

COMP6237 Data Mining

Lecture 4: Embedding Data (Dimensionality Reduction II)

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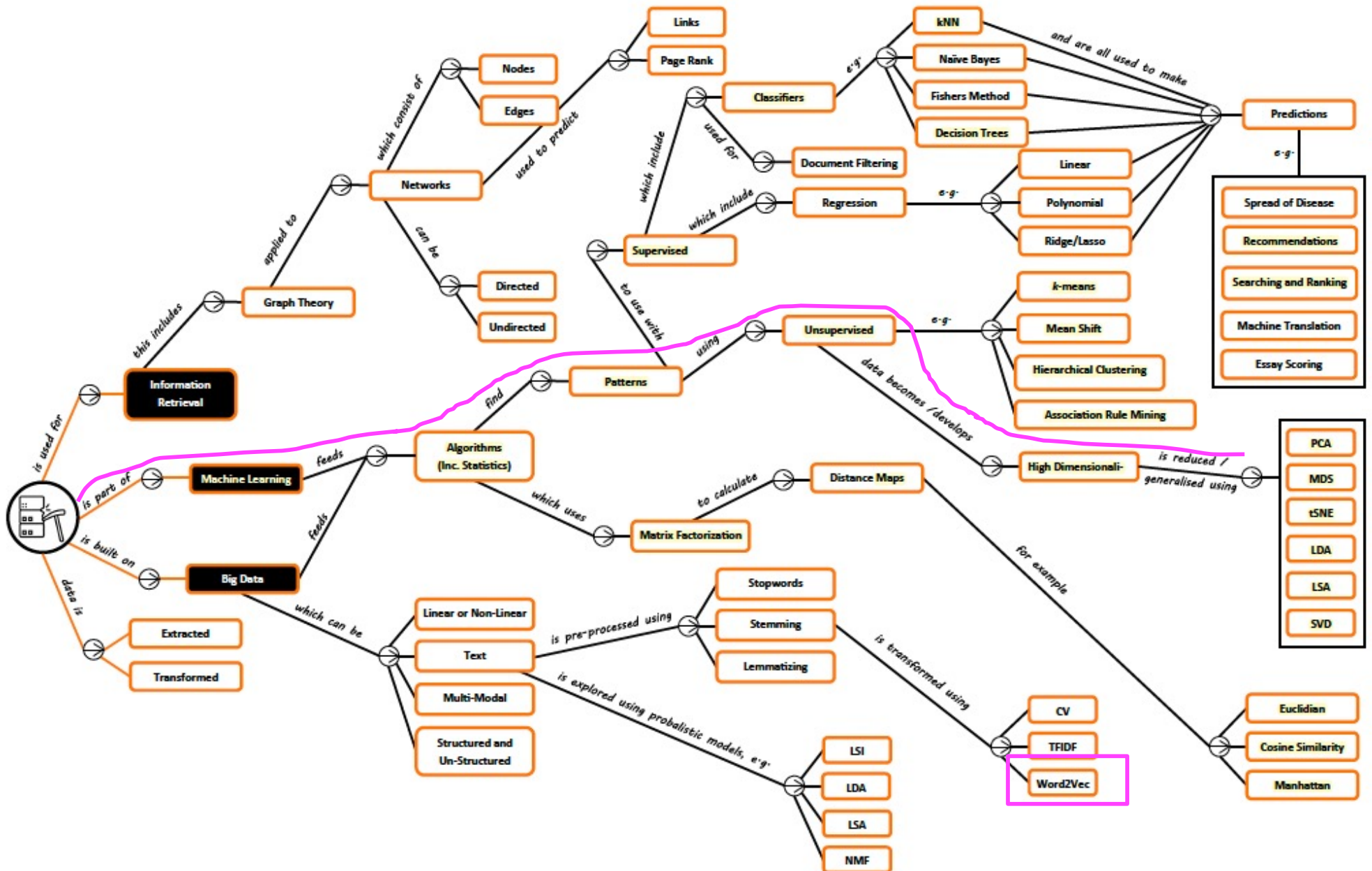
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Lecture slides available here:

<http://comp6237.ecs.soton.ac.uk/zh.html>

(Thanks to Prof. Jonathon Hare and Dr. Jo Grundy for providing the lecture materials used to develop the slides.)

Dimensionality Reduction II – Roadmap



Dimensionality Reduction II – Textbook

CHAPTER 7

Dimensionality Reduction

We saw in Chapter 6 that high-dimensional data has some peculiar characteristics, some of which are counterintuitive. For example, in high dimensions the center of the space is devoid of points, with most of the points being scattered along the surface of the space or in the corners. There is also an apparent proliferation of orthogonal axes. As a consequence high-dimensional data can cause problems for data mining and analysis, although in some cases high-dimensionality can help, for example, for nonlinear classification. Nevertheless, it is important to check whether the dimensionality can be reduced while preserving the essential properties of the full data matrix. This can aid data visualization as well as data mining. In this chapter we study methods that allow us to obtain optimal lower-dimensional projections of the data.

- ▶ Data Mining and Machine Learning: Fundamental Concepts and Algorithms M. L. Zaki and W. Meira

Chapter 11

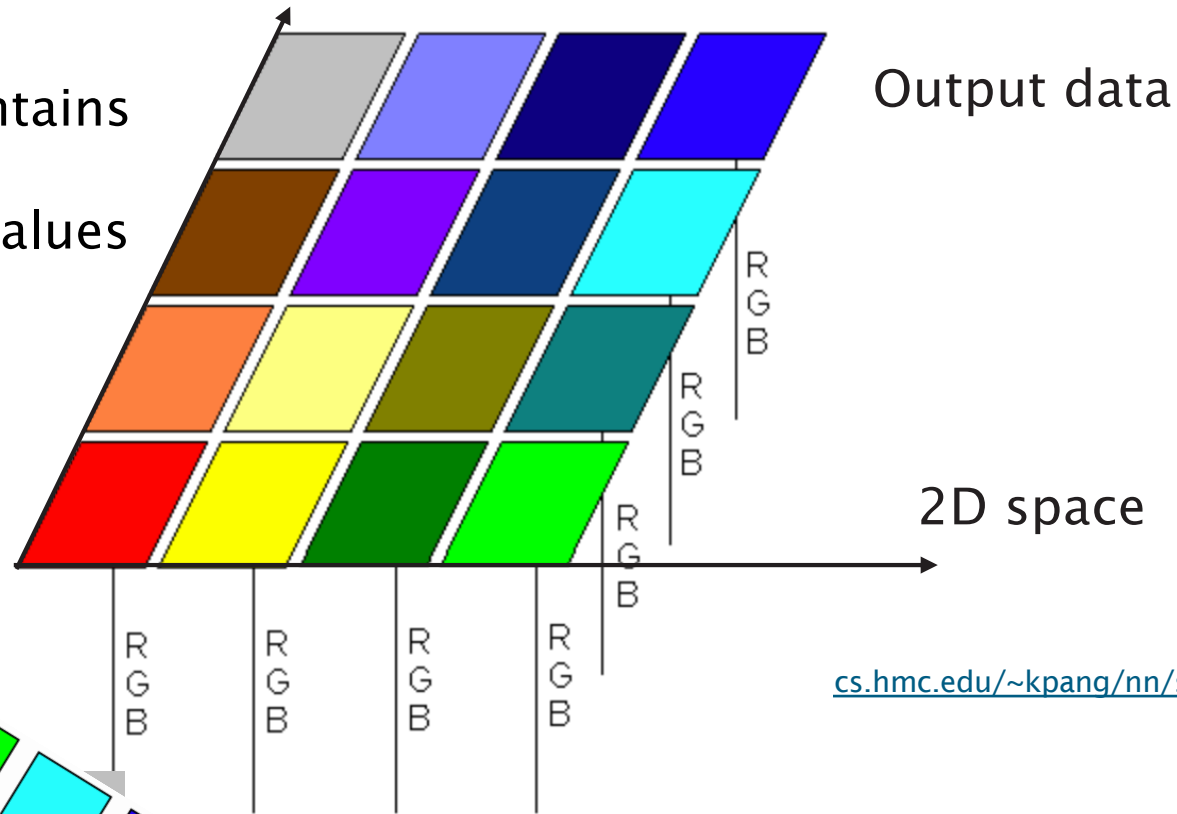
Dimensionality Reduction

There are many sources of data that can be viewed as a large matrix. We saw in Chapter 5 how the Web can be represented as a transition matrix. In Chapter 9, the utility matrix was a point of focus. And in Chapter 10 we

- ▶ Mining of Massive Datasets J. Leskovec *et al*

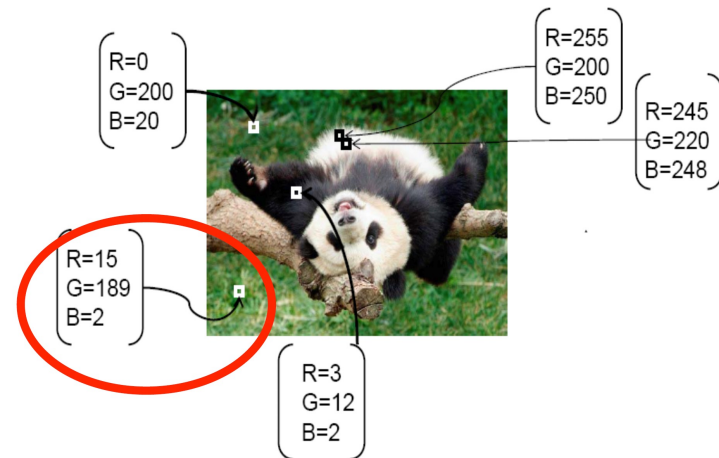
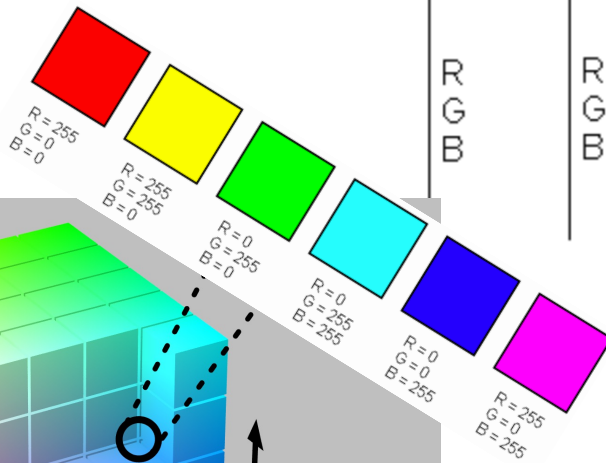
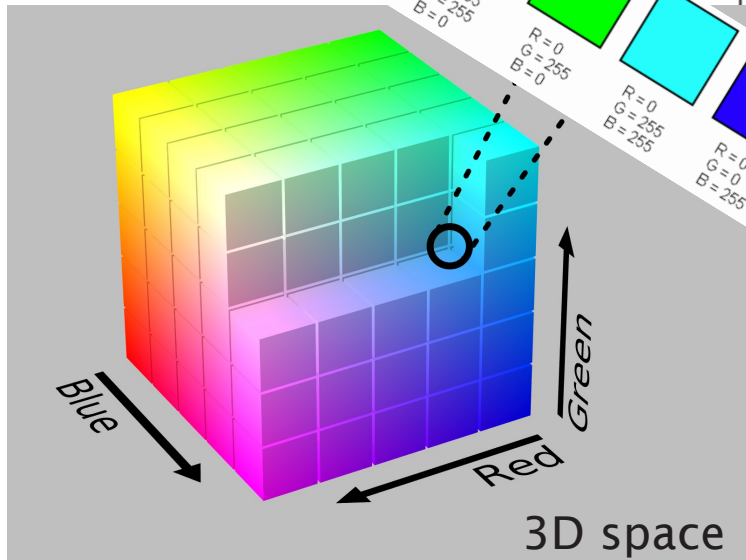
Dimensionality Reduction II – Overview (1/2)

Each neuron/node contains a **3D weight vector** representing its RGB values and a **2D geometric location** in the grid.

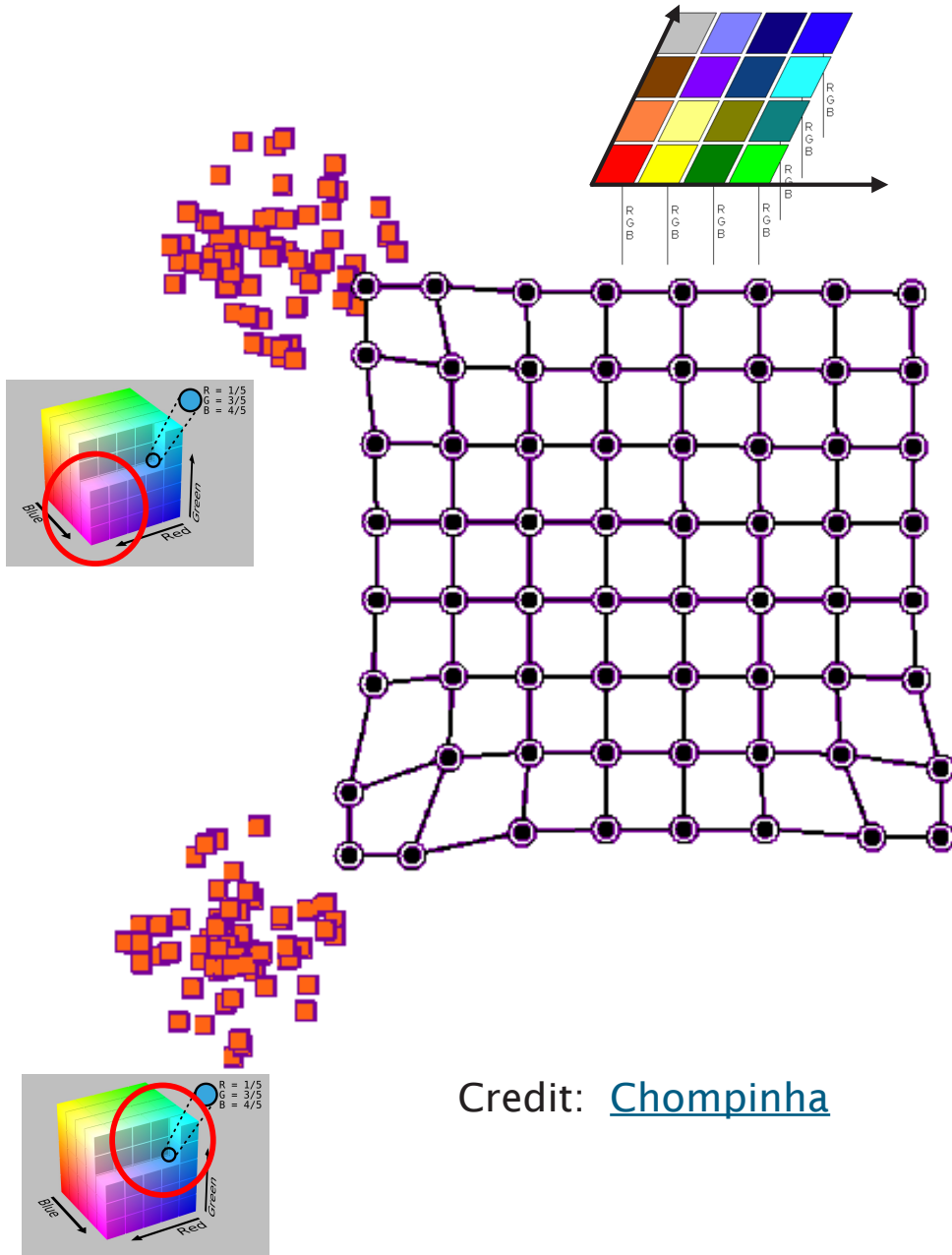


cs.hmc.edu/~kpang/nn/som.html

Input data



Dimensionality Reduction II – Overview (2/2)



- Input data at a high-D space
- ⊙ Output data at a low-D space

Generate a reduced-dimensional representation, usually two-dimensional, of a higher-dimensional dataset while **maintaining the topological structure of the data.**

Credit: [Chompinha](#)

Dimensionality Reduction II – Learning Outcomes

- **LO1:** understand the basic idea of all the learned dimensionality reduction methods (exam)
 - ❖ E.g., describe the basic idea of the suggested algorithm
 - ❖ E.g., understand the pros and cons of the learned algorithms
 - ❖ E.g., given a dataset, be prepared to follow the selected algorithm to calculate its key steps on the given data
- **LO2:** Implement the learned algorithms for dimensionality reduction (course work)

Assessment hints: Multi-choice Questions (single answer: concepts, calculation etc)

- *Textbook Exercises: textbooks (Programming + Mining)*
- *Other Exercises: <https://www-users.cse.umn.edu/~kumar001/dmbook/sol.pdf>*
- *ChatGPT or other AI-based techs*

Embedding Data

Understanding large data sets is *hard*

Especially when the data are highly dimensional

It would help if we knew:

- ▶ which data items are similar
- ▶ which features are similar

Embedding Data

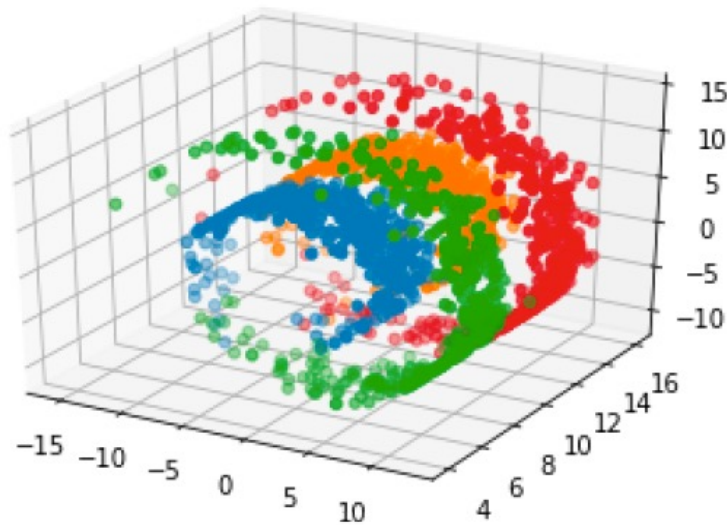
With 2D data, we can plot it to easily visualise relationships

This is not possible with highly dimensional data

However: PCA can reduce the dimensionality to 2, based on the first and second principle axes

Embedding Data - PCA

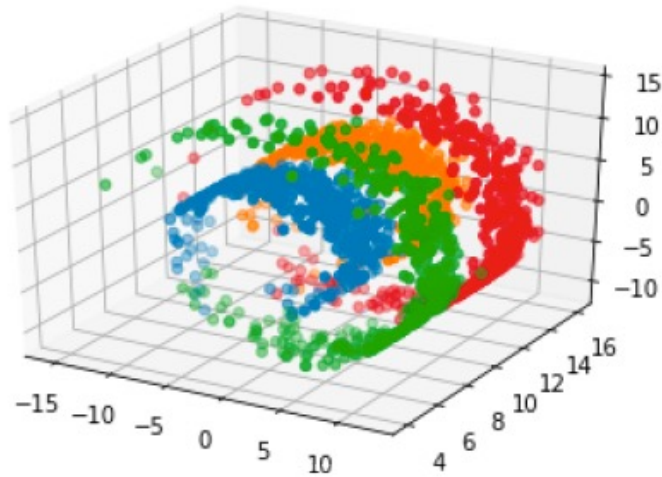
We can use the so called 'Swiss Roll' data set to exemplify this:



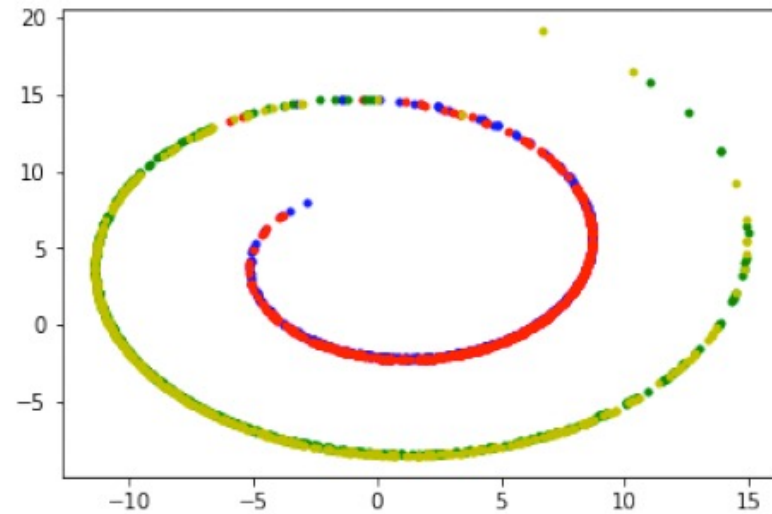
The data in 3D has 4 clearly separate groups

Embedding Data - PCA

We can use the so called 'Swiss Roll' data set to exemplify this:



The data in 3D has 4 clearly separate groups



Using PCA, it does not separate the data well at all

Embedding Data - PCA

Unfortunately there is no control over the distance measure

Using axes of greatest variance does not mean similar things appear close together.

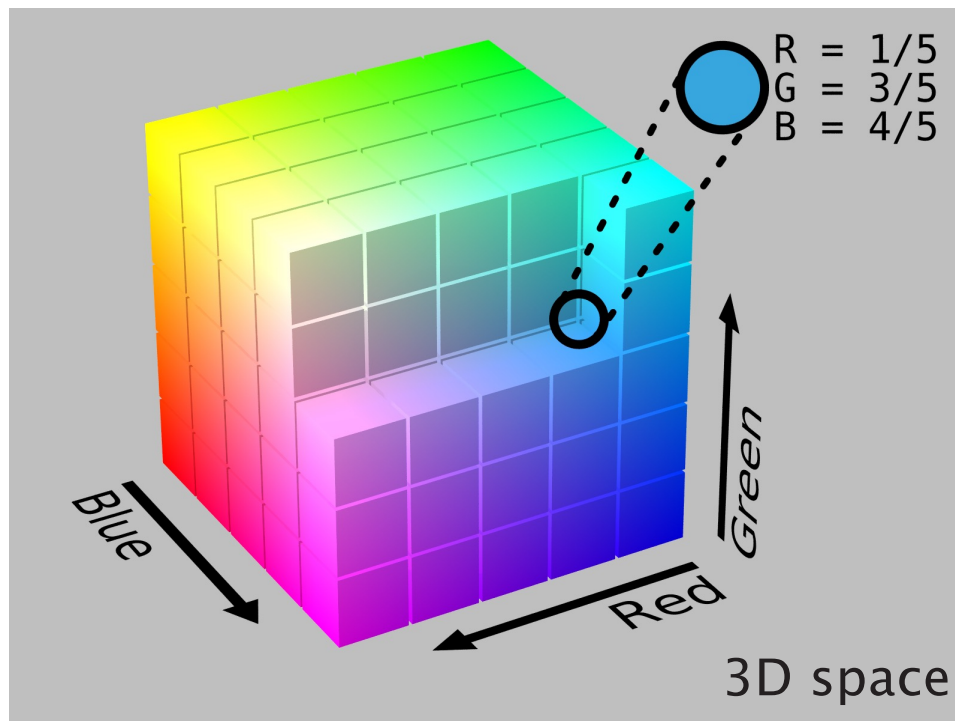
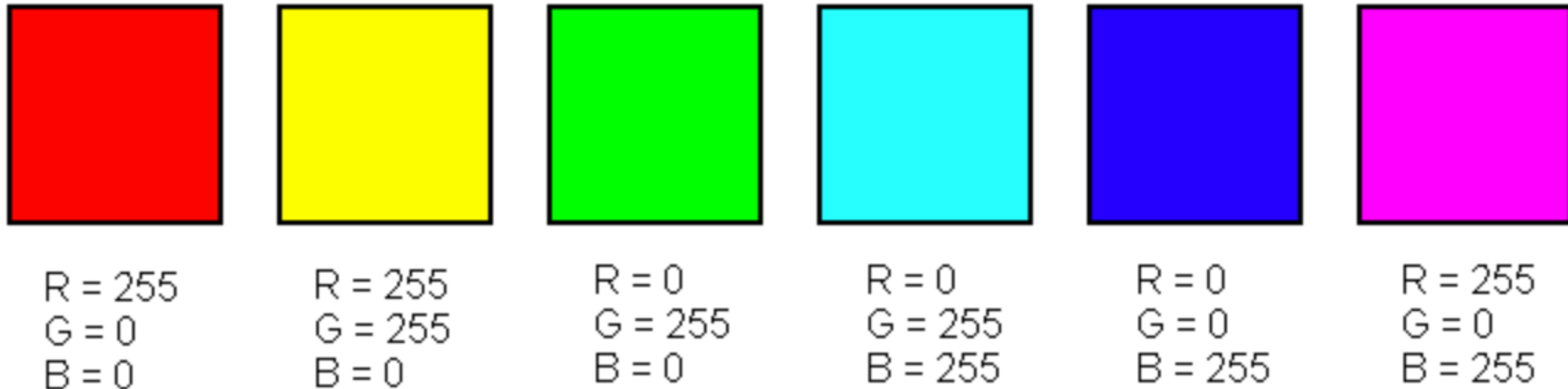
PCA is only rotation of original space, followed by removal of less significant dimensions

Embedding Data – Self-Organising Maps

Kohonen 1982: Self-Organising Maps (SOM)

- ▶ Inspired by neural networks
- ▶ 2D n by m array of nodes
- ▶ Units close to each other are considered to be neighbours
- ▶ Maps high dimensional vectors to unit with coordinates closest (Euclidean Distance)
- ▶ This is the *best matching unit*

Embedding Data – Self-Organising Maps



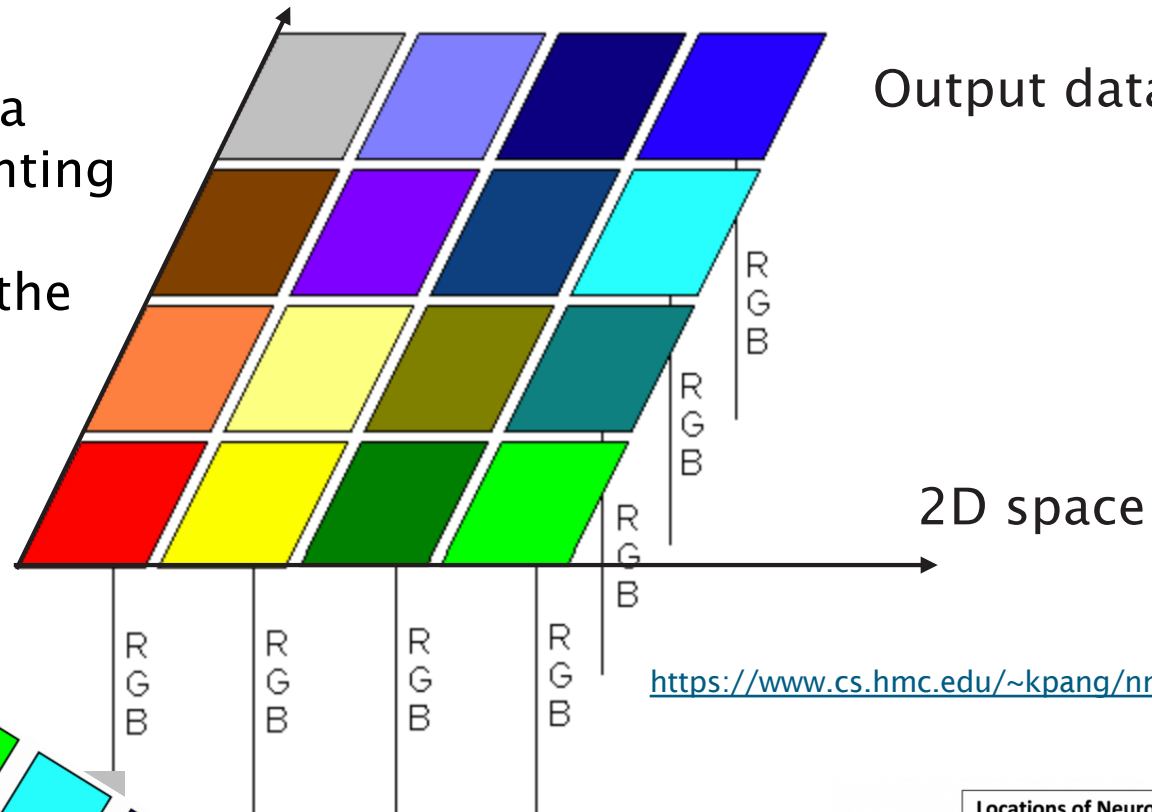
A good mapping would group the red, green, and blue colors far away from one another and place the intermediate colors between their base colors (e.g. yellow close to red and green, teal close to green and blue, etc)

<https://www.cs.hmc.edu/~kpanq/nn/som.html>

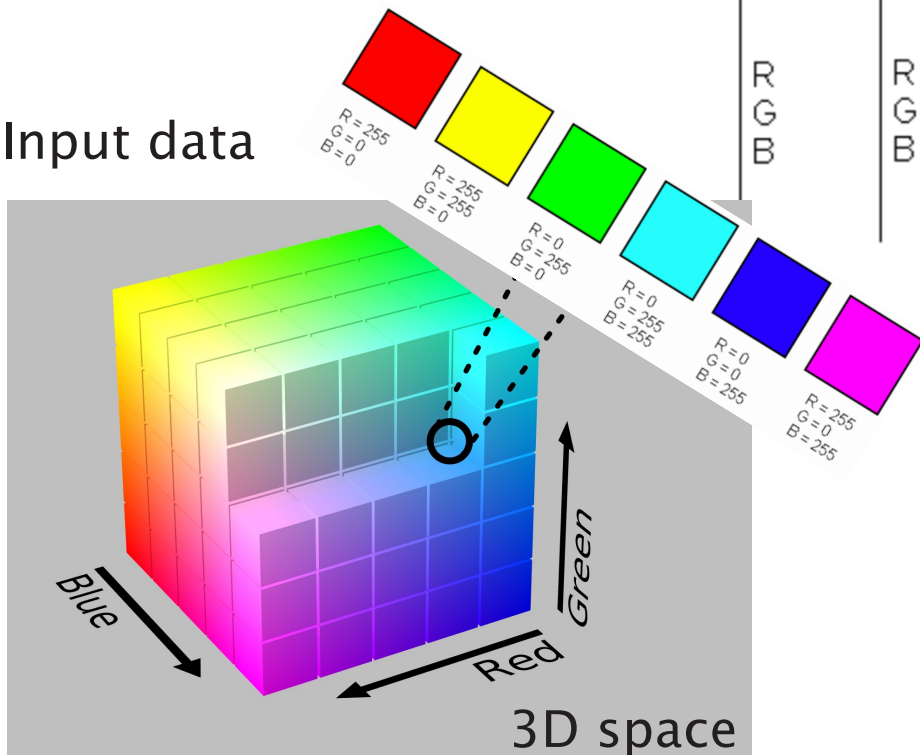
<https://en.wikipedia.org>

Embedding Data – Self-Organising Maps

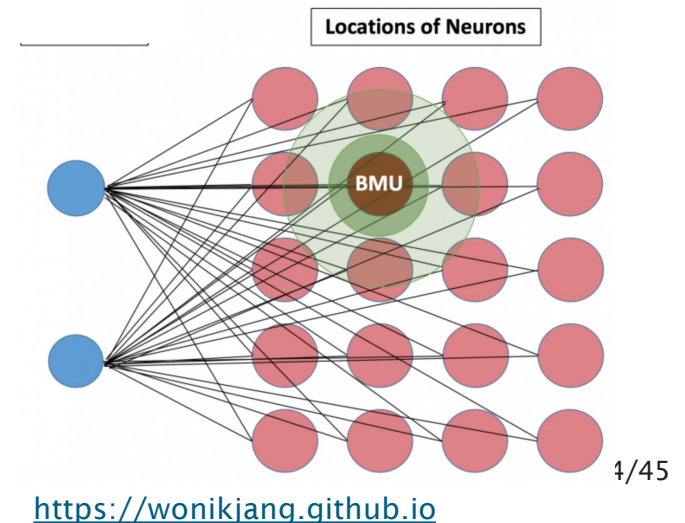
Each neuron contains a weight vector representing its RGB values and a geometric location in the grid.



Input data



<https://www.cs.hmc.edu/~kpang/nn/som.html>



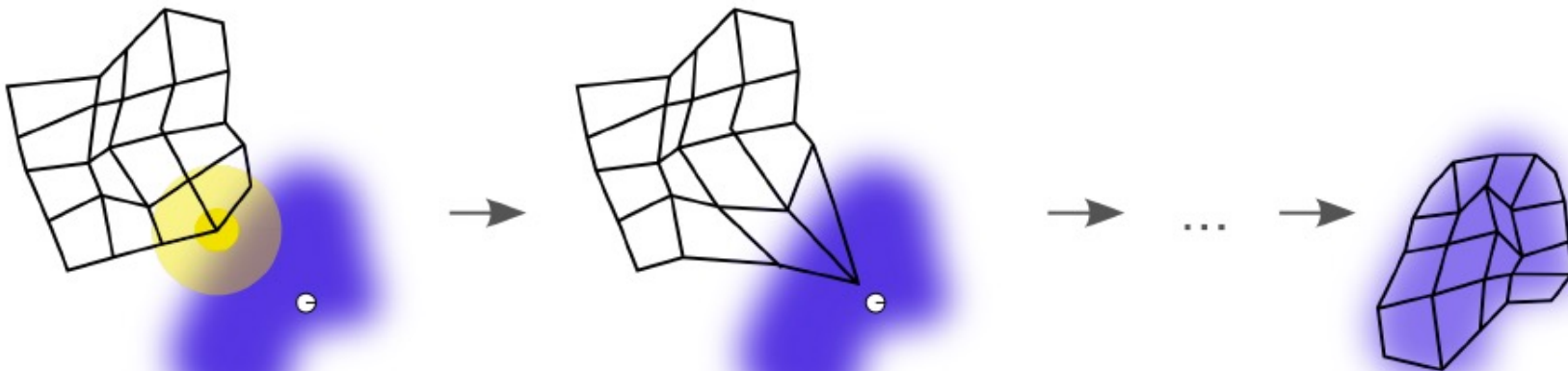
Embedding Data – Self-Organising Maps

SOMs: two phases

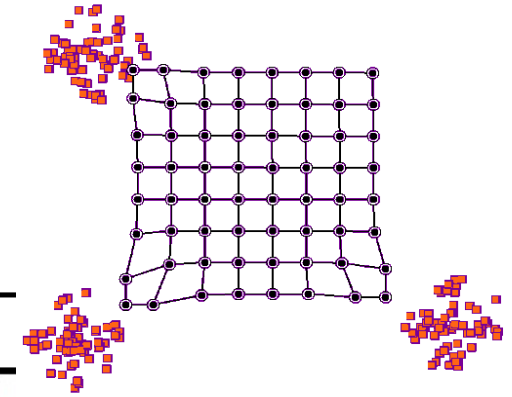
- ▶ training
- ▶ mapping

To start training, the set of nodes each has a random starting position defined in the feature space.

This is then updated by taking one feature vector, finding which unit is the *best matching unit* (BMU) then moving that unit and, to a lesser extent, its neighbours, closer to that data point



Embedding Data – Self-Organising Maps



Algorithm 1: Training Self-Organising Maps

Data: N data points with d dimensional feature vectors X_i
 $i = 1 \dots N$, number of iterations λ

w = randomly initialise $n \times m$ units with weight vector ;

$t = 0$;

while $t < \lambda$ **do**

for each x_i **do**

$BMU = w_{nm}$ with min distance;

 Update BMU and its neighbours by moving closer to x_i ;

end

$t = t + 1$

end

Credit: [Chompinha](#)

Embedding Data – Self-Organising Maps

To update the weight vector \mathbf{w}

$$\mathbf{w}(t + 1) = \mathbf{w}(t) + \theta(u, v, t)\alpha(t)(x_i - \mathbf{w}(t))$$

where θ is the neighbourhood weighting function (usually Gaussian)

$$\text{e.g., } \theta = \begin{cases} 1 & \text{if the node is the winner} \\ 0.5 & \text{if the node is immediate neighbour of the winner} \\ 0 & \text{otherwise} \end{cases}$$

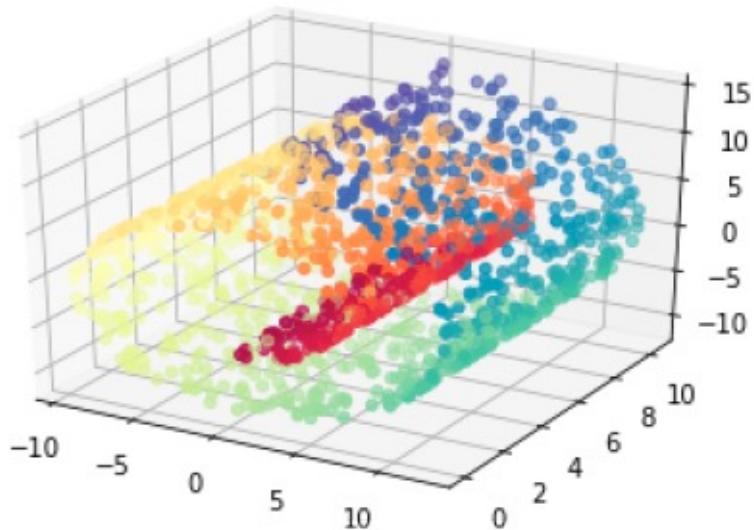
α is the learning rate

x_i is the input vector

u is the winner node, and v is the node whose weight is \mathbf{w}

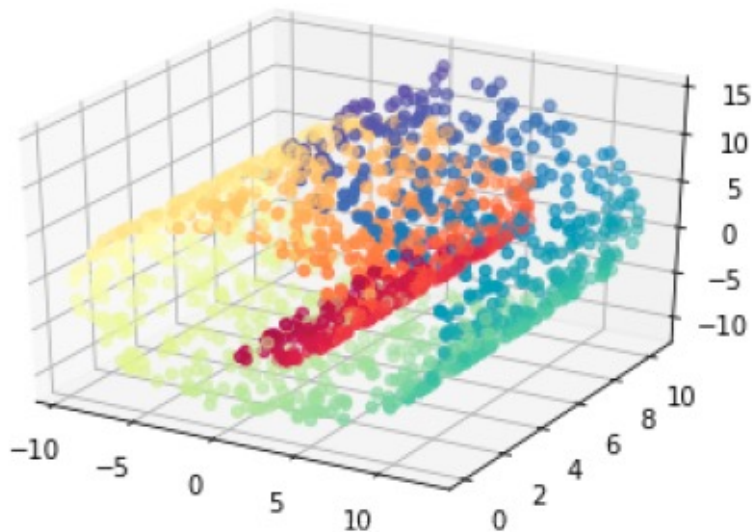
Both the learning rate and the neighbourhood weighting function get smaller over time

Embedding Data – Self-Organising Maps

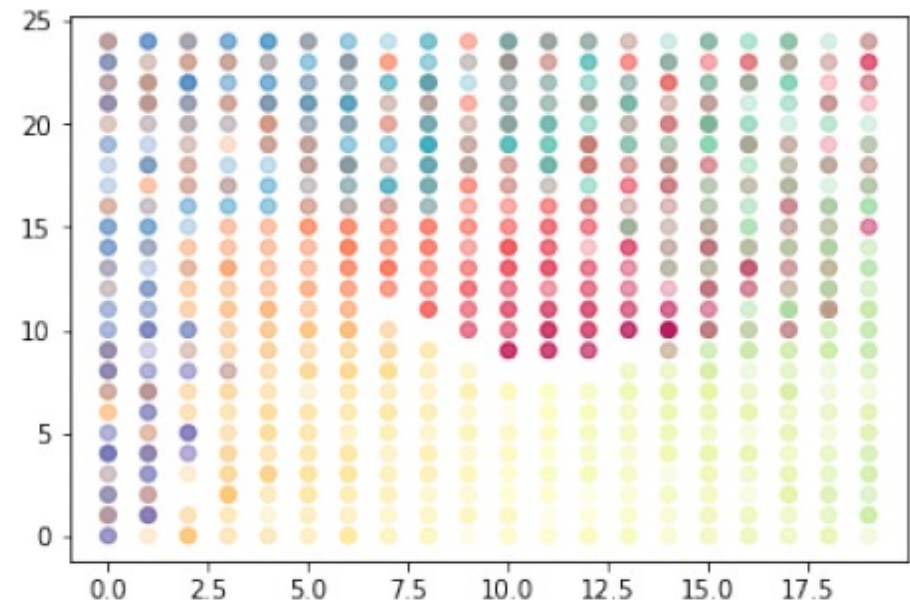


The data in 3D has a clear structure

Embedding Data – Self-Organising Maps



The data in 3D has a clear structure

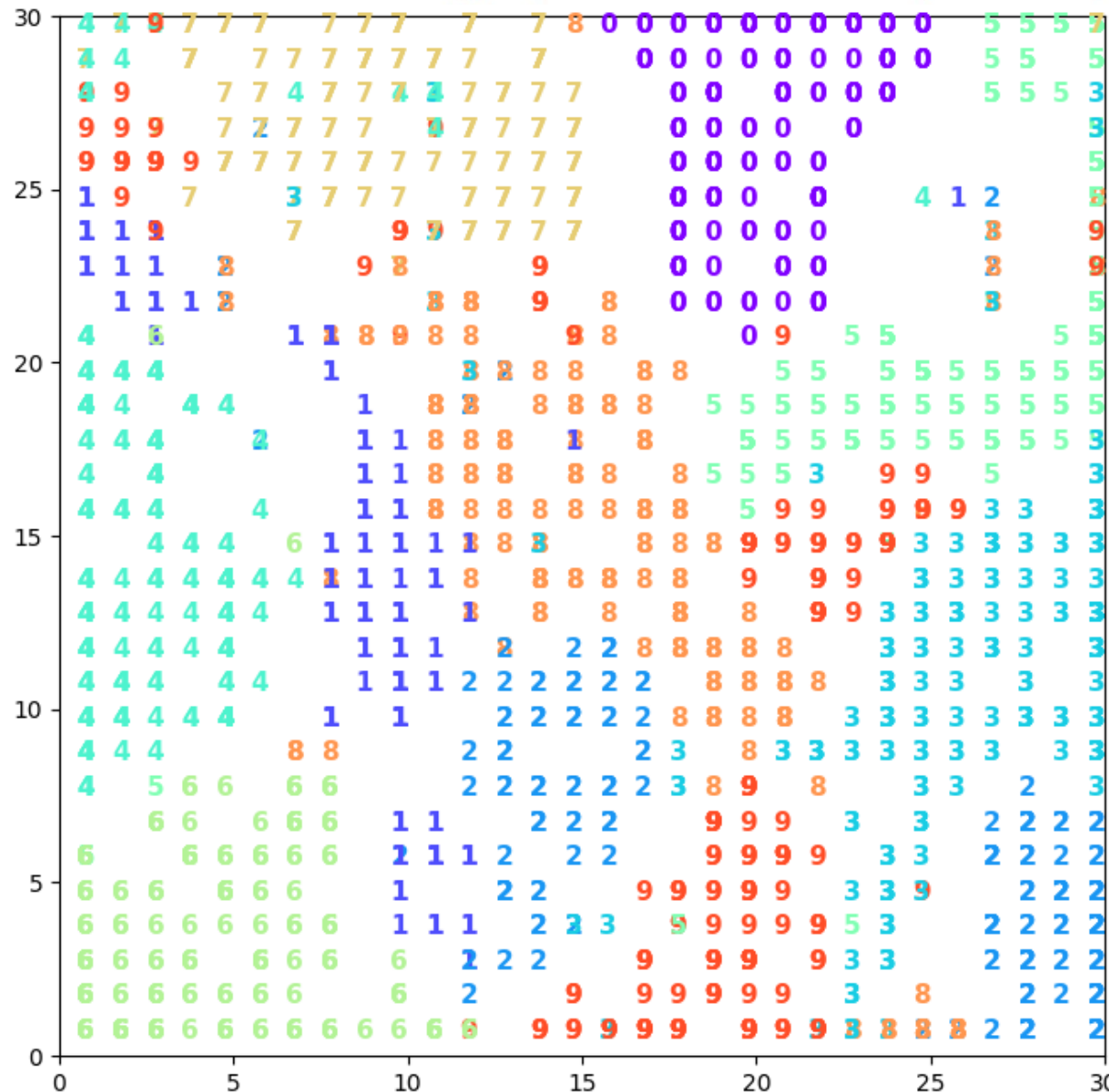


Using SOM, it can much better group the data

[ipynb mean centering demo](https://github.com/zhiwu-huang/Data-Mining-Demo-Code-18-19) <https://github.com/zhiwu-huang/Data-Mining-Demo-Code-18-19>

Embedding Data – Self-Organising Maps

With the MNIST digits, the results are quite impressive



ipython mean centering demo

<https://github.com/zhiwu-huang/Data-Mining-Demo-Code-18-19>

Embedding Data – Multidimensional Scaling

Multi Dimensional Scaling involves:

- ▶ Start with data in a high dimensional space and a set of corresponding points in a lower dimensional space
- ▶ Optimise the positions of points in lower dimensional space so their Euclidean distances are *like* the distances between the high dimensional points
- ▶ Can use any distance measure in the high D space

Embedding Data – Multidimensional Scaling

There are two main sorts of multidimensional scaling:

- ▶ Metric MDS - Tries to match distances
- ▶ Non-metric MDS - tries to match rankings

Only requires distances between items as input

Unlike PCA and SOM, there is no explicit mapping

Both metric and non-metric measure goodness of fit between two spaces

They try to minimise a *stress function*

Embedding Data – Multidimensional Scaling

Stress functions:

- ▶ Least-squares scaling / Kruskal-Shepard scaling
- ▶ Shepard-Kruskal non-metric scaling
- ▶ Sammon Mapping

Sammon Mapping is given by:

$$S(z_1, z_2, \dots, z_n) = \sum_{i \neq j} \frac{(\delta_{ij} - \|z_i - z_j\|)^2}{\delta_{ij}}$$

where $\delta_{i,j}$ is the distance in high dimensional space
and $\|z_i - z_j\|$ is the distance in low dimensional space
Looks at all combinations of points with all different points.

Embedding Data – Multidimensional Scaling

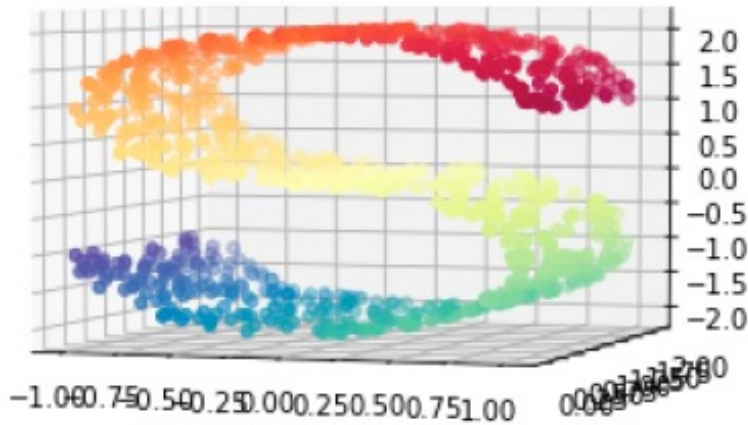
For non-linear, need to use gradient descent
start at arbitrary point, take steps in direction of gradient, with
step size a proportion of the gradient magnitude.

$$z_j(k + 1) = z_j(k) - \gamma_k \Delta_{z_j} S(z_1(k), z_2(k), \dots, z_n(k))$$

Where the derivative of the Sammon stress is:

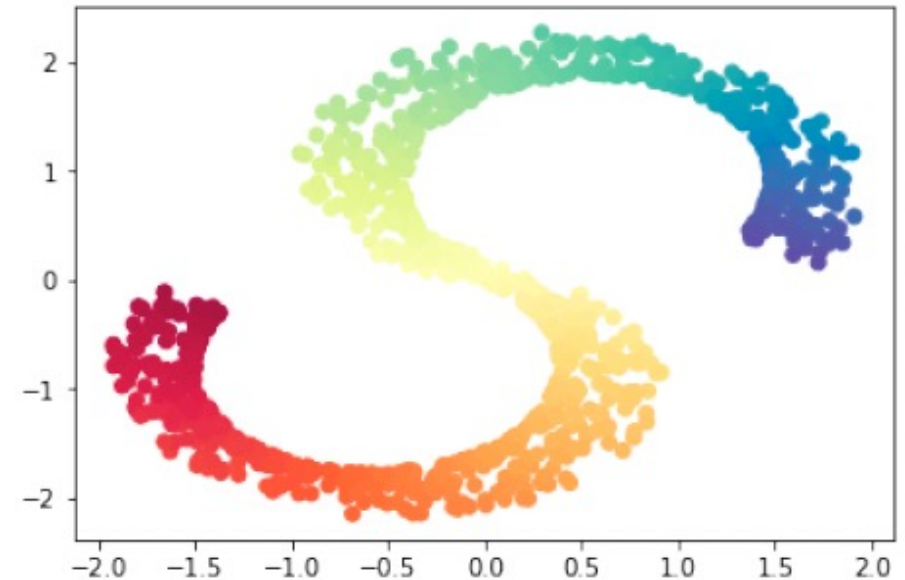
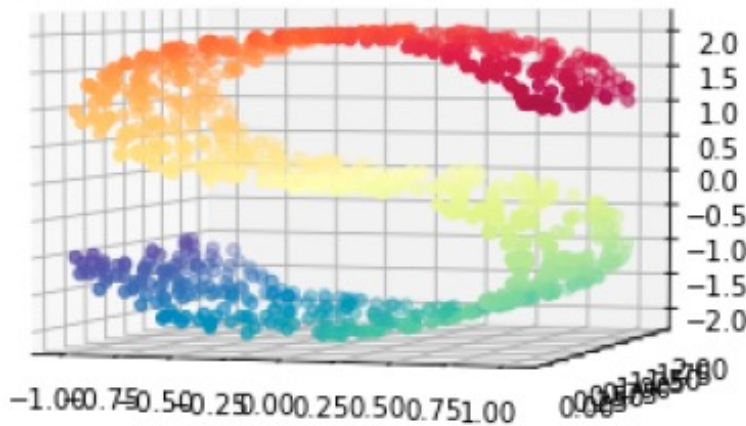
$$\Delta_{z_j} S() = 2 \sum_{i \neq j} \left(\frac{\|z_i(k) - z_j(k)\| - \delta_{ij}}{\delta_{ij}} \right) \left(\frac{z_j(k) - z_i(k)}{\|z_i(k) - z_j(k)\|} \right)$$

Embedding Data – Multidimensional Scaling



The data in 3D has a clear structure.
MDS gives..

Embedding Data – Multidimensional Scaling



The data in 3D has a clear structure.
MDS gives..

Not particularly brilliant, has red the same distance from yellow as orange and green
Has however preserved the structure from 3D in to 2D

Embedding Data – Stochastic Neighbour Embedding

Stochastic Neighbour Embedding (SNE)

Works in a similar way to MDS

MDS optimises distances, SNE optimises the distribution of data

Aims to make the distribution of the projected data in low dimensional space close to the actual distribution in high dimensional space

Embedding Data – Stochastic Neighbour Embedding

To calculate the source distribution:

We define a conditional probability that high-dimensional x_i would pick x_j as a neighbour if the neighbours were picked in proportion to their probability density under a Gaussian centred at x_i

$$p_{j|i} = \frac{\exp^{-\|x_i - x_j\|^2 / 2\sigma_i^2}}{\sum_{k \neq i} \exp^{-\|x_i - x_k\|^2 / 2\sigma_i^2}}$$

The SNE algorithm chooses σ for each data point such that smaller σ is chosen for points in dense parts of the space, and larger σ is chosen for points in sparse parts

Embedding Data – Stochastic Neighbour Embedding

To calculate the target distribution:

Define a conditional probability that low-dimensional y_i would pick y_j as a neighbour if the neighbours were picked in proportion to their probability density under a Gaussian centred at y_i

$$q_{j|i} = \frac{\exp(-\|y_i - y_j\|^2)}{\sum_{k \neq i} \exp(-\|y_i - y_k\|^2)}$$

In this space we assume the variance of all Gaussians is $1/\sqrt{2}$ in this space

Embedding Data – Stochastic Neighbour Embedding

To measure the difference between two probability distributions we use the *Kullback-Leibler* (KL) Divergence:

$$D_{KL}(P|Q) = \sum_i P(i) \log \frac{P(i)}{Q(i)}$$

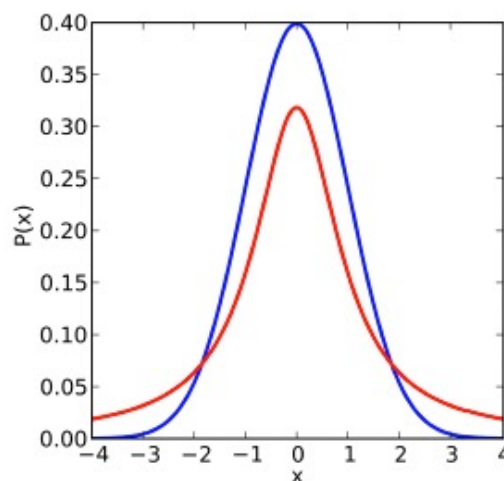
The cost function is the KL divergence summed over all data points

$$C = \sum_i \sum_j p_{i|j} \log \frac{p_{j|i}}{q_{j|i}}$$

C can be minimised using *gradient descent* - but..
difficult to optimise, leads to crowded visualisations, big clumps of data together in the center

Embedding Data – t- Stochastic Neighbour Embedding

In 2008 Maaten and Hinton came up with a way to improve SNE, by replacing the Gaussian distribution for the lower dimensional space with a Student's t distribution.



Student's t distribution in red, Gaussian distribution in blue

The Student's t distribution has a much longer tail, helps avoid clumping in the middle.

Embedding Data – t- Stochastic Neighbour Embedding

The cost function is also modified, making gradients simpler, so faster to compute

$$p_{ij} = \frac{p_{j|i} + p_{i|j}}{2N}$$

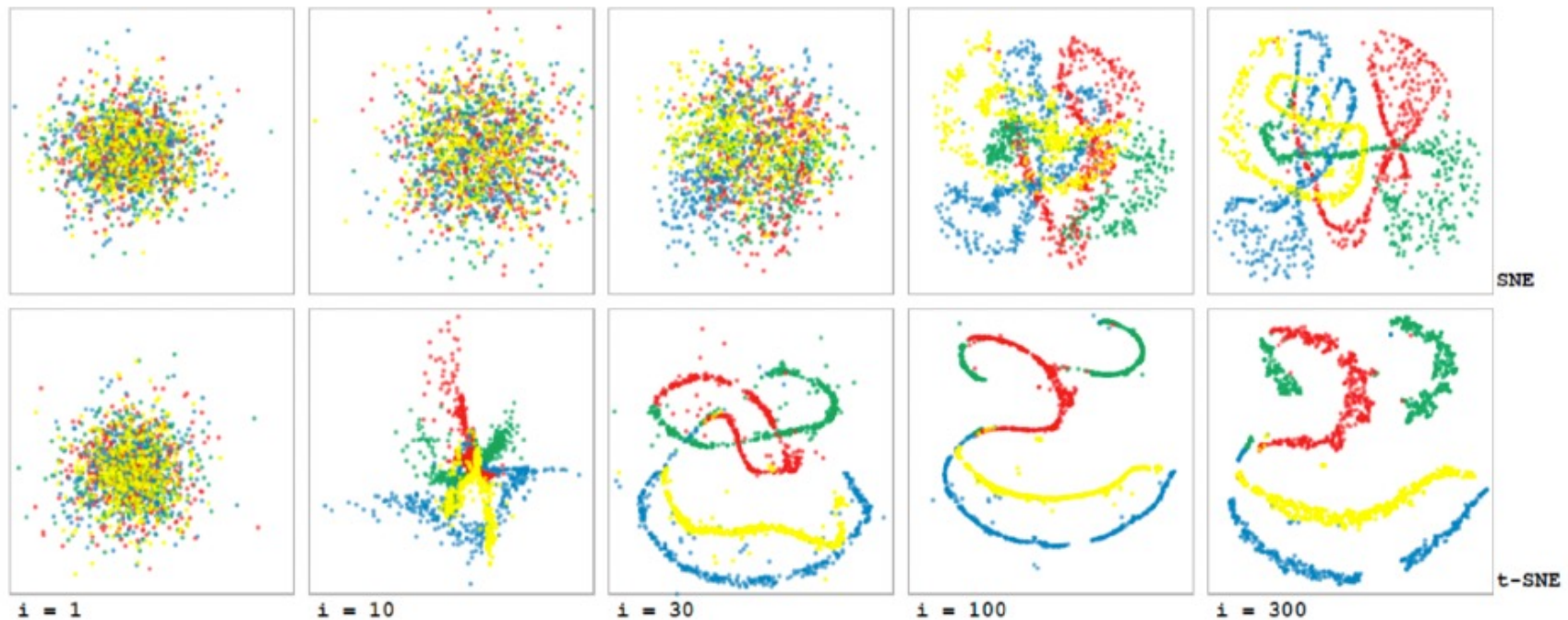
To alleviate crowding using Student's t distribution in the lower dimensional space:

$$q_{ij} = \frac{(1 + \|x_i - x_j\|)^{-1}}{\sum_{k \neq i} (1 + \|x_i - x_k\|)^{-1}}$$

we use 1 degree of freedom with the Student's t distribution, equivalent to a Cauchy distribution

Embedding Data – t- Stochastic Neighbour Embedding

For the Swiss Roll data:



Lorenzo Amabili, <http://lorenzoamabili.github.io>

Embedding Data

Instead of projecting high dimensional data down in to 2 or 3 dimensions, we can use a medium dimensionality, keeping the useful information, capturing the key distinguishing features

This is called an *embedding*

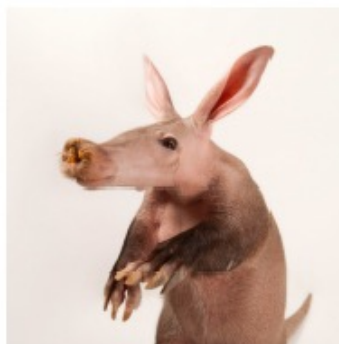
For example: **word2vec**

Embedding Data – One Hot Encoding

For example: Documents

We use a 'Bag of Words', where each word is a vector:

- ▶ $a \rightarrow [1, 0, 0, 0, 0, 0, 0, 0, \dots, 0]$
- ▶ $aa \rightarrow [0, 1, 0, 0, 0, 0, 0, 0, \dots, 0]$
- ▶ $aardvark \rightarrow [0, 0, 1, 0, 0, 0, 0, 0, \dots, 0]$
- ▶ $aardwolf \rightarrow [0, 0, 0, 1, 0, 0, 0, 0, \dots, 0]$



This is called *One Hot Encoding*

Embedding Data – One Hot Encoding

What problems does this encoding have?

- ▶ all vectors are *orthogonal*, i.e. unrelated

But we know that many words in English are related, e.g. ‘rain’, ‘drizzle’, ‘downpour’, ‘shower’, ‘squall’ all mean pretty much the same thing.

- ▶ vectors are very long and *very sparse*

English has over 250,000 words (depending on how you count them) so each vector that could fully describe a document should be 250,000 long for every word.

Embedding Data – word2vec

Boiling this data down to a manageable vector size, while still retaining meaning is not a simple task

word2vec was proposed in 2013 by Mikolov *et al* (although the paper was rejected by the ICLR conference!)

Published at NeurIPS 2013 and cited over 40,000 times, the work introduced the seminal word embedding technique word2vec. Demonstrating the power of learning from large amounts of unstructured text, the work catalyzed progress that marked the beginning of a new era in natural language processing.

NeurIPS 2023 ‘Test of Time Award’

<https://blog.neurips.cc/2023/12/11/announcing-the-neurips-2023-paper-awards/>

Embedding Data – word2vec

For example:

“the quick brown fox jumps over the lazy dog”

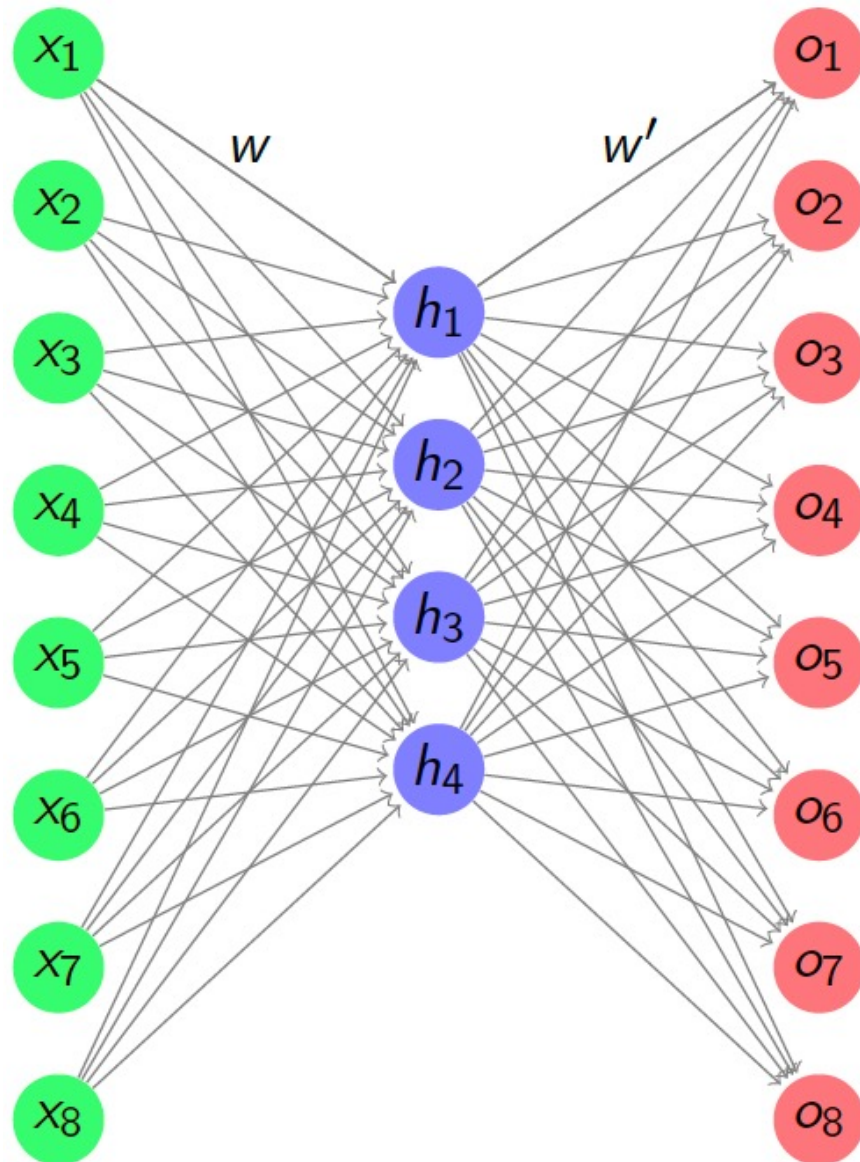
The word ‘brown’ has the words ‘the’, ‘quick’, ‘fox’ and ‘jumps’ close by

This gives training samples:

- ▶ ‘the’, ‘brown’
- ▶ ‘quick’, ‘brown’
- ▶ ‘fox’, ‘brown’
- ▶ ‘jumps’, ‘brown’

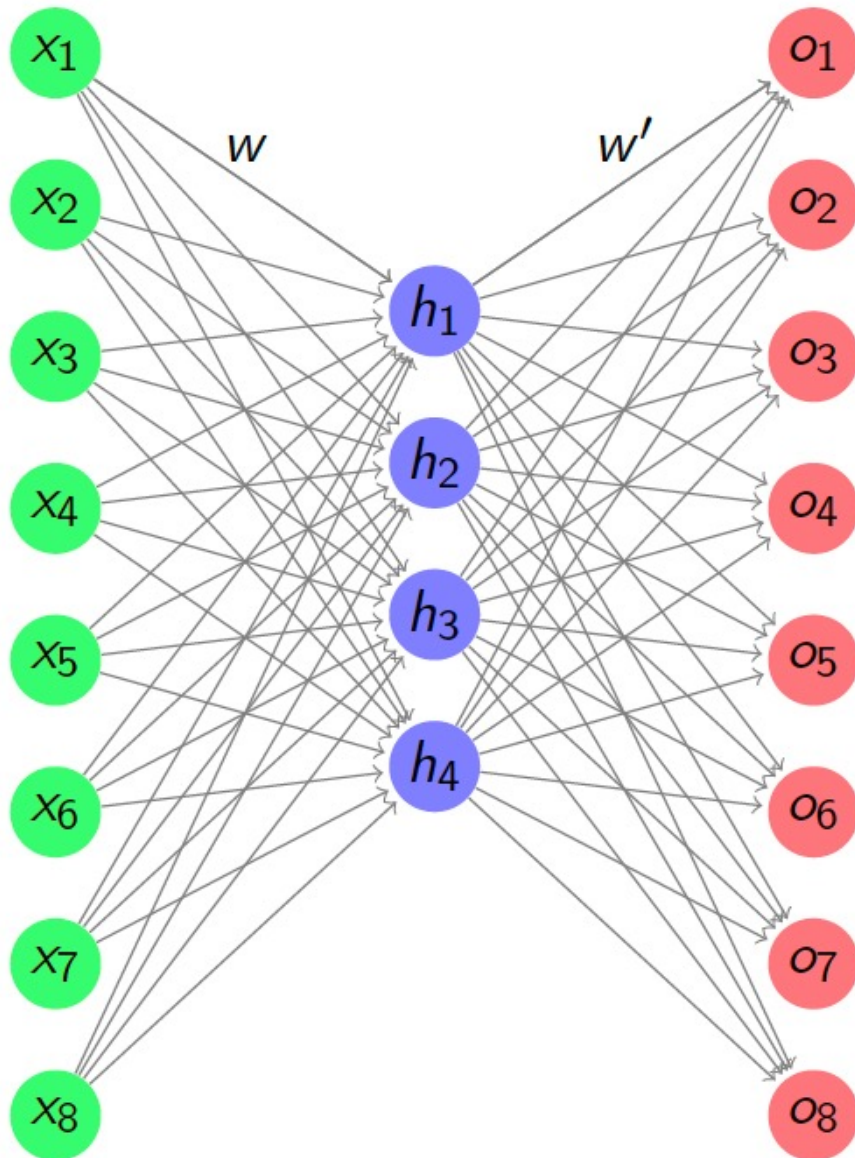
which we train a simple neural network with.

Embedding Data – word2vec



input vector is the 'one hot'
encoding of the word in question

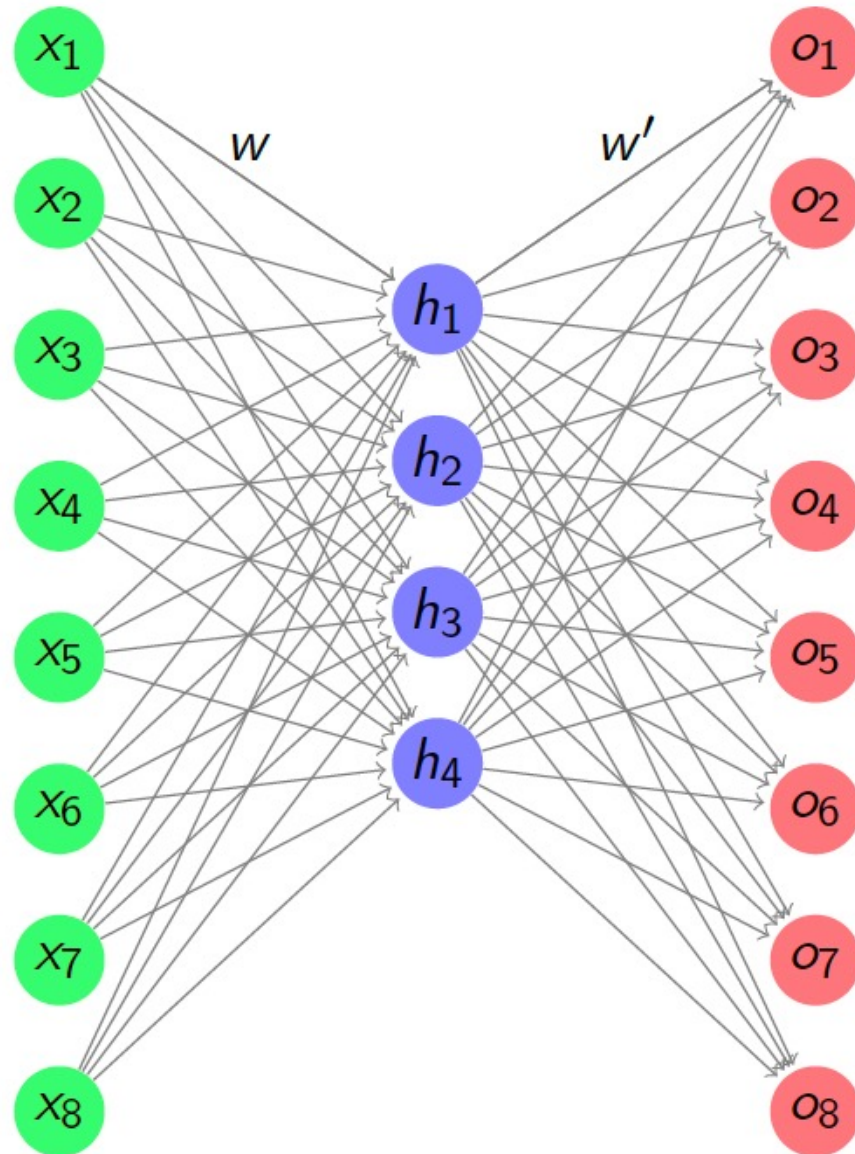
Embedding Data – word2vec



input vector is the 'one hot' encoding of the word in question

the output vector is the vector for the predicted word.

Embedding Data – word2vec

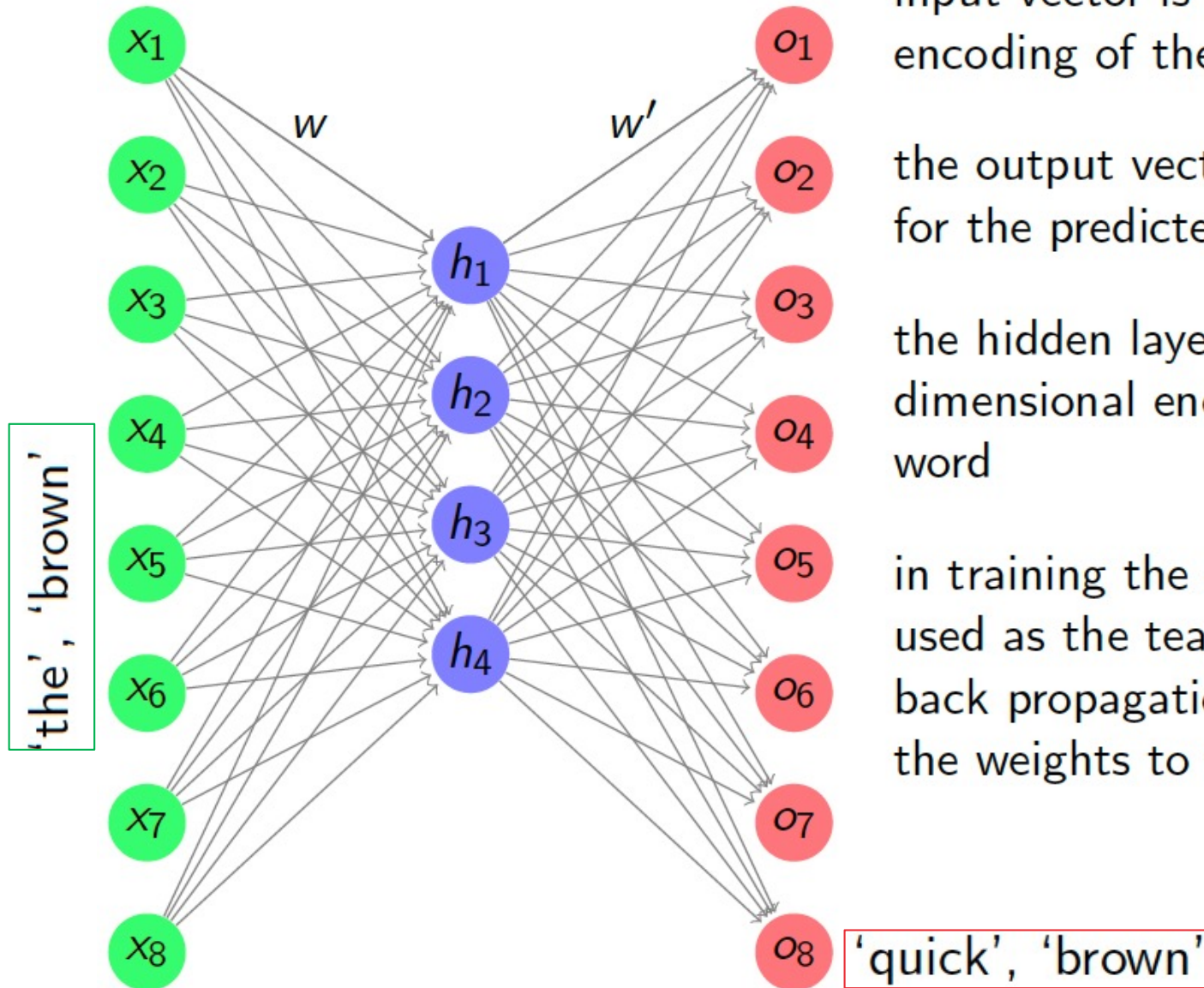


input vector is the 'one hot' encoding of the word in question

the output vector is the vector for the predicted word.

the hidden layer learns a lower dimensional encoding of each word

Embedding Data – word2vec



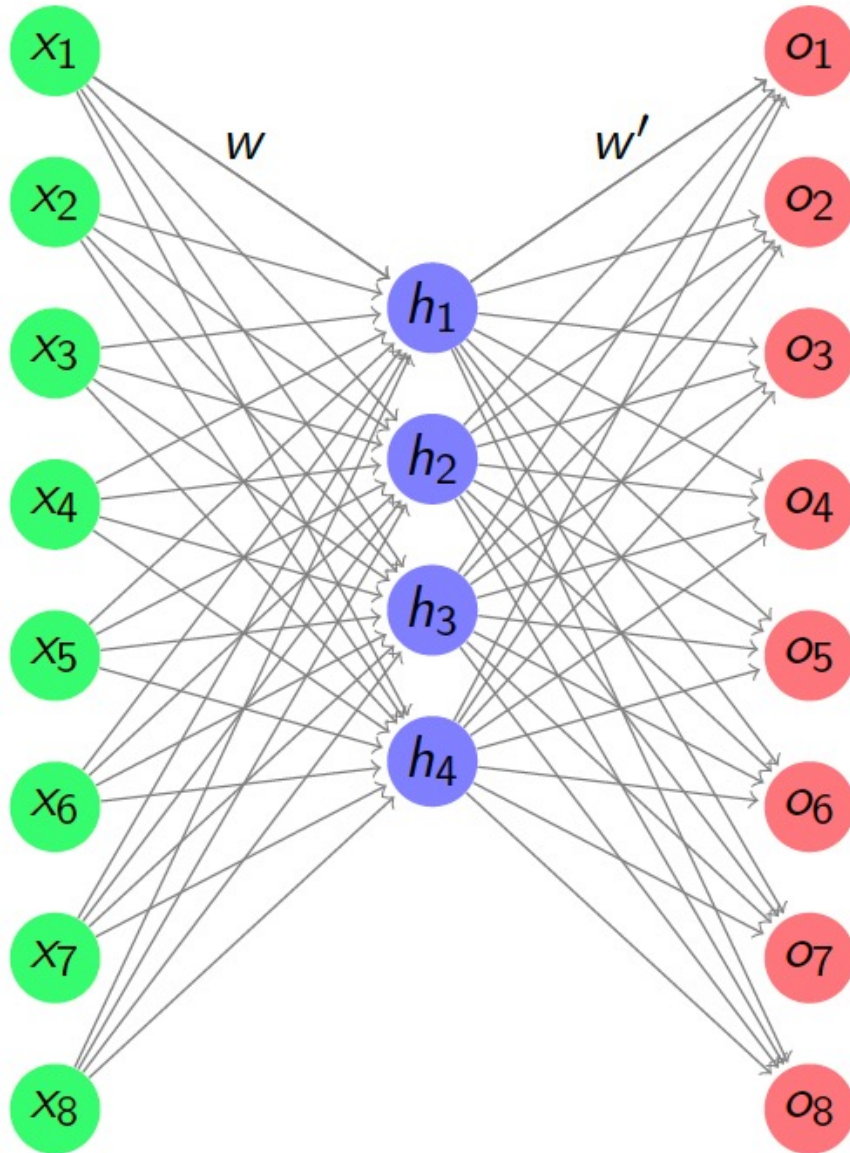
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in training the correct word is used as the teaching signal, and back propagation is used to learn the weights to the hidden layer.

Embedding Data – word2vec



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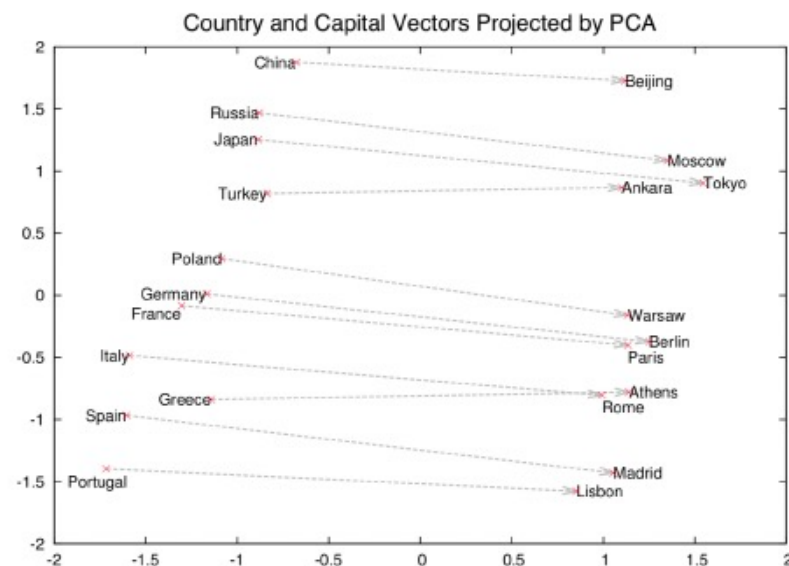
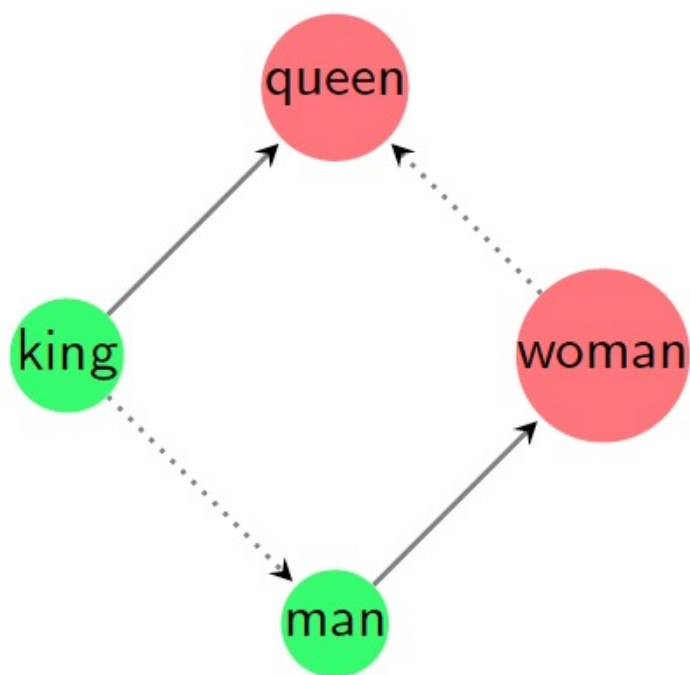
in training the correct word is used as the teaching signal, and back propagation is used to learn the weights to the hidden layer.

the output from the hidden layer after training is used as the lower dimensional representation

Embedding Data – word2vec

These lower dimensional representations can include a good deal of semantic meaning

i.e. $\text{vector}(\text{king}) - \text{vector}(\text{man}) + \text{vector}(\text{woman}) \approx \text{vector}(\text{queen})$



From Mikolov *et al* NIPS 2013

Embedding Data – word2vec

Dimensionality reduction and visualisation is key to understanding the data

Useful for your **coursework**

There are many ways to do this, with the `sklearn.manifold` library:

<https://scikit-learn.org/stable/modules/manifold.html>

