A Geometric Approach for Separating Several Speech Signals^{*}

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Abstract. In this paper a new geometrical approach for separating speech signals is presented. This approach can be directly applied to separate more than two speech signals. It is based on clustering the observation points, and then fitting a line (hyper-plane) onto each cluster. The algorithm quality is shown to be improved by using DCT coefficients of speech signals, as opposed to using speech samples.

1 Introduction

Blind Source Separation (BSS) or Independent Component Analysis (ICA) consists in retrieving unknown statistically independent signals from their observed mixtures, assuming there is no information about the original source signals, or about the mixing system (hence the term *Blind*).

For linear instantaneous mixtures $\mathbf{x}(t) = \mathbf{As}(t)$, where the sources $\mathbf{s}(t) \triangleq (s_1(t), \ldots, s_N(t))^T$ are (unknown) statistically independent signals, the observation signals are denoted $\mathbf{x}(t) \triangleq (x_1(t), \ldots, x_N(t))^T$, and \mathbf{A} is the $N \times N$ (unknown) mixing matrix. In this paper, the number of observations and sources are assumed to be equal. The problem is then to estimate the source vector $\mathbf{s}(t)$ only by knowing the observation vector $\mathbf{x}(t)$.

One approach to solve the problem is to determine a separating matrix **B** such that the outputs $\mathbf{y}(t) \triangleq \mathbf{Bx}(t)$ become statistically independent. This independence insures the estimation of the sources, up to a scale and a permutation indeterminacy [1].

Another approach is the geometric source separation algorithm, which has been first introduced in [2]. In this approach (for 2-dimensional case), it is

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Fig. 1. Distribution of a) source samples, and b) observation samples.



Fig. 2. Distribution of a) two speech samples, and (b) their mixtures.

first noted that because of the independence of source signals, $p_{s_1s_2}(s_1, s_2) = p_{s_1}(s_1)p_{s_2}(s_2)$, where p stands for the Probability Density Function (PDF). Consequently, for bounded sources, the points (s_1, s_2) will be distributed in a rectangular region (Fig. 1-a). Now, because of the scale indeterminacy, the mixing matrix is assumed to be of the form (normalized with respect to diagonal elements):

$$\mathbf{A} = \begin{pmatrix} 1 & a \\ b & 1 \end{pmatrix} \tag{1}$$

Under the transformation $\mathbf{x} = \mathbf{As}$, the rectangular region of the *s*-plane will be transformed into a parallelogram (Fig. 1-b), and the slopes of the borders of this parallelogram are 1/a and *b*. In other words, for estimating the mixing matrix, it is sufficient to determine the slopes of the borders of the distribution of the observation samples.

Although this approach is not easily generalized to higher dimensions, it is successful in separating two sources, provided that their distributions allow a good estimation of the borders of the parallelogram (*e.g.* uniform and sinusoidal sources). However, this technique cannot be used in separating speech signals because the PDF of a speech is mostly concentrated about zero. This comes from the fact that in a speech signal, there are many low energy (silence or unvoiced) sections. Consequently, as it can be seen in Fig. 2, it is practically impossible to find the borders of the parallelogram when the sources are speech signals. This is explained by a probabilistic manner in [3]: the probability of having a point in the borders of the parallelogram is very low.

Although for speech signals the borders of the parallelogram are not visible in Fig. 2, there are two visible "axes", corresponding to lines $s_1 = 0$ and $s_2 = 0$ in

the s-plane (throughout the paper, it is assumed that the sources and hence the observations have zero-means). The slopes of these axes, too, determine a and b in (1). In other words, for speech signals, instead of finding the borders, we try to find these axes. This idea is used in [3] for separating speech signals by utilizing an "angular" histogram for estimating these axes. In this method, the resolution of the histogram cannot be too fine (requires more data points), and cannot be too coarse (bad estimation of the mixing matrix). Moreover, this approach cannot be easily generalized to mixtures of more than two speech signals.

In this paper, we propose another approach for estimating these "axes" based on line (or hyper-plane) fitting. The main idea is to fit two lines on the scatter plot of observations, which will be the required axes. This approach does not suffer from the problem of the resolution of a histogram. Moreover, we will see that this approach can be directly used in higher dimensions.

2 Speech Separation by Line Fitting

2.1 Two Dimensional Case

As it is explained in the previous section, the main idea of our method is to estimate the slopes of two axes of the scatter plot of observations (Fig. 2-b). These axes corresponds to the lines $s_1 = 0$ and $s_2 = 0$ in the scatter plot of sources. The existence of these lines is a result of many low-energy sections of a speech signal. For example, the points with small s_1 and different values for s_2 will be concentrated about the axis $s_1 = 0$.

However, we do not use (1) as a model for mixing matrix, because it has two restrictions. Firstly, in this model, it is implicitly assumed that the diagonal elements of the actual mixing matrix are not zero, otherwise infinite values for a and b may be encountered (this situation corresponds to vertical axes in the x-plane). Secondly, this approach is not easy to be generalized to higher dimensions.

Instead of model (1), let us consider a general "separating matrix" $\mathbf{B} = [b_{ij}]_{2\times 2}$. Under the transformation $\mathbf{y} = \mathbf{B}\mathbf{x}$, one of the axes must be transformed to $y_1 = 0$, and the other to $y_2 = 0$. In other words, for every (x_1, x_2) on the first axis:

$$\begin{pmatrix} 0\\y_2 \end{pmatrix} = \begin{pmatrix} b_{11} & b_{12}\\b_{21} & b_{22} \end{pmatrix} \begin{pmatrix} x_1\\x_2 \end{pmatrix} \Rightarrow b_{11}x_1 + b_{12}x_2 = 0$$
(2)

That is, the equation of the first axis is $b_{11}x_1 + b_{12}x_2 = 0$. In a similar manner, the second axis will be $b_{21}x_1 + b_{22}x_2 = 0$. Consequently, for estimating the separating matrix, the equations of the two axes must be found in the form of $\alpha_1x_1 + \alpha_2x_2 = 0$, and then each row of the separating matrix is composed of the coefficients of one of the axes. For finding the axes we suggest is to "fit" two straight lines on the scatter plot of the observations.

It is seen that by this approach, we are not restricted to non-vertical axes (non-zero diagonal elements of the mixing matrix). More interestingly, this approach can be directly used in higher dimensions, as stated below.

2.2 Higher Dimensions

The approach stated above can be directly generalized to higher dimensions. For example, for 3 speech signals and 3 sources, the low-energy (silence and unvoiced) values of s_1 with different values of s_2 and s_3 will form the plane $s_1 = 0$ in the 3-dimensional scatter plot of sources. Hence, in this 3-dimensional scatter plot, there are 3 visible planes: $s_1 = 0$, $s_2 = 0$ and $s_3 = 0$. These planes will be transformed to three main planes in the scatter plot of observations. With calculations similar as (2), it is seen that each row of the separating matrix is composed of the coefficients of one of these main planes in the form of $\alpha_1 x_1 + \alpha_2 x_2 + \alpha_3 x_3 = 0$.

Consequently, for N-dimensional case, N (hyper-)planes in the form of $\alpha_1 x_1 + \cdots + \alpha_N x_N = 0$ must be first "fitted" onto the scatter plot of observations. Then, each row of the separating matrix is the coefficients $(\alpha_1, \ldots, \alpha_N)$ of one of these (hyper-)planes.

3 Line Fitting

To use the idea of the previous section, we need a method for fitting two lines (or N hyper-planes) onto the scatter plot of observations.

3.1 Fitting a Straight Line onto a Set of Points

First of all, consider the problem of fitting a line onto K data points $(x_i, y_i)^T$, i = 1...K. In the traditional least squares method, this is done by finding the line y = mx + h which minimizes $\sum_{i=1}^{K} (y - y_i)^2 = \sum_{i=1}^{K} (mx_i + h - y_i)^2$. This is equivalent to minimizing the "vertical" distances between the line and the data points, as shown in Fig. 3-a. This technique is mainly used in linear regression analysis where there are errors in y_i 's, but not in x_i 's.

However, in our application of fitting a line onto a set of points, a better measure is minimizing the sum of "orthogonal distances" between the points and the line, as shown in Fig. 3-b. Moreover, as discussed in the previous sections, we are seeking a line in the form ax + by = 0. Consequently, the best fitted line is determined by minimizing $\sum_{i=1}^{K} d_i^2$, where d_i is the orthogonal distance between the *i*-th point and the line:

$$d_i = \frac{|ax_i + by_i|}{\sqrt{a^2 + b^2}} \tag{3}$$

However, ax + by = 0 is not uniquely determined by a pair (a, b), because (ka, kb) represents the same line. To obtain a unique solution, the coefficients are normalized such that $a^2 + b^2 = 1$. To summarize, the best fitted line ax + by = 0 is obtained by minimizing $\sum_{i=1}^{K} (ax_i + by_i)^2$ under the constraint $a^2 + b^2 = 1$.

N-Dimensional Case. In a similar manner, an *N*-dimensional hyper-plane $\alpha_1 x_1 + \alpha_2 x_2 + \cdots + \alpha_N x_N = 0$ is fitted onto a set of *K* data points $\mathbf{x}_i = (x_1^{(i)}, x_2^{(i)}, \dots, x_N^{(i)})^T$, $i = 1, \dots, K$ by minimizing the cost function:



Fig. 3. a) Least squares line fitting, b) Orthogonal line fitting.

$$\mathcal{C}(\alpha_1,\ldots,\alpha_N) = \sum_{i=1}^K \left(\alpha_1 x_1^{(i)} + \cdots + \alpha_N x_N^{(i)}\right)^2 \tag{4}$$

under the constraint $g(\alpha_1, \ldots, \alpha_N) \equiv \alpha_1^2 + \cdots + \alpha_N^2 - 1 = 0.$

Solution. Using Lagrange multipliers, the optimum values for $\alpha_1, \ldots, \alpha_N$ satisfy $\nabla C = \lambda \nabla g$. After a few algebraic calculations, this equation is written in the matrix form:

$$\mathbf{R}_{\mathbf{x}}\boldsymbol{\alpha} = \frac{\lambda}{K}\boldsymbol{\alpha} \tag{5}$$

where $\boldsymbol{\alpha} \triangleq (\alpha_1, \ldots, \alpha_N)^T$ and $\mathbf{R}_{\mathbf{x}} \triangleq \frac{1}{K} \sum_{i=1}^K \mathbf{x}_i \mathbf{x}_i^T$ is the correlation matrix of data points. Equation (5) shows that λ/K and $\boldsymbol{\alpha}$ are eigen value and eigen vector of the correlation matrix $\mathbf{R}_{\mathbf{x}}$, respectively. Moreover:

$$\mathcal{C} = \sum_{i=1}^{K} \left(\boldsymbol{\alpha}^{T} \mathbf{x}_{i} \right)^{2} = \sum_{i=1}^{K} \boldsymbol{\alpha}^{T} \mathbf{x}_{i} \mathbf{x}_{i}^{T} \boldsymbol{\alpha} = K \boldsymbol{\alpha}^{T} \mathbf{R}_{\mathbf{x}} \boldsymbol{\alpha} = \lambda \boldsymbol{\alpha}^{T} \boldsymbol{\alpha} = \lambda$$

and hence for minimizing the cost function, λ must be minimum. Consequently, the solution of the hyper-plane fitting problem is given by the eigen vector of the correlation matrix which corresponds to its minimum eigen value.

Discussion. It is interesting to think about the conjunction of the above approach to Principal Component Analysis (PCA). Note that $\boldsymbol{\alpha}$ is the vector perpendicular to the plane $\alpha_1 x_1 + \cdots + \alpha_N x_N = 0$, and the above theorem states that this vector must be chosen in the direction with the minimum spread of data points, which is compatible with our heuristic interpretations of plane (line) fitting. This method has old foundations in mathematics [4], and somewhat called Principal Component Regression (PCR) [5].

3.2 Fitting 2 Straight Lines (N Hyper-planes)

However, as stated in Section 2, for 2 mixtures of 2 sources our problem is to fit 2 lines onto the observation points, not just 1 line. In other words, as it is seen in Fig. 2, we need to divide the data points into 2 clusters, and then to fit a line onto the points of each cluster. The extension to N mixtures of N sources

- Initially distribute the points into clusters S_1, \ldots, S_N (*e.g.* random initialization).
- Loop:
 - 1. Fit a line (hyper-plane) onto each set of points S_i (we call it l_i).
 - 2. Recalculate the clusters: Let S_i be the set of all points which are closer to line (hyper-plane) l_i than other lines (hyper-planes), that is:

$$\mathcal{S}_i = \{ \mathbf{x} \mid d(\mathbf{x}, l_i) < d(\mathbf{x}, l_j), \forall j \neq i \}$$

- Repeat until convergence.

Fig. 4. Algorithm of fitting two lines (N hyper-planes) onto a set of points.

is straightforward: we need to divide the data into N clusters, and then to fit a hyper-plane onto the points of each cluster. Mathematically, this is equivalent to minimizing the following cost function (for the N-dimensional case):

$$\mathcal{C} = \sum_{\mathbf{x}_i \in \mathcal{S}_1} d^2(\mathbf{x}_i, l_1) + \sum_{\mathbf{x}_i \in \mathcal{S}_2} d^2(\mathbf{x}_i, l_2) + \dots + \sum_{\mathbf{x}_i \in \mathcal{S}_N} d^2(\mathbf{x}_i, l_N)$$
(6)

where S_j is the *j*-th cluster of points and $d^2(\mathbf{x}_i, l_j)$ denotes the perpendicular distance of the *i*-th point from the *j*-th plane.

Having divided the points into clusters S_1, \ldots, S_N , the previous section gives us the best line fitted onto the points of each cluster. For clustering the data points, we use the algorithm stated in Fig. 4, which is inspired from the *k*-means (or Lloyd) algorithm for data clustering [6]. Its difference with *k*-means is that in *k*-means, each cluster is mapped onto a point (point \rightarrow point), but in our algorithm each cluster is mapped onto a line or hyper-plane (point \rightarrow line).

The following theorem is similar to a corresponding theorem for the k-means algorithm [6].

Theorem 1. The algorithm of Fig. 4 converges in a finite number of iterations.

Proof. At each iteration, the cost function (6) cannot be increased. This is because in the first step (fitting hyper-planes onto the clusters) the cost function is either decreased or does not change. In the second step, too, the redistribution of the points in the clusters is done such that it decreases the cost function or does not change it. Moreover, there is a finite number of possible clustering of finite number of points. Consequently, the algorithm must converge in a finite number of iterations. \Box

Initialization. The fact that the cost-function is non-increasing in the algorithm, shows that the algorithm may get trapped in a local minimum. This is one of major problems of the k-means algorithm, too. It depends on the initialization of the algorithm, and become more severe when the dimensionality increases. In k-means, one approach is to run the algorithm with several randomly chosen initializations, and then to take the result which produces the minimum cost-function.



Fig. 5. Distribution of a) the observations, and b) their DCT coefficients (right).

4 Final Algorithm, and Its Improvement by Using DCT

The final separation algorithm is now evident. First, run the algorithm of Fig. 4. After convergence, there are N lines (hyper-planes) $l_i : \alpha_{i1}x_1 + \cdots + \alpha_{iN}x_N = 0$, $i = 1 \dots N$. Then, the *i*-th row of the separating matrix is $(\alpha_{i1}, \dots, \alpha_{iN})$.

However, the separation quality of the algorithm can be improved, with a simple trick. Recall that the success of the algorithm is because of the existence of two visible "axes" in Fig. 2. These axes were formed because of the small-valued (low-energy) parts of one speech and other parts of the second one. Now, recall that the Discrete Cosine Transform (DCT) coefficients of a speech frame (10-20 msec) contain a lot of nearly zero values. Moreover, DCT is a linear transformation, and hence, the DCT coefficients of the observations are a mixture of the DCT coefficients of the original speeches with the same mixing matrix. Therefore, it seems that it is a good idea to apply the algorithm on the DCT coefficients of observations, and that of their DCT coefficients. It is seen visually that the "axes" are more visible in the scatter plot of DCT coefficients. Consequently, one expects to get better results by applying the algorithm on the DCT coefficients of the observations, as is confirmed by our experiments, too.

5 Experimental Results

Many simulations have been conducted to separate 2, 3 or 4 sources. In all these simulations, typically less than 30 iterations are needed to achieve separation. The experimental study shows that local minima depends on the initialization phase of the algorithm and on the number of sources (local minima have been never encountered in separating two sources).

Here, the simulation results of 4 typical speech signals (sampled with 8KHz sampling rate) are presented. In all the experiments, the diagonal elements of the mixing matrix are 1, while all other elements are 0.5. For each simulation, 10 random initializations are used, and then the matrix which creates minimum cost-function is taken as the answer.

To measure the performance of the algorithm, let $\mathbf{C} \triangleq \mathbf{B}\mathbf{A}$ be the global mixing-separating matrix. Then, we define the Signal to Noise Ratio by (assuming no permutation) $\mathrm{SNR}_i(\mathrm{in} \,\mathrm{dB}) \triangleq 10 \log_{10} \frac{c_{ii}^2}{\sum_{j \neq i} c_{ij}^2}$. This criterion shows



Fig. 6. Separation result in separating N speech signals, a) N = 2, b) N = 3, c) N = 4.

how much the global matrix **C** is close to the identity matrix. As a performance criterion of the algorithm, we take the average of the SNR's of all outputs: $\text{SNR} = \frac{1}{N} \sum_{i} \text{SNR}_{i}$. To virtually create different source signals, each speech signals is shifted randomly in time (more precisely, each speech signal is shifted 128k samples, where k is a randomly chosen integer). This results in a completely different source scatter plot, and virtually creates a new set of source signals. Then, for each experiment, the algorithm is run 50 times (with 50 different random shifts), and the averaged SNR is calculated.

Figure 6 shows this averaged SNR's with respect to number of samples, for separating 2, 3 and 4 speech signals. The figure clearly shows the ability of the algorithm for speech separation, and the advantage obtained by using DCT coefficients. Moreover, it is seen that when the number of sources increases, more data samples are required to reach a given separation quality. This was expected, because the algorithm is based on the "sparsity" of the speech signals. In other words, for forming the planes, it is required that one speech signal is low-energy (silence/unvoiced), and the others are not. If p is the probability of being in a low energy state, the probability of sparsity is $p(1-p)^{(N-1)}$, which decreases exponentially with N.

6 Conclusion

In this paper, a geometrical approach for separating several speech signals has been presented. It has been shown that for speech signals (or other sources whose PDF's are concentrated about zero), the ICA can be accomplished by a clustering of observation samples and then applying a PCA on each cluster and taking the smallest principal component. Although this approach was based on geometric interpretations, its final algorithm is completely algebraic.

Initialization is the main problem of this algorithm. Finding better initialization approaches is currently under study.

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