Big Data Fundamentals and Applications

Principal Component Principal Component<

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Outlines

- 1. Review
- 2. Why Do We Need Dimension Reduction?
- 3. PCA Mathematical Explanation
- 4. PCA Programming in Python
- 5. Part VIII Basic numerical methods (Scipy) Part IX Advanced numerical methods (Scipy)
- 6. Notice



 Before we explain PCA, we need to review the mathematical meaning of three basic descriptive statistics, including expectation, variance, and covariance.

$$\mu = \frac{1}{n} \sum_{i=1}^{n} x_i$$

$$var(x) = \sigma^2 = \left(\sqrt{\frac{\sum_{i=1}^{n} (x_i - \mu)^2}{n}}\right)^2 = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu)^2$$

$$Cov(x, y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_x) (y_i - \mu_y)$$

 In previous courses or your understanding, these three parameters usually perform as above equations.

$$\mu = \frac{1}{n} \sum_{i=1}^{n} x_i \mid var(x) = \sigma^2 = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu)^2 \mid Cov(x, y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_x) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} ($$





$$\mu = \frac{1}{n} \sum_{i=1}^{n} x_i \mid var(x) = \sigma^2 = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu)^2 \mid Cov(x, y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_x) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y)$$



$$\mu = \frac{1}{n} \sum_{i=1}^{n} x_i \mid var(x) = \sigma^2 = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu)^2 \mid Cov(x, y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_x) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_y) (y_i - \mu_y) (y_i - \mu_y)$$

- Pearson's correlation coefficient
- Given two parameters x_i and y_i , ⁰³ where *i* ranges from 1 to *n*. ⁰⁴ Then, Pearson's correlation ⁰⁵ coefficient could be defined as follows. ¹⁰

$$\rho = \frac{\sum_{i=1}^{n} (x_i - \mu_x) (y_i - \mu_y)}{\sqrt{\sum_{i=1}^{n} (x_i - \mu_x)^2 \sum_{i=1}^{n} (y_i - \mu_y)^2}}$$



Question 1

If x is highly correlated with y, and then what do you expect from their covariance and standard deviations?

Expectation

$$E(X) = \sum_{x} xP(X = x) = \mu$$

Variance

$$var(X) = E([X - \mu]^2)$$

= $E(X^2 - 2\mu X + \mu^2)$
= $E(X^2) - 2\mu E(X) + \mu^2$
= $E(X^2) - 2\mu^2 + \mu^2$
= $E(X^2) - \mu^2$
= $E(X^2) - \mu^2$

Characteristics of Expectation

E(aX + bY) = aE(X) + bE(Y), a, b $\in \mathbb{R}$ X and Y can be independent or dependent.

E(XY) = E(X)E(Y)Where cov(X,Y) = 0

- Covariance
- If x and y are independent... var(X + Y) = var(X) + var(Y)
- If x and y are dependent... $var(X + Y) = E([(X + Y) - E(X + Y)]^2)$ $= E\left(\left[(X+Y) - \left(E(X) + E(Y)\right)\right]^2\right)$ $= E\left(\left[\left(X - E(X)\right) + \left(Y - E(Y)\right)\right]^2\right)$ $= E\left((X - E(X))^{2} + 2(X - E(X))(Y - E(Y)) + (Y - E(Y))^{2} \right)$ $= E\left[(X - E(X))^{2} \right] + E\left[(Y - E(Y))^{2} \right] + 2E\left[(X - E(X))(Y - E(Y)) \right]$ = var(X) + var(Y) + 2cov(X,Y)

 $E(aX + bY) = aE(X) + bE(Y), a, b \in \mathbb{R}$ X and Y can be independent or dependent. E(XY) = E(X)E(Y)Where cov(X,Y) = 0

Why Do We Need Dimension Reduction?

- Here comes the first question into your mind.
 - Why do we need dimension reduction?
 - What's the importance of dimension reduction?
 - Can we directly import all datasets into your model without dimension reduction?
- Statistical models (e.g., linear regression) have several assumptions when you adopt them. One of them is "all variables have to be linearly independent," indicating no collinearity.
- To achieve this goal, various methods were developed for orthogonalizing parameters and reducing the dimension of the dataset, such as PCA, LDA, LLE, and Laplacian Eigenmaps.

PCA – Math $\mu = \frac{1}{n} \sum_{i=1}^{n} x_i | var(x) = \sigma^2 = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu)^2 | Cov(x, y) = \frac{1}{n} \sum_{i=1}^{n} (x_i - \mu_x) (y_i - \mu_y)$





• Given a point "x" and project onto a vector "v".

$$\|x_i\|\cos\theta = \|x_i\| \frac{x_i^T \cdot \nu}{\|x_i\| \|\nu\|} = \frac{x_i^T \cdot \nu}{\|\nu\|}$$

$$if \ v \ is \ unit \ vector \ \dots \ \|\nu\| = 1$$

$$= \frac{x_i^T \cdot \nu}{\|\nu\|} = x_i^T \cdot \nu$$

 $cos\theta = \frac{x_i^T \cdot v}{\|x_i\| \|v\|}$

$$X = \begin{bmatrix} | & | & \cdots & | \\ x_1 & x_2 & \cdots & x_n \\ | & | & \cdots & | \end{bmatrix} \to X^T = \begin{bmatrix} - & x_1 & - \\ - & x_2 & - \\ - & \vdots & - \\ - & x_n & - \end{bmatrix}$$

$$P = \begin{bmatrix} x_1^T \\ x_2^T \\ \vdots \\ x_n^T \end{bmatrix} = X^T u \implies solve P$$

$$x_{i}$$

$$x_{i}$$

$$x_{i}$$

$$x_{i}$$

$$x_{i}$$

$$x_{i}^{T} \cdot v$$

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$$J(u) = ||P^{2}|| = P^{T}P = (X^{T}u)^{T}(X^{T}u) = u^{T}XX^{T}u$$

$$\underset{u}{\operatorname{argmax}} J(u) = u^T X X^T u, where \ subject \ to \ u^T u = 1$$

when u is the eigen vector $J(u) = ||P^2|| = u^T X X^T u = u^T \lambda u = \lambda u^T u = \lambda$

Given an eigenvector, the total square of projected values is the eigenvalue = λ

 $XX^T u = \lambda u$

Eigenvector is a symmetry matrix u is an unit vector $uu^T = u^T u = 1$



A is a square symmetric matrix has orthogonal eigenvectors with different eigenvalues.

 $[x_{1}, \lambda_{1}], [x_{2}, \lambda_{2}]$ $\begin{cases}
Ax_{1} = \lambda_{1}x_{1} \\
Ax_{2} = \lambda_{2}x_{2} \\
x_{1}^{T}Ax_{2} = x_{1}^{T}\lambda_{2}x_{2} = \lambda_{2}x_{1}^{T}x_{2} \\
x_{1}^{T}A^{T}x_{2} = (Ax_{1})^{T}x_{2} = (\lambda_{1}x_{1})^{T}x_{2} = \lambda_{1}x_{1}^{T}x_{2} \\
(\because A \in symmetric matrix, \therefore A = A^{T})
\end{cases}$

Eigenvector is a symmetry matrix u is an unit vector

 $XX^T u = \lambda u$

 $uu^T = u^T u = 1$

$$\lambda_2 x_1^T x_2 = \lambda_1 x_1^T x_2$$
$$\underline{x_1^T x_2} (\lambda_2 - \lambda_1) = 0$$

Orthogonal All eigenvalues $x_1^T x_2 = 0$ are different

 Conversion between orthogonal bases $u_{i} \cdot u_{j} = u_{i}^{T} \cdot u_{j} = \begin{cases} 1, if \ i = j \\ 0, otherwise \end{cases}$ $U = \begin{vmatrix} | & | & | \\ u_1 & u_2 & \dots & u_d \\ | & | & | \end{vmatrix} \Rightarrow U^T U = I = U^{-1} = U^T$ $x = y_1 u_1 + y_2 u_2 + \dots + y_d u_d = \begin{bmatrix} | & | & | & | \\ u_1 & u_2 & \dots & u_d \\ | & | & | & | \end{bmatrix} \begin{vmatrix} y_1 \\ y_2 \\ \vdots \\ \vdots \\ y_d \end{vmatrix} = Uy$ $\Rightarrow y = U^{-1}x = U^Tx$

 $XX^T v = \lambda v \Rightarrow AV = \lambda V$ where *V* is eigenvector and λ is eigenvalue. **Principal component** *PC* $AV = \lambda V$

From Singular Vector Decomposition (SVD)

$$A = U\Sigma V^{T} \Rightarrow AV = U\Sigma \Rightarrow \lambda = \frac{\Sigma^{2}}{N}, \therefore AV = U\Sigma$$

Variable loading L (A onto PC)
Standardize PC = $\sqrt{N}U$
 $cov(A, PC) = \frac{A^{T}\sqrt{N}U}{N} = \frac{V\Sigma U^{T}U}{\sqrt{N}} = V\frac{\Sigma}{\sqrt{N}} = V\sqrt{E} = L$

PCA – Math

 $J(u) = ||P^{2}|| = P^{T}P = (X^{T}u)^{T}(X^{T}u) = u^{T}XX^{T}u$

argmax $J(u) = u^T X X^T u$, where subject to $u^T u = 1$ Add Lagrange multiplier argmax $J(u, \lambda) = u^T X X^T u - \lambda (1 - u^T u)$ $\stackrel{u,\lambda}{\nabla_u J(u, \lambda)} = \nabla_u (u^T X X^T u - \lambda (1 - u^T u)) = 0$ $\Rightarrow 2X X^T u - 2\lambda u = 0$ $\Rightarrow X X^T u = \lambda u \rightarrow \text{eigenvector} \qquad Au = \lambda u$

The steps of PCA

- 1. Find the sample mean $\mu = \frac{1}{n} \sum_{i=1}^{n} x_i$
- 2. Subtract mean
- 3. Compute covariance matrix $C = \frac{1}{n}XX^T = \frac{1}{n}\sum_{i=1}^n (x_i \mu)(x_i \mu)$
- 4. Find the eigenvalues of C and arrange them into descending order

$$\lambda_1 > \lambda_2 > \dots > \lambda_d, \{u_1, u_2, \dots, u_d\}$$

5. The transformation is $y = U^T X$.

PCA – Python

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learn

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scikit-learn 1.1.2 Other versions

Please **cite us** if you use the software.

sklearn.decomposition.PCA
Examples using
sklearn.decomposition.PCA

sklearn.decomposition.PCA

class sklearn.decomposition.PCA(n_components=None, *, copy=True, whiten=False, svd_solver='auto', tol=0.0, iterated_power='auto', n_oversamples=10, power_iteration_normalizer='auto', random_state=None)

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Principal component analysis (PCA).

Linear dimensionality reduction using Singular Value Decomposition of the data to project it to a lower dimensional space. The input data is centered but not scaled for each feature before applying the SVD.

It uses the LAPACK implementation of the full SVD or a randomized truncated SVD by the method of Halko et al. 2009, depending on the shape of the input data and the number of components to extract.

It can also use the scipy.sparse.linalg ARPACK implementation of the truncated SVD.

Notice that this class does not support sparse input. See TruncatedSVD for an alternative with sparse data.

Read more in the User Guide.

Parameters:: n_components : int, float or 'mle', default=None

Number of components to keep. if n_components is not set all components are kept:

n_components == min(n_samples, n_features)

If n_components == 'mle' and svd_solver == 'full', Minka's MLE is used to guess the dimension. Use of n_components == 'mle' will interpret svd_solver == 'auto' as svd_solver == 'full'.

If $0 < n_{components} < 1$ and $svd_{solver} == 'full'$, select the number of components such that the amount of variance that needs to be explained is greater than the percentage specified by n_components.

If svd_solver == 'arpack', the number of components must be strictly less than the minimum of n_features and n_samples.

PCA – Python

linear algebra from sklearn.decomposition import PCA # call PCA pca = PCA(n_components=2) # number of preserved components pca.fit(X) # show results print(pca.mean_) # mean print(pca.explained_variance_) # eigenvalues print(pca.components_) # eigenvectors

Notice

- Advantages of PCA
 - Easy to use
 - Accelerate model fitting
 - Avoid overfitting
- Disadvantage of PCA
 - Low interpretability
 - The trade-off between information loss and features (dimensions)

- Limitations of PCA
 - Linear relationship between features
 - Correlation between features
 - Sensitive to the scale of data
 - Not robust to outliers
 - Not accept missing values

Question Time

If you have any questions, please do not hesitate to ask me.

References:

張智星 科學計算講義 Sklearn



Big Data Fundamentals and Applications Principal Component Analysis – PCA

The End Thank you for your attention))

