

ICAR: INDEPENDENT COMPONENT ANALYSIS USING REDUNDANCIES

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ABSTRACT

A new Blind Source Separation (BSS) algorithm, called ICAR and using only Fourth Order (FO) statistics of the data, is proposed. The latter method is compared by computer experiments with the well-known methods COM1, COM2, JADE, FastICA and SOBI. Since ICAR has given very good convergence results and has performed the source separation in the presence of a Gaussian noise with unknown spatial correlation, it appears as being one of the most attracting BSS algorithms.

1. INTRODUCTION

Blind Source separation (BSS) or Independent Component Analysis (ICA) have lately raised great interest. These problems find their place in numerous applications including telecommunications, speech processing, or biomedical engineering. For instance, in antenna array processing, if several sources impinge of an array of discrete sensors, and if the channel time spread associated with every source and every sensor is negligible compared to the symbol period, then the signals received can be modeled as a static mixture of the transmitted sources. BSS aims at extracting the sources from the sole observation of the mixtures received on the array. Several techniques are available, depending on the assumptions made. In this paper, it will be assumed that sources are mutually statistically independent. Contrary to Principal Component Analysis (PCA), which exploits only statistics of order 2, ICA resorts to statistics of higher order, and is thus able to impose some stronger statistical independence than just a mere decorrelation. This is made possible if sources are not Gaussian, and made easier if there are at least as many sensors as sources.

Since the early works of Jutten [11], the concept of ICA has evolved, and most of the material has been presented in the seminal paper of Comon [7], where it is proposed to maximize contrast functionals; an algorithm is proposed there, which shall be referred to as COM2, for Contrast-based Maximization of squared fourth order cumulants. Some years later, he also proposed a simple algorithm [8], which will be called COM1 in the sequel, maximizing a

squares-free criterion [13]. On the other hand, Cardoso and Souloumiac have proposed [4] slightly later an efficient algorithm performing Joint Approximate Diagonalization of Eigen cumulant matrices (JADE). Later, Hyvarinen and others propose the so-called FastICA fixed point algorithm, which extracts one source at a time, first for real signals [12], then for complex [3]. All these methods can be sensitive to the presence of additive Gaussian noise, of unknown spatial coherence; such a noise is frequent in applications such as HF radiocommunications. In fact, they perform as a preliminary stage an exact second-order decorrelation by a preliminary “spatial whitening”; this operation is sub-optimal in several respects, because it puts too much emphasis on second order statistics.

The solution (also sub-optimal though) proposed in this paper consists of getting rid of this whitening stage, and of using exclusively higher order statistics, namely fourth order cumulants. More precisely, the redundancy theoretically present in the quadricovariance of the observations is exploited. This concept can be extended to statistics of order strictly higher than 4, allowing for instance to address the case of *underdetermined mixtures*, in which the number of sources present exceeds in permanence the number of sensors. Such extensions to order 6, or more generally to order $m=2q$ ($q \geq 2$), have been recently proposed by the authors under the names of BIRTH [2] and BIOME [1].

2. NOTATIONS AND ASSUMPTIONS

Let N sensors be available, and denote $x_n(k)$ the noisy observations received on these sensors, $1 \leq n \leq N$. The vector of observations $\mathbf{x}(k) \stackrel{\text{def}}{=} [x_1(k) \ x_2(k) \ \cdots \ x_N(k)]^T$ can be modeled in the form below:

$$\mathbf{x}(k) = \mathbf{A} \mathbf{s}(k) + \boldsymbol{\nu}(k) = \sum_{p=1}^P \mathbf{a}_p s_p(k) + \boldsymbol{\nu}(k) \quad (1)$$

where $s_p(k)$ denotes the unknown sources, and $\mathbf{A} \stackrel{\text{def}}{=} [\mathbf{a}_1 \ \mathbf{a}_2 \ \cdots \ \mathbf{a}_P]$, $\mathbf{s}(k) \stackrel{\text{def}}{=} [s_1(k) \ s_2(k) \ \cdots \ s_P(k)]^T$, $\boldsymbol{\nu}(k) \stackrel{\text{def}}{=} [\nu_1(k) \ \nu_2(k) \ \cdots \ \nu_N(k)]^T$ denote respectively the $N \times P$ mixing matrix, and the source and noise random vectors,

assumed to be independent. Also define for very index k the entries of the 4-th order cumulant tensor, C_x , of a random process $x(k)$, stationary and ergodic up to order 4:

$$C_{de,x}^{fg} \stackrel{\text{def}}{=} \text{Cum}\{x_d(k), x_e(k), x_f(k)^*, x_g(k)^*\} \quad (2)$$

It is then possible to store every entry of tensor C_x in a matrix Q_x , sometimes called the *quadricovariance*:

$$\forall 1 \leq d, e, f, g \leq N, Q_x(N(d-1)+g, N(f-1)+e) = C_{de,x}^{fg} \quad (3)$$

where $Q_x(r, q)$ corresponds to the (r, q) -th component of Q_x . Note that this arrangement is of course not unique, but this is the one that yields the best performances in terms of identifiability [6]. Matrix Q_x can be denoted more generally as $C_{4,x}^1$ [1].

Let us add that the expression of 4-th order cumulants of zero-mean complex variables as a function of moments appears in many papers; see for instance [10] for non zero mean variables. We also assume the following:

- H1.** For every index k , the source vector $s(k)$ is stationary and ergodic (extension to cyclostationary and cycloergodic is straightforward), with complex values, mutually independent at order 4 (i.e. all cross cumulants of order 4 are null);
- H2.** For every k , the noise vector $\nu(k)$ is stationary and ergodic;
- H3.** For every k , $s(k)$ and $\nu(k)$ are statistically independent;
- H4.** Source *kurtoses* (standardized autocumulants of order 4), $\kappa_{ppp,s}^{ppp} = C_{ppp,s}^{ppp} / \pi_p^2$, are non zero and have all the same sign ($\pi_p \stackrel{\text{def}}{=} E[|s_p(k)|^2]$ denotes the power of the p -th source);
- H5.** The number of sources is smaller than the number of sensors: $P \leq N$. The mixture is then referred to as *overdetermined*;
- H6.** The mixing matrix A is full rank and does not contain any null entry.

The goal is to determine a separating matrix, W , such that

$$y(n) \stackrel{\text{def}}{=} W^H x(n) \quad (4)$$

is an estimate of vector $s(k)$ up to a multiplicative *trivial* matrix (i.e. of the form $\Lambda\Pi$ where Λ is invertible diagonal and Π is a permutation). In ICAR, as in numerous other BSS algorithms, the construction of W requires the identification of mixture A .

3. THE CORE OF THE ICAR METHOD

3.1. Properties of the quadricovariance

The multilinearity property enjoyed by cumulants allows to write

$$Q_x = [A \otimes A^*] Q_s [A \otimes A^*]^H \quad (5)$$

where Q_x and Q_s denote quadricovariance matrices of $x(k)$ and $s(k)$, of size $N^2 \times N^2$ and $P^2 \times P^2$, respectively. Because sources are independent, Q_s is diagonal. However, it is not full rank. Another decomposition,

$$Q_x = A_Q Q_s A_Q^H \quad (6)$$

exhibits two new matrices. The first one, $Q_s \stackrel{\text{def}}{=} \text{diag}([C_{11,s}^{11}, C_{22,s}^{22}, \dots, C_{PP,s}^{PP}])$, is diagonal of size $P \times P$, and invertible from **(H4)**. The second one, A_Q , of size $N^2 \times P$, is also of full rank from **(H6)** and [1, prop. 1], can be written as:

$$\begin{aligned} A_Q &\stackrel{\text{def}}{=} [a_1 \otimes a_1^* \ a_2 \otimes a_2^* \ \dots \ a_P \otimes a_P^*] \\ &= [[A^* \Phi_1]^T \ [A^* \Phi_2]^T \ \dots \ [A^* \Phi_N]^T]^T \end{aligned} \quad (7)$$

where

$$\Phi_n \stackrel{\text{def}}{=} \text{diag}([A(n, 1) \ A(n, 2) \ \dots \ A(n, P)]) \quad (8)$$

Entries $A(n, \cdot)$ of the diagonal matrix Φ_n , of size $P \times P$, form the n -th line of A . (Matrices Q_s and A_Q are referred to as $\zeta_{4,s}$ et $\mathcal{A}_2^1$ in [1]).

3.2. Principle of ICAR

the principle is similar to that described in [2] with the difference that statistics of order 4 are utilized instead of 6. This difference is important because 4-th order statistics do not allow any more to address *underdetermined mixtures*, at least in the present framework. Since sources have the same kurtosis sign, it is possible to determine a unique real square root matrix of Q_x or $-Q_x$. Without restricting the generality, assume source kurtoses are all positive. The square root of Q_x can be computed by an Eigen Value Decomposition (EVD), from (6):

$$Q_x^{1/2} = E_s L_s^{1/2} = A_Q Q_s^{1/2} V^H \quad (9)$$

where L_s ($L_s^{1/2}$ denotes a square root of L_s) is the $P \times P$ diagonal matrix containing the P non-zero eigenvalues of Q_x , and E_s is the $N^2 \times P$ matrix of associated normalized eigenvectors. Because A_Q is full rank, it can be shown that **(H4)** amounts to assuming that diagonal elements of L_s are non zero and have the same sign [1, prop. 2], here positive. Moreover, (9) establishes a link between $Q_x^{1/2}$ and A_Q , where V is a unitary matrix, uniquely defined once L_s and E_s are fixed. Next, (9) and (7) relate $Q_x^{1/2}$ and A such that:

$$\begin{aligned} Q_x^{1/2} &= [[A^* \Phi_1 Q_s^{1/2} V^H]^T \ \dots \ [A^* \Phi_N Q_s^{1/2} V^H]^T]^T \\ &\stackrel{\text{def}}{=} [\Gamma_1^T \ \Gamma_2^T \ \dots \ \Gamma_N^T]^T \end{aligned} \quad (10)$$

where $\Gamma_n = \mathbf{A}^* \Phi_n \mathbf{Q}_s^{1/2} \mathbf{V}^H$ is the n -th matrix block of $\mathbf{Q}_x^{1/2}$, of size $N \times P$. Consequently, matrix $\mathbf{V}$ diagonalizes the $N(N-1)$ matrices Θ_{n_1, n_2} given by

$$\forall 1 \leq n_1 \neq n_2 \leq N, \quad \Theta_{n_1, n_2} = \Gamma_{n_1}^\# \Gamma_{n_2} \quad (11)$$

where $\#$ denotes pseudo-inversion. In fact, from the expression of Γ_n yielded by (10), and under assumptions (H5)-(H6), the pseudo-inverse of Γ_n can be written as $\Gamma_n^\# \stackrel{\text{def}}{=} (\Gamma_n^H \Gamma_n)^{-1} \Gamma_n^H$. As a consequence, matrices Θ_{n_1, n_2} can be rewritten in the form below:

$$\forall 1 \leq n_1 \neq n_2 \leq N, \quad \Theta_{n_1, n_2} = \mathbf{V} D_{n_1, n_2} \mathbf{V}^H \quad (12)$$

where matrices $D_{n_1, n_2} \stackrel{\text{def}}{=} \Phi_{n_1}^{-1} \Phi_{n_2}$ are diagonal. Denote $\mathbf{V}_{sol}$ the unitary matrix that jointly diagonalizes matrices Θ_{n_1, n_2} . Then $\mathbf{V}_{sol} = \mathbf{V} \mathbf{T}$ with $\mathbf{T}$ unitary. Thus, from (9), $\mathbf{V}_{sol}$ allows to identify $\mathcal{A}_Q$ up to a unitary matrix:

$$\mathbf{Q}_x^{1/2} \mathbf{V}_{sol} = \mathcal{A}_Q \mathbf{Q}_s^{1/2} \mathbf{T} \stackrel{\text{def}}{=} \widehat{\mathbf{A}}_Q \quad (13)$$

On the other hand, from assumptions (H5)-(H6), for every pair $(p_1, p_2)_{p_1 \neq p_2}$ belonging to $\{1, 2, \dots, P\}^2$, there exists at least a pair $(n_1, n_2)_{n_1 \neq n_2}$ belonging to $\{1, 2, \dots, N\}^2$ such that $D_{n_1, n_2}(p_1, p_1) \neq D_{n_1, n_2}(p_2, p_2)$. This implies that $\mathbf{T}$ is trivial. Then from (13), matrix $\widehat{\mathbf{A}}_Q$ is consequently an estimate of $\mathcal{A}_Q$, up to a trivial matrix.

Next from (10), equation (13) can be rewritten as

$$\begin{aligned} \mathbf{Q}_x^{1/2} \mathbf{V}_{sol} &= \left[[\mathbf{A}^* \Phi_1 \mathbf{Q}_s^{1/2} \mathbf{T}]^\top \dots [\mathbf{A}^* \Phi_N \mathbf{Q}_s^{1/2} \mathbf{T}]^\top \right]^\top \\ &\stackrel{\text{def}}{=} [\Sigma_1^\top \Sigma_2^\top \dots \Sigma_N^\top]^\top = \widehat{\mathbf{A}}_Q \end{aligned} \quad (14)$$

The matrix block Σ_1 , formed of the N first rows of $\mathbf{Q}_x^{1/2} \mathbf{V}_{sol}$, corresponds to $\mathbf{A}^*$, up to a trivial matrix:

$$\Sigma_1 = \mathbf{A}^* \Phi_1 \mathbf{Q}_s^{1/2} \mathbf{T} \quad (15)$$

where $\mathbf{Q}_s^{1/2}$ and Φ_n are diagonal for every n , $1 \leq n \leq N$. This method is named ICAR1. One could also think of several improvements, for instance by averaging the N blocks Σ_n for estimating $\mathbf{A}^*$, giving rise to ICAR2.

3.3. Refinement of the method

In order to fully exploit the information contained in matrix $\widehat{\mathbf{A}}_Q$ (14), namely its redundancies of $\mathcal{A}_Q$ in (7), it is possible to mimic the last step of FOBUM [9]. Indeed, from (13) and (7):

$$\begin{aligned} \widehat{\mathbf{A}}_Q &= \left[\lambda_{\xi(1)} [\mathbf{a}_{\xi(1)} \otimes \mathbf{a}_{\xi(1)}^*] \dots \lambda_{\xi(P)} [\mathbf{a}_{\xi(P)} \otimes \mathbf{a}_{\xi(P)}^*] \right] \\ &\stackrel{\text{def}}{=} [\mathbf{b}_{\xi(1)} \dots \mathbf{b}_{\xi(P)}] \end{aligned} \quad (16)$$

where $|\lambda_p| = |C_{ppp, s}^{ppp}|^{1/2}$, and ξ is a permutation mapping on $\{1, 2, \dots, P\}$. It is then possible to associate every vector

$\mathbf{b}_{\xi(p)}$ of size $N \times 1$ with a matrix $\mathbf{B}_{\xi(p)}$ of size $N \times N$, whose columns are precisely the N successive vectors $\mathbf{b}_{\xi(p)}$ formed of N values:

$$\mathbf{B}_{\xi(p)} = \lambda_{\xi(p)} \left[\mathbf{a}_{\xi(p)} \mathbf{a}_{\xi(p)}^H \right]^* \quad (17)$$

A mere diagonalization of matrices $\mathbf{B}_{\xi(p)}^*$ allows to yield the P directional vectors $\mathbf{a}_p$, by retaining each time the eigenvector associated with the dominant eigenvalue, and up to a permutation and a scalar multiplicative factor. The method with the latter improvement is named ICAR3.

4. SIMULATIONS

Two computer experiments show the performances of ICAR and some efficient BSS techniques (COM1, JADE, FastICA, SOBI). $P = 4$ statistically independent sources, i.e. 2 BPSK and 2 QPSK, all with a raised cosine pulse shape of roll-off equal to 0.25, are received by a UCA of $N = 4$ identical sensors of radius R such that $R/\lambda = 0.55$ (λ : wavelength). The four sources, assumed synchronized, have the same input SNR (Signal to Noise Ratio) of 20dB and the noise is spatially and temporally white Gaussian. The symbol period T_1 associated with the first BPSK is equal to three times the sample period T_e . The other sources have a symbol period equal to twice the sample period. The directions of arrival of the sources are such that the source steering vectors are orthogonal and the associated carrier residus are such that $f_{c1} T_e = 0$, $f_{c2} T_e = 0.3$, $f_{c3} T_e = 0.2$ and $f_{c4} T_e = 0.1$. The performance measure assumed to evaluate the quality of the extraction of source p is the maximal signal to interference plus noise ratio associated with source p , denoted $SINRM_p$ [5]. It can be compared to the optimal $SINRM_p$ computed using the exact mixing matrix instead of the estimated one and denoted by *Optimum SMF*. These are precisely the comparisons that are drawn now. Figure 1

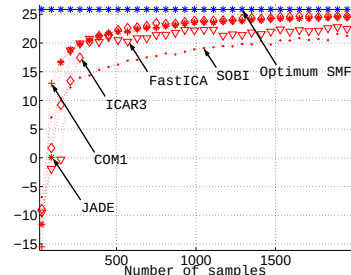


Fig. 1. $SINRM_3$ for a SNR of 20 dB

displays the $SINRM$ of source 3 averaged over 200 realizations, associated with ICAR3, and BSS techniques JADE, COM1, FastICA and SOBI, as a function of the number of samples. Figure 1 shows the good performance of the ICAR3 algorithm facing the other methods. Note that the SOBI method gives in this simulation good results since sources have been chosen with different spectral densities,

especially taking different carrier residus. Similar results are observed for the other sources, and are not reported here.

The influence of the noise spatial coherence coefficient ρ is next studied. Now we have 3 sources, i.e. 2 BPSK and 1 QPSK, all with a raised cosine pulse shape of roll-off equal to 0.25, are assumed to be received by a UCA of $N = 5$ identical sensors of radius R such that $R/\lambda = 0.55$. Their symbol periods are equal to $T_1 = 2T_e$, $T_2 = 3T_e$ and $T_3 = 4T_e$ respectively. Their carrier residus are chosen equal to zero. The source steering vectors are built orthogonal, the SNR is 0 dB and 1500 samples are used for separation in this scenario. We apply the COM1, COM2, JADE, SOBI and ICAR3 methods, and the SINRM associated with each source is computed and averaged still over 200 realizations. The Gaussian noise is modeled as a sum of two noises, $\nu_{in}(k)$ and $\nu_{out}(k)$, of covariance $\mathbf{R}_\nu^{in}$ and $\mathbf{R}_\nu^{out}$ respectively, such that:

$$\mathbf{R}_\nu^{in}(r, q) \stackrel{\text{def}}{=} \sigma^2 \delta(r - q) / 2 \quad \mathbf{R}_\nu^{out}(r, q) \stackrel{\text{def}}{=} \sigma^2 \rho^{|r - q|} / 2 \quad (18)$$

where σ^2 is the global noise variance per sensor. Note that $\mathbf{R}_\nu(r, q) \stackrel{\text{def}}{=} \mathbf{R}_\nu^{in}(r, q) + \mathbf{R}_\nu^{out}(r, q)$ is the (r, q) -th component of the global noise covariance. Contrary to COM1, COM2, JADE and SOBI, algorithm ICAR3 is totally insensitive to the increase in coefficient ρ . In fact, the classical methods such as COM1, COM2, JADE, FastICA and SOBI require a prior spatial whitening based on second order moments. This stage theoretically needs the perfect knowledge of the noise covariance. If this is not the case, a whitening of the observed data is performed instead, which is biased. ICAR3 does not suffer from this drawback, since it uses only 4-th order cumulants, which are (asymptotically) insensitive to Gaussian noise, regardless of its space/time color. Again, similar results have been observed for sources 1 and 2.

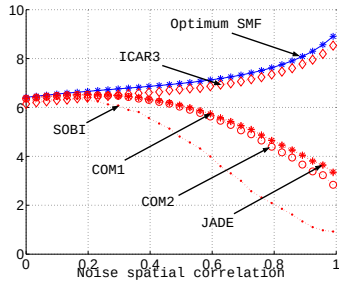


Fig. 2. SINRM_3 for a SNR of 0 dB

5. CONCLUSION

The new algorithm proposed in this paper, referred to as ICAR, utilizes only fourth order statistics of observations, and seems to be much more attractive than previous (classical) BSS techniques needing prior second order decorrelation, according to our computer simulations. We currently work on another solution to the problem presented in section 3.3, in other words, on another technique to extract the mixing matrix $\mathbf{A}$ contained in $\mathcal{A}_Q(7)$.

6. REFERENCES

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