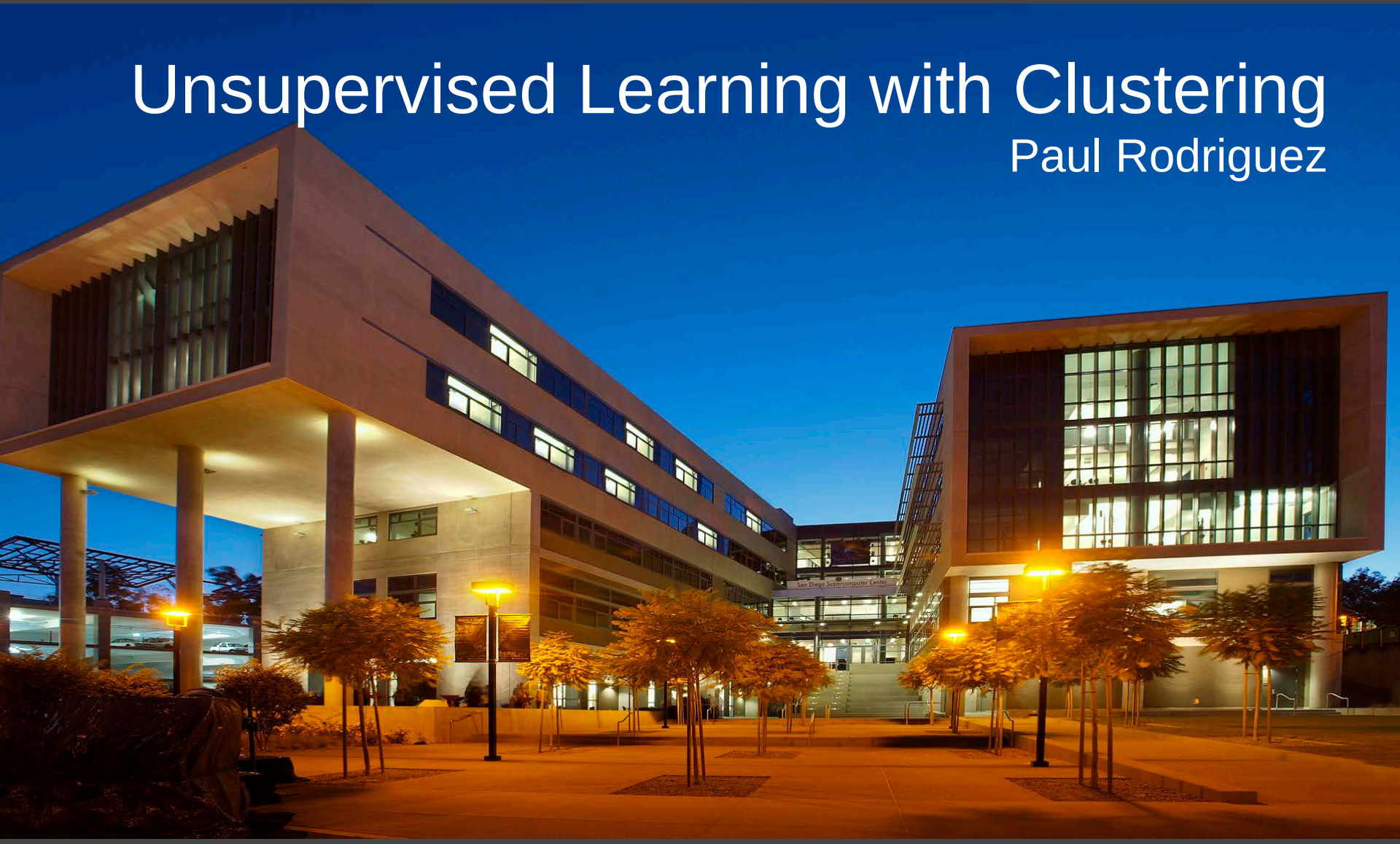
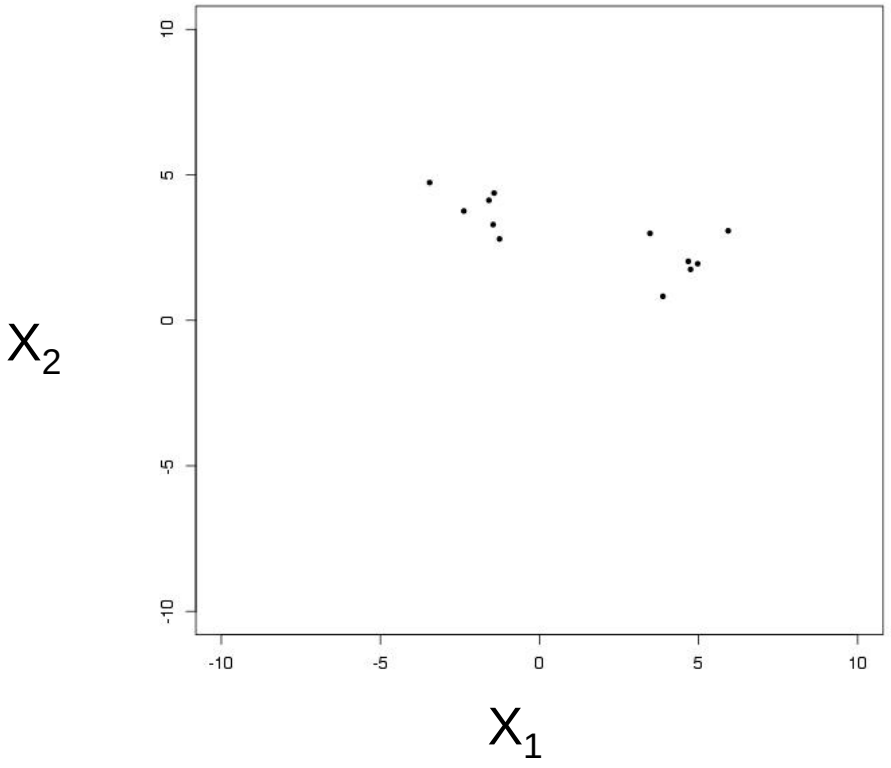
Unsupervised Learning with Clustering

Paul Rodriguez



Clustering Idea

- **Given a set of data can we find a natural grouping?**



Essential R commands:

```
D     =rnorm(12,0,1) #generate 12
                     #random normal
X1    =matrix(D,6,2) #put into 6x2 matrix
X1[,1]=X1[,1]+4;     #shift center
X1[,2]=X1[,2]+2;

#repeat for another set of points

#bind data points and plot
plot(rbind(X1,X2),
          xlim=c(-10,10),ylim=c(-10,10));
```

Why Clustering

- A good grouping implies some structure

Why Clustering

- **A good grouping implies some structure**
- **In other words, given a good grouping, we can then:**
  - Interpret and label clusters
  - Identify important features
  - Characterize new points by the closest cluster
  - Use the cluster assignments as a summary of the data

Clustering Objective

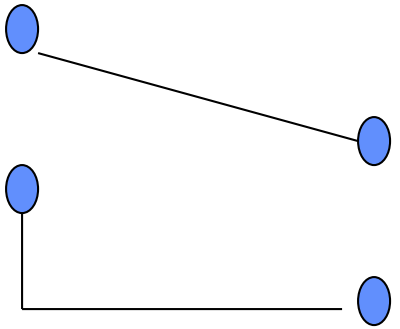
- **Objective: find subsets that are similar within cluster and dissimilar between clusters**

Clustering Objective

- **Objective: find subsets that are similar within cluster and dissimilar between clusters**
- **Similarity defined by distance measures**

Clustering Objective

- **Objective: find subsets that are similar within cluster and dissimilar between clusters**
- **Similarity defined by distance measures**
  - Euclidean distance
  - Manhattan distance



Clustering Objective

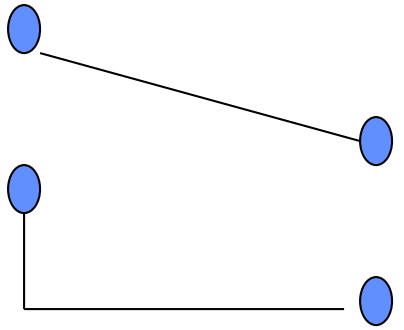
- **Objective: find subsets that are similar within cluster and dissimilar between clusters**
- **Similarity defined by distance measures**

  - Euclidean distance =

    $\text{sqrt}[(a1 - b1)^2 + (a2 - b2)^2 + \ldots)]$

  - Manhattan distance

    $[|a1 - b1| + |a2 - b2| + \ldots)]$

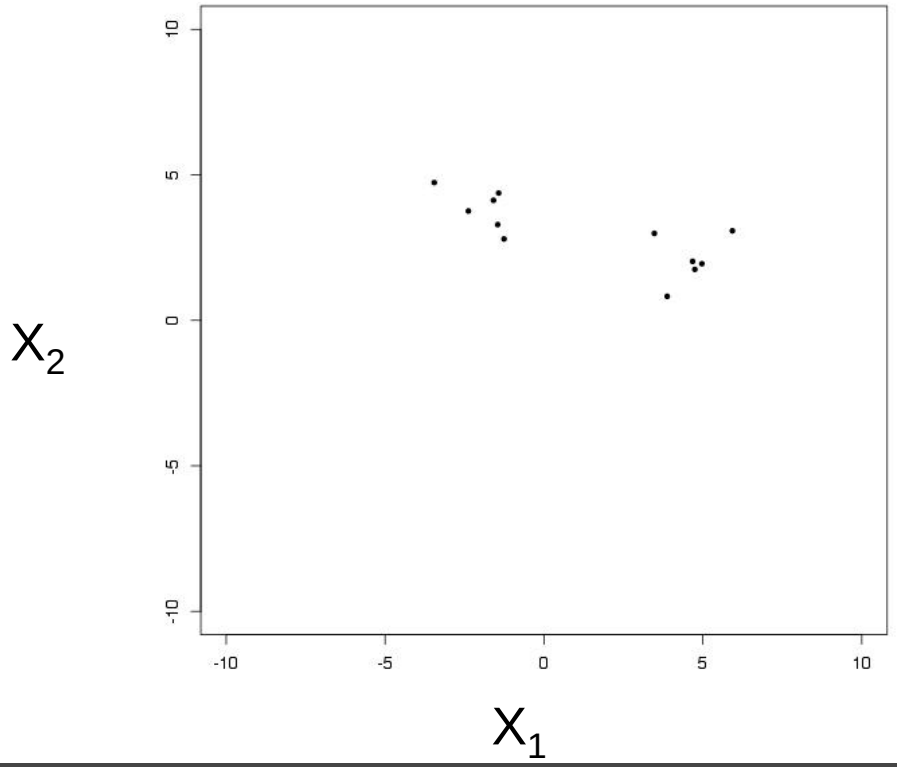


Kmeans Clustering

- **A simple, effective, and standard method**
  - Start with K initial cluster centers
  - Loop:
    - Assign each data point to nearest cluster center
    - Calculate mean of cluster for new center
  - Stop when assignments don't change
- **Issues:**
  - How to choose K?
  - How to choose initial centers?
  - Will it always stop?

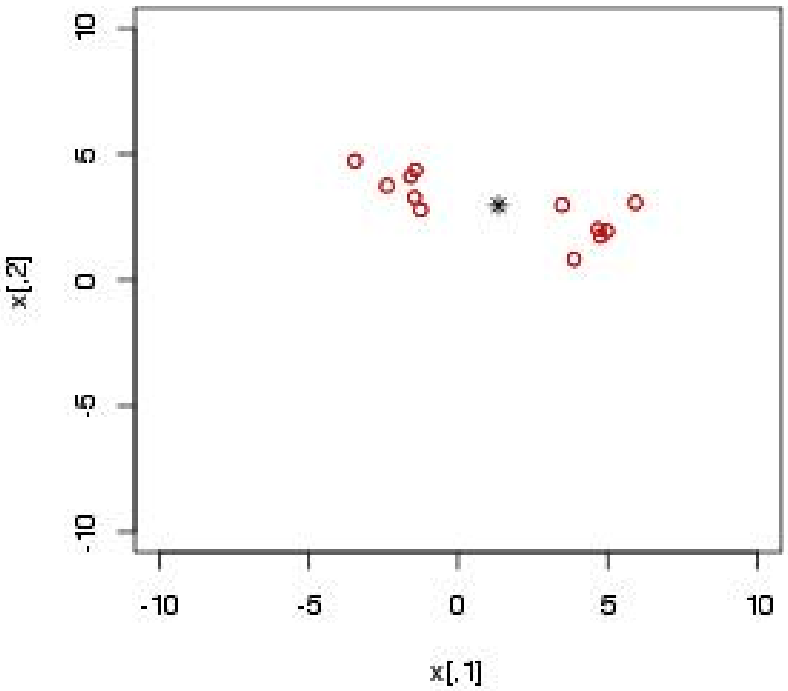
Kmeans Example

- **For K=1, using Euclidean distance, where will the cluster center be?**



Kmeans Example

For K=1, the overall mean minimizes Sum Squared Error (SSE), aka Euclidean dist. squared

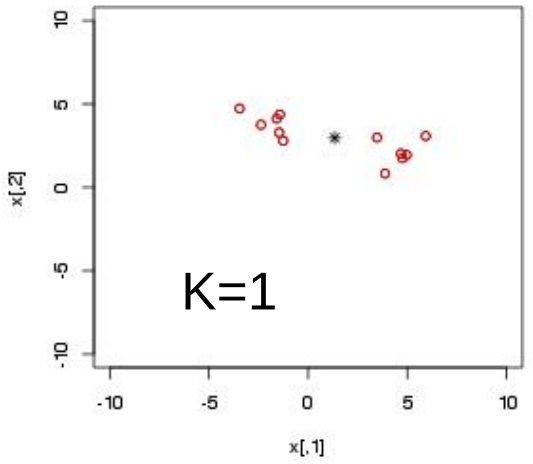


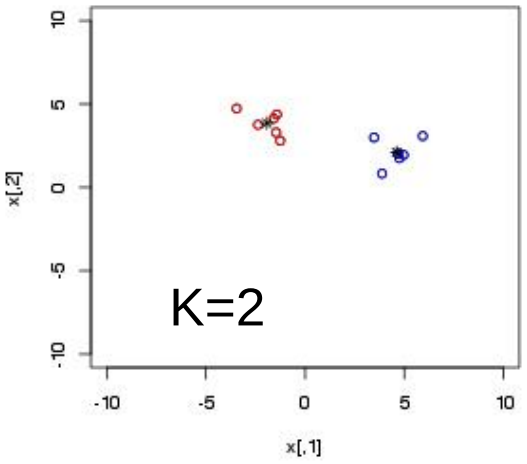
Essential R commands:
Kresult = kmeans(X,1,10,1)

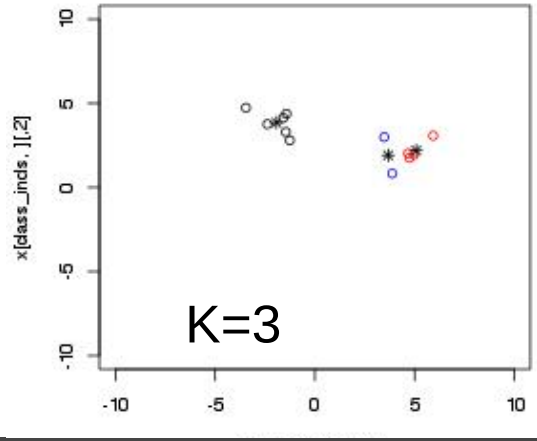
#choose 1 data point as initial K centers
#10 is max loop iterations
#1 is number of initial sets to try

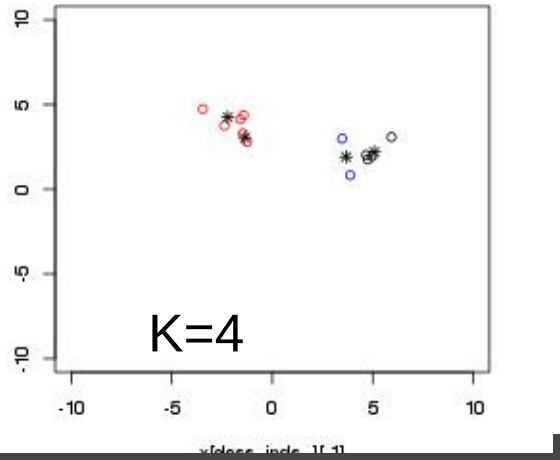
#Kresult is an R object with subfields
Kresult$cluster #cluster assignments
Kresults$tot.withinss # tot within SSE

Kmeans Example







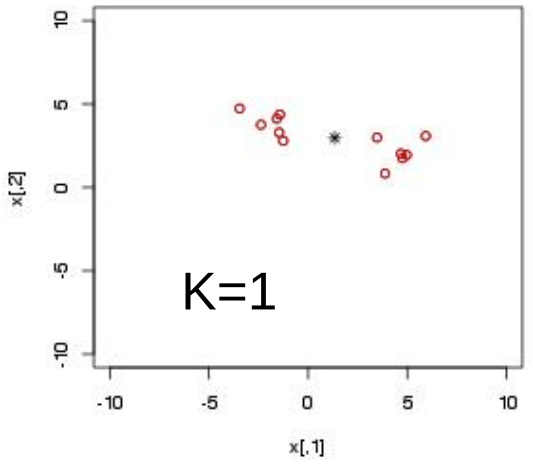


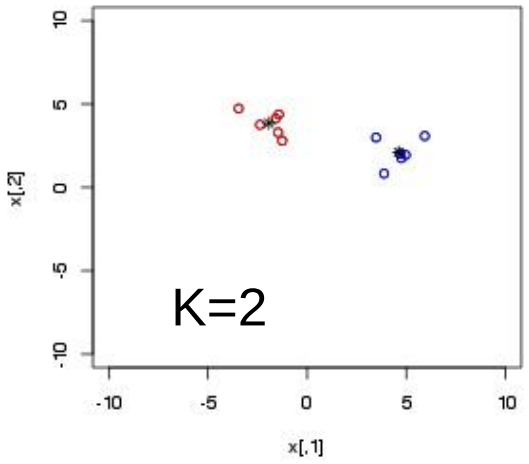
Essential R commands:
inds=which(Kresult$cluster==K)

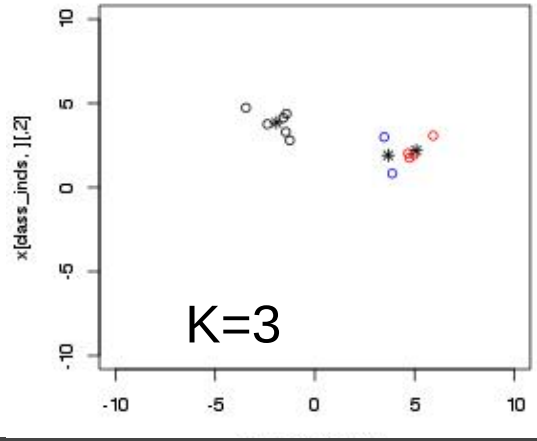
plot(X[inds,],col2use="red");

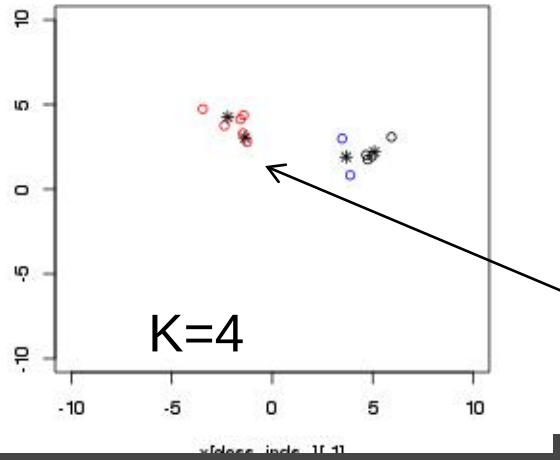
…

Kmeans Example









Essential R commands:
inds=which(Kresult$cluster==K)

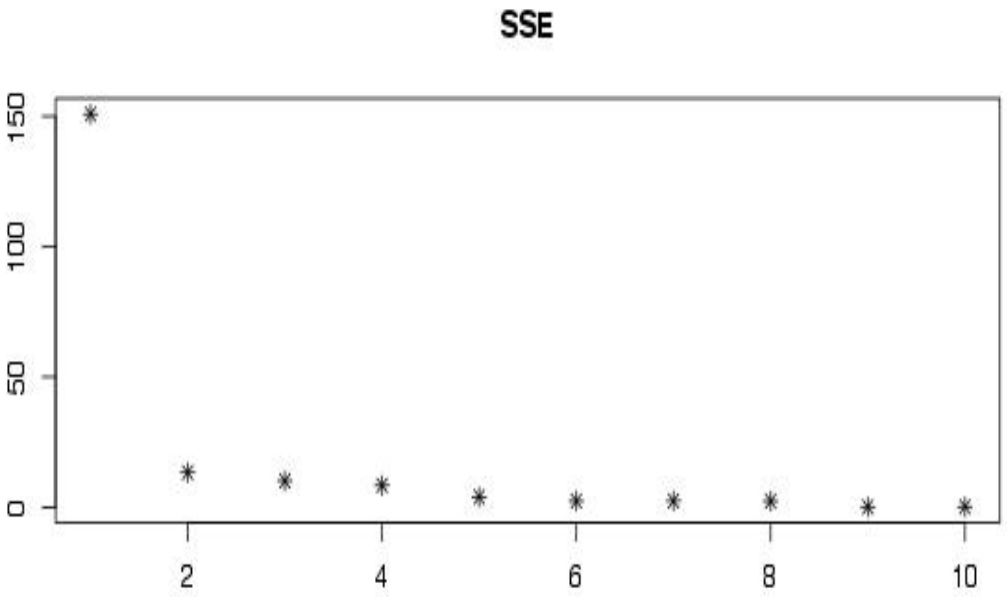
plot(X[inds,],col2use="red");

…

As K increases individual points get a cluster

Choosing K for Kmeans

Total Within Cluster SSE



K=1 to 10

Essential R commands:

```
for (num_k in 1:10) {
Kres=kmeans(X,num_k,10,1);
```

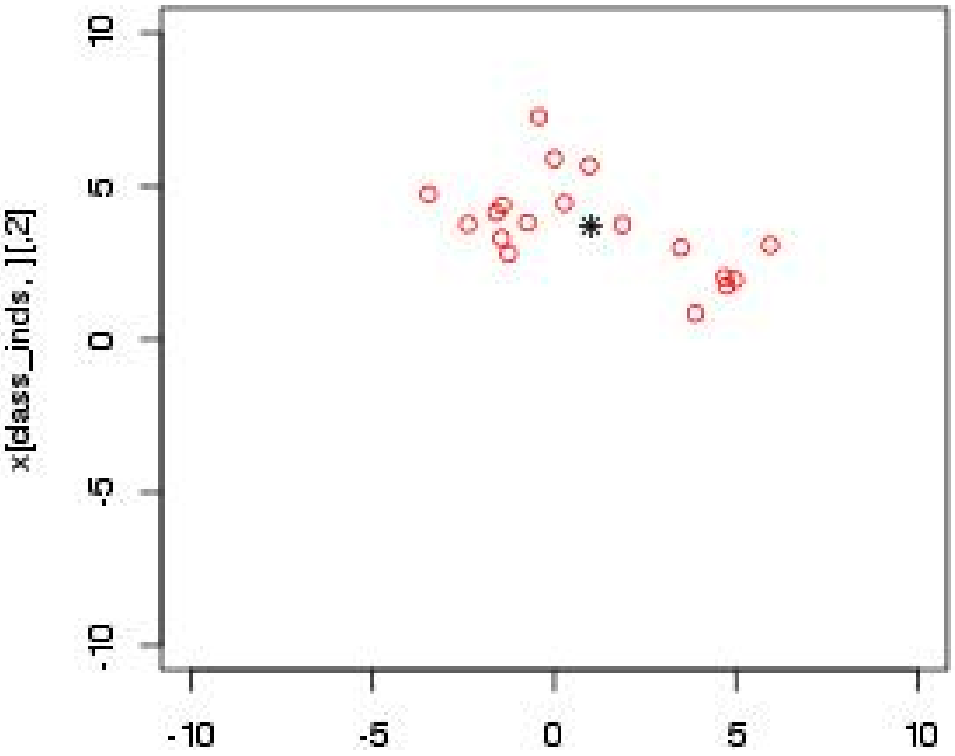
Save and then plot
Kres$tot.withinss

…

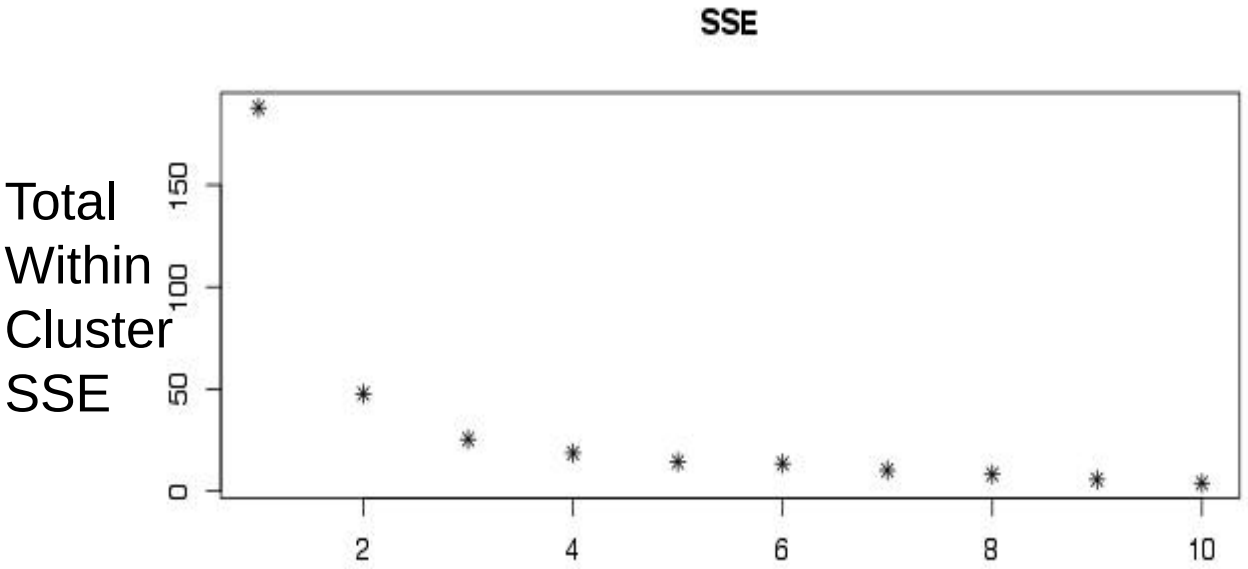
- Not much improvement after K=2 (“elbow”)

Kmeans Example – more points

How many clusters should there be?



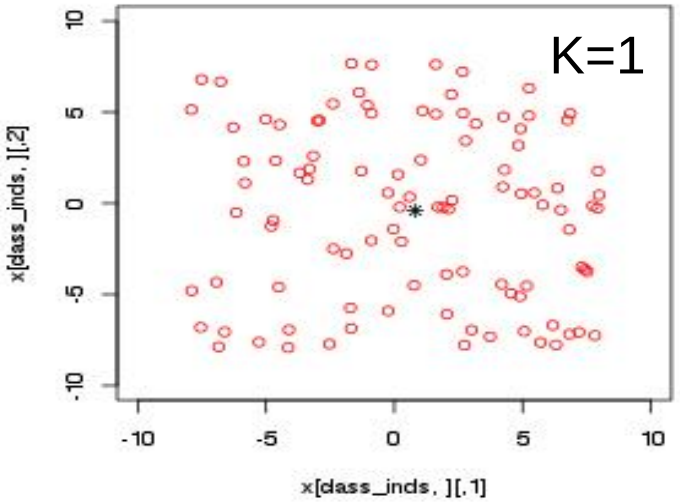
Choosing K for Kmeans

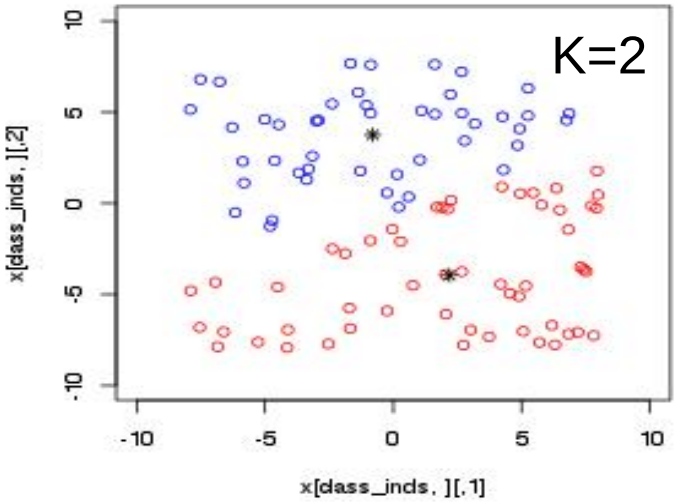


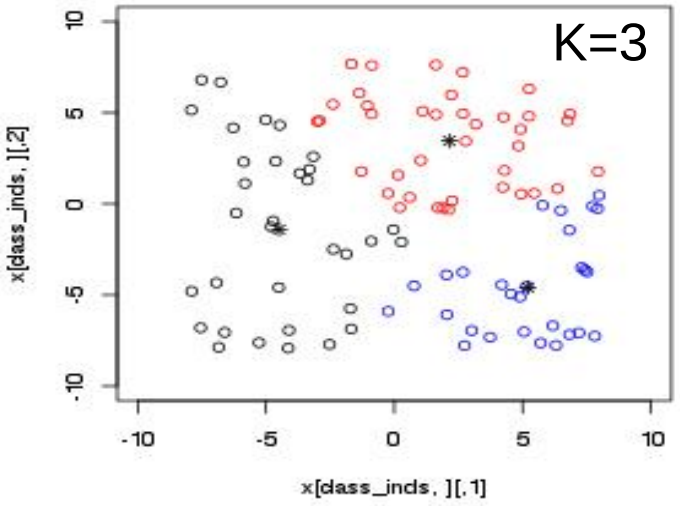
K=1 to 10

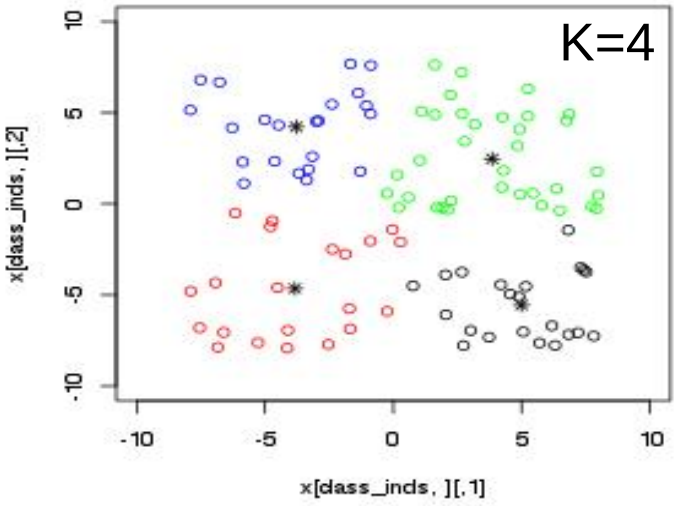
- Smooth decrease at K ≥ 2, harder to choose
- In general, smoother decrease => less structure

Kmeans Example: uniform dist.



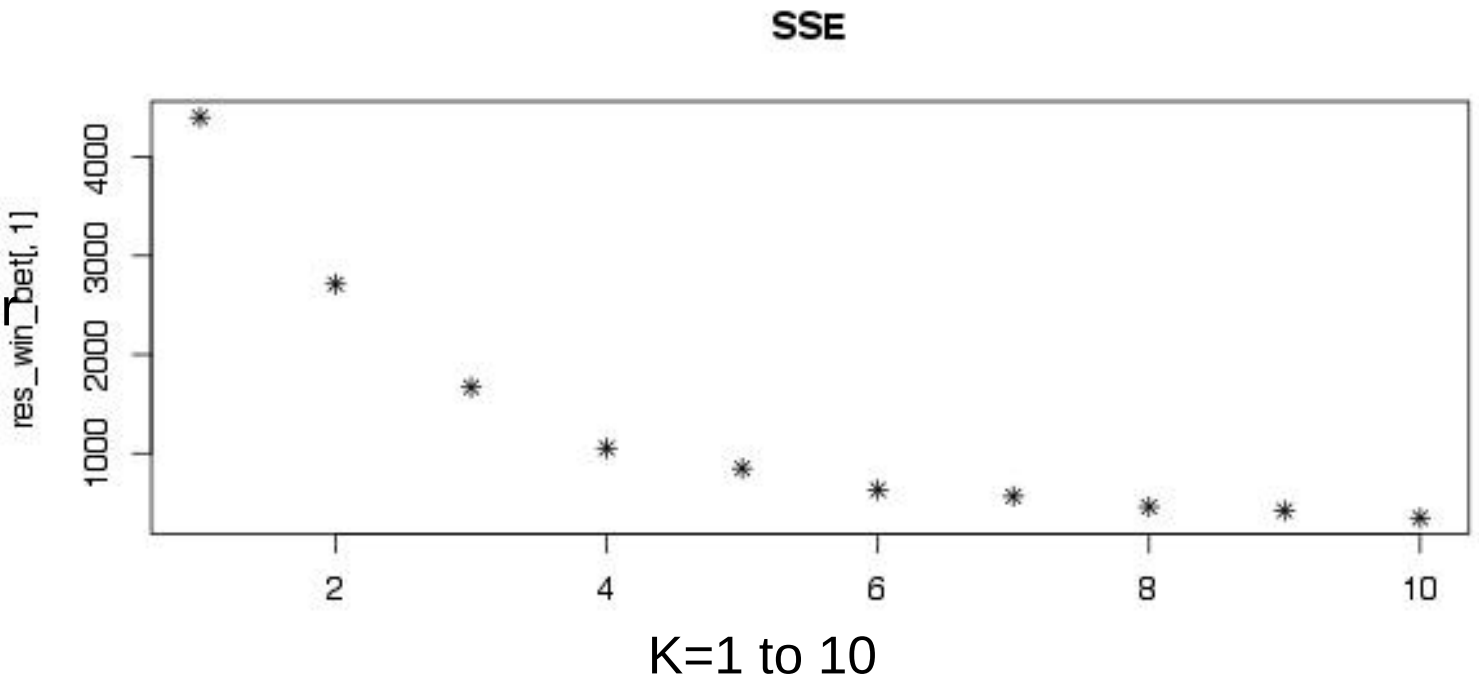






Choosing K - uniform

Total Within Cluster SSE



K=1 to 10

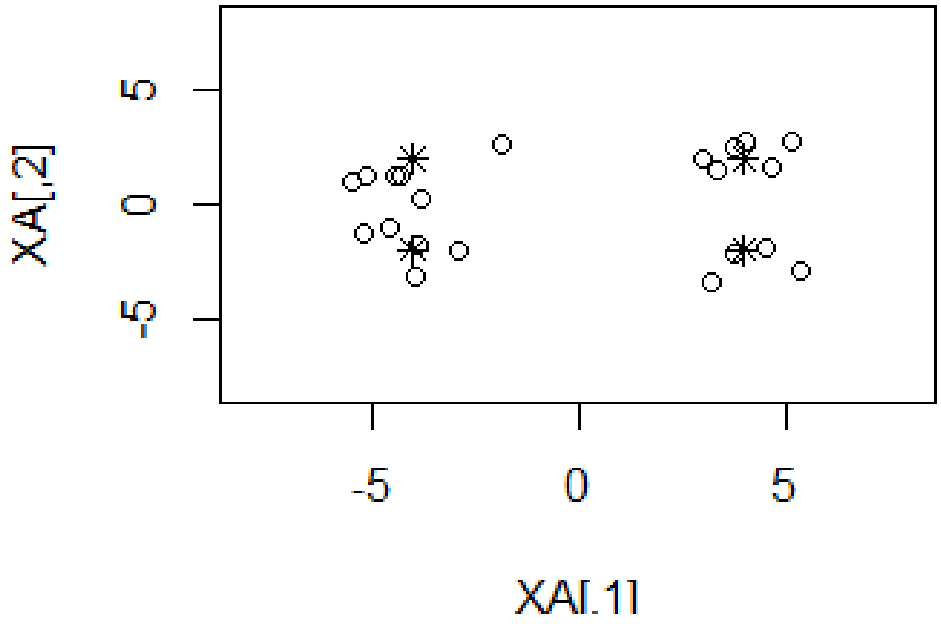
- Smooth decrease across K => less structure

Choosing K methods

- **Sum Squared Error Within Clusters**
  - Take the “elbow” in the decrease as K=1…N
- **“Silhouette”: mean distances within a cluster vs distances to next best cluster**
  - Take the maximum value overall mean as K=1…N
- **“Gap” value of SSE-within-cluster of data vs uniform distribution**
  - Take maximum value as K=1…N

Choosing K methods

- **Example, 4 clusters normal distribution, small sample**

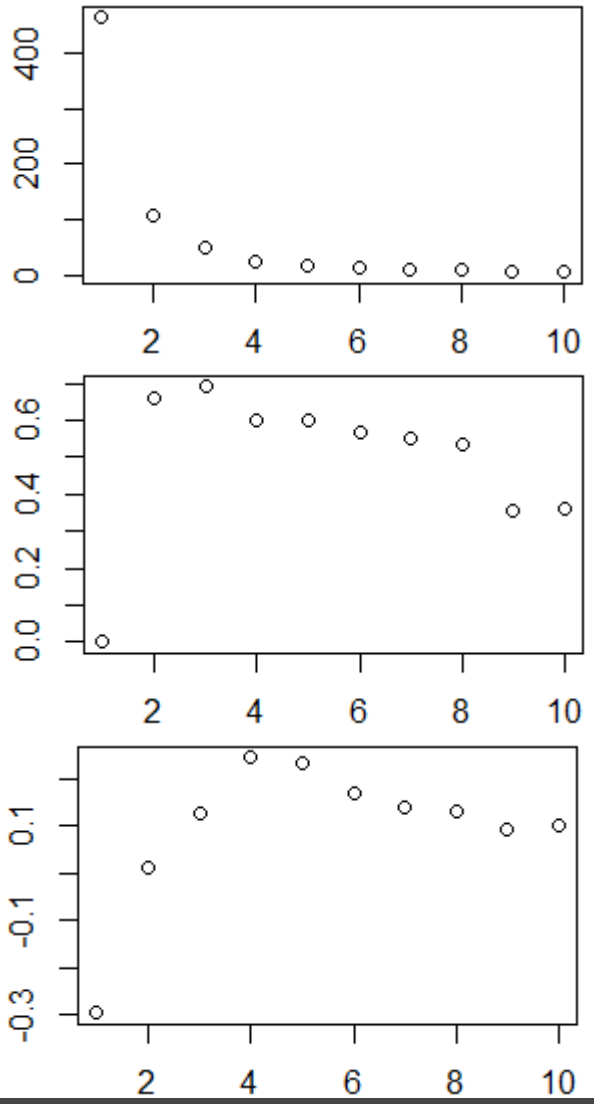


Choosing K methods

SSE within cluster (elbow)

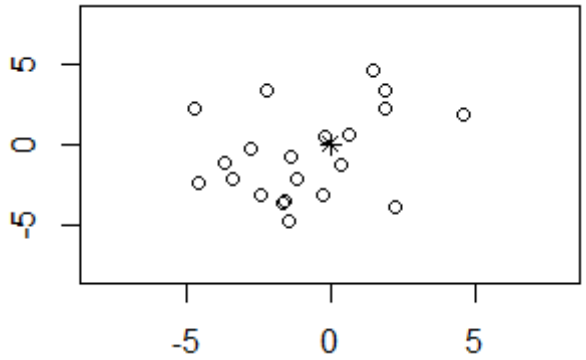
Silhouette intra vs next best

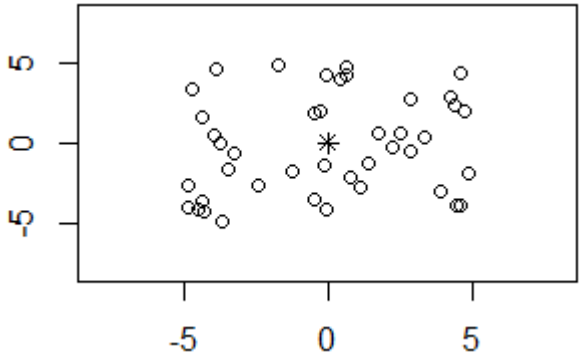
Gap SSE within cluster vs uniform data baseline

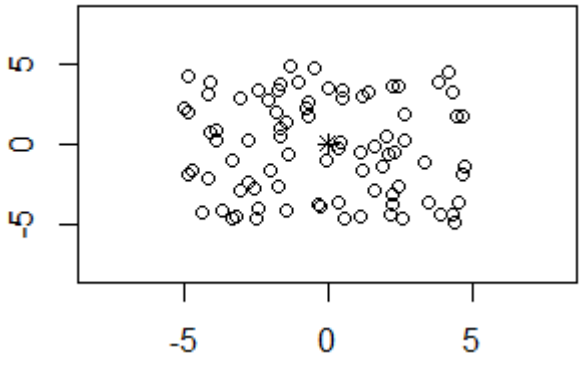


Choosing K methods

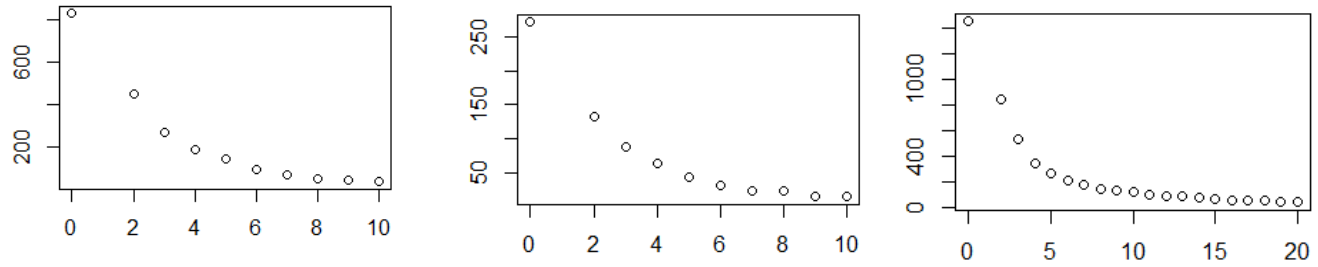
- **Example, random uniform data, small, med, larger samples**



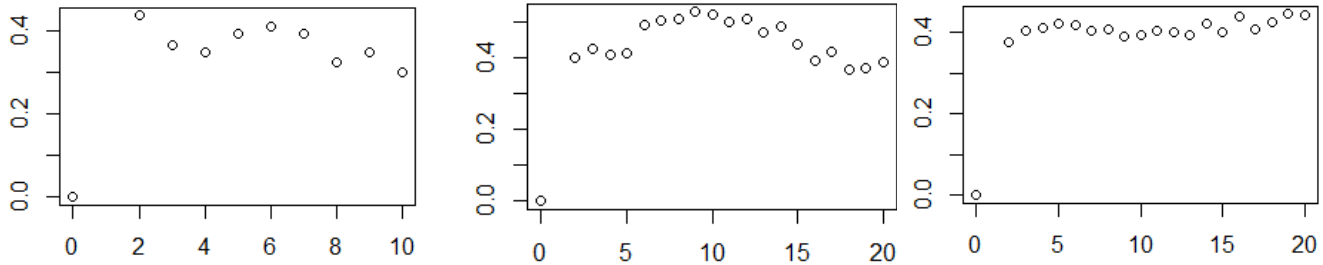




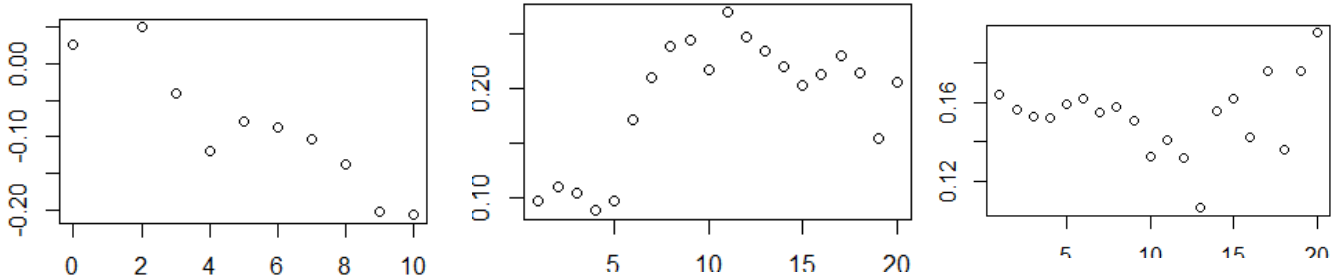
SSE within cluster (elbow)



Silhouette intra vs next best



Gap SSE within cluster vs uniform data baseline



small to large data sample size →

Kmeans Guidelines

- **Initial starting points and convergence:**

  Kmeans++ algorithm selects good initial clusters that also helps convergence

  10 iterations often good

  Can also run several times and choose best result

Kmeans Clustering Issues

- **Scale:**
  - Dimensions with large numbers may dominate distance metrics (so can be good to normalize or scale data)

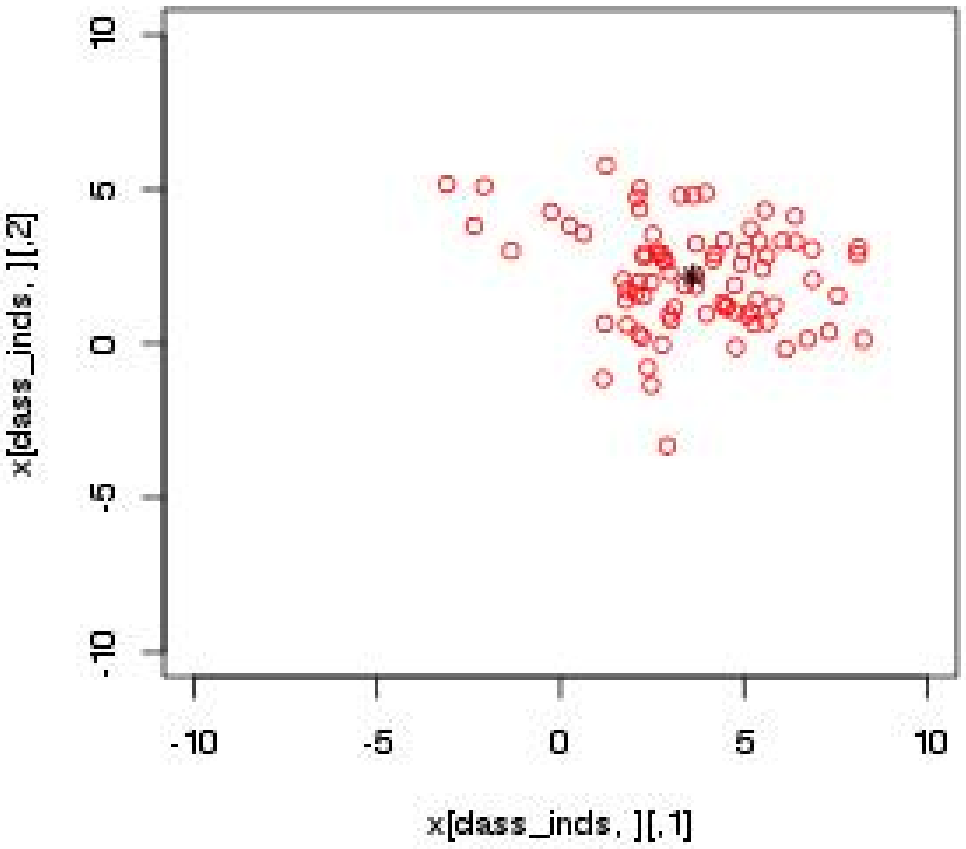
- **Outliers:**
  - Outliers can pull cluster mean

    (K-mediods uses median instead of mean)

Soft Clustering Methods

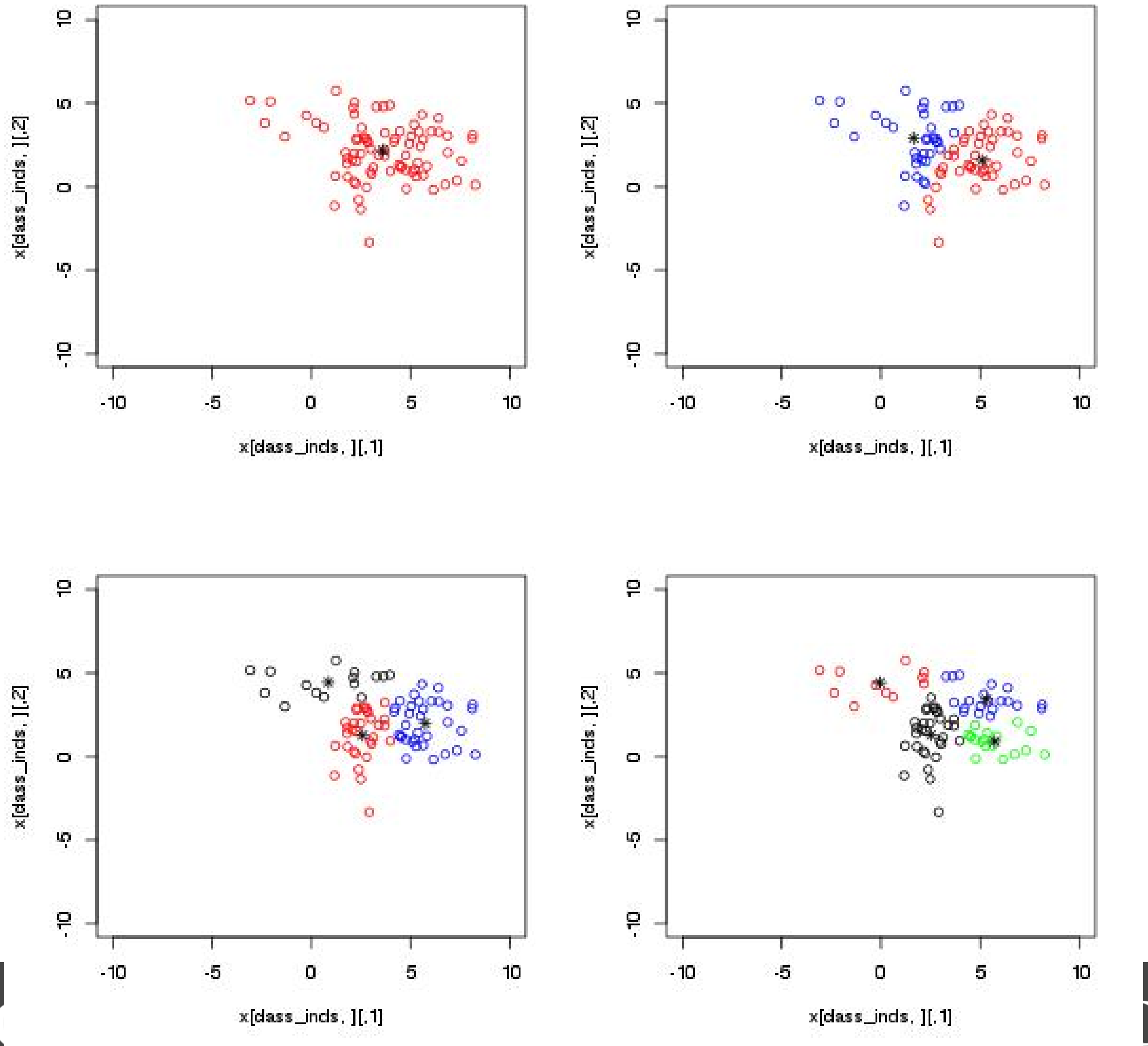
- **Fuzzy Clustering**
  - Kmeans with weighted assignments to all clusters
  - Weights depend on relative distance
  - Find min weighted SSE
- **Expectation-Maximization:**
  - Initialize a mixture of multivariate Gaussian distributions
  - Find means, variances, and mixture weights that maximize probability of data

Kmeans with unequal cluster variance and/or size

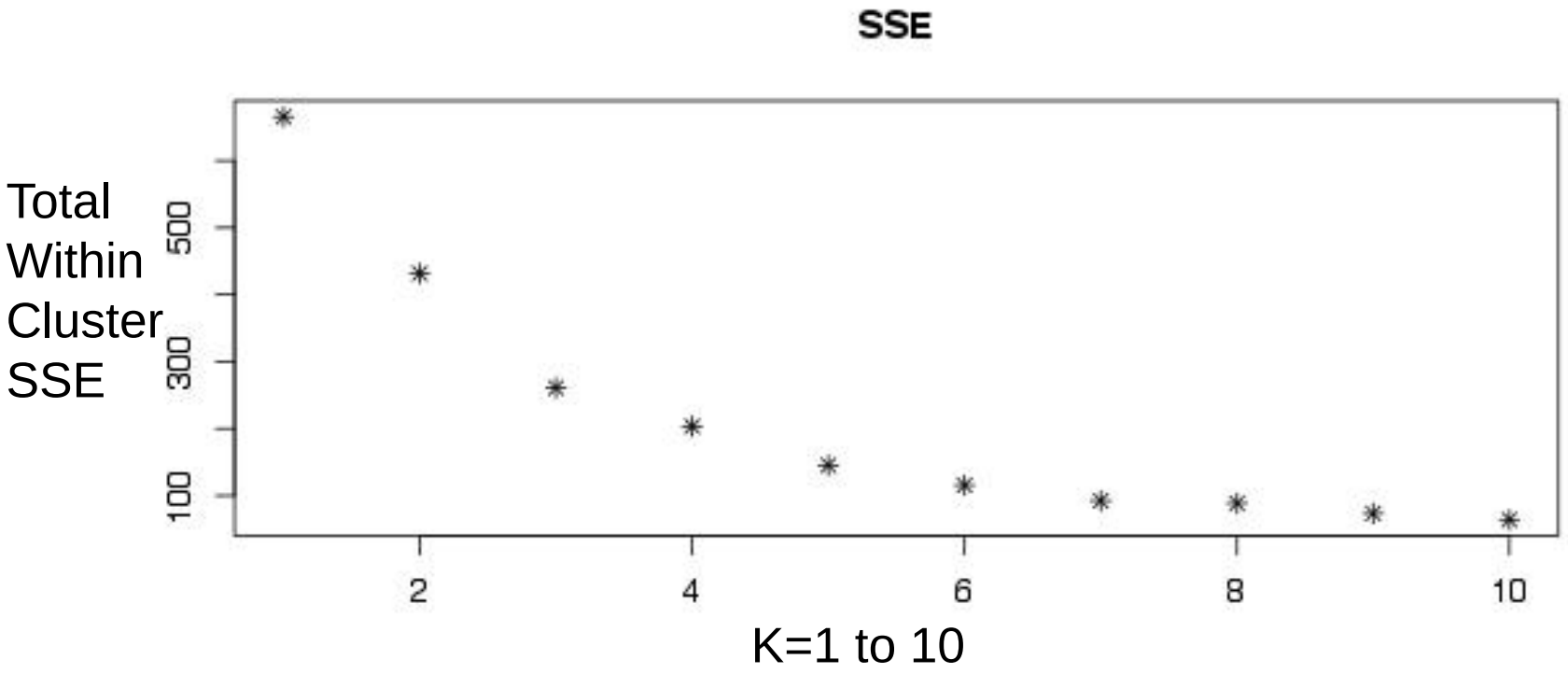


Can you guess K?

Kmeans – unequal cluster variance

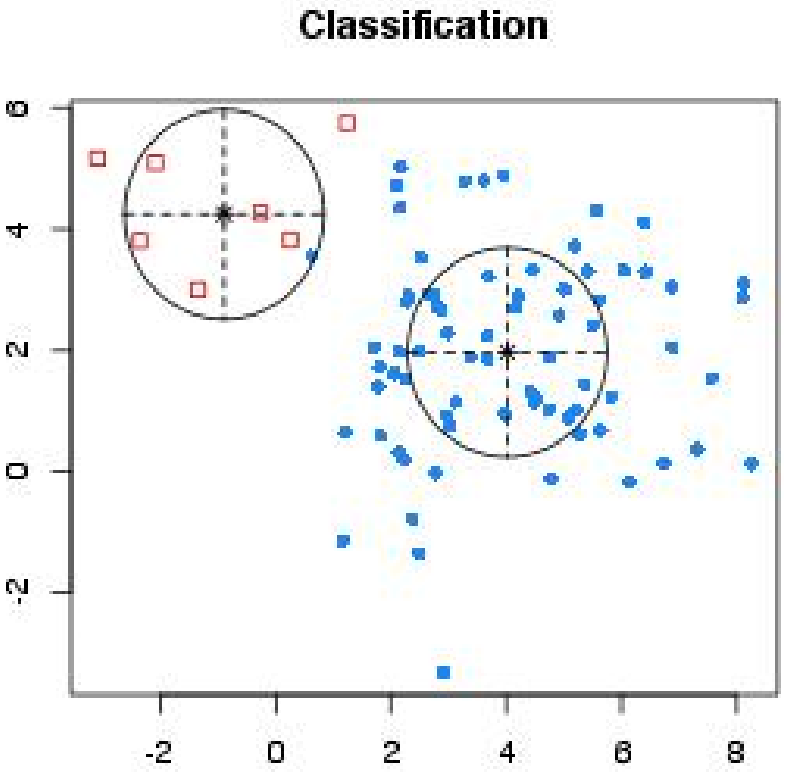


Choosing K – unequal distributions



- Smooth decrease across K => less structure (try Silhouette or Gap?)

EM clustering



- Selects K=2
  (either by Information Criterion=
  min of SSE+ K*logN,
  Or by cross-validation)

- Handles unequal variance and/or size

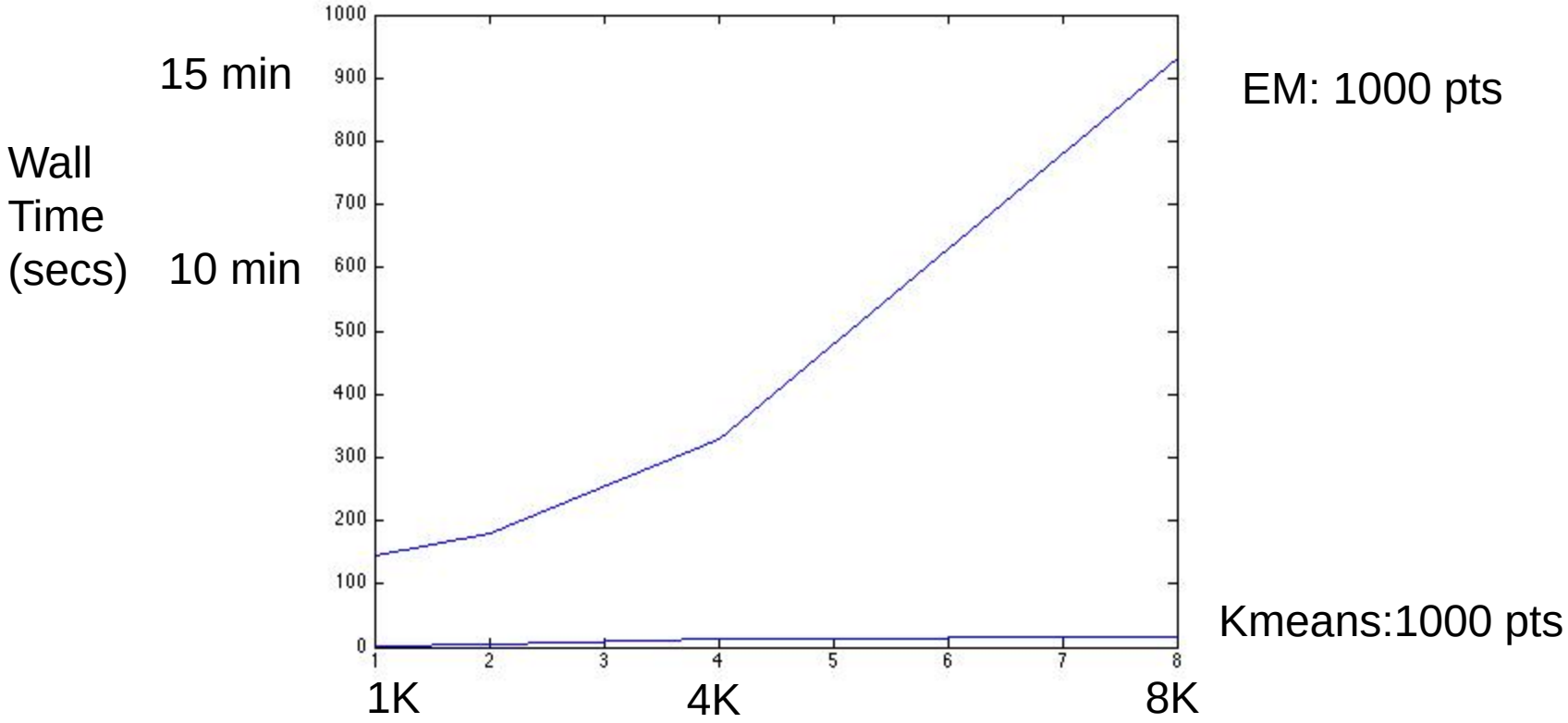
R:
library('mclust')
em_fit=Mclust(x);
plot(em_fit);

Kmeans computations

- **Distance of each point to each cluster center**
  - For *N* points, *D* dimensions: each loop requires *N*D*K* operations
- **Update Cluster centers**
  - only track points that change, get change in cluster center
- **But for EM errors to each cluster center update a probability function**

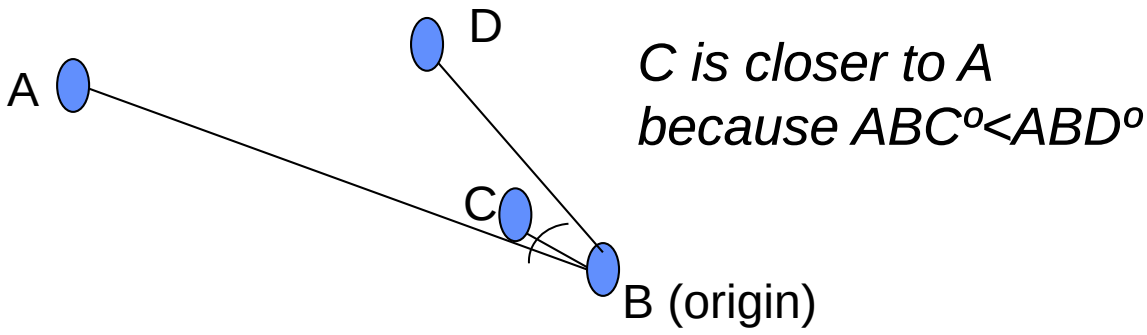
Kmeans vs EM performance

1 Gordon compute node, normal random matrices
R: system.time(Mclust())



Other distance measures

- Cosine: take angle difference (good for sparse vectors)

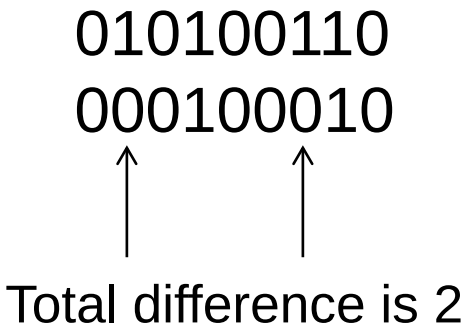


- Mahalanobis: dimensions rescaled by variance

- Jaccard (over sets A,B): 1- (|A∩B| / |AUB|)
  (e.g. sets of words in documents)

Other distance measures

- Hamming distance: count 1 if values different
  e.g. appropriate for binary strings

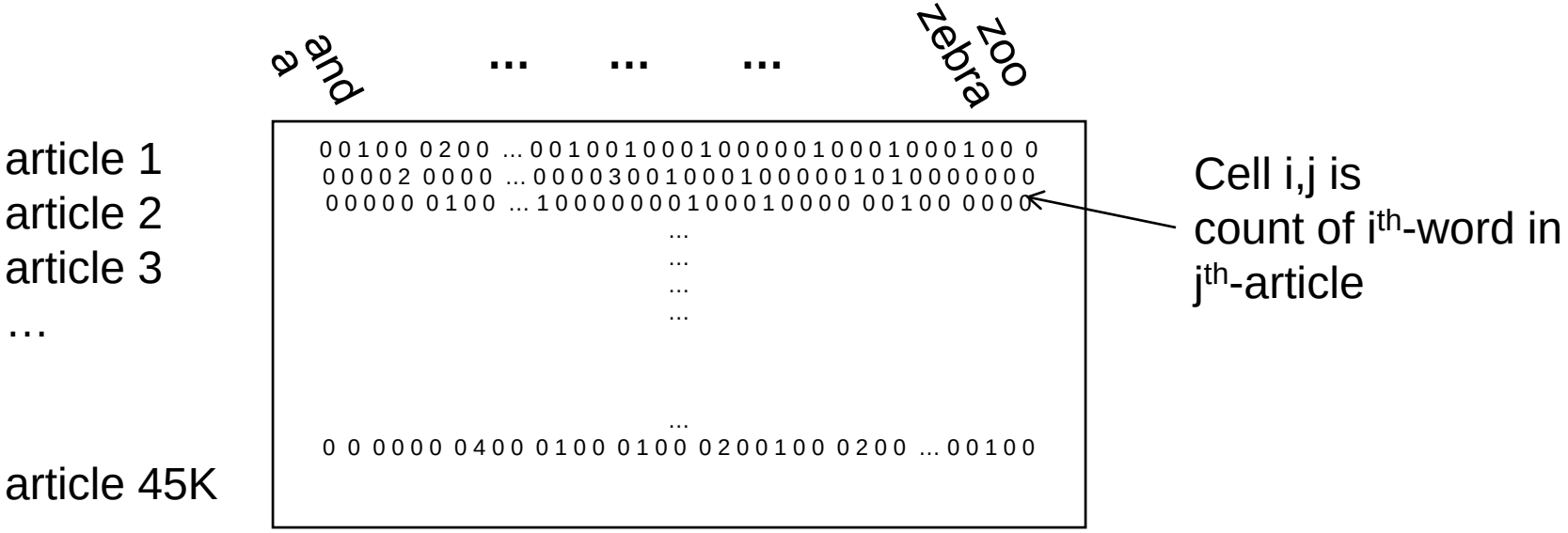


Kmeans big data example

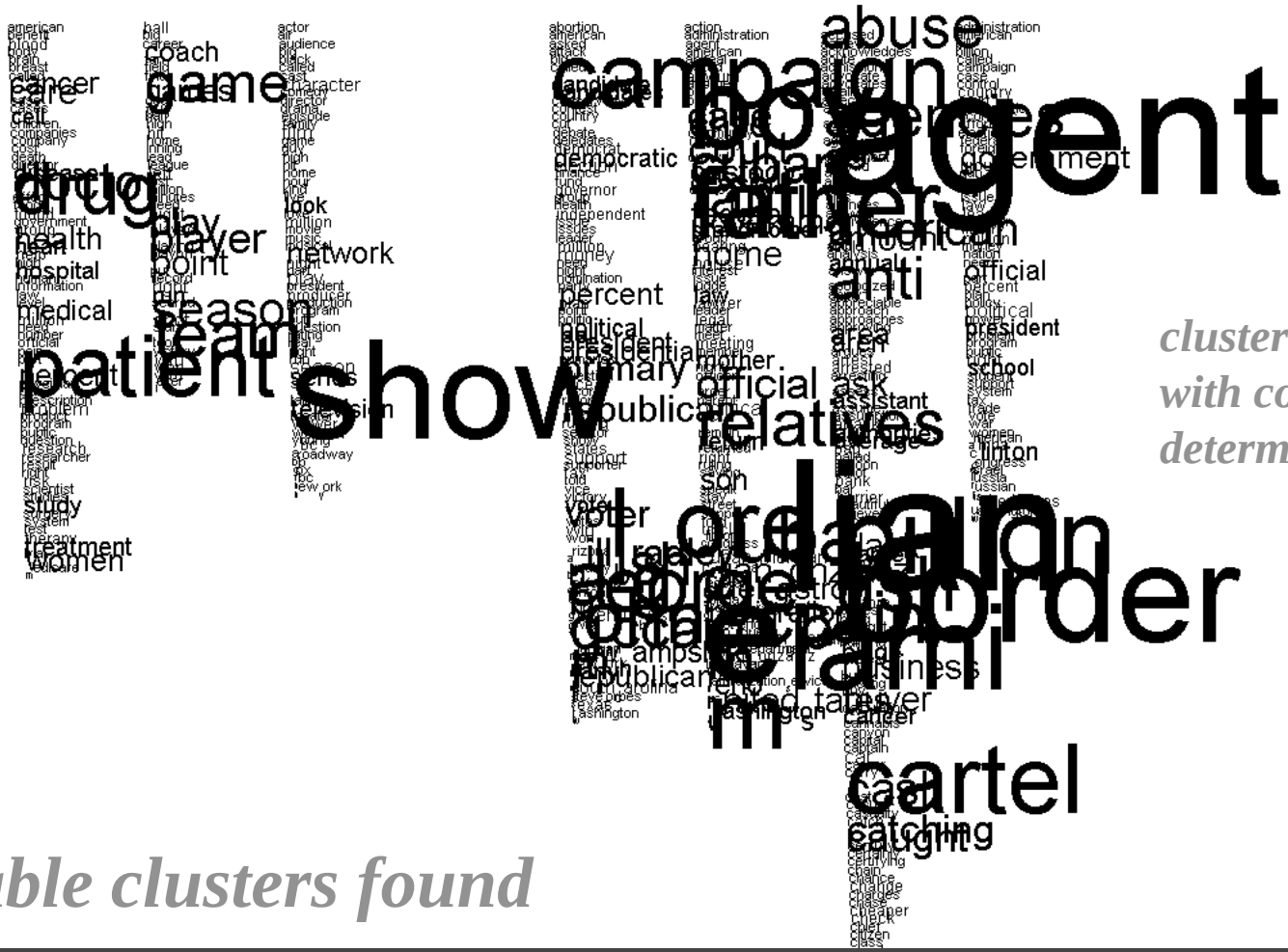
- **45,000 NYTimes articles, 102,000 unique words**

  (UCI Machine Learning repository)

- **Full Data Matrix: 45Kx102K ~ 40Gb**



Kmeans results



cluster means shown with coordinates determining fontsize

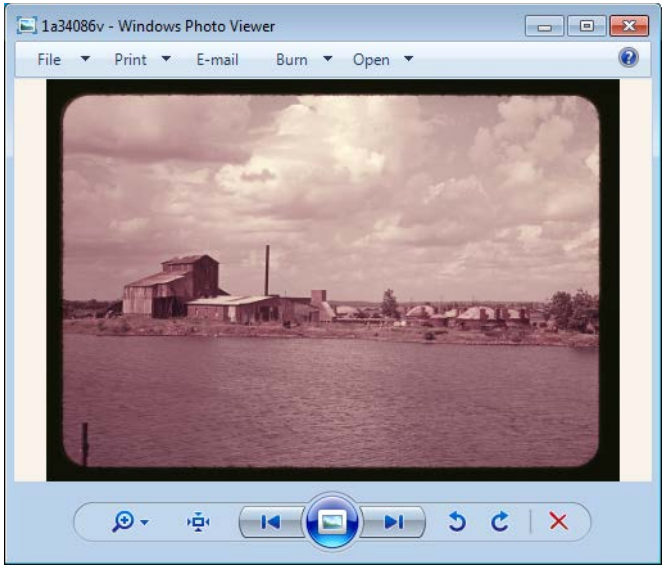
7 viable clusters found

Kmeans for image segmentation

R snippet

get packages
read 1024X718X3 RGB image
convert to matrix 1024*718 X 3
Choose K by trial and error
run Kmeans and display

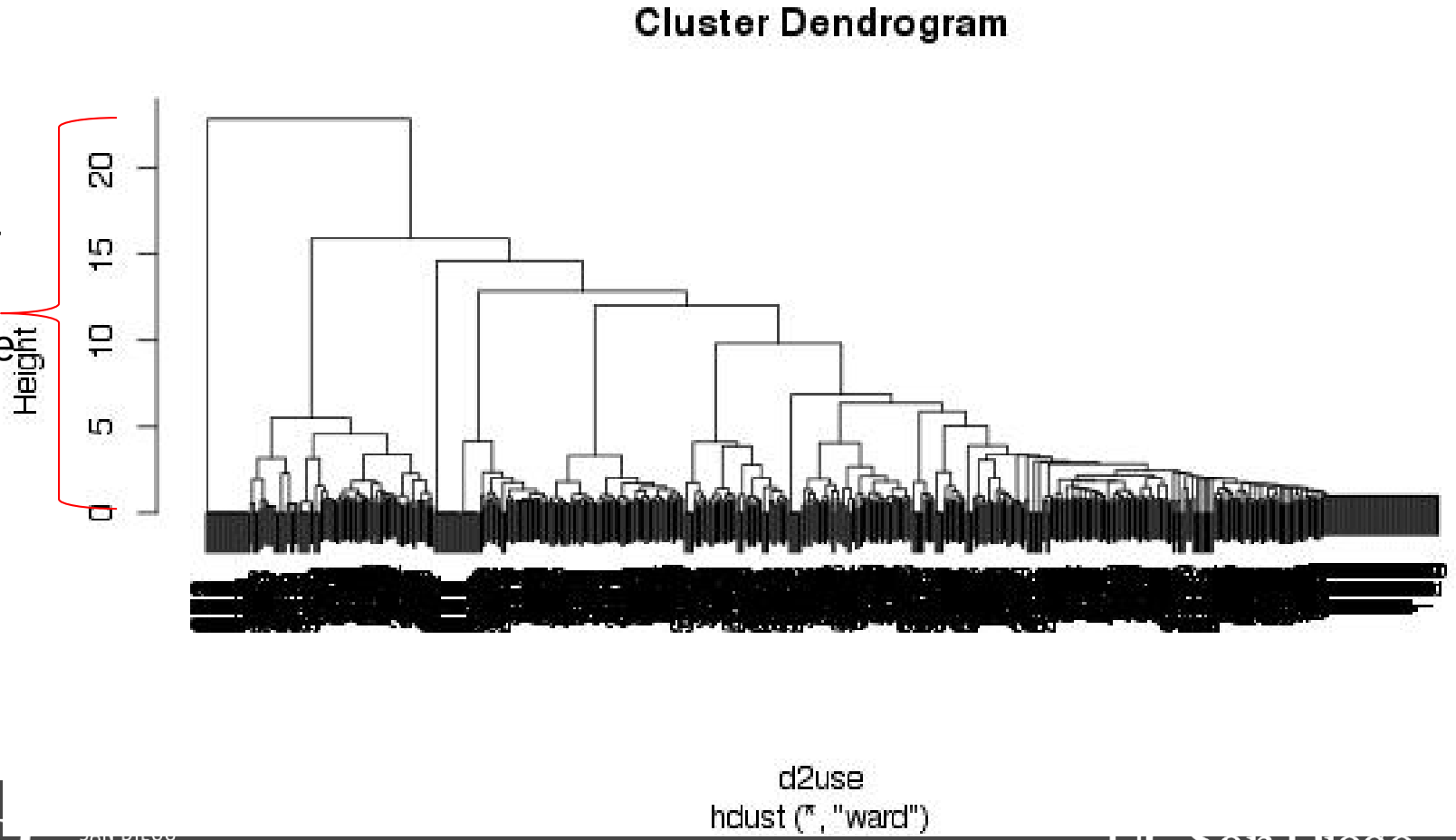
```
install.packages('ripa')
library('ripa')

source("http://bioconductor.org/biocLite.R")
biocLite()
biocLite("EBImage")

library('EBImage')
im=readImage('1a34086v.jpg')

library('ripa')
img=rgb2grey(im, coefs=c(0.30, 0.59, 0.11))

imgx1  =as.vector(img)
numk=8
km_imx1=kmeans(imgx1,numk,50,1);
img_km_mat =matrix(km_imx1$cluster,dim(im)[1],dim(im)[2])

display(img_km_mat/numk)
```



Other Clustering Methods

Hierarchical Clustering

- **hclust with “Ward” distance gives spherical clusters**

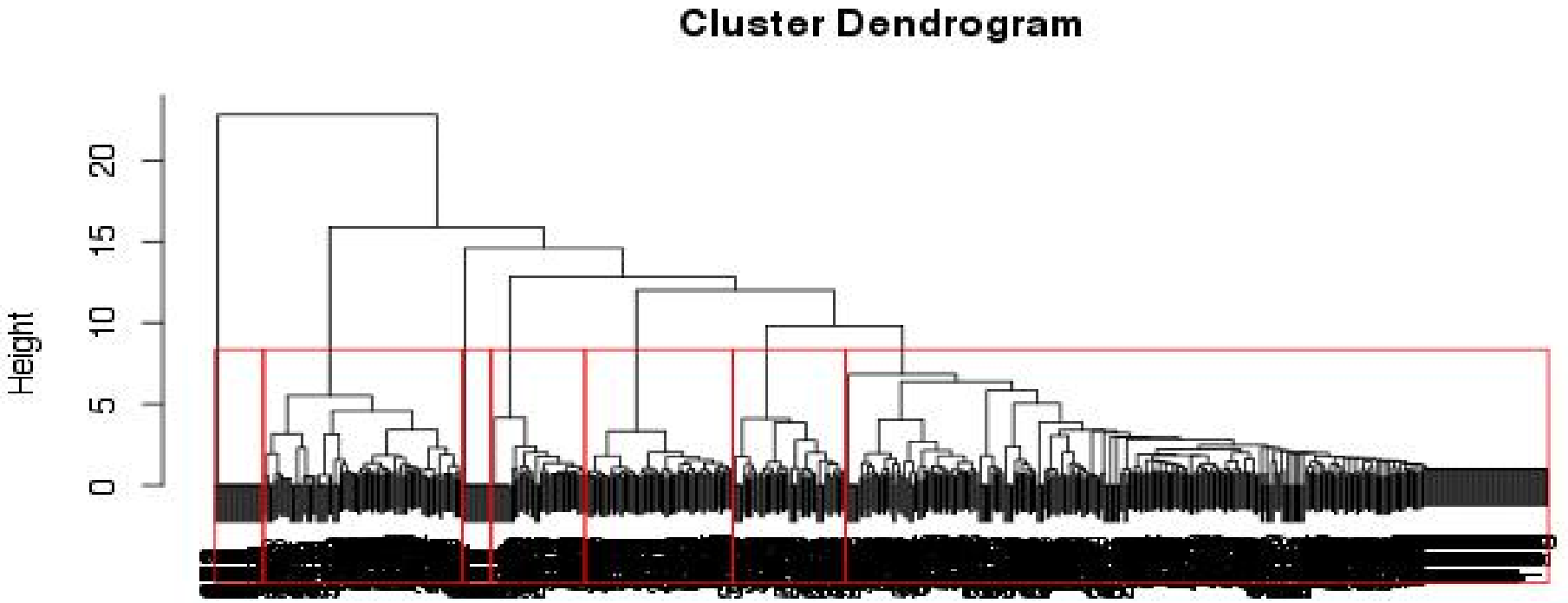


the cluster distance when 2 are combined

Hierarchical Clustering

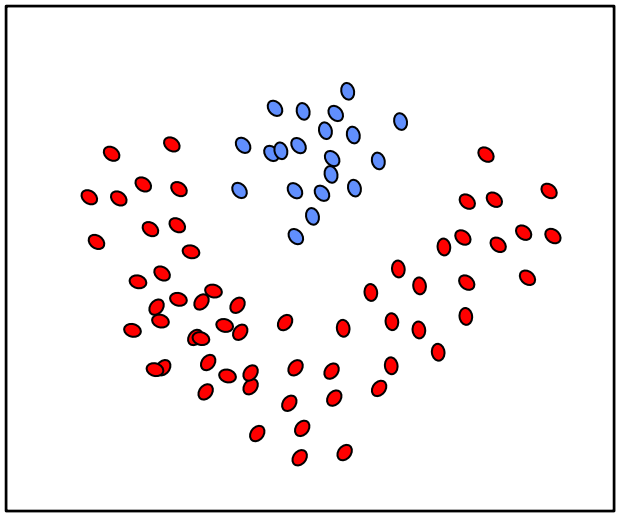
- **Where height change looks big, cut off tree**

```
groups <- cutree(fit, k=7)
rect.hclust(fit, k=7, border="red")
```



Other Clustering

- **Density based clustering**
  - build neighborhoods around seed points
  - link neighborhoods
  - Results in arbitrary cluster shapes, good for image and spatial clustering



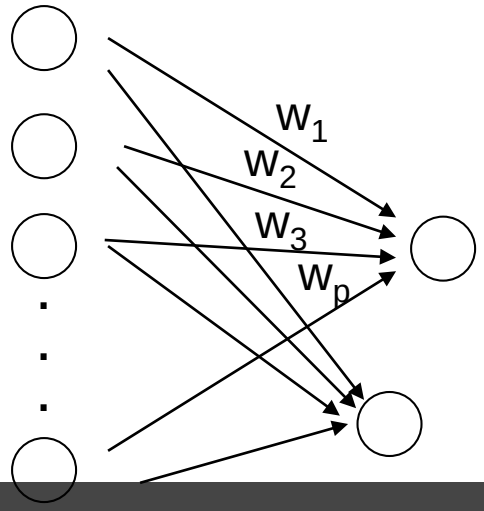
Other Clustering

- **Neural Network Based (e.g.)**

  initialize weights to coordinate values for a seed point

  set input nodes to data points

  get best match to seed for each data point and adjust weights toward the data point

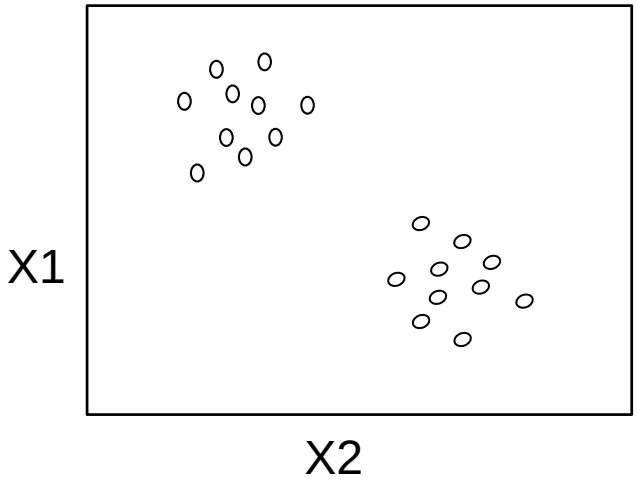


Target node(s), starts as a seed point and ends up as a cluster mean

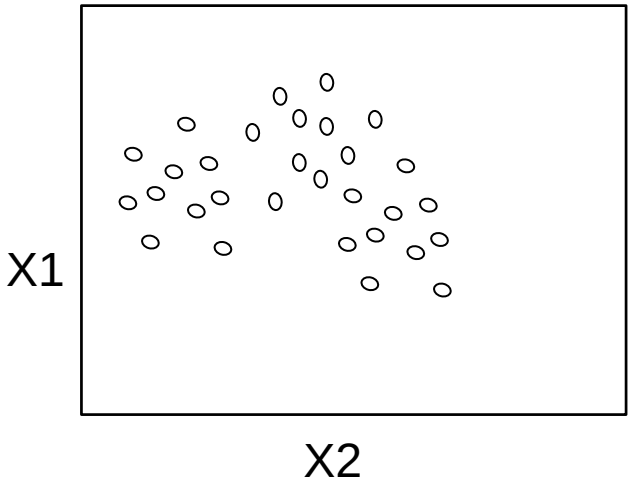
Pause

Imagine these 2 dimensional input spaces:
Which of these is easy or hard to cluster? (no class labels)

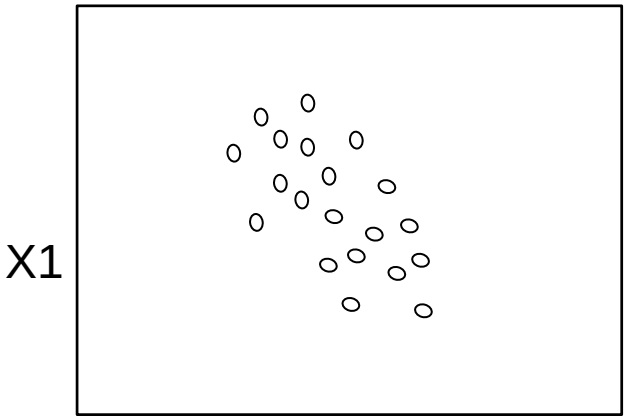
A



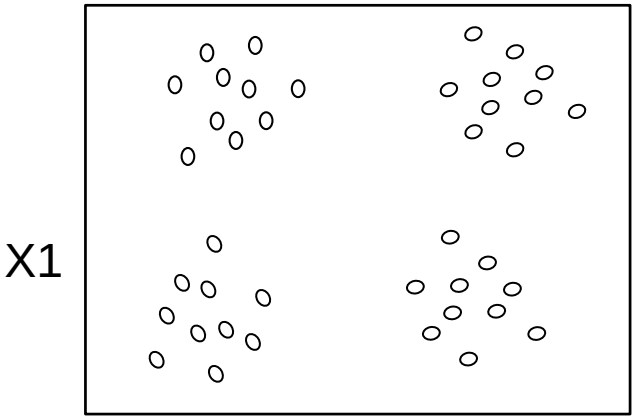
C



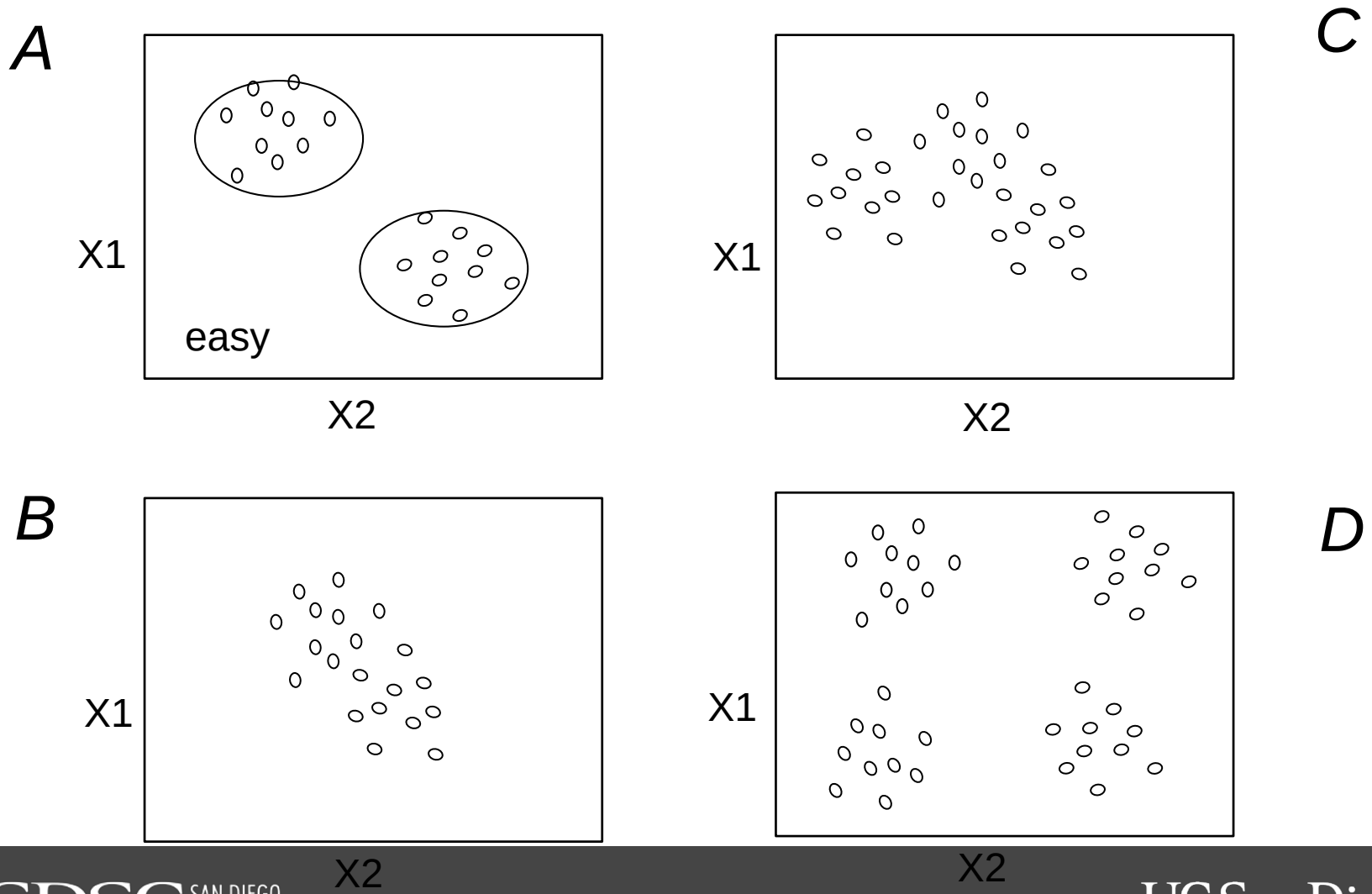
B



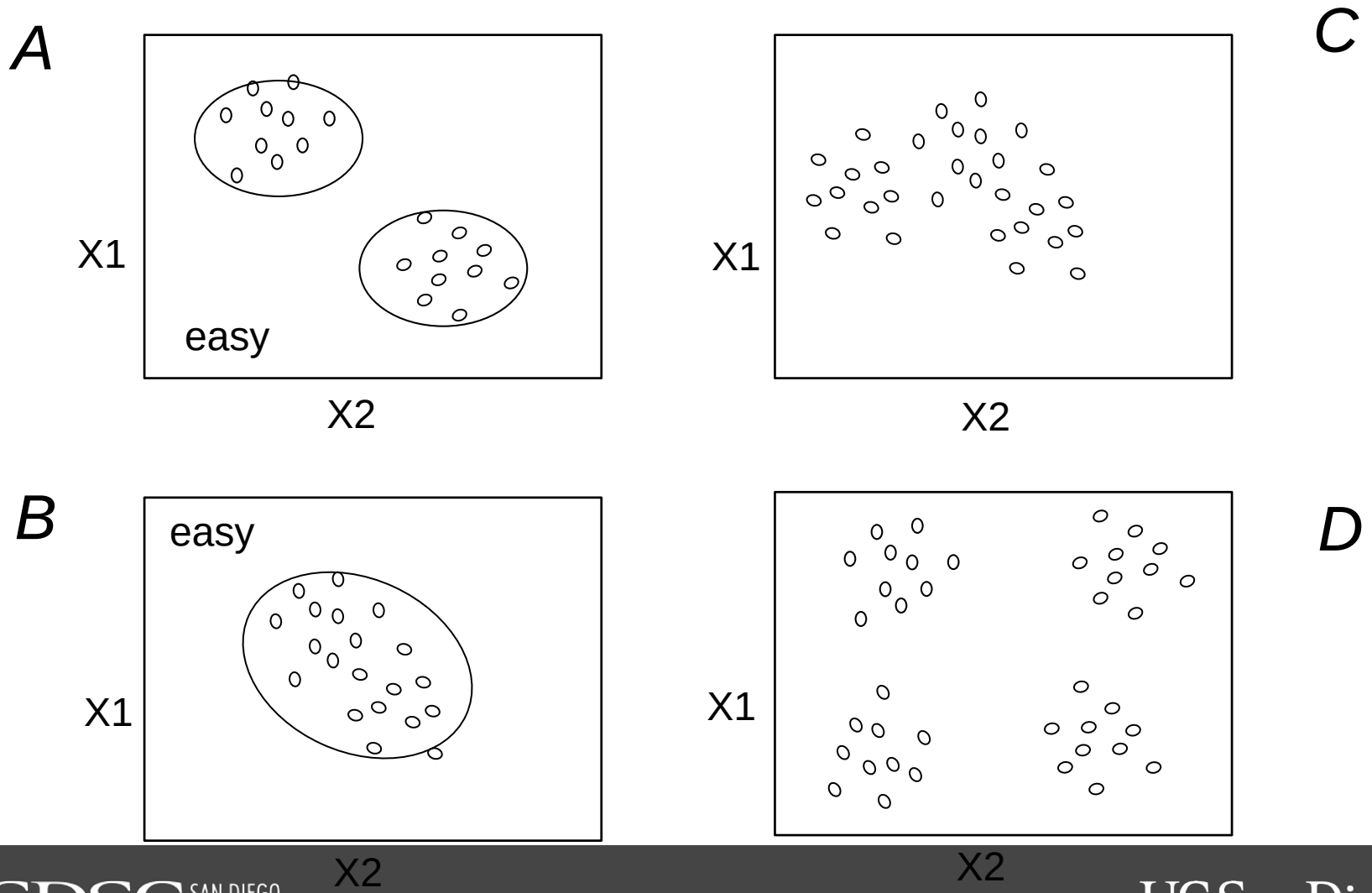
D



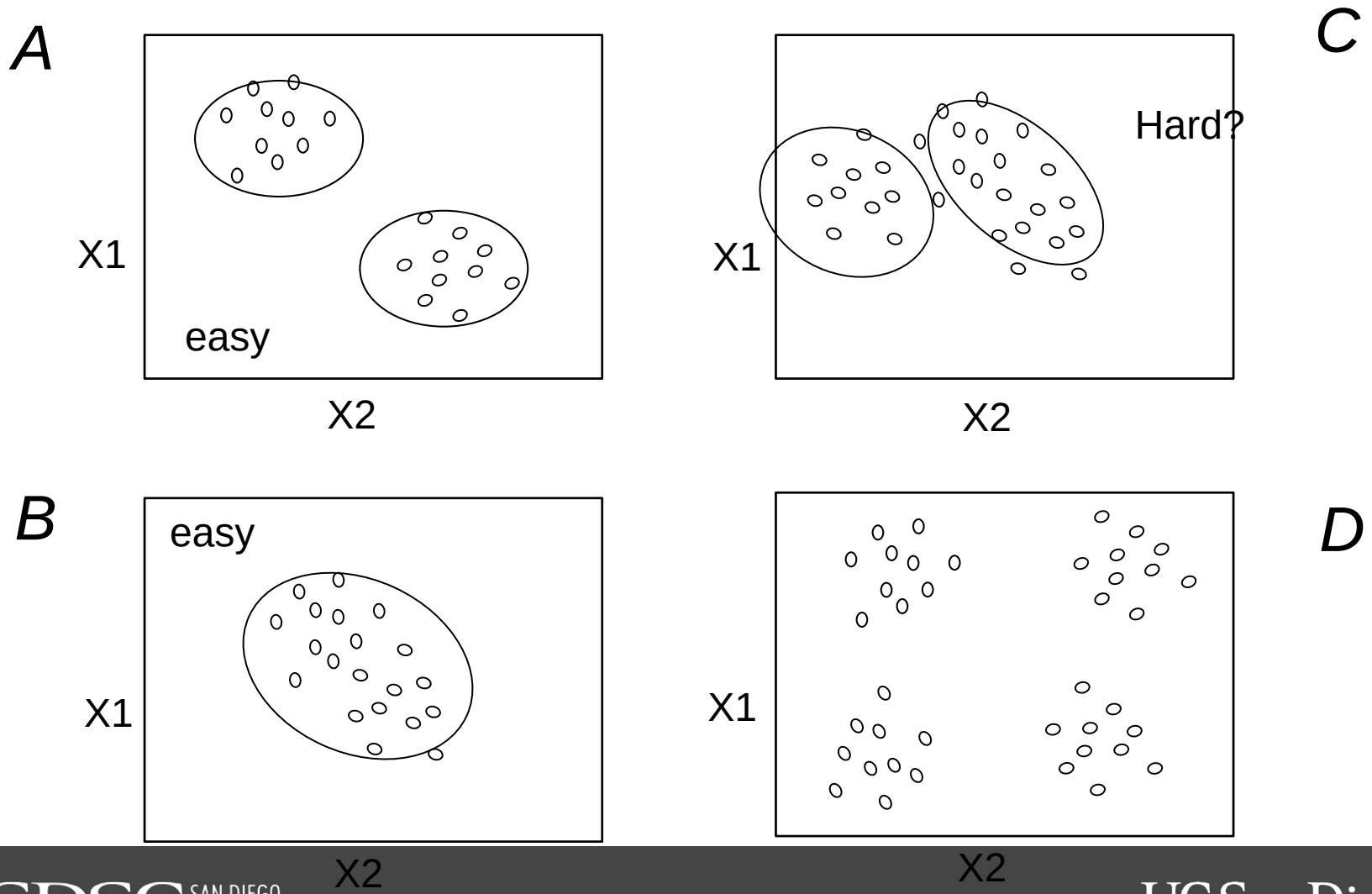
Potential clusters



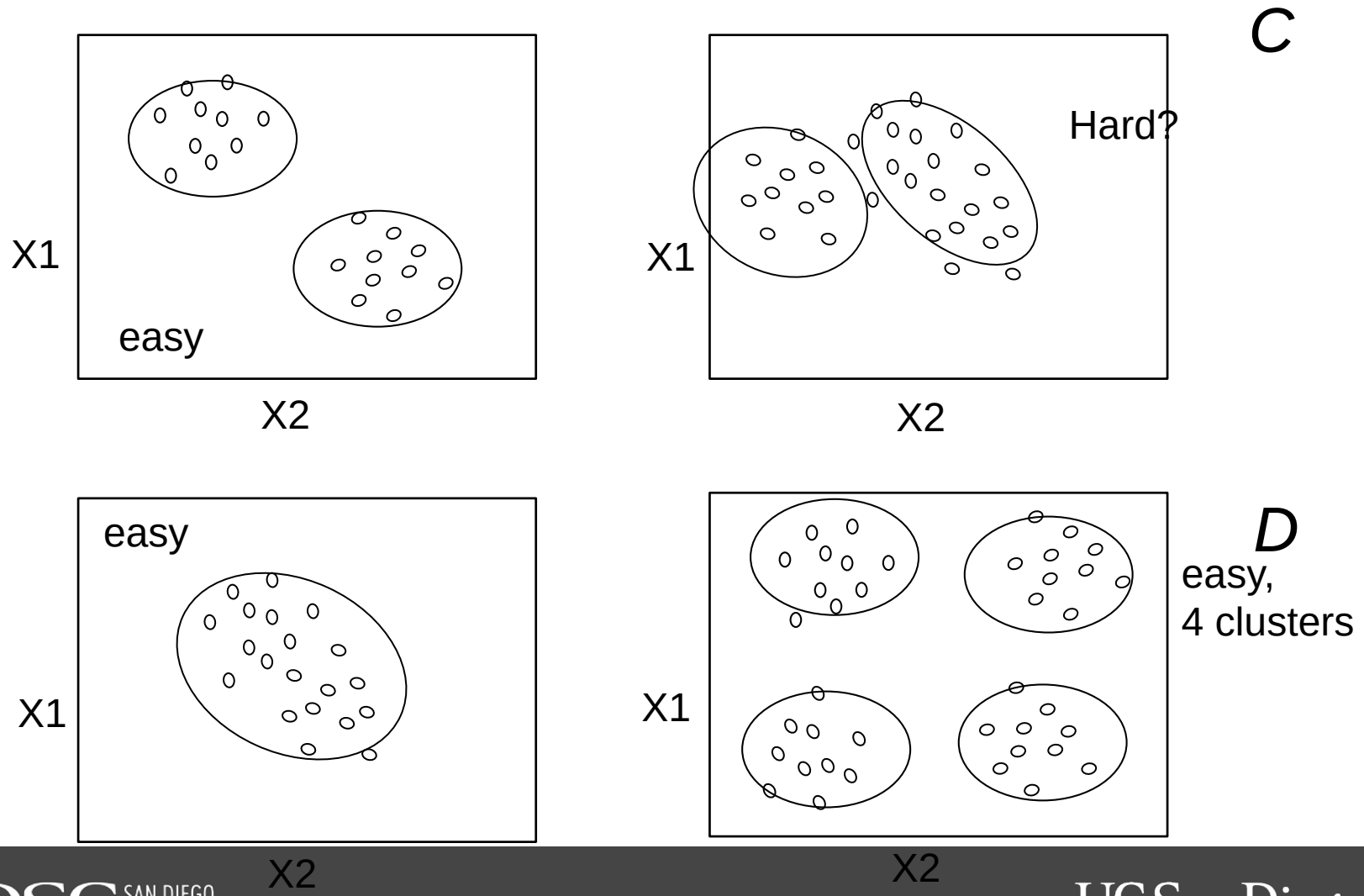
Potential clusters



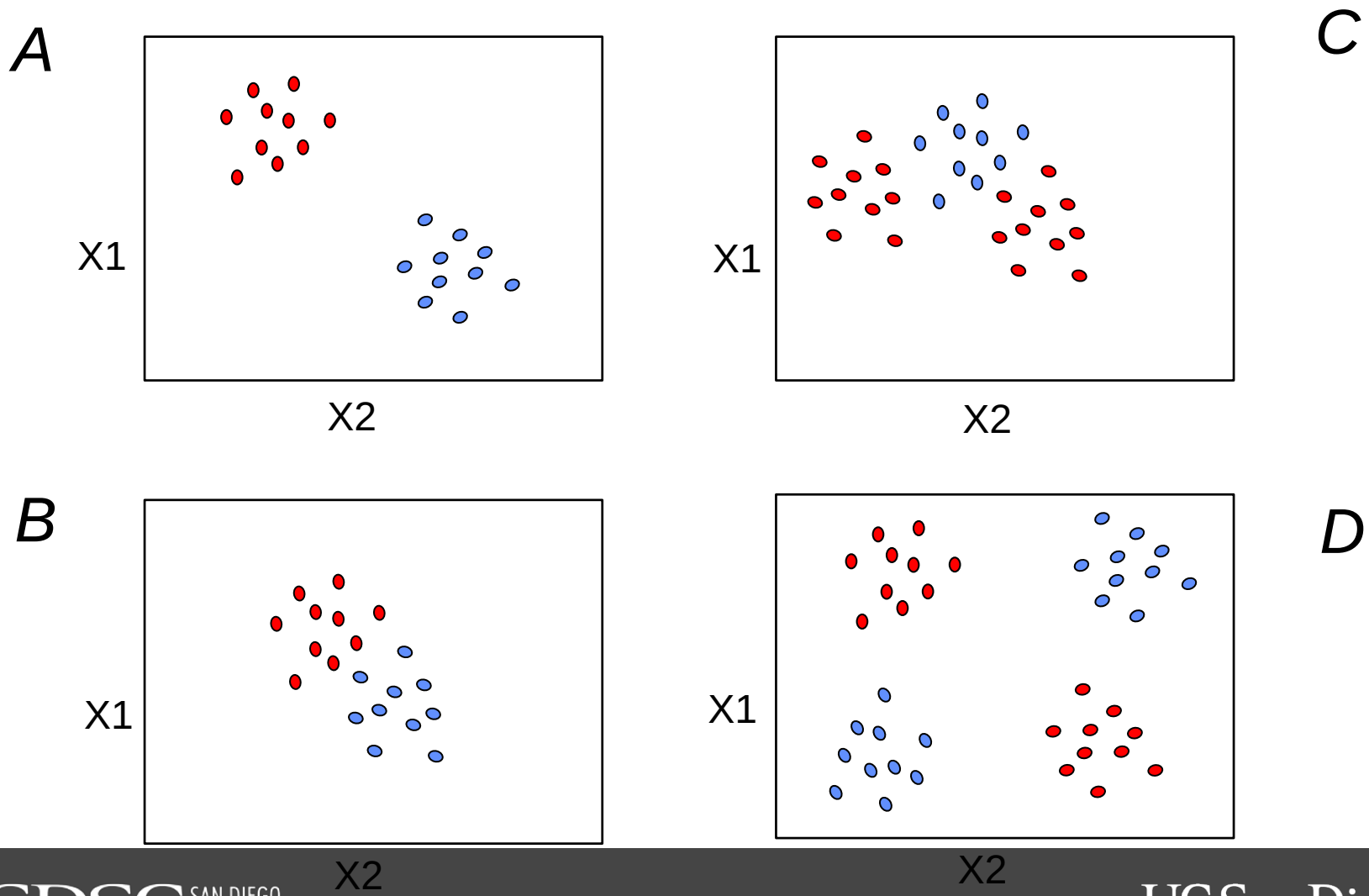
Potential clusters



Potential clusters

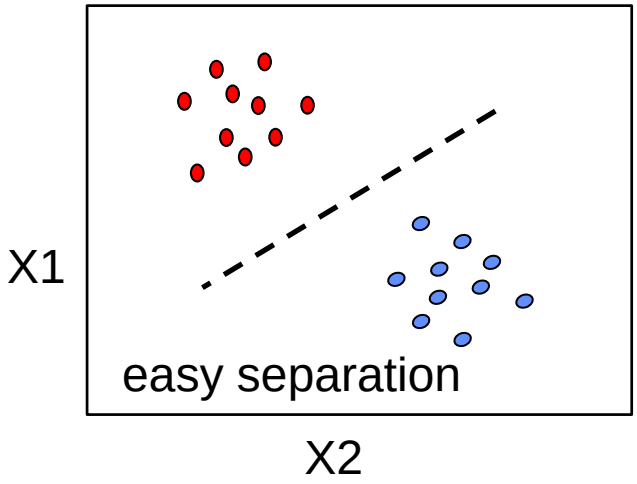


Now imaging there are two classes

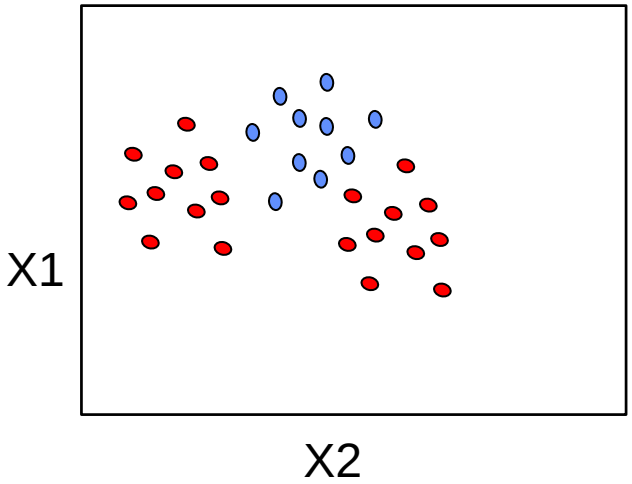


Which are easy or hard to classify?
(ie separate red or blue with lines)

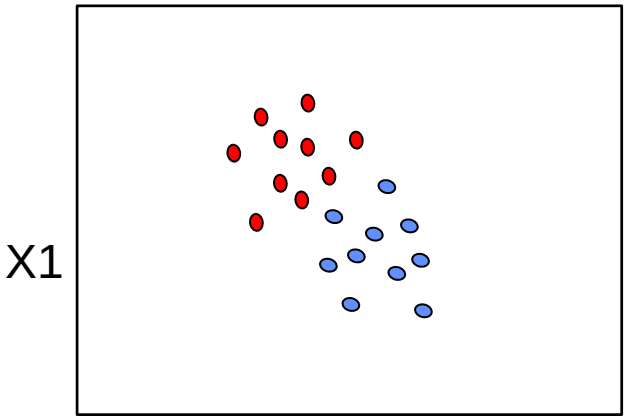
A



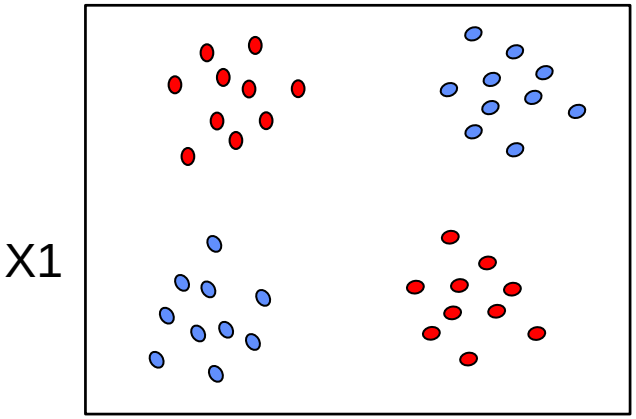
C



B

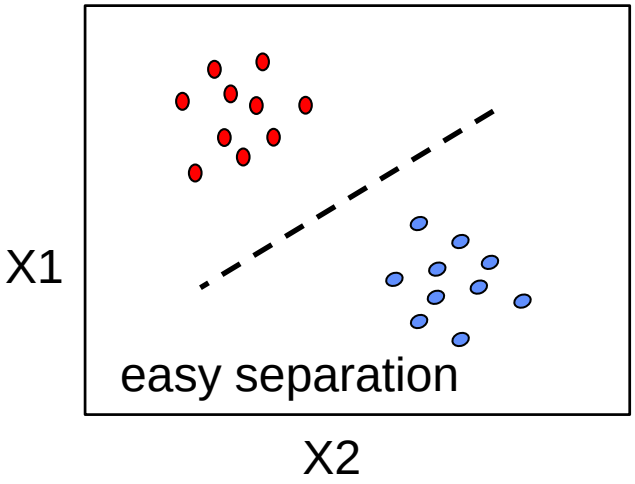


D

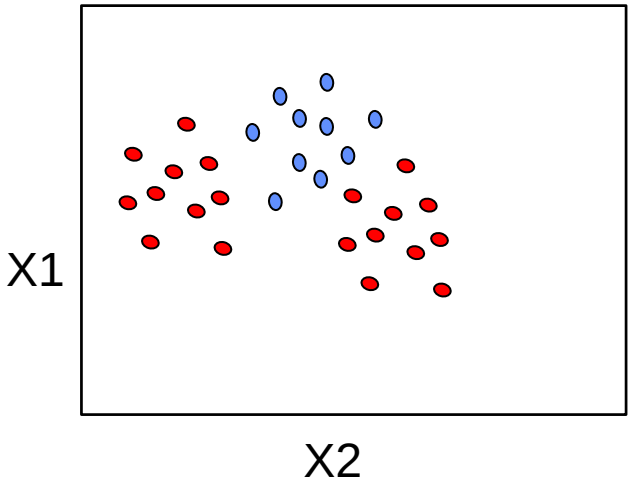


Which are easy or hard to classify?
(ie separate red or blue with lines)

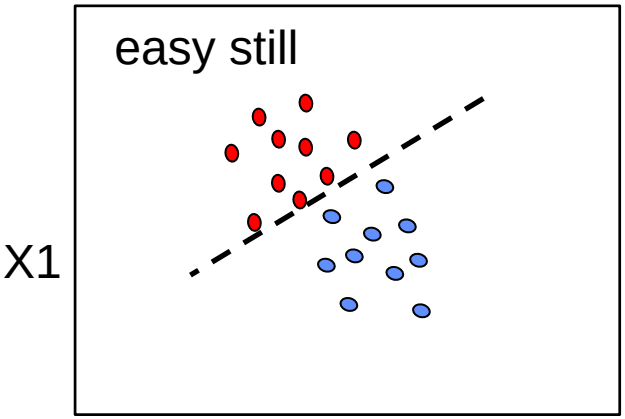
A



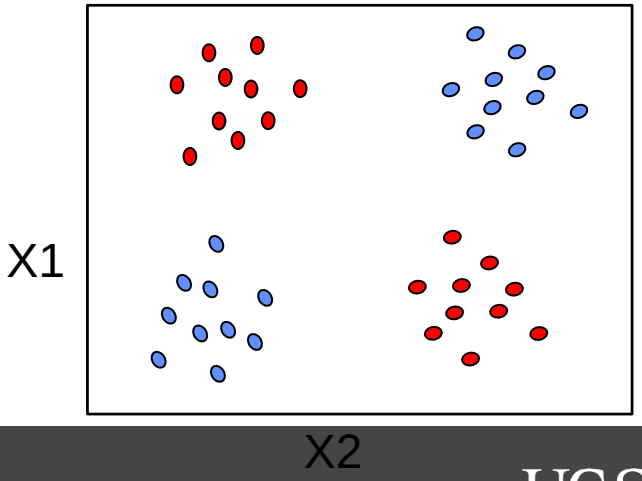
C



B

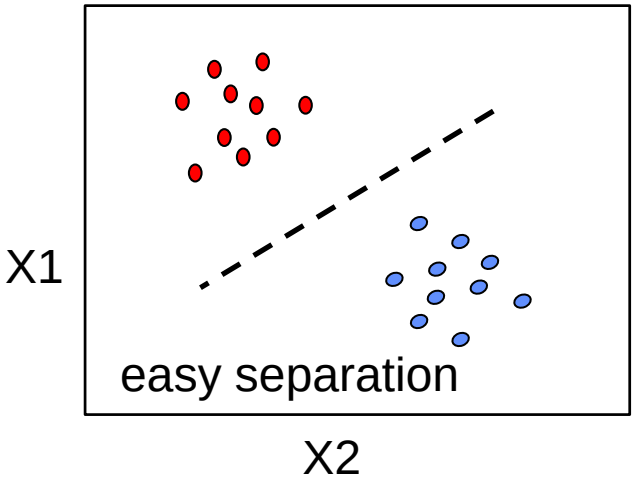


D

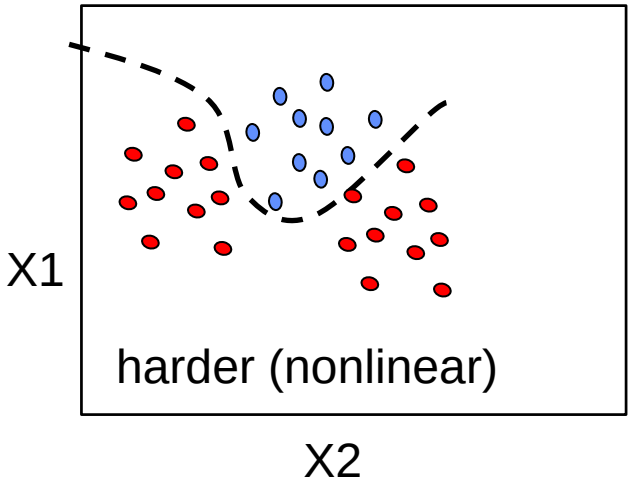


Which are easy or hard to classify?
(ie separate red or blue with lines)

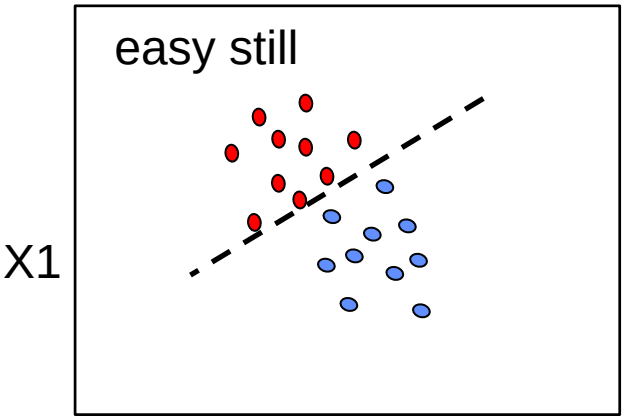
A



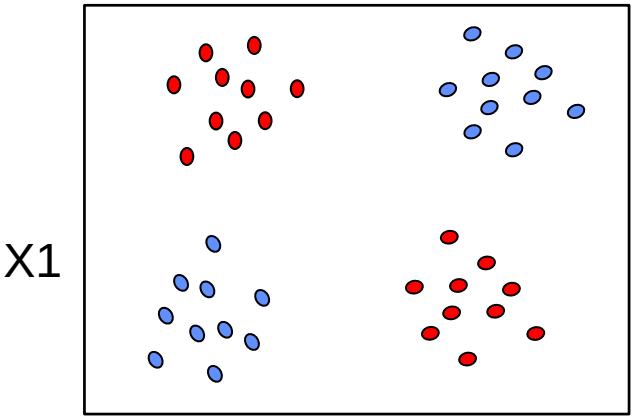
C



B

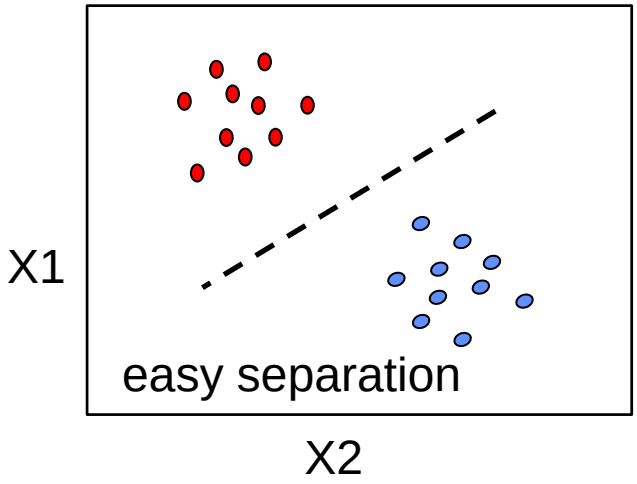


D

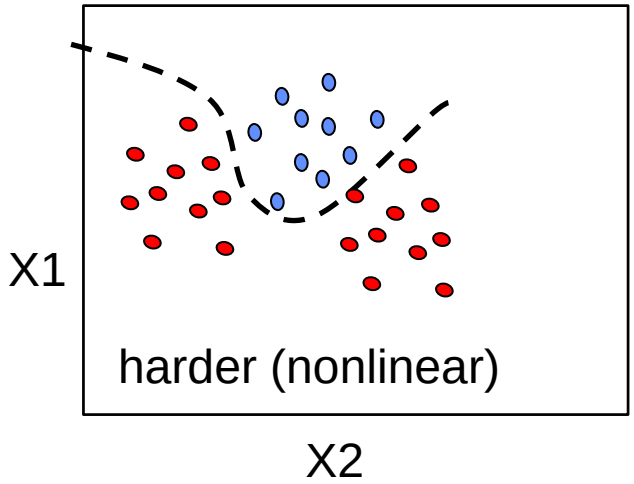


Which are easy or hard to classify?
(ie separate red or blue with lines)

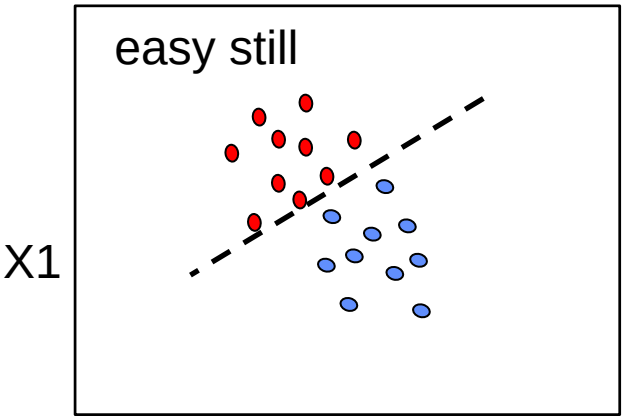
A



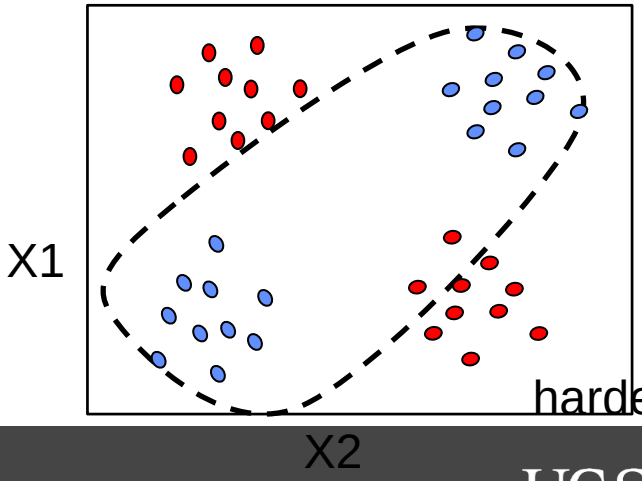
C



B

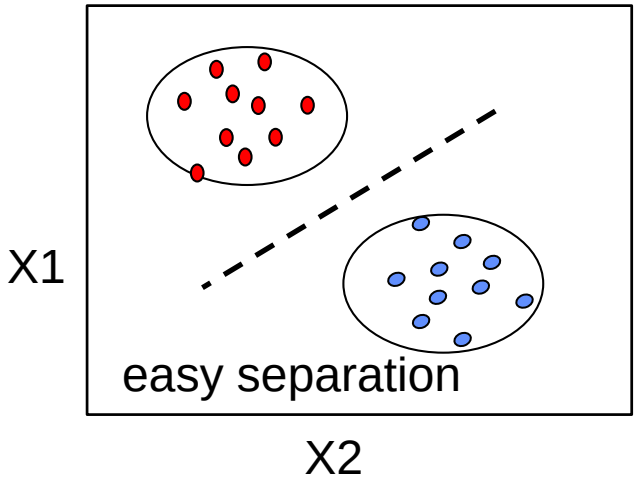


D

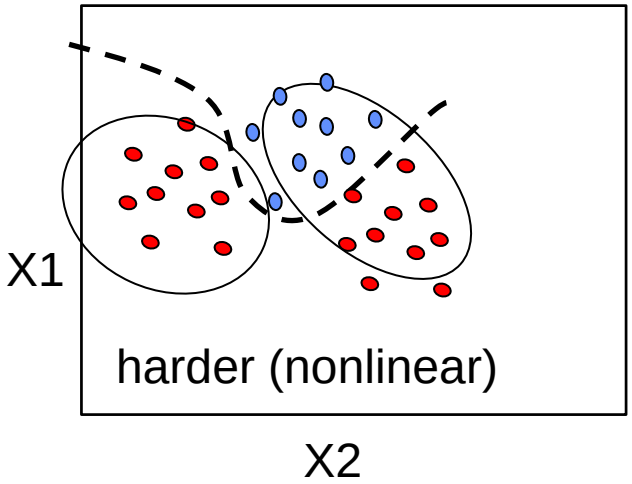


Which are easy or hard to classify?
(ie separate red or blue with lines)

A

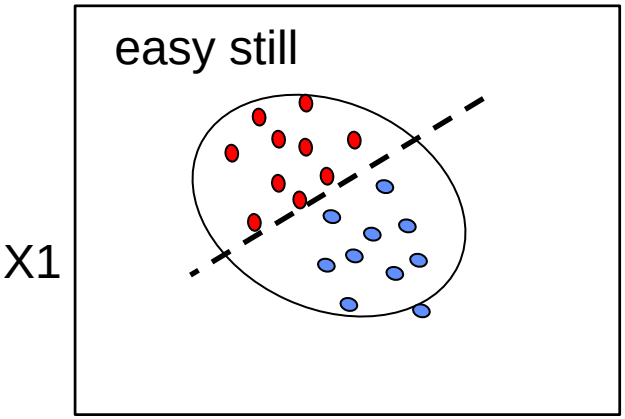


C

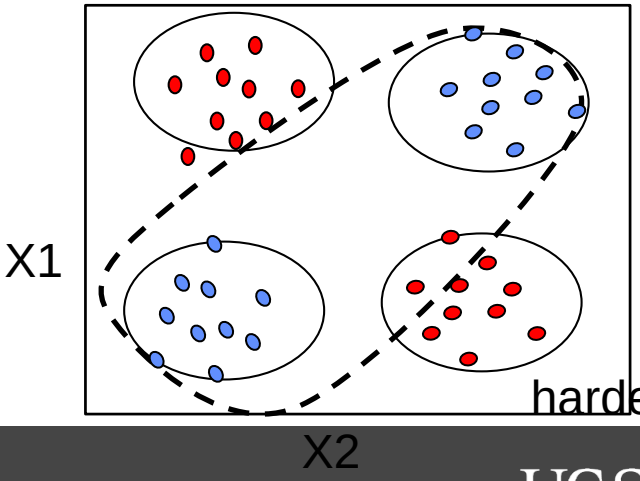


In summary:
No easy relationship between clusters and classification

B



D



Pause

Exercise, Kmeans and visualization in 2-D using SVD

- Using same data from SVD exercise, and SVD reduced matrix
- Run Kmeans
- Project data points onto first 2 SVD factors and plot them colored by cluster

(Which 2 factors?

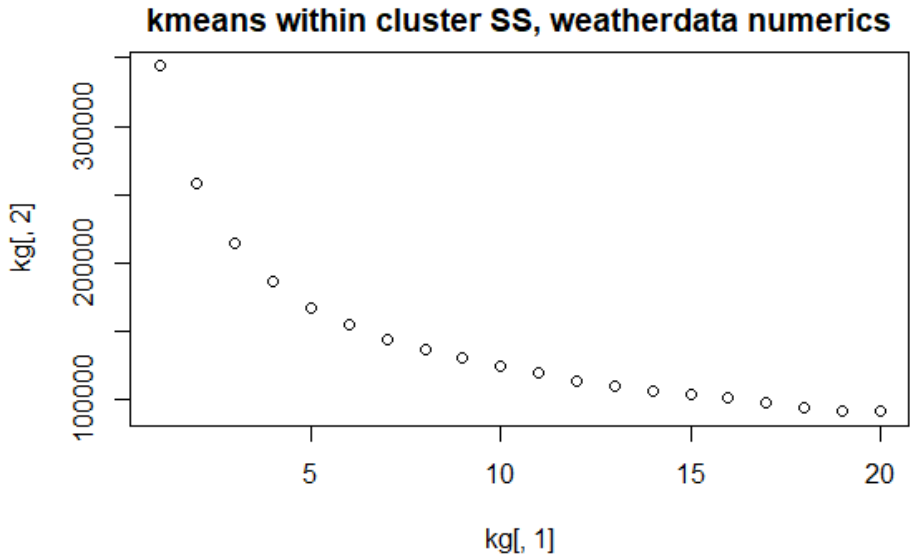
Keep in mind that:

Wsvd$U is N rows x P cols

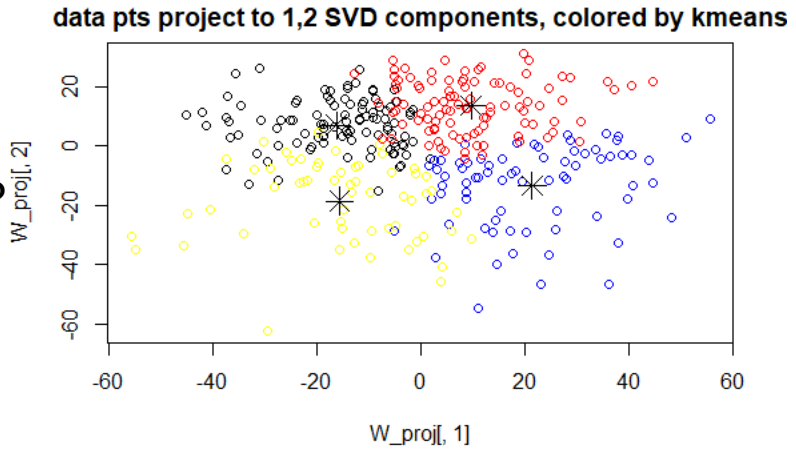
Wsvd$V is PxP,

Data point is Px1 in original space and 2x1 in reduced space)

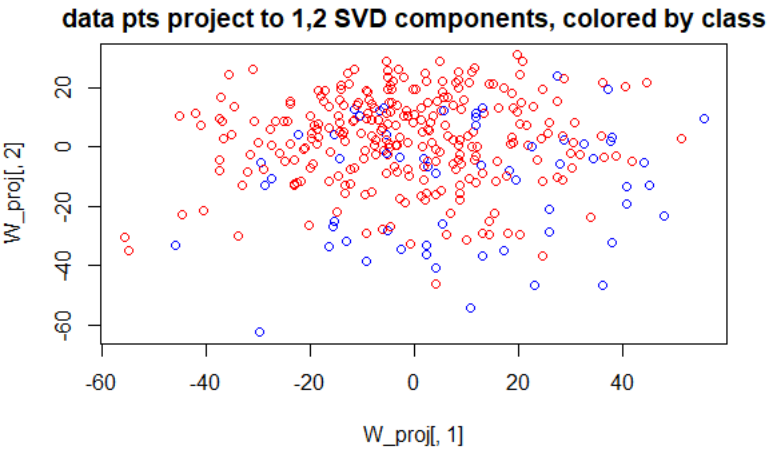
See clustering_exercise Rmd file
Rerun the SVD exercise if you need to, to get the W_mncntr and Wsvd matrices
Run Kmeans



Run Kmeans with k=4
Plot data points onto 2-D space, use colors determined by Kmeans
Plot cluster centers



Are points well separated in 2-D projection?
Should they be?



Plot using colors determined by class (raintomorrow)

Are classes well separated?

Principle Components vs Clustering

- PCA, SVD reduces dimensions, Clustering reduces to categorical groups
- In some cases, *k* PCs ⇔ *k* clusters
- It is also useful to visualize clusters in PC space

Summary

- **Having no label doesn’t stop you from finding structure in data**
- **Unsupervised methods are somewhat related**