Nonlinear Embeddings

-Statistical Machine Learning-

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Lower dimesional (metric) embeddings

Spectral connectivity analysis (SCA) is a general process for finding lower dimensional structure in the data

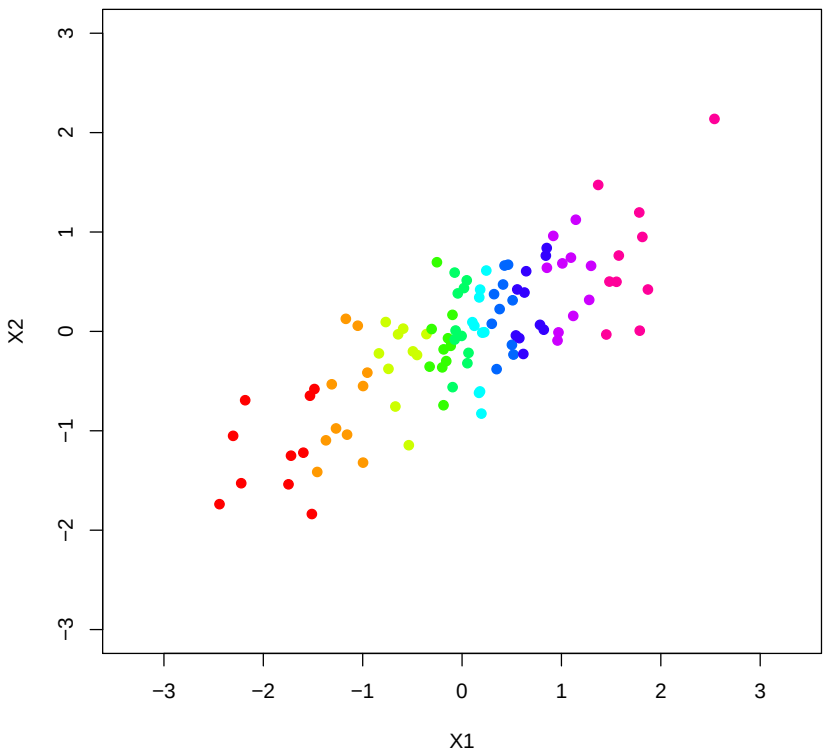
It can be...

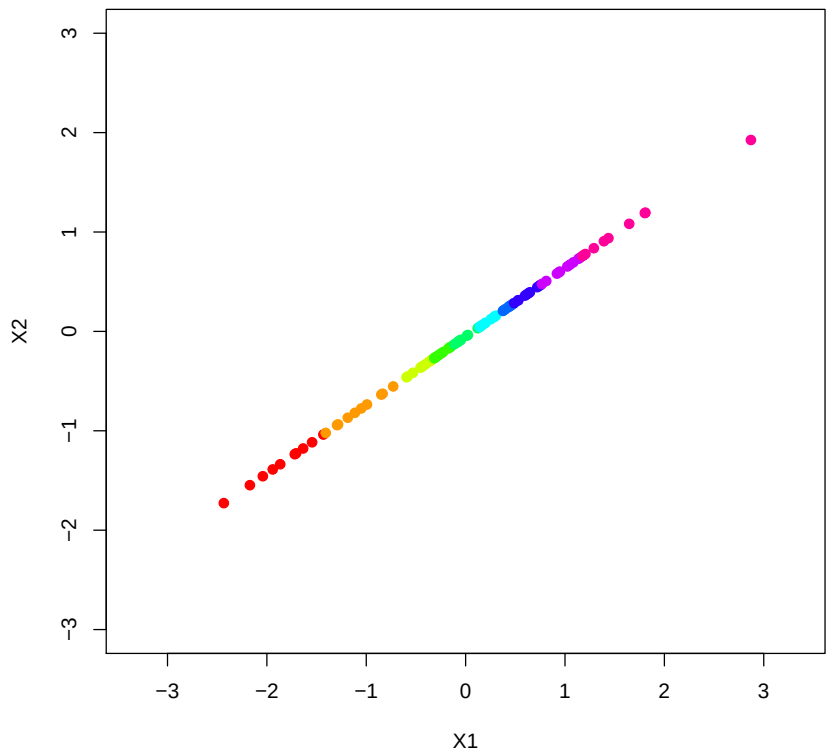
- Linear or nonlinear
- Used for dimension reduction or feature creation
- PCA, Fisher discriminant analysis, Locally linear embeddings, Hessian eigenmaps, Laplacian eigenmaps, kernel PCA
- Useful as an input to classification, clustering, and regression approaches

Let's take one last look at PCA before proceding

When PCA Works Well

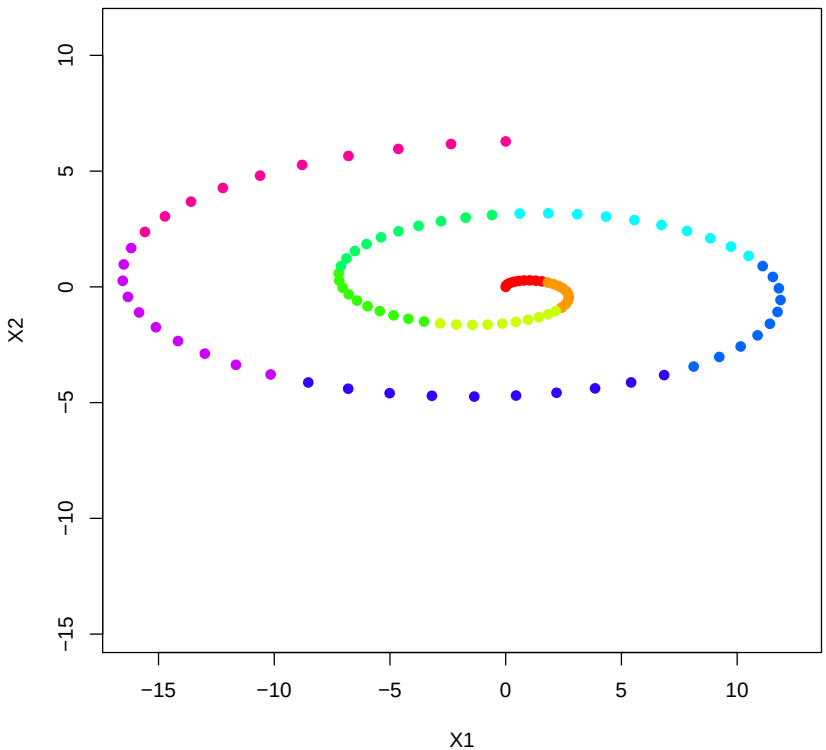
PCA can do effective dimension reduction (that is, explain most of the data with $m < p$ components) as long as the data can be efficiently represented as 'lines' (or planes, or hyperplanes). So, in two dimensions:

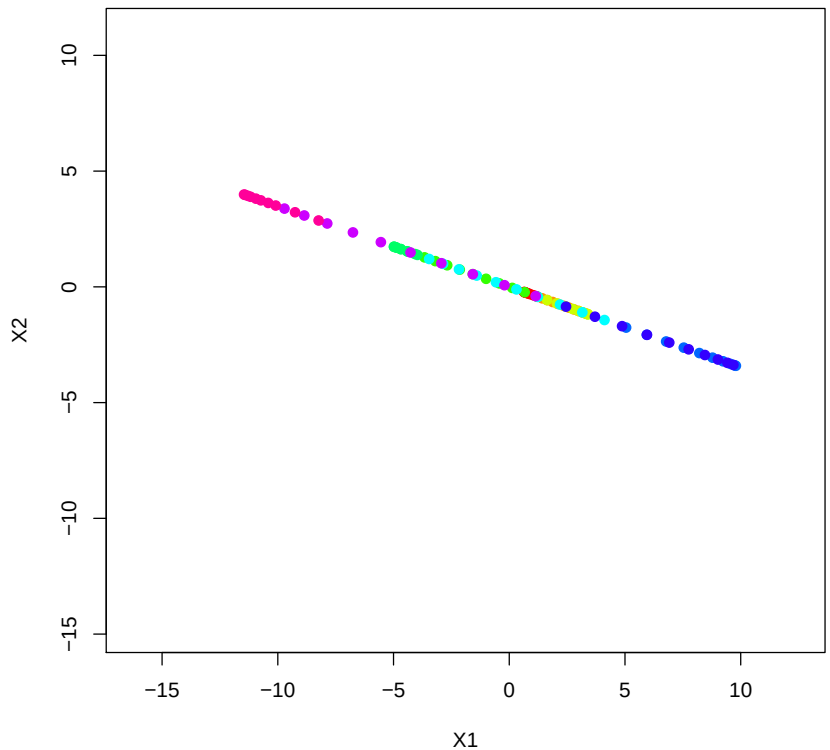




When PCA doesn't work well

What about other data structures? Again in two dimensions



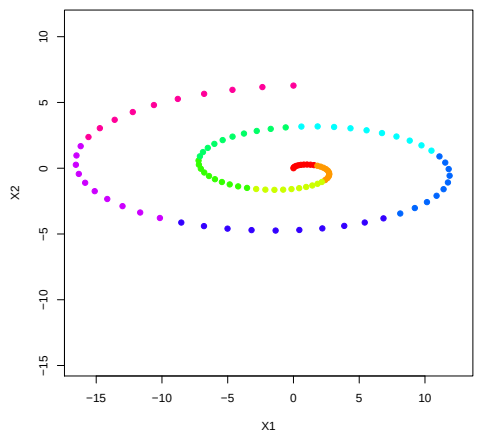


Here, we have failed miserably.

Explanation

- PCA wants to minimize distances (equivalently maximize variance). This means it slices through the data at the meatiest point, and then the next one, and so on. If the data are 'curved' this is going to induce artifacts.
- PCA also looks at things as being close if they are near each other in a Euclidean sense

  [this is essentially all covariance is].
- On the spiral, our intuition says that things are 'close' only if the distance is constrained to go along the curve. In other words, purple and blue are close, blue and red are not.



PCA and covariance

PCA: Find the directions of greatest variance. This doesn't on its face seem like it maintains correlations, but observe:

$$var([a,b]^\top X) = var(ax_1 + bx_2) = a^2 Var(x_1) + b^2 Var(x_2) + 2abCov(x_1, x_2)$$

If we standardize the matrix, then this reduces to

$$var(ax_1 + bx_2) = a^2 + b^2 + 2abCov(x_1, x_2)$$

This gets maximized over $a^2 + b^2 = 1$.

- If $Cov(x_1, x_2) \approx 0$, then this gets maximized by any $a^2 + b^2 = 1$ (it doesn't matter)
- If $Cov(x_1, x_2) \approx 1$, then this gets maximized by setting $a = b = 1/\sqrt{2}$

So, in either case, we are really maintaining correlations

Correlation is fundamentally a linear phenomenon

Graphical example of the phenomenon

```
library(mvtnorm)
sigma   = matrix(c(1,sig,sig,1),nrow=2)
nsweep  = 1000
outcome = matrix(0,nrow=nsweep,ncol=2)
for(sweep in 1:nsweep){
  x                = rmvnorm(200,c(0,0),sigma)
  out.pca          = prcomp(x,center=T,scale=F)
  outcome[sweep,] = out.pca$rotation[,1]
}
plot(outcome,xlab='PC1',ylab='PC2')
```

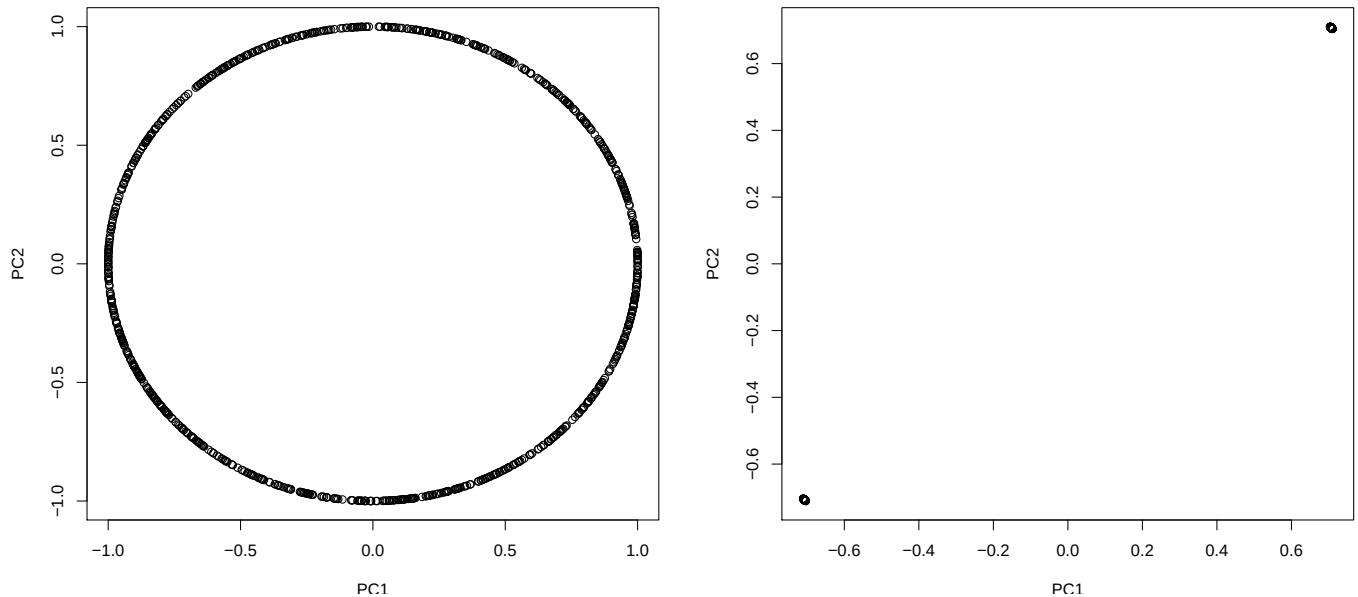


Figure: Left: `sig = 0`. Right: `sig = .999`

Nonlinear embeddings

Kernel PCA (KPCA)

Classical PCA comes from $\tilde{\mathbb{X}} = \mathbb{X} - M\mathbb{X} = UDV^\top$, where $M = \mathbf{1}\mathbf{1}^\top/n$ and $\mathbf{1} = (1, 1, \ldots, 1)^\top$

However, we can just as easily get it from the outer product

$$\mathbb{K} = \tilde{\mathbb{X}}\tilde{\mathbb{X}}^\top = (I - M)\mathbb{X}\mathbb{X}^\top(I - M) = UD^2U^\top$$

The intuition behind KPCA is that $\mathbb{K}$ is an expansion into a kernel space, where

$$\mathbb{K}_{i,i'} = k(\tilde{X}_i, \tilde{X}_{i'}) = \langle \tilde{X}_i, \tilde{X}_{i'} \rangle$$

Reminder: Anytime we see an inner product, we can kernelize it

Kernel PCA

Following this intuition, the approach is simple:

1. Specify a kernel k
   (e.g. $k(X, X') = \exp\{-\gamma^{-1} ||X - X'||_2^2\}$)
2. Form $K_{i,i'} = k(X_i, X_{i'})$
3. Standardize and get eigenvector decomposition

$$\mathbb{K} = (I - M)K(I - M) = UD^2U^\top$$

This implicitly finds the inner product:

$$k(X_i, X_{i'}) = \langle \phi(X_i), \phi(X_{i'}) \rangle$$

However, we need only specify the kernel

Kernel PCA

The scores are still $Z = UD$

The q^{th} KPCA score is (up to centering)

$$Z_{iq} = \sum_{i'=1}^{n} \beta_{i'q} k(X_i, X_{i'})$$

where $\beta_{i',q} = u_{i'q}/d_q$

Note: As we don't explicitly generate the feature map, there are no loadings

Small detour

Reproducing kernel Hilbert space

Reminder: Mercer's theorem assures us that

$$k(X, X') = \sum_{j=1}^{\infty} \theta_j \phi_j(X) \phi_j(X')$$

Here, the system $(\phi_j)_{j=1}^{\infty}$ spans a space $\mathcal{H}_k$

The function space $\mathcal{H}_k$ is known as a reproducing kernel Hilbert space (RKHS)

It can also be thought of as roughly

$$\mathcal{H}_k = \{f : f(X) = \sum_{i=1}^{n} \beta_i k(X, X_i)\}$$

Which has a special inner product

$$\langle f, f \rangle_{\mathcal{H}_k} = ||f||^2_{\mathcal{H}_k} = \sum_{j=1}^{\infty} f_j^2 / \theta_j < \infty$$

Reproducing kernel Hilbert space

Writing

$$f(X) = \sum_{i=1}^{n} \beta_i k(X, X_i)$$

The terms $k(X, X_i)$ are the representers, as

$$\langle k(\cdot, X), f \rangle_{\mathcal{H}_k} = f(X)$$

and $\mathcal{H}_k$ is called a reproducing kernel Hilbert space (RKHS) as

$$\langle k(\cdot, X), k(\cdot, X') \rangle_{\mathcal{H}_k} = k(X, X')$$

Note: For kernel methods, we are generalizing the finite dimensional Euclidean inner product

$$\langle X, X' \rangle = X^\top X'$$

Kernel methods via regularization

After specifying a kernel function k, we can define an estimator via

$$\min_{f \in \mathcal{H}_k} \hat{\mathbb{P}} \ell_f + \lambda ||f||^2_{\mathcal{H}_k}$$

This is a (potentially) infinite dimensional optimization problem

(hard, especially with a computer)

It can be shown that the solution has the form

$$\hat{f}(X) = \sum_{i=1}^{n} \beta_i k(X, X_i)$$

(This is known as the representer theorem)

Back to KPCA

Kernel PCA

Reminder: To get the first PC in classical PCA, we want to solve

$$\max_{\alpha} \mathbb{V}\alpha^\top X \quad \text{subject to} \quad ||\alpha||_2^2 = 1$$

Translate this into the kernel setting, and we are trying to solve

$$\max_{g \in \mathcal{H}_k} \mathbb{V}g(X) \text{ subject to } ||g||_{\mathcal{H}_k} = 1$$

The representer theorem states that a solution to this problem is

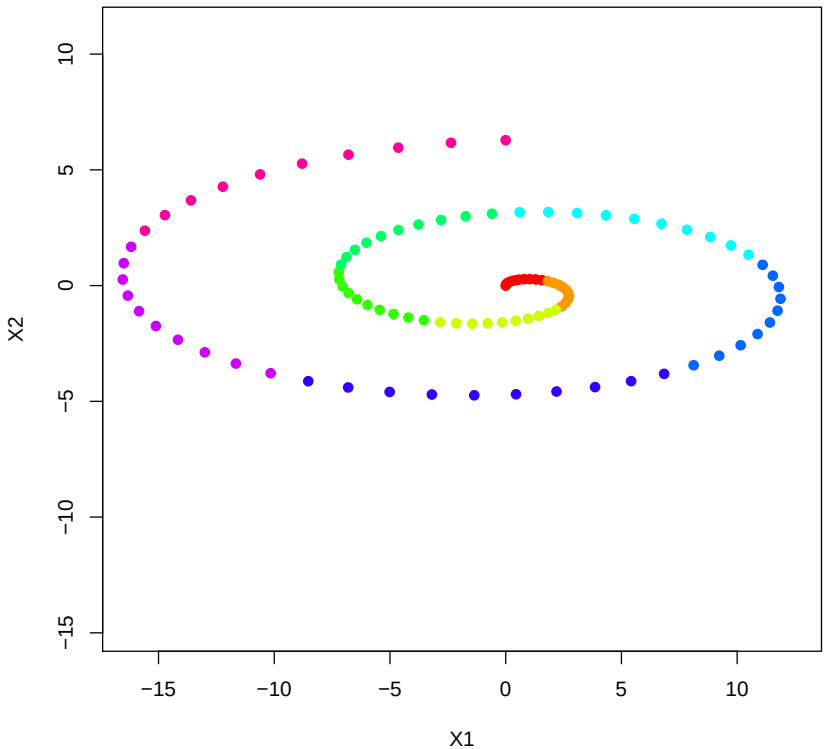
$$\hat{g}(X) = \sum_{i=1}^{n} \beta_i k(X, X_i)$$

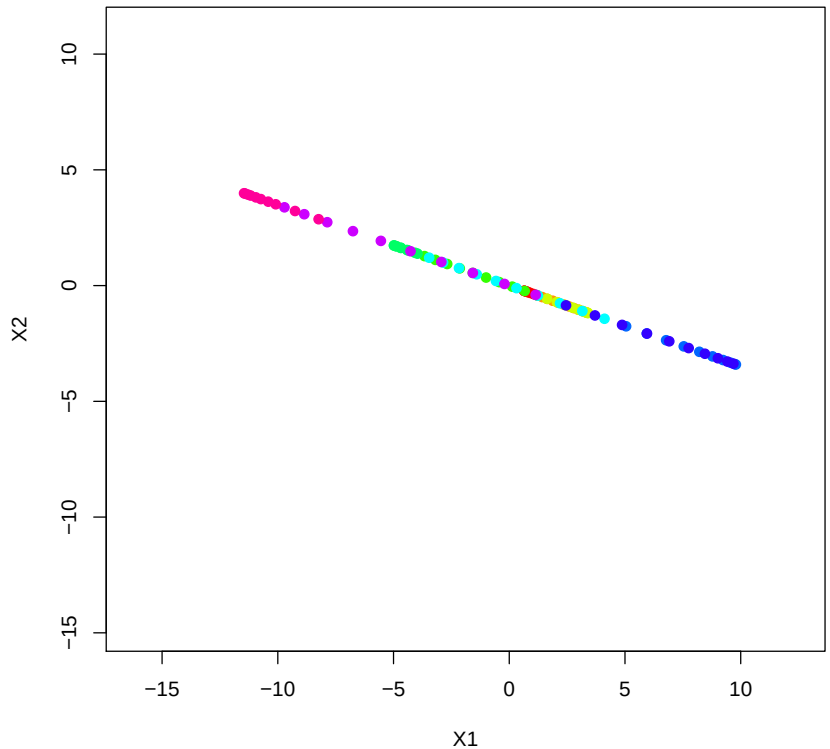
Compare

$$Z_{iq} = \sum_{i'=1}^{n} \beta_{i'q} k(X_i, X_{i'})$$

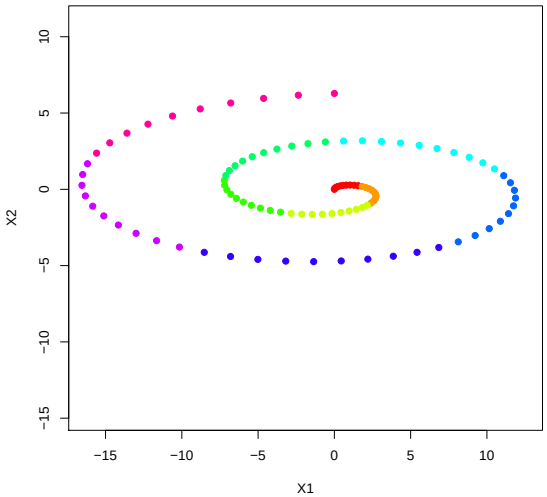
where $\beta_{i',q} = u_{i'q}/d_q$

KPCA: some results

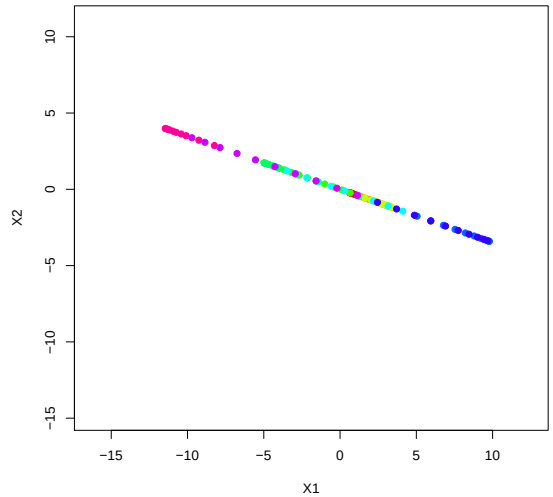




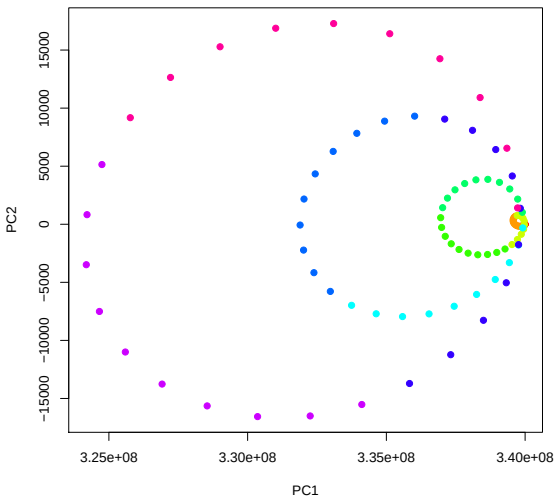
KPCA: SOME RESULTS



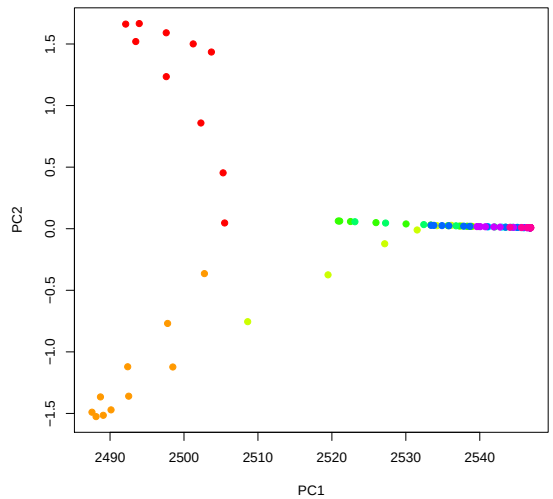
Data



PCA



KPCA: polynomial(2)



KPCA: gaussian(10)

Semisupervised learning in practice

Looking at:

$$Z_{iq} = \sum_{i'=1}^{n} \beta_{i'q} k(X_i, X_{i'})$$

this is only defined at our observed features

Write

- $\mathcal{D}_{train} = \{(X_1, Y_1), \ldots, (X_n, Y_n)\}$
- $\mathcal{D}_{test} = \{(X_1^*, Y_1^*), \ldots, (X_{n^*}^*, Y_{n^*}^*)\}$

Two common scenarios are

1. We are given $\mathcal{D}_{train}$ and $X_1^*, \ldots, X_{n^*}^*$ to build $\hat{f}$
2. We are given only $\mathcal{D}_{train}$ to build $\hat{f}$

Case 1

We are given $\mathcal{D}_{train}$ and $X_1^*, \ldots, X_{n^*}^*$ to build $\hat{f}$

Then we can just use straight forward KPCA

(Or any unsupervised learning step)

1. Form $\mathbb{K}$ on $\mathcal{D}_{train}$ and $X_1^*, \ldots, X_{n^*}^*$
2. Get UD
3. Pass $Z_q = UD[, 1 : q]$ to train $\hat{f}$
4. Get $\hat{Y} = \hat{f}(Z_q)$

Case 2

We are given only $\mathcal{D}_{train}$ to build $\hat{f}$

Now, we don't know the coordinates of $X_1^*, \ldots, X_{n^*}^*$ in the representation space

To get a new observation X^* embedded into this representation:

$$Z_0 = D^{-1}U^\top(I - M)[k^* - K\mathbf{1}/n]$$

where $k^* = [k(X^*, X_1), \ldots, k(X^*, X_n)]^\top$

Then we compute:

1. Form $\mathbb{K}$ on $\mathcal{D}_{train}$
2. Get UD
3. Pass $Z_q = UD[, 1 : q]$ to train $\hat{f}$
4. Form Z_q^* for all $X_1^*, \ldots, X_{n^*}^*$
5. Get $\hat{Y}_{test} = \hat{f}(Z_q^*)$

KPCA: SUMMARY

Kernel PCA seeks to generalize the notion of similarity using a kernel map

This can be interpreted as finding smooth, orthogonal directions in a RKHS

This can allow us to start picking up nonlinear (in the original feature space) aspects of our data

This new representation can be passed to a supervised method to form a semisupervised learner

Laplacian Eigenmaps

In order to use the intuitive distance, we need to know the geometry of the data. This needs to be estimated.

We can get an estimate of the distance in the unknown geometry that the data come from (known as a manifold) by altering the usual Euclidean distance.

Some notes:

- The name Laplacian Eigenmaps comes from getting the eigenvector decomposition of the Laplacian restricted to the manifold (which is the second derivative version of the gradient).
- If the manifold is smooth, then local Euclidean distance is an approximation to the distance on the manifold.

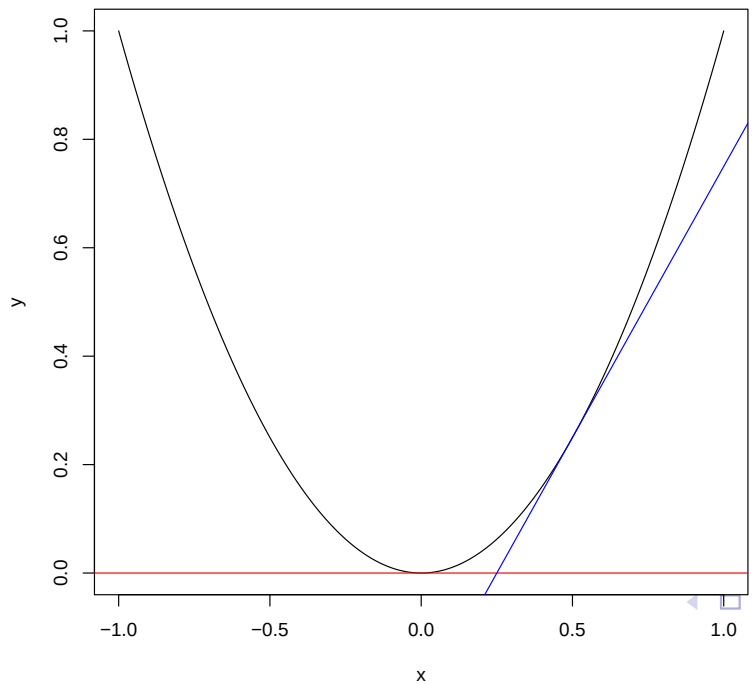
What is a manifold?

How good of an approximation is Euclidean distance?

This question is equivalent to how asking: how quickly does the tangent (space) change?

In 1-D, the tangent space is just the first derivative at that point:

$$f(x) = x^2 \Rightarrow f'(x) = 2x.$$



What is a Manifold?

Therefore, the quality of the (local) Euclidean distance, depends on the second derivative

(ie: how fast does the first derivative change?)

In higher dimensions, the second derivative is known as the Laplacian:

$$\sum_j \frac{\partial^2 f}{\partial x_j^2}$$

(Note: This is also known as the divergence of the gradient)

What are Laplacian Eigenmaps, then?

Imagine the operator $\mathbb{L}$ that performs this operation:

$$\mathbb{L}f = \sum_j \frac{\partial^2 f}{\partial x_j^2}$$

Then $\mathbb{L}$ is the Laplacian, mapping a function to the divergence of its gradient

Key Idea: We can get the eigenvectors/eigenvalues of $\mathbb{L}$. Analogously to PCA, we can now do inference with these eigenvectors.

Note: There is a substantial overlap with KPCA, the difference being the centering of K and the row sum versus column sum normalization

Laplacian Eigenmaps

Collect data: $X_1, \ldots, X_n$ where $X_i \in \mathbb{R}^p$.

1. Form the distance matrix $\Delta_{ij} = ||X_i - X_j||_2^2$.
2. Compute
$$\mathbb{K} = \exp\left(-\frac{\Delta}{\gamma}\right)$$
3. Form the Laplacian $\mathbb{L} = \mathbb{I} - \mathbb{M}^{-1}\mathbb{K}$,
$$\mathbb{M} = \texttt{diag}(\texttt{rowSums}(\mathbb{K}))$$
4. Compute the spectrum: $\mathbb{L} = U\Sigma U^\top$.
5. Return U_d, where U_d corresponds to the smallest d (nontrivial) eigenvalues of $\mathbb{L}$

   (Note that the eigenvectors of $\mathbb{L}$ and $\mathbb{M}^{-1}\mathbb{K}$ are the same but the order of the eigenvalues are reversed)

Deeper investigation

1. Form the distance matrix Δ.

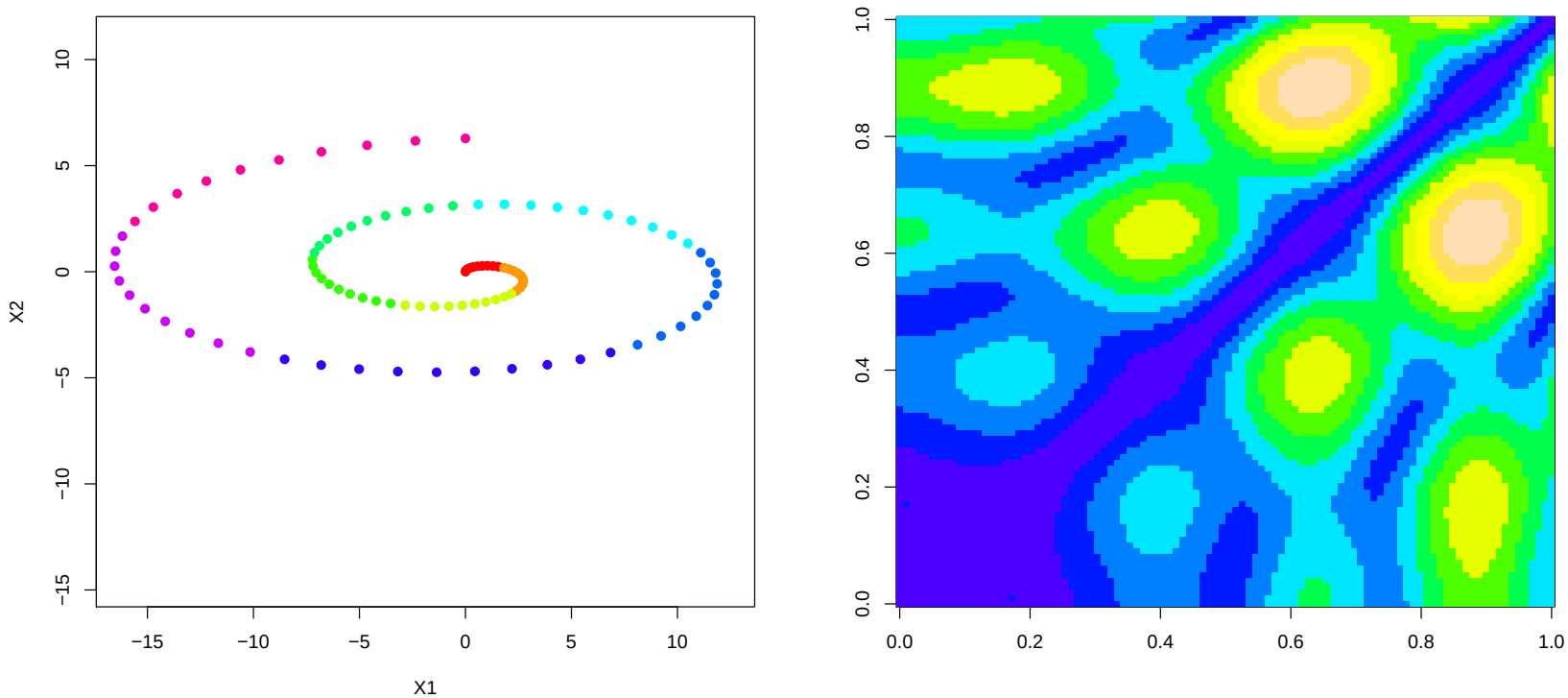


Figure: If we think about the center as 0 and the last blue circle as 1, then each entry the plot on the Right is the Euclidean distance between each data point on the plot on the Left (that is, Δ). The color on the Right plot goes from purple (small distance) to beige/pink (large distance).

Deeper investigation

```
Delta = as.matrix(dist(X,diag=TRUE,upper=TRUE))

image(Delta,col=topo.colors(10))
```

Deeper investigation

2. Exponentiate $-\Delta/\gamma$ to form $\mathbb{K}$ for some fixed γ.

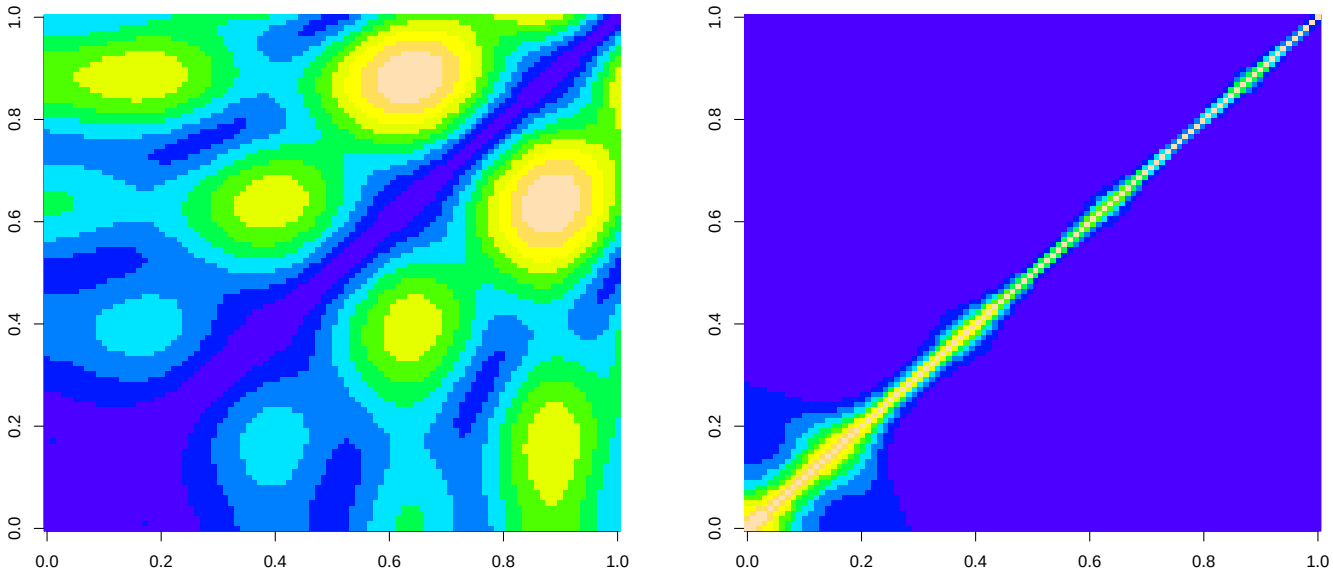
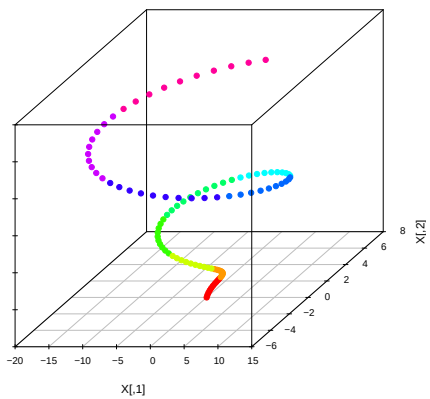


Figure: The Left plot is Δ and the Right plot is $\mathbb{K}$ for $\gamma = 0.95$.

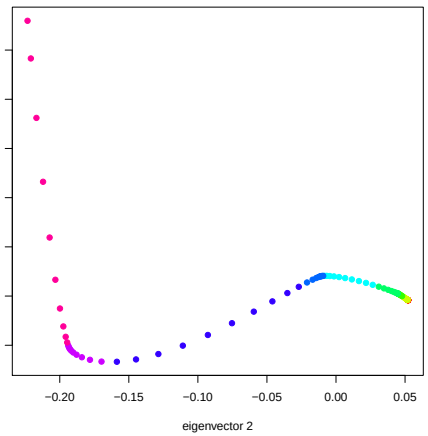
Deeper investigation

```
gamma = 0.95
Wgamma = exp(-Delta/gamma)
image(Wgamma,col=topo.colors(10))
```

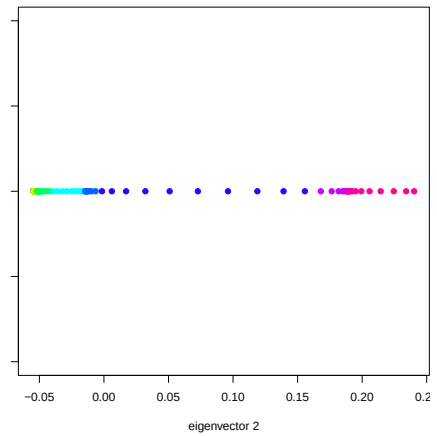
Spiral in $\mathbb{R}^3$



Original data

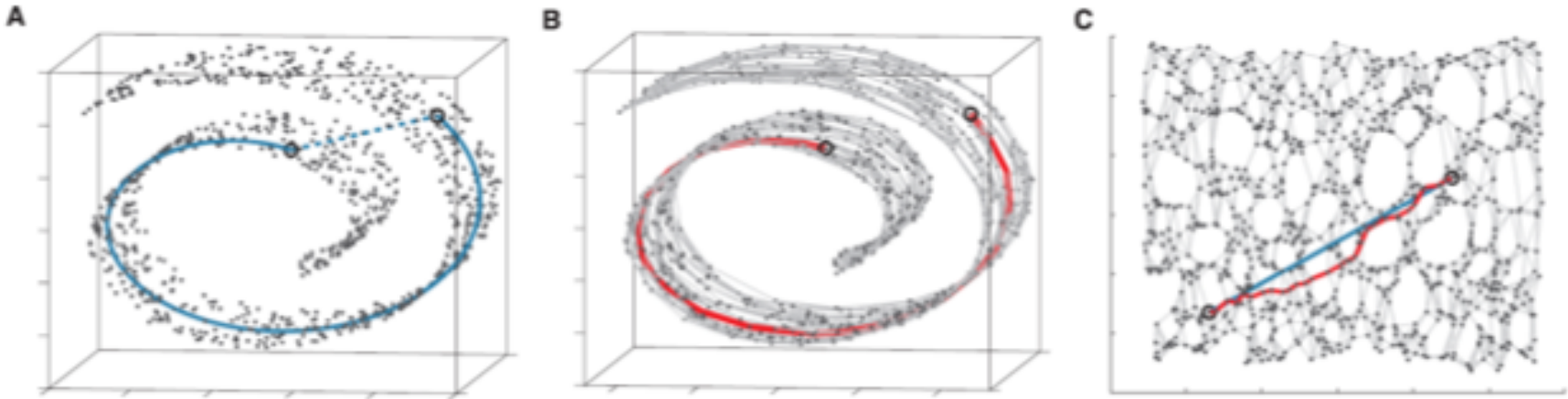


1^{st} & 2^{nd} nontrivial eigenvectors



1-dimensional

Local Euclidean distance approximates the geodesic



The red line is the local Euclidean path between the two points, while the blue line is the path along the manifold.

James, Witten, Hastie, Tibshirani (2013)