



# Computer Vision

## *---Face Recognition Using Principal Components Analysis (PCA)*

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2020.10.30

# Face recognition

☞ What makes face detection and recognition hard?



Hilary



NCC, SSD,.....?

# Face recognition

☞ What makes face detection and recognition hard?

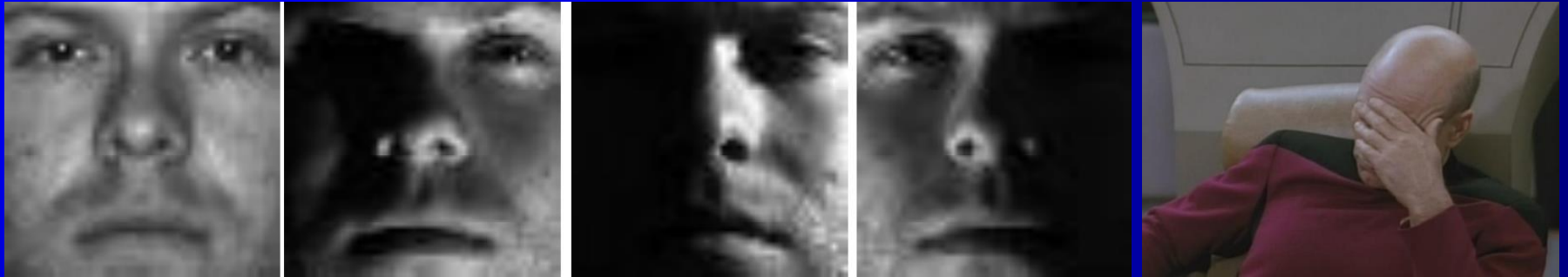


Variation in appearance: can't match a single face template and expect it to work.



# Face recognition

☞ What makes face detection and recognition hard?



Lighting

Occlusion



Viewpoint



Do this image  
contain faces?  
Where?

# Face recognition

## Simple Idea for Face Detection

- Treat each window in the image like a vector



$x$

whether  $x$  matches some  $y_j$  in the database



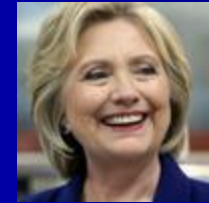
$$\text{SSD: } (y_j - x)^2$$

$$\text{Cross-correlation: } y_j \cdot x$$

NCC, zero-mean NCC...

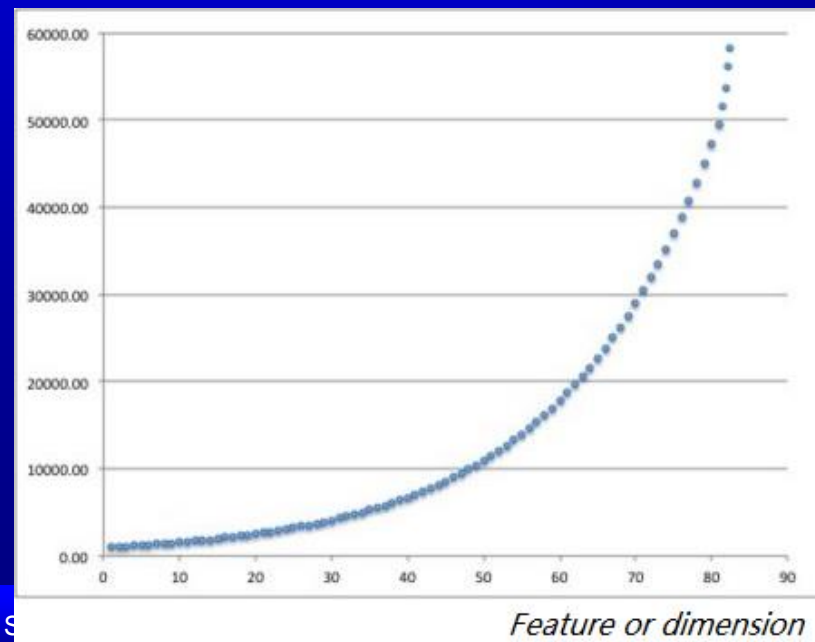
# Face recognition

- ➡ When viewed as vectors of pixel values, face images are extremely high-dimensional
  - 100x100 image = 10,000 dimensions
  - Slow and lots of storage
- ➡ But very few 10,000-dimensional vectors are valid face images.
- ➡ We want to effectively model the subspace of face images.



# Principal Component Analysis (PCA)

- ➡ Pattern recognition in high-dimensional spaces
  - Problems arise when performing recognition in a high-dimensional space (**curse of dimensionality**--维数灾难).
  - As the number of **features or dimensions** grows, the amount of data we need to generalize accurately grows **Exponentially!**



# Principal Component Analysis (PCA)

- ☞ Pattern recognition in high-dimensional spaces
- Problems arise when performing recognition in a high-dimensional space (**curse of dimensionality**--维数灾难).
  - Significant improvements can be achieved by first mapping the data into a lower-dimensional sub-space.

$$x = \begin{bmatrix} a_1 \\ a_2 \\ \dots \\ a_N \end{bmatrix} \longrightarrow \text{reduce dimensionality} \longrightarrow y = \begin{bmatrix} b_1 \\ b_2 \\ \dots \\ b_K \end{bmatrix} \quad (K \ll N)$$

- The goal of PCA is to reduce the dimensionality of the data while retaining **as much information as possible** in the original dataset.



# Principal Component Analysis (PCA)

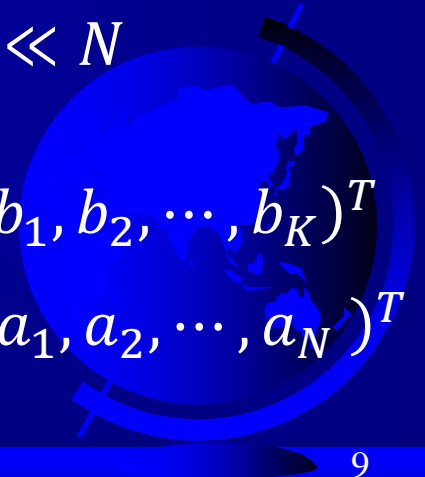
## ☞ Dimensionality reduction

- PCA allows us to compute a linear transformation that maps data from a high dimensional space to a lower dimensional sub-space.

$$y = Tx \text{ where } T = \begin{bmatrix} t_{11} & t_{12} & \dots & t_{1N} \\ t_{21} & t_{22} & \dots & t_{2N} \\ \dots & \dots & \dots & \dots \\ t_{K1} & t_{K2} & \dots & t_{KN} \end{bmatrix} \quad K \times N \quad K \ll N$$

$$\begin{aligned} b_1 &= t_{11}a_1 + t_{12}a_2 + \dots + t_{1N}a_N \\ b_2 &= t_{21}a_1 + t_{22}a_2 + \dots + t_{2N}a_N \\ &\dots \\ b_K &= t_{K1}a_1 + t_{K2}a_2 + \dots + t_{KN}a_N \end{aligned}$$

$$\begin{aligned} y &= (b_1, b_2, \dots, b_K)^T \\ x &= (a_1, a_2, \dots, a_N)^T \end{aligned}$$



# Principal Component Analysis (PCA)

- Lower dimensionality basis
  - Approximate vectors by finding a basis in an appropriate lower dimensional space.

(1) Higher-dimensional space representation:

$$x = a_1 v_1 + a_2 v_2 + \cdots + a_N v_N$$

$v_1, v_2, \dots, v_N$  is a basis of the  $N$ -dimensional space

(2) Lower-dimensional space representation:

$$\hat{x} = b_1 u_1 + b_2 u_2 + \cdots + b_K u_K$$

$u_1, u_2, \dots, u_K$  is a basis of the  $K$ -dimensional space

- Note: if both bases have the same size ( $N = K$ ), then  $x = \hat{x}$ )



# Principal Component Analysis (PCA)

- Information loss

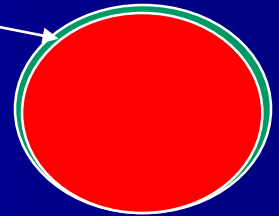
- Dimensionality reduction implies information loss!
- PCA preserves as much information as possible, that is, it minimizes the error:

$$\|x - \hat{x}\|$$

- How should we determine the best lower dimensional subspace?

*The best low-dimensional space can be determined by the "best" eigenvectors of the covariance matrix of  $x$  (i.e., the eigenvectors corresponding to the "largest" eigenvalues -- also called "principal components").*

Information loss



# Principal Component Analysis (PCA)

- Methodology
  - Suppose  $x_1, x_2, \dots, x_M$  are  $N \times 1$  vectors

$$\frac{1}{M} \left[ \begin{matrix} \vdots \\ x_1 \\ \vdots \end{matrix} + \begin{matrix} \vdots \\ x_2 \\ \vdots \end{matrix} + \dots + \begin{matrix} \vdots \\ x_M \\ \vdots \end{matrix} \right]_{N \times M} \quad \begin{matrix} \vdots \\ \bar{x} \\ \vdots \end{matrix}_{N \times 1}$$

Step1:  $\bar{x} = \frac{1}{M} \sum_{i=1}^M x_i$

Step2: subtract the mean:  $\Phi_i = x_i - \bar{x}$  (i.e. center at zero)

Step3: from the matrix  $A = [\Phi_1 \Phi_2 \dots \Phi_M]$  ( $N \times M$  matrix), then compute

$$C = \frac{1}{M} \sum_{n=1}^M \Phi_n \Phi_n^T = \frac{1}{M} A A^T$$

(sample covariance matrix,  $N \times N$ , characterizes the scatter of the data)

Step4: compute the eigenvalue of C:  $\lambda_1 > \lambda_2 > \dots > \lambda_N$   $AA^T u_i = \lambda_i u_i$

Step5: compute the eigenvectors of C:  $u_1, u_2, \dots, u_N$

# Principal Component Analysis (PCA)

- Methodology – cont.

- Since  $C$  is symmetric,  $u_1, u_2, \dots, u_N$  form a basis, (i.e., any vector  $x$  or actually  $(x - \bar{x})$ , can be written as a linear combination of the eigenvectors):

$$x - \bar{x} = b_1 u_1 + b_2 u_2 + \dots + b_N u_N = \sum_{i=1}^N b_i u_i \quad b_i = u_i^T (x - \bar{x})$$

Step 6: (dimensionality reduction step) keep only the terms corresponding to the  $K$  largest eigenvalues:

$$\hat{x} - \bar{x} = \sum_{i=1}^K b_i u_i \text{ where } K \ll N \quad u_i - \text{dimension } N$$

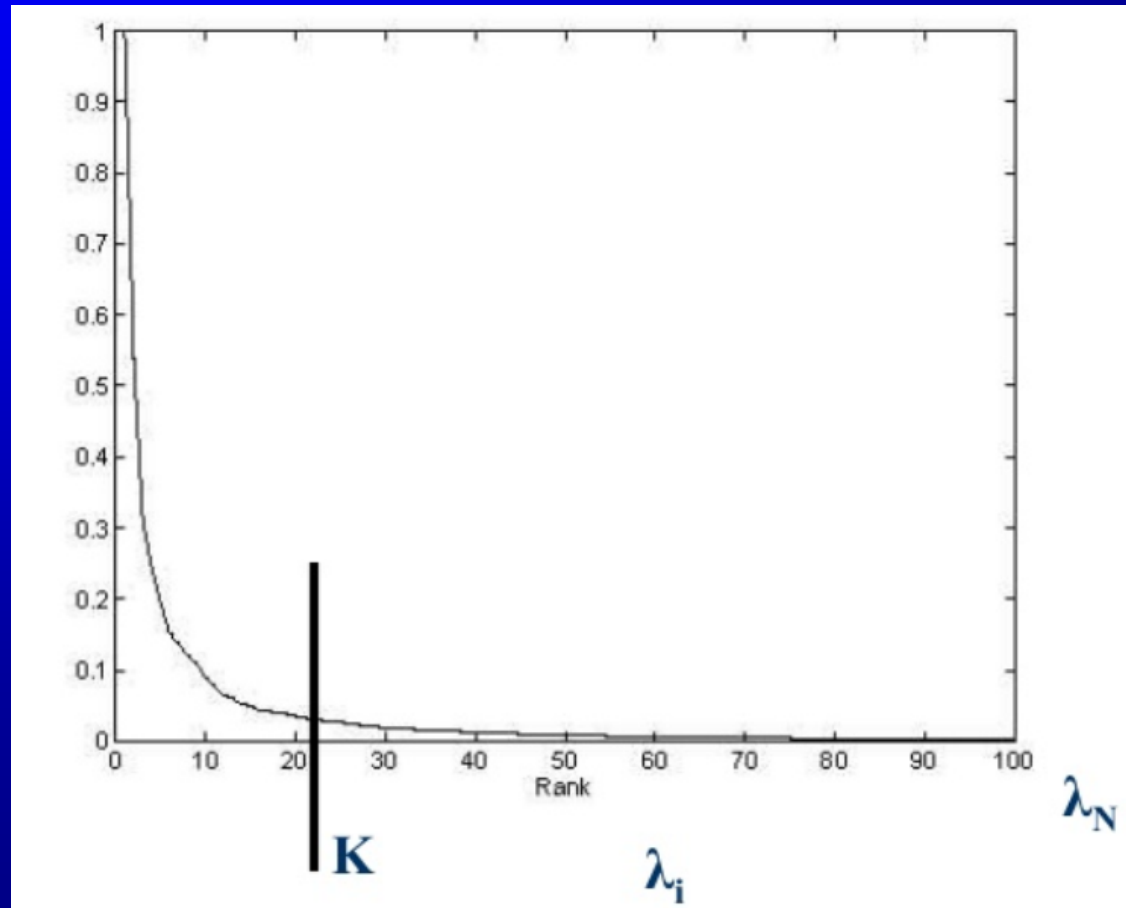
- The representation of  $\hat{x} - \bar{x}$  into the basis  $u_1, u_2, \dots, u_K$  is thus

$$\begin{bmatrix} b_1 \\ b_2 \\ \dots \\ b_K \end{bmatrix}$$



# Principal Component Analysis (PCA)

- Eigenvalue spectrum.



# Principal Component Analysis (PCA)

- Linear transformation implied by PCA
  - The linear transformation  $R^N \rightarrow R^K$  that performs the dimensionality reduction is:

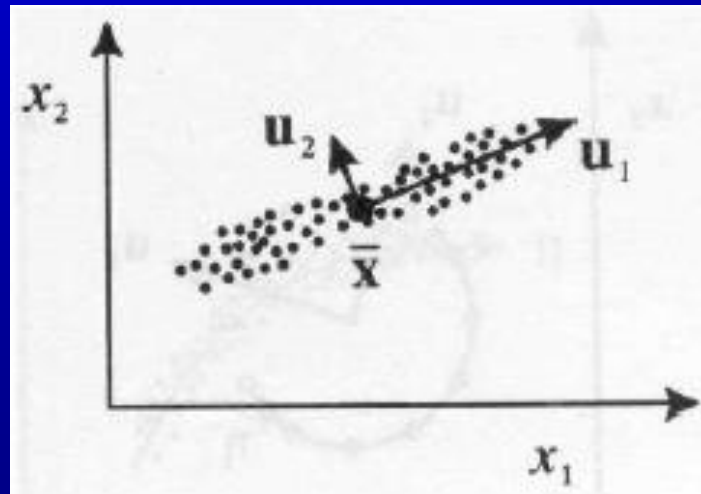
$$\begin{bmatrix} b_1 \\ b_2 \\ \dots \\ b_K \end{bmatrix} = \begin{bmatrix} u_1^T \\ u_2^T \\ \dots \\ u_K^T \end{bmatrix} (x - \bar{x}) = U^T (x - \bar{x})$$

(i.e., simply computing coefficients of linear expansion)



# Principal Component Analysis (PCA)

- Geometric interpretation
  - PCA projects the data along the directions where the data varies the most.
  - These directions are determined by the eigenvectors of the covariance matrix corresponding to the largest eigenvalues.
  - The magnitude of the eigenvalues corresponds to the variance of the data along the eigenvector directions.



# Principal Component Analysis (PCA)

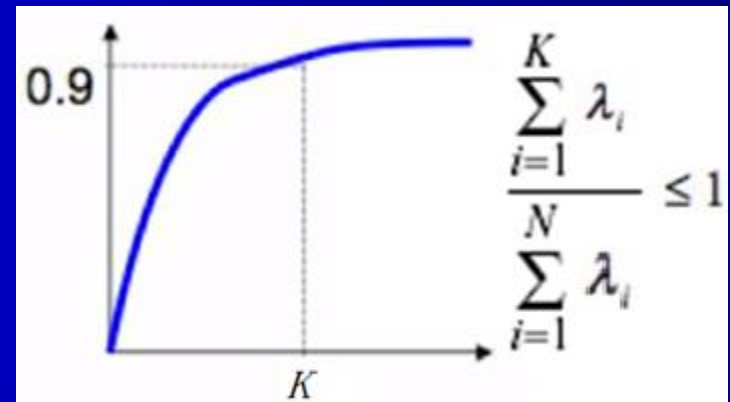
- How to choose  $K$  (i.e., number of principal components) ?
  - To choose  $K$ , use the following criterion:



# Principal Component Analysis (PCA)

- How to choose  $K$  (i.e., number of principal components) ?
  - To choose  $K$ , use the following criterion:

$$\frac{\sum_{i=1}^K \lambda_i}{\sum_{i=1}^N \lambda_i} > \text{Threshold} \quad (\text{e.g., } 0.9 \text{ or } 0.95)$$



- In this case, we say that we “preserve” 90% or 95% of the information in our data.
- If  $K=N$ , then we “preserve” 100% of the information in our data.





# Principal Component Analysis (PCA)

- What is the error due to dimensionality reduction?
  - The original vector  $x$  can be reconstructed using its principal components:

$$\hat{x} - \bar{x} = \sum_{i=1}^K b_i u_i \text{ or } \hat{x} = \sum_{i=1}^K b_i u_i + \bar{x}$$

- PCA minimizes the reconstruction error:

$$e = \|x - \hat{x}\|$$

- It can be shown that the error is equal to:

$$e = 1/2 \sum_{i=K+1}^N \lambda_i$$

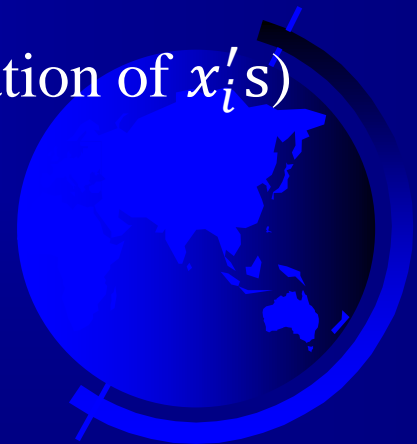


# Principal Component Analysis (PCA)

- Standardization

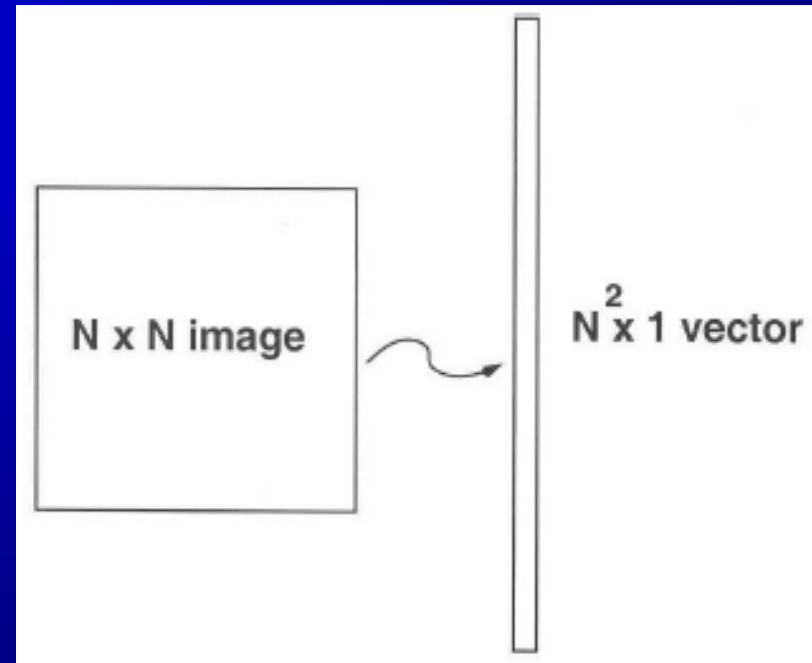
- The principal components are dependent on the *units* used to measure the original variables as well as on the *range* of values they assume.
- You should always **standardize** the data prior to using PCA.
- A common standardization method is to transform all the data to have zero mean and unit standard deviation:

$$\frac{x_i - \mu}{\sigma} \quad (\mu \text{ and } \sigma \text{ are the mean and standard deviation of } x_i\text{'s})$$



# Application to Faces

- Case Study: Eigenfaces for Face Detection/Recognition
  - M. Turk, A. Pentland, "Eigenfaces for Recognition", *Journal of Cognitive Neuroscience*, vol. 3, no. 1, pp. 71-86, 1991.
- Face Recognition
  - The simplest approach is to think of it as a template matching problem
  - Problems arise when performing recognition in a high-dimensional space.
  - Significant improvements can be achieved by first mapping the data into a *lower dimensionality* space.

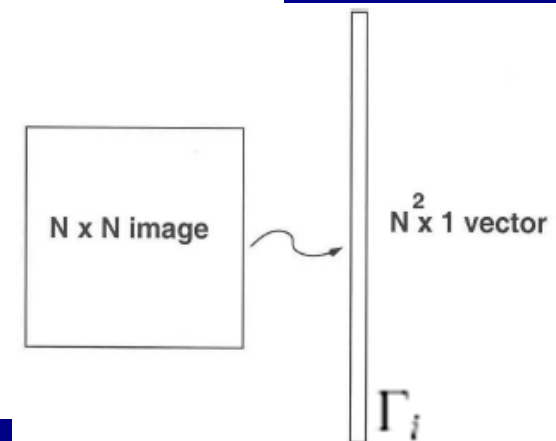
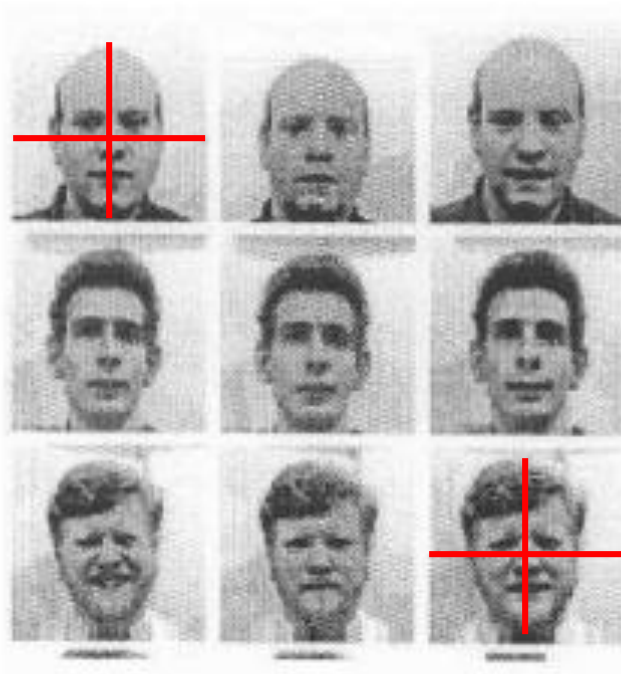


# Application to Faces

- Computation of low-dimensional basis (i.e., eigenfaces):

Step 1: obtain face images  $I_1, I_2, \dots, I_M$  (training faces)

(**very important:** the face images must be *centered* and of the same *size*)



Step 2: represent every image  $I_i$  as a vector  $\Gamma_i$

# Application to Faces

- Computation of the eigenfaces – cont.

$$\frac{1}{M} \left\{ \begin{matrix} \text{column 1} \\ \text{column 2} \\ \vdots \\ \text{column } M \end{matrix} \right\} + \dots + \begin{matrix} \text{column } N^2 \end{matrix}$$

1                      M                       $N^2$

Step 3: compute the average face vector  $\Psi$ :

$$\Psi = \frac{1}{M} \sum_{i=1}^M \Gamma_i$$

Step 4: subtract the mean face:

$$\Phi_i = \Gamma_i - \Psi$$

Step 5: compute the covariance matrix  $C$ :

$$C = \frac{1}{M} \sum_{n=1}^M \Phi_n \Phi_n^T = \frac{1}{M} A A^T \quad (N^2 \times N^2 \text{ matrix})$$

$$\text{where } A = [\Phi_1 \ \Phi_2 \ \dots \ \Phi_M] \quad (N^2 \times M \text{ matrix})$$





# Application to Faces

- Computation of the Eigen faces – cont.

Step 6: compute the eigenvectors  $u_i$  of  $AA^T \rightarrow AA^T u_i = \lambda_i u_i$   $C \boxed{u_i} = \lambda_i \boxed{u_i}$

The matrix  $AA^T$  is very large --> not practical !!

Step 6.1: consider the matrix  $A^T A$  ( $M \times M$  matrix)

Step 6.2: compute the eigenvectors  $v_i$  of  $A^T A$

$$A^T A v_i = \mu_i v_i$$

What is the relationship between  $u_i$  and  $v_i$ ?

$$A^T A v_i = \mu_i v_i \Rightarrow AA^T A v_i = \mu_i A v_i \Rightarrow C \boxed{A v_i} = \mu_i \boxed{A v_i}$$

$$\rightarrow u_i = A v_i \quad \text{and} \quad \lambda_i = \mu_i$$

Thus,  $AA^T$  and  $A^T A$  have the same eigenvalues and their eigenvectors are related as follows:  $u_i = A v_i$  !!

$$A^T A \quad \mu_i = \lambda_i \quad AA^T$$

Eigenvector of  $C$

6.4276	0	0	0	0
0	0.9822	0	0	0
0	0	0.3178	0	0
0	0	0	0.0000	0
0	0	0	0	0.0000

6.4276	0	0	0	0
0	0.9822	0	0	0
0	0	0.3178	0	0

$AA^T$   
 $\lambda_i$  eigenvalue

$u_i$  eigenvector

$A^T A$   
 $\mu_i$  eigenvalue

$v_i$  eigenvector

$AA^T$

$A^T A$

# Application to Faces

- Computation of the eigenfaces – cont.

Note 1:  $AA^T$  can have up to  $N^2$  eigenvalues and eigenvectors.

Note 2:  $A^T A$  can have up to  $M$  eigenvalues and eigenvectors.

Note 3: The  $M$  eigenvalues of  $A^T A$  (along with their corresponding eigenvectors) correspond to the  $M$  *largest* eigenvalues of  $AA^T$  (along with their corresponding eigenvectors).

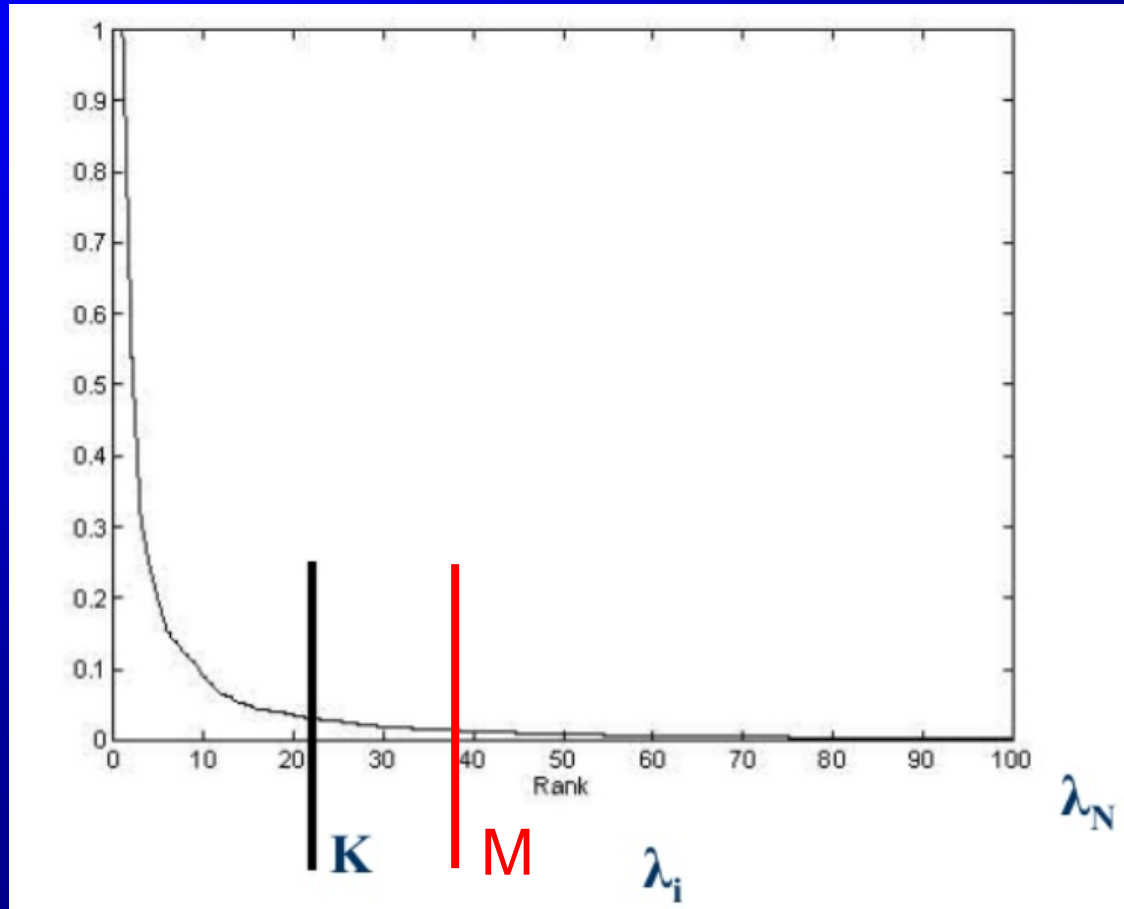
Step 6.3: compute the  $M$  best eigenvectors of  $AA^T$ :  $u_i = Av_i$

(**important:** normalize  $u_i$  such that  $\|u_i\| = 1$ )

Step 7: keep only  $K$  eigenvectors (corresponding to the  $K$  largest eigenvalues)

# Application to Faces

- Computation of the eigenfaces – cont.



# Eigenfaces example

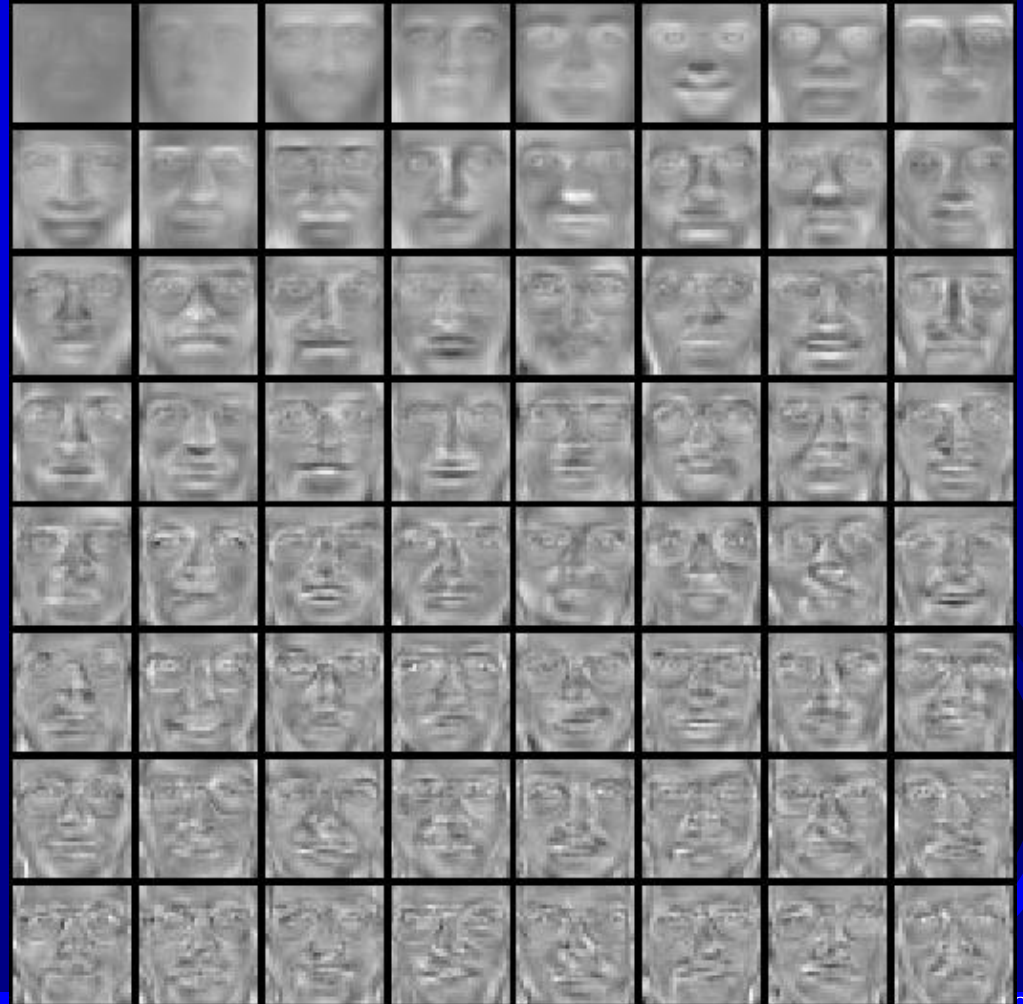
Training  
images



# Eigenfaces example

Top eigenvectors:  $u_1, \dots, u_k$

Mean:  $\mu$





# Application to Faces

- Representing faces onto this basis

- Each face (minus the mean)  $\Phi_i$  in the training set can be represented as a linear combination of the best  $K$  eigenvectors:

$$\hat{\Phi}_i - mean = \sum_{j=1}^K w_j u_j, \quad (w_j = u_j^T \Phi_i)$$

(we call the  $u_j$ 's *eigenfaces*)



Face  
reconstruc  
tion:

# Eigenfaces

- Face Recognition Using Eigenfaces

- Given an unknown face image  $\Gamma$  (centered and of the same size like the training faces) follow these steps:

Step 1: normalize  $\Gamma$ :  $\Phi = \Gamma - \Psi$

Step 2: project on the eigenspace

$$\hat{\Phi} = \sum_{i=1}^K w_i u_i \quad (w_i = u_i^T \Phi) \quad (\text{where } \|u_i\| = 1)$$

Step 3: represent  $\Phi$  as:  $\Omega = \begin{bmatrix} w_1 \\ w_2 \\ \dots \\ w_K \end{bmatrix}$

Step 4: find  $e_r = \min_k \|\Omega - \Omega^k\|$ , where  $\Omega^k$  is the  $k$ th face vector.

Step 5: if  $e_r < T_r$ , then  $\Gamma$  is recognized as face  $l$  from the training set.

# Eigenfaces

- Face Recognition Using Eigenfaces – cont.
  - The distance  $e_r$  is called distance within face space (difs)
  - The *Euclidean distance* can be used to compute  $e_r$ , however, the *Mahalanobis distance (马氏距离)* has shown to work better:

$$\|\Omega - \Omega^k\| = \sum_{i=1}^K (w_i - w_i^k)^2 \quad \text{Euclidean distance}$$



*Mahalanobis distance*

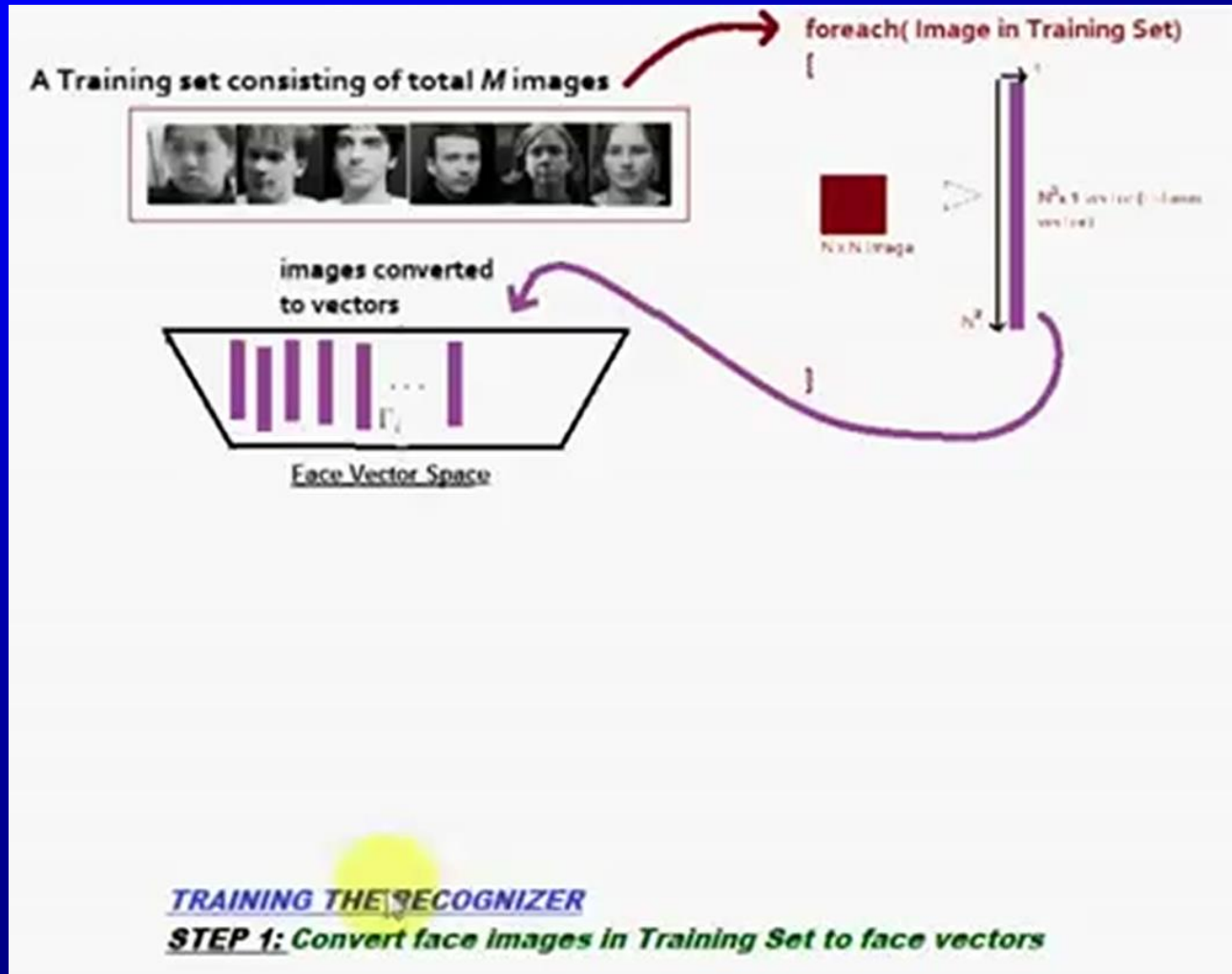
$$\|\Omega - \Omega^k\| = \sum_{i=1}^K \frac{1}{\lambda_i} (w_i - w_i^k)^2$$

(variations along all axes are treated as equally significant)



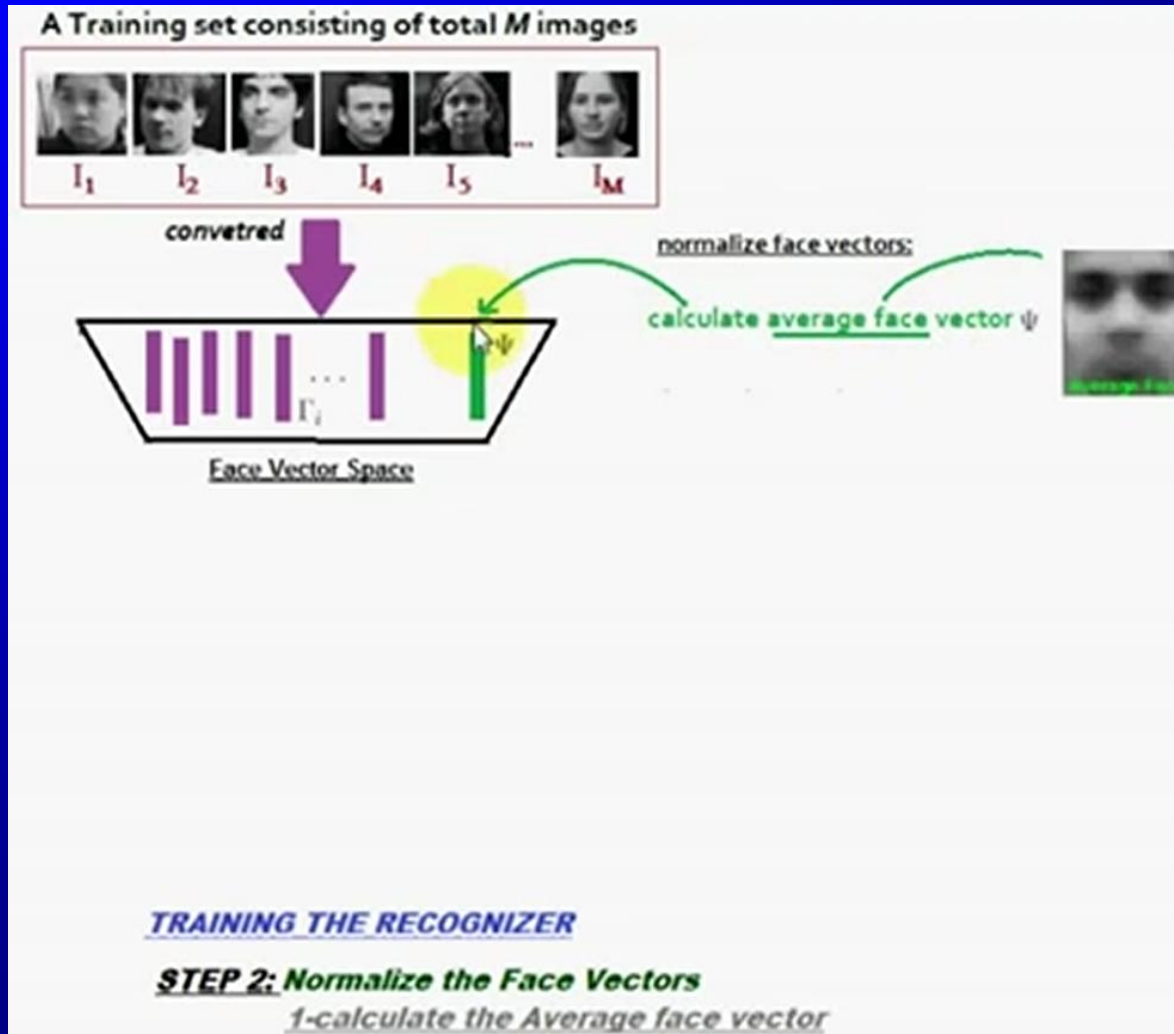
# Eigenfaces

- Face Recognition Using Eigenfaces



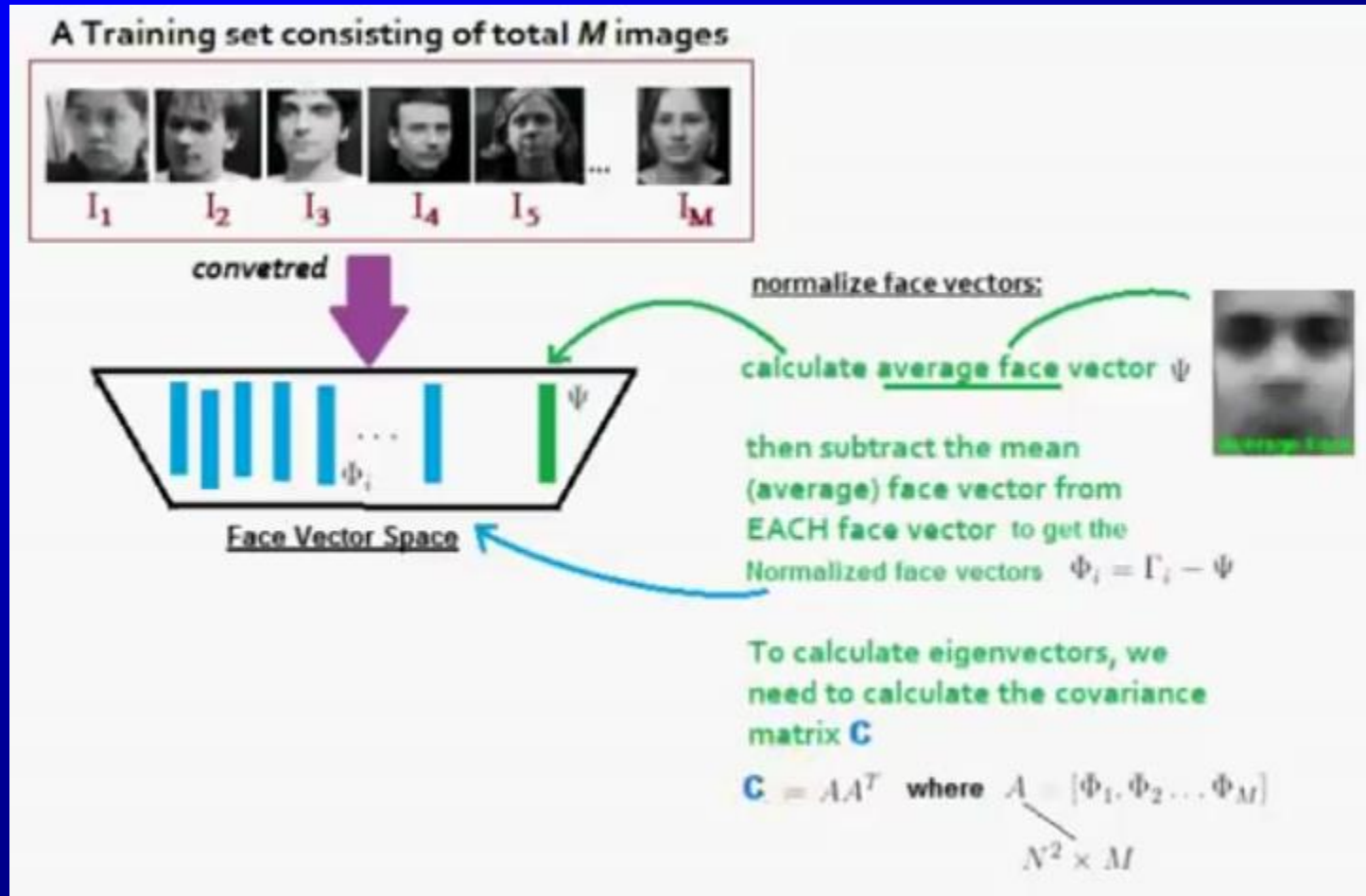
# Eigenfaces

- Face Recognition Using Eigenfaces



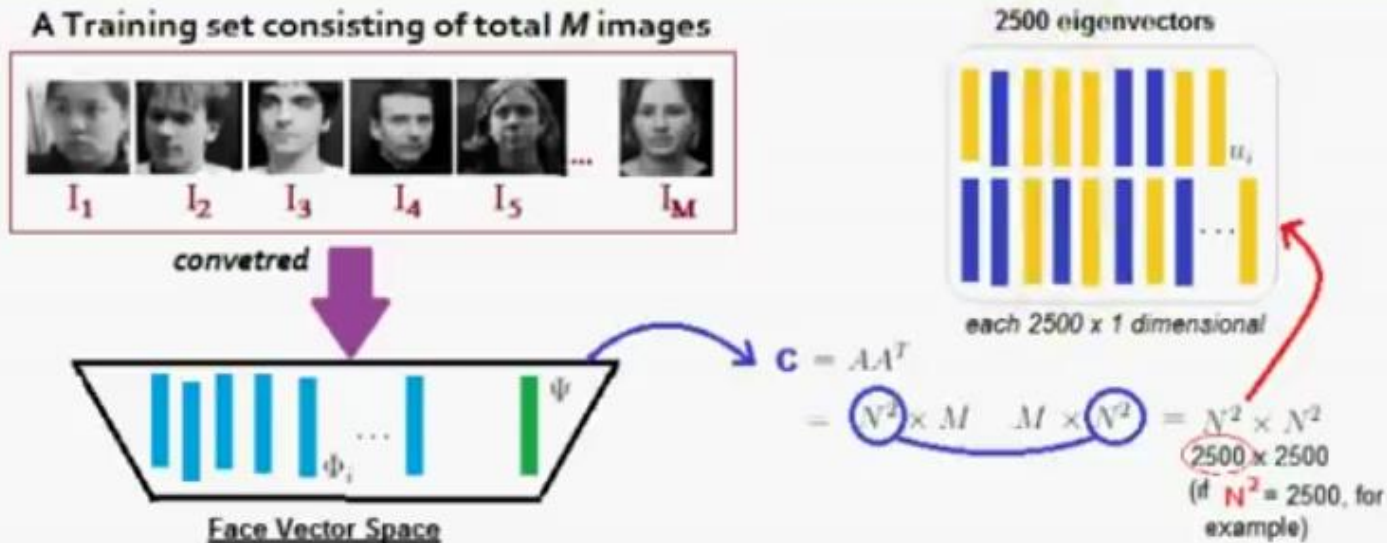
# Eigenfaces

- Face Recognition Using Eigenfaces



# Eigenfaces

- Face Recognition Using Eigenfaces



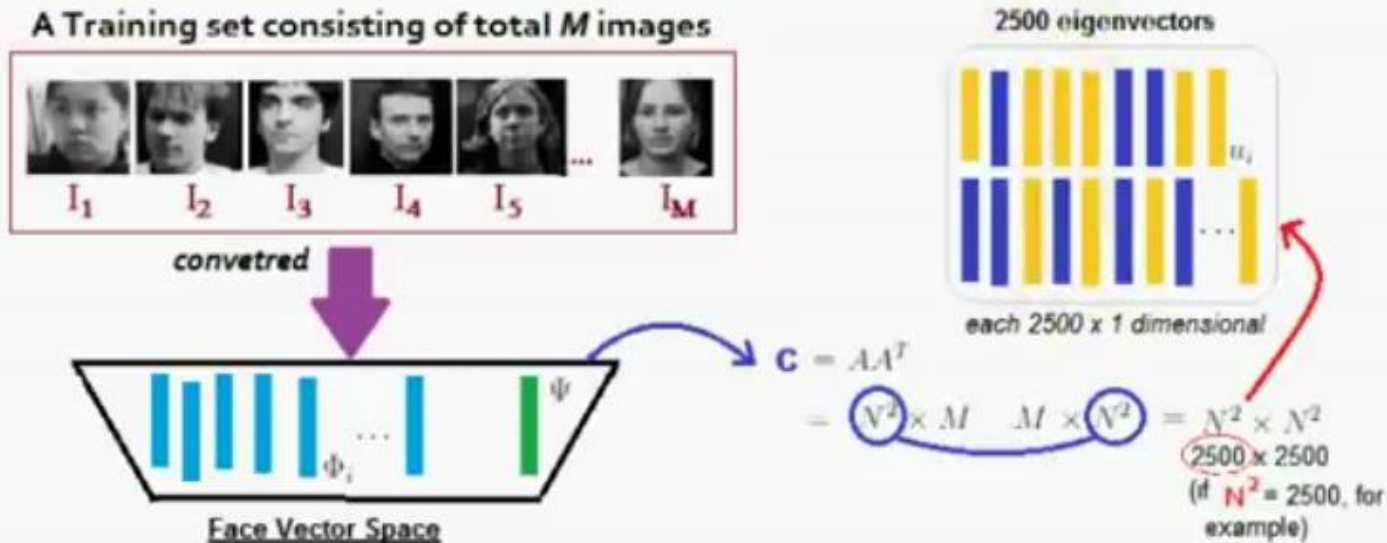
WARNING!!!

System may slow down terribly or run out of memory! Computations required are HUGE.



# Eigenfaces

- Face Recognition Using Eigenfaces



Solution!!!

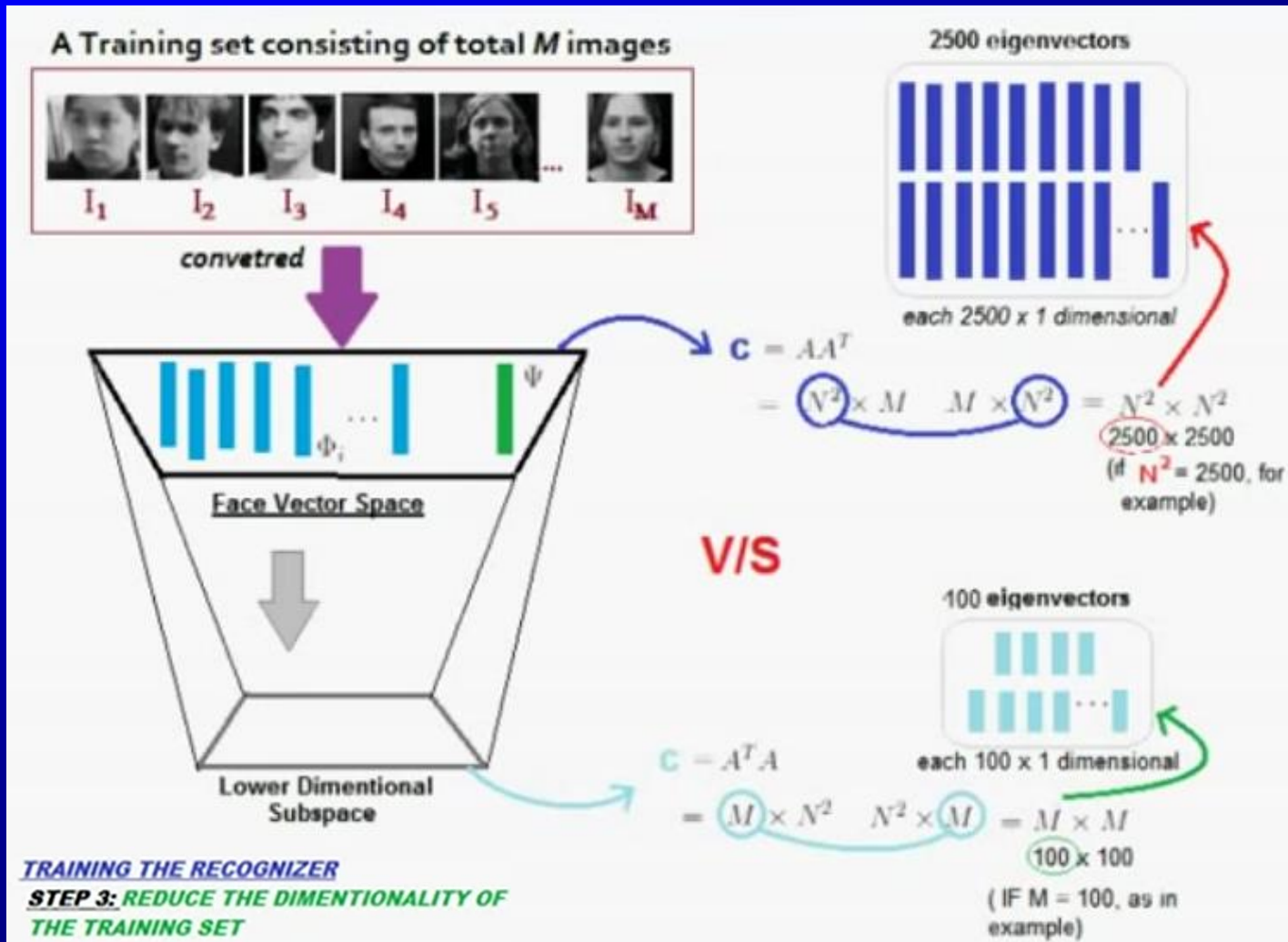
Dimensionality Reduction

Calculate eigenvectors  
from a Covariance with  
reduced dimensionality.



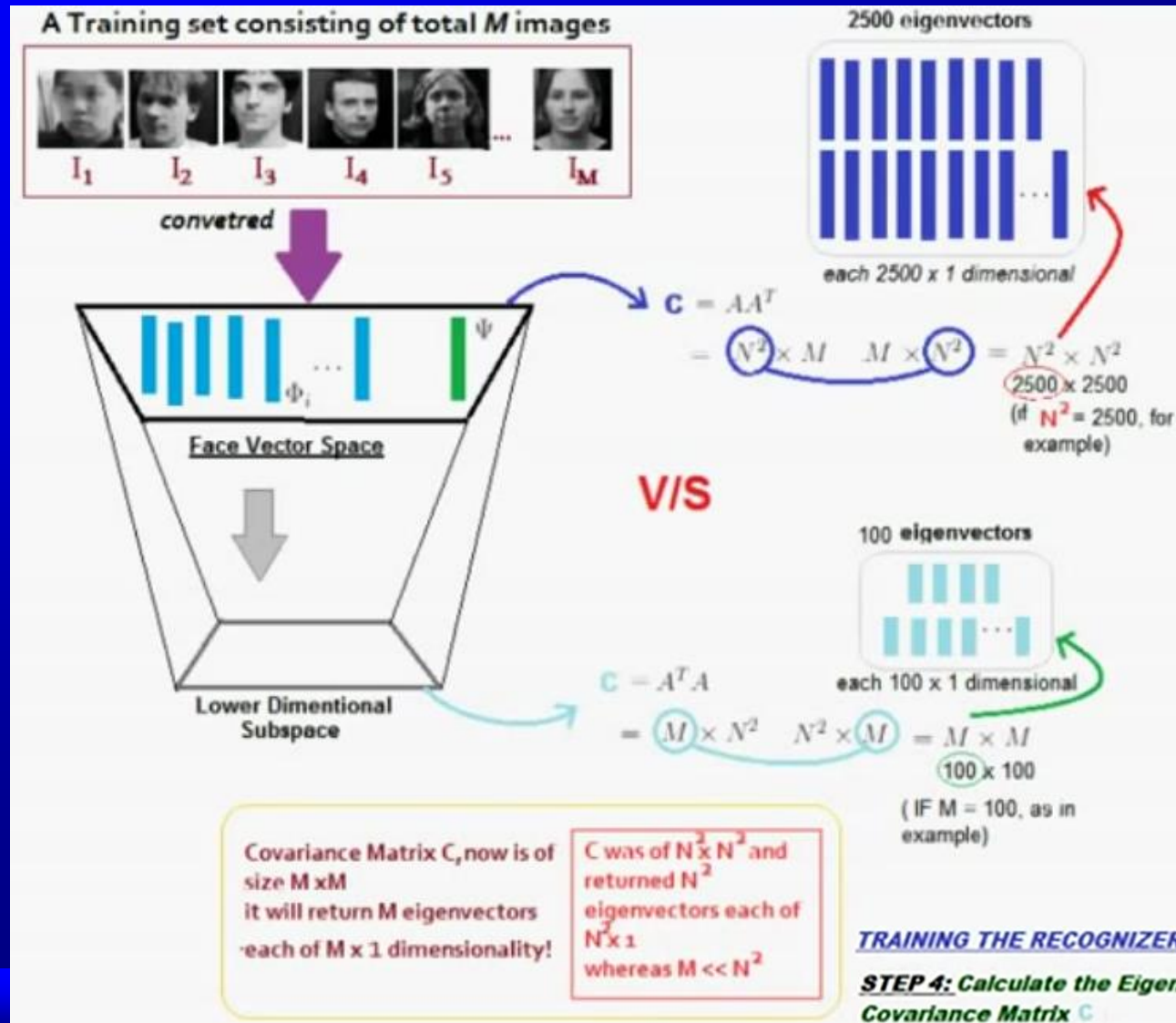
# Eigenfaces

- Face Recognition Using Eigenfaces



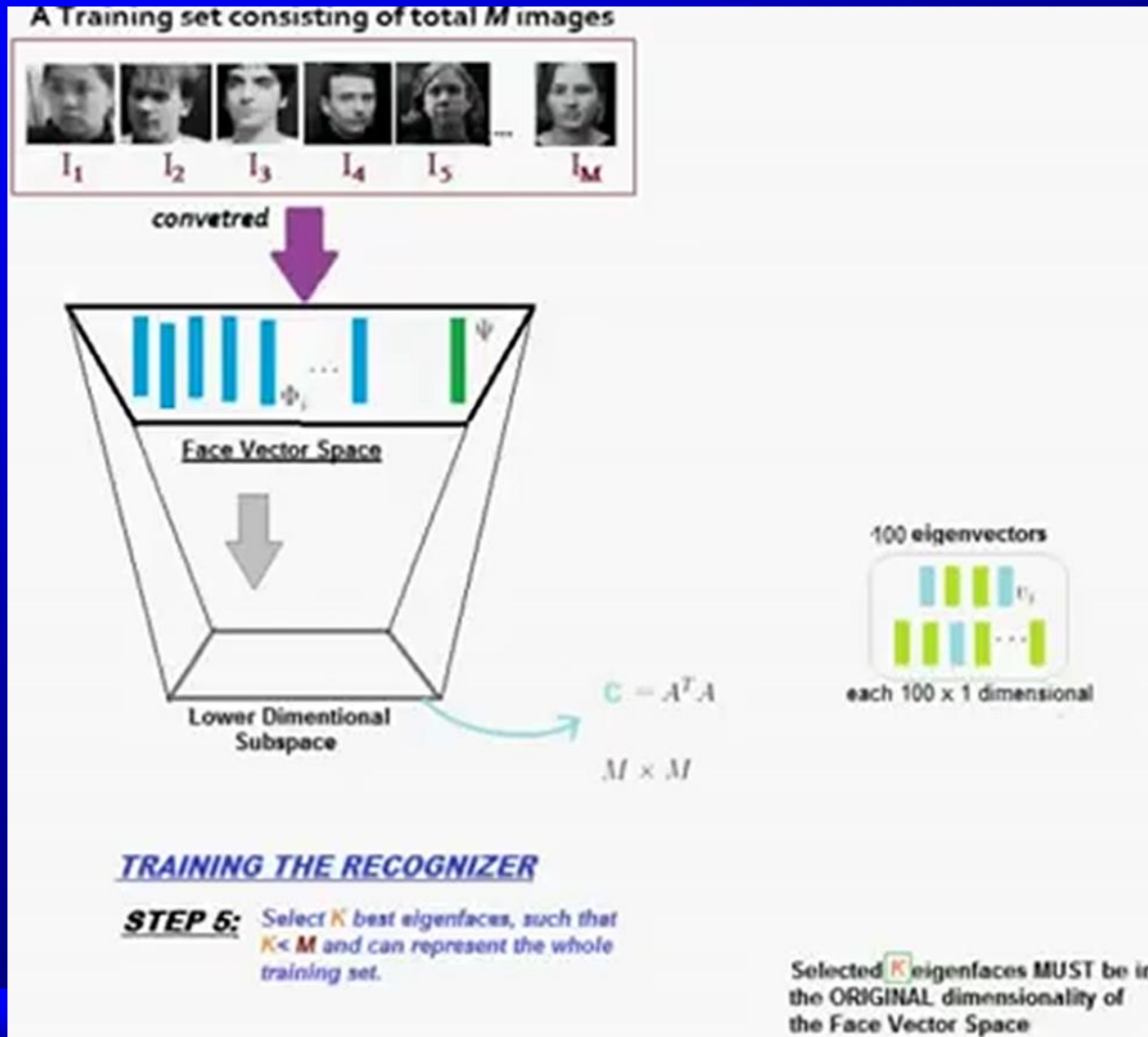
# Eigenfaces

- Face Recognition Using Eigenfaces



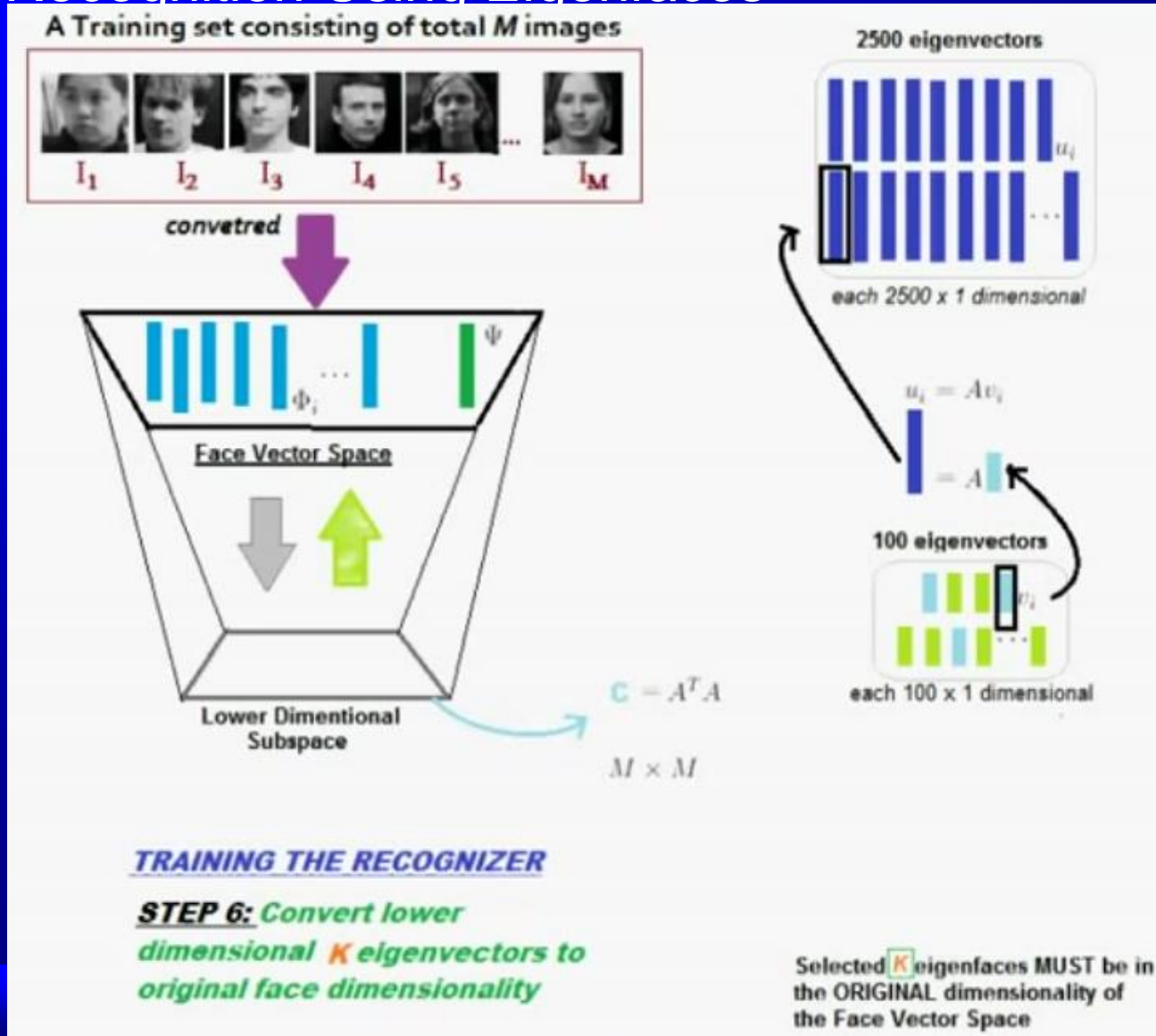
# Eigenfaces

- Face Recognition Using Eigenfaces



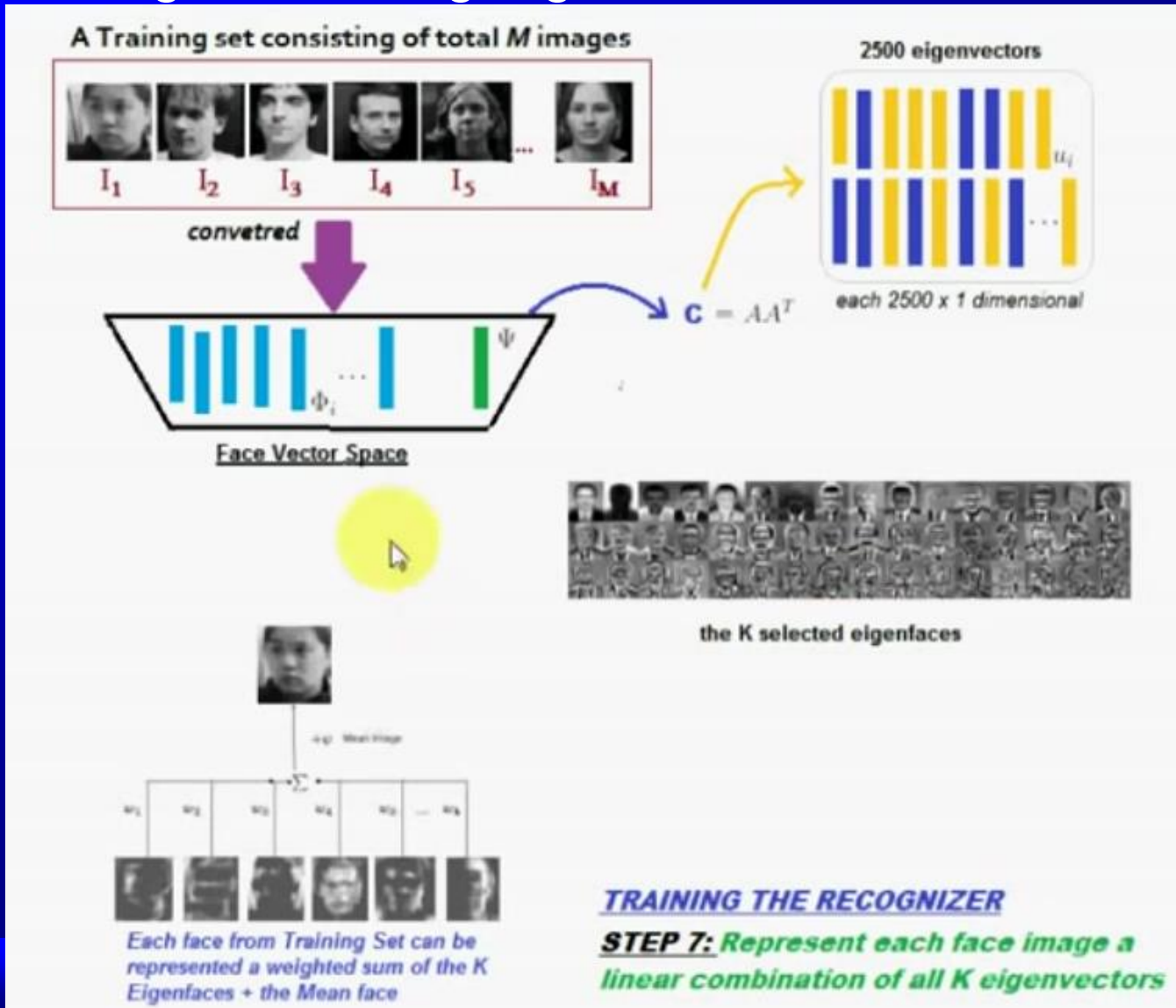
# Eigenfaces

- Face Recognition Using Eigenfaces



# Eigenfaces

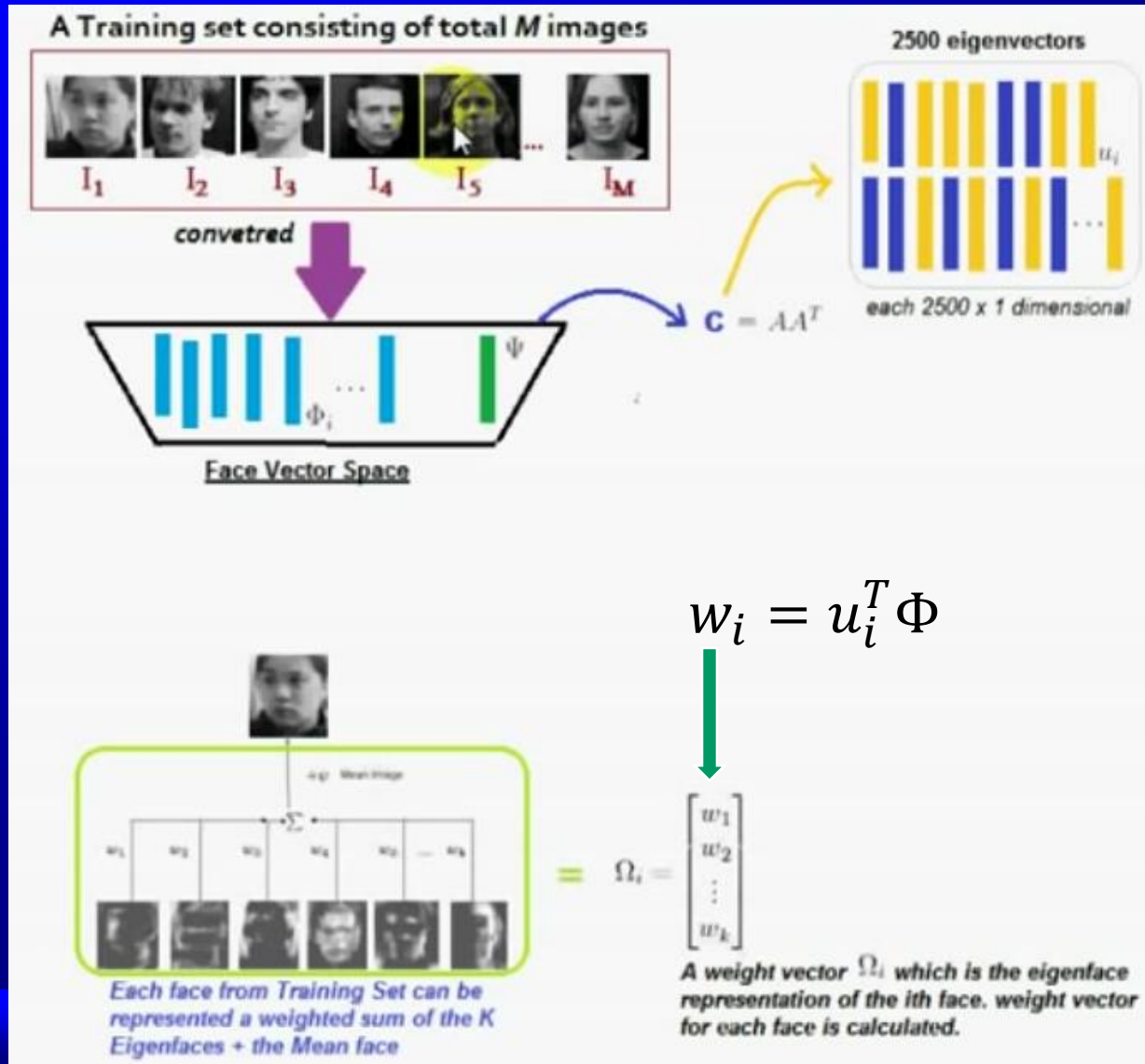
- Face Recognition Using Eigenfaces





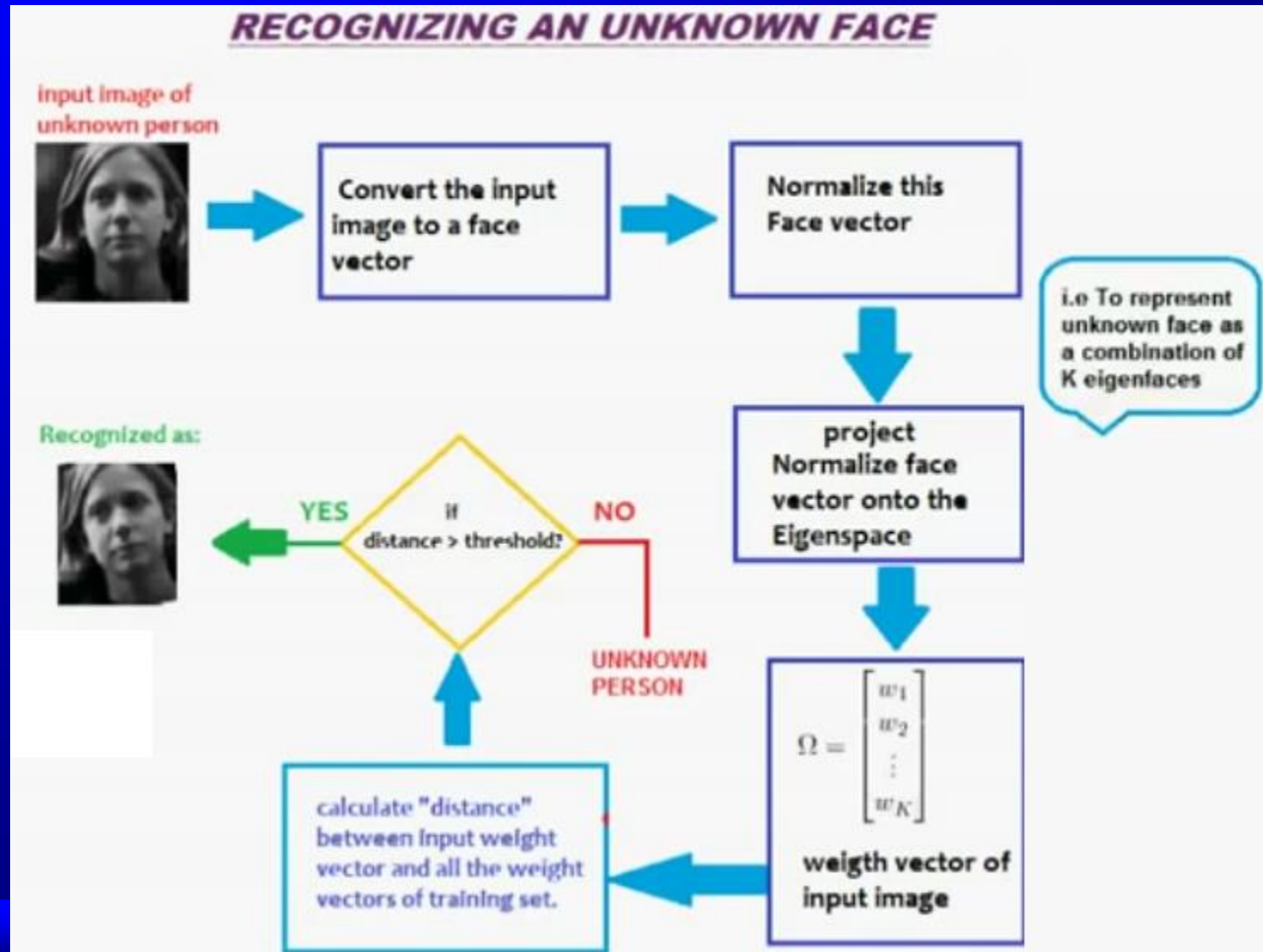
# Eigenfaces

- Face Recognition Using Eigenfaces

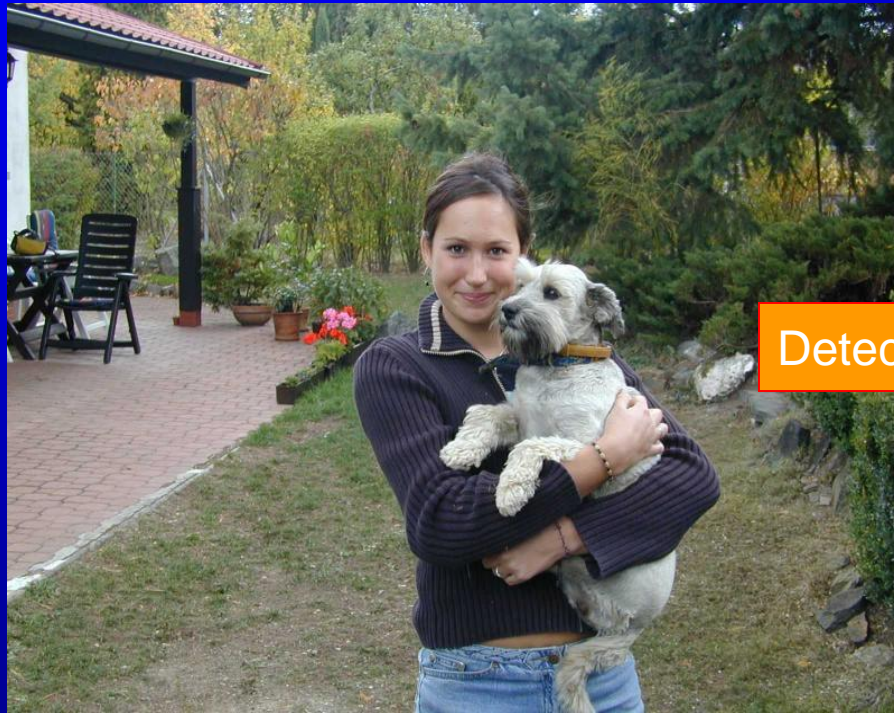


# Eigenfaces

- Face Recognition Using Eigenfaces



# Face detection and recognition



Detection



Recognition

"Sally"





# Eigenfaces

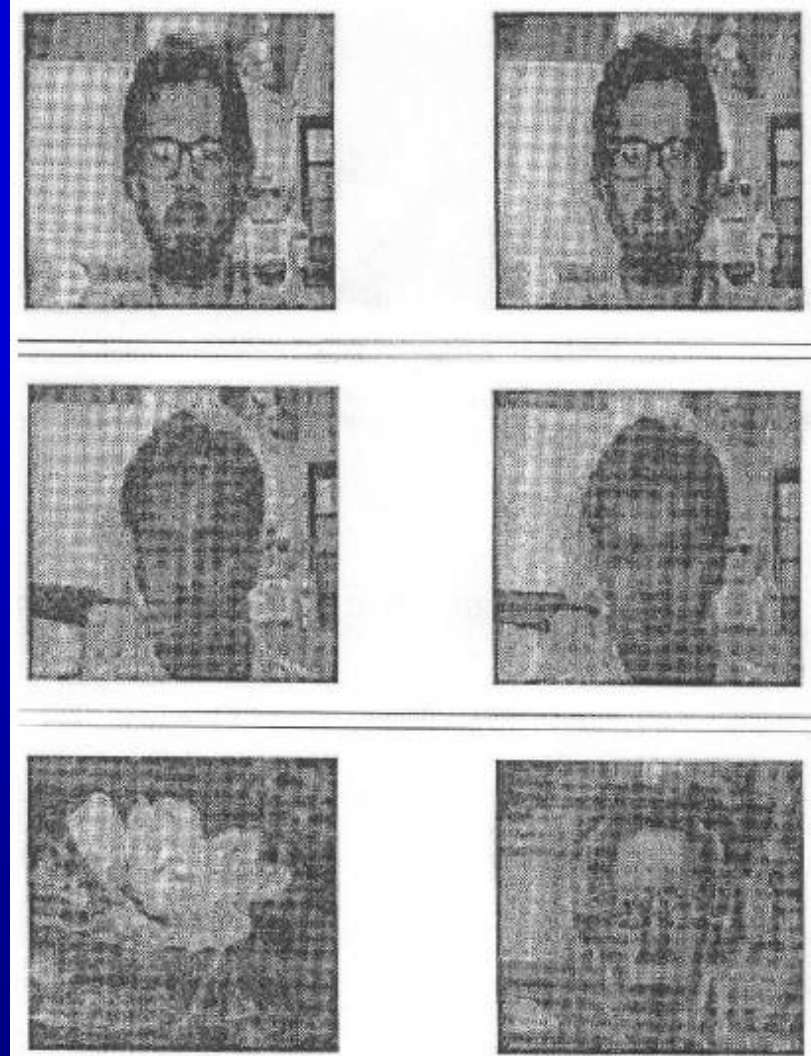
- Reconstruction of faces and non-faces

Reconstructed face looks like a face.

Reconstructed non-face looks like a face again!

Input

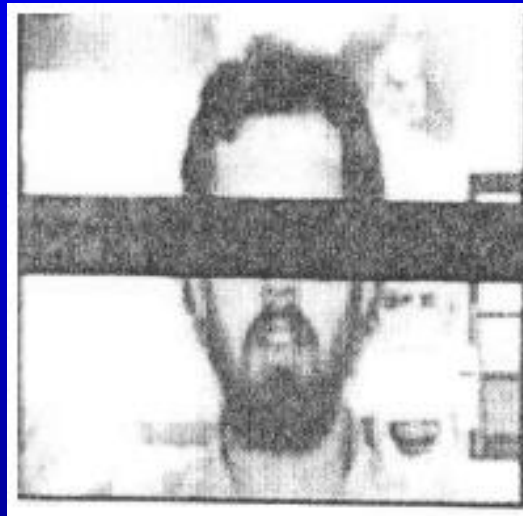
Reconstructed



# Reconstruction using partial information

- Robust to partial face occlusion.

Input



Reconstructed



# Limitations

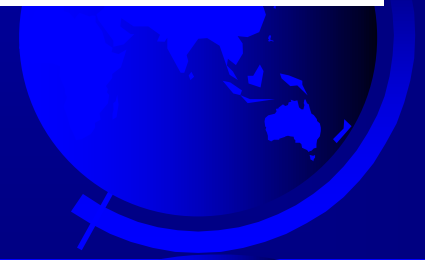


- **Background** changes cause problems
  - De-emphasize the outside of the face (e.g., by multiplying the input image by a 2D Gaussian window centered on the face).
- **Light** changes degrade performance
  - Light normalization helps.
- Performance decreases quickly with changes to face size
  - Multi-scale eigenspaces.
  - Scale input image to multiple sizes.
- Performance decreases with changes to **face orientation** (but not as fast as with scale changes)
  - Plane rotations are easier to handle.
  - Out-of-plane rotations are more difficult to handle.



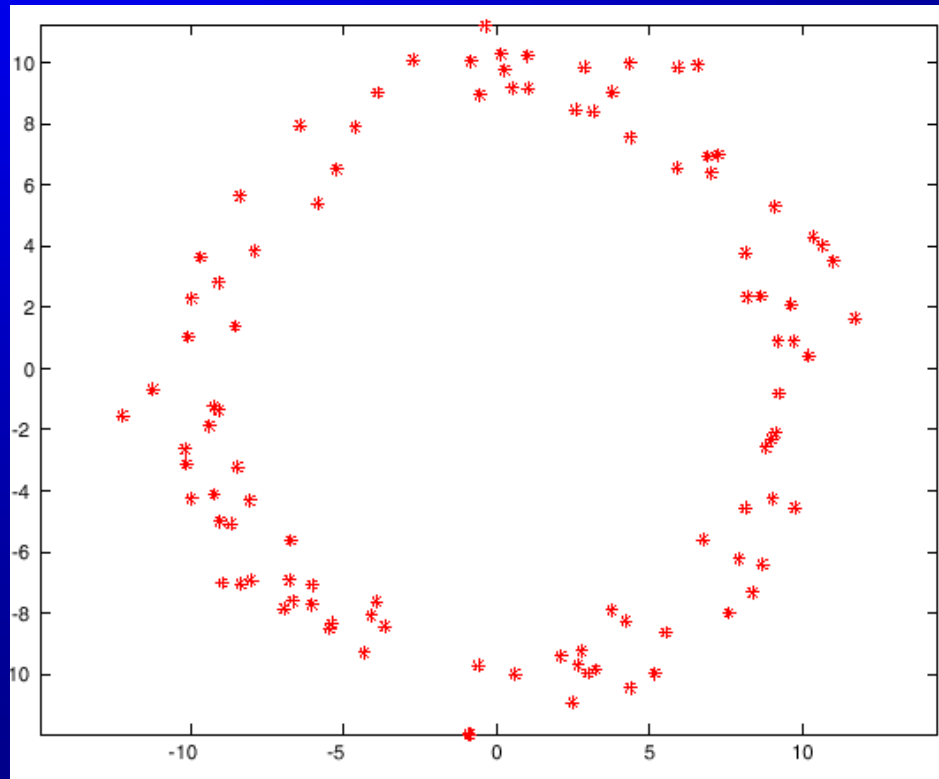
# Limitations

☞ Not robust to misalignment

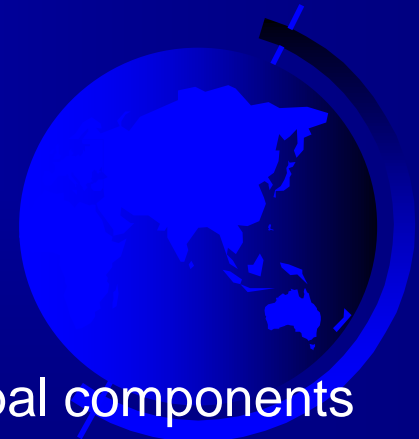


# Limitations

- ✎ PCA assumes that the data follows a Gaussian distribution (mean  $\mu$ , covariance matrix  $\Sigma$ )

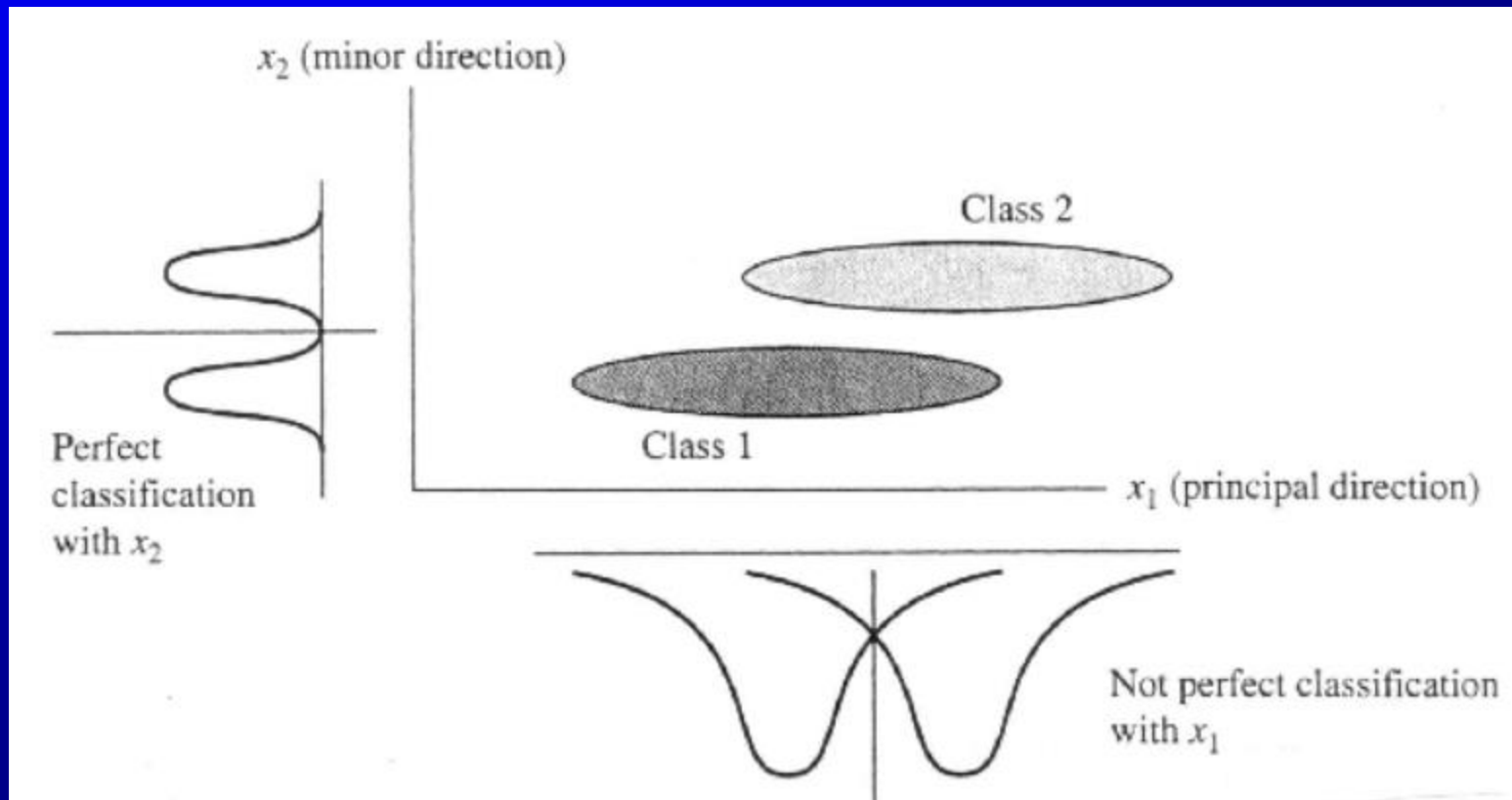


The shape of this dataset is not well described by its principal components



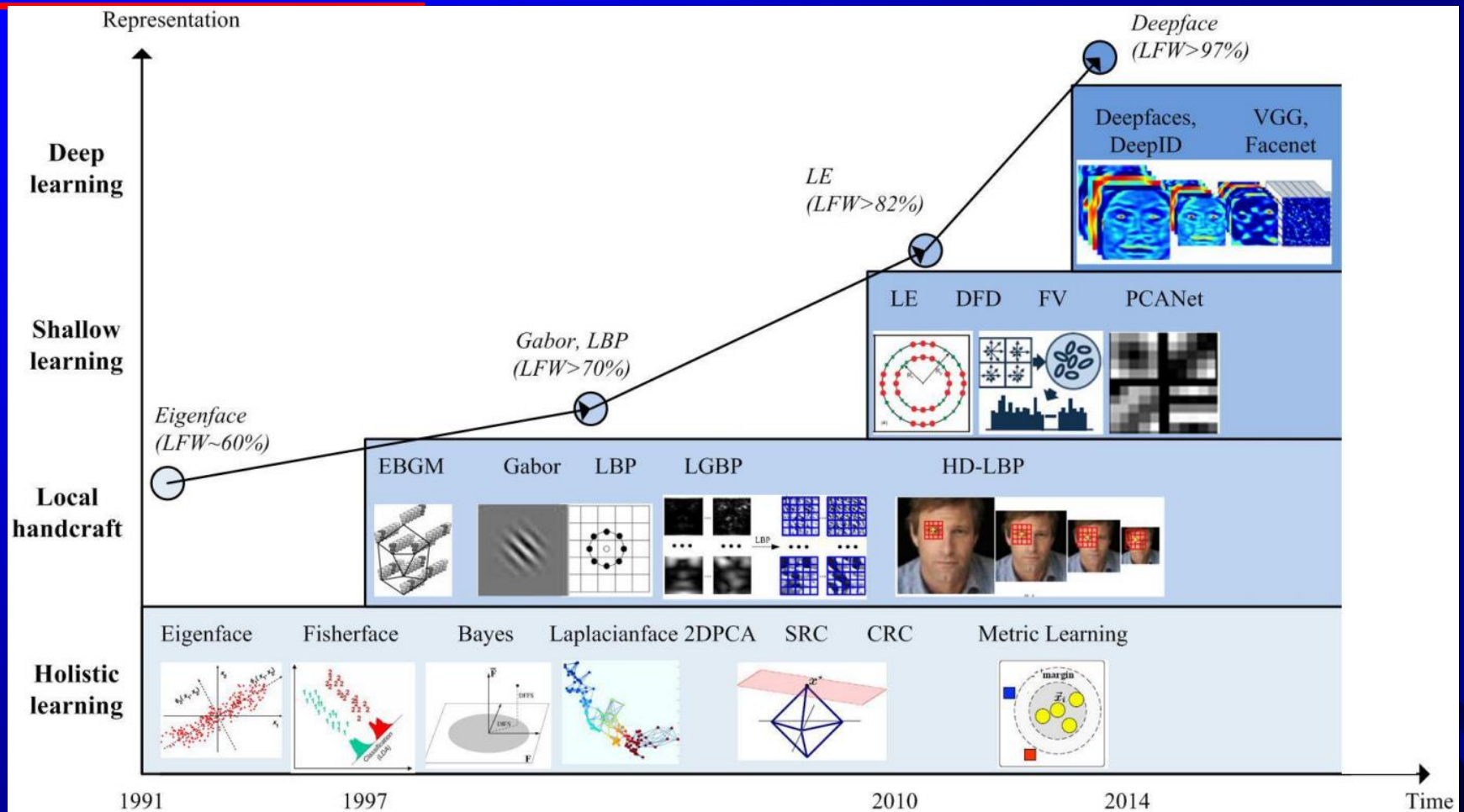
# Limitations

- PCA is not always an optimal dimensionality-reduction procedure for classification purposes:





# Face Recognition-State-of-the-art



Milestones of feature representation for FR. The holistic approaches dominated the FR community in the 1990s and 2000s. In the early 2000s and 2010s, Local-feature-based FR and learning-based local descriptors were introduced successively. In 2014, DeepFace and DeepID achieved state-of-the-art accuracy, and research focus has shifted to deep-learning-based approaches. As the representation pipeline becomes deeper and deeper, the LFW (Labeled Face in-the-Wild) performance steadily improves from around 60% to above 90%, while deep learning boosts the performance to 99.80% in only three years.

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# See You

