

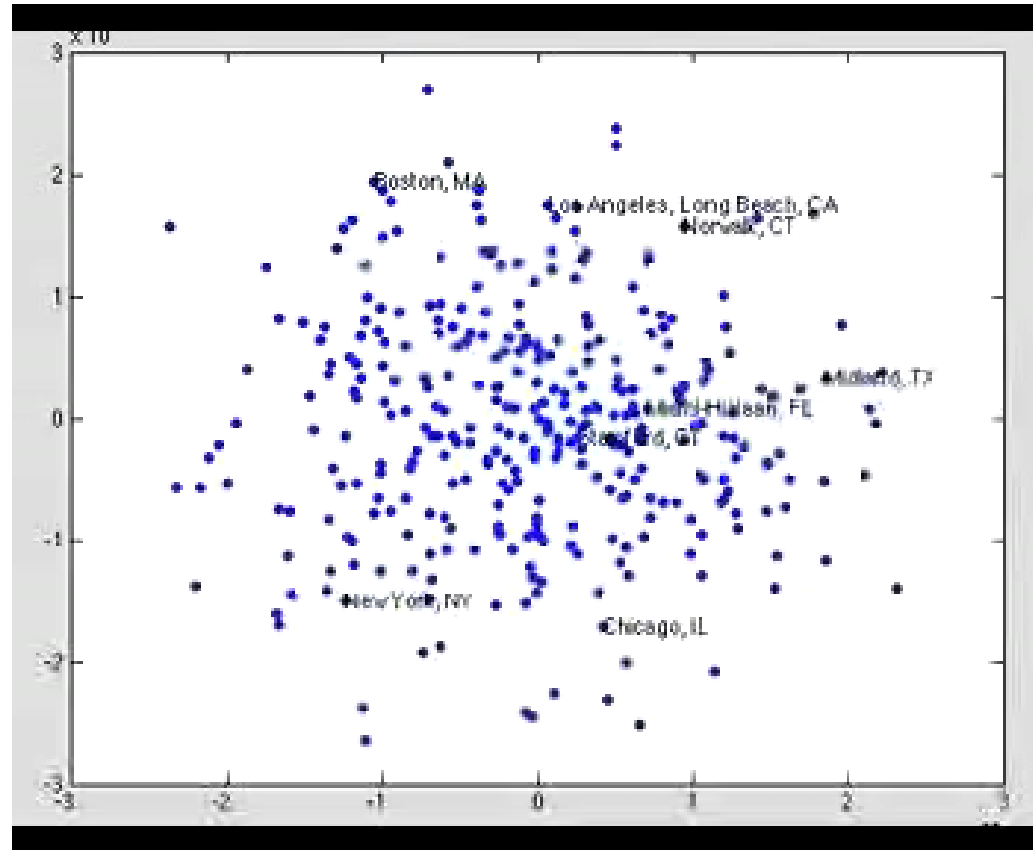
CPSC 340: Machine Learning and Data Mining

**Multi-Dimensional Scaling
Summer 2021**

Admin

- Assignment 6 out, due Friday 11:55pm
- Today is final exam coverage cut-off
- **Final exam is next Wednesday (June 23)**
 - Prep materials go up soon
- **Course evaluation is open.**
 - Please give me an honest feedback! How did I do?

Last Time: Multi-Dimensional Scaling



$$f(Z) = \sum_{i=1}^n \sum_{j=i+1}^n (\underbrace{\|z_i - z_j\|}_{\text{distance in scatterplot}} - \underbrace{\|x_i - x_j\|}_{\text{Distance between points in original 'd' dimensions}})^2$$

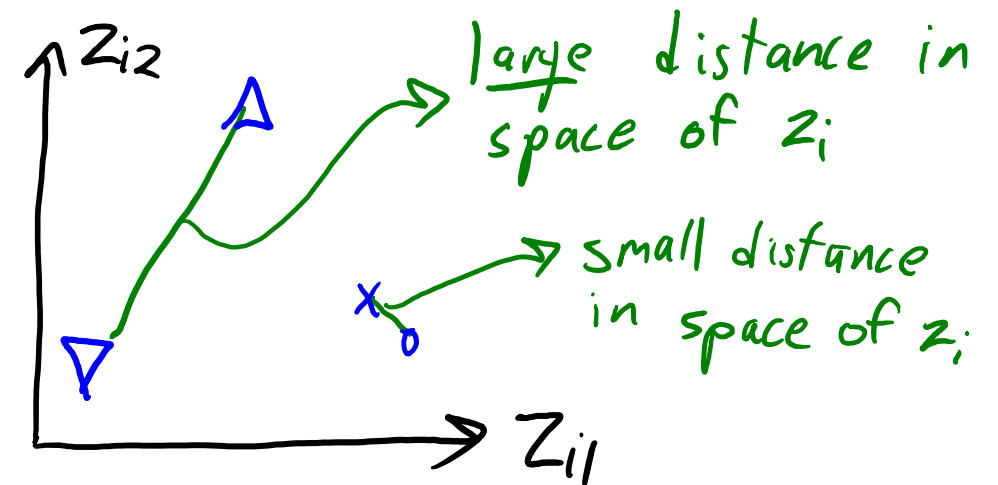
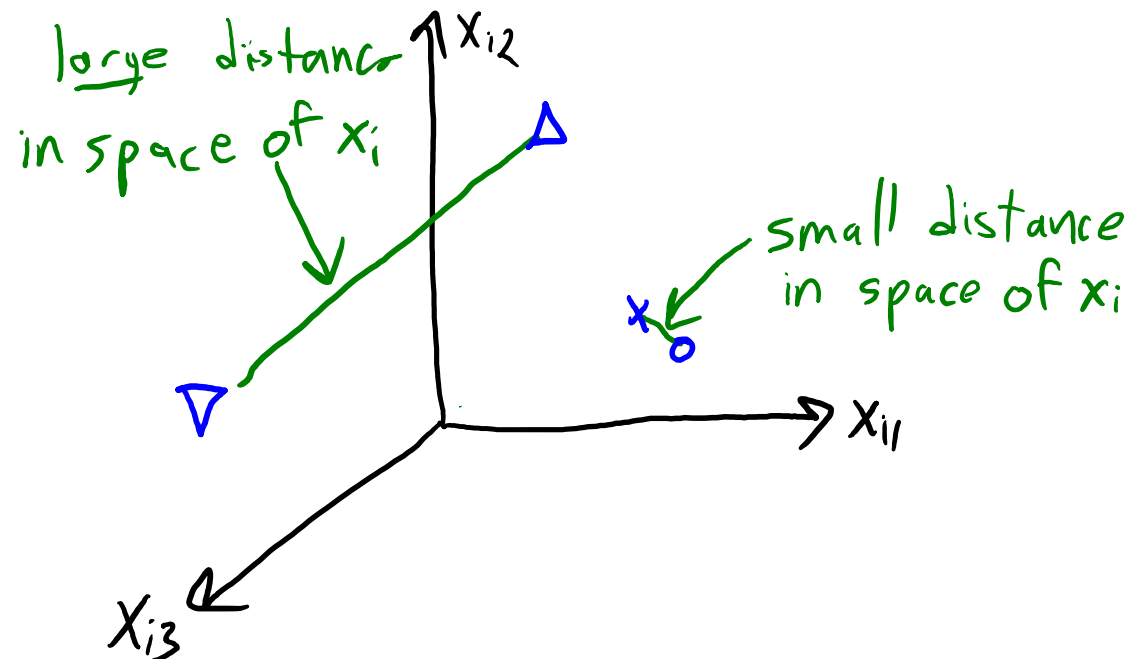
Try to make scatterplot distances match high-dimensional distance

sum over pairs of examples

Multi-Dimensional Scaling

- Multi-dimensional scaling (MDS):
 - Optimize the final locations of the z_i values.

$$f(z) = \sum_{i=1}^n \sum_{j=i+1}^n (\|z_i - z_j\| - \|x_i - x_j\|)^2$$

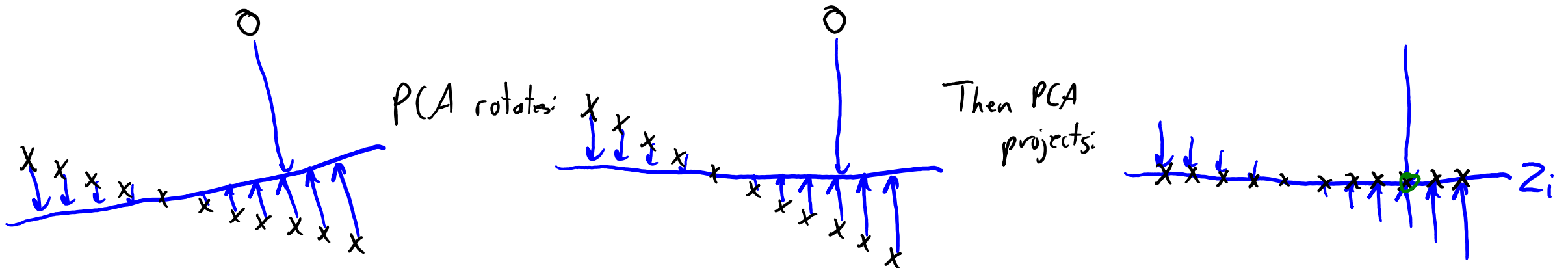


Multi-Dimensional Scaling

- Multi-dimensional scaling (MDS):
 - Optimize the final locations of the z_i values.

$$f(z) = \sum_{i=1}^n \sum_{j=i+1}^n (\|z_i - z_j\| - \|x_i - x_j\|)^2$$

- Non-parametric dimensionality reduction and visualization:
 - No 'W': just trying to make z_i preserve high-dimensional distances between x_i .

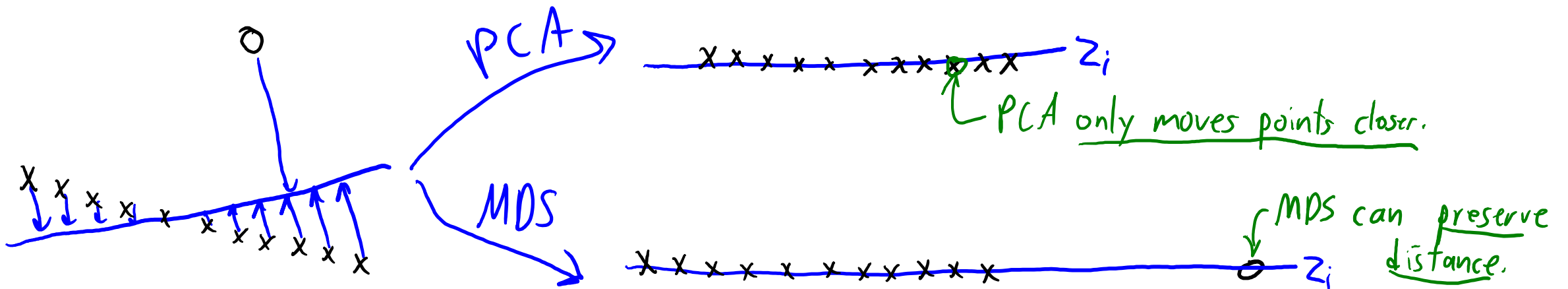


Multi-Dimensional Scaling

- Multi-dimensional scaling (MDS):
 - Optimize the final locations of the z_i values.

$$f(z) = \sum_{i=1}^n \sum_{j=i+1}^n (\|z_i - z_j\| - \|x_i - x_j\|)^2$$

- Non-parametric dimensionality reduction and visualization:
 - No 'W': just trying to make z_i preserve high-dimensional distances between x_i .

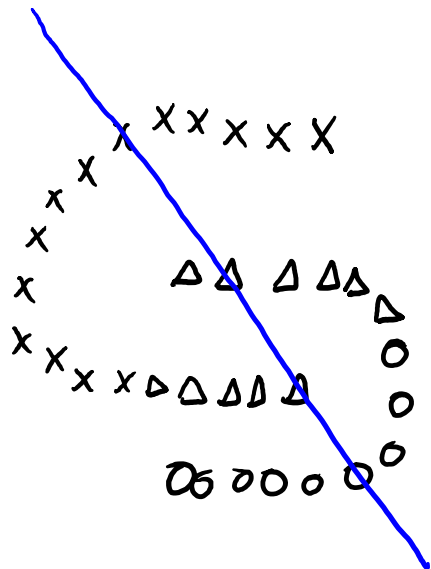


Multi-Dimensional Scaling

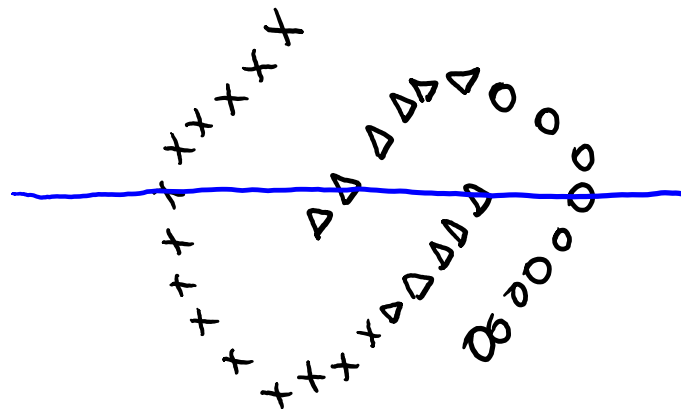
- Multi-dimensional scaling (MDS):
 - Optimize the final locations of the z_i values.

$$f(z) = \sum_{i=1}^n \sum_{j=i+1}^n (\|z_i - z_j\| - \|x_i - x_j\|)^2$$

- Non-parametric dimensionality reduction and visualization:
 - No 'W': just trying to make z_i preserve high-dimensional distances between x_i .



PCA rotation:



PCA projection:

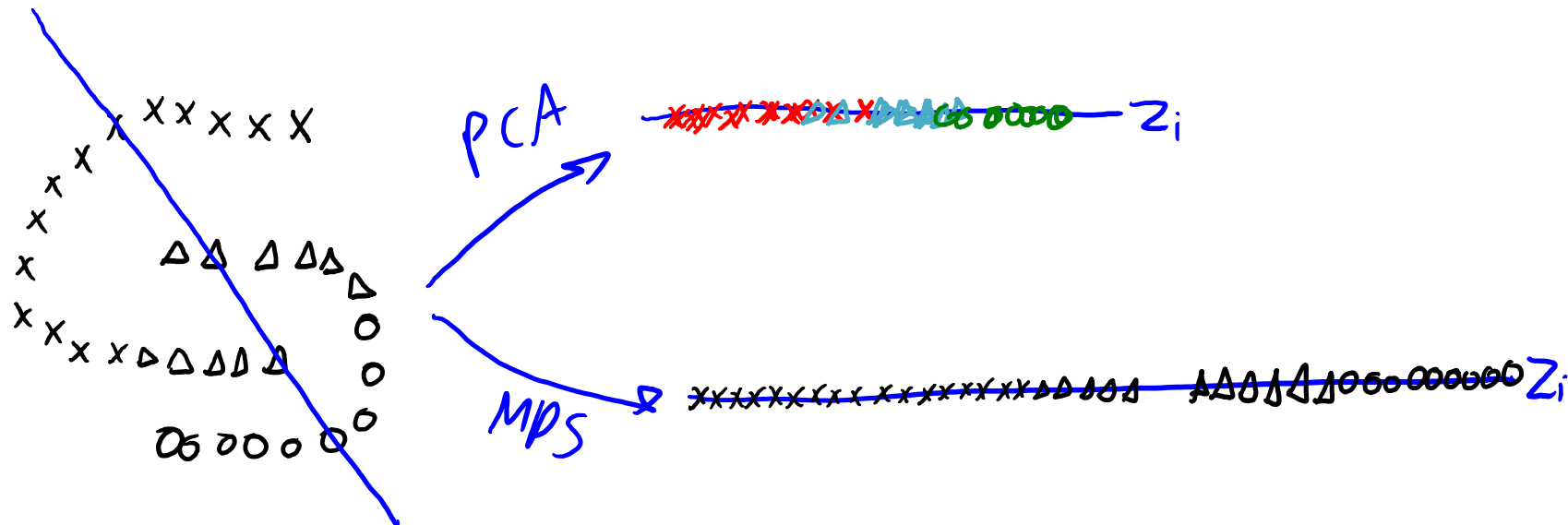


Multi-Dimensional Scaling

- Multi-dimensional scaling (MDS):
 - Optimize the final locations of the z_i values.

$$f(z) = \sum_{i=1}^n \sum_{j=i+1}^n (\|z_i - z_j\| - \|x_i - x_j\|)^2$$

- Non-parametric dimensionality reduction and visualization:
 - No 'W': just trying to make z_i preserve high-dimensional distances between x_i .



Multi-Dimensional Scaling

- Multi-dimensional scaling (MDS):
 - Optimize the final locations of the z_i values.

$$f(z) = \sum_{i=1}^n \sum_{j=i+1}^n (\|z_i - z_j\| - \|x_i - x_j\|)^2$$

- Cannot use SVD to compute solution:
 - Instead, do gradient descent on the z_i values.
 - You “learn” a scatterplot that tries to visualize high-dimensional data.
 - Not convex and sensitive to initialization.
 - And solution is not unique due to various factors like translation and rotation.

In This Lecture

1. Multi-Dimensional Scaling

- Euclidean MDS
- Sammon Mapping
- Geodesic MDS (ISOMAP)

2. Latent Factors for Language

Coming Up Next

EUCLIDEAN MDS VARIANTS

Different MDS Cost Functions

- **MDS** default objective: squared difference of Euclidean norms:

$$f(z) = \sum_{i=1}^n \sum_{j=i+1}^n (\|z_i - z_j\| - \|x_i - x_j\|)^2$$

Q: How many distance functions are involved here?

Q: Can we generalize this to other measures of distance?

Different MDS Cost Functions

- **MDS** default objective function with **general distances/similarities**:

$$f(Z) = \sum_{i=1}^n \sum_{j=i+1}^n d_3(d_2(z_i, z_j) - d_1(x_i, x_j))$$

– Functions are **not necessarily the same**:

- d_1 := high-dimensional distance we want to match.

$$d_1 : \mathbb{R}^d \times \mathbb{R}^d \rightarrow \mathbb{R}$$

- d_2 := low-dimensional distance we can control.

$$d_2 : \mathbb{R}^2 \times \mathbb{R}^2 \rightarrow \mathbb{R}$$

- d_3 := how we compare high-/low-dimensional distances.

$$d_3 : \mathbb{R} \times \mathbb{R} \rightarrow \mathbb{R}$$

Different MDS Cost Functions

- **MDS** default objective function with **general distances/similarities**:

$$f(Z) = \sum_{i=1}^n \sum_{j=i+1}^n d_3(d_2(z_i, z_j) - d_1(x_i, x_j))$$

- **“Classic” MDS:**

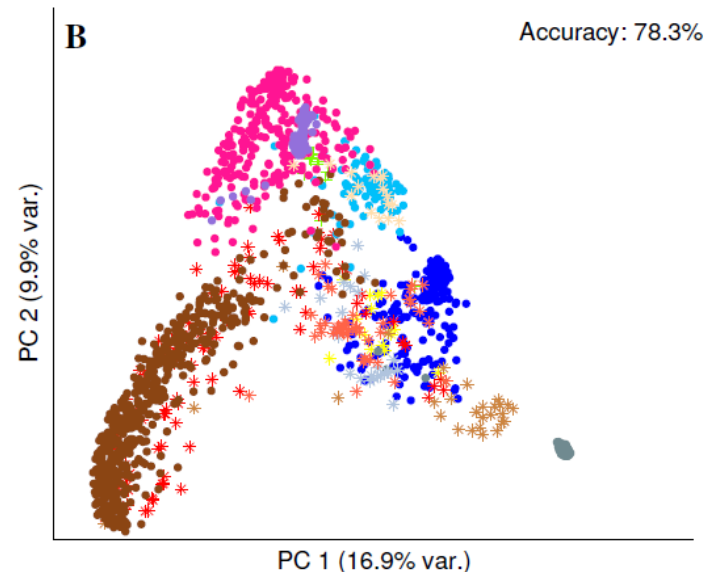
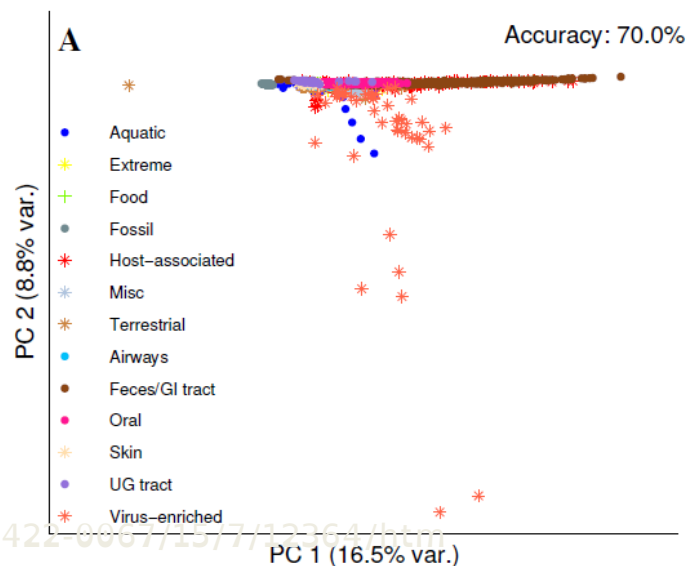
- $d_1(x_i, x_j) = x_i^T x_j$, $d_2(z_i, z_j) = z_i^T z_j$, $d_3(a, b) = (a - b)^2$
- This is a factorless version of _____.
- Not a great choice because it's _____.

Different MDS Cost Functions

- **MDS** default objective function with **general distances/similarities**:

$$f(z) = \sum_{i=1}^n \sum_{j=i+1}^n d_3(d_2(z_i, z_j) - d_1(x_i, x_j))$$

- Another possibility: $d_1(x_i, x_j) = \|x_i - x_j\|_1$ and $d_2(z_i, z_j) = \|z_i - z_j\|$.
 - z_i approximates high-dimensional L_1 -norm distances.



Sammon's Mapping

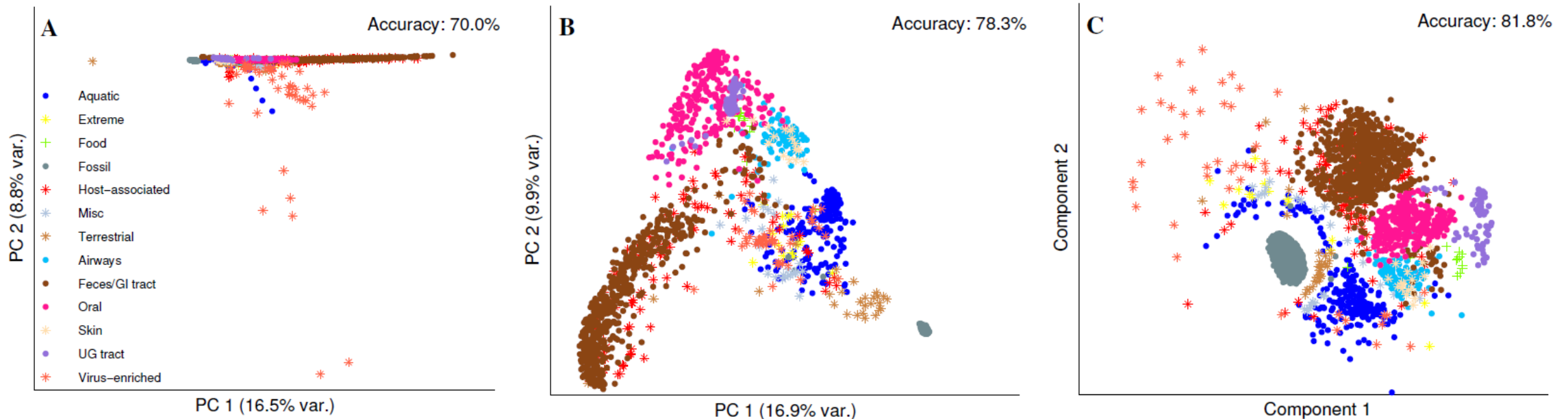
- Challenge for most MDS models: they **focus on** _____.
 - Leads to “crowding” effect like with PCA.
- Early attempt to address this is **Sammon's mapping**:
 - **Weighted MDS** so large/small distances are more comparable.

$$f(Z) = \sum_{i=1}^n \sum_{j=i+1}^n \left(\frac{d_2(z_i, z_j) - d_1(x_i, x_j)}{d_1(x_i, x_j)} \right)^2$$

- Denominator **reduces focus on large distances**.

Sammon's Mapping

- Challenge for most MDS models: they **focus on large distances**.
 - Leads to “crowding” effect like with PCA.
- Early attempt to address this is **Sammon's mapping**:
 - **Weighted MDS** so large/small distances are more comparable.

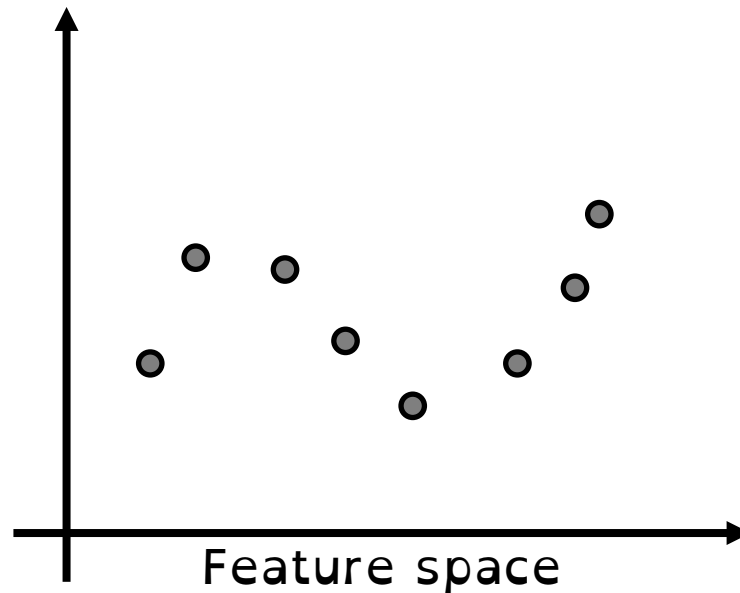


Coming Up Next

MANIFOLDS

“Manifold”

- “Manifold” := non-Euclidean subspace of feature space where datapoints live

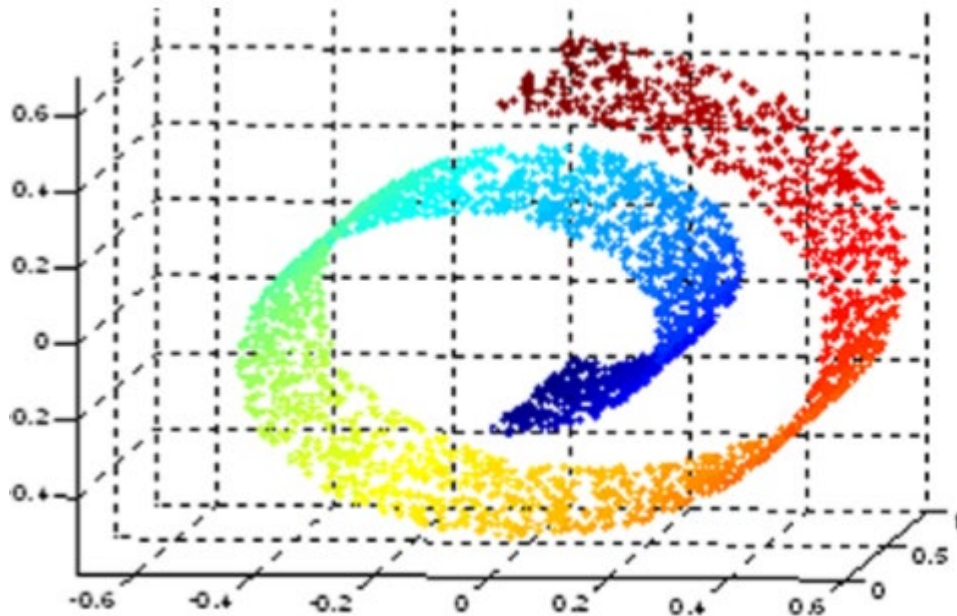


- Assumption: most data live on a manifold, not a true Euclidean feature space!

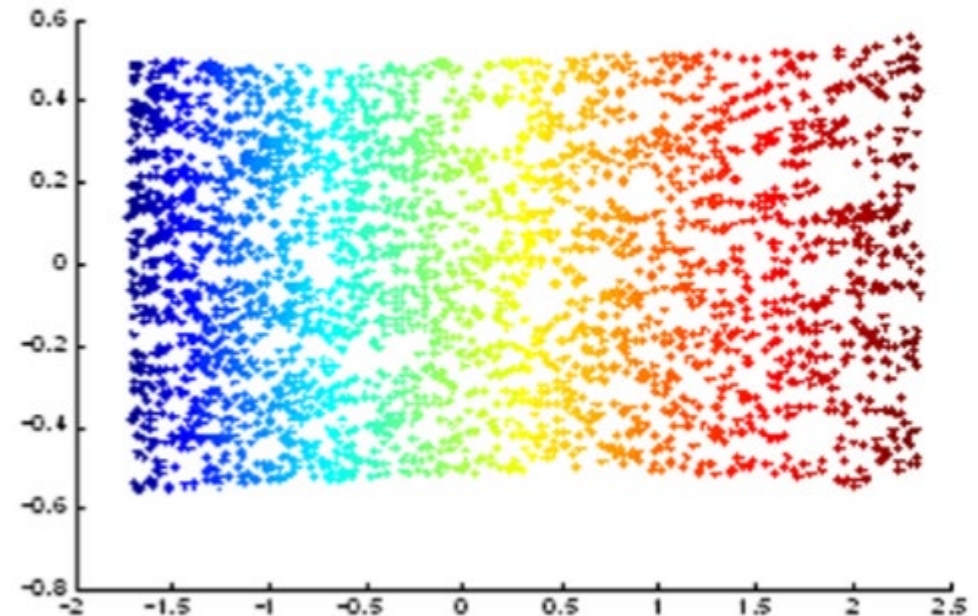
Learning Manifolds

- Consider data that lives on a **low-dimensional “manifold”**.
- e.g. ‘Swiss roll’:

Original data

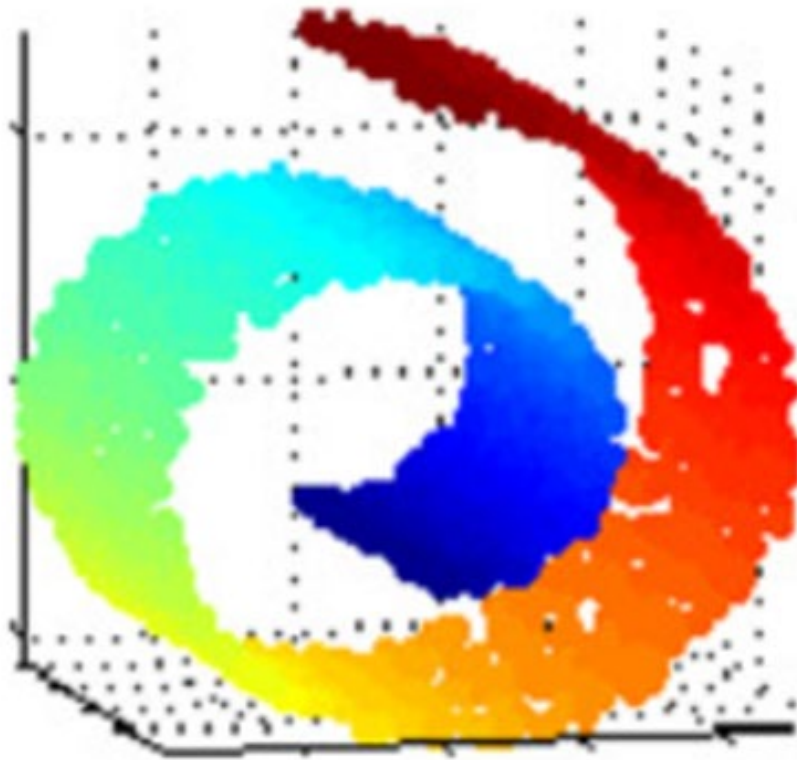


Two-dimensional manifold

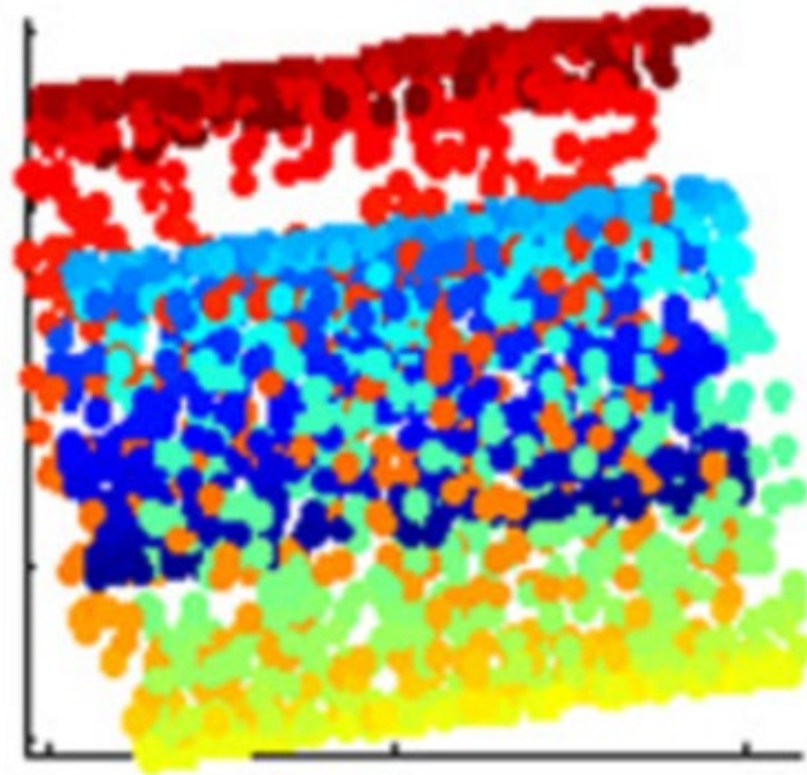


Learning Manifolds

- Consider data that lives on a **low-dimensional “manifold”**.
 - With usual distances, **PCA/MDS will not discover non-linear manifolds.**



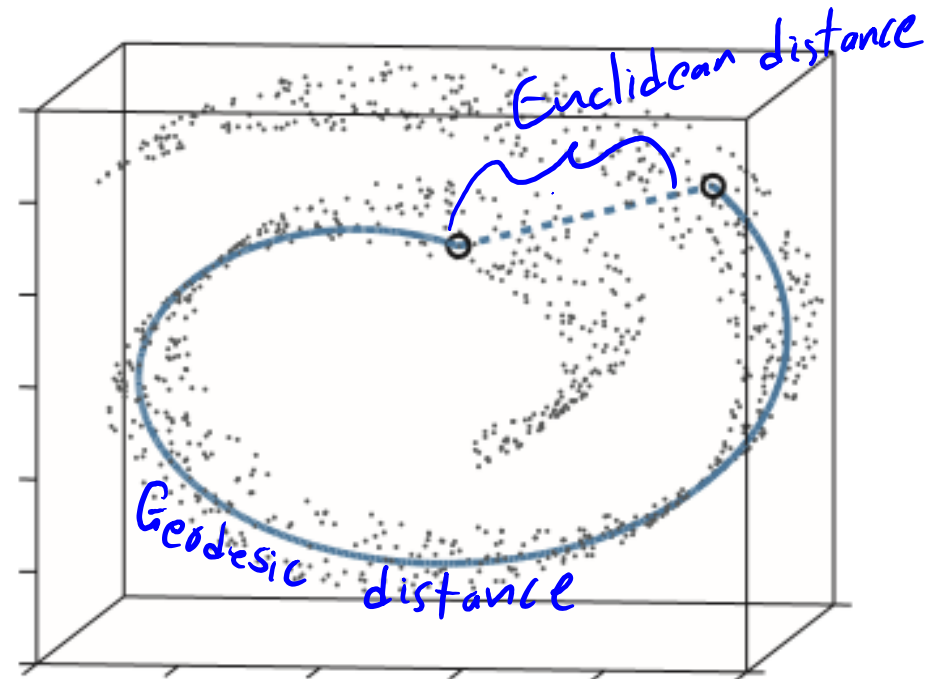
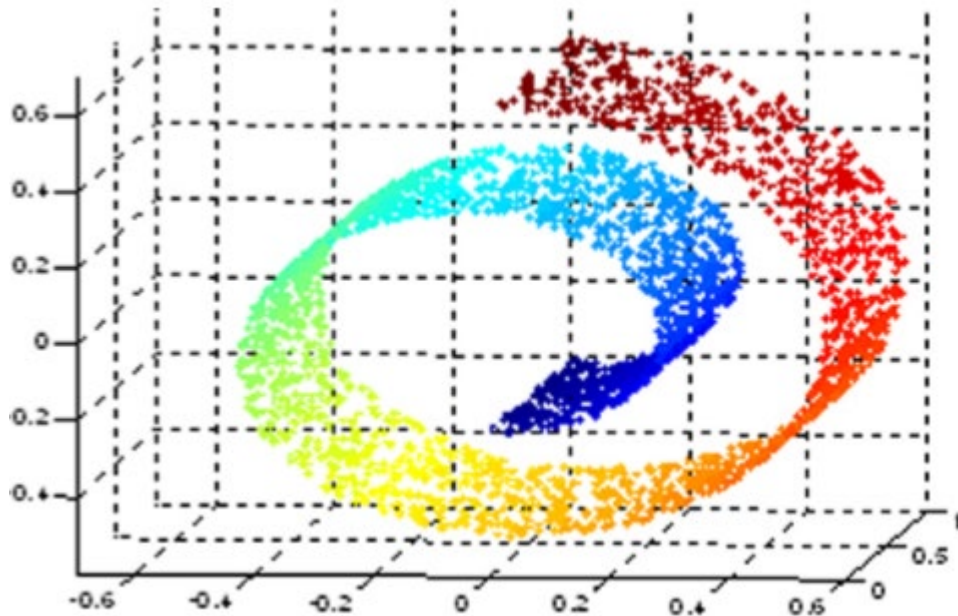
Original data



PCA

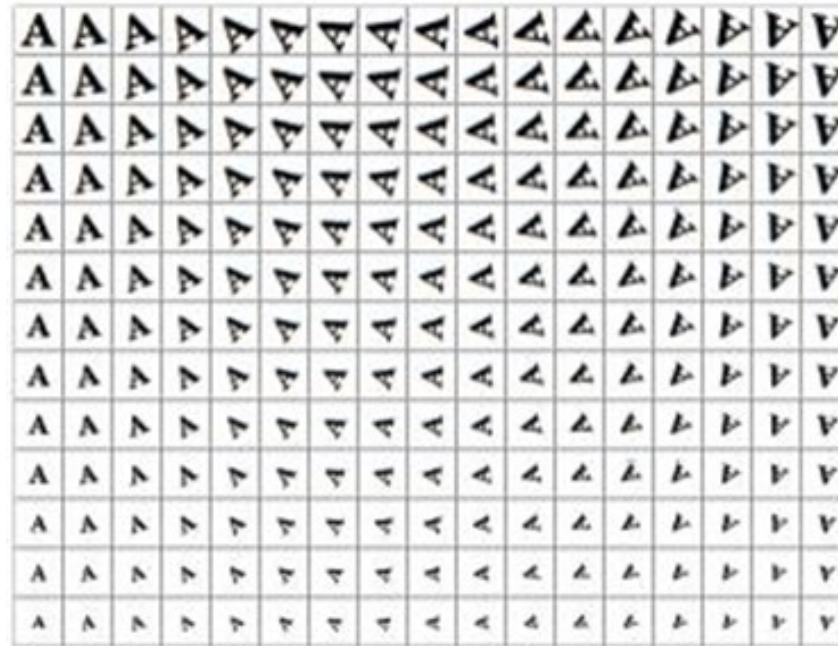
Learning Manifolds

- Consider data that lives on a **low-dimensional “manifold”**.
 - With usual distances, **PCA/MDS will not discover non-linear manifolds**.
- We need **geodesic distance**: the _____.



Manifolds in Image Space

- Consider slowly-varying transformation of image:



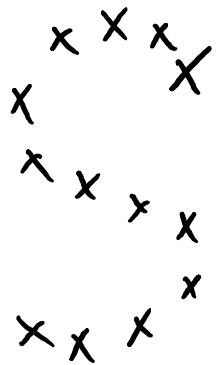
- **Images are on a manifold** in the high-dimensional space.
 - Euclidean distance **doesn't reflect manifold structure**.
 - **Geodesic distance** is **distance through space of rotations/resizings**.

Coming Up Next

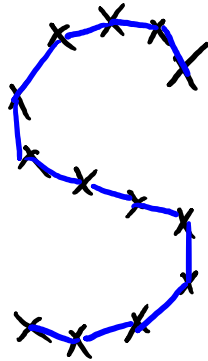
ISOMAP

ISOMAP

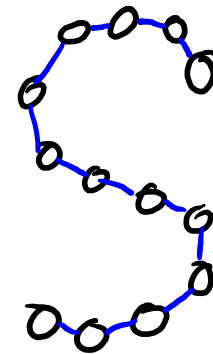
- ISOMAP is MDS on manifolds:



find "neighbours"
of each point



Represent points
and neighbours
as a weighted
graph.



"weight" on each
edge is distance
between points

Approximate geodesic distance
by shortest path through
graph.

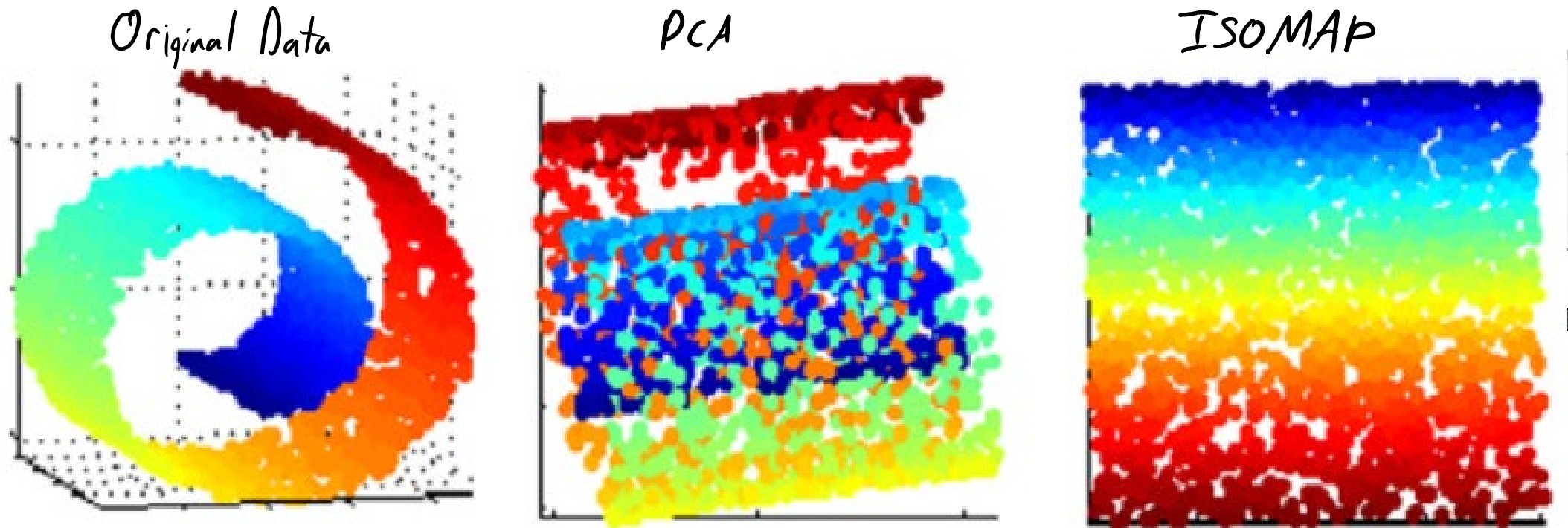
$$D = \begin{bmatrix} 0 & 1 & 2 & 3 & \dots \\ 1 & 0 & 1 & 2 & \dots \\ 2 & 1 & 0 & 1 & \dots \\ 3 & 2 & 1 & 0 & \dots \\ \vdots & \vdots & \vdots & \vdots & \ddots \end{bmatrix}$$

ISOMAP z_i values in 1D or 2D

Run MDS
with these
approximate geodesic distances.

ISOMAP

- ISOMAP can “unwrap” the roll:
 - _____ are approximations to geodesic distances.



- Sensitive to having _____ :
 - Points off of manifold and gaps in manifold cause problems.

Constructing Neighbour Graphs

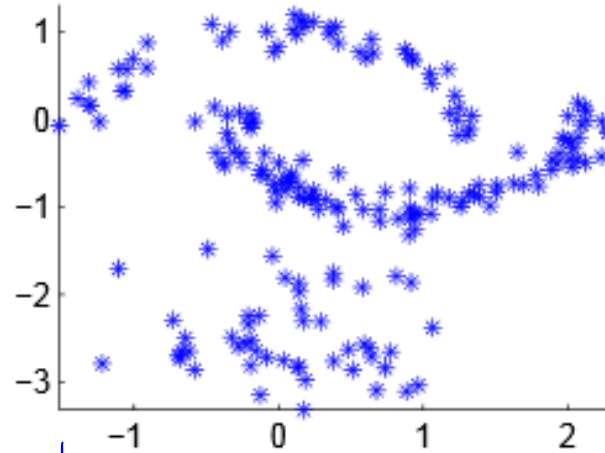
- Sometimes you can **define the graph/distance without features**:
 - Facebook friend graph.
 - Connect YouTube videos if one video tends to follow another.
- But we can also **convert from features x_i to a “neighbour” graph (A6)**:
 - Approach 1 (“**epsilon graph**”): connect x_i to all x_j within some threshold ϵ .
 - Like we did with density-based clustering.
 - Approach 2a (“**KNN graph**”): connect x_i to x_j if:
 - x_j is a KNN of x_i **OR** x_i is a KNN of x_j .
 - Approach 2b (“**mutual KNN graph**”): connect x_i to x_j if:
 - x_j is a KNN of x_i **AND** x_i is a KNN of x_j .

Converting from Features to Graph

add edge
if $\|x_i - x_j\| \leq 0.3$

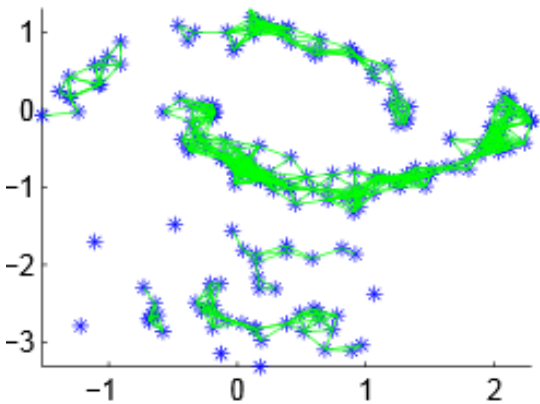
add edge if
 i is 5-NN
of j or
 j is
5-NN
of i

Data points

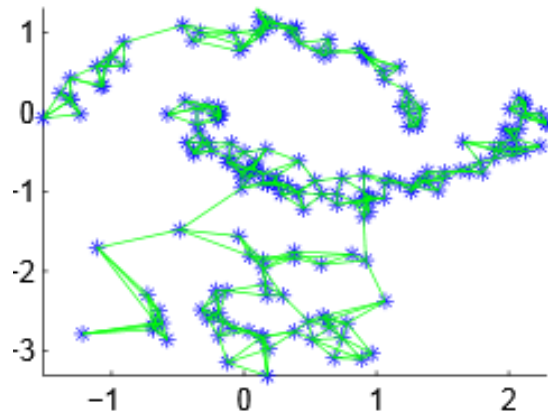


add edge if
 i and j
are kNNs
of each other.

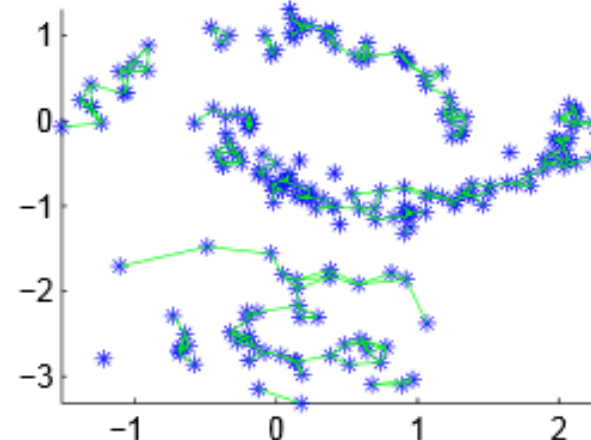
epsilon-graph, epsilon=0.3



kNN graph, k = 5

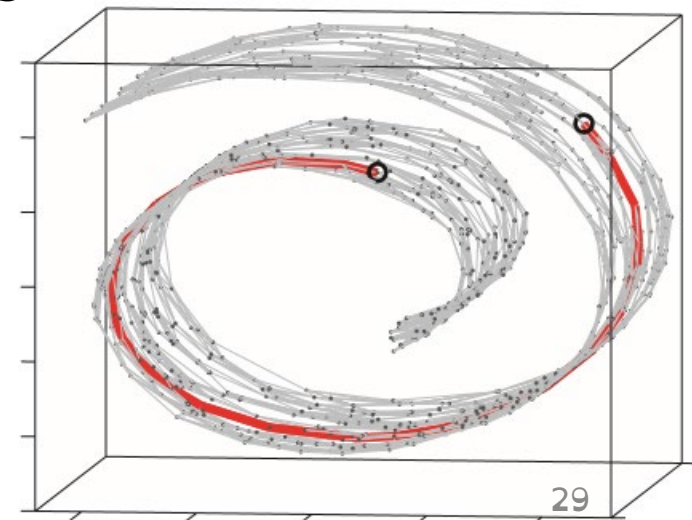


Mutual kNN graph, k = 5

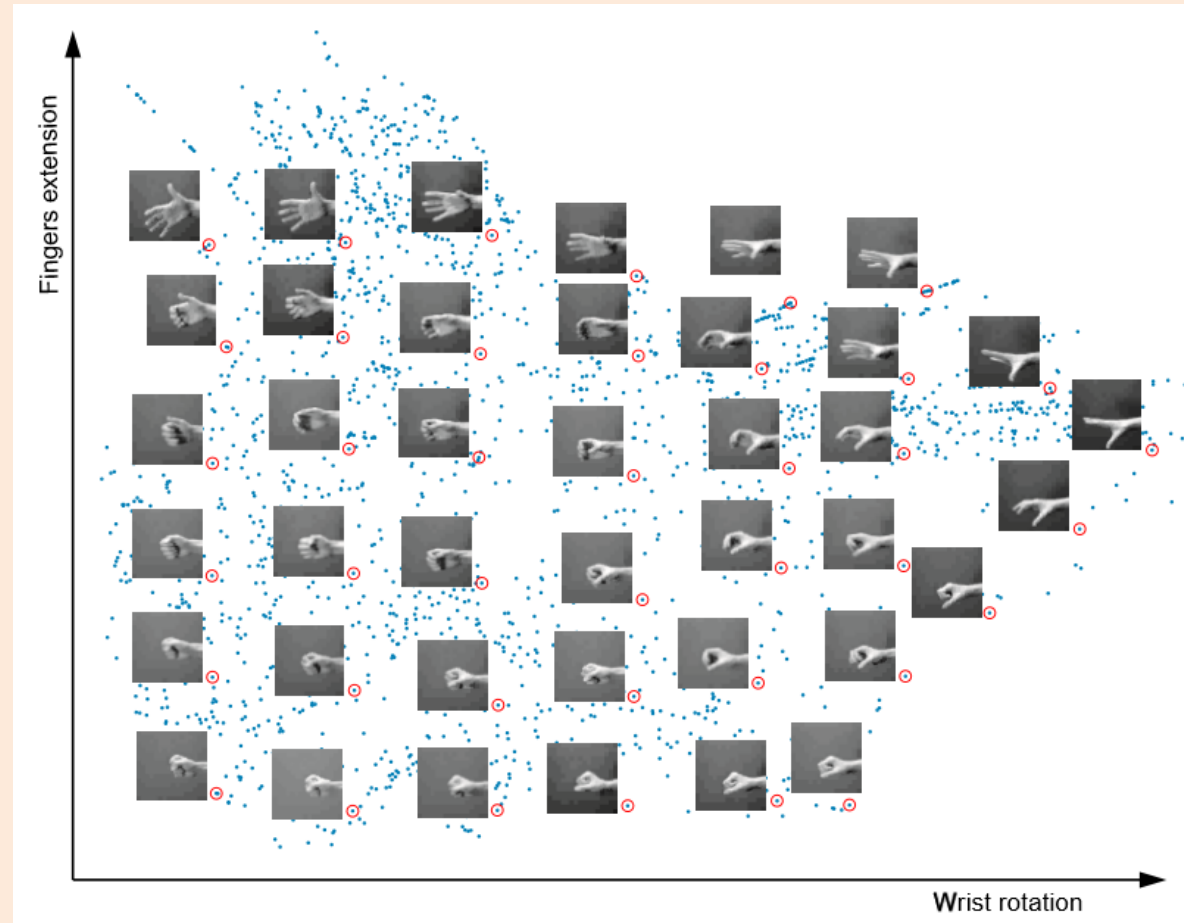


ISOMAP

- **ISOMAP** is latent-factor model for visualizing data on manifolds:
 1. Find the **neighbours** of each point.
 - Usually “k-nearest neighbours graph”, or “epsilon graph”.
 2. Compute **edge weights**:
 - Usually distance between neighbours.
 3. Compute **weighted shortest path** between all points
 - Dijkstra or other shortest path algorithm.
 4. Run **MDS** using these distances.



ISOMAP on Hand Images



- Related method is “local linear embedding”.

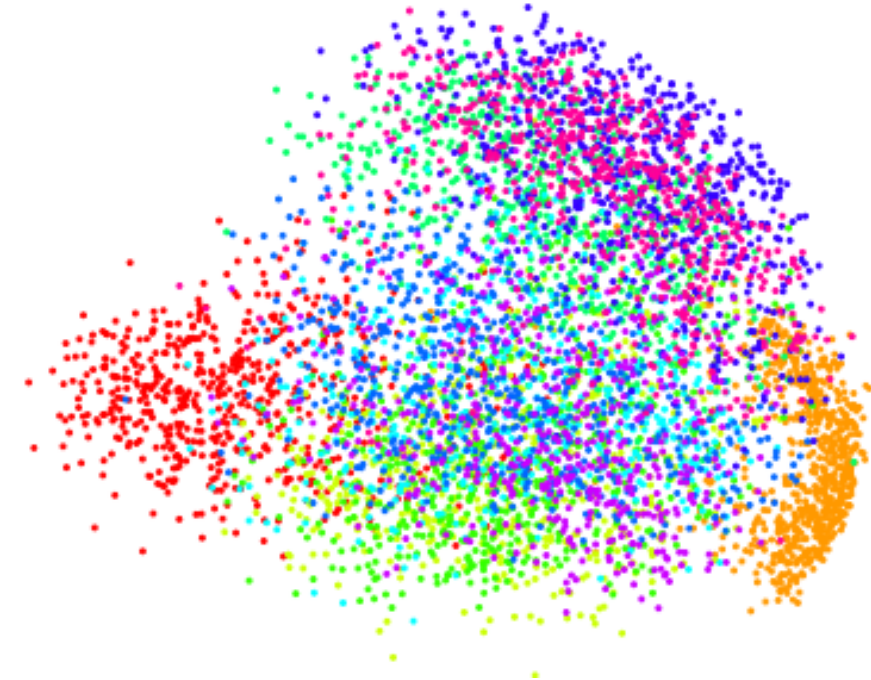
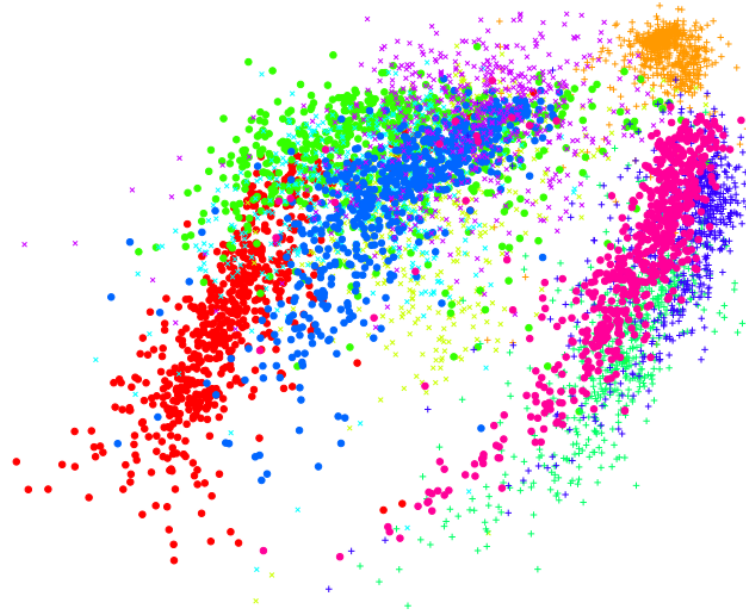
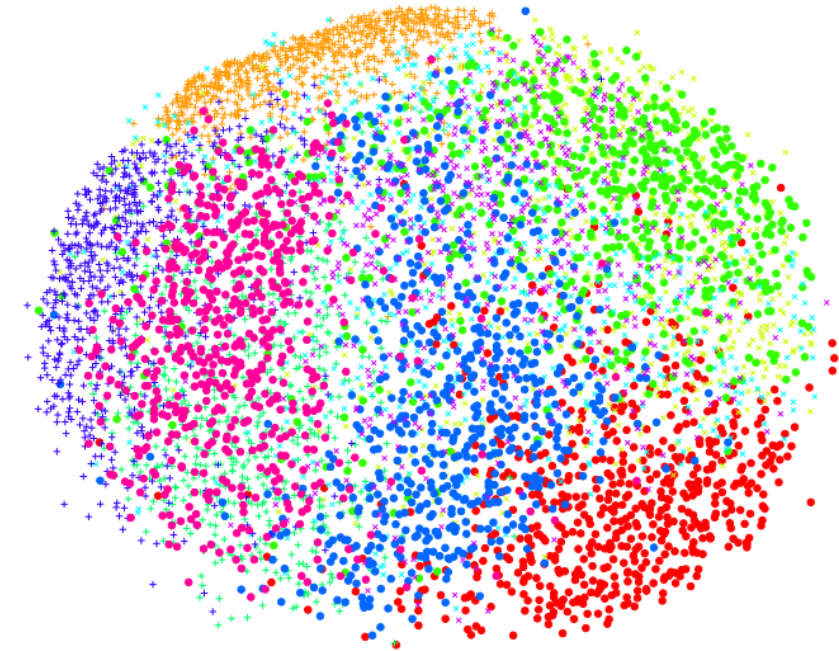
Sammon's Map vs. ISOMAP vs. PCA



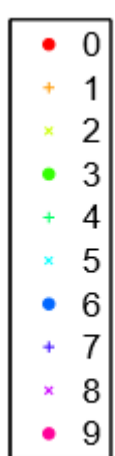
Sammon Map

ISOMAP

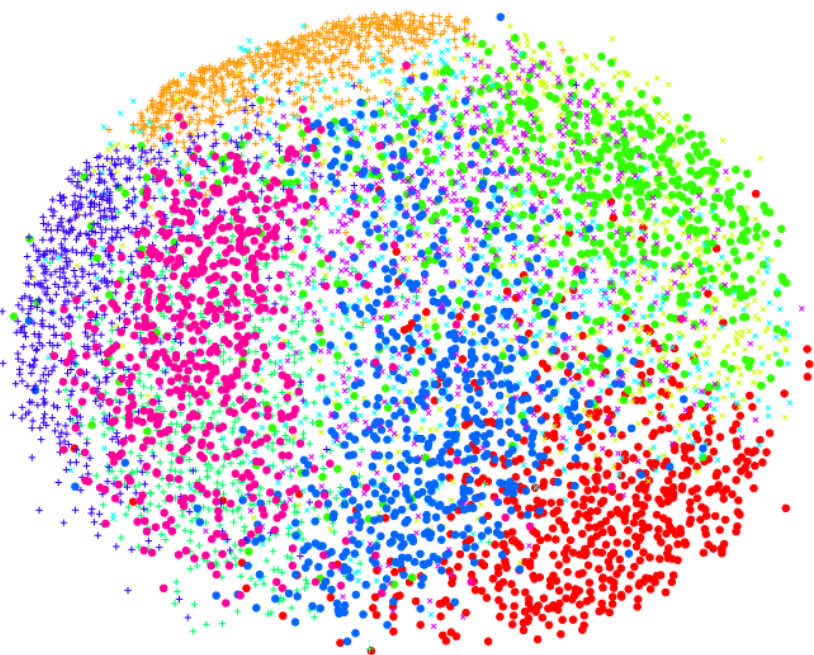
PCA



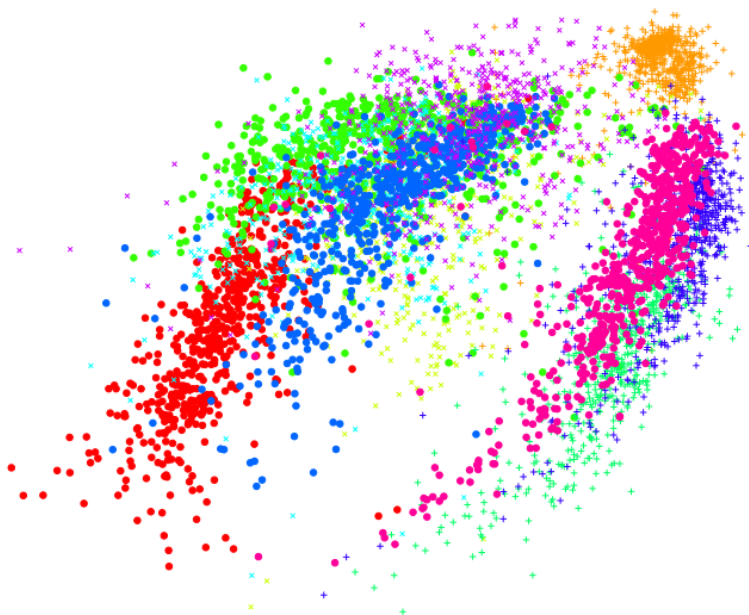
Sammon's Map vs. ISOMAP vs. t-SNE



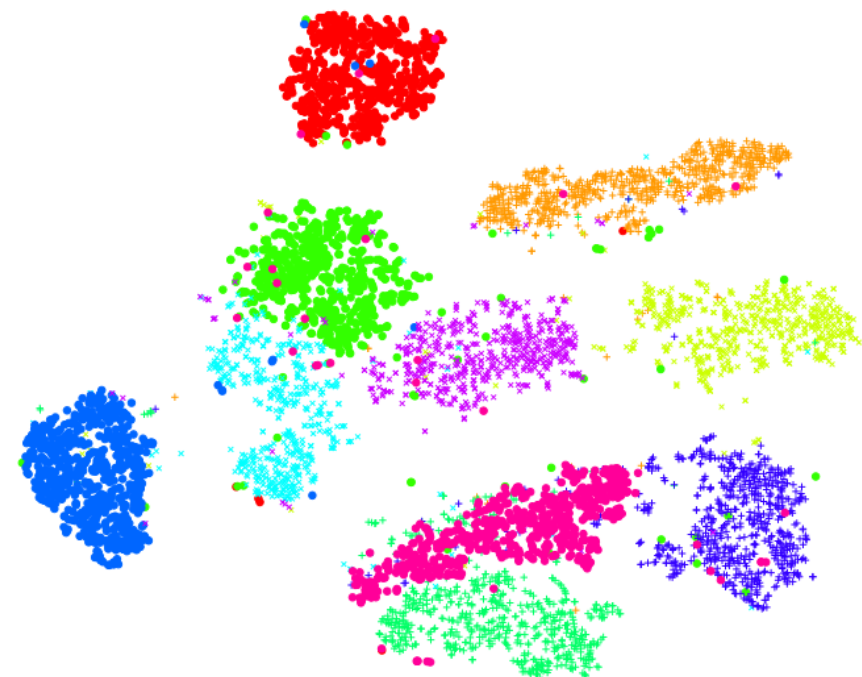
Sammon Map



ISOMAP



t-SNE



Sammon's Map vs. ISOMAP vs. t-SNE



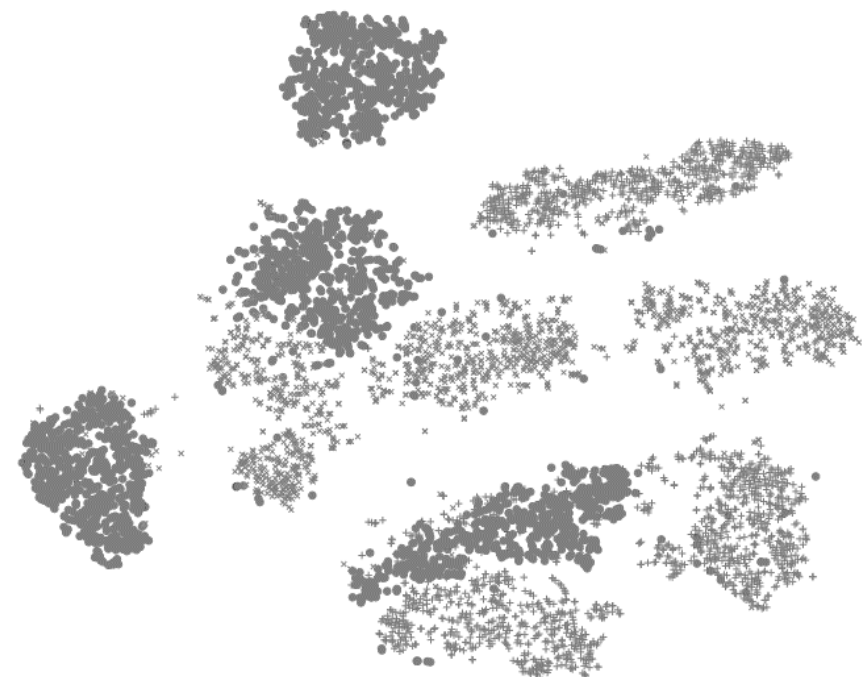
Sammon Map



ISOMAP



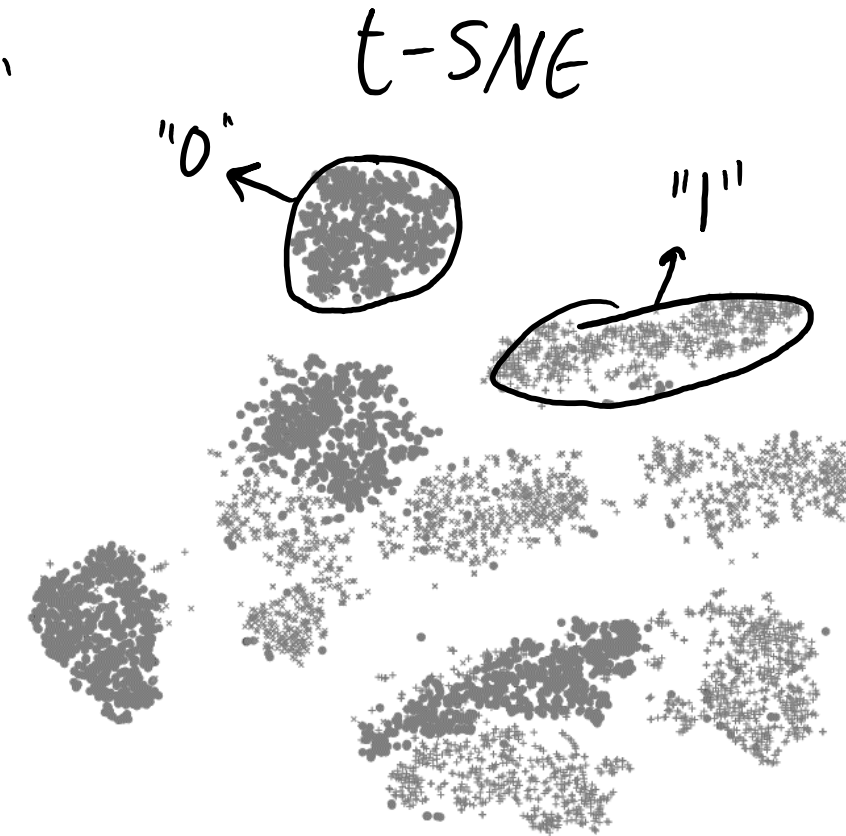
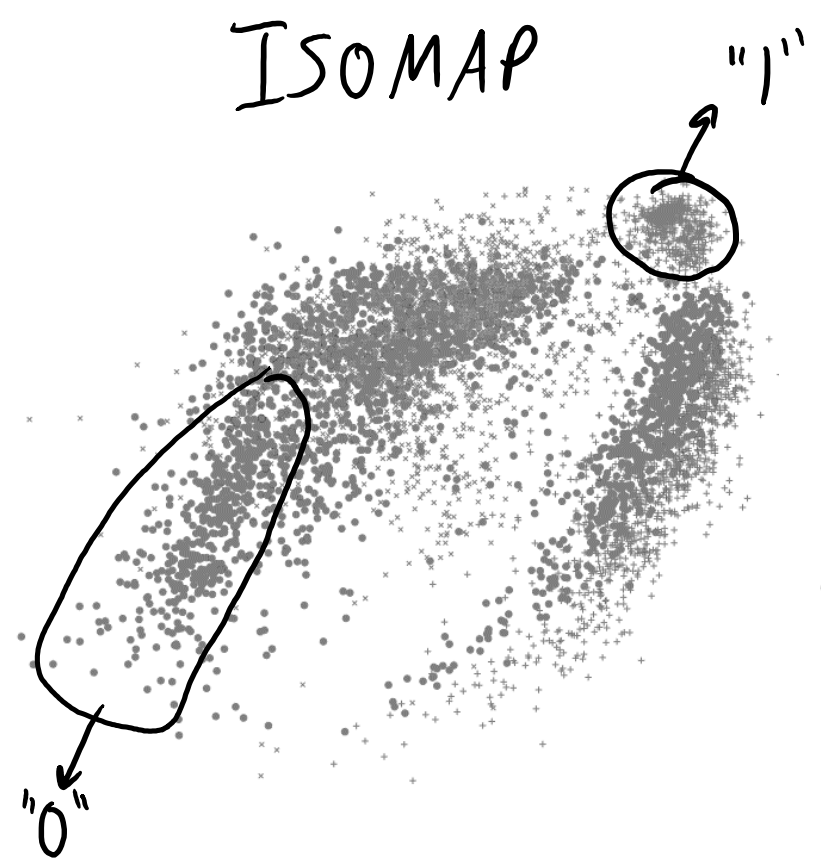
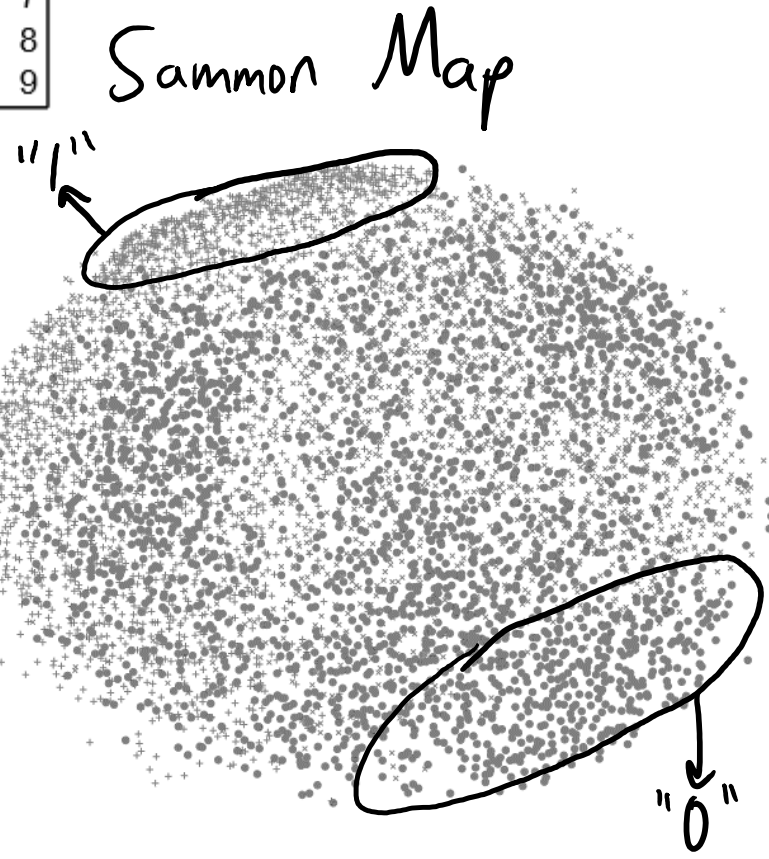
t-SNE



Remember this is unsupervised, algorithms do not know the labels.

Sammon's Map vs. ISOMAP vs. t-SNE

- 0
- 1
- 2
- 3
- 4
- 5
- 6
- 7
- 8
- 9

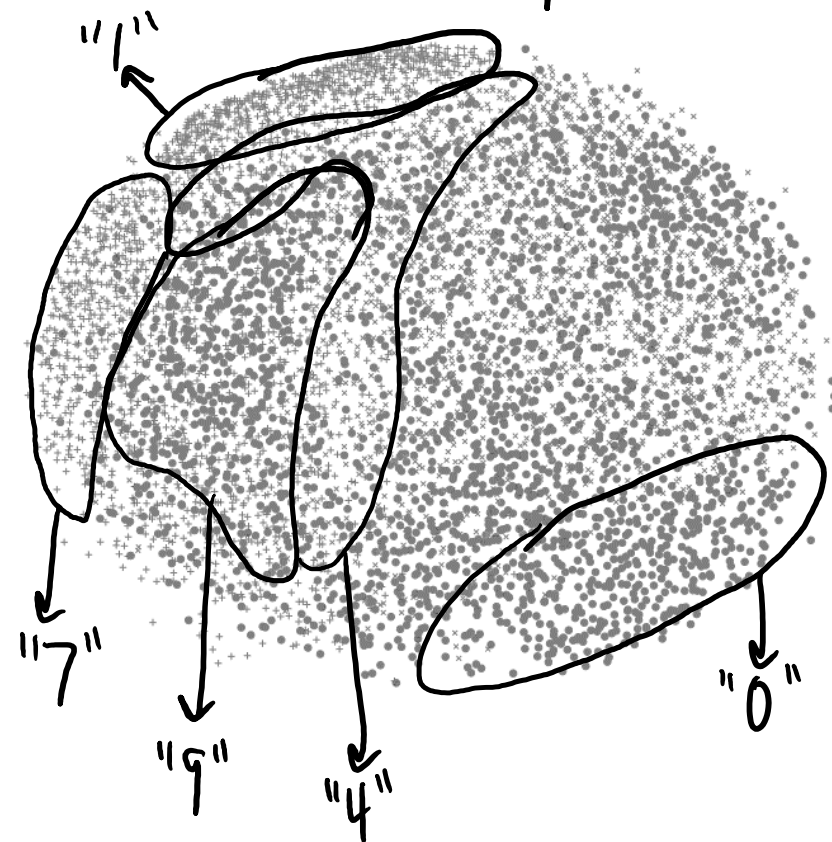


Remember this is unsupervised, algorithms do not know the labels.

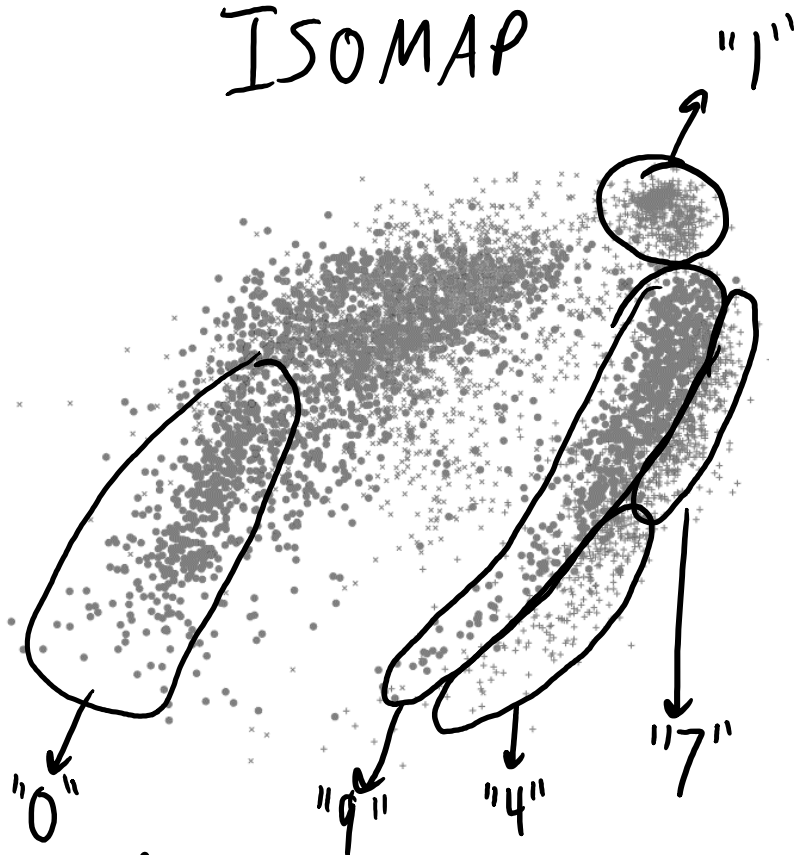
Sammon's Map vs. ISOMAP vs. t-SNE

- 0
- 1
- 2
- 3
- 4
- 5
- 6
- 7
- 8
- 9

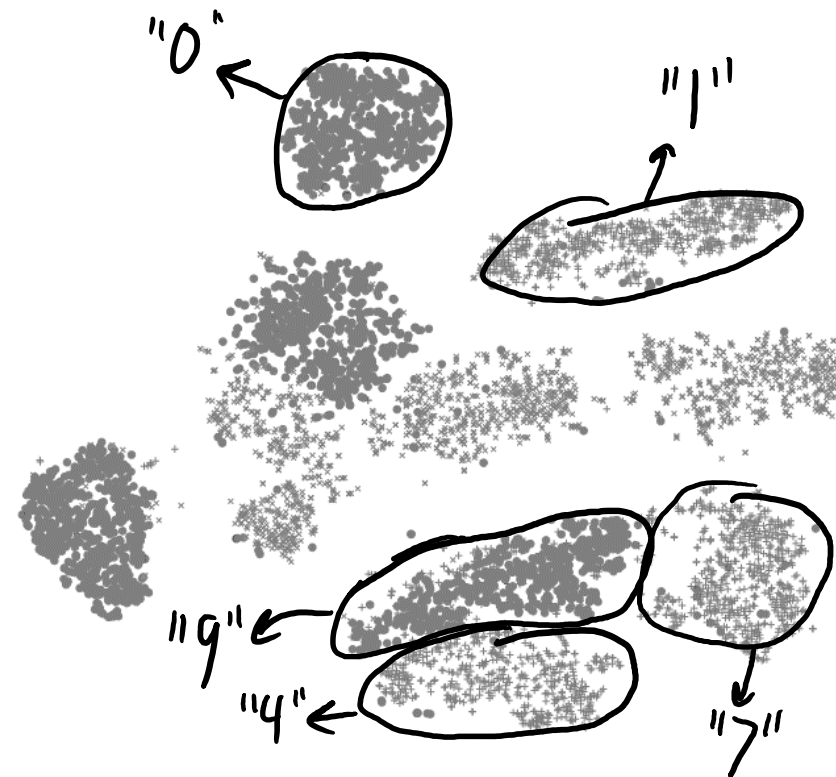
Sammon Map



ISOMAP



t-SNE

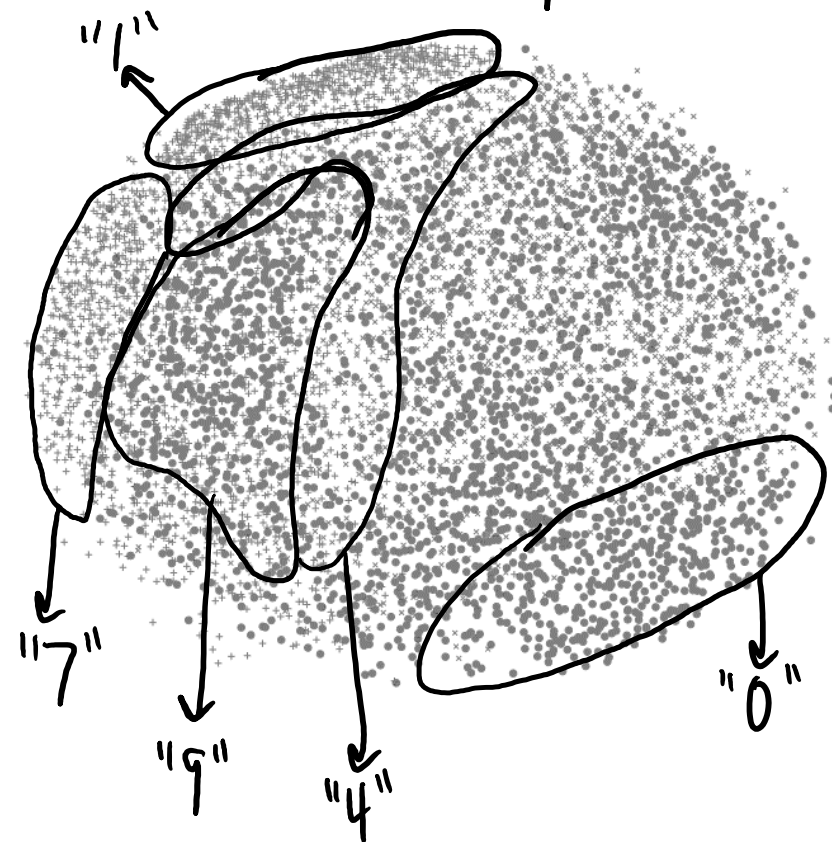


Remember this is unsupervised, algorithms do not know the labels.

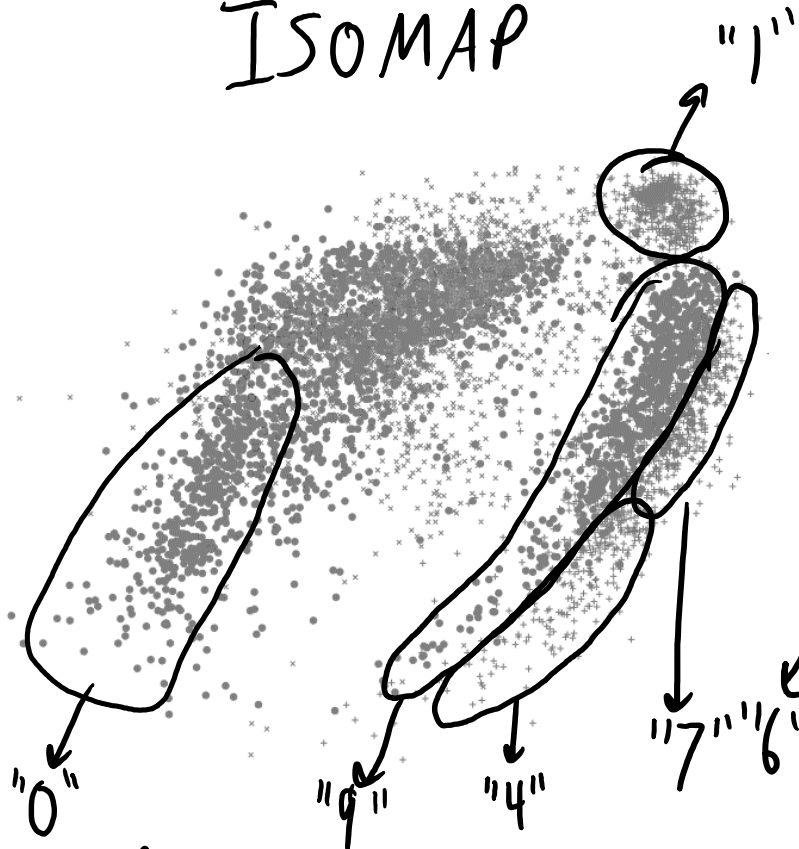
Sammon's Map vs. ISOMAP vs. t-SNE

- 0 ●
- 1 +
- 2 ×
- 3 ●
- 4 +
- 5 ×
- 6 ●
- 7 +
- 8 ×
- 9 ●

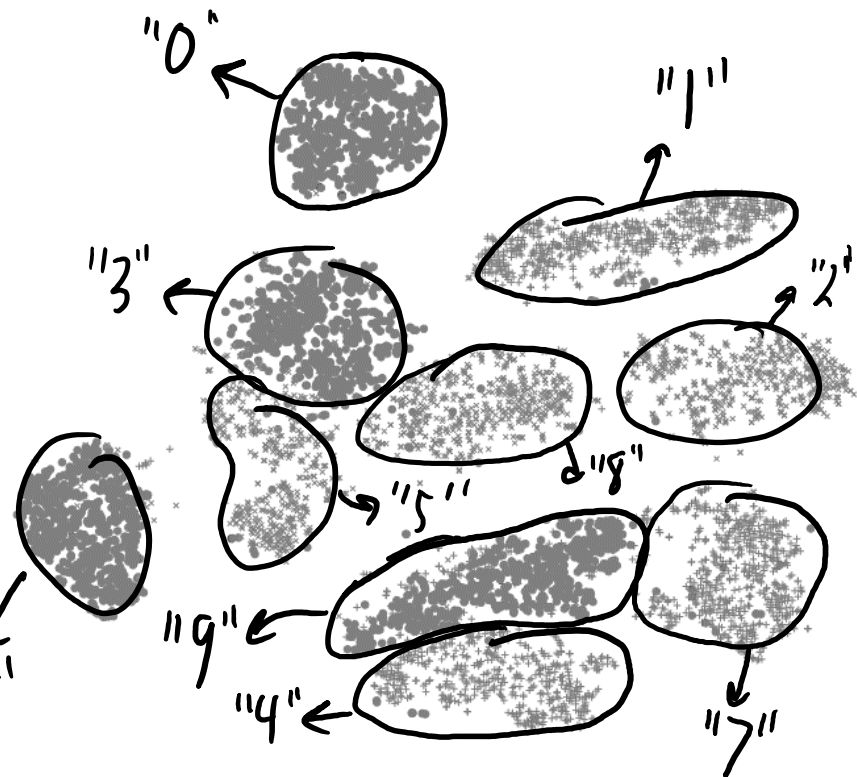
Sammon Map



ISOMAP



t-SNE



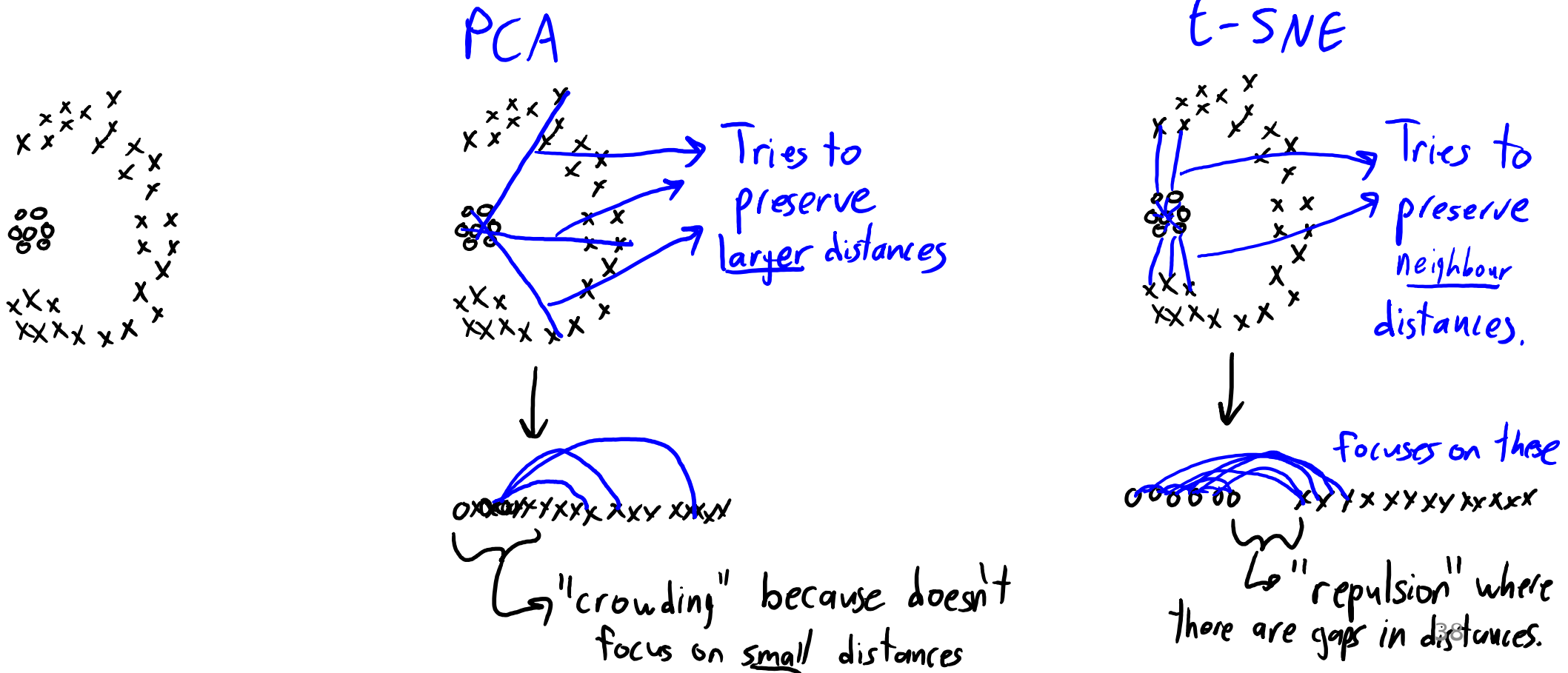
Remember this is unsupervised, algorithms do not know the labels.

Coming Up Next

T-SNE

t-Distributed Stochastic Neighbour Embedding

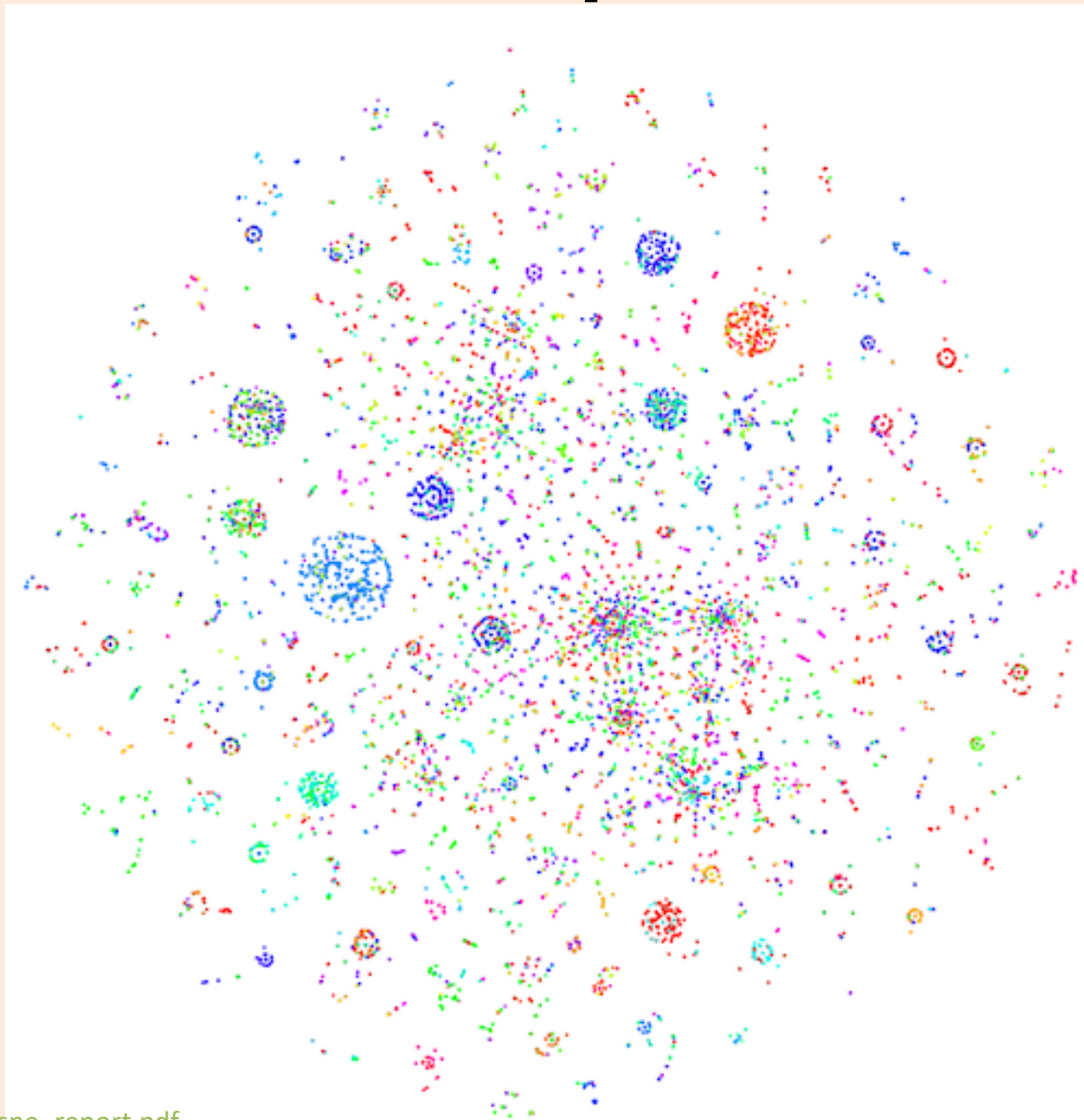
- One key idea in **t-SNE**:
 - Focus on distance to “neighbours”
(allow large variance in other distances)



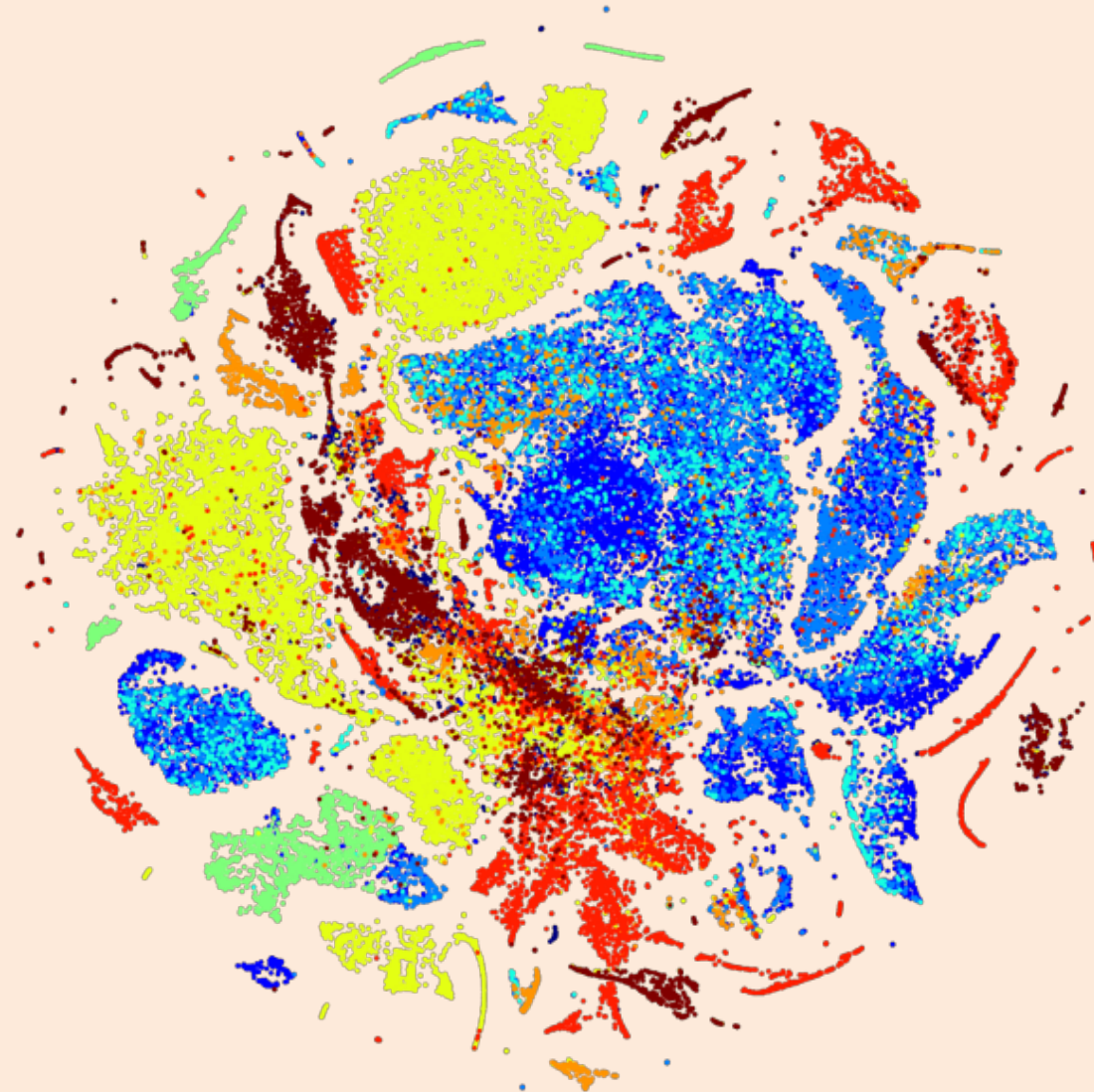
t-Distributed Stochastic Neighbour Embedding

- **t-SNE** is a special case of MDS (specific d_1 , d_2 , and d_3 choices):
 - d_1 : for each x_i , compute 'neighbour-ness' of each x_j
 - Computation is similar to k-means++, but most weight to close points (Gaussian).
 - Doesn't require explicit graph.
 - d_2 : for each z_i , compute 'neighbour-ness' of each z_j .
 - Similar to above, but use student's t (grows really slowly with distance).
 - Avoids 'crowding', because you have a huge range that large distances can fill.
 - d_3 : Compare x_i and z_i using an entropy-like measure:
 - How much 'randomness' is in probabilities of x_i if you know the z_i (and vice versa)?
- Interactive demo: <https://distill.pub/2016/misread-tsne>

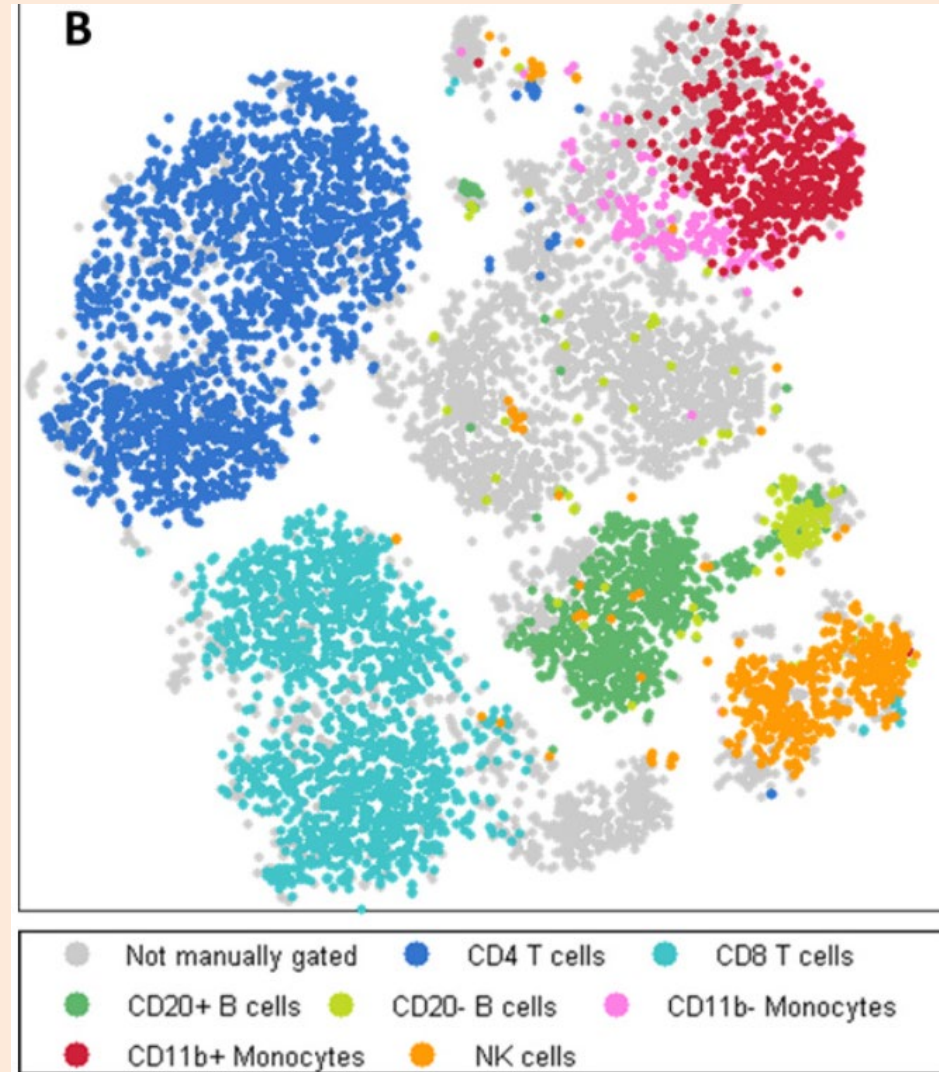
t-SNE on Wikipedia Articles



t-SNE on Product Features



t-SNE on Leukemia Heterogeneity



End of Part 4: Latent Factor Models

End of Part 4: Key Concepts

- We discussed **linear latent-factor models**:

$$\begin{aligned} f(W, Z) &= \sum_{i=1}^n \sum_{j=1}^d (\langle w_j, z_i \rangle - x_{ij})^2 \\ &= \sum_{i=1}^n \|W^T z_i - x_i\|^2 \\ &= \|ZW - X\|_F^2 \end{aligned}$$

- Represent 'X' as linear combination of **latent factors 'w_c'**.
 - **Latent features 'z_i'** give a lower-dimensional version of each 'x_i'.
 - When k=1, finds **direction that minimizes squared orthogonal distance**.
- Applications:
 - Outlier detection, dimensionality reduction, data compression, features for linear models, visualization, factor discovery, filling in missing entries.

End of Part 4: Key Concepts

- We discussed **linear latent-factor models**:

$$f(W, z) = \sum_{i=1}^n \sum_{j=1}^d (\langle w_j, z_i \rangle - x_{ij})^2$$

- **Principal component analysis (PCA)**:
 - Often uses **orthogonal factors** and fits them **sequentially** (via **SVD**).
- **Non-negative matrix factorization**:
 - Uses **non-negative** factors giving sparsity.
 - Can be minimized with **projected gradient**.
- Many variations are possible:
 - Different regularizers (**sparse coding**) or loss functions (**robust/binary PCA**).
 - Missing values (**recommender systems**) or change of basis (**kernel PCA**).

End of Part 4: Key Concepts

- We discussed **multi-dimensional scaling (MDS)**:
 - **Non-parametric** method for high-dimensional **data visualization**.
 - Tries to match distance/similarity in high-/low-dimensions.
 - “Gradient descent on scatterplot points”.
- Main **challenge in MDS methods is “crowding”** effect:
 - Methods focus on large distances and lose local structure.
- Common solutions:
 - **Sammon mapping**: use weighted cost function.
 - **ISOMAP**: approximate geodesic distance using via shortest paths in graph.
 - **T-SNE**: give up on large distances and focus on neighbour distances.

Summary

- Different MDS distances/losses/weights usually gives better results.
- Manifold learning focuses on low-dimensional curved structures.
- ISOMAP is most common approach:
 - Approximates geodesic distance by shortest path in weighted graph.
- t-SNE is promising new data MDS method.

- Next time: deep learning.

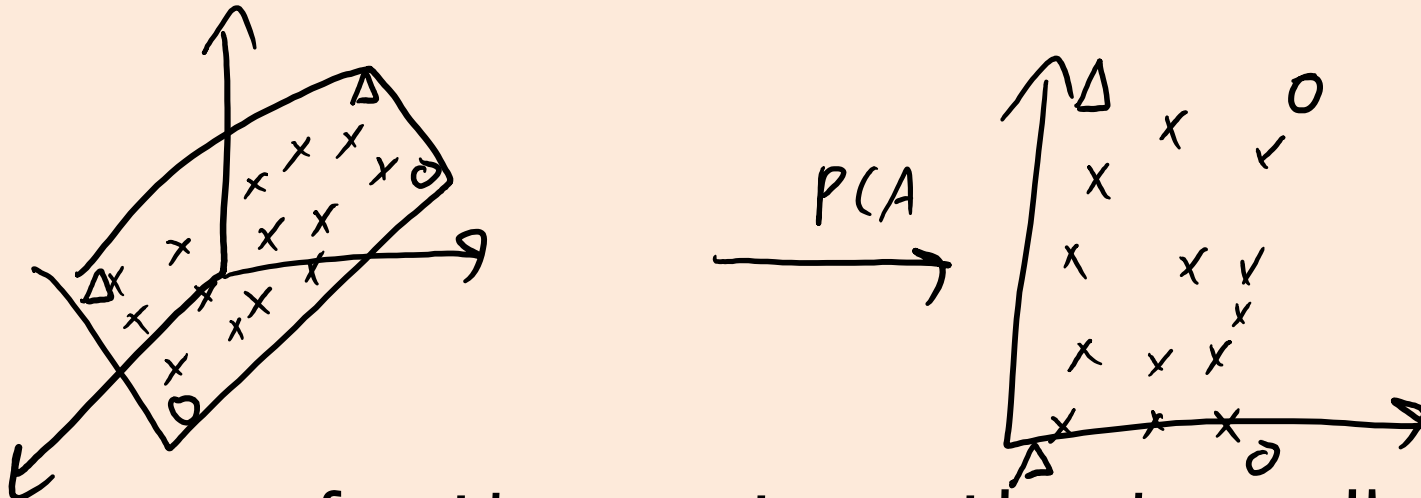
Please Do Course Evaluation!

Review Questions

- Q1: Is MDS sensitive to initialization? Why?
- Q2: What is the problem with using linear dimensionality reduction for data on manifold?
- Q3: How does ISOMAP compute pair-wise distances among examples?
- Q4: What is the key idea behind t-SNE in terms of preserving distances in 2D?

Does t-SNE always outperform PCA?

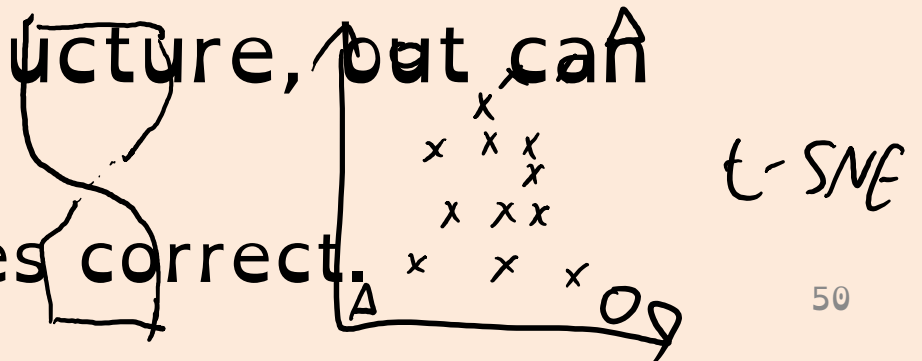
- Consider 3D data living on a 2D hyper-plane:



- PCA can perfectly capture the low-dimensional structure.

- T-SNE can capture the local structure, but can “twist” the plane.

– It doesn't try to get long distances correct.



Graph Drawing

- A closely-related topic to MDS is **graph drawing**:
 - Given a graph, how should we display it?
 - Lots of interesting methods: https://en.wikipedia.org/wiki/Graph_drawing



Bonus Slide: Multivariate Chain Rule

- Recall the **univariate chain rule**: $\frac{d}{dw} [f(g(w))] = f'(g(w)) g'(w)$
- The **multivariate chain rule**: $\underbrace{\nabla [f(g(w))]}_{1 \times 1} = \underbrace{f'(g(w))}_{1 \times 1} \underbrace{\nabla g(w)}_{d \times 1}$
- Example:

$$\nabla \left[\frac{1}{2} (w^T x_i - y_i)^2 \right]$$

$$= \nabla [f(g(w))]$$

with $g(w) = w^T x_i - y_i$

and $f(r_i) = \frac{1}{2} r_i^2$

$$\nabla g(w) = x_i$$

$$f'(r_i) = r_i$$

$$\nabla [f(g(w))] = r_i x_i$$

$$= (w^T x_i - y_i) x_i$$

Bonus Slide: Multivariate Chain Rule for MDS

- General **MDS** formulation:

$$\operatorname{argmin}_{Z \in \mathbb{R}^{n \times k}} \sum_{i=1}^n \sum_{j=i+1}^n g(d_1(x_i, x_j), d_2(z_i, z_j))$$

- Using **multivariate chain rule** we have:

$$\nabla_{z_i} g(d_1(x_i, x_j), d_2(z_i, z_j)) = g'(d_1(x_i, x_j), d_2(z_i, z_j)) \nabla_{z_i} d_2(z_i, z_j)$$

- **Example:**

If $d_1(x_i, x_j) = \|x_i - x_j\|$ and $d_2(z_i, z_j) = \|z_i - z_j\|$ and $g(d_1, d_2) = \frac{1}{2}(d_1 - d_2)^2$

$$\nabla_{z_i} g(d_1(x_i, x_j), d_2(z_i, z_j)) = \underbrace{-(d_1(x_i, x_j) - d_2(z_i, z_j))}_{g'(d_1, d_2)} \left[\underbrace{-\frac{(z_i - z_j)}{2\|z_i - z_j\|}}_{\text{(how distance changes in } z \text{ space)}} \right] \rightarrow \nabla_{z_i} d_2(z_i, z_j)$$

↳ Assuming $z_i \neq z_j$ (move distances closer)

Latent-Factor Representation of Words

- For natural language, we often **represent words by an index**.
 - E.g., “cat” is word 124056 among a “bag of words”.
- But this may be inefficient:
 - Should “cat” and “kitten” **share parameters** in some way?

Latent-Factor Representation of Words

- **Latent-factor representation** of individual words:
 - Closeness in latent space should indicate similarity.
 - Distances could represent meaning?
- Recent alternative to PCA/NMF is **word2vec...**

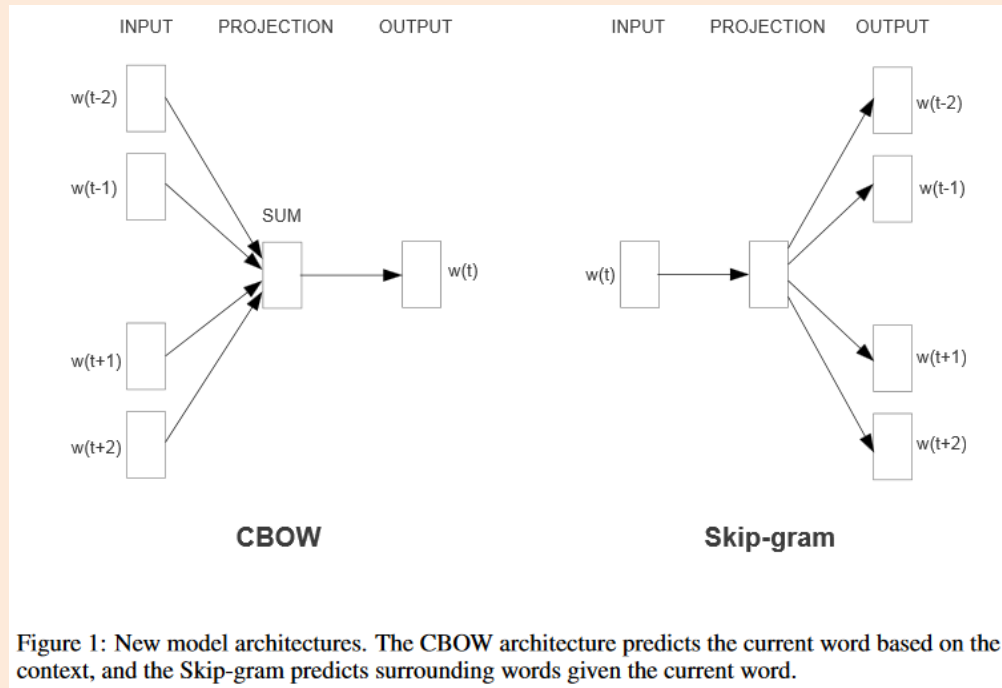
Using Context

- Consider these phrases:
 - “the cat purred”
 - “the kitten purred”

 - “black cat ran”
 - “black kitten ran”
- Words that occur in the same context likely have similar meanings.
- **Word2vec** uses this insight to design an **MDS distance function**.

Word2Vec

- Two common **word2vec** approaches:
 1. Try to **predict word from surrounding words** (continuous bag of words).
 2. Try to **predict surrounding words from word** (skip-gram).



- Train **latent-factors** to solve one of these supervised learning tasks.

Word2Vec

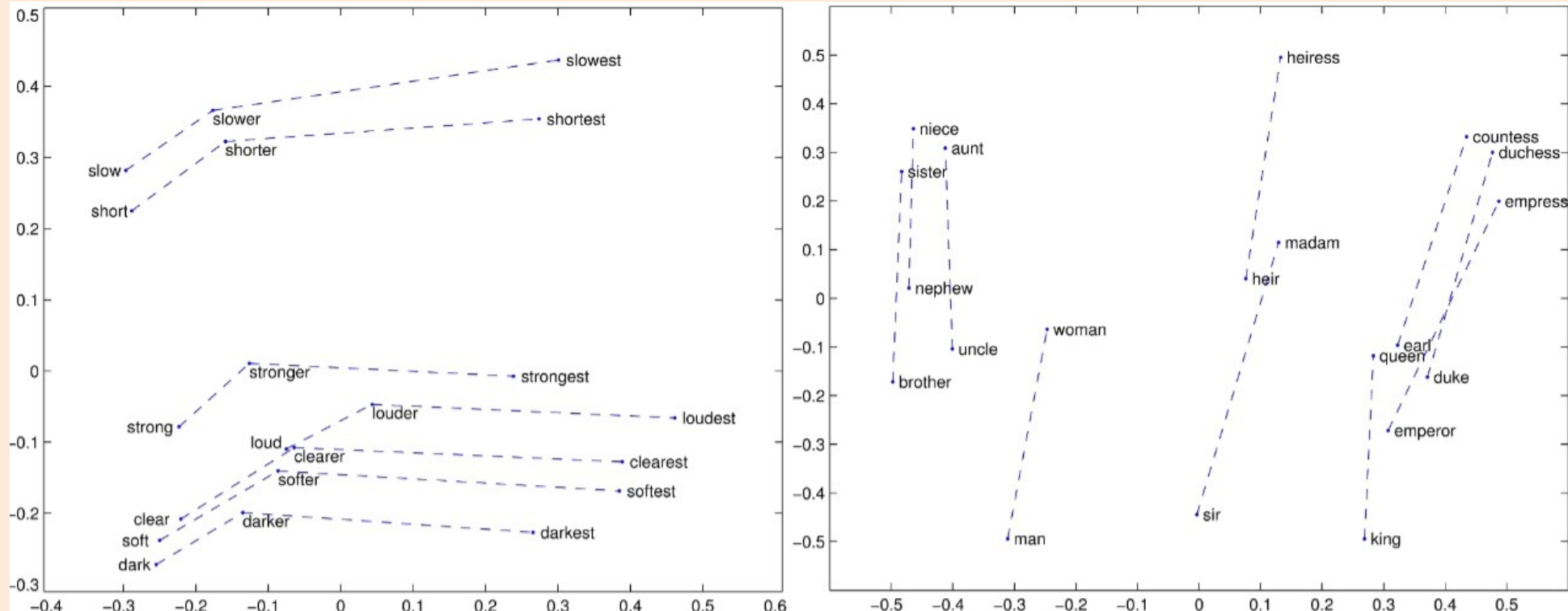
- In both cases, each word 'i' is represented by a vector z_i .
- In **continuous bag of words (CBOW)**, we optimize the following likelihood:

$$p(x_i | x_{\text{surround}}) = \prod_{j \in \text{surround}} p(x_i | x_j) \quad (\text{independence assumption})$$
$$= \prod_{j \in \text{surround}} \frac{\exp(z_i^T z_j)}{\sum_{c=1}^K \exp(z_c^T z_j)} \quad (\text{softmax over all words})$$

- Apply gradient descent to logarithm:
 - Encourages $z_i^T z_j$ to be big for words in same context (making z_i close to z_j).
 - Encourages $z_i^T z_j$ to be small for words not appearing in same context (makes z_i and z_j far).
- For **CBOW**, denominator sums over all words.
- For **skip-gram** it will be over **all possible surrounding words**.
 - Common trick to speed things up: sample terms in denominator (“negative sampling”).

Word2Vec Example

- MDS visualization of a set of related words:



- Distances between vectors might represent semantics.

Word2Vec

- Subtracting word vectors to find related vectors.

Table 8: *Examples of the word pair relationships, using the best word vectors from Table 4 (Skip-gram model trained on 783M words with 300 dimensionality).*

Relationship	Example 1	Example 2	Example 3
France - Paris	Italy: Rome	Japan: Tokyo	Florida: Tallahassee
big - bigger	small: larger	cold: colder	quick: quicker
Miami - Florida	Baltimore: Maryland	Dallas: Texas	Kona: Hawaii
Einstein - scientist	Messi: midfielder	Mozart: violinist	Picasso: painter
Sarkozy - France	Berlusconi: Italy	Merkel: Germany	Koizumi: Japan
copper - Cu	zinc: Zn	gold: Au	uranium: plutonium
Berlusconi - Silvio	Sarkozy: Nicolas	Putin: Medvedev	Obama: Barack
Microsoft - Windows	Google: Android	IBM: Linux	Apple: iPhone
Microsoft - Ballmer	Google: Yahoo	IBM: McNealy	Apple: Jobs
Japan - sushi	Germany: bratwurst	France: tapas	USA: pizza

Table 8 shows words that follow various relationships. We follow the approach described above: the relationship is defined by subtracting two word vectors, and the result is added to another word. Thus for example, $Paris - France + Italy = Rome$. As it can be seen, accuracy is quite good, although

- Word vectors for 157 languages [here](#).

Multiple Word Prototypes

- What about **homonyms** and **polysemy**?
 - The word vectors would **need to account for all meanings.**
- More recent approaches:
 - Try to **cluster the different contexts** where words appear.
 - Use **different vectors for different contexts.**

$$X_{jaguar} \approx \begin{bmatrix} \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \\ \cdot & \cdot & \cdot & \cdot & \cdot & \cdot & \cdot \end{bmatrix} \begin{matrix} z_{j1} \\ z_{j2} \\ z_{j3} \end{matrix}$$

61

