

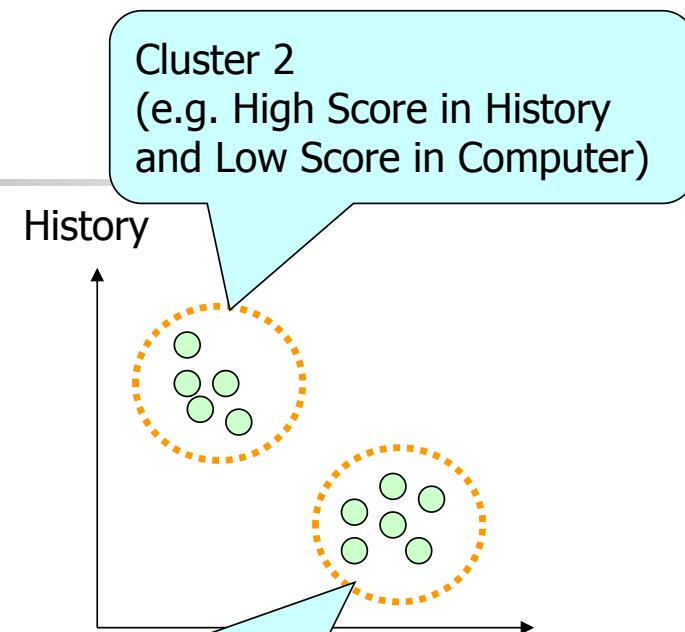


# Subspace Clustering

Prepared by Raymond Wong  
Presented by Raymond Wong  
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# Clustering

	Computer	History
Raymond	100	40
Louis	90	45
Wyman	20	95
...	...	...



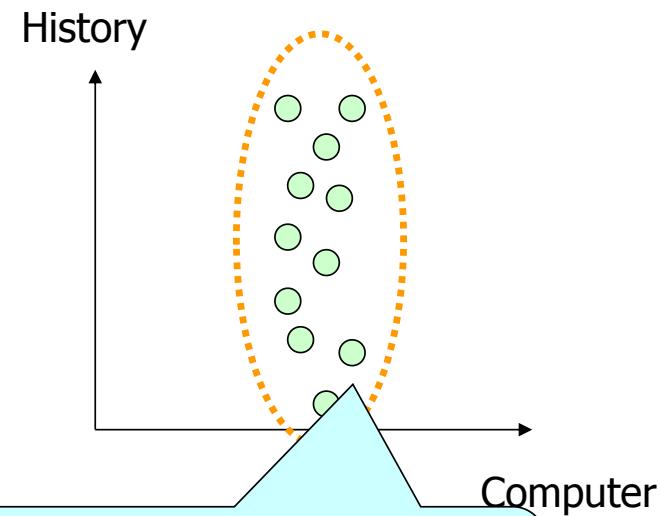
Cluster 2  
(e.g. High Score in History  
and Low Score in Computer)

Problem: to find all clusters

This kind of clustering considers only FULL space (i.e.  
computer and history)!

# Subspace Clustering

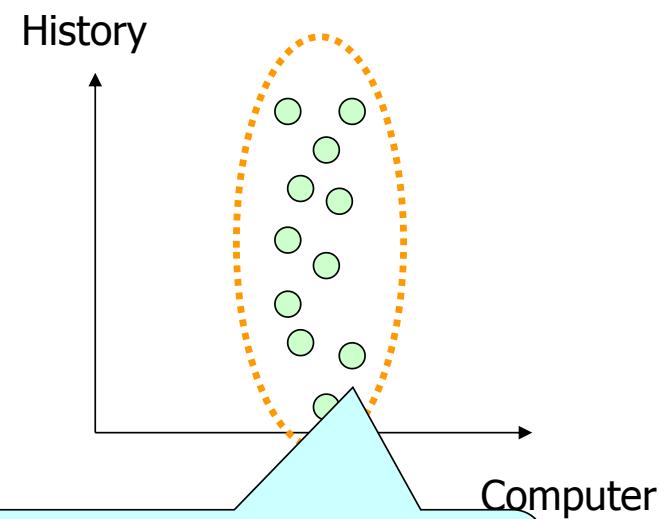
	Computer	History
Raymond	50	40
Louis	60	45
Wyman	40	95
...	...	...



Cluster 1  
(e.g. Middle Score in Computer  
and Any Score in History)

# Subspace Clustering

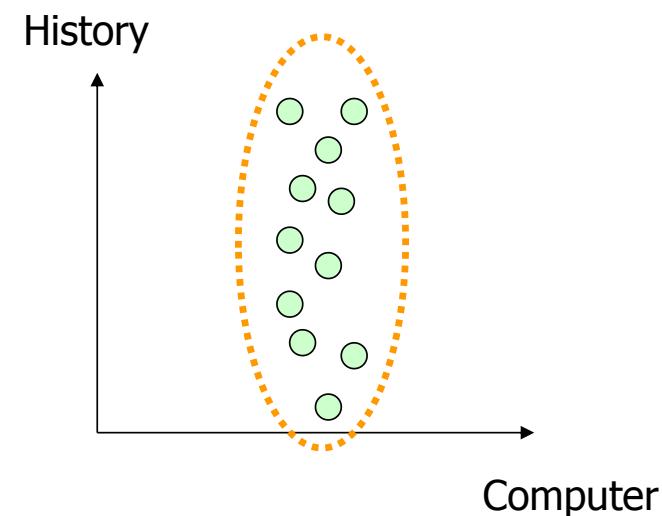
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Cluster 1  
(e.g. Middle Score in Computer  
and History)

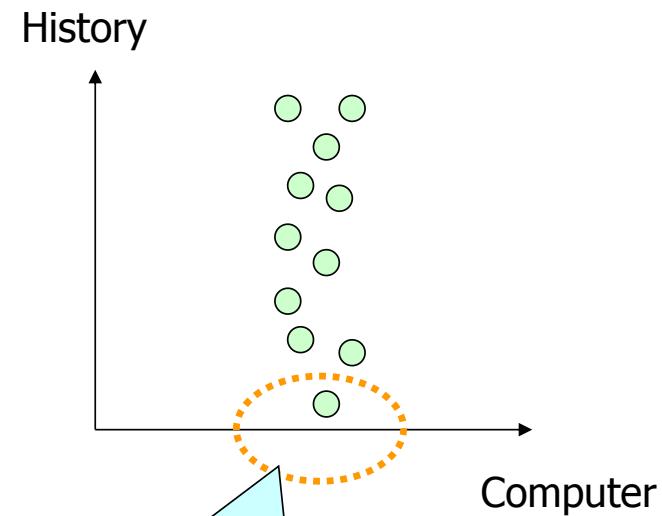
# Subspace Clustering

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Raymond	50	40
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# Subspace Clustering

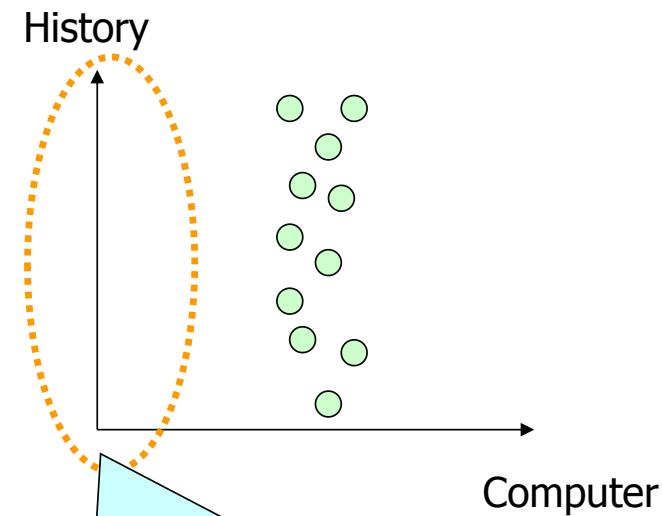
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Cluster 1  
(e.g. Middle Score in Computer)

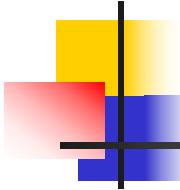
# Subspace Clustering

	Computer	History
Raymond	50	40
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Wyman	40	95
...	...	...



No Cluster!  
The data points span along history dimension.

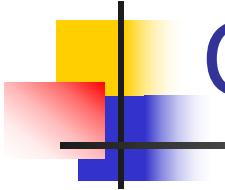
Problem: to find all clusters in the subspace (i.e. some of the dimensions)



# Why Subspace Clustering?

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- Clustering for Understanding
  - Applications
    - Biology
      - Group different species
    - Psychology and Medicine
      - Group medicine
    - Business
      - Group different customers for marketing
    - Network
      - Group different types of traffic patterns
    - Software
      - Group different programs for data analysis



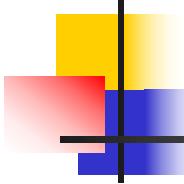
# Curse of Dimensionality

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- When the number of dimensions increases,
  - the distance between any two points is nearly the same

Surprising results!

This is the reason why we need to study subspace clustering



# Subspace Clustering Methods

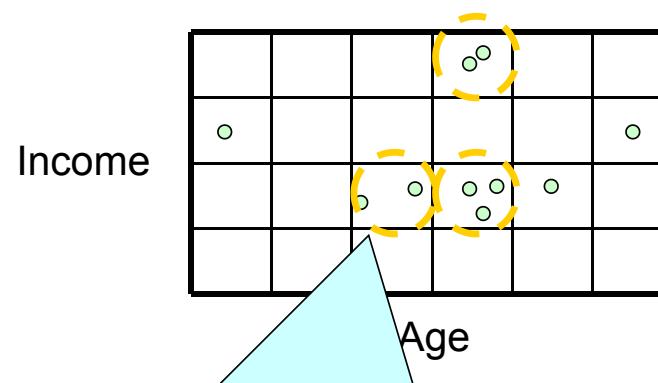
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- Dense Unit-based Method
- Entropy-Based Method
- Transformation-Based Method

# Dense Unit-based Method for Subspace Clustering

Density

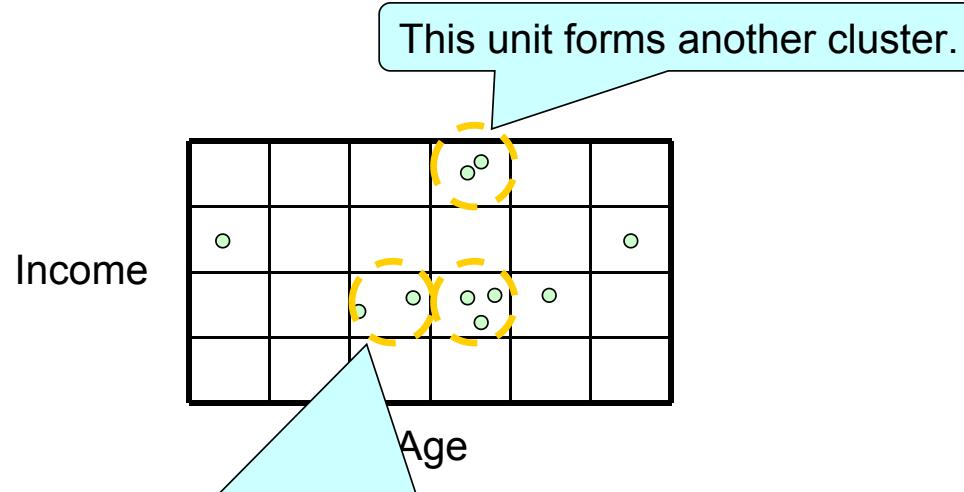
**Dense unit:** a unit if the fraction of data points contained in it is at least a threshold,  $T$



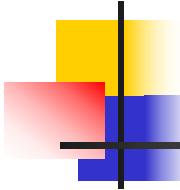
If  $T = 20\%$ , these three units are dense.

# Dense Unit-based Method for Subspace Clustering

**Cluster:** a maximal set of connected dense units in k-dimensions



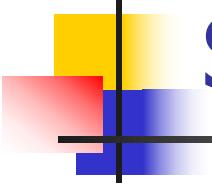
The problem is to find which **sub-spaces** contain dense units.  
The second problem is to find clusters from each sub-space containing dense units



# Dense Unit-based Method for Subspace Clustering

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- **Step 1:** Identify sub-spaces that contain dense units
- **Step 2:** Identify clusters in each sub-spaces that contain dense units



# Step 1

---

Suppose we want to find all dense units (e.g.,  
dense units with density  $\geq 20\%$ )

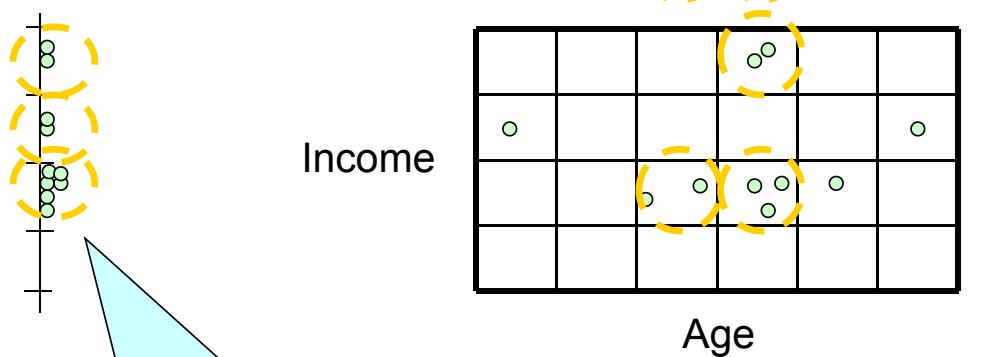
## ■ Property

- If a set  $S$  of points is a cluster in a  $k$ -dimensional space, then  $S$  is also part of a cluster in any  $(k-1)$ -dimensional projections of the space.

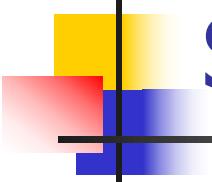
# Step 1

Suppose we want to find all dense units (e.g., dense units with density  $\geq 20\%$ )

If  $T = 20\%$ , these two units are dense.



If  $T = 20\%$ , these three units are dense.



## Step 1

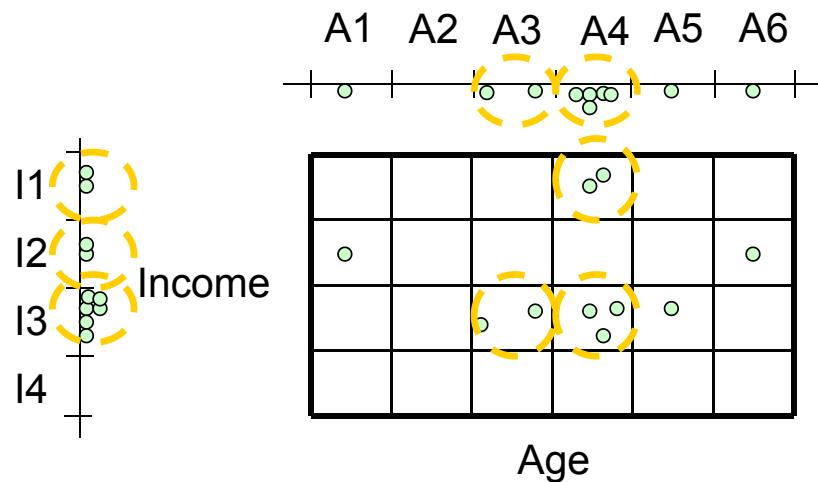
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Suppose we want to find all dense units (e.g.,  
dense units with density  $\geq 20\%$ )

- We can make use of apriori approach to solve the problem

# Step 1

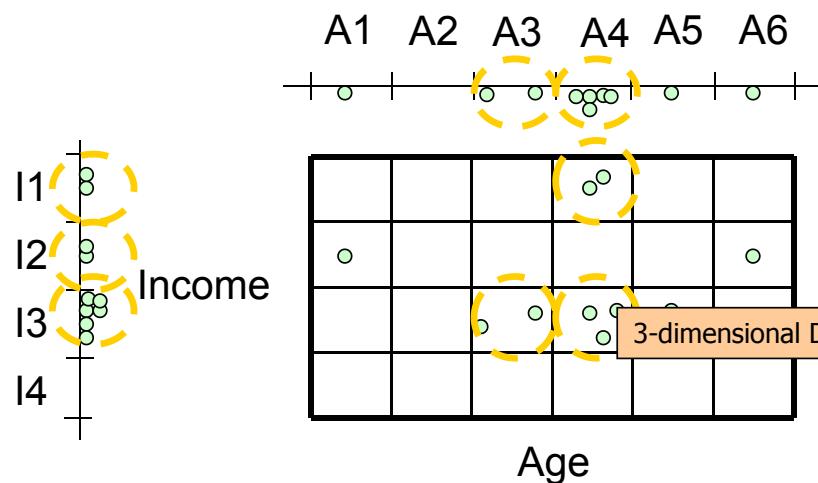
Suppose we want to find all dense units (e.g.,  
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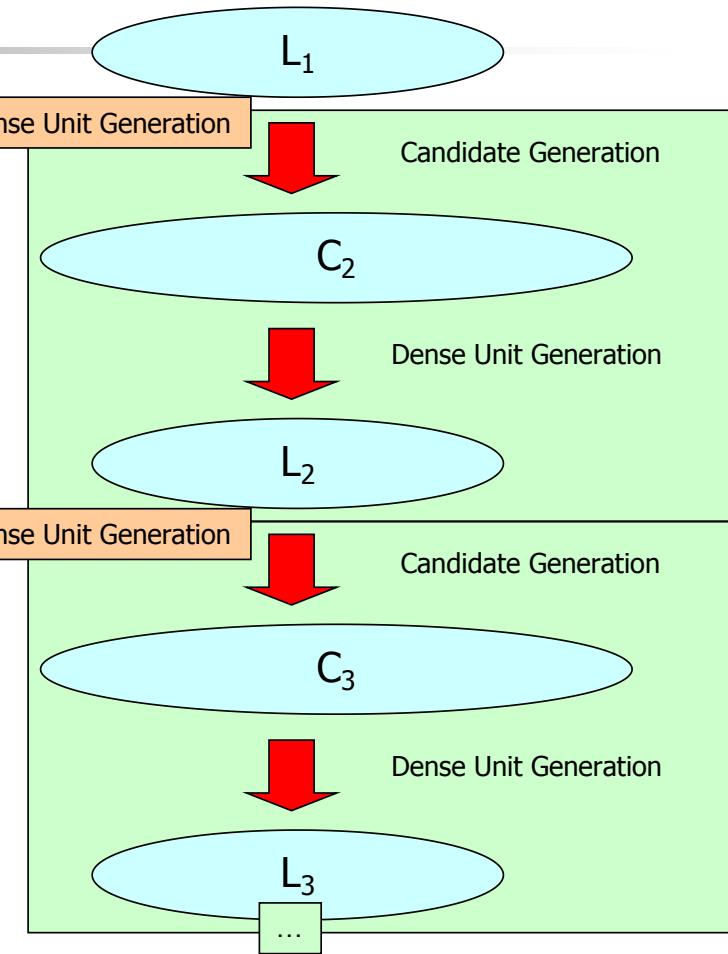
With respect to dimension Age,  
A3 and A4 are dense units.  
With respect to dimension Income,  
I1, I2 and I3 are dense units

# Apriori

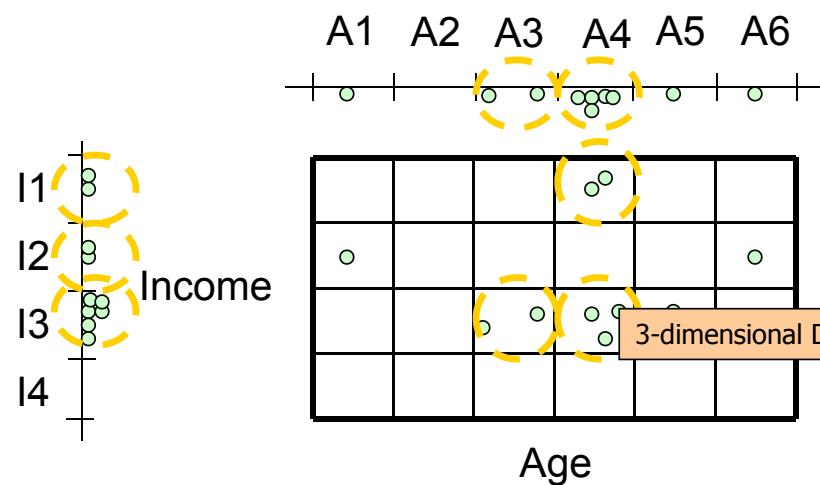
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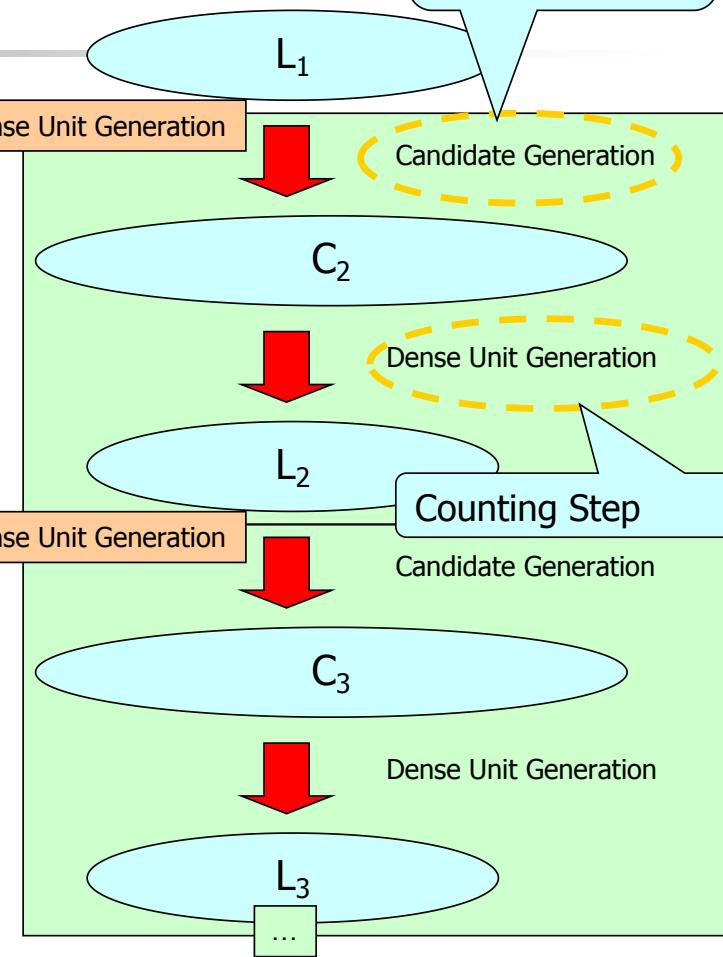
# Apriori

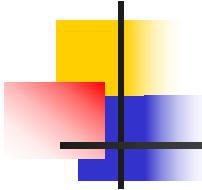


With respect to dimension Age,  
A3 and A4 are dense units.  
With respect to dimension Income,  
I1, I2 and I3 are dense units

Suppose we want to find all dense units  
(dense units with density  $\geq 20\%$ )

1. Join Step
2. Prune Step





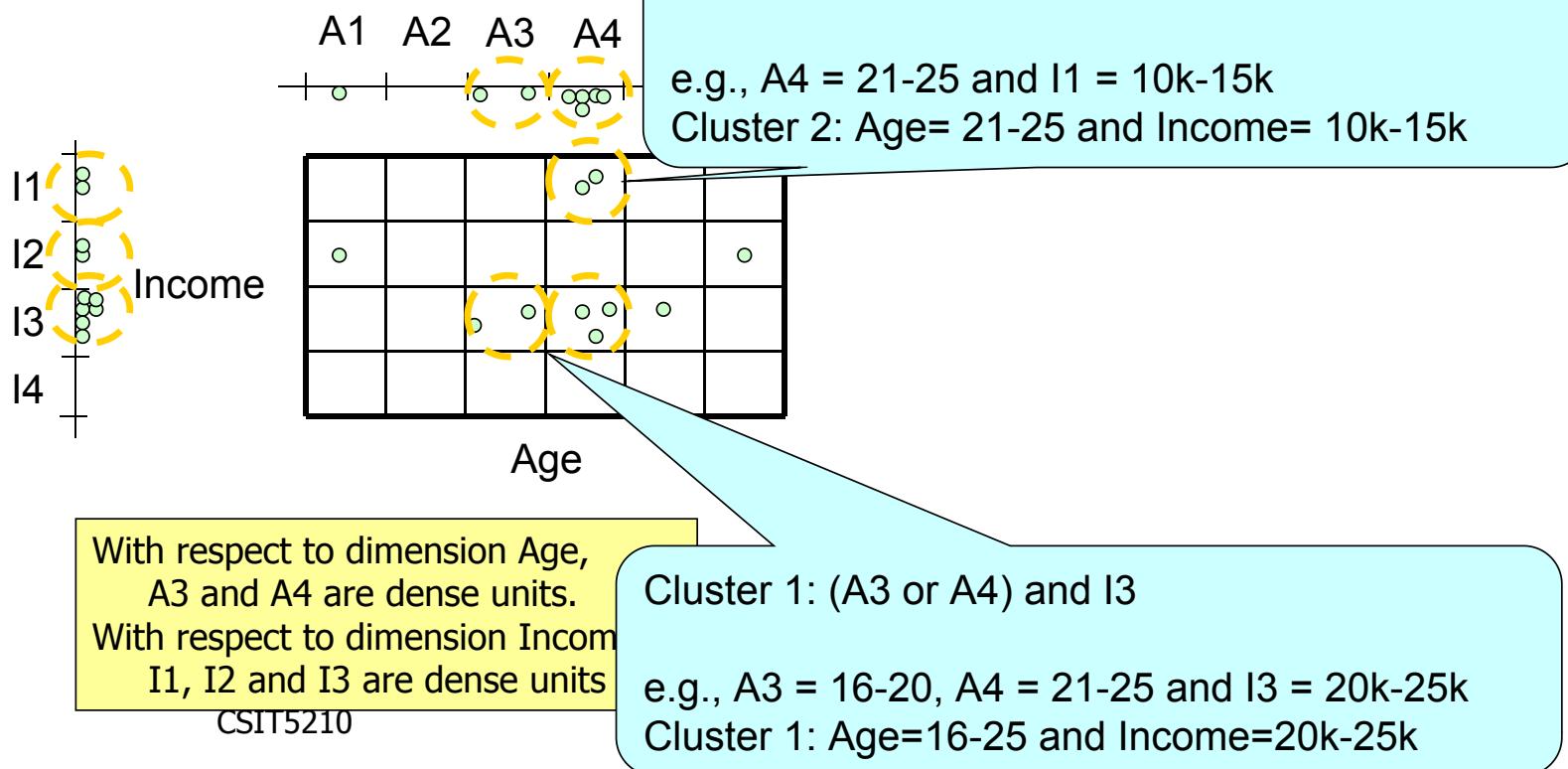
# Dense Unit-based Method for Subspace Clustering

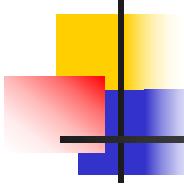
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- **Step 1:** Identify sub-spaces that contain dense units
- **Step 2:** Identify clusters in each sub-spaces that contain dense units

## Step 2

Suppose we want to find all dense units (e.g.,  
dense units with density  $\geq 20\%$ )

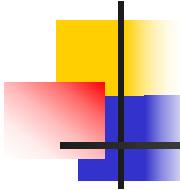




# Subspace Clustering Methods

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- Dense Unit-based Method
- Entropy-Based Method
- Transformation-Based Method



# Entropy-Based Method for Subspace Clustering

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- **Entropy**
- Problem
  - Good subspace Clustering
- Algorithm
  - Property
  - Apriori

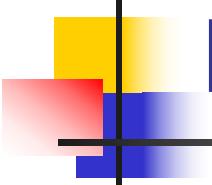


# Entropy

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- Example

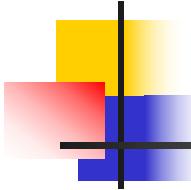
- Suppose we have a horse race with eight horses taking part.
- Assume that the probabilities of winning for the eight horses are
- $\frac{1}{2}, \frac{1}{4}, \frac{1}{8}, \frac{1}{16}, \frac{1}{64}, \frac{1}{64}, \frac{1}{64}, \frac{1}{64}$



# Entropy

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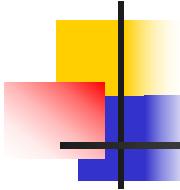
- Suppose we want to send a message to another person indicating which horse won the race. One method is to send a 3 bit string to denote the index of the winning horse



# Entropy

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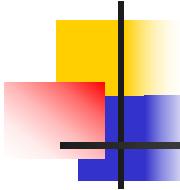
- Another method is to use a variable length coding set (i.e. 0, 10, 110, 1110, 111100, 111101, 111110, 111111) to represent the eight horses.
- The average description length is
  - $\frac{1}{2} \log \frac{1}{2} - \frac{1}{4} \log \frac{1}{4} - \frac{1}{8} \log \frac{1}{8} - \frac{1}{16} \log \frac{1}{16} - 4 \times \frac{1}{64} \log \frac{1}{64}$
  - = 2 bits



# Entropy

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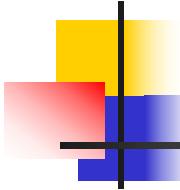
- The **entropy** is a way to measure the amount of information.
- The smaller the entropy (viewed as the average length of description length in the above example), the more informative we have.



# Entropy

---

- Assume that the probabilities of winning for the eight horses are
- $(1/8, 1/8, 1/8, 1/8, 1/8, 1/8, 1/8, 1/8)$
- Entropy of the horse race:
- $$\begin{aligned} H(X) &= -(1/8 \log 1/8) \times 8 \\ &= 3 \text{ bits} \end{aligned}$$



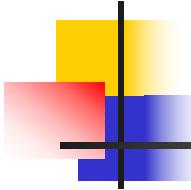
# Entropy

---

- Assume that the probabilities of winning for the eight horses are
- $(1, 0, 0, 0, 0, 0, 0, 0)$
- Entropy of the horse race:
- $$\begin{aligned} H(X) &= -1 \log 1 - 7(0 \log 0) \\ &= 0 \text{ bits} \end{aligned}$$

We use the convention that  $0 \log 0 = 0$

justified by continuity since  $x \log x \rightarrow 0$  as  $x \rightarrow 0$



# Entropy

---

- Let  $A$  be the set of possible outcomes of random variable  $X$ .
- Let  $p(x)$  be the probability mass function of the random variable  $X$ .
- The entropy  $H(X)$  of a discrete random variable  $X$  is

$$H(X) = - \sum_{x \in A} p(x) \log p(x)$$

Unit: bit

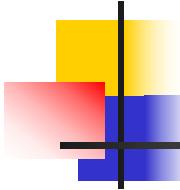
CSI If the base of log is 2, the unit for entropy is bit.



# Entropy

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- $H(X) \geq 0$
- Because  $0 \leq p(x) \leq 1$

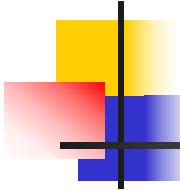


## More variables

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- When there are more than one variable, we can calculate the **joint entropy** to measure their uncertainty
- $X_i$ : the i-th random variable
- $A_i$ : the set of possible outcomes of  $X_i$
- Entropy:

$$H(X_1, \dots, X_n) = - \sum_{x_1 \in A_1} \dots \sum_{x_n \in A_n} p(x_1, \dots, x_n) \log p(x_1, \dots, x_n)$$



## More variables

---

$X_1 \setminus X_2$	1	2
1	$1/4$	$1/2$
2	0	$1/4$

$$p(1, 1) = \frac{1}{4}$$

$$p(1, 2) = \frac{1}{2}$$

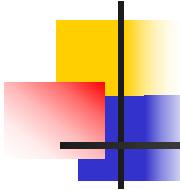
$$p(2, 1) = 0$$

$$p(2, 2) = \frac{1}{4}$$

$$H(X_1, X_2)$$

$$= -\frac{1}{4} \log \frac{1}{4} - \frac{1}{2} \log \frac{1}{2} - 0 \log 0 - \frac{1}{4} \log \frac{1}{4}$$

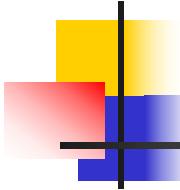
$$= 1.5 \text{ bits}$$



# Entropy-Based Method for Subspace Clustering

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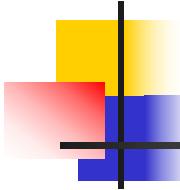
- Entropy
- Problem
  - Good subspace Clustering
- Algorithm
  - Property
  - Apriori



# Subspace Clustering

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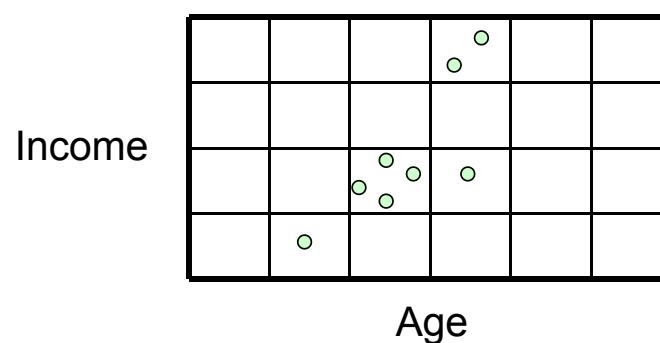
- We divide each dimension into intervals of equal length  $\Delta$ , so the subspace is partitioned into a grid.



# Subspace Clustering

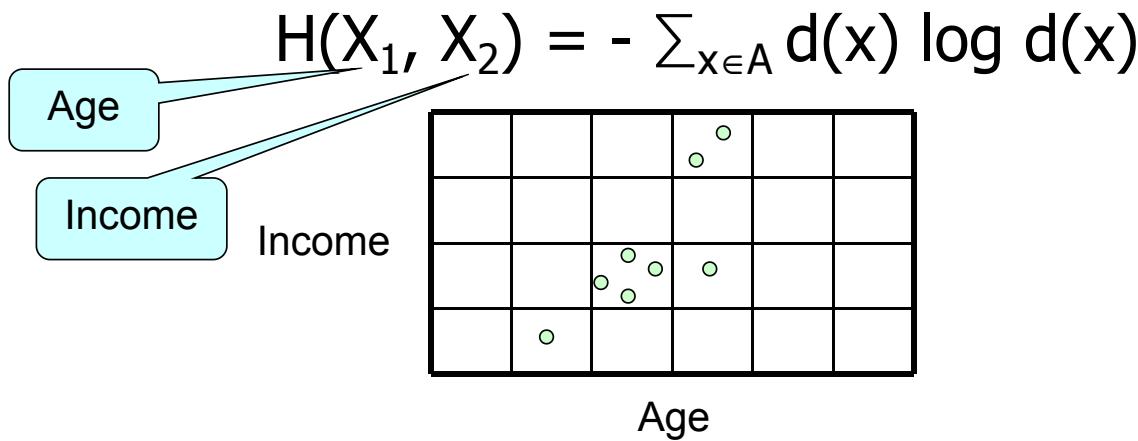
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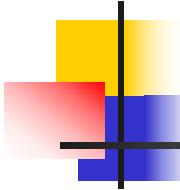
- We divide each dimension into intervals of equal length  $\Delta$ , so the subspace is partitioned into a grid.



# Subspace Clustering

- Let  $A$  be the set of all cells.
- $d(x)$  be the density of a cell  $x$  in terms of the percentage of data contained in  $x$ .
- We can define the entropy to be:





# Subspace Clustering

---

- Let  $A$  be the set of all cells.
- $d(x)$  be the density of a cell  $x$  in terms of the percentage of data contained in  $x$ .
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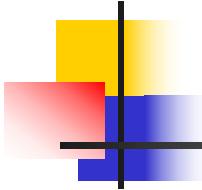
$$H(X_1, X_2) = - \sum_{x \in A} d(x) \log d(x)$$

Age

Income

Given a parameter  $\omega$ ,  
k dimensions (or random variables) are said  
to have **good clustering** if

$$H(X_1, X_2, \dots, X_k) \leq \omega$$



# Subspace Clustering

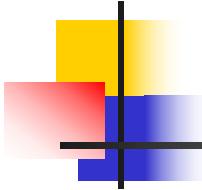
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- **Problem:** We want to find all subspaces with good clustering.

e.g., we want to find sub-spaces with entropy  $\leq 0.2$

Given a parameter  $\omega$ ,  
k dimensions (or random variables) are said  
to have **good clustering** if

$$H(X_1, X_2, \dots, X_k) \leq \omega$$

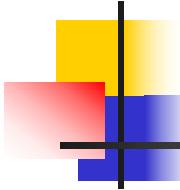


# Conditional Entropy

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- The conditional entropy  $H(Y|X)$  is defined as

$$H(Y|X) = \sum_{x \in A} p(x)H(Y|X = x)$$



# Conditional Entropy

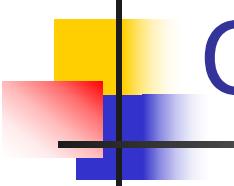
X\Y	1	2
1	0	3/4
2	1/8	1/8

$$H(Y|X=1) = 0 \text{ bit}$$

$$H(Y|X=2) = 1 \text{ bit}$$

$$\begin{aligned} H(Y|X) &= \frac{3}{4} \times H(Y|X=1) + \frac{1}{4} \times H(Y|X=2) \\ &= 0.25 \text{ bit} \end{aligned}$$

$$H(Y|X) = - \sum_{x \in A} \sum_{y \in B} p(x, y) \log p(y|x)$$



# Conditional Entropy

---

- A : a set of possible outcomes of random variable X
- B : a set of possible outcomes of random variable Y
- $H(Y|X) = - \sum_{x \in A} \sum_{y \in B} p(x, y) \log p(y|x)$

$$H(Y|X) = - \sum_{x \in A} \sum_{y \in B} p(x, y) \log p(y|x)$$

# Conditional Entropy

X\Y	1	2
1	0	3/4
2	1/8	1/8

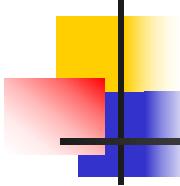
$$p(Y = 1 | X = 1) = 0$$

$$p(Y = 2 | X = 1) = 1$$

$$p(Y = 1 | X = 2) = \frac{1}{2}$$

$$p(Y = 2 | X = 2) = \frac{1}{2}$$

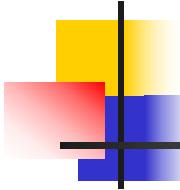
$$\begin{aligned} H(Y|X) &= -0 \log 0 - \frac{3}{4} \log 1 - 1/8 \log \frac{1}{2} - 1/8 \log \frac{1}{2} \\ &= 0.25 \text{ bit} \end{aligned}$$



## Chain Rule

---

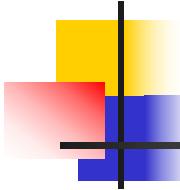
- $H(X, Y) = H(X) + H(Y | X)$



## Chain Rule

---

- $H(X, Y) = H(X) + H(Y | X)$
- $$\begin{aligned} H(X_1, \dots, X_{k-1}, X_k) \\ = H(X_1, \dots, X_{k-1}) + H(X_k | X_1, \dots, X_{k-1}) \end{aligned}$$



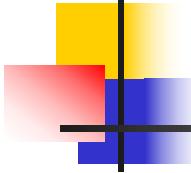
# Entropy-Based Method for Subspace Clustering

---

- Entropy
- Problem
  - Good subspace Clustering

## Algorithm

- Property
- Apriori



# Property

Given a parameter  $\omega$ ,  
k dimensions (or random variables) are said  
to have **good clustering** if

$$H(X_1, X_2, \dots, X_k) \leq \omega$$

- **Lemma:** If a k-dimensional subspace  $X_1, \dots, X_k$  has good clustering,  
then each of the  $(k-1)$ -dimensional  
projections of this space has also good  
clustering.

Proof: Since the subspace  $X_1, \dots, X_k$  has good clustering,

$$H(X_1, \dots, X_k) \leq \omega$$

Consider a  $(k-1)$ -dimensional projections, say  $X_1, \dots, X_{k-1}$  :

$$\begin{aligned} H(X_1, \dots, X_{k-1}) &\leq H(X_1, \dots, X_{k-1}) + H(X_k | X_1, \dots, X_{k-1}) \\ &= H(X_1, \dots, X_k) \end{aligned}$$

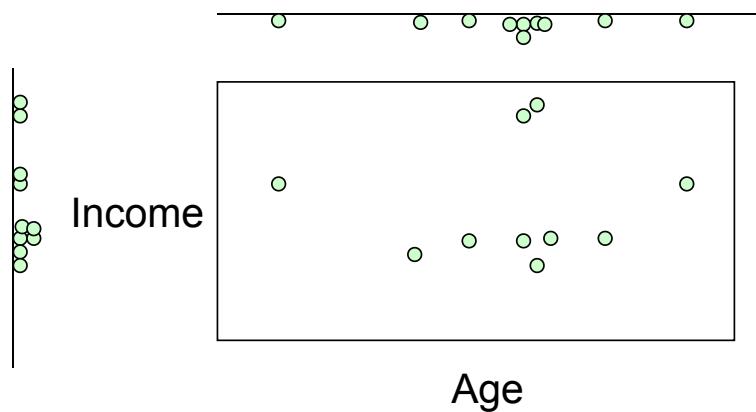


Suppose we want to find sub-spaces with entropy  $\leq 0.2$

- We can make use of apriori approach to solve the problem

# Apriori

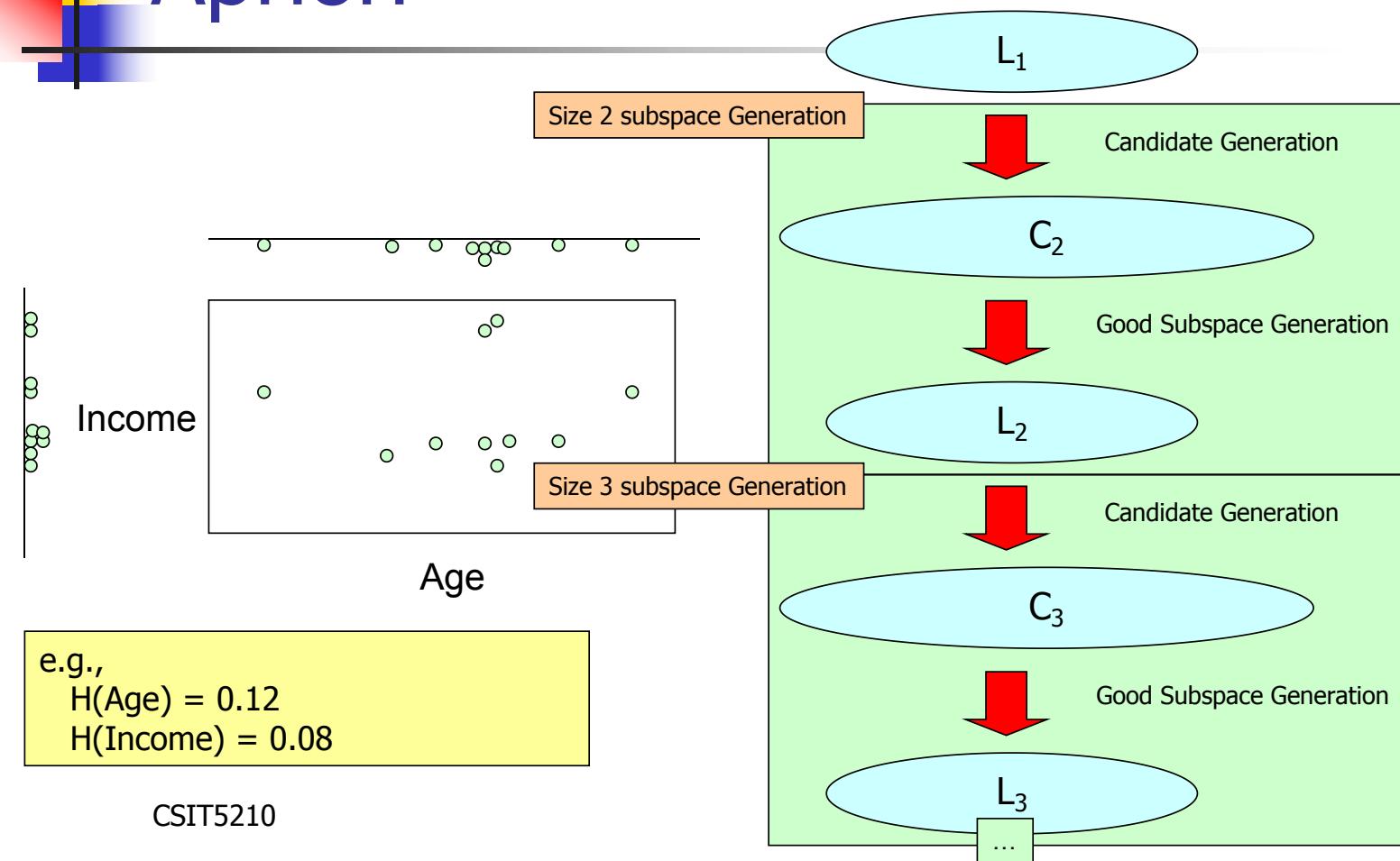
Suppose we want to find sub-spaces with entropy  $\leq 0.2$



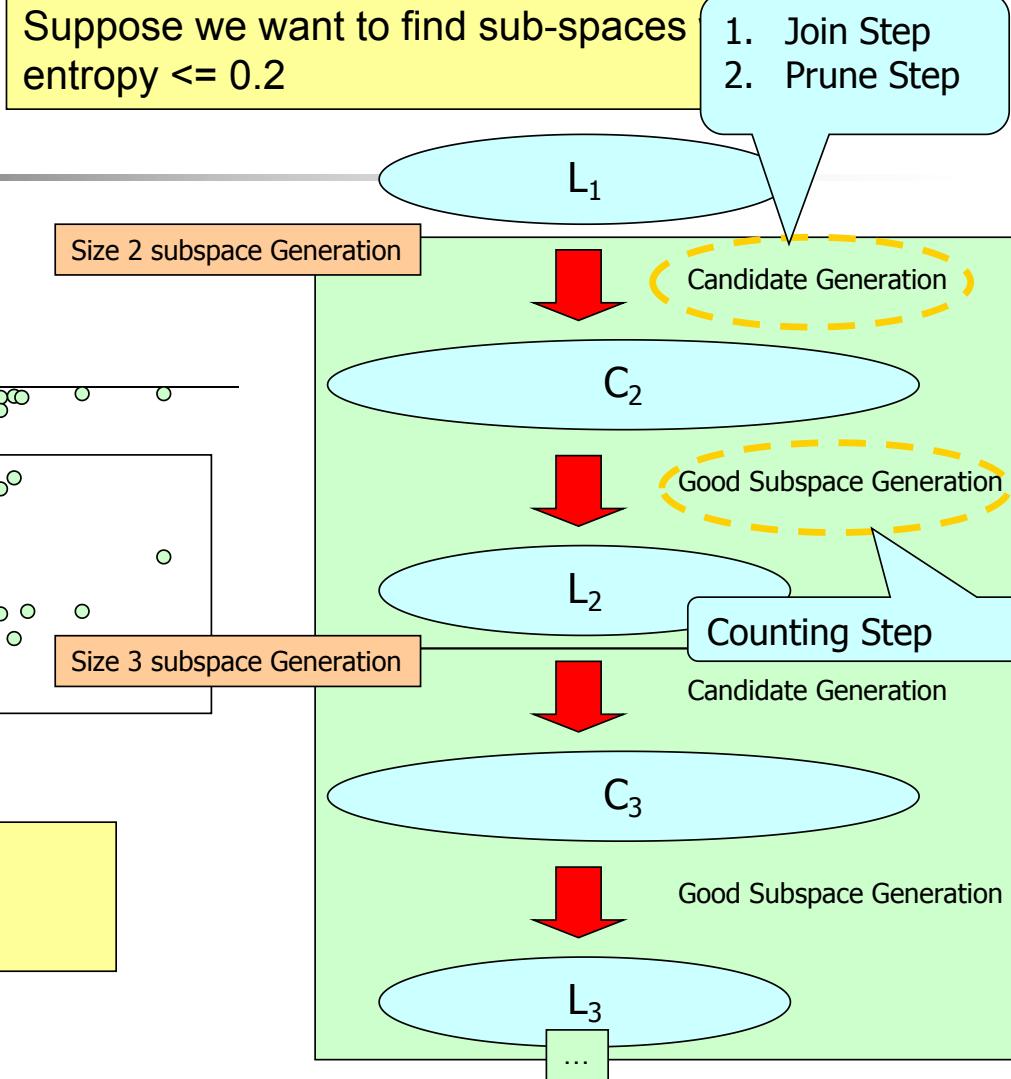
e.g.,  
 $H(\text{Age}) = 0.12$   
 $H(\text{Income}) = 0.08$

# Apriori

Suppose we want to find sub-spaces with entropy  $\leq 0.2$



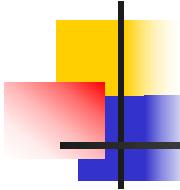
# Apriori





Suppose we want to find sub-spaces with entropy  $\leq 0.2$

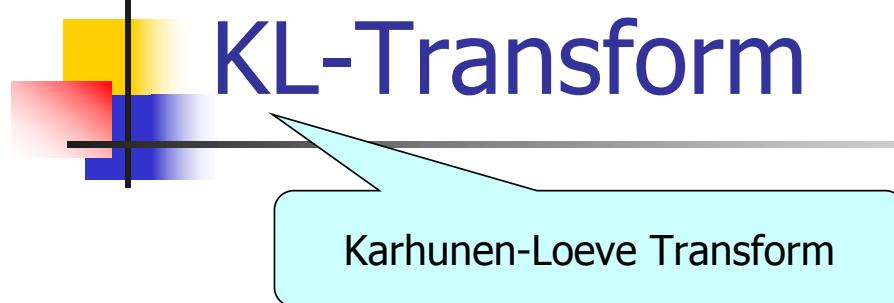
- After finding the subspaces with entropy  $\leq 0.2$ ,
- We can find the real clusters by existing methods (e.g., k-mean) in each of the subspaces found.



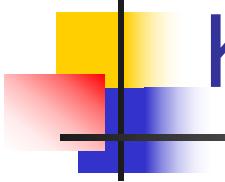
# Subspace Clustering Methods

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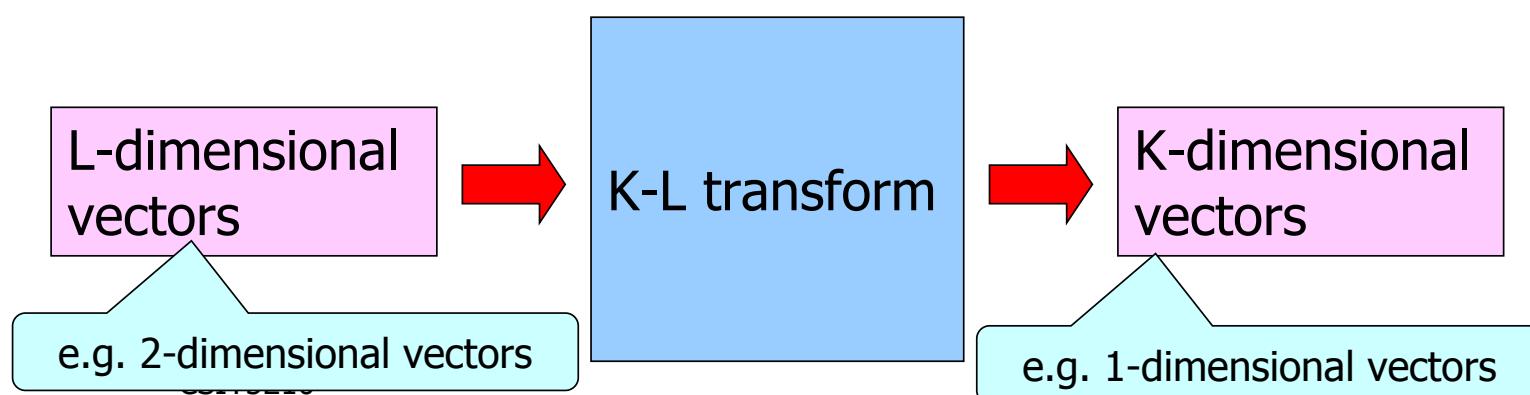
- Dense Unit-based Method
- Entropy-Based Method
- Transformation-Based Method

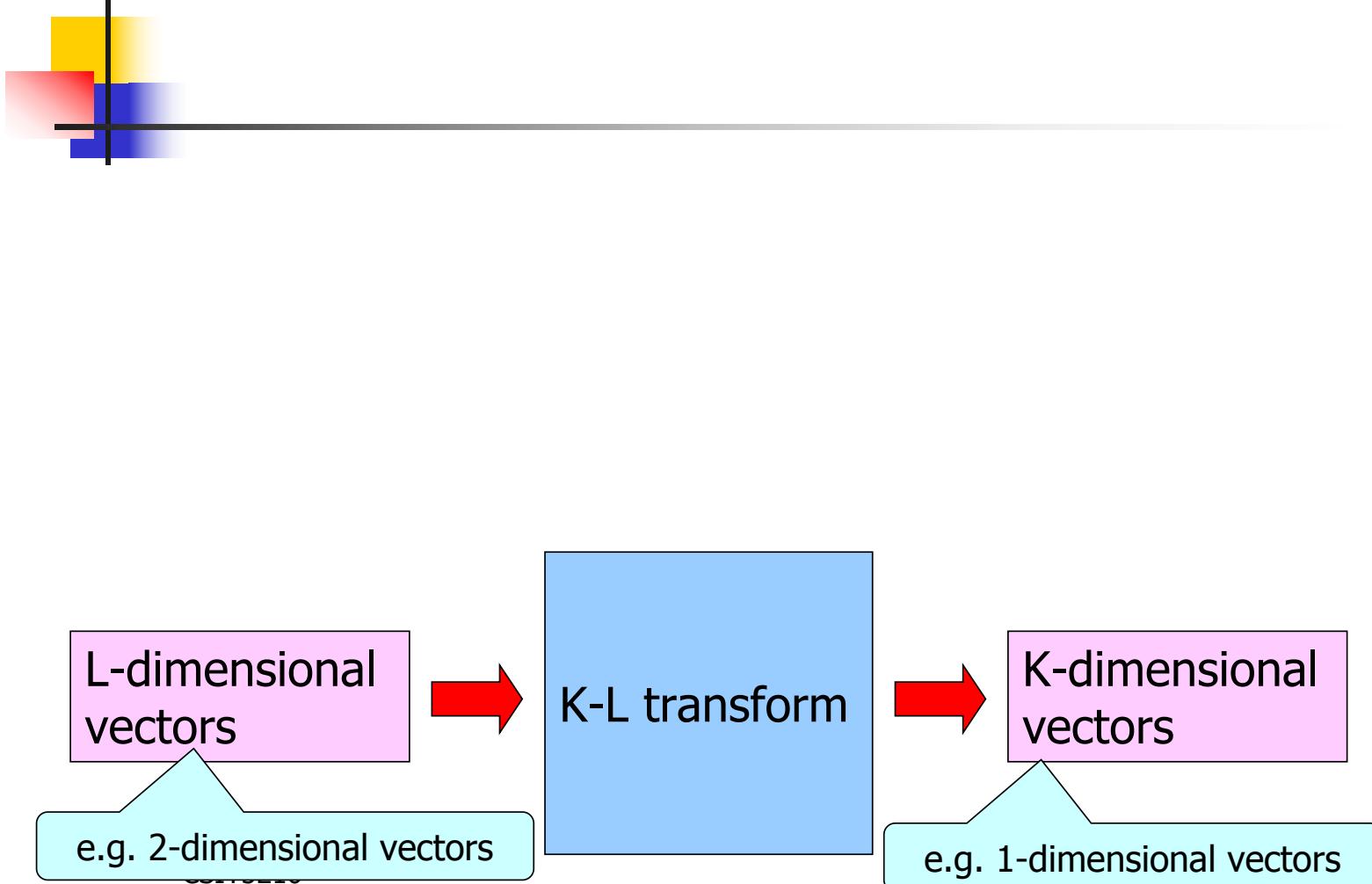


- The two previous approaches find the sub-space in the original dimensions
- KL-Transform “transforms” the data points from the original dimensions into other dimensions



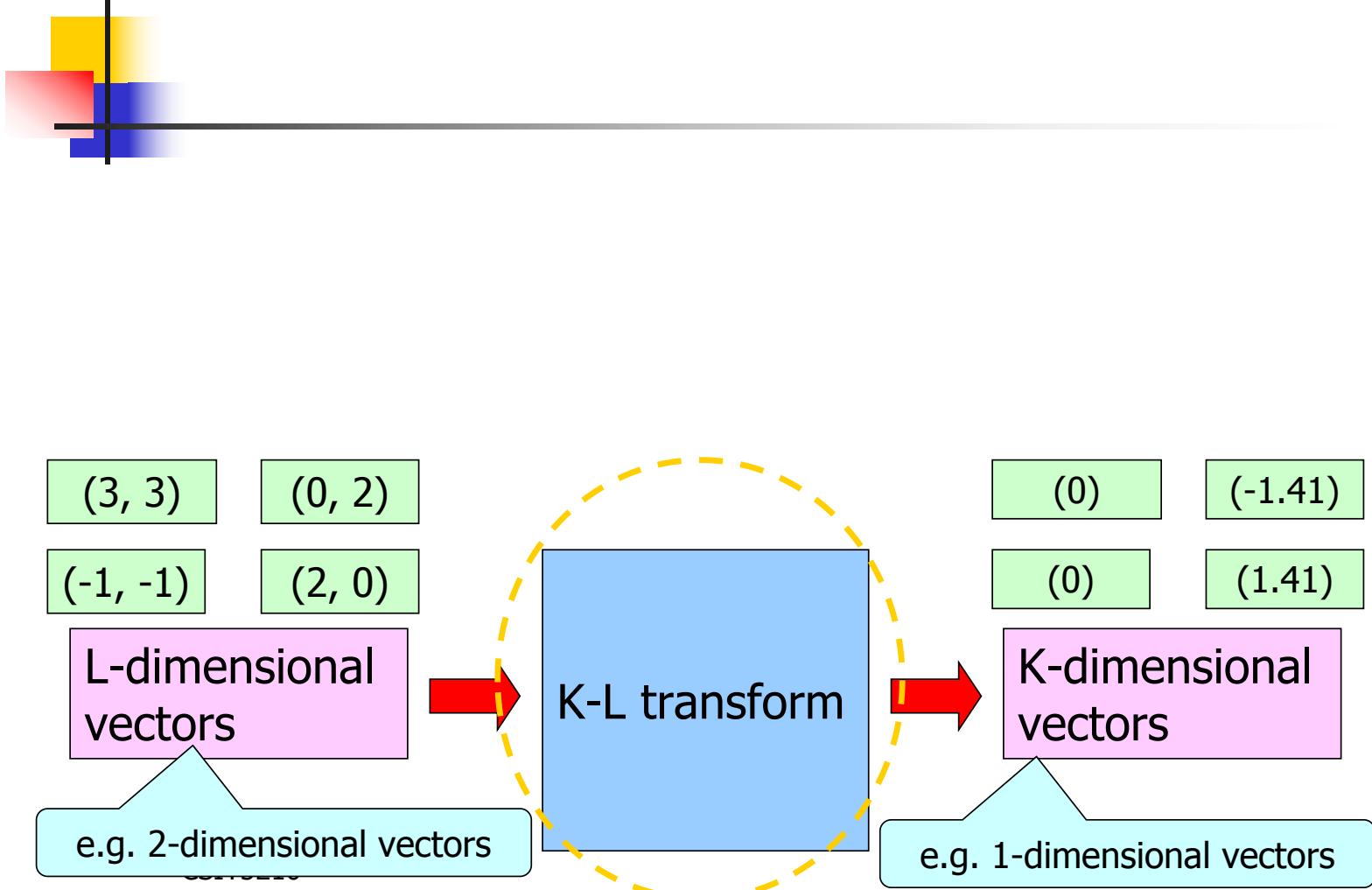
# KL-Transform

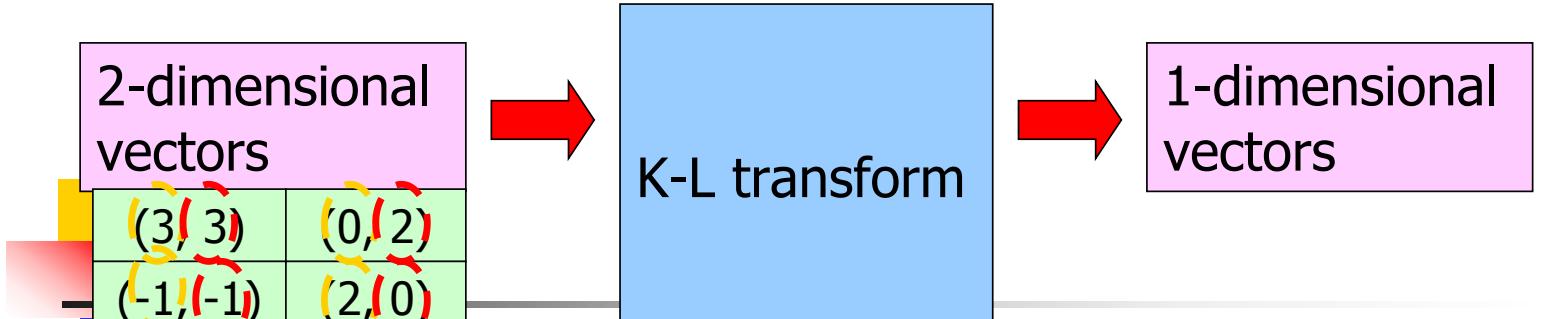






<b>Step 1</b>	For each dimension $i$ , calculate the mean $e_i$ (expected value) For each L-dimensional data $\{x_1, x_2, \dots, x_L\}$ find $\{x_1 - e_1, x_2 - e_2, \dots, x_L - e_L\}$
<b>Step 2</b>	Obtain the covariance matrix $\Sigma$
<b>Step 3</b>	Find the eigenvalues and eigenvectors of $\Sigma$ Choose the eigenvectors of unit lengths
<b>Step 4</b>	Arrange the eigenvectors in descending order of the eigenvalues
<b>Step 5</b>	Transform the given L-dimensional vectors by eigenvector matrix
<b>Step 6</b>	For each “transformed” L-dimensional vector, keep only the K values $\{y_1, y_2, \dots, y_K\}$ corresponding to the smallest K eigenvalues.





### Step 1

For each dimension  $i$ ,  
calculate the mean  $e_i$  (expected value)  
For each  $L$ -dimensional data  $\{x_1, x_2, \dots, x_L\}$   
find  $\{x_1 - e_1, x_2 - e_2, \dots, x_L - e_L\}$

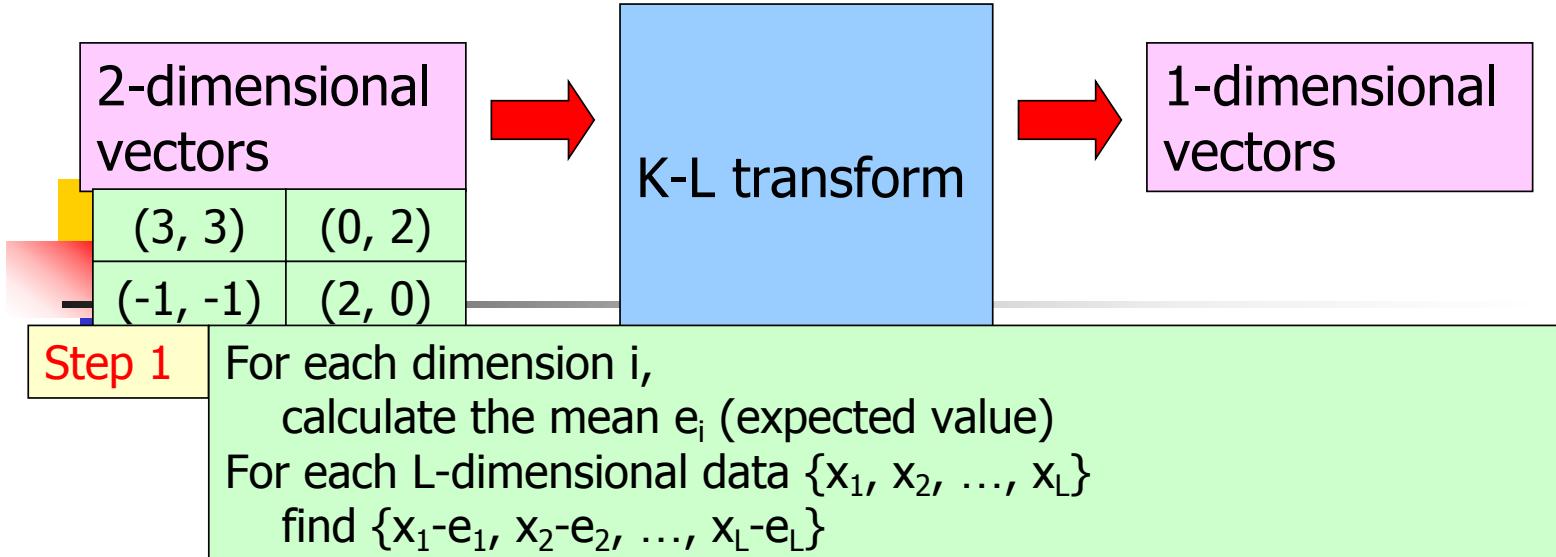
For **dimension 1**,

$$\text{mean} = (3 + 0 + (-1) + 2)/4 = 4/4 = 1$$

For **dimension 2**,

$$\text{mean} = (3 + 2 + (-1) + 0)/4 = 4/4 = 1$$

$$\text{mean vector} = (1, 1)$$



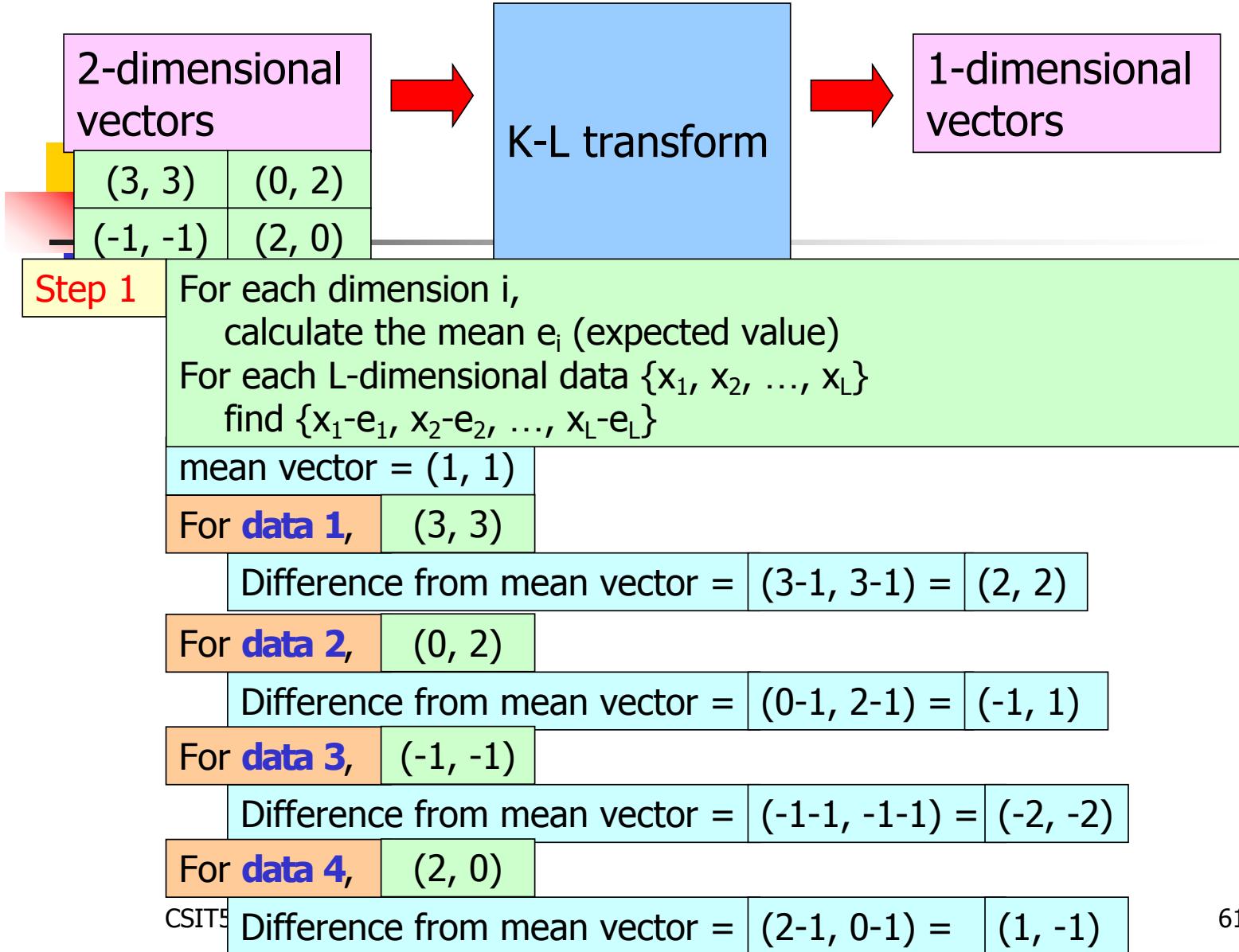
For **dimension 1**,

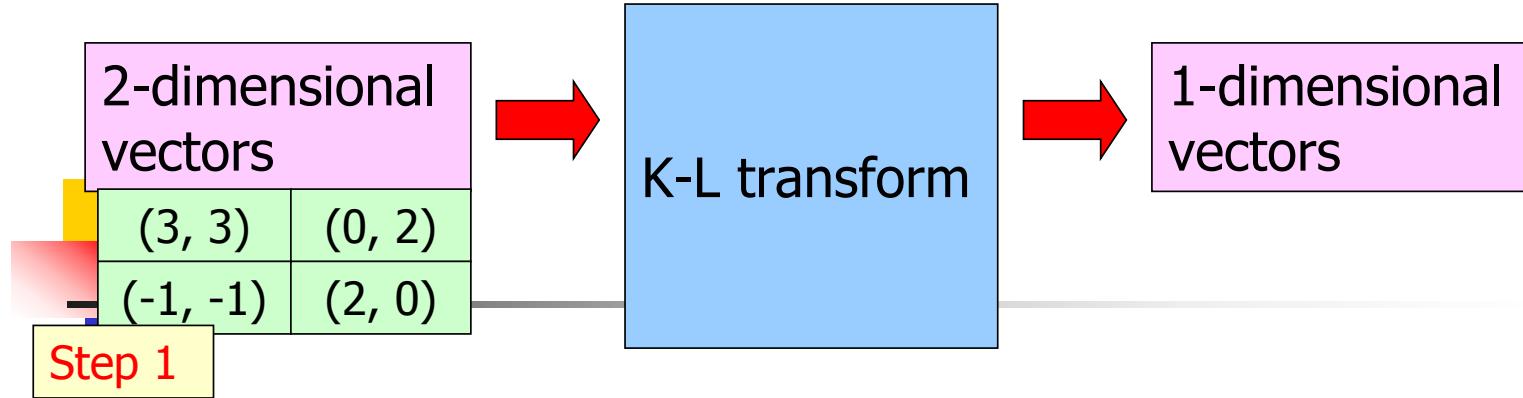
$$\text{mean} = (3 + 0 + (-1) + 2)/4 = 4/4 = 1$$

For **dimension 2**,

$$\text{mean} = (3 + 2 + (-1) + 0)/4 = 4/4 = 1$$

$$\text{mean vector} = (1, 1)$$





mean vector = (1, 1)

For **data 1**, (3, 3)

Difference from mean vector = (2, 2)

For **data 2**, (0, 2)

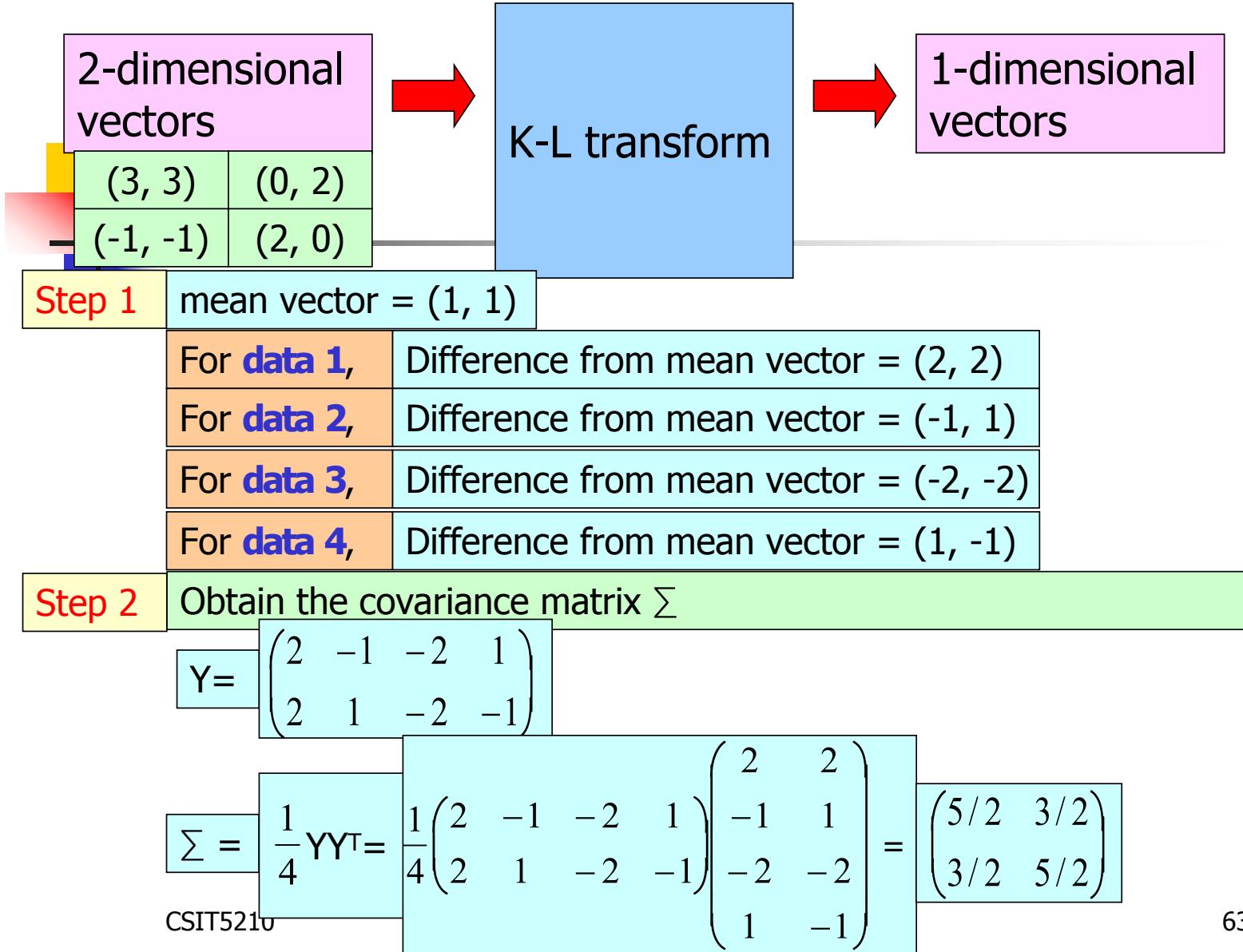
Difference from mean vector = (-1, 1)

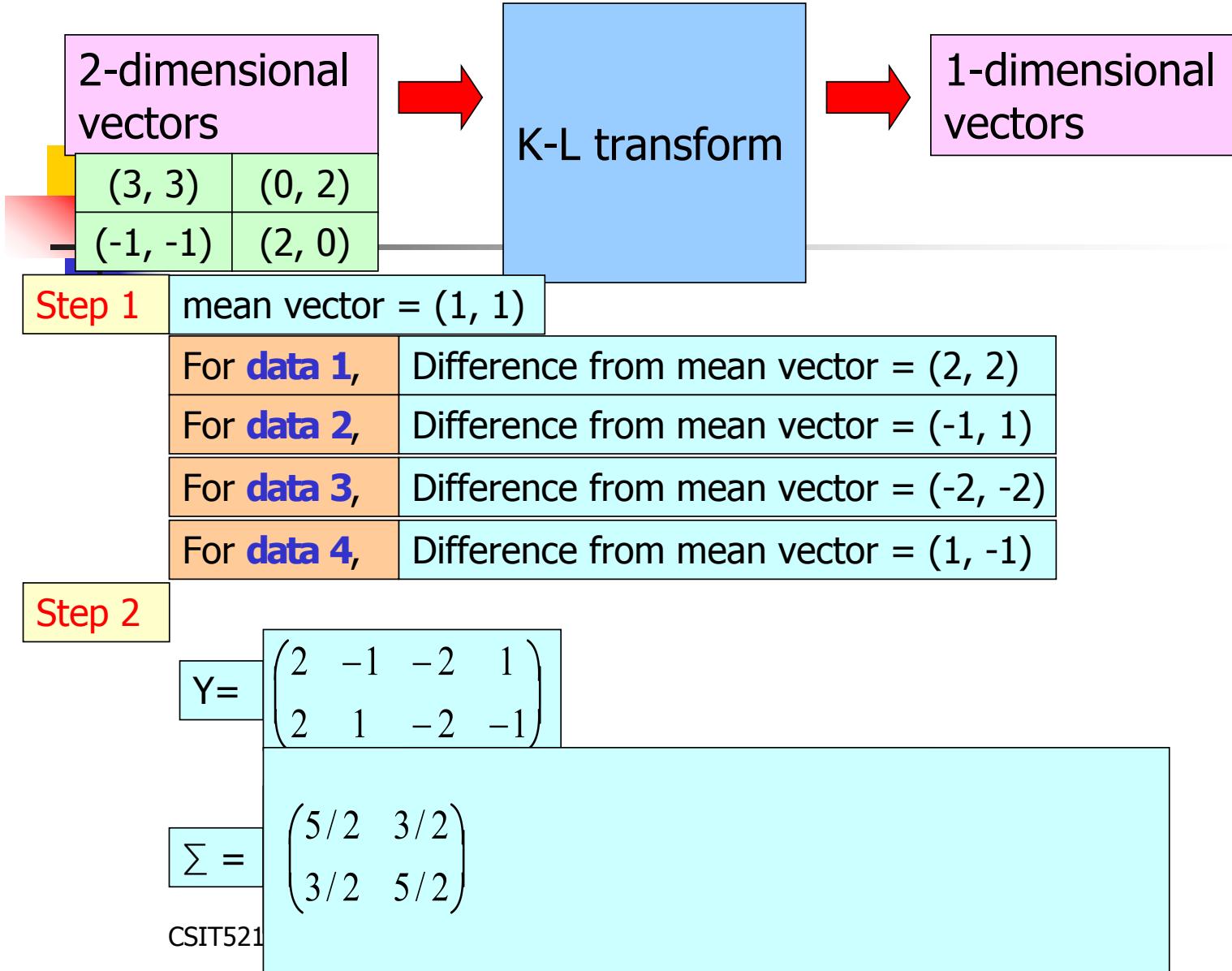
For **data 3**, (-1, -1)

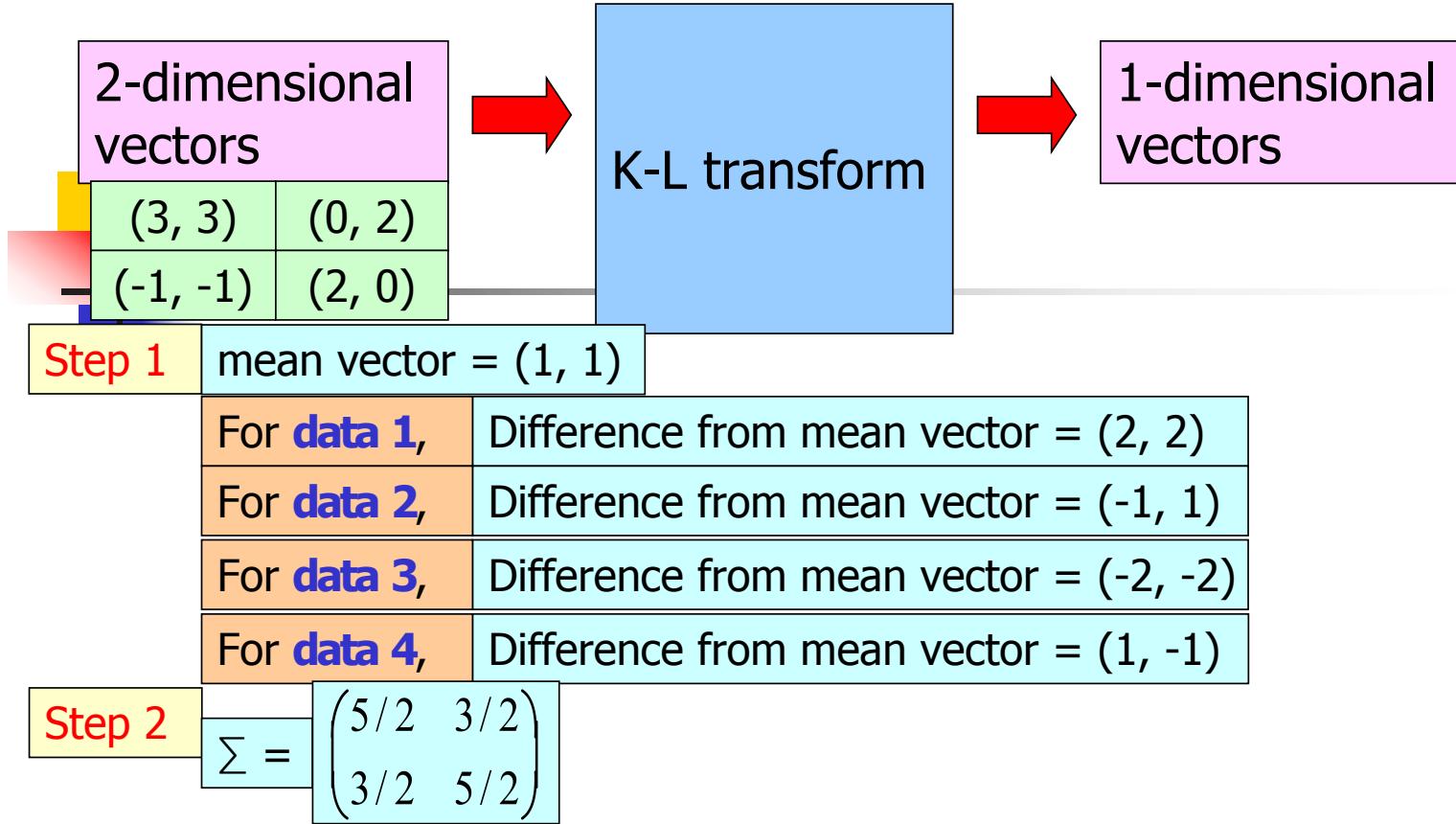
Difference from mean vector = (-2, -2)

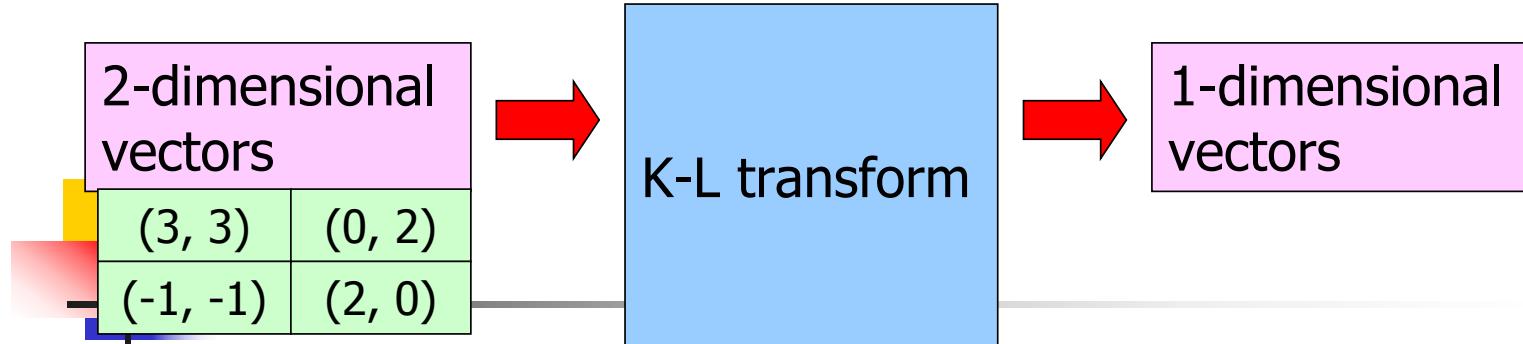
For **data 4**, (2, 0)

Difference from mean vector = (1, -1)









Step 2

$$\Sigma = \begin{pmatrix} 5/2 & 3/2 \\ 3/2 & 5/2 \end{pmatrix}$$



Step 2

$$\Sigma = \begin{pmatrix} 5/2 & 3/2 \\ 3/2 & 5/2 \end{pmatrix}$$

Step 3

Find the eigenvalues and eigenvectors of  $\Sigma$   
Choose the eigenvectors of unit lengths

$$\begin{vmatrix} 5/2 - \lambda & 3/2 \\ 3/2 & 5/2 - \lambda \end{vmatrix} = 0$$

$$(5/2 - \lambda)^2 - (3/2)^2 = 0$$

$$25/4 - 5\lambda + \lambda^2 - 9/4 = 0$$

$$4 - 5\lambda + \lambda^2 = 0$$

$$(\lambda - 4)(\lambda - 1) = 0$$

$$\lambda = 4 \quad \text{or} \quad \lambda = 1$$



Step 2

$$\Sigma = \begin{pmatrix} 5/2 & 3/2 \\ 3/2 & 5/2 \end{pmatrix}$$

Step 3

$$\begin{vmatrix} 5/2 - \lambda & 3/2 \\ 3/2 & 5/2 - \lambda \end{vmatrix} = 0$$

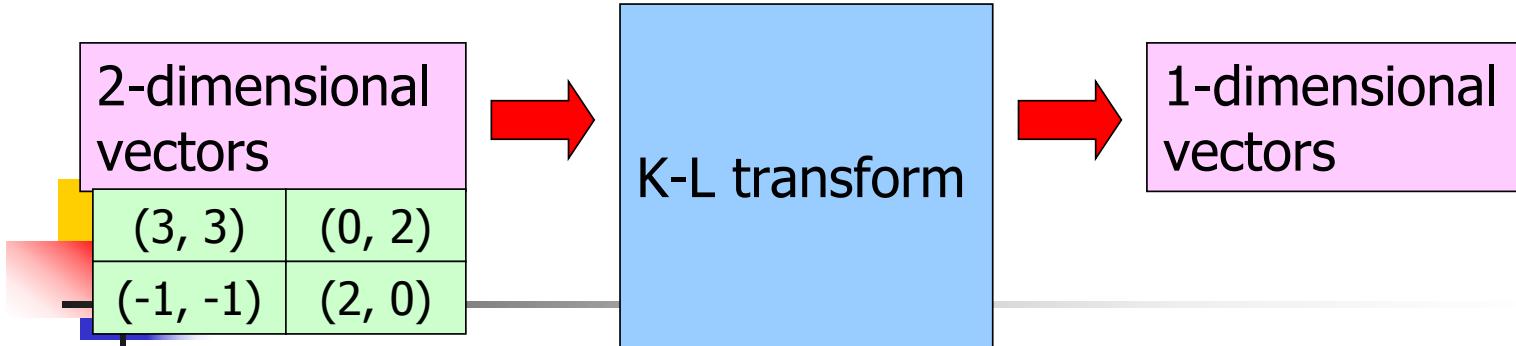
$$(5/2 - \lambda)^2 - (3/2)^2 = 0$$

$$25/4 - 5\lambda + \lambda^2 - 9/4 = 0$$

$$4 - 5\lambda + \lambda^2 = 0$$

$$(\lambda - 4)(\lambda - 1) = 0$$

$$\lambda = 4 \quad \text{or} \quad \lambda = 1$$



Step 2

$$\Sigma = \begin{pmatrix} 5/2 & 3/2 \\ 3/2 & 5/2 \end{pmatrix}$$

Step 3

$$\lambda = 4 \quad \text{or} \quad \lambda = 1$$

When  $\lambda = 4$

$$\begin{pmatrix} 5/2-4 & 3/2 \\ 3/2 & 5/2-4 \end{pmatrix} \begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} 0 \\ 0 \end{pmatrix}$$

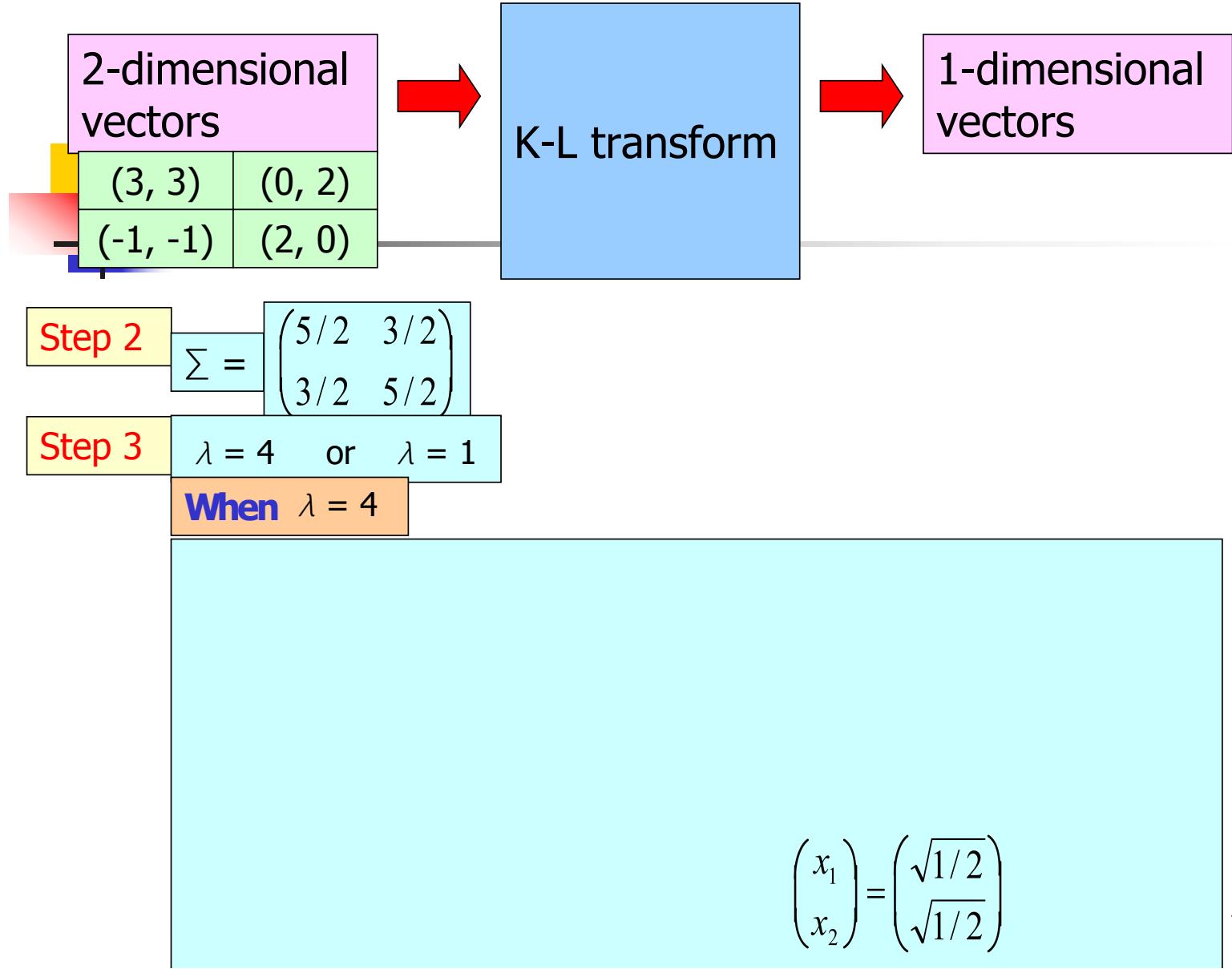
$$\begin{pmatrix} -3/2 & 3/2 \\ 3/2 & -3/2 \end{pmatrix} \begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} 0 \\ 0 \end{pmatrix}$$

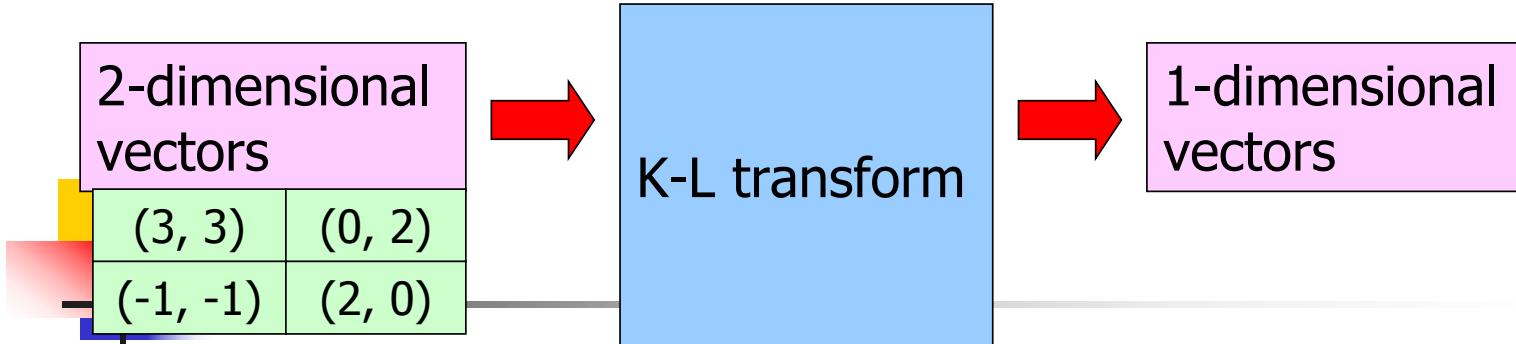
$$\begin{pmatrix} -1 & 1 \\ 1 & -1 \end{pmatrix} \begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} 0 \\ 0 \end{pmatrix}$$

$$\begin{aligned} x_1 - x_2 &= 0 \\ x_1 &= x_2 \\ \begin{pmatrix} x_1 \\ x_2 \end{pmatrix} &= \begin{pmatrix} a \\ a \end{pmatrix} \quad \text{where } a \in R \end{aligned}$$

We choose the eigenvector of unit length

$$\begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} \sqrt{1/2} \\ \sqrt{1/2} \end{pmatrix}$$





Step 2

$$\Sigma = \begin{pmatrix} 5/2 & 3/2 \\ 3/2 & 5/2 \end{pmatrix}$$

Step 3

$$\lambda = 4 \quad \text{or} \quad \lambda = 1$$

When  $\lambda = 4$

$$\begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} \sqrt{1/2} \\ \sqrt{1/2} \end{pmatrix}$$

When  $\lambda = 1$

$$\begin{pmatrix} 5/2-1 & 3/2 \\ 3/2 & 5/2-1 \end{pmatrix} \begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} 0 \\ 0 \end{pmatrix}$$

$$\begin{pmatrix} 3/2 & 3/2 \\ 3/2 & 3/2 \end{pmatrix} \begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} 0 \\ 0 \end{pmatrix}$$

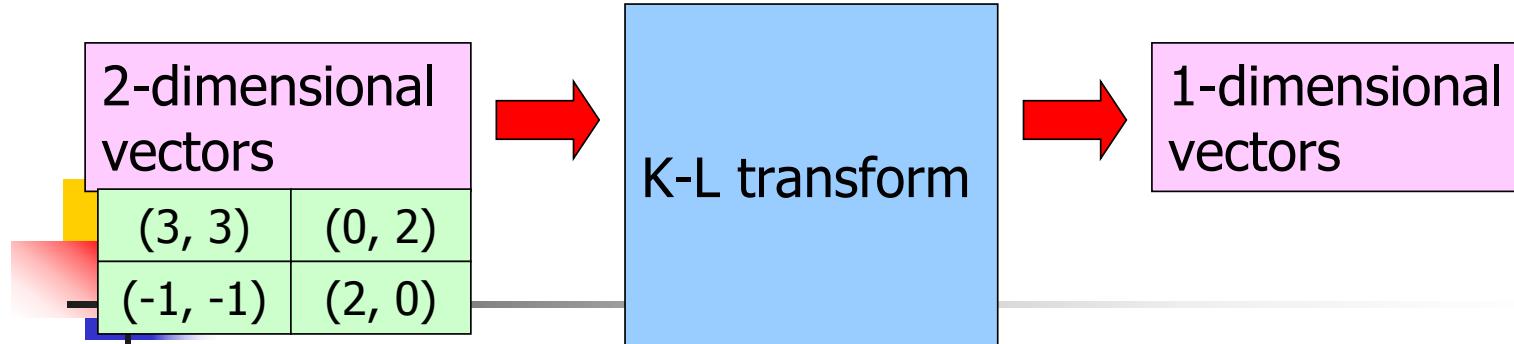
$$\begin{pmatrix} 1 & 1 \\ 1 & 1 \end{pmatrix} \begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} 0 \\ 0 \end{pmatrix}$$

$$\begin{aligned} x_1 + x_2 &= 0 \\ x_1 &= -x_2 \end{aligned}$$

$$\begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} a \\ -a \end{pmatrix} \quad \text{where } a \in R$$

We choose the eigenvector of unit length

$$\begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} \sqrt{1/2} \\ -\sqrt{1/2} \end{pmatrix}$$



Step 2

$$\Sigma = \begin{pmatrix} 5/2 & 3/2 \\ 3/2 & 5/2 \end{pmatrix}$$

Step 3

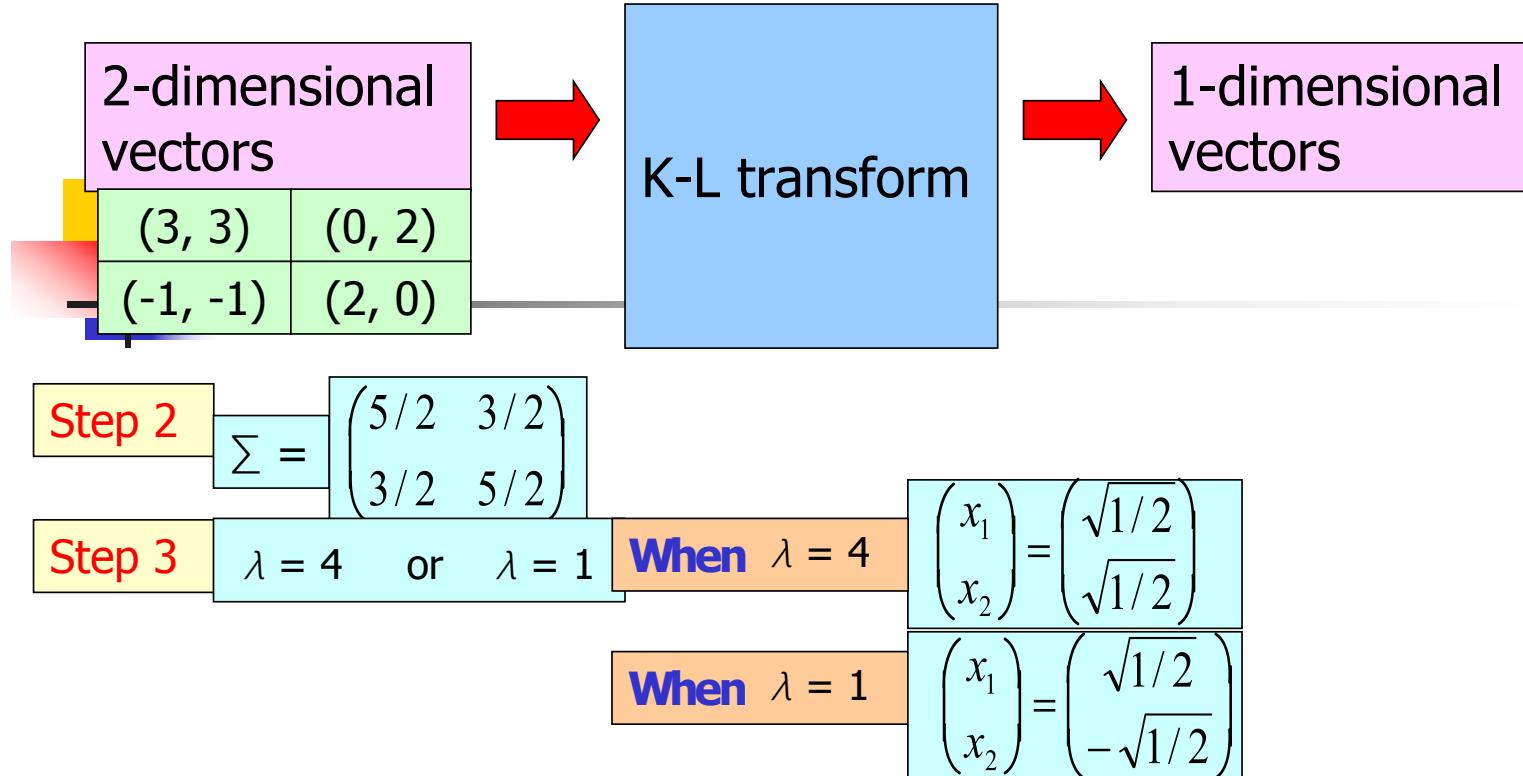
$\lambda = 4$  or  $\lambda = 1$

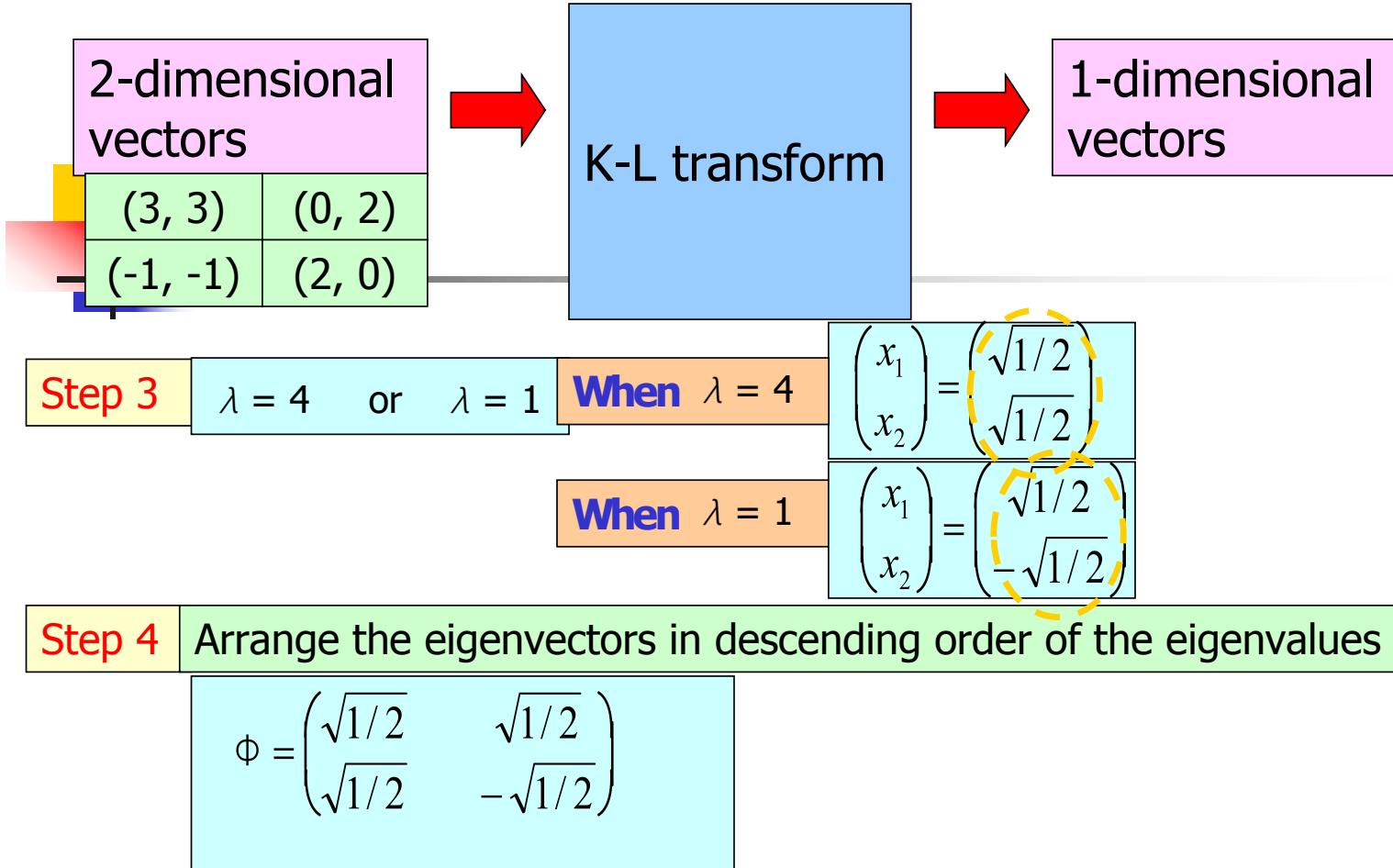
When  $\lambda = 4$

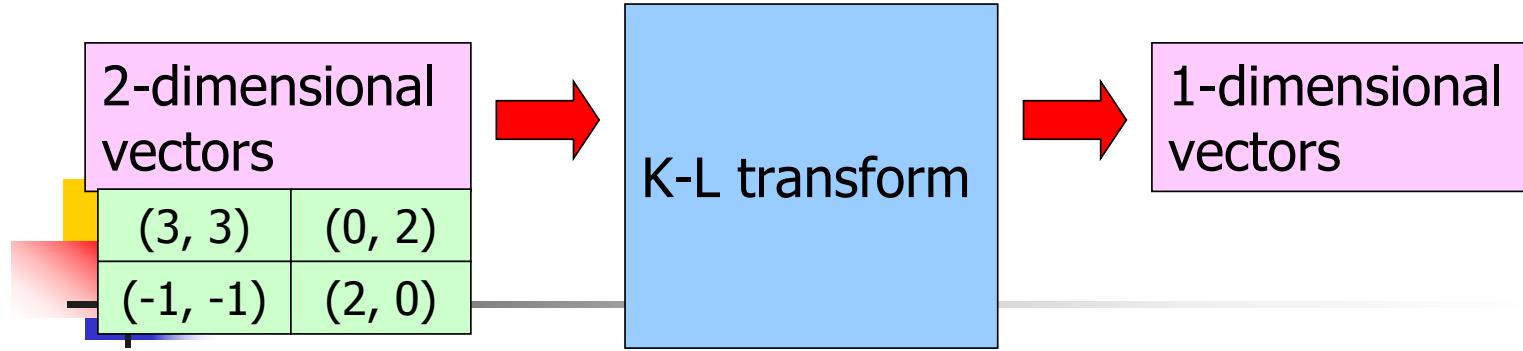
$$\begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} \sqrt{1/2} \\ \sqrt{1/2} \end{pmatrix}$$

When  $\lambda = 1$

$$\begin{pmatrix} x_1 \\ x_2 \end{pmatrix} = \begin{pmatrix} \sqrt{1/2} \\ -\sqrt{1/2} \end{pmatrix}$$



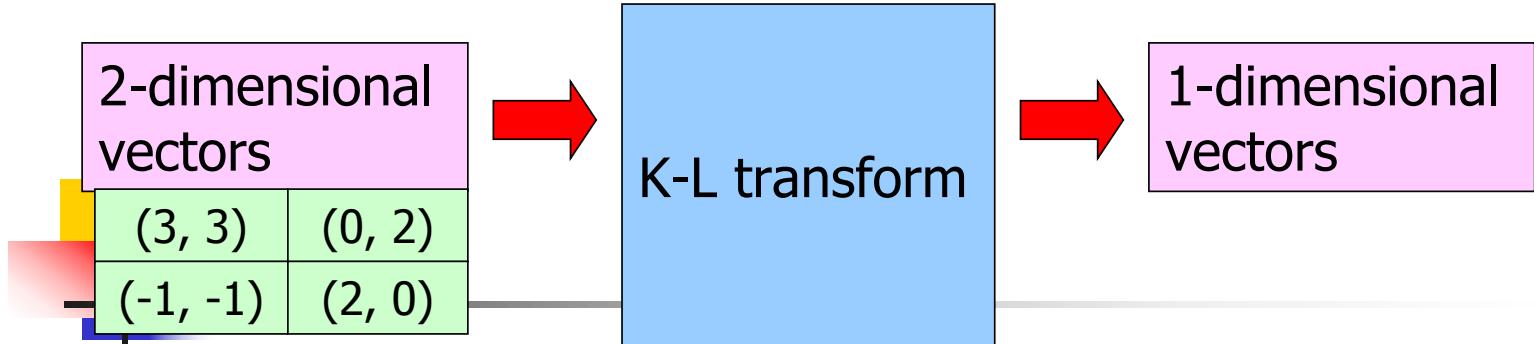




Step 4

Arrange the eigenvectors in descending order of the eigenvalues

$$\phi = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix}$$



Step 4

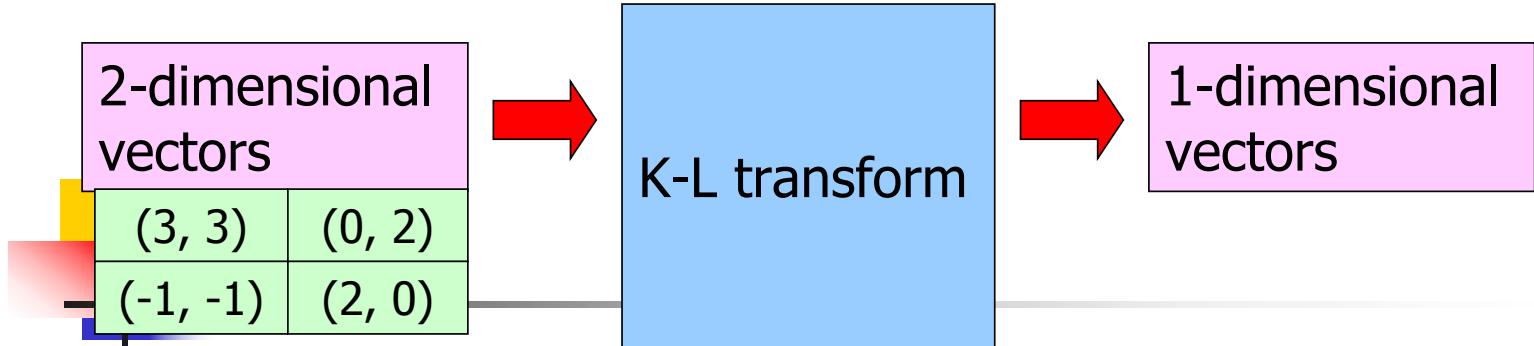
$$\Phi = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix}$$

Step 5

Transform the given L-dimensional vectors by eigenvector matrix

$$Y = \Phi^T X$$

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} X$$



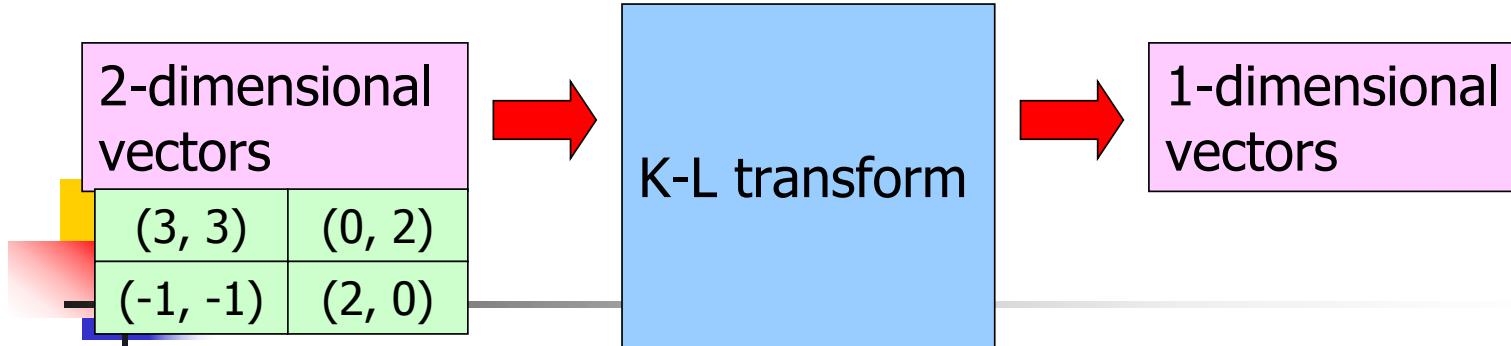
Step 4

$$\Phi = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix}$$

Step 5

Transform the given L-dimensional vectors by eigenvector matrix

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} X$$



Step 4

$$\Phi = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix}$$

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} X$$

Step 5

Transform the given L-dimensional vectors by eigenvector matrix



Step 4

$$\Phi = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix}$$

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} X$$

Step 5

Transform the given L-dimensional vectors by eigenvector matrix

For **data 1**,

(3, 3)

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} \begin{pmatrix} 3 \\ 3 \end{pmatrix} = \begin{pmatrix} 4.24 \\ 0 \end{pmatrix}$$

For **data 3**,

(-1, -1)

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} \begin{pmatrix} -1 \\ -1 \end{pmatrix} = \begin{pmatrix} -1.41 \\ 0 \end{pmatrix}$$

For **data 2**,

(0, 2)

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} \begin{pmatrix} 0 \\ 2 \end{pmatrix} = \begin{pmatrix} 1.41 \\ -1.41 \end{pmatrix}$$

For **data 4**,

(2, 0)

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} \begin{pmatrix} 2 \\ 0 \end{pmatrix} = \begin{pmatrix} 1.41 \\ 1.41 \end{pmatrix}$$



Step 4

$$\Phi = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix}$$

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} X$$

Step 5

Transform the given L-dimensional vectors by eigenvector matrix

For data 1,

(3, 3)

$$= \begin{pmatrix} 4.24 \\ 0 \end{pmatrix}$$

For data 3,

(-1, -1)

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} \begin{pmatrix} -1 \\ -1 \end{pmatrix} = \begin{pmatrix} -1.41 \\ 0 \end{pmatrix}$$

For data 2,

(0, 2)

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} \begin{pmatrix} 0 \\ 2 \end{pmatrix} = \begin{pmatrix} 1.41 \\ -1.41 \end{pmatrix}$$

For data 4,

(2, 0)

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} \begin{pmatrix} 2 \\ 0 \end{pmatrix} = \begin{pmatrix} 1.41 \\ 1.41 \end{pmatrix}$$



Step 4

$$\Phi = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix}$$

$$Y = \begin{pmatrix} \sqrt{1/2} & \sqrt{1/2} \\ \sqrt{1/2} & -\sqrt{1/2} \end{pmatrix} X$$

Step 5

Transform the given L-dimensional vectors by eigenvector matrix

For **data 1**,

(3, 3)



(4.24, 0)

For **data 2**,

(0, 2)



(1.41, -1.41)

For **data 3**,

(-1, -1)



(-1.41, 0)

For **data 4**,

(2, 0)



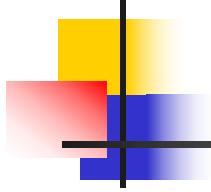
(1.41, 1.41)



**Step 6**

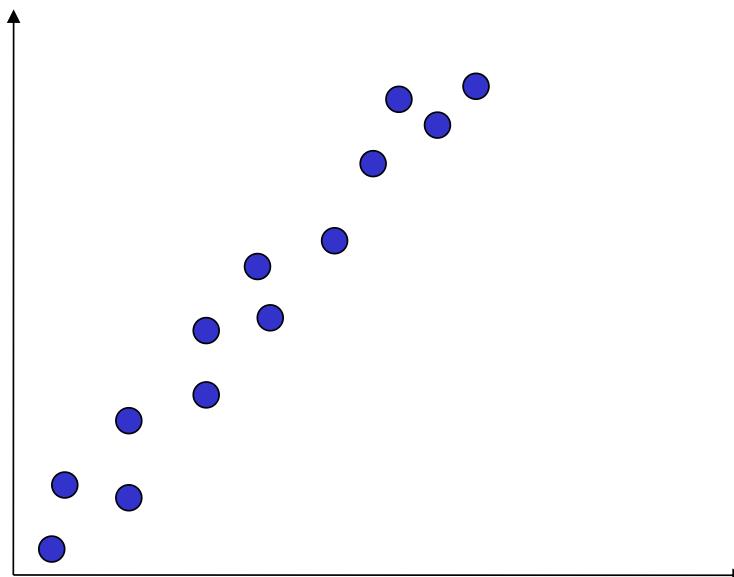
For each “transformed” L-dimensional vector, keep only the K values  $\{y_1, y_2, \dots, y_K\}$  corresponding to the smallest k eigenvalues.

For <b>data 1</b> ,	(3, 3)	$(4.24, 0)$	(0)
For <b>data 2</b> ,	(0, 2)	$(1.41, -1.41)$	(-1.41)
For <b>data 3</b> ,	(-1, -1)	$(-1.41, 0)$	(0)
For <b>data 4</b> ,	(2, 0)	$(1.41, 1.41)$	(1.41)

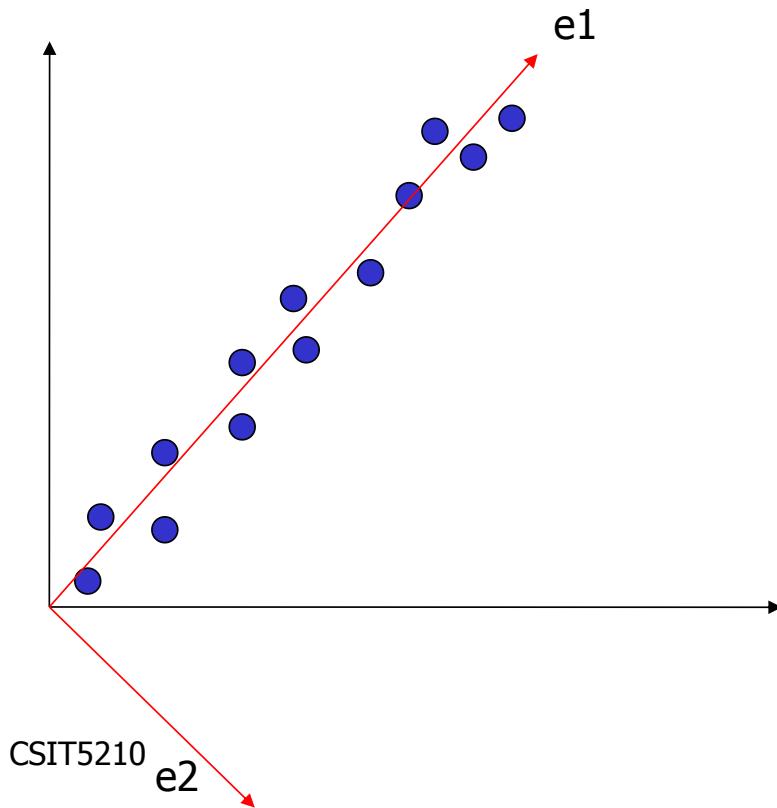


- 
- Why do we need to do this KL-transform?
  - Why do we choose the eigenvectors with the smallest eigenvalues?

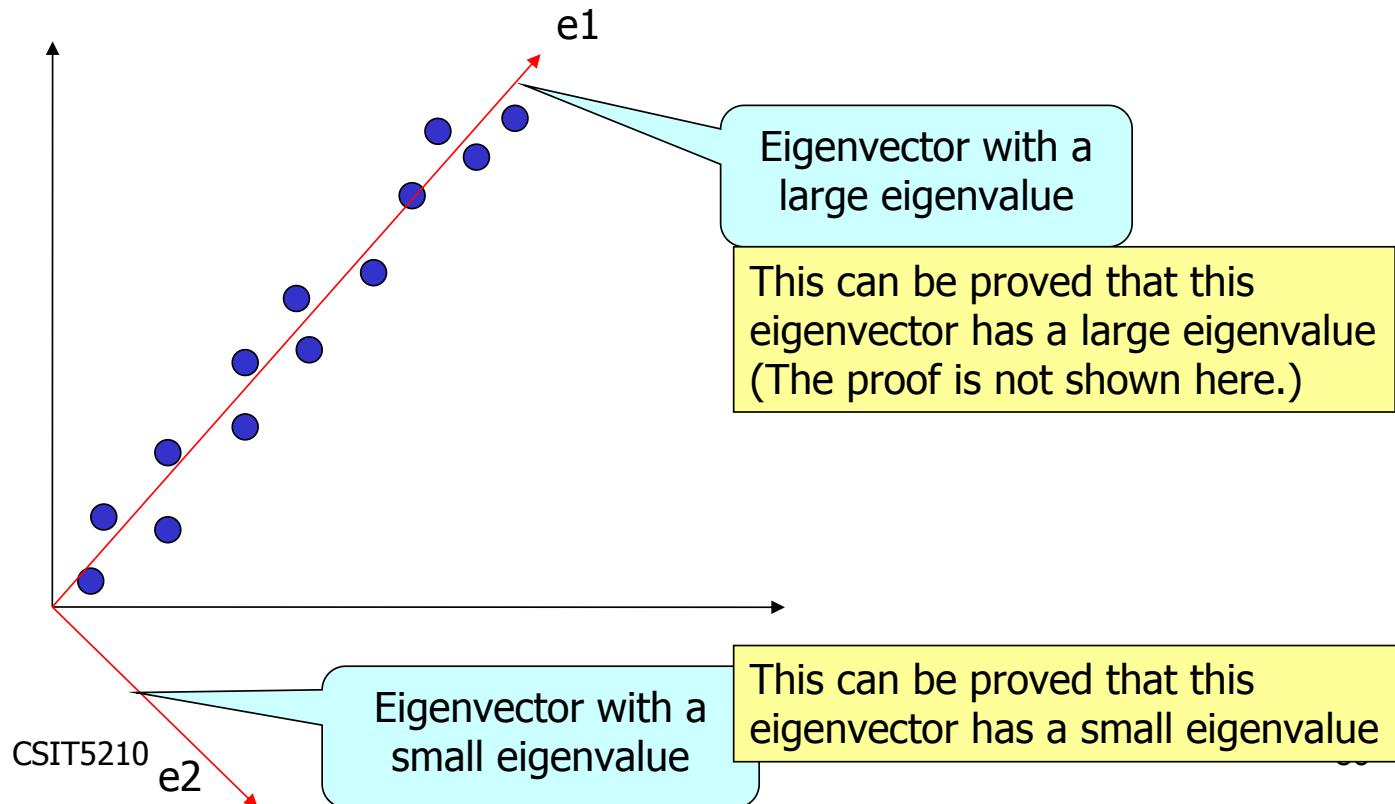
Suppose we have the  
following data set

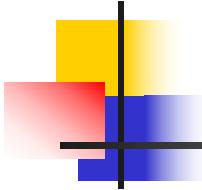


According to the data, we find the following eigenvectors (marked in red)



According to the data, we find the following eigenvectors (marked in red)





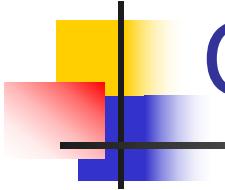
# Consider two cases

---

- Case 1
  - Consider that the data points are projected on  $e_1$
- Case 2
  - Consider that the data points are projected on  $e_2$

It corresponds to dimension reduction.

It corresponds to subspace clustering.

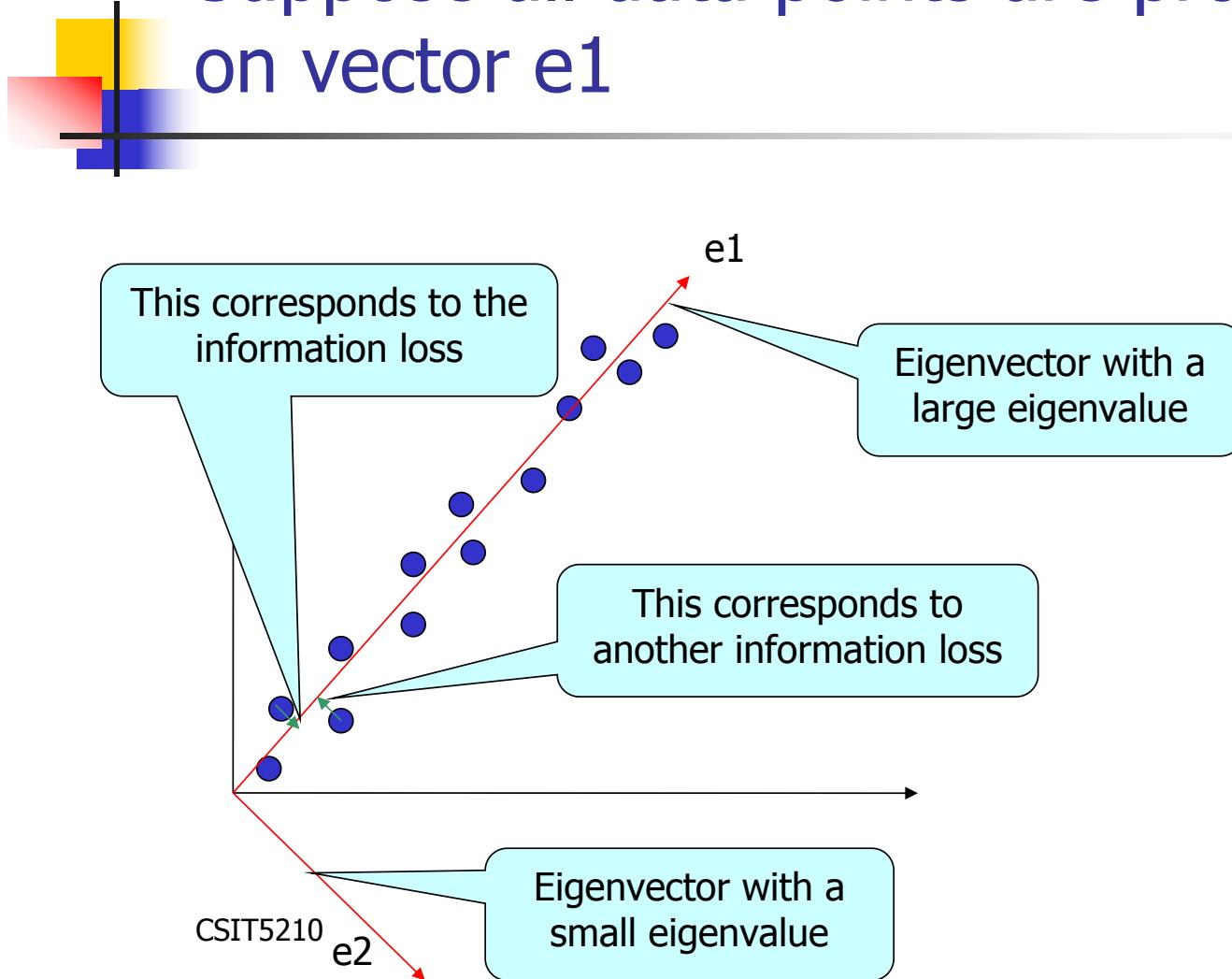


## Case 1

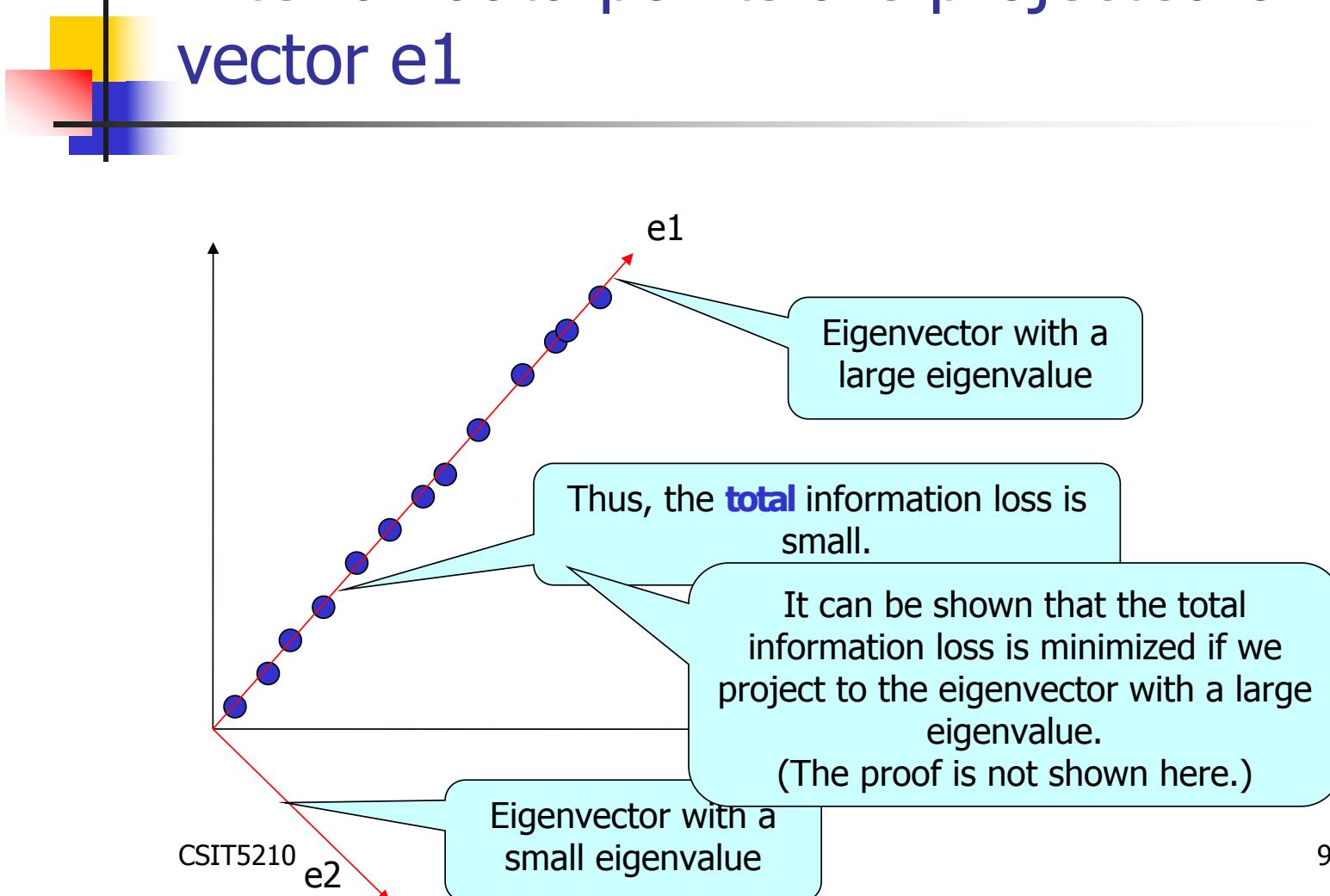
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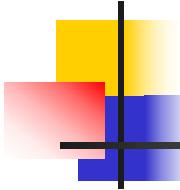
- Consider that the data points are projected on  $e_1$

# Suppose all data points are projected on vector $e_1$



## After all data points are projected on vector $e_1$

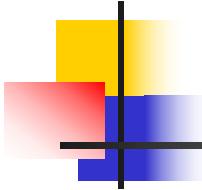




# Objective of Dimension Reduction

---

- The objective of Dimension Reduction
  - To **reduce** the total number of dimensions
  - At the same time, we want to keep information as much as possible.  
(i.e., **minimize** the information loss)
- In our example,
  - We reduce from two dimensions to one dimension
  - The eigenvector with a large eigenvector corresponds to this dimension
  - After we adopt this dimension, we can minimize the information loss

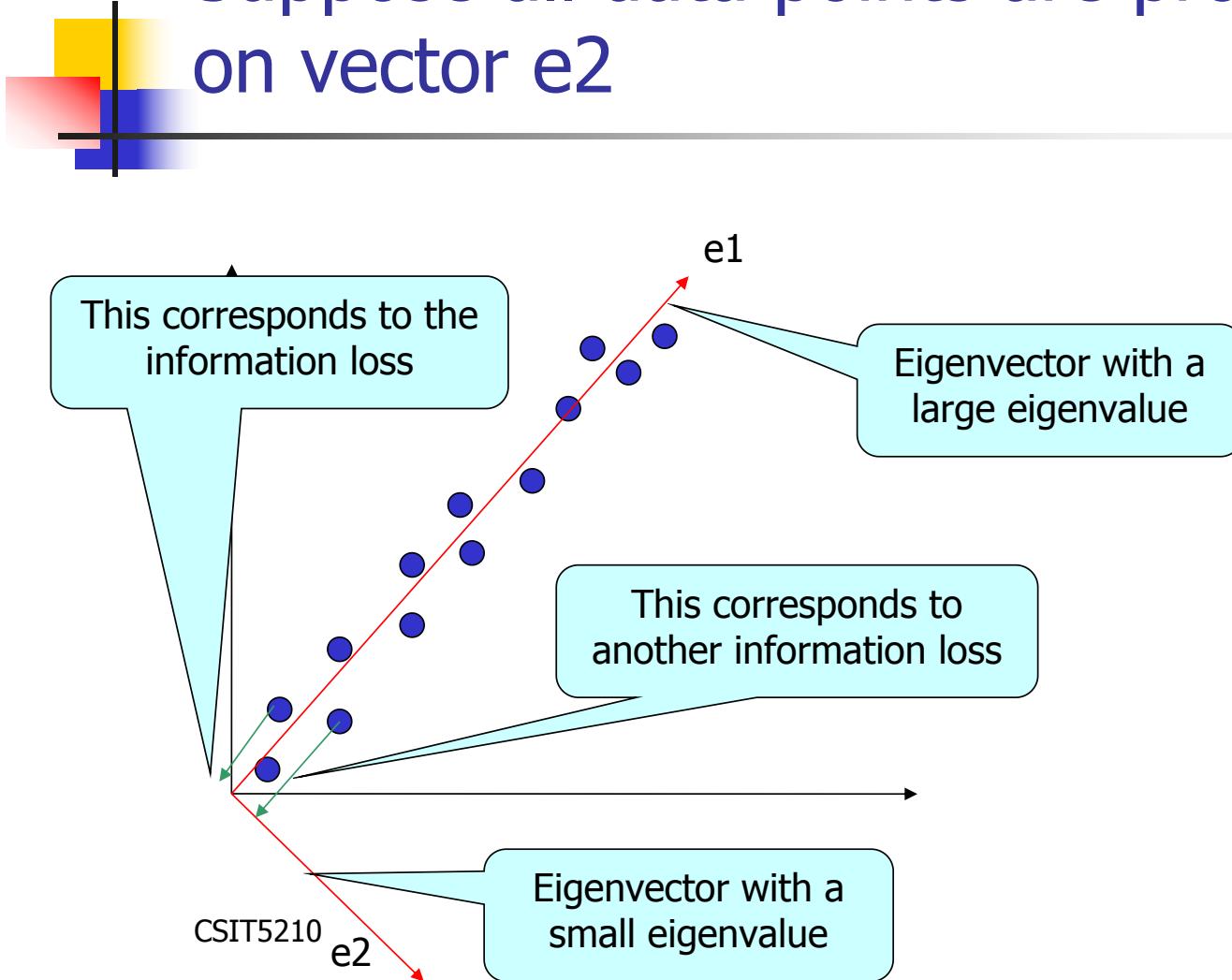


## Case 2

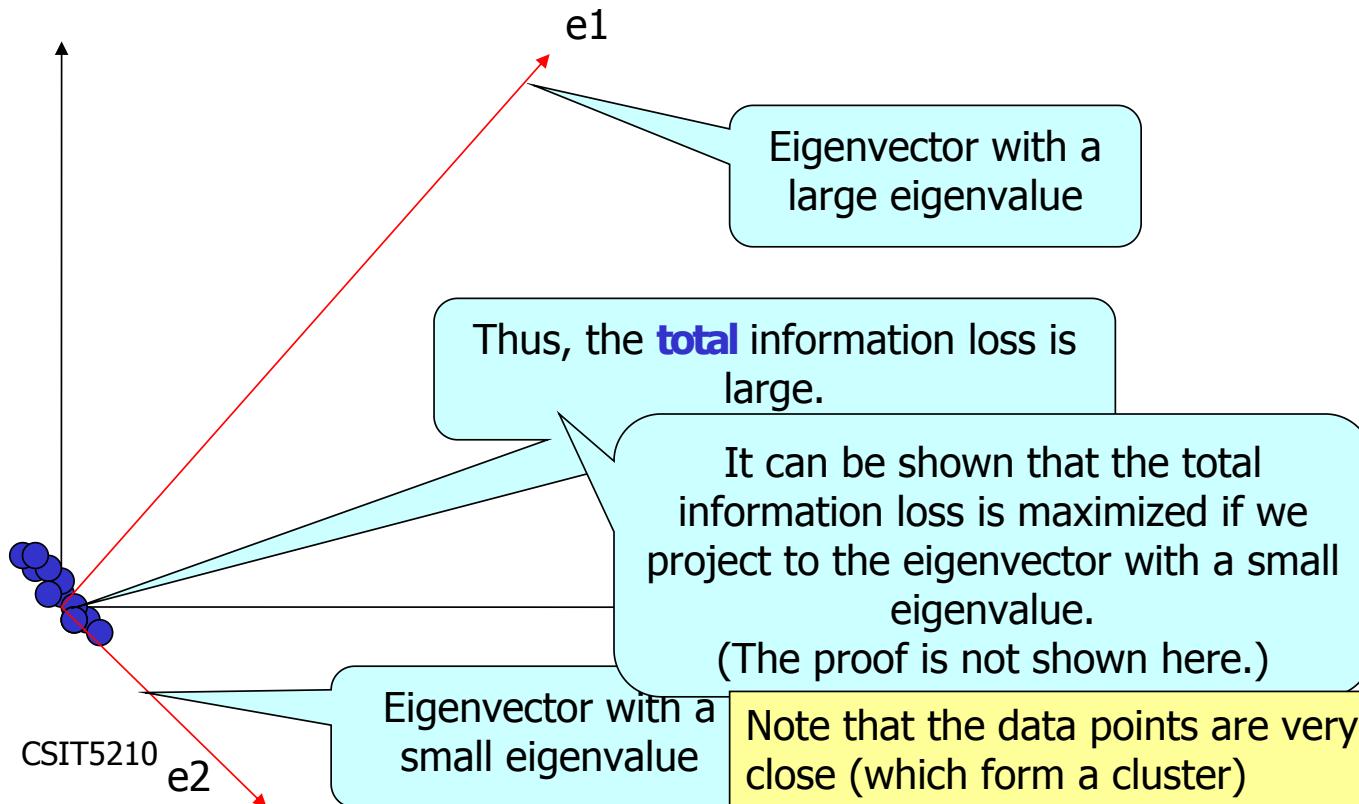
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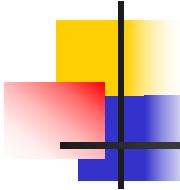
- Consider that the data points are projected on  $e_2$

# Suppose all data points are projected on vector $e_2$



## After all data points are projected on vector $e_2$





# Objective of Subspace Clustering (KL-Transform)

- The objective of Subspace clustering
  - To **reduce** the total number of dimensions
  - At the same time, we want to find a cluster
    - A cluster is a group of “close” data points
    - This means that, after the data points are transformed, the data points are very close.
    - In KL-transform, you can see that the information loss is **maximized**.
- In our example,
  - We reduce from two dimensions to one dimension
  - The eigenvector with a small eigenvalue corresponds to this dimension
  - After we adopt this dimension, we can maximize the information loss