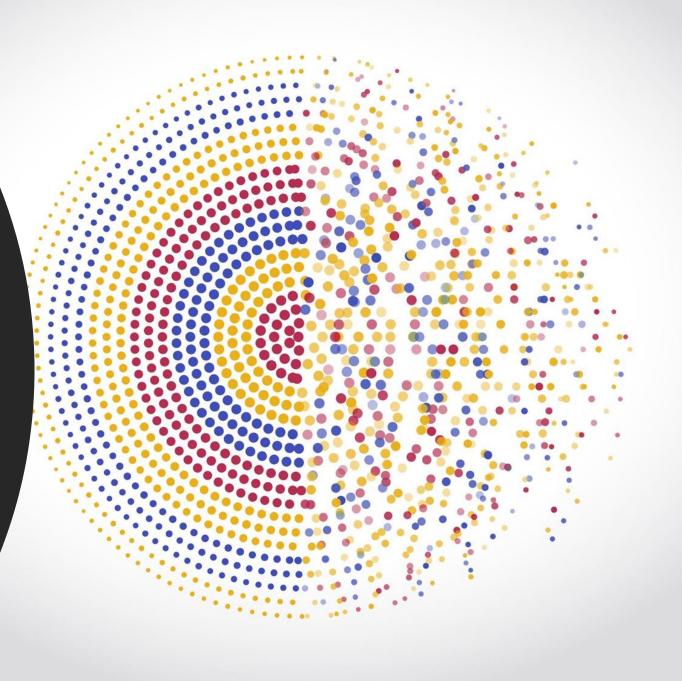
Data Preprocessing

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Data Preprocessing

- Dimensionality Reduction
- Feature Selection
- Feature Creation

Dimensionality Reduction

- Dimensionality reduction is a technique that transforms high-dimensional data into a lower-dimensional space:
 - · Principal Component Analysis (PCA)
 - Linear Discriminant Analysis
 - · t-Distributed Stochastic Neighbor Embedding (t-SNE)
 - · UMAP (Uniform Manifold Approximation and Projection)
 - · Random Projection

• PCA is a dimensionality reduction technique that transforms high-dimensional data into a lower-dimensional space while capturing the directions of maximum variance. This is widely used for tasks such as data compression, noise reduction, and visualization.

- Before performing PCA, we center the data by subtracting the mean of each feature. This ensures that the principal components align with the directions of maximum variance, and that the covariance matrix accurately represents the relationships between features.
- For a dataset $X \in R^{n \times d}$, where n is the number of samples and d is the number of features, we compute the mean vector $\mu \in R^{1 \times d}$ for each feature. Subtracting the mean results in the centered data matrix $X_{centered} \in R^{n \times d}$:

$$X_{centered} = X - \mu$$

• The covariance matrix $\Sigma \in \mathbb{R}^{d \times d}$ is then computed as:

$$\Sigma = \frac{1}{n-1} X_{centered}^T X_{centered}$$

Each element Σ_{ij} represents the covariance between features i and j, with the diagonal elements representing the variance of individual features.

- Eigenvalue Decomposition
- PCA finds the principal components by solving the eigenvalue decomposition of the covariance matrix:

$$\Sigma v_i = \lambda_i v_i$$

Where:

- $v_i \in \mathbb{R}^{d \times 1}$ are the eigenvectors (principal components).
- $-\lambda_i \in R$ are the eigenvalues, representing the variance captured by each component.
- The top k eigenvectors form the matrix $W_k \in \mathbb{R}^{d \times k}$, which projects the original data into a lower-dimensional subspace:

$$Y = X_{centered}W_k$$
, where $Y \in \mathbb{R}^{n \times k}$.

- Singular Value Decomposition (SVD)
- Alternatively, PCA can be performed using Singular Value Decomposition (SVD). SVD decomposes the original data matrix $X \in \mathbb{R}^{n \times d}$ as:

$$X = U \Sigma V^T$$

Where:

- $U \in \mathbb{R}^{\{n \times n\}}$ contains the left singular vectors.
- $V \in \mathbb{R}^{\{d \times d\}}$ contains the right singular vectors (the principal components).
- $-\Sigma \in \mathbb{R}^{\{n \times d\}}$ contains the singular values.

- Connection Between Maximizing Variance and Minimizing Error
- The objective of PCA is to find a projection matrix $W_k \in R^{\{d \times k\}}$ that maximizes the variance of the projected data. The variance of the projected data $Y = X_{\{centered\}} W_k$ can be expressed as:

$$Var(Y) = \frac{1}{n-1} Y^T Y = \frac{1}{n-1} W_k^T X_{\{centered\}}^T X_{\{centered\}} W_k = W_k^T \Sigma W_k$$

To maximize the variance, we solve the optimization problem:

 $max_W tr(W^T \Sigma W)$, subject to $W^T W = I$.

- Connection Between Maximizing Variance and Minimizing Error
- Minimizing the reconstruction error is an alternative interpretation of PCA. After projecting the data onto the lower-dimensional subspace, we attempt to reconstruct the original data, minimizing the Frobenius norm of the difference between the original data and its reconstruction:

$$min_W ||X_{\{centered\}} - X_{\{centered\}} W W^T||_F^2$$
.

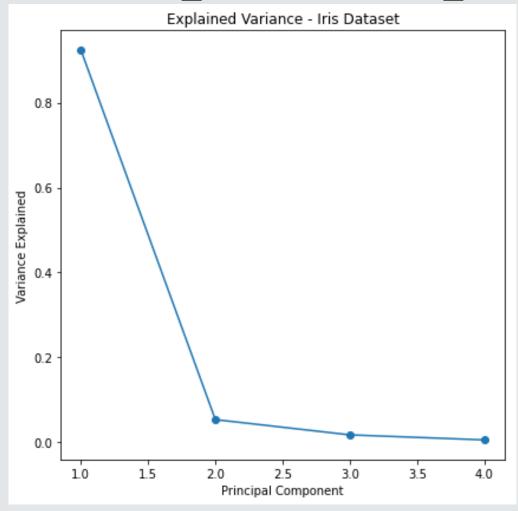
The expanded Frobenius norm is:

$$||X_{\{centered\}} - X_{\{centered\}} W W^T||_F^2$$

$$= tr(X_{\{centered\}}^T X_{\{centered\}}) - 2tr(W^T X_{\{centered\}}^T X_{\{centered\}} W)$$

$$+ tr(W^T X_{\{centered\}}^T X_{\{centered\}} W W^T W).$$

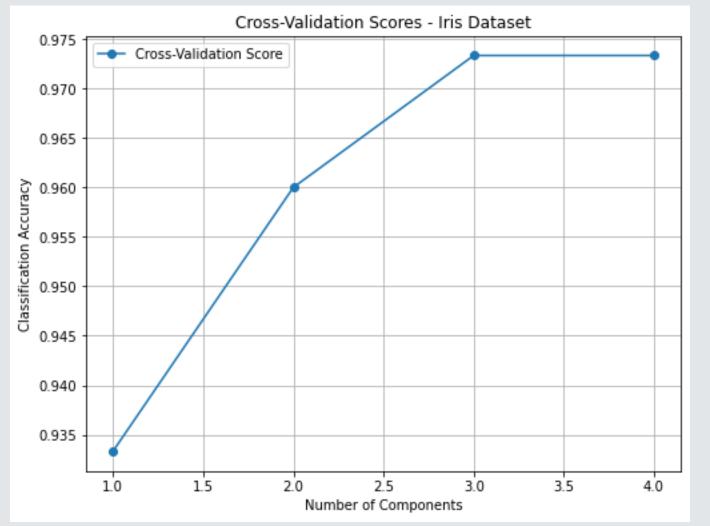
- Choosing the Number of Principal Components
- The following methods help to determine the optimal number of principal components:
 - 1. Elbow Method.
 - 2. Cumulative Explained Variance.
 - 3. Cross-Validation.

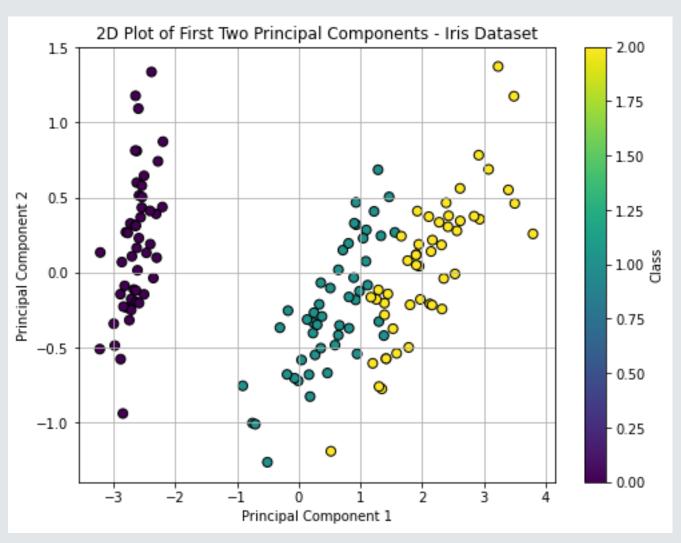


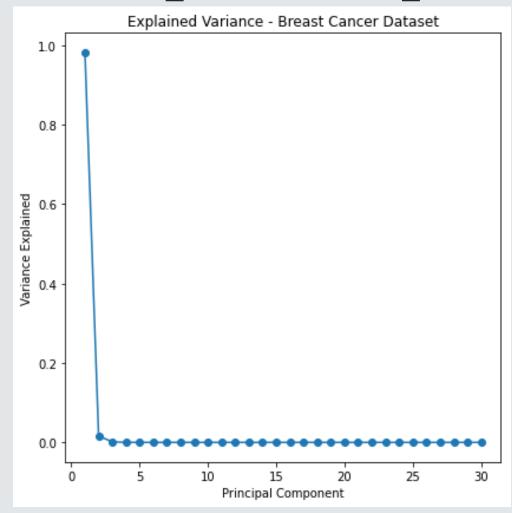
Cumulative Variance - Iris Dataset Cumulative Variance --- 95% Variance Threshold 0.99 0.98 Cumulative Variance 0.97 0.96 0.94 0.93 1.5 2.5 3.5 2.0 3.0 4.0 1.0 Principal Component

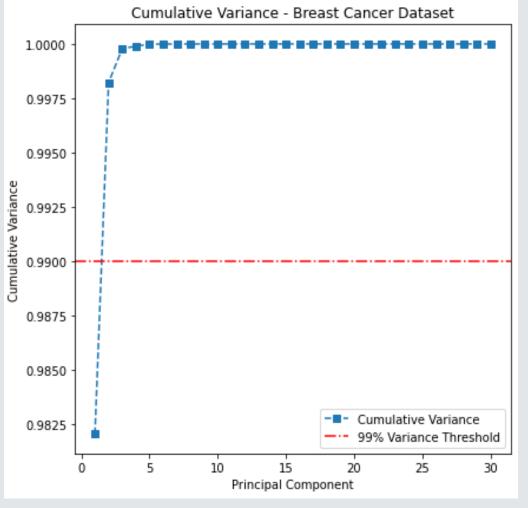
Elbow Method

Cumulative Explained Variance.



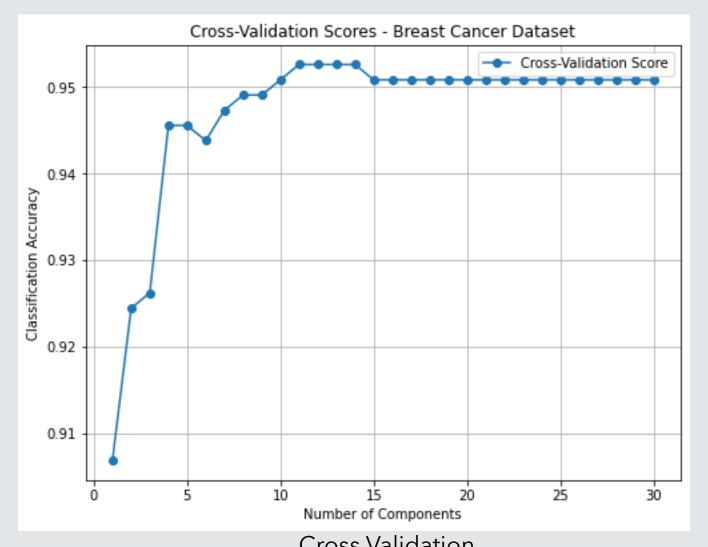


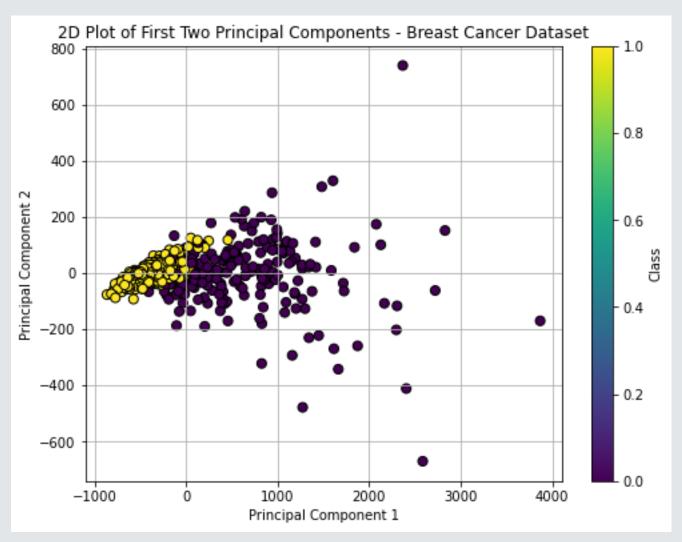




Elbow Method

Cumulative Explained Variance.





Limitations

- Linearity: PCA assumes that the principal components are linear combinations of the original features. It fails to capture non-linear relationships in the data.
- Interpretability: The principal components are often linear combinations of many features, making it difficult to interpret their meaning.
- Variance-Based: PCA focuses on maximizing variance, which may not always capture the most important features for specific tasks such as classification.
- Sensitivity to Scaling: PCA is sensitive to the scaling of features. Features with higher magnitude may dominate the principal components unless the data is appropriately scaled.

• Feature selection is the process of selecting the most relevant features from a dataset to improve the performance of a machine learning model. It helps by reducing the dimensionality of the feature space, improving model accuracy, and decreasing computational costs. There are three main types of feature selection methods: **filter methods**, wrapper methods, and embedded methods.

- Filter Methods
- Filter methods are statistical techniques that rank and select features based on their properties, independent of any specific machine learning model. They are fast and efficient, making them suitable for pre-processing steps.
 - · Variance Threshold
 - · Correlation Coefficient
 - · Mutual Information

Filter Method - Variance Threshold

- Variance threshold removes features that have a variance below a specified threshold. Features with low variance typically carry little information.
- **Mathematics**: For a feature $X_i \in \mathbb{R}^n$, the variance is:

$$Var(X_j) = \frac{1}{n} \sum (X_{ij} - \mu_j)^2$$

Features with variance below the threshold are removed.

Filter Method - Correlation Coefficient

- The Pearson correlation coefficient measures the linear relationship between a feature and the target variable. Features with low correlation to the target are considered irrelevant.
- Mathematics: The Pearson correlation coefficient between a feature X_j and target y is:

$$\rho(X_j, y) = \frac{Cov(X_j, y)}{\sigma_X \sigma_y}$$

is the covariance between X_j and y, and σ_{X_j} and σ_y are the standard deviations.

Filter Method - Mutual Information

- Mutual information (MI) measures the dependency between a feature and the target variable. It captures both linear and non-linear relationships.
- Mathematics: The mutual information between a feature X_i and target y is:

$$I(X_j; y) = \sum_{x \in X_j, \alpha \in y} p(x, \alpha) \log(\frac{p(x, \alpha)}{p(x)p(\alpha)})$$

Higher MI indicates a stronger relationship.

- Wrapper Methods
- Wrapper methods evaluate subsets of features using a machine learning model to determine which features contribute the most to model performance. These methods are computationally expensive but often more accurate than filter methods.
 - · Recursive Feature Elimination
 - · Forward Feature Selection

Wrapper Method - Recursive Feature Elimination

 Recursive Feature Elimination (RFE) ranks features based on their importance and recursively eliminates the least important ones.

Process:

- Train a model with all features.
- Rank the features based on their importance (e.g., model coefficients).
- Remove the least important feature.
- Repeat until the desired number of features remains.

Wrapper Method - Forward Feature Selection

• Forward selection starts with no features and adds features one by one, selecting the feature that improves model performance the most at each step.

Embedded Methods

- Embedded methods perform feature selection during the model training process. Lasso regression is one of the most common embedded methods, as it automatically selects features by penalizing the coefficients of less important features.
- Lasso Regression (L1 Regularization)
- Lasso regression adds an L_1 regularization term to the objective function, which forces some feature coefficients to become zero, effectively selecting a subset of features.
- Mathematics: Lasso regression solves the following optimization problem:
- $\min_{w} \frac{1}{2n} || Xw y ||_{2}^{2} + \lambda || w ||_{1}$
- where λ controls the strength of regularization. Features corresponding to zero coefficients are removed.