: Condensed Nearest Neighbor



1. Randomly select k data points into $\mathcal{S}$.
2. Classify all data according to $\mathcal{S}$.
3. Let $\mathbf{x}_*$ be an inconsistent point and y_* its class w.r.t. $\mathcal{D}$.
4. Add the closest point to $\mathbf{x}_*$ not in $\mathcal{S}$ that has class y_*.
5. Iterate until $\mathcal{S}$ classifies all points consistently with $\mathcal{D}$.

Consider the solid blue point:

i. blue w.r.t. selected points
ii. red w.r.t. $\mathcal{D}$

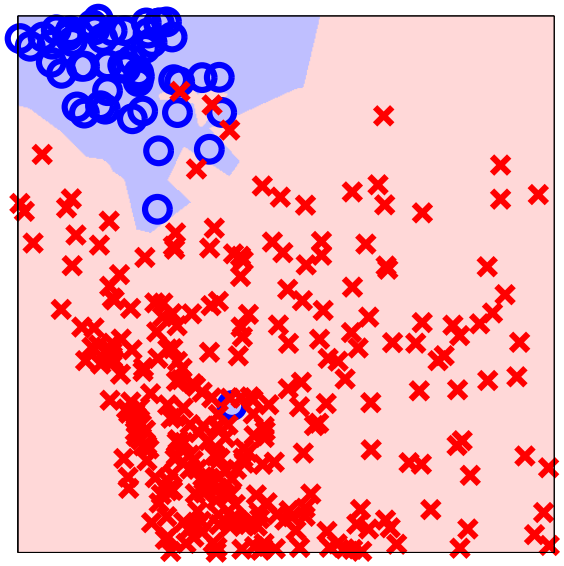
Add a red point:

i. not already selected
ii. closest to the inconsistent point

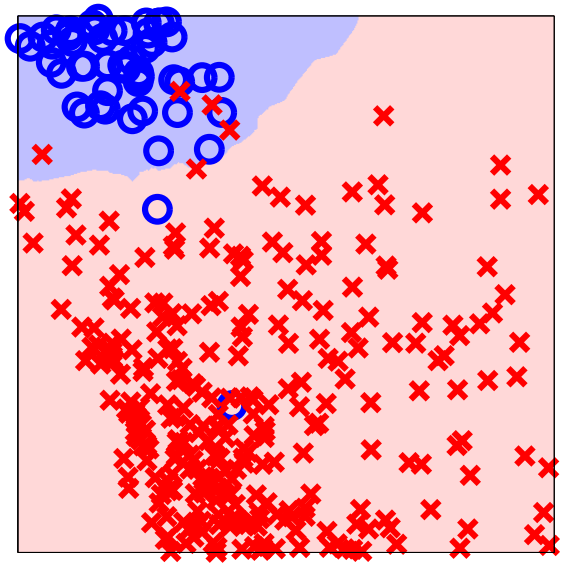
Minimum consistent set (MCS)? ← NP-hard

Nearest Neighbor on Digits Data

1-NN rule

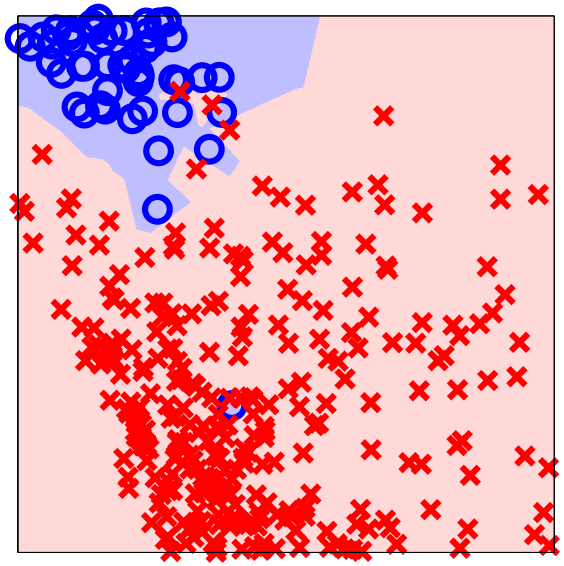


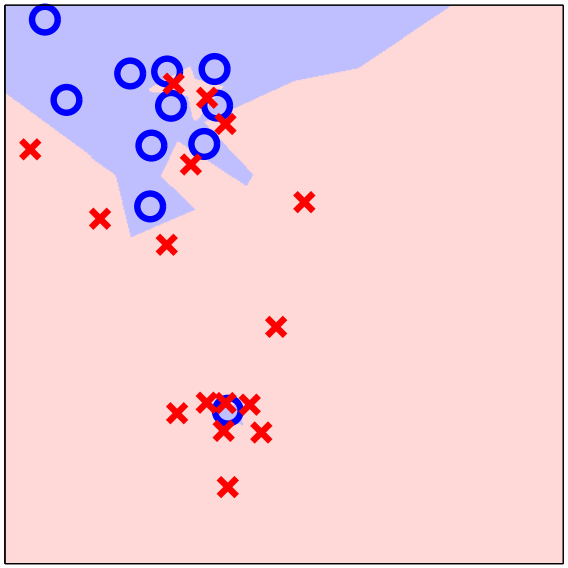
21-NN rule



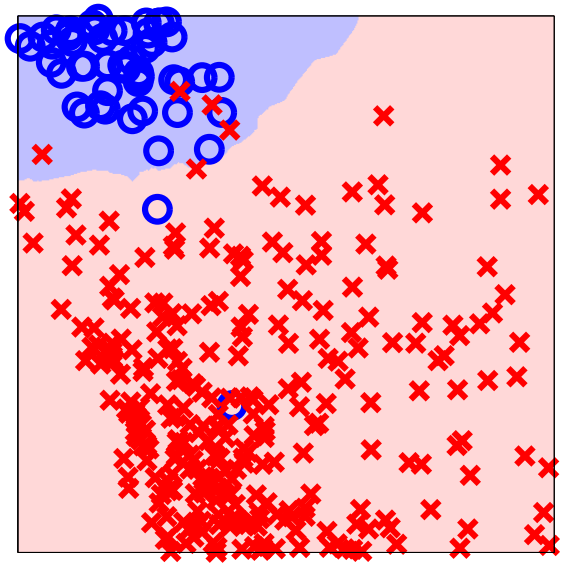
Condensing the Digits Data

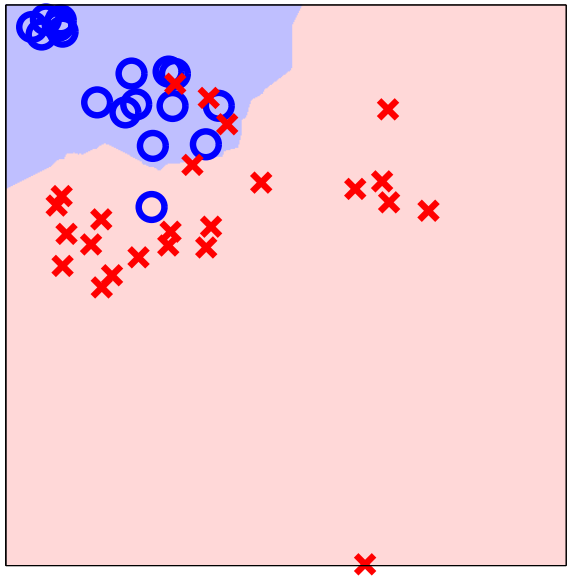
1-NN rule





21-NN rule





Finding the Nearest Neighbor

1. S_1, S_2 are 'clusters' with centers $\boldsymbol{\mu}_1, \boldsymbol{\mu}_2$ and radii r_1, r_2.

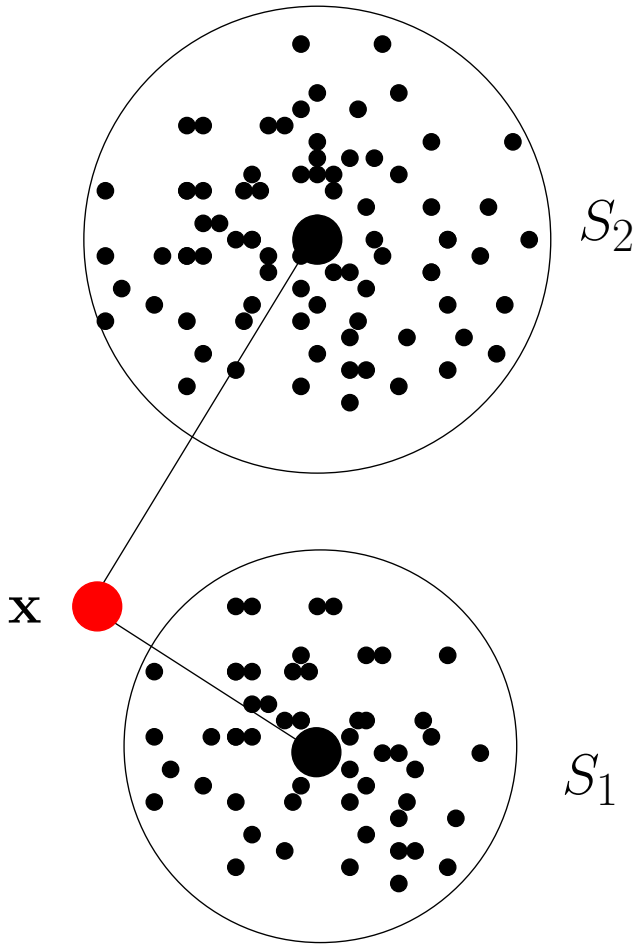
2. **[Branch]** Search S_1 first $\rightarrow \hat{\mathbf{x}}_{[1]}$.

3. The distance from $\mathbf{x}$ to any point in S_2 is at least

$$\| \mathbf{x} - \boldsymbol{\mu}_2 \| - r_2$$

4. **[Bound]** So we are done if

$$\| \mathbf{x} - \hat{\mathbf{x}}_{[1]} \| \leq \| \mathbf{x} - \boldsymbol{\mu}_2 \| - r_2$$



A *branch and bound* algorithm

Can be applied *recursively*

When Does the Bound Hold?

Bound condition: $\| \mathbf{x} - \hat{\mathbf{x}}_{[1]} \| \leq \| \mathbf{x} - \boldsymbol{\mu}_2 \| - r_2$.

$$\| \mathbf{x} - \hat{\mathbf{x}}_{[1]} \| \leq \| \mathbf{x} - \boldsymbol{\mu}_1 \| + r_1$$

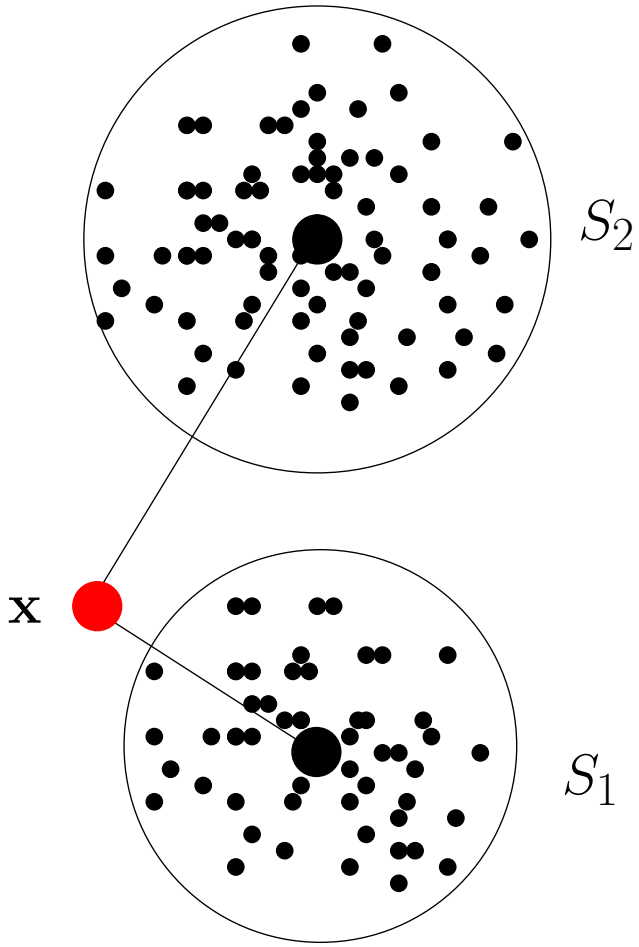
So, it suffices that

$$r_1 + r_2 \leq \| \mathbf{x} - \boldsymbol{\mu}_2 \| - \| \mathbf{x} - \boldsymbol{\mu}_1 \|.$$

$\| \mathbf{x} - \boldsymbol{\mu}_1 \| \approx 0$ means $\| \mathbf{x} - \boldsymbol{\mu}_2 \| \approx \| \boldsymbol{\mu}_2 - \boldsymbol{\mu}_2 \|$.

It suffices that

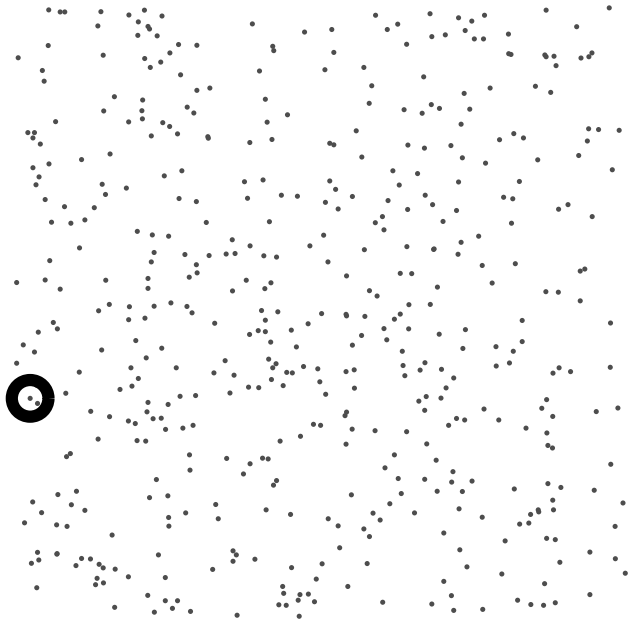
$$r_1 + r_2 \leq \| \boldsymbol{\mu}_2 - \boldsymbol{\mu}_1 \|.$$



within cluster spread should be less than *between cluster spread*

Finding Clusters – Lloyd's Algorithm

1. Pick well separated centers for each cluster.



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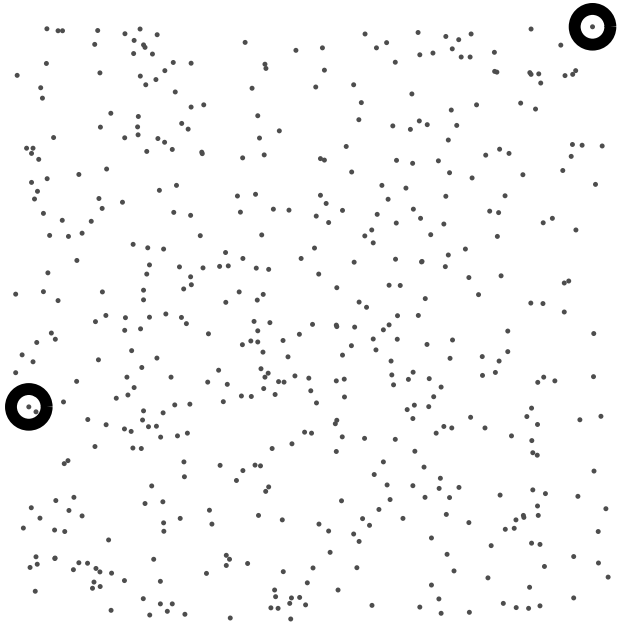

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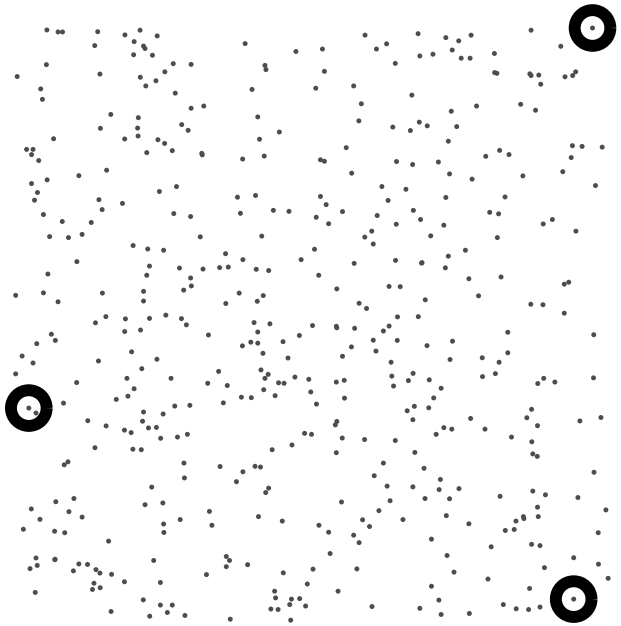

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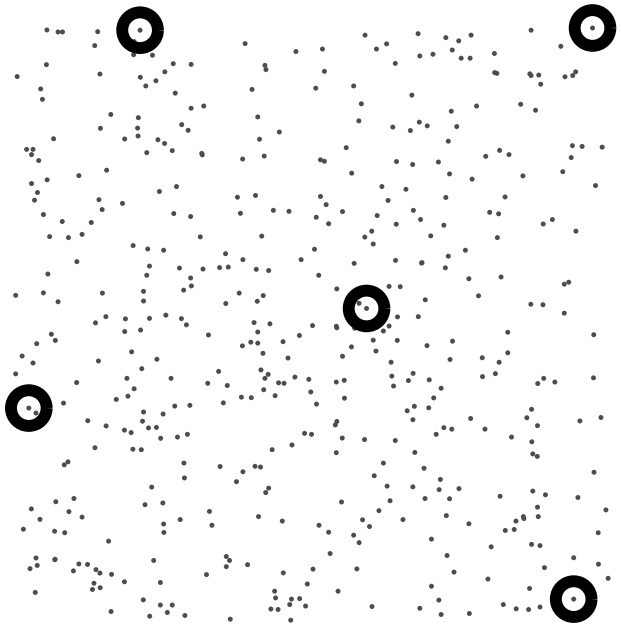
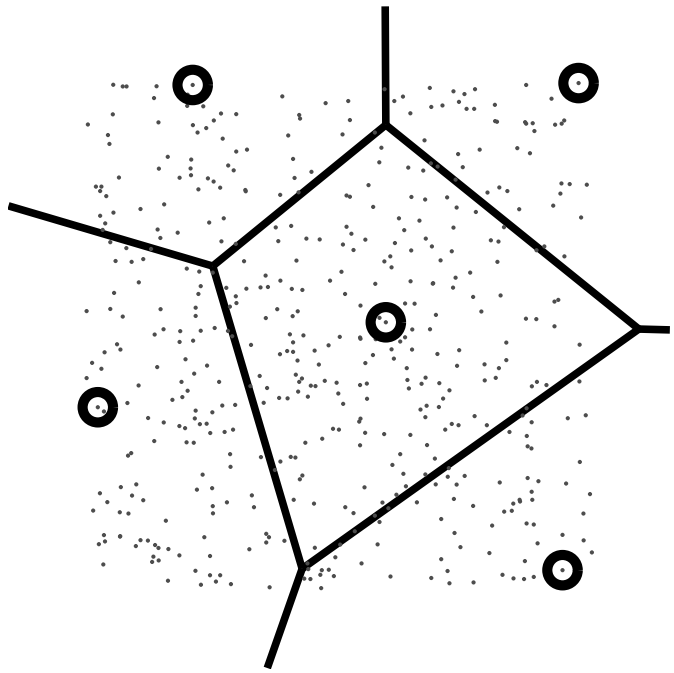


Finding Clusters – Lloyd's Algorithm

1. Pick well separated centers for each cluster.

2. Compute Voronoi regions as the clusters.

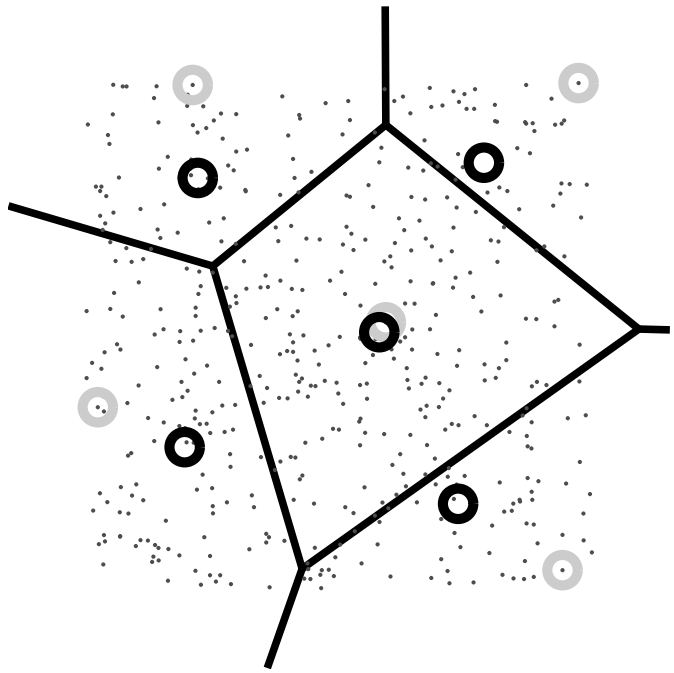


Finding Clusters – Lloyd's Algorithm

1. Pick well separated centers for each cluster.

2. Compute Voronoi regions as the clusters.

3. Update the Centers.



Finding Clusters – Lloyd's Algorithm

1. Pick well separated centers for each cluster.

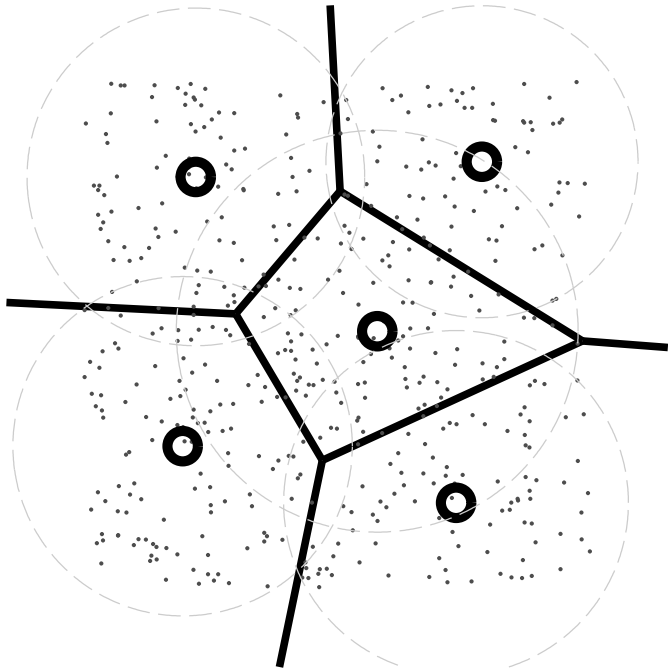
2. Compute Voronoi regions as the clusters.

3. Update the Centers.

4. Update the Voronoi regions.

5. Compute centers and radii:

$$\boldsymbol{\mu}_j = \frac{1}{|S_j|} \sum_{\mathbf{x}_n \in S_j} \mathbf{x}_n; \qquad r_j = \max_{\mathbf{x}_n \in S_j} \| \mathbf{x}_n - \boldsymbol{\mu}_j \|.$$



Radial Basis Functions (RBF)

k-Nearest Neighbor: Only considers k-nearest neighbors.

each neighbor has equal weight

What about using *all* data to compute $g(\mathbf{x})$?

RBF: Use all data.

data further away from $\mathbf{x}$ have less weight.