

## Quick RECAP OF LAST CLASS

In the last class, our discussion was as follows:

Supposing we have  $N$  realizations of random variable vectors

$x \in \mathbb{R}^n$  (random variable, is a collection of all possible images

of different kinds of dogs in Michigan state), the mean or

expected value based on the  $N$  realizations, or assuming only  $N$  realizations exist

$$\mu = \sum_{i=1}^N P(X = \vec{x}_i) x_i = \frac{1}{N} \sum_{i=1}^N x_i$$

The covariance matrix expresses the expectation of how any two components (or pixels) are related, i.e.

$$\Sigma = \frac{1}{N} \sum_{i=1}^N (x_i - \mu)(x_i - \mu)^T$$

We showed that the eigen vectors of  $\Sigma$  are orthonormal to one another because it is PSD (and Hermitian); & the eigenvalues are real & non-negative.

We noted that these eigenvectors can serve a new coordinate system or basis for our random variable space; and more importantly the e.vectors are ordered, in terms of relative importance of how much they express distinction or variance in the data.

We order the eigenvectors according to corresponding e. values

Note unbiased Covariance uses  $\frac{1}{n-1}$

$$T = [v_1, v_2, \dots, v_{n-1}]$$

Principal components

$$y = T^H x \quad \leftarrow \text{PCA transformation}$$

which is the projection of the data  
onto the eigen vector space

Further  $\Sigma = T \Lambda T^H$

$\Lambda$  is a Diagonal Matrix of eigenvalues

## Variance of Transformed Data

Notice that

$$\Sigma = \frac{1}{N} \sum_{i=1}^N (\vec{x}_i - \mu)(\vec{x}_i - \mu)^T$$

is equivalent to

$$\Sigma = \frac{1}{n} (X - \vec{\mu} \vec{1}^T) (X - \vec{\mu} \vec{1}^T)^T$$

where  $X = \begin{bmatrix} \vec{x}_1 & \vec{x}_2 & \dots & \vec{x}_n \\ 1 & 1 & & 1 \end{bmatrix}$        $\vec{1} = \underbrace{[1, 1, \dots, 1]^T}_n$

$$\Sigma = \frac{1}{n} X_m X_m^T$$

where  $X_m$  is the mean-subtracted data.

① In the PCA-transformed space, the data has zero covariance.

$$Y_m = T_n^H X_m$$

$$\begin{aligned} \Sigma_y &= \frac{1}{n} Y_m Y_m^T = \frac{1}{n} T_n^H X_m X_m^H T_n \\ &= T_n^H \Sigma_x T_n \end{aligned}$$

But recall that  $\Sigma_x = T_n \Lambda T_n^H$

$$\therefore \Sigma_y = T_n^H T_n \Lambda T_n^H T_n = \Lambda : \text{Covariance is diagonal}$$

so just Variance

## Singular Value Decomposition

Every complex-valued matrix  $A \in \mathbb{C}^{n \times m}$  can be decomposed as

$$A = U S V^H$$

where  $U \in \mathbb{C}^{n \times n}$  and  $V \in \mathbb{C}^{m \times m}$  are unitary matrices with orthonormal columns, and  $S \in \mathbb{R}^{n \times m}$  is a matrix with real non-negative entries on the diagonals and zeros off the diagonal.

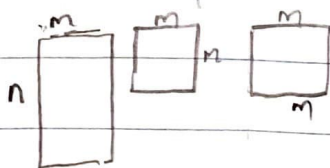
When  $n \geq m$ , the matrix  $S$  has at most  $m$  non-zero elements on the diagonal and may be written as

$$S = \begin{bmatrix} \hat{S} \\ 0 \end{bmatrix}$$

Therefore, we can exactly represent the matrix  $A$  using 'economy' SVD:

$$A = U S V^H = \begin{bmatrix} \hat{U} & \hat{U}^\perp \end{bmatrix} \begin{bmatrix} \hat{S} \\ 0 \end{bmatrix} V^H = \hat{U} \hat{S} V^H$$

where the columns of  $\hat{U}^\perp$  span a vector space that is complementary and orthogonal to that spanned by  $\hat{U}$ .





The columns of  $U$  are called left singular vectors; those of  $V$  are right singular vectors. The diagonal elements of  $\hat{S}$  are called singular values and are ordered from largest to smallest. The rank of matrix  $A$  is equal to the number of non-zero singular values.

We can write the SVD for the mean subtracted data as

$$X_m = U S V^H$$

Recall that the covariance matrix is

$$\Sigma = \frac{1}{n} X_m X_m^H = \frac{1}{n} U S^2 U^H = T \Lambda T^H$$

Hence, the principal components of  $T$  are also the left singular vectors of  $X_m$ . The variance of the data in the new coordinates

$$\lambda_k = \frac{\sigma^2}{n}$$