

Advanced Data Analysis: Kernel PCA

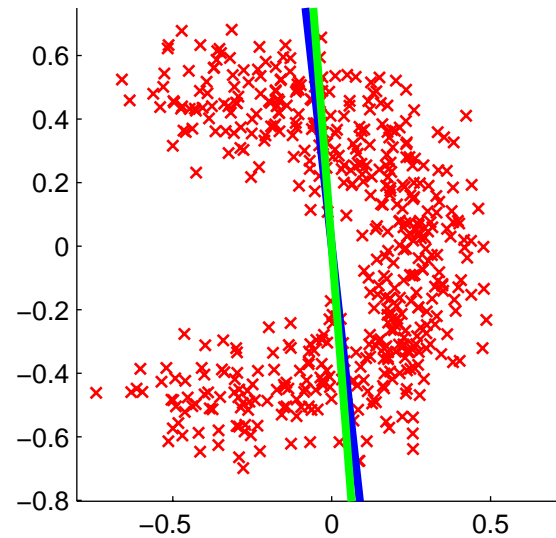
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Data with Curved Structures

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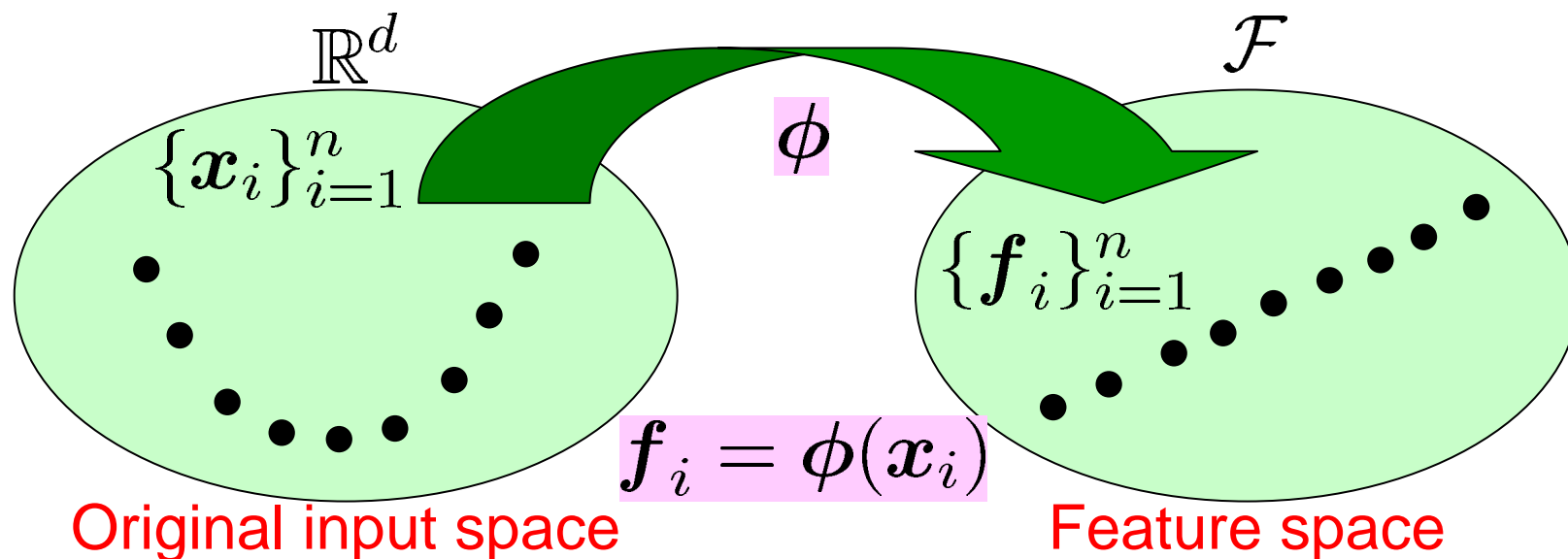


- If the data cloud is bent, **any linear methods** can not find the **curved** structure.

➡ Limitation of linear method!

Non-Linearizing Linear Methods⁸⁴

- A simple non-linear extension of linear methods while keeping computational advantages of linear methods:
 - Map the original data to a **feature space** by a **non-linear transformation**
 - Run **linear algorithm** in the feature space

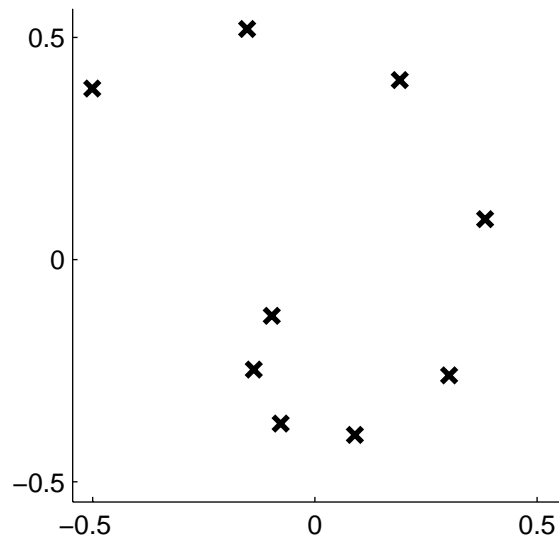


Example

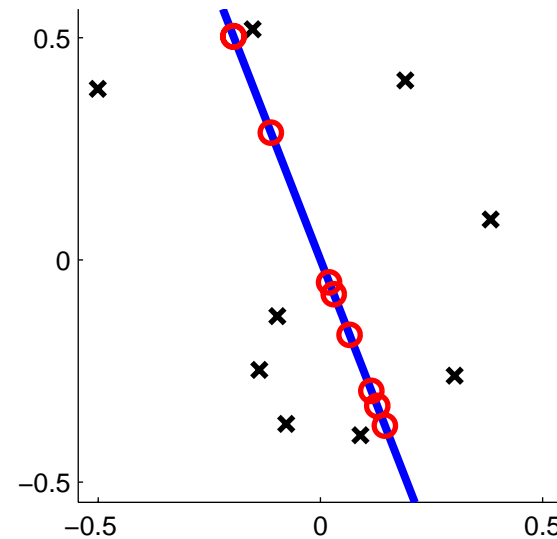
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■ $d = 2$

Centered data
in input space



Linear PCA



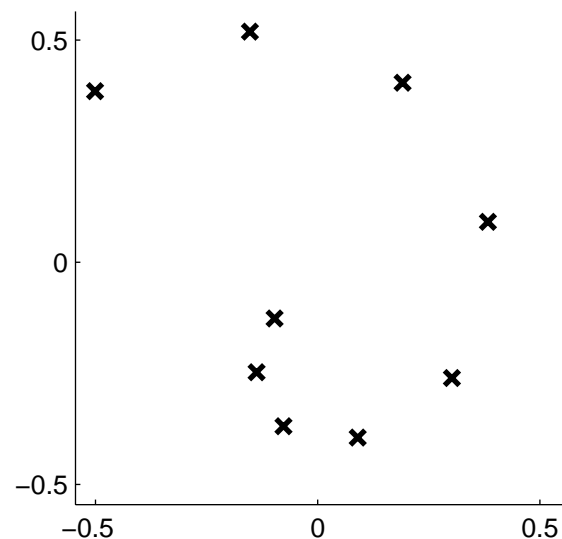
Example (cont.)

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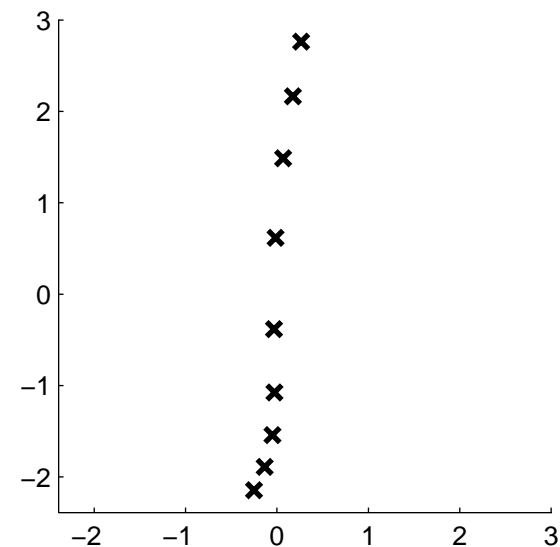
■ Polar coordinate:

$$\mathbf{x} = \begin{pmatrix} a \\ b \end{pmatrix} \longrightarrow \mathbf{f} = \begin{pmatrix} r \cos \theta \\ r \sin \theta \end{pmatrix}$$

Centered data
in input space



Centered data
in feature space

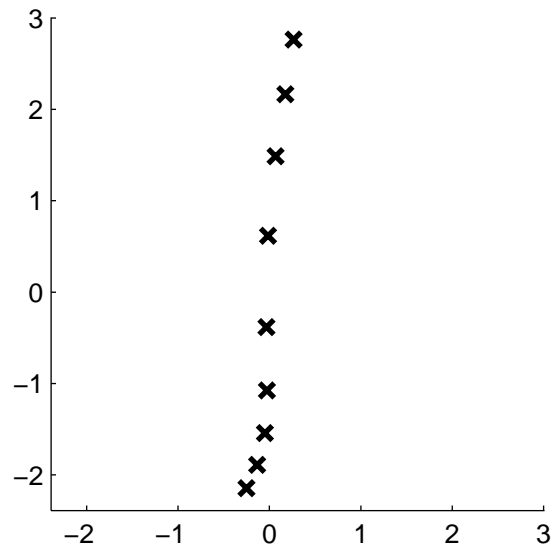


Example (cont.)

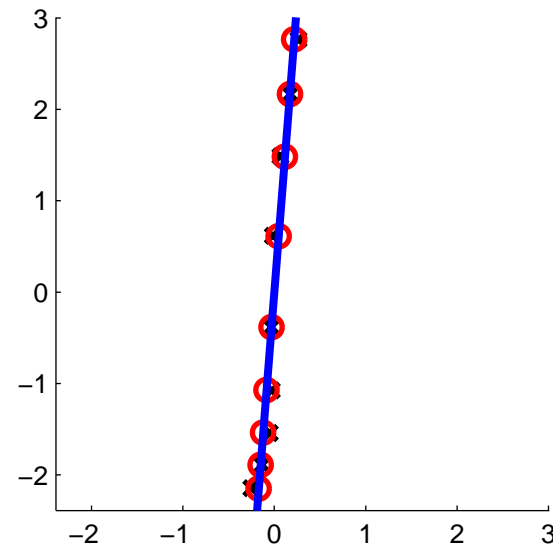
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- Run PCA in feature space.

Centered data
in feature space



PCA projection
in feature space

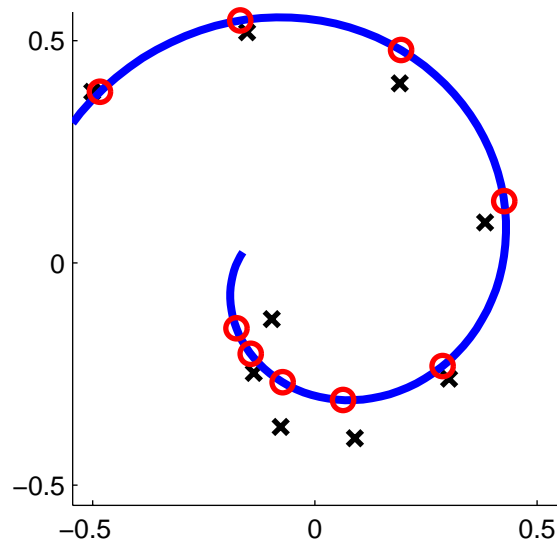


Example (cont.)

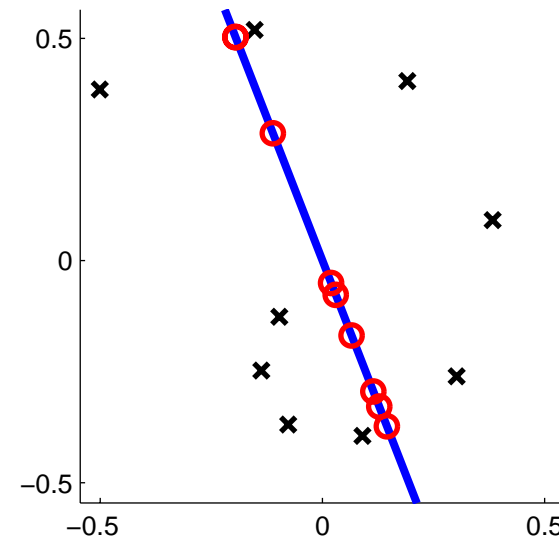
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- Pull the results back to input space.

Non-linear PCA



Linear PCA



- Non-linear PCA describes the original data much better than linear PCA.

Centering in Feature Space

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- PCA requires centered samples, thus we need to center samples by

$$\bar{f}_i = f_i - \frac{1}{n} \sum_{j=1}^n f_j$$

- In matrix form,

$$\bar{F} = FH$$

$$F = (f_1 | f_2 | \cdots | f_n)$$

$$\bar{F} = (\bar{f}_1 | \bar{f}_2 | \cdots | \bar{f}_n)$$

$$H = I_n - \frac{1}{n} \mathbf{1}_{n \times n}$$

I_n : n -dimensional identity matrix

$\mathbf{1}_{n \times n}$: $n \times n$ matrix with all ones

PCA in Feature Space

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$$\overline{C}\psi = \lambda\psi$$

$$\overline{C} = \overline{F} \overline{F}^\top$$

■ PCA solution:

$$\mathbf{B}_{PCA} = (\psi_1 | \psi_2 | \cdots | \psi_m)^\top$$

- $\{\lambda_i, \psi_i\}_{i=1}^m$: Sorted eigenvalues and normalized eigenvectors of $\overline{C}\psi = \lambda\psi$

$$\langle \psi_i, \psi_j \rangle = \delta_{i,j}$$

$$\lambda_1 \geq \lambda_2 \geq \cdots \geq \lambda_\mu$$

■ PCA embedding of a sample f :

$$\overline{g} = \mathbf{B}_{PCA} \left(f - \frac{1}{n} \mathbf{F} \mathbf{1}_n \right)$$

$$\mu = \dim(\mathcal{F})$$

$\mathbf{1}_n$: n -dimensional vector with all ones

PCA in High-Dimensional Feature Space

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$$\mu = \dim(\mathcal{F})$$

- If μ is high,
 - Description ability of non-linear PCA will increase.
 - However, computational cost increases since the dimension of $\overline{\mathcal{C}}$ is μ .
- We may be able to reduce the computational cost since

$$\text{rank}(\overline{\mathcal{C}}) = \min(\mu, n) \leq \mu$$

$$\overline{\mathcal{C}} = \overline{\mathbf{F}} \overline{\mathbf{F}}^\top$$

$$\overline{\mathbf{F}} = (\overline{\mathbf{f}}_1 | \overline{\mathbf{f}}_2 | \cdots | \overline{\mathbf{f}}_n)$$

Dual Formulation

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$$(A) \quad \overline{C}\psi = \lambda\psi$$

$$\overline{C} = \overline{F} \overline{F}^\top$$

$$(B) \quad \overline{K}\alpha = \lambda\alpha$$

$$\overline{K} = \overline{F}^\top \overline{F}$$

■ Solution of (A) can be obtained from (B).

• **Proof:** If α is a solution of (B), it holds that

$$\overline{C} \overline{F}\alpha = \overline{F}\overline{F}^\top \overline{F}\alpha = \overline{F}\overline{K}\alpha = \lambda\overline{F}\alpha$$

This implies that $\psi = \overline{F}\alpha$ is a solution of (A).

■ **Note:** solution of (B) can also be obtained from (A).

■ Given \overline{K} , solving (B) is faster than (A) when $\mu > n$ since

$$\text{rank}(\overline{C}) = n < \mu$$

Renormalization of Eigenvectors⁹³

$$\overline{K}\alpha = \lambda\alpha$$

- Standard eigensolvers output an orthonormal eigenvectors.

$$\langle \alpha_i, \alpha_j \rangle = \delta_{i,j}$$

- However, PCA requires the primal eigenvectors $\{\psi_i\}_{i=1}^m$ to be orthonormal.

- Since $\langle \psi_i, \psi_j \rangle = \langle \overline{K}\alpha_i, \alpha_j \rangle = \lambda_i \delta_{i,j}$, we need to renormalize $\{\psi_i\}_{i=1}^m$ by

$$\psi_i \leftarrow \frac{\psi_i}{\|\psi_i\|} = \frac{1}{\sqrt{\lambda_i}} \overline{F}\alpha_i$$

$$\psi_i = \overline{F}\alpha_i$$

$$\overline{K}\alpha_i = \lambda_i\alpha_i$$

PCA in Feature Space (Dual) ⁹⁴

■ PCA embedding of a sample f :

$$\bar{g} = \Lambda^{-\frac{1}{2}} A^\top H \left(k - \frac{1}{n} K \mathbf{1}_n \right) \quad (\text{Homework})$$

- $\{\lambda_i, \alpha_i\}_{i=1}^m$: Sorted eigenvalues and normalized eigenvectors of $\bar{K} \alpha = \lambda \alpha$

$$\lambda_1 \geq \lambda_2 \geq \dots \geq \lambda_n \quad \langle \alpha_i, \alpha_j \rangle = \delta_{i,j}$$

$$\Lambda = \text{diag}(\lambda_1, \lambda_2, \dots, \lambda_m)$$

$$A = (\alpha_1 | \alpha_2 | \dots | \alpha_m)$$

$$\bar{K} = H K H \quad K = F^\top F$$

$$H = I_n - \frac{1}{n} \mathbf{1}_{n \times n} \quad k = F^\top f$$

I_n : n -dimensional identity matrix

$\mathbf{1}_{n \times n}$: $n \times n$ matrix with all ones

$\mathbf{1}_n$: n -dimensional vector with all ones

PCA in Feature Space (Dual) ⁹⁵

$$\mu = \dim(\mathcal{F})$$

- In the dual formulation, the computational complexity depends not on μ but only on n , if K and k are given.
- However, the computation of K and k still depends on μ .

$$K = F^\top F$$

$$k = f^\top F$$

- **Note:** K and k depend on μ only through the inner product between samples.

$$K_{i,j} = \langle f_i, f_j \rangle$$

$$k_i = \langle f, f_i \rangle$$

- For some transformation $\phi(x)$ ($= f$), there exists a bivariate function $K(x, x')$ such that

$$K_{i,j} = \langle f_i, f_j \rangle = K(x_i, x_j)$$

- Such implicit mapping $\phi(x)$ exists if
 - K is symmetric: $K^\top = K$
 - K is positive semi-definite: $\forall y, \langle Ky, y \rangle \geq 0$
- Such $K(x, x')$ is called the **reproducing kernel**.
- Rather than directly defining $\phi(x)$, we implicitly specify $\phi(x)$ by a reproducing kernel.

Examples of Kernels

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■ Polynomial kernel:

$$\mu = \dim(\mathcal{F})$$

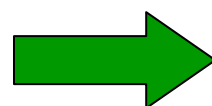
$$K(\mathbf{x}, \mathbf{x}') = \langle \mathbf{x}, \mathbf{x}' \rangle^c \quad c \in \mathbb{N}$$

- When $d = 2$ and $c = 2$,

$$\langle \mathbf{x}, \mathbf{x}' \rangle^2 = (ss' + tt')^2$$

$$\mathbf{x} = \begin{pmatrix} s \\ t \end{pmatrix}$$

$$= sss's' + 2ss'tt' + ttt't'$$



$$\mathbf{f}_i = \phi(\mathbf{x}_i) = \begin{pmatrix} s^2 \\ \sqrt{2}st \\ t^2 \end{pmatrix}$$

$$\mu = 3$$

- In general,

$$\mu = {}_{c+d-1}C_c$$

Examples of Kernels (cont.)

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■ Gaussian kernel:

$$K(\mathbf{x}, \mathbf{x}') = \exp(-\|\mathbf{x} - \mathbf{x}'\|^2 / c^2)$$

$$c > 0$$

Note: $\mu = \infty$!

$$\mu = \dim(\mathcal{F})$$

Kernel PCA: Summary

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■ Kernel PCA embedding of a sample f is

$$\bar{g} = \Lambda^{-\frac{1}{2}} A^\top H \left(k - \frac{1}{n} K \mathbf{1}_n \right)$$

- $\{\lambda_i, \alpha_i\}_{i=1}^m$: Sorted eigenvalues and normalized eigenvectors of $H K H \alpha = \lambda \alpha$

$$\lambda_1 \geq \lambda_2 \geq \dots \geq \lambda_n$$

$$\langle \alpha_i, \alpha_j \rangle = \delta_{i,j}$$

$$\Lambda = \text{diag}(\lambda_1, \lambda_2, \dots, \lambda_m)$$

$$A = (\alpha_1 | \alpha_2 | \dots | \alpha_m)$$

$$H = I_n - \frac{1}{n} \mathbf{1}_{n \times n}$$

$$k = (K(x, x_1), K(x, x_2), \dots, K(x, x_n))^\top$$

I_n : n -dimensional identity matrix

$\mathbf{1}_{n \times n}$: $n \times n$ matrix with all ones

$\mathbf{1}_n$: n -dimensional vector with all ones

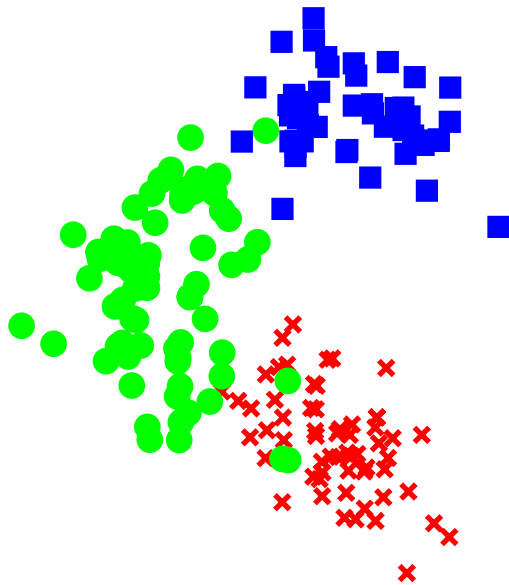
$$K_{i,j} = K(x_i, x_j)$$

Examples

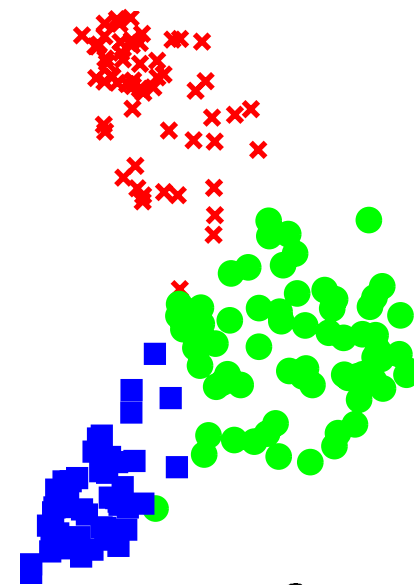
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- Wine data (UCI): 13-dim, 178 samples

$$K(x, x') = \exp(-\|x - x'\|^2 / c^2)$$



Linear PCA



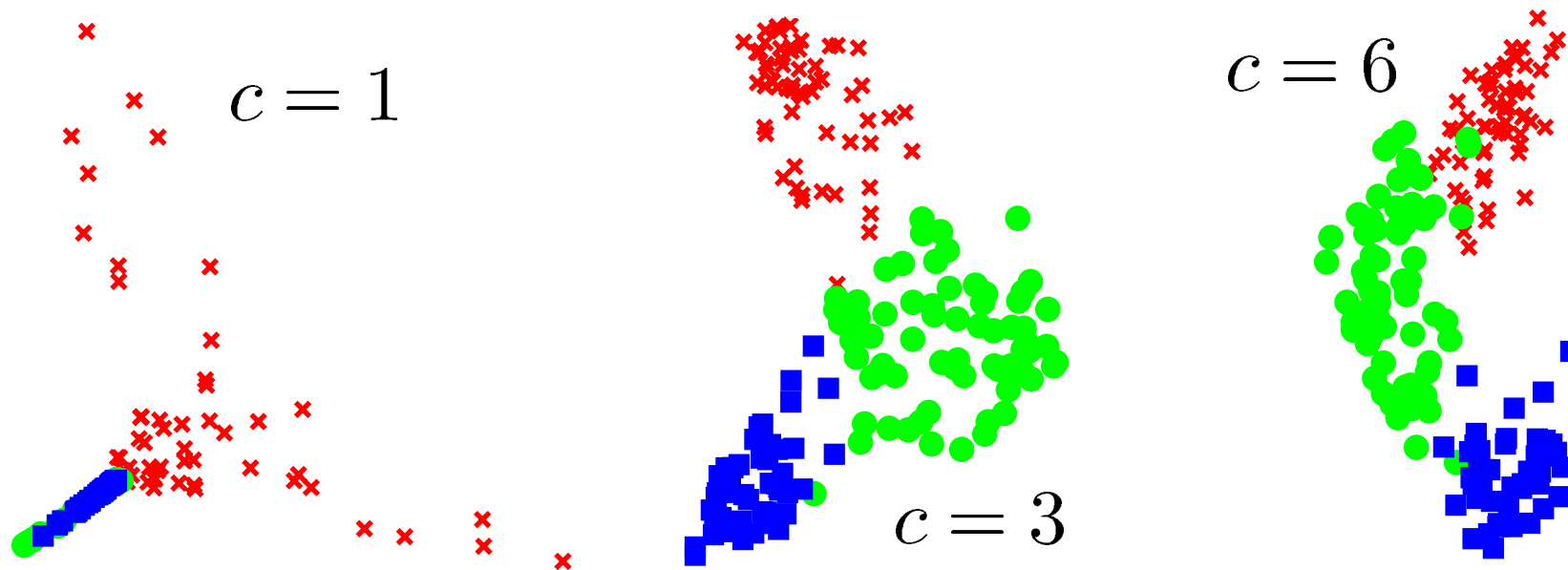
$c = 3$

Gaussian KPCA

Examples (cont.)

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$$K(\mathbf{x}, \mathbf{x}') = \exp(-\|\mathbf{x} - \mathbf{x}'\|^2 / c^2)$$



- Choice of kernels (type and parameter) depends on the result.
- Appropriately choosing kernels is not straightforward in practice.

Homework

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1. Implement kernel PCA with Gaussian kernels and reproduce the embedding result of the Wine data set.

<http://sugiyama-www.cs.titech.ac.jp/~sugi/data/DataAnalysis>

Test kernel PCA with your own (artificial or real) data and analyze the characteristics of kernel PCA.

2. Prove that kernel PCA embedding of a sample f is given by

$$\bar{g} = \Lambda^{-\frac{1}{2}} A^{\top} H \left(k - \frac{1}{n} K \mathbf{1}_n \right)$$

Suggestion

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- Read the following article for the next class:
 - M. Belkin & P. Niyogi: Laplacian eigenmaps for dimensionality reduction and data representation, Neural Computation, 15(6), 1373-1396, 2003.

<http://neco.mitpress.org/cgi/reprint/15/6/1373.pdf>