Independent Component Analysis(ICA)

Independent Component Analysis (ICA): A Deep Dive

Problem Context: The Cocktail Party Problem

The cocktail party problem is a classic example of source separation. Imagine you're in a room where multiple people are speaking simultaneously, and you have several microphones recording the mixed audio signals. The goal is to separate the voices of the individual speakers from the recorded mixtures.

The Mathematical Framework

1. Observed Data (Mixed Signals)

Let the observed signals be represented as: $X = A \cdot S$

- X: A matrix of observed signals (mixed signals), where each column corresponds to a signal recorded by a microphone.
- ullet A: A mixing matrix, which combines the independent source signals.
- S: A matrix of independent source signals (e.g., voices of different speakers).

The goal of ICA is to recover S (the independent source signals) using only X.

2. Assumptions of ICA

ICA relies on three key assumptions:

1. Linearity:

The observed signals are linear mixtures of the source signals.

2. Statistical Independence:

The source signals S are statistically independent.

3. Non-Gaussianity:

 The source signals are non-Gaussian. This is crucial for separation since Gaussian signals cannot be separated based solely on statistical independence.

3. ICA Objective

ICA seeks a demixing matrix W such that: $S_{\text{estimated}} = W \cdot X$ where $S_{\text{estimated}}$ approximates the original independent sources S.

Key Steps in ICA

Step 1: Centering

The data is centered to ensure zero mean for each feature: $X_{\rm centered} = X - \mu$ This simplifies subsequent calculations by aligning the data with the origin.

Step 2: Whitening

Whitening removes correlations between features and scales the data to unit variance:

$$X_{ ext{whitened}} = W_{ ext{whitening}} \cdot X_{ ext{centered}}$$

ullet Whitening ensures that the covariance matrix of $X_{
m whitened}$ is the identity matrix:

$$Cov(X_{whitened}) = I$$

Step 3: Iterative Optimization

The weight matrix W is iteratively optimized to maximize the statistical independence of the estimated components. This involves:

- 1. Applying a **non-linear function** to amplify non-Gaussianity: $g(x) = \tanh(x)$
- 2. Updating W using gradient-based optimization and normalizing it after each update to maintain stability.

Step 4: Decorrelation

To ensure orthogonality of the weight matrix W, it is decorrelated after each iteration: $W_{\text{decorrelated}} = Q \cdot D \cdot Q^{\top} \cdot W$ where Q and D are derived from the eigenvalue decomposition of WW^{\top} .

Step 5: Projection

Once W is optimized, the independent components are extracted by projecting the whitened data: $S=W\cdot X_{ ext{whitened}}$

How ICA Solves the Cocktail Party Problem

1. Input:

ullet The observed signals X are mixtures of the original independent sources.

2. Processing:

ullet ICA learns a demixing matrix W that inverts the mixing process.

3. Output:

ullet The independent components $S_{
m estimated}$ are the separated source signals (e.g., the voices of individual speakers).

Applications of ICA

1. Audio Signal Processing:

Separating voices or musical instruments from mixed audio recordings.

2. Neuroscience:

• Analyzing EEG or fMRI data to identify independent brain activity patterns.

3. Financial Analysis:

• Identifying independent factors driving market behaviors.

Summary

ICA is a powerful algorithm for blind source separation. By leveraging the principles of statistical independence and non-Gaussianity, ICA can decompose mixed signals into their independent components, solving problems like the cocktail party problem.

Step 1: Initializing Parameters

Parameters in the ICA Class

1. n_components

- Definition: The number of independent components to extract from the data.
- **Purpose**: Determines the size of the weight matrix and the number of independent signals to recover.
- **Typical Use Case**: If the input data has $n_{\rm features}$, n_components is usually less than or equal to $n_{\rm features}$.
- Initialization: $self. n_components = n_components$

2. max_iter

- Definition: The maximum number of iterations the algorithm will run before stopping.
- Purpose: Acts as a safeguard to ensure the algorithm terminates even if convergence isn't achieved.
- **Typical Use Case**: Values like 1000 or higher are common for complex datasets.
- Initialization: $self. max_i ter = max_i ter$

3. tol

- **Definition**: The convergence tolerance, controlling when the algorithm stops iterating.
- **Purpose**: If the change between iterations falls below this threshold, the algorithm assumes convergence and stops.
- Typical Use Case: Smaller values (e.g., 1e-5) enforce stricter convergence but may increase runtime.
- Initialization: self. tol = tol

Summary

In the __init__ method of the ICA class, these parameters are stored as attributes for later use in the algorithm.

Step 2: Centering the Data

Goal

The purpose of centering is to ensure that the dataset has a mean of zero along each feature. This is a critical preprocessing step in ICA and other statistical methods.

Mathematical Explanation

Given a dataset X of shape $(n_{\rm samples}, n_{\rm features})$, the mean of each feature is computed as: $\mu = \frac{1}{n_{\rm samples}} \sum_{i=1}^{n_{\rm samples}} X[i]$ The centered dataset is obtained by subtracting this mean from each feature: $X_{\rm centered} = X - \mu$

Importance of Centering

- 1. **Statistical Independence**: Centering removes the bias introduced by non-zero means, which is crucial for identifying independent components.
- Simplifies Whitening: Whitening transforms the covariance matrix into an identity

matrix. Centering ensures that the covariance matrix is well-defined and unbiased.

3. **Numerical Stability**: Subtracting the mean can improve numerical stability in subsequent steps, like eigenvalue decomposition.

Example Dataset

Consider the dataset:
$$X=\begin{bmatrix}1&2\\3&4\\5&6\end{bmatrix}$$
 After centering, each column should have a mean of zero.

Step 3: Whitening the Data

Goal

Whitening transforms the dataset so that its variables are decorrelated and have unit variance. This simplifies the optimization in ICA by standardizing the data and aligning it to a "spherical" feature space.

Mathematical Foundation

1. Covariance Matrix

The covariance matrix measures the linear relationships between features: $Cov(X) = \frac{1}{n-1}X^{\top}X$ In practice, we compute the covariance matrix using: Cov(X) = np.cov(X, rowvar=False)

• Why rowvar=False ? By default, np.cov assumes rows are variables and columns are observations. Setting rowvar=False ensures rows represent samples, and columns represent features.

2. Eigenvalue Decomposition

The covariance matrix is symmetric, so it can be decomposed into eigenvalues and eigenvectors: $\mathrm{Cov}(X) = Q\Lambda Q^{\top}$

- Q: Orthogonal matrix of eigenvectors.
- Λ : Diagonal matrix of eigenvalues.
- Eigenvalues represent the variance in the direction of the corresponding eigenvectors.

3. Whitening Transformation

To whiten the data, we construct a whitening matrix W using the eigenvalues and eigenvectors: $W=Q\Lambda^{-1/2}Q^{\top}$

- $\Lambda^{-1/2}$: Diagonal matrix where each eigenvalue λ is replaced by $1/\sqrt{\lambda}$.
- ullet Q: Rotates the data to decorrelate it, and $\Lambda^{-1/2}$ scales it to unit variance.

The whitened data is: $X_{
m whitened} = XW$

Implementation Steps

- 1. **Covariance Matrix**: Compute the covariance matrix of X using: covariance = $\operatorname{np.cov}(X, \operatorname{rowvar} = \operatorname{False})$
- 2. **Eigenvalue Decomposition**: Perform eigenvalue decomposition on the covariance matrix: eigvals, eigvecs = np.linalg.eigh(covariance)
 - eigvals : Eigenvalues (variances along principal axes).
 - eigvecs: Eigenvectors (principal axes).
- 3. Whitening Matrix: Construct the diagonal matrix from the eigenvalues: $D = \operatorname{diag}(1/\sqrt{\operatorname{eigvals}})$ Compute the whitening matrix: $W = \operatorname{eigvecs}^{\top}$
- 4. **Transform the Data**: Apply the whitening matrix to the dataset: $X_{ ext{whitened}} = X \cdot W$

Why Whitening is Important

- 1. **Decorrelates Variables**: After whitening, the covariance matrix of the data becomes the identity matrix: $Cov(X_{whitened}) = I$
- Equalizes Variance: Ensures all features have unit variance, simplifying the ICA optimization process.
- 3. **Numerical Stability**: Reduces sensitivity to differences in scaling and correlations, making the algorithm more robust.

Summary

Whitening prepares the data for ICA by removing correlations and standardizing variance. It ensures that the independent component extraction focuses on higher-order statistics, as lower-order correlations are eliminated.

Step 4: Non-Linear Function (_g)

Purpose

The non-linear function in ICA is used to approximate statistical independence by maximizing non-Gaussianity. It operates on the transformed data during optimization and plays a key role in the weight update step.

Mathematical Explanation

Why Non-Linear?

- **Gaussian Distributions**: The central limit theorem states that a sum of independent random variables tends to be Gaussian, even if the original variables are not.
- Non-Gaussianity as a Measure: ICA leverages the idea that independent components
 are less Gaussian than mixtures. A non-linear function enhances the separation of
 independent signals by amplifying non-Gaussian features.

The Hyperbolic Tangent Function (tanh)

One common choice for the non-linear function is the hyperbolic tangent: $g(x) = \tanh(x)$

- **Definition**: $anh(x) = \frac{\sinh(x)}{\cosh(x)} = \frac{e^x e^{-x}}{e^x + e^{-x}}$
- Range: (-1,1)
- **Shape**: A smooth, S-shaped curve that saturates as $x \to \pm \infty$.

Why Use tanh in ICA?

- 1. **Amplifies Non-Gaussianity**: It is non-linear, emphasizing differences in the distribution shape between Gaussian and non-Gaussian signals.
- 2. **Robustness**: Smooth and bounded, preventing extreme values that could destabilize optimization.
- 3. **Mathematical Simplicity**: Its derivative (used in weight updates) is straightforward: $g'(x) = 1 (\tanh(x))^2$

Implementation Details

- ullet Input: A single variable or NumPy array x containing the data to transform.
- Output: The transformed data after applying tanh(x).
- Key Step: Use np.tanh(x) for element-wise application of the hyperbolic tangent.

Summary

The non-linear function tanh(x) is essential for ICA to extract independent components. It enhances non-Gaussianity and stabilizes the optimization process, making it an ideal choice for this step.

Step 5: Derivative of the Non-Linear Function (_g_prime)

Purpose

The derivative of the non-linear function is critical for the optimization process in ICA. It provides the gradient information necessary for updating the weights in the algorithm.

Mathematical Explanation

The Non-Linear Function

The non-linear function used in ICA is: $g(x) = \tanh(x)$

• Recall that tanh(x) is defined as:

$$anh(x)=rac{e^x-e^{-x}}{e^x+e^{-x}}$$

Derivative of tanh(x)

Using calculus, the derivative of tanh(x) can be derived as: $g'(x) = 1 - (tanh(x))^2$

Why?

- 1. From the definition of tanh(x): $tanh(x) = \frac{\sinh(x)}{\cosh(x)}$
 - The derivative of $\sinh(x)$ is $\cosh(x)$.
 - The derivative of $\cosh(x)$ is $\sinh(x)$.
- 2. Applying the quotient rule: $rac{d}{dx} anh(x)=rac{\cosh(x)\cosh(x)-\sinh(x)\sinh(x)}{\cosh^2(x)}$
 - ullet Simplify the numerator using the hyperbolic identity: $\cosh^2(x) \sinh^2(x) = 1$
 - Thus: $g'(x)=rac{1}{\cosh^2(x)}$
- 3. Using the identity: $\frac{1}{\cosh^2(x)} = 1 (\tanh(x))^2$

• The result is: $g'(x) = 1 - (\tanh(x))^2$

Why This Formula is Useful

1. Efficiency:

• The derivative is computed using only the value of tanh(x), avoiding recomputation of cosh(x) or sinh(x).

2. Optimization:

• The derivative scales the weight updates in ICA, guiding the algorithm to converge faster and more accurately.

Implementation Details

- ullet Input: A single variable or NumPy array x containing the data.
- Output: The derivative g'(x), computed element-wise as $1 (\tanh(x))^2$.
- Key Step: Use np.tanh(x) to compute the tangent hyperbolic and square it to obtain the derivative.

Summary

The derivative of the non-linear function, $g'(x) = 1 - (\tanh(x))^2$, is vital for ICA as it enables gradient-based optimization. Its simplicity and efficiency make it a natural choice for this process.

Step 6: Fitting the ICA Model (fit)

Purpose

The fit method trains the ICA model on the given dataset. It finds the unmixing matrix (W) that separates mixed signals into statistically independent components.

Key Steps in fit

1. Preprocessing

Before starting the ICA optimization, the input data (X) must be:

1. **Centered**: Remove the mean of each feature to ensure zero-mean data.

$$X = \operatorname{self.} \setminus \operatorname{center}(X)$$

2. **Whitened**: Decorrelate and scale the data to unit variance. $X = \text{self.} \setminus \text{whiten}(X)$

2. Initialize Parameters

• **Weight Matrix**: Initialize the unmixing matrix (W) with random values:

```
W = \text{np.random.randn}(\text{self.n} \setminus \text{components}, n\_features)
```

Shape Parameters: Determine the number of samples and features:

```
n\_samples, n\_features = X. shape
```

1. Preprocessing

Before starting the ICA optimization, the data must be prepared:

1. Centering:

- Subtract the mean from each feature to ensure the data has zero mean.
- This step is critical because ICA assumes the input data is centered.
- Use the _center method you implemented earlier: $X = \text{self.} \setminus \text{_center}(X)$

2. Whitening:

- Decorrelate and normalize the data so that each feature has unit variance.
- This step simplifies the ICA optimization by aligning the data into a spherical space.
- Use the _whiten method: $X = \text{self.} \setminus \text{whiten}(X)$

2. Initialization

The following parameters must be initialized:

1. Number of Samples and Features:

ullet Determine the dimensions of the input data (X): $n_{
m samples}, n_{
m features} = X. {
m shape}$

2. Weight Matrix:

- Initialize the unmixing matrix (W) with random values:
 - $W = \text{np.random.randn}(\text{self.n} \setminus \text{components}, n_{\text{features}})$
- This matrix will be iteratively updated to find the independent components.

Summary of Missing Steps

- 1. Preprocess the input data by centering and whitening it.
- 2. Initialize the necessary variables: the number of samples, features, and the random weight matrix.

These steps prepare the data and parameters for the iterative optimization process, which is already included in the provided code.

3. Iterative Optimization

- 1. **Iterate**: Run the optimization loop for up to self.max_iter iterations: for i in range(self.max_iter):
- 2. **Update Each Component**: For each component (j), calculate:
 - (wx = X \cdot W[j]): The projection of (X) onto the (j)-th weight vector.
 - $(g_wx = g(wx))$: Apply the non-linear function.
 - (g_prime_wx = g'(wx)): Compute its derivative.
 - Update (W[j]) using the formula: $W[j] = \frac{X^{ op} \cdot g wx \text{mean}(g prime wx) \cdot W[j]}{n samples}$
 - Normalize (W[j]) to have unit length: $W[j]/=\|\|W[j]\|\|$
- 3. **Decorrelate the Weight Matrix**: After updating all components, decorrelate (W) to ensure orthogonality: $W = \text{self.} \setminus \text{decorrelate}(W)$
- 4. **Check Convergence**: Compute the maximum absolute change in the diagonals of (W \cdot W_{\text{old}}^\top):

```
\max_abs_cange = np.\max(np.abs(np.diag(W @ W_old.T)) - 1)) If the change is less than self.tol , stop the iterations.
```

4. Finalize

- ullet Store the final weight matrix as the learned components: $\mathrm{self.components} ackslash = W$
- Return the fitted model instance.

Summary

The fit method optimizes the weight matrix (W) to separate independent components by:

- 1. Preprocessing the data (centering and whitening).
- 2. Iteratively updating (W) using gradient-based optimization and ensuring orthogonality.
- 3. Stopping when the convergence criterion is met or the maximum iterations are reached.

Step 7: Decorrelating the Weight Matrix (_decorrelate)

Purpose

The goal of this step is to ensure that the rows of the weight matrix W remain orthogonal during the ICA optimization process. Orthogonality is critical because independent components should be extracted without overlapping or redundancy.

Mathematical Explanation

Why Decorrelate?

1. Orthogonality:

 In ICA, the weight matrix W should ideally represent independent directions in the feature space. This requires its rows (or components) to be orthogonal to each other.

2. Stability:

• Without decorrelation, the optimization process might converge to suboptimal or unstable solutions due to redundancy between components.

Matrix Orthogonality

A matrix W is orthogonal if: $WW^{\top} = I$ where I is the identity matrix.

Steps to Decorrelate

1. Compute WW^{\top} :

 \bullet The product WW^{\top} gives a symmetric matrix representing the correlation between rows of W .

2. Eigenvalue Decomposition:

- ullet Perform eigenvalue decomposition on $WW^ op\colon WW^ op = Q\Lambda Q^ op$
 - Q: Orthogonal matrix of eigenvectors.
 - Λ : Diagonal matrix of eigenvalues.

3. Construct Normalization Matrix:

• Create a diagonal matrix D that scales each eigenvalue by its inverse square root: $D={
m diag}(1/\sqrt{\lambda})$ where λ are the eigenvalues.

4. Transform W:

ullet Normalize W using the eigenvectors and the normalization matrix:

$$W_{ ext{decorrelated}} = QDQ^{ op}W$$

ullet This ensures that $W_{ ext{decorrelated}}W_{ ext{decorrelated}}^ op=I.$

Implementation Steps

- 1. Compute WW^{\top} .
- 2. Perform eigenvalue decomposition on $WW^{ op}$ to obtain eigenvalues and eigenvectors.
- 3. Construct the normalization matrix ${\cal D}$ from the eigenvalues.
- 4. Apply the transformation to W using the eigenvectors and D.

Why This Works

- 1. The eigenvalue decomposition aligns the matrix to an orthogonal basis.
- 2. Normalizing by the eigenvalues ensures unit-length scaling in each direction, resulting in an orthogonal matrix.

Summary

The _decorrelate method ensures that the weight matrix remains orthogonal, a critical property for extracting independent components. This step stabilizes and improves the convergence of the ICA algorithm.

Your Task

- 1. Implement the _decorrelate method based on the explanation.
- 2. Test it on a sample matrix W to ensure it returns an orthogonal matrix.

Step 8: Transforming the Data (transform)

Purpose

The transform method applies the trained ICA model to a new dataset X. It projects the input data onto the learned independent components to extract their representations.

Mathematical Explanation

1. Preprocessing

Before applying the learned components, the input data X must be:

1. Centered:

- ullet Remove the mean of each feature: $X_{
 m centered} = X \mu$
- This ensures that the data aligns with the assumptions of the trained ICA model.

2. Whitened:

- ullet Decorrelate and scale the data to unit variance: $X_{
 m whitened} = W_{
 m whitening} \cdot X_{
 m centered}$
- Whitening simplifies the subsequent projection step by normalizing the feature space.

2. Projection onto Independent Components

Once the data is preprocessed, it is projected onto the learned independent components:

$$S = X_{\text{whitened}} \cdot W_{\text{components}}^{\top}$$

- ullet $W_{
 m components}$: The weight matrix learned during the fit step, which contains the directions of the independent components.
- S: The extracted independent components of the input data.

Key Steps in transform

1. Center the Data:

Use the _center method to ensure zero-mean input.

2. Whiten the Data:

Use the _whiten method to decorrelate and normalize the input.

3. Project onto Components:

Multiply the preprocessed data by the transpose of self.components_.

Why This Works

- The learned matrix self.components_ captures the directions of the independent components.
- 2. By projecting the preprocessed data onto these directions, we recover the independent components present in the new data.

Summary

The transform method extracts independent components from a new dataset by:

- 1. Centering and whitening the data.
- 2. Projecting it onto the learned independent components.

Step 9: Combining Fitting and Transformation (fit_transform)

Purpose

The fit_transform method combines two key steps:

- 1. Fit:
 - Trains the ICA model on the dataset X to learn the independent components.
- 2. Transform:
 - ullet Projects the input data X onto the learned independent components to extract their representations.

This method is useful for workflows where the dataset X is both the training set and the data to be transformed.

Mathematical Explanation

1. Fitting the ICA Model

The fit step involves:

- 1. Centering:
 - Remove the mean of each feature in X.
- 2. Whitening:
 - Decorrelate and normalize the features to unit variance.
- 3. Optimization:
 - Iteratively update the weight matrix (W) to maximize the statistical independence of the components.
- 4. Learned Components:
 - Store the final weight matrix (W) as self.components_.

2. Transforming the Data

After fitting the model, the transform step:

- 1. Preprocesses the input data (centering and whitening).
- 2. Projects it onto the independent components: $S = X_{ ext{whitened}} \cdot W_{ ext{components}}^{ op}$
- 3. Returns the independent components (S).

Implementation Steps

1. Call the fit Method:

- ullet Train the ICA model on X to learn the independent components.
- This step populates self.components_ with the weight matrix (W).

2. Call the transform Method:

• Use the trained model to extract independent components from X.

3. Return the Transformed Data:

• The method returns the independent components as the output.

Why fit_transform Works

1. Combines Steps:

 By combining fit and transform, the method provides a seamless way to preprocess, train, and extract independent components in one step.

2. Reusability:

 After fit_transform is called, the model is trained and can be reused to transform new datasets.

Summary

The fit_transform method first trains the ICA model using fit and then extracts independent components from the same dataset using transform. It combines these steps into a single, convenient method.

```
import numpy as np

class ICA:
    def __init__(self, n_components, max_iter=1000, tol=1e-5):
        self.n_components = n_components
```

```
self.max_iter = max_iter
    self.tol = tol
def _center(self, X):
    """Centers the dataset (zero mean)."""
    self.mean = X.mean(axis=0) # Average of the columns
    X = X-self.mean # Centering the dataset
    return X
def _whiten(self, X):
    """Whitens the dataset (decorrelates the variables)."""
    covariance = np.cov(X, rowvar=False) # Covariance matrix of the matrix X, r
    eig_vals, eig_vecs = np.linalg.eigh(covariance) #Eigenvalue decomp
    self.whitening_matrix = eig_vecs@np.diag(1.0 / np.sqrt(eig_vals))@eig_vecs.
    X_whitened = X@self.whitening_matrix # The altered x matrix
    return X_whitened
def _g(self, x):
    """Non-linear function (logistic)."""
    return np.tanh(x)
def _g_prime(self, x):
    """Derivative of the non-linear function."""
    return 1-np.tanh(x)**2
def fit(self, X):
    X= self._center(X)
    X = self._whiten(X)
    n_samples, n_features = X.shape
    W= np.random.randn(self.n_components,n_features)
    print("Starting ICA fitting...")
    for i in range(self.max_iter):
        W_old = W_copy()
        for j in range(self.n_components):
            wx = np.dot(X, W[j])
            g_wx = self._g(wx)
            g_prime_wx = self._g_prime(wx)
            W[j] = (X.T @ g_wx - np.mean(g_prime_wx) * W[j]) / n_samples
            W[j] /= np.linalg.norm(W[j])
        # Decorrelate weights
        W = self._decorrelate(W)
        # Check convergence
        max_abs_change = np.max(np.abs(np.abs(np.diag(W @ W_old.T)) - 1))
        print(f"Iteration {i}, change: {max_abs_change}")
        if max_abs_change < self.tol:</pre>
            print("Converged!")
            break
    self.components_ = W
    print("ICA fitting completed.")
    return self
def decorrelate(self, W):
```

```
"""Ensures that the matrix W remains orthogonal."""
        matrix = W@W.T #Multiply W and its transpose
        eig_vals, eig_vecs = np.linalg.eigh(matrix) #Eigenvalue decomp
        return eig_vecs@np.diag(1.0 / np.sqrt(eig_vals))@eig_vecs.T@W
    def transform(self, X):
        X = self._center(X) #centering the data
        X= self._whiten(X) #Whitening the data
        return X@self.components_.T #Multiply components by altered X
    def fit_transform(self, X):
        self.fit(X)
        return self.transform(X)
# Example usage:
# X is your dataset with shape (n_samples, n_features)
# n_components is the number of independent components you want to extract
# Generate a random dataset for demonstration purposes
# Create synthetic sources
np.random.seed(0)
n_samples = 1000
time = np.linspace(0, 8, n_samples)
# Create two independent signals
s1 = np.sin(2 * time) # Sinusoidal signal
s2 = np.sign(np.sin(3 * time)) # Square signal
# Stack signals into a matrix
S = np.c_[s1, s2]
# Mix the signals with a random mixing matrix
A = np.array([[1, 1], [0.5, 2]]) # Mixing matrix
X = S @ A.T # Generate observations by mixing the signals
# Apply ICA to separate the sources
ica = ICA(n_components=2)
S_ = ica.fit_transform(X)
print("Original Signals (first 5 samples):")
print(S[:5])
print("\nRecovered Signals (first 5 samples):")
print(S_[:5])
import matplotlib.pyplot as plt
plt.figure(figsize=(12, 6))
# Original signals
plt.subplot(2, 1, 1)
plt.plot(S[:100], label="Original Signals")
plt.title("Original Signals")
plt.legend(loc='upper right')
```

```
# Recovered signals
 plt.subplot(2, 1, 2)
 plt.plot(S_[:100], label="Recovered Signals")
 plt.title("Recovered Signals")
 plt.legend(loc='upper right')
 plt.tight_layout()
 plt.show()
Starting ICA fitting...
Iteration 0, change: 1.3055495001726078
Iteration 1, change: 0.0012452944914410624
Iteration 2, change: 0.0005997391255762041
Iteration 3, change: 0.0002734536808511878
Iteration 4, change: 0.0001215131822759341
Iteration 5, change: 5.337125746307514e-05
Iteration 6, change: 2.3320898175915872e-05
Iteration 7, change: 1.0166962624569287e-05
Iteration 8, change: 4.427914454985249e-06
Converged!
ICA fitting completed.
Original Signals (first 5 samples):
[[0.
              0.
 [0.01601533 1.
 [0.03202655 1.
 [0.04802956 1.
                         1
 [0.06402025 1.
                         11
Recovered Signals (first 5 samples):
[[-0.17517734 -0.03858102]
 [-0.20119239 0.96408661]
 [-0.1779273
                0.96322394]
 [-0.15467414 0.96236171]
 [-0.13143888 0.96150014]]
                                          Original Signals
                                                                                  Original Signals
                                                                                 Original Signals
0.8
0.6
0.4
0.0
                                       40
                                                                                       100
                                         Recovered Signals
1.2
                                                                                Recovered Signals
                                                                                Recovered Signals
1.0
0.8
0.6
0.2
0.0
                                                                                       100
```