Blind Signal Separation: Statistical Principles

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Invited Paper

Blind signal separation (BSS) and independent component analysis (ICA) are emerging techniques of array processing and data analysis that aim to recover unobserved signals or "sources" from observed mixtures (typically, the output of an array of sensors), exploiting only the assumption of mutual independence between the signals. The weakness of the assumptions makes it a powerful approach, but it requires us to venture beyond familiar second-order statistics. The objectives of this paper are to review some of the approaches that have been recently developed to address this exciting problem, to illustrate how they stem from basic principles, and to show how they relate to each other.

Keywords—Adaptive arrays, adaptive signal processing, array signal processing, asymptotic stability, blind source separation, higher order statistics, independent component analysis, maximum likelihood estimation, minimum entropy methods, signal analysis, signal processing antennas, signal representations, signal restoration, signal separation, source separation, unsupervised learning.

I. INTRODUCTION

Blind signal separation (BSS) consists of recovering unobserved signals or "sources" from several observed mixtures. Typically the observations are obtained at the output of a set of sensors, where each sensor receives a different combination of the source signals. The adjective "blind" stresses the fact that 1) the source signals are not observed and 2) no information is available about the mixture. This is a sound approach when modeling the transfer from the sources to the sensors is too difficult: it is unavoidable when no prior information is available about the transfer. The lack of prior knowledge about the mixture is compensated by a statistically strong but often physically plausible assumption of independence between the source signals. The so-called blindness should not be understood negatively; the weakness of the prior information is precisely the strength of the BSS model, making it a versatile tool for exploiting the spatial diversity provided by an array of sensors. Promising applications can already be found in the processing of communications signals, e.g., [23], [63],

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[67], [5], in biomedical signals¹ like ECG [30] and EEG [50], [46], in monitoring [37], [35], or as an alternative to principal component analysis (e.g., see [46], [9], [52], and [6]).

The simplest BSS model assumes the existence of n independent signals $s_1(t), \dots, s_n(t)$ and the observation of as many mixtures $x_1(t), \dots, x_n(t)$, these mixtures being linear and instantaneous, i.e., $x_i(t) = \sum_{j=1}^n a_{ij} s_j(t)$ for each i = 1, n.

This is represented compactly by the mixing equation

$$\mathbf{x}(t) = A\mathbf{s}(t) \tag{1}$$

where $\mathbf{s}(t) = [s_1(t), \cdots, s_n(t)]^{\dagger}$ is an $n \times 1$ column vector collecting the source signals, vector $\mathbf{x}(t)$ similarly collects the n observed signals, and the square $n \times n$ "mixing matrix" A contains the mixture coefficients. Here, as in the following, † denotes transposition. The BSS problem exists in recovering the source vector $\mathbf{s}(t)$ using only the observed data $\mathbf{x}(t)$, the assumption of independence between the entries of the input vector $\mathbf{s}(t)$, and possibly some prior information about the probability distribution of the inputs. It can be formulated as the computation of an $n \times n$ separating matrix B whose output

$$\mathbf{y}(t) = B\mathbf{x}(t) \tag{2}$$

is an estimate of the vector s(t) of the source signals.

Fig. 2 shows an example of adaptive separation of (real) digital communications signals. A two-sensor array collects complex-valued noisy mixtures of two source signals which both have a constant modulus envelope. Successful separation upon adaptation is evidenced by the restoration of the constant modulus at each output.

In Fig. 2, the underlying BSS algorithm optimizes a cost function composed of two penalty terms: one for correlation between outputs and one for deviation of the modulus from a constant value. This example introduces several points to be developed below.

 A penalty term involving only pairwise decorrelation (second-order statistics) would not lead to separa-

¹ See http://www.cnl.salk.edu/~tewon/ica_cnl.html, the ICA page of the CNL group, for several biomedical applications.

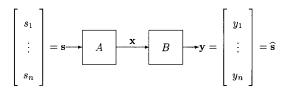


Fig. 1. Mixing and separating. Unobserved signals s; observations x, estimated source signals y.

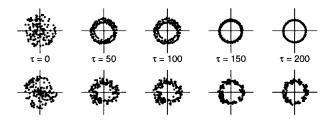


Fig. 2. Outputs $y_1(t)$ (top row) and $y_2(t)$ (bottom row) when using the separating matrix obtained after adaptation based on 0, 50, 100, 150, 200 samples of a 2×2 mixture of constant modulus signals. Each subplot is in the complex plane. The clustering around circles shows the restoration of the constant modulus property.

- tion. Source separation must go beyond second-order statistics (see Section II).
- Source separation can be obtained by optimizing a contrast function, i.e., a scalar measure of some distributional property of the output y. The constant modulus property is very specific; more general contrast functions are based on other measures, such as entropy, mutual independence, high-order decorrelations, divergence between the joint distribution of y and some model, etc. Contrast functions are discussed in Section III, where it is seen how they relate to each other and can be derived from the maximum likelihood principle.
- Fast adaptation is possible, even with simple algorithms (see Sections IV and V), and blind identification can be accurate even with a small number of samples (see Section VI on performance analysis).

The basic BSS model can be extended in several directions. Considering, for instance, more sensors than sources, noisy observations, and complex signals and mixtures, one obtains the standard narrow band array processing/beamforming model. Another extension is to consider convolutive mixtures: this results in a multichannel blind deconvolution problem. These extensions are of practical importance, but this paper is restricted to the simplest model (i.e., real signals, as many sensors as sources, nonconvolutive mixtures, and noise-free observations) because it captures the essence of the BSS problem and because our objective is to present the basic statistical ideas, focusing on principles. Nonetheless, some references to papers addressing more general models are provided in the last section.

The paper is organized as follows: Section II discusses blind identifiability; Sections III and IV present contrast functions and estimating functions, starting from information—theoretic ideas and moving to suboptimal high-

order approximations; adaptive algorithms are described in Section V; Section VI addresses some performance issues.

II. CAN IT BE DONE? MODELING AND IDENTIFIABILITY

When is source separation possible? To what extent can the source signals be recovered? What are the properties of the source signals allowing for partial or complete blind recovery? These issues are addressed in this section.

A. The BSS Model

Source separation exploits primarily spatial diversity, that is, the fact that different sensors receive different mixtures of the sources. Spectral diversity, if it exists, could also be exploited, but the approach of source separation is essentially spatial, meaning that it looks for structure across the sensors, not across time. The consequence of ignoring any time structure is that the information contained in the data is represented exhaustively by the sample distribution of the observed vector \boldsymbol{x} (as depicted graphically in Fig. 3). Thus, BSS becomes the problem of identifying the probability distribution of a vector $\boldsymbol{x} = A\boldsymbol{s}$, given a sample distribution. In this perspective, the statistical model has two components: the mixing matrix A and the probability distribution of the source vector \boldsymbol{s} .

— Mixing matrix. The mixing matrix A is the parameter of interest. Its columns are assumed to be linearly independent so that it is invertible (see [13] for the discussion of a more general case).

There is something special about having an invertible matrix as the unknown parameter, because matrices represent linear transformations. Indeed, model (1) is a particular instance of a transformation model. Furthermore, the set of all $n \times n$ invertible matrices forms a multiplicative group. This simple fact has a profound impact on source separation because it allows for the design of algorithms with uniform performance, i.e., whose behavior is completely independent of the particular mixture (Sections V-A and VI-C).

— Source distribution. The probability distribution of each source is a "nuisance parameter," meaning that we are not primarily interested in it even though knowing or estimating these distributions is necessary to estimate efficiently the parameter of interest. Even if we say nothing about the distribution of each source, we say a lot about their joint distribution by the key assumption of mutual source independence. If each source i=1,n is assumed to have a probability density function (pdf) denoted $q_i(\cdot)$, the independence assumption has a simple mathematical expression. The (joint) pdf q(s) of the source vector s is

$$q(\mathbf{s}) = q_1(s_1) \times \dots \times q_n(s_n) = \prod_{i=1,n} q_i(s_i)$$
 (3)

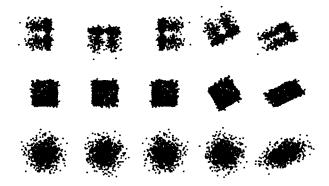


Fig. 3. Sample distributions of (x_1, x_2) when $\mathbf{x} = A\mathbf{s}$ for five different transformation matrices A and three pairs of distributions for (s_1, s_2) . From left to right: the identity transform, permutation of the sources, sign change, a $\pi/3$ rotation, and a generic linear transform.

i.e., it is the product of the densities for all sources (the marginal densities). Source separation techniques differ widely by the (explicit or implicit) assumptions made on the individual distributions of the sources. There is a whole range of options.

- 1) The source distributions are known in advance.
- 2) Some features are known (moments, heavy tails, bounded support, etc.).
- 3) They belong to a parametric family.
- 4) No distribution model is available.

A priori, the stronger the assumption, the narrower the applicability. However, well-designed approaches are in fact surprisingly robust even to gross errors in modeling the source distributions, as shown below. For ease of exposition, zero mean sources are assumed throughout

$$Es = 0$$
, i.e., $Es_i = 0$ $1 \le i \le n$. (4)

B. Blind Identifiability

The issue of blind identifiability is to understand to which extent matrix A is determined from the sole distribution of the observed vector $\mathbf{x} = A\mathbf{s}$. The answer depends on the distribution of \mathbf{s} and on what is known about it.

A square matrix is said to be nonmixing if it has one and only one nonzero entry in each row and each column. If C is nonmixing then $\mathbf{y} = C\mathbf{s}$ is a copy of \mathbf{s} , i.e., its entries are identical to those of \mathbf{s} up to permutations and changes of scales and signs. Source separation is achieved if such a copy is obtained. When the distribution of \mathbf{s} is unknown, one cannot expect to do any better than signal copy, but the situation is a bit different if some prior information about the distribution of \mathbf{s} is available. If the sources have distinct distributions, a possible permutation can be detected; if the scale of a given source is known, the amplitude of the corresponding column of A can be estimated, etc.

Some intuition about identifiability can be gained by considering simple examples of 2×2 mixing. Each row of Fig. 3 shows (sample) distributions of a pair (s_1, s_2) of independent variables after various linear transforms. The

columns successively show (s_1, s_2) , (s_2, s_1) , $(-s_1, -s_2)$, and the effect of a $\pi/3$ rotation and of a nondescript linear transform.

Visual inspection of the transformed distribution compared to the original one gives a feeling of how well the transform matrix A can be identified based on the observation of a mixture. The first row of Fig. 3 shows a case where the second column of A can be identified only up to sign because s_2 is symmetrically distributed about the origin (and therefore has the same distribution as $-s_2$). The second row shows a more severe indetermination. Here, s_1 and s_2 have the same symmetric distribution, and the transform can be determined only up to arbitrary changes of sign and a permutation. The last row shows the most severe case. Here, s_1 and s_2 are normally distributed with equal variance so that their joint distribution is invariant under rotation.

These simple examples suggest that A can be blindly identified—possibly up to some indeterminations induced by the symmetries in the distribution of the source vector—in the case of known source distributions. However, this knowledge is not necessary; the eye certainly can capture the distortion in the last columns of Fig. 3, even without reference to the undistorted shapes in first column. This is because the graphical "signature of independence" (the pdf shape in the first column) clearly appears as distorted in the last colum. This intuition is supported by the following statement adapted from [25] after a theorem of Darmois (see also [13]).

For a vector \mathbf{s} of independent entries with at most one Gaussian entry and for any invertible matrix C, if the entries of $\mathbf{y} = C\mathbf{s}$ are independent, then \mathbf{y} is a copy of \mathbf{s} (C is nonmixing).

Thus, unless a linear transform is nonmixing, it turns a vector of independent entries (at most one being Gaussian) into a vector whose entries are not independent. This is a key result because it means that blind signal separation can be achieved by restoring statistical independence. This is not only a theoretical result about blind identifiability, but it also suggests that BSS algorithms could be devised by maximizing the independence between the outputs of a separating matrix. Section III shows that the maximum likelihood principle does support this idea and leads to a specific measure of independence.

1) Independence and Decorrelation: Blind separation can be based on independence but independence cannot be reduced to the simple decorrelation conditions that $Ey_iy_j = 0$ for all pairs $1 \le i \ne j \le n$. This is readily seen from the fact that there are, by symmetry, only n(n-1)/2 such conditions (one for each pair of sources) while there are n^2 unknown parameters.

Second-order information (decorrelation), however, can be used to reduce the BSS problem to a simpler form. Assume, for simplicity, that the source signals have unit variance so that their covariance matrix is the identity matrix $\mathbf{E} \boldsymbol{s} \boldsymbol{s}^\dagger = I$; vector \boldsymbol{s} is said to be spatially white. Let W be a whitening matrix for \boldsymbol{x} , that is, $\boldsymbol{z} \stackrel{\triangle}{=} W \boldsymbol{x}$ is spatially white.

The composite transform WA is necessarily a rotation matrix because it relates two spatially white vectors \mathbf{s} and $\mathbf{z} = WA\mathbf{s}$. Therefore, whitening or sphering the data reduces the mixture to a rotation matrix. It means that a separating matrix can be found as a product B = UW, where W is a whitening matrix and U is a rotation matrix. Note that any further rotation of \mathbf{z} into $\mathbf{y} = U\mathbf{z}$ preserves spatial whiteness, so that two equivalent approaches to exploiting source decorrelation are 1) find B as B = UW with W a spatial whitener and U a rotation or 2) find B under the whiteness constraint $E\mathbf{y}\mathbf{y}^{\dagger} = I$. For further reference, we write the whiteness constraint as

$$EH_{w}(\mathbf{y}) = 0$$
 where $H_{w}(\mathbf{y}) \stackrel{\Delta}{=} \mathbf{y}\mathbf{y}^{\dagger} - I$. (5)

Spatial whiteness imposes n(n+1)/2 constraints, leaving n(n-1)/2 unknown (rotation) parameters to be determined by other than second-order information. Second-order information is able to do "about half of the BSS job."

The prewhitening approach is sensible from an algorithmic point of view but it is not necessarily statistically efficient (see Section VI-B). Actually, enforcing the whiteness constraint amounts to believing that second-order statistics are "infinitely more reliable" than any other kinds of statistics. This is, of course, untrue.

C. Likelihood

This section examines the likelihood of source separation models in a simple, graphical way. The likelihood in a given model is the probability of a data set as a function of the parameters of the model. The simple model $\boldsymbol{x} = A\boldsymbol{s}$ for vector \boldsymbol{x} discussed in Section II-A is parameterized by the pair (A,q), which is made of the mixing matrix A and the density q for the source vector \boldsymbol{s} . The density of $\boldsymbol{x} = A\boldsymbol{s}$ for a given pair (A,q) is classically given by

$$p(\mathbf{x}; A, q) = |\det A|^{-1} q(A^{-1}\mathbf{x}).$$
 (6)

If T samples $X_{1:T} \stackrel{\triangle}{=} [\boldsymbol{x}(1), \cdots, \boldsymbol{x}(T)]$ of \boldsymbol{x} are modeled as independent, then $p(X_{1:T}) = p(\boldsymbol{x}(1)) \times \cdots \times p(\boldsymbol{x}(T))$. Thus, the normalized (i.e., divided by T) log likelihood of $X_{1:T}$ for the parameter pair (A,q) is

$$\frac{1}{T}\log p(X_{1:T}; A, q) = \frac{1}{T} \sum_{t=1}^{T} \log q(A^{-1}\boldsymbol{x}(t)) - \log|\det A|.$$
(7)

Figs. 5–7 show the likelihood landscape when A is varied while q is kept fixed. For each figure, T=1000 independent realizations of $\mathbf{s}=[s_1,s_2]^\dagger$ are drawn according to some pdf $r(\mathbf{s})=r_1(s_1)r_2(s_2)$ and are mixed with a 2×2 matrix \underline{A} to produce T samples of \mathbf{x} . Therefore, this data set $X_{1:T}$ follows exactly model (1) with a true mixing matrix \underline{A} and a true source distribution $r(\mathbf{s})$. The figures show the log likelihood when A is varied around its true value \underline{A} while model density $q(\mathbf{s})$ is kept fixed. These figures illustrate the impact of the choice of a particular model density.

In each of these figures, the matrix parameter A is varied in two directions in matrix space according to A =

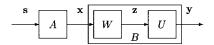


Fig. 4. Decorrelation leaves an unknown rotation.

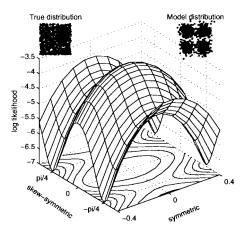


Fig. 5. Log likelihood with a slightly misspecified model for source distribution. Maximum is reached close to the true value.

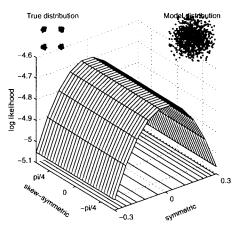


Fig. 6. Log likelihood with a Gaussian model for source distribution. There is no contrast in the skew-symmetric direction.

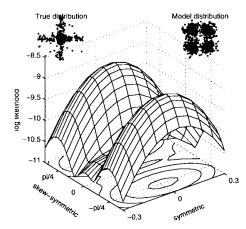


Fig. 7. Log likelihood with a widely misspecified model for source distribution; maximum is reached for a mixing system.

 $\underline{A}M(u,v)$, where M(u,v) is the matrix

$$M(u,v) \stackrel{\Delta}{=} \begin{bmatrix} \cosh u & \sinh u \\ \sinh u & \cosh u \end{bmatrix} \cdot \begin{bmatrix} \cos v & -\sin v \\ \sin v & \cos v \end{bmatrix}. \quad (8)$$

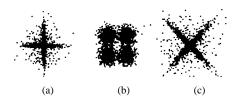


Fig. 8. How the maximum likelihood estimator is misled: (a) true distribution, (b) hypothesized distribution, and (c) estimated distribution.

This is just a convenient way to generate a neighborhood of the identity matrix. For small u and v

$$M(u,v) \approx I + u \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix} + v \begin{bmatrix} 0 & -1 \\ 1 & 0 \end{bmatrix}.$$
 (9)

Therefore, u and v are called symmetric and skew-symmetric parameters, respectively. Each one controls a particular deviation of M(u,v) away from the identity.

In Fig. 5, the true source distributions r_1 and r_2 are uniform on [-1, +1] but the model takes q_1 and q_2 to be each a mixture of two normal distributions with same variance but different means (as in the second column of Fig. 11). True and hypothesized sample distributions of $\mathbf{s} = (s_1, s_2)$ are displayed in the upper left and right corners of the plot. Even though an incorrect model is used for the source distribution $q \neq r$, the figure shows that the likelihood is maximal around (u, v) = (0, 0), i.e., the most likely mixing matrix given the data and the model is close to \underline{A} .

In Fig. 6, the true sources are almost binary (see upper left corner) but a Gaussian model is used: $q_1(s) = q_2(s) \propto \exp{-s^2/2}$. The figure shows that the likelihood of $A = \underline{A}M(u,v)$ does not depend on the skew-symmetric parameter v, again evidencing the insufficiency of Gaussian modeling.

In Fig. 7, the sources are modeled as in Fig. 5, but the true (and identical) source distributions r_1 and r_2 are now mixtures of normal distributions with the same mean but different variances (as in the second column of Fig. 11). A disaster happens: the likelihood is no longer maximum for A in the vicinity of A.

Actually, if the value \hat{A} of A maximizing the likelihood is used to estimate the source signals as $\mathbf{y} = \hat{s} = \hat{A}^{-1}\mathbf{x}$, one obtains maximally mixed sources. This is explained in Section III-A and Fig. 8.

The main point of this informal study is the necessity of non-Gaussian modeling (Fig. 6), the possibility of using only an approximate model of the sources (Fig. 5), and the existence of a limit to the misspecification of the source model (Fig. 7). How wrong can the source distribution model be? This is quantified in Section VI-A.

III. CONTRAST FUNCTIONS

This section introduces contrast functions, which are objective functions for source separation. The maximum likelihood principle is used as a starting point, suggesting several information—theoretic objective functions (Section III-A) which are then shown to be related to another class

of objective functions based on high-order correlations (Section III-B).

Minimum contrast estimation is a general technique of statistical inference [57] which encompasses several techniques like maximum likelihood or least squares. It is relevant for blind deconvolution (see [36] and [11]) and has been introduced in the related BSS problem by Comon in [25]. In both instances, a contrast function is a real function of a probability distribution. To deal with such functions, a special notation is useful: for \boldsymbol{x} , a given random variable, $f(\boldsymbol{x})$ generically denotes a function of \boldsymbol{x} , while $f[\boldsymbol{x}]$ denotes a function of the distribution of \boldsymbol{x} . For instance, the mean value of \boldsymbol{x} is denoted $m[\boldsymbol{x}] \stackrel{\triangle}{=} \mathbf{E}\boldsymbol{x}$.

Contrast functions for source separation (or simply "contrasts," for short) are generically denoted $\phi[y]$. They are real-valued functions of the distribution of the output y = Bx and they serve as objectives: they must be designed in such a way that source separation is achieved when they reach their minimum value. In other words, a valid contrast function should, for any matrix C, satisfy $\phi[Cs] \ge \phi[s]$ with equality only when y = Cs is a copy of the source signals. Since the mixture can be reduced to a rotation matrix by enforcing the whiteness constraint $Eyy^{\dagger} = I$ (Section II-B), one can also consider orthogonal contrast functions. These are denoted as $\phi^{\circ}[y]$ and must be minimized under the whiteness constraint $Eyy^{\dagger} = I$.

A. Information Theoretic Contrasts

The maximum likelihood (ML) principle leads to several contrasts which are expressed via the Kullback divergence. The Kullback divergence between two probability density functions f(s) and g(s) on \mathbb{R}^n is defined as

$$K(f|g) \stackrel{\Delta}{=} \int_{\mathbf{s}} f(\mathbf{s}) \log \left(\frac{f(\mathbf{s})}{g(\mathbf{s})} \right) d\mathbf{s}$$
 (10)

whenever the integral exists [27]. The divergence between the distributions of two random vectors \boldsymbol{w} and \boldsymbol{z} is concisely denoted $\boldsymbol{K}[\boldsymbol{w}|\boldsymbol{z}]$. An important property of \boldsymbol{K} is that $\boldsymbol{K}[\boldsymbol{w}|\boldsymbol{z}] \geq 0$ with equality if and only if \boldsymbol{w} and \boldsymbol{z} have the same distribution. Even though \boldsymbol{K} is not a distance (it is not symmetric), it should be understood as a statistical way of quantifying the closeness of two distributions.

1) Matching Distributions—Likelihood and Infomax: The likelihood landscapes displayed in Figs. 5–7 assumes a particular pdf $q(\cdot)$ for the source vector. Denoting \underline{s} a random vector with distribution q, simple calculus shows that

$$\frac{1}{T}\log p(X_{1:T}; A, q) \xrightarrow{T \to \infty} -\mathbf{K}[A^{-1}\mathbf{x}|\underline{\mathbf{s}}] + \text{cst.} \quad (11)$$

Therefore, Figs. 5–7 approximately display (up to a constant term) minus the Kullback divergence between the distribution of $\mathbf{y} = A^{-1}\mathbf{x}$ and the hypothesized distribution of the sources. This shows that the maximum likelihood principle is associated with a contrast function

$$\phi_{ML}[\mathbf{y}] = \mathbf{K}[\mathbf{y}|\underline{\mathbf{s}}] \tag{12}$$

and the normalized log likelihood can be seen, via (11), as an estimate of $-K[y|\underline{s}]$ (up to a constant). The ML principle thus says something very simple when applied to the BSS problem, that is, find matrix A such that the distribution of $A^{-1}x$ is as close as possible (in the Kullback divergence) to the hypothesized distribution of the sources.

The instability problem illustrated by Fig. 7 may now be understood as follows: in this figure, the likelihood is maximum when M(u,v) is a $\pm \pi/4$ rotation, because the true source distribution is closer to the hypothesized source distribution after it is rotated by $\pm \pi/4$. As Fig. 8 shows, after such a rotation the areas of highest density of \boldsymbol{y} correspond to the points of highest probability of the hypothesized source model.

A different approach to derive the contrast function (12) is very popular among the neural network community. Denote $g_i(\cdot)$ the distribution function

$$g_i(s) \stackrel{\Delta}{=} \int_{-\infty}^s q_i(t) dt \in [0,1] \qquad 1 \le i \le n$$
 (13)

so that $g'_i = q_i$, and denote $g(\mathbf{s}) = [g_1(s_1), \dots, g_n(s_n)]^{\dagger}$. An interpretation of the infomax principle (see [8], [14], and references therein) suggests the contrast function

$$\phi_{IM}[\boldsymbol{y}] \stackrel{\Delta}{=} -\boldsymbol{H}[g(\boldsymbol{y})] \tag{14}$$

where $\boldsymbol{H}[\cdot]$ denotes the Shannon entropy (for a random vector \boldsymbol{u} with density $p(\boldsymbol{u})$, this is $\boldsymbol{H}[\boldsymbol{u}] = -\int p(\boldsymbol{u}) \log p(\boldsymbol{u}) \ d\boldsymbol{u}$ with the convention $0 \log 0 = 0$). This idea can be understood as follows: on one hand, $g(\boldsymbol{s})$ is uniformly distributed on $[0,1]^n$ if \boldsymbol{s} has pdf q; on the other hand, the uniform distribution has the highest entropy among all distributions on $[0,1]^n$ [27]. Therefore, $g(C\underline{\boldsymbol{s}})$ has the highest entropy when C=I. The infomax idea, however, yields the same contrast as the likelihood $\phi_{IM}[\boldsymbol{y}] = \phi_{ML}[\boldsymbol{y}]$. The connection between maximum likelihood and infomax was noted by several authors (see [56], [18], [49]).

2) Matching the Structure—Mutual Information: The simple likelihood approach described above is based on a fixed hypothesis about the distribution of the sources. This becomes a problem if the hypothesized source distributions differ too much from the true ones, as illustrated by Figs. 7 and 8. This remark suggests that the observed data should be modeled by adjusting both the unknown system and the distributions of the sources. In other words, one should minimize the divergence $K[y|\underline{s}]$ with respect to A (via the distribution of \underline{s} . The last minimization problem has a simple and intuitive theoretical solution. Denote \tilde{y} a random vector with 1) independent entries and 2) each entry distributed as the corresponding entry of y. A classic property (e.g., see [27]) of \tilde{y} is that

$$K[y|\underline{s}] = K[y|\tilde{y}] + K[\tilde{y}|\underline{s}]$$
(15)

for any vector \underline{s} with independent entries. Since $K[y|\tilde{y}]$ does not depend on \underline{s} , (15) shows that $K[y|\underline{s}]$ is minimized in \underline{s} by minimizing its second term, i.e., $K[\tilde{y}|\underline{s}]$; this is

simply achieved by taking $\underline{s} = \tilde{y}$ for which $K[\tilde{y}|\underline{s}] = 0$, so that $\min_{\underline{s}} K[y|\underline{s}] = K[y|\tilde{y}]$. Having minimized the likelihood contrast $K[y|\underline{s}]$ with respect to the source distribution leading to $K[y|\tilde{y}]$, our program is completed if we minimize the latter with respect to y, i.e., if we minimize the contrast function

$$\phi_{MI}[\boldsymbol{y}] \stackrel{\Delta}{=} \boldsymbol{K}[\boldsymbol{y}|\tilde{\boldsymbol{y}}]. \tag{16}$$

The Kullback divergence $K[y|\tilde{y}]$ between a distribution and the closest distribution with independent entries is traditionally called the mutual information (between the entries of y). It satisfies $\phi_{MI}[y] \geq 0$ with equality if and only if y is distributed as \tilde{y} . By definition of \tilde{y} , this happens when the entries of y are independent. In other words, $\phi_{MI}[y]$ measures the independence between the entries of y. Thus, mutual information apears as the quantitative measure of independence associated to the maximum likelihood principle.

Note further that $K[\tilde{y}|\underline{s}] = \sum_{i=1}^{n} K[y_i|\underline{s}_i]$ (because both \tilde{y} and \underline{s} have independent entries). Therefore

$$\phi_{ML}[\mathbf{y}] = \phi_{MI}[\mathbf{y}] + \sum_{i=1}^{n} \mathbf{K}[y_i|\underline{s}_i]$$
 (17)

so that the decomposition (15) or (17) of the global distribution matching criterion $\phi_{ML}[\mathbf{y}] = \mathbf{K}[\mathbf{y}|\mathbf{s}]$ should be understood as

$$\binom{\text{Total}}{\text{mismatch}} = \binom{\text{Deviation from}}{\text{independence}} + \binom{\text{Marginal}}{\text{mismatch}}.$$

Therefore, maximizing the likelihood with fixed assumptions about the distributions of the sources amounts to minimizing a sum of two terms: the first term is the true objective (mutual information as a measure of independence) while the second term measures how far the (marginal) distributions of the outputs y_1, \dots, y_n are from the assumed distributions.

3) Orthogonal Contrasts: If the mixing matrix has been reduced to a rotation matrix by whitening as explained in Section II-B, contrast functions like ϕ_{ML} or ϕ_{MI} can still be used. The latter takes an interesting alternative form under the whiteness constraint $\mathbf{E} y y^\dagger = I$. In this case, one can show that $\phi_{MI}[y]$ is, up to a constant term, equal to the sum of the Shannon entropies of each output. Thus, under the whiteness constraint, minimizing the mutual information between the entries of y is equivalent to minimizing the sum of the entropies of the entries of y, and we define

$$\phi_{MI}^{\circ}[\boldsymbol{y}] \stackrel{\Delta}{=} \sum_{i} \boldsymbol{H}[y_{i}]. \tag{18}$$

There is a simple interpretation: mixing the entries of s tends to increase their entropies. It seems natural to find separated source signals as those with minimum marginal entropies. It is also interesting to notice that $-H[y_i]$ is (up to a constant) the Kullback divergence between the distribution of y_i and the zero-mean unit-variance normal distribution. Therefore, minimizing the sum of the marginal

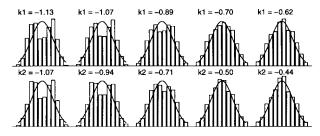


Fig. 9. Gaussianization by mixing. Histograms of y_1 (top row) and y_2 (bottom row) when \boldsymbol{y} is rotated by $\alpha\pi/4$ for $\alpha=0,1/4,1/2,3/4,1$. Each subplot also shows the estimated kurtosis k_1 and k_2 [defined at (21)] decreasing (in absolute value) upon mixing.

entropies is also equivalent to driving the marginal distributions of y as far away as possible from normality. Again, the interpretation is that mixing tends to Gaussianize the marginal distributions so that a separating technique should go in the opposite direction. Fig. 9 is a visual illustration of the tendency towards normality by mixing.

The first column shows histograms for two independent variables s_1 and s_2 with a bimodal distribution and (superimposed to it as a solid line) the best Gaussian approximation. The following columns show the histograms after rotations by steps of $\pi/16$, going from 0 to $\pi/4$ where mixing is maximal. The tendency towards normality is very apparent.

The entropic form (18) of the mutual information was used as starting point by Comon [25]; it remains a valid contrast under the weaker constraint that B is a volume preserving transformation [55].

4) Discussion: The canonical contrast for source separation is the mutual information ϕ_{MI} because it expresses the key property of source independence and nothing else. It does not include any explicit or implicit assumption about the distributions of the sources. On the other hand, if the source distributions are known, ϕ_{ML} is more appropriate because it directly expresses the fit between data and model. Also, ϕ_{ML} is easier to minimize because its gradient is easily estimated [see (31)] while estimating the gradient of ϕ_{MI} is computationally demanding [59]. Even when the source distributions are unknown, one may use ϕ_{ML} with hypothesized source distributions which only need to be close enough to the true distributions (Section II-C provides a qualitative explanation, Section VI-A provides a quantitative statement, and Section V-B discusses adapting the model distributions). Another approach is to approximate the Kullback-based contrasts using high-order statistics, as examined next.

B. High-Order Approximations

High-order statistics can be used to define contrast functions which are simple approximations to those derived from the ML approach. High-order information is most simply expressed by using cumulants. The discussion being limited to cumulants of orders two and four, only the following definitions are needed. For zero-mean random variables a, b, c, d, second-order cumulants are identical to

second-order moments $\operatorname{Cum}[a,b] \stackrel{\Delta}{=} \operatorname{E}\!ab$ and fourth-order cumulants are

$$\operatorname{Cum}[a, b, c, d] \stackrel{\Delta}{=} \operatorname{E}abcd - \operatorname{E}abEcd - \operatorname{E}acEbd - \operatorname{E}adEbc.$$
(19)

Whenever the random variables a,b,c,d can be split into two mutually independent groups, their cumulant is zero. Therefore, independence beyond second-order decorrelation can be tested easily using high-order cumulants.

For simplicity, the following notation for the cumulants of the elements of a given vector y is used throughout:

$$C_{ij}[\mathbf{y}] \stackrel{\triangle}{=} \operatorname{Cum}[y_i, y_j] \quad C_{ijkl}[\mathbf{y}] \stackrel{\triangle}{=} \operatorname{Cum}[y_i, y_j, y_k, y_l].$$

Since the source vector s has independent entries, all its cross-cumulants vanish

$$C_{ij}[\mathbf{s}] = \sigma_i^2 \delta_{ij} \quad C_{ijkl}[\mathbf{s}] = k_i \delta_{ijkl}$$
 (20)

where δ is the Kronecker symbol, and where we have defined the variance σ_i^2 and the kurtosis k_i of the *i*th source as the second- and fourth-order autocumulants of s_i

$$\sigma_i^2 \stackrel{\Delta}{=} \mathcal{C}_{ii}[\mathbf{s}] = \mathrm{E}s_i^2 \quad k_i \stackrel{\Delta}{=} \mathcal{C}_{iiii}[\mathbf{s}] = \mathrm{E}s_i^4 - 3\mathrm{E}^2s_i^2.$$
 (21)

The likelihood contrast $\phi_{ML}[y] = K[y|\underline{s}]$ is the measure of mismatch between output distribution and a model source distribution. A cruder measure can be defined from the quadratic mismatch between the cumulants

$$\phi_{2}[\mathbf{y}] \stackrel{\triangle}{=} \sum_{ij} (C_{ij}[\mathbf{y}] - C_{ij}[\mathbf{s}])^{2} = \sum_{ij} (C_{ij}[\mathbf{y}] - \sigma_{i}^{2} \delta_{ij})^{2}$$

$$\phi_{4}[\mathbf{y}] \stackrel{\triangle}{=} \sum_{ijkl} (C_{ijkl}[\mathbf{y}] - C_{ijkl}[\mathbf{s}])^{2}$$

$$= \sum_{ijkl} (C_{ijkl}[\mathbf{y}] - k_{i} \delta_{ijkl})^{2}.$$

Are ϕ_2 and ϕ_4 contrast functions, as introduced in the beginning of this section? Clearly ϕ_2 is not a contrast because $\phi_2[\boldsymbol{y}] = 0$ only expresses the decorrelation between the entries of \boldsymbol{y} . On the contrary, one can show that $\phi_4[\boldsymbol{y}]$ is a contrast if all the sources have known nonzero kurtosis. Even though fourth order information is by itself sufficient to solve the BSS problem, it is interesting to use ϕ_2 and ϕ_4 in conjunction because they jointly provide an approximation to the likelihood contrast; if \boldsymbol{s} and \boldsymbol{y} are symmetrically distributed with distributions that are close to normal, then

$$K[y|s] \approx \phi_{24}[y] \stackrel{\Delta}{=} \frac{1}{48} (12\phi_2[y] + \phi_4[y]).$$
 (22)

Room is lacking to discuss the validity of this approximation (which stems from an Edgeworth expansion, see Section V-B). The point, however, is not to determine how closely $\phi_{24}[\boldsymbol{y}]$ approximates $\boldsymbol{K}[\boldsymbol{y}|\boldsymbol{s}]$, but rather to follow the suggestion that second- and fourth-order information could be used jointly.

1) Orthogonal Contrasts: We consider cumulant-based orthogonal contrasts. The orthogonal approach, which enforces whiteness, i.e., $\phi_2[\boldsymbol{y}] = 0$, corresponds to replacing the factor 12 in (22) by an infinite weight (optimal weighting is considered in [19]; see also Section V-B) or is equivalent to minimizing $\phi_4[\boldsymbol{y}]$ under the whiteness constraint $\phi_2[\boldsymbol{y}] = 0$. Simple algebra shows that if $\phi_2[\boldsymbol{y}] = 0$, then $\phi_4[\boldsymbol{y}]$ is equal (up to a constant additive term) to

$$\phi_4^{\circ}[\boldsymbol{y}] \stackrel{\Delta}{=} -2 \sum_{i=1}^n k_i \mathcal{C}_{iiii}[\boldsymbol{y}] = \mathbf{E} f_4(\boldsymbol{y})$$
 (23)

where we have defined $f_4(\mathbf{y}) \stackrel{\triangle}{=} -2 \sum_{i=1}^n k_i (y_i^4 - 3)$. Since this contrast function is the expectation of a function of \mathbf{y} , it is particularly simple to estimate by a sample average.

Recall that the contrast function ϕ_{ML} defined in (12) depends on a source model, i.e., it is defined using a hypothetical density $q(\cdot)$ for the source distribution. Similarly, the fourth-order approximation $\phi_4^{\rm o}$ requires a hypothesis about the sources, but it is only a fourth-order hypothesis in the sense that only the kurtosis k_i for each source must be specified in definition (23). In the same manner that minimizing ϕ_{ML} over the source distribution yields the mutual information contrast ϕ_{MI} , minimizing $\phi_4^{\rm o}[\boldsymbol{y}]$ (which approximates ϕ_{ML}) over the kurtosis k_i of each source yields an approximation to the mutual information which is

$$\phi_{\text{ICA}}^{\circ}[\boldsymbol{y}] = \sum_{ijkl \neq iiii} C_{ijkl}^{2}[\boldsymbol{y}] = -\sum_{i} C_{iiii}^{2}[\boldsymbol{y}] + \text{cst.} \quad (24)$$

This was first obtained by Comon [26] (along a slightly different route) and by Gaeta *et al.* [40] by approximating the likelihood by a Gram–Charlier expansion. This contrast is similar to ϕ_{MI} also in that its first form involves only terms measuring the (fourth-order) independence between the entries of \boldsymbol{y} . Its second form stems from the fact that $\Sigma_{ijkl} \, \mathcal{C}^2_{ijkl}[\boldsymbol{y}]$ is constant if $\phi_2[\boldsymbol{y}] = 0$ holds (e.g., see [25]). It is also similar to (23) when $k_i \approx \mathcal{C}_{iiii}[\boldsymbol{y}]$, which is indeed the case close to separation.

One benefit of considering fourth-order orthogonal contrasts like $\phi_{\rm ICA}^{\rm o}$ is that they can be optimized by the Jacobi technique: the unknown rotation (Section II-B) can be found as a sequence of 2×2 rotations applied in sequence to all pairs (y_i,y_j) , for $i\neq j$ with the optimal angle at each step often being available in close form. Comon [25] has such a formula for $\phi_{\rm ICA}$ in the case of real signals.

Independence can also be tested on a smaller subset of cross-cumulants with

$$\phi_{\text{JADE}}^{\circ}[\boldsymbol{y}] \stackrel{\Delta}{=} \sum_{ijkl \neq ijkk} C_{ijkl}^{2}[\boldsymbol{y}].$$
 (25)

The motivation for using this specific subset is that ϕ_{JADE} also is a joint diagonalization criterion, implying that it can be optimized by the Jacobi technique for which the rotation angles can be found in close form, even in the complex case [22]. A similar technique is described in [31].

Simpler contrasts can be used if the kurtosis of the sources are known. For instance (23) suggests, for negative

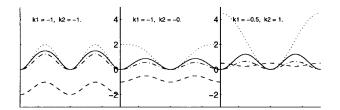


Fig. 10. Variation of orthogonal contrast functions (solid: $\phi_{\rm ICA}^{\circ}$, dash-dots: $\phi_{\rm JADE}^{\circ}$, dashes: $\phi_{\rm m}^{\circ}$, dots: $\phi_{\rm d}^{\circ}$) when two sources with kurtosis k_1 and k_2 are rotated between $-\pi/2$ and $\pi/2$. Left: $(k_1, k_2) = (-1, -1)$; center: $(k_1, k_2) = (-1, 0)$; right: $(k_1, k_2) = (-0.5, 1)$.

kurtosis $(k_i < 0 \ \forall i)$, a very simple contrast

$$\phi_m^{\circ}[\mathbf{y}] = \sum_{i=1}^n \mathbf{E} y_i^4 \tag{26}$$

(see [21], [53], [47], and [44]). Actually, the condition that k_i+k_j0 for all pairs of sources is sufficient for the stationary points of this orthogonal contrast function to be locally stable (see Section VI-A).

Some properties of the fourth-order contrasts discussed above are illustrated by Fig. 10, which displays the variation of some orthogonal contrast functions in the two-source case. A 2×1 source vector \mathbf{s} with kurtosis (k_1,k_2) is rotated into \mathbf{y} by an angle $\theta\in(-\pi/2,\pi/2)$. On the left panel, the sources have identical kurtosis $k_1=k_2=-1$; all the four contrasts are minimized at integer multiples of $\pi/2$. On the center panel, one source is Gaussian $(k_2=0)$; the contrasts show smaller variations except for ϕ_4^{o} . Note that ϕ_4 "knows" that one source has zero kurtosis, thus distinguishing between even and odd multiplies of $\pi/2$. On the right panel, $k_1=-0.5$ and $k_2=1$, so that $k_1+k_2>0$ which violates the condition for ϕ_m^{o} to be a contrast: its minima become maxima and vice versa. This is the same phenomenon illustrated by Fig. 7.

IV. ESTIMATING FUNCTIONS

By design, all valid contrast functions reach their minima at a separating point when the model holds; in this sense, no one is better than another. In practice, however, contrasts are only estimated from a finite data set. Sample-based contrasts depend not on the distribution of \boldsymbol{y} but on its *sample* distribution. Estimation from a finite data set introduces stochastic errors depending on the available samples and also on the contrast function. Thus, a statistical characterization of the minima of sample-based contrast functions is needed and will provide a basis for comparing contrast functions. For this purpose, the notion of estimating function is introduced. It is also closely related to gradient algorithms for BSS (Section V-A).

A. Relative Gradient

The variation of a contrast function $\phi[y]$ under a linear transform of y may be expressed by defining a "relative gradient." This specific notion builds on the fact that the parameter of interest is a square matrix.

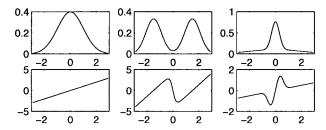


Fig. 11. Some densities and their associated scores.

Definition: An infinitesimal transform of y is $y \rightarrow (I + \mathcal{E})y = y + \mathcal{E}y$, where \mathcal{E} is a "small" matrix

$$oldsymbol{y}
ightarrow oldsymbol{I+\mathcal{E}}
ightarrow oldsymbol{y+\mathcal{E}y}.$$

If ϕ is smooth enough, $\phi[y + \mathcal{E}y]$ can be expanded as

$$\phi[\mathbf{y} + \mathcal{E}\mathbf{y}] = \phi[\mathbf{y}] + \sum_{i,j=1}^{n} G_{ij}\mathcal{E}_{ij} + o(||\mathcal{E}||)$$
 (27)

with G_{ij} being the partial derivative of $\phi[y + \mathcal{E}y]$ with respect to \mathcal{E}_{ij} at $\mathcal{E} = 0$. These coefficients form an $n \times n$ matrix, denoted as $\nabla \phi[y]$, and called the "relative gradient" [21] of $\phi[y]$ at [y]. In matrix form, expansion (27) reads

$$\phi[\mathbf{y} + \mathcal{E}\mathbf{y}] = \phi[\mathbf{y}] + \langle \nabla \phi[\mathbf{y}] | \mathcal{E} \rangle + o(||\mathcal{E}||)$$
 (28)

where $\langle \cdot | \cdot \rangle$ is the Euclidean scalar product between matrices $\langle M | N \rangle = \operatorname{trace}(MN^{\dagger}) = \sum_{i,j=1}^{n} M_{ij} N_{ij}$.

Note that the relative gradient is defined without explicit reference to the possible dependence of \boldsymbol{y} on B as $\boldsymbol{y} = B\boldsymbol{x}$. Thus, it actually characterizes the first order variation of the contrast function itself. It is of course possible to relate $\nabla \phi[\boldsymbol{y}]$ to a regular gradient with respect to B if $\boldsymbol{y} = B\boldsymbol{x}$. Elementary calculus yields

$$\nabla \phi[B\mathbf{x}] = B^{\dagger} \frac{\partial \phi[B\mathbf{x}]}{\partial B}.$$
 (29)

The notion of "natural gradient" was independently introduced by Amari [1]. In general, it is distinct from the relative gradient in that the latter is defined in any continuous group of transformation while the former is defined in any smooth statistical model. However, for the BSS model, which as a statistical transformation model combines both features, the two ideas yield the same class of algorithms (Section V-A).

1) Score Functions: The source densities q_1, \dots, q_n , used in (3) and (7) to define the likelihood of a BSS model, enter into the estimating function via their log-derivatives, the so-called score functions $\varphi_1, \dots, \varphi_n$, defined as

$$\varphi_i \stackrel{\Delta}{=} -(\log q_i)' \quad \text{or} \quad \varphi_i(\cdot) = -\frac{q_i(\cdot)'}{q_i(\cdot)}.$$
 (30)

Fig. 11 displays some densities and their associated score functions.

Note that the score for the most basic distribution is the most basic function: if s is a zero-mean unit-variance Gaussian variable $q(s) = (2\pi)^{-1/2} \exp{-(s^2/2)}$, then the associated score is $\varphi(s) = s$. Actually, Gaussian densities

are precisely these densities associated with linear score functions. Thus, the necessity of non-Gaussian modeling (recall Section II) translates into the necessity of considering nonlinear score functions.

2) Relative Gradient of the Likelihood Contrast: At the core of the BSS contrast functions is $\phi_{ML}[y]$, which is associated with the likelihood given the source densities q_1, \dots, q_n . Its relative gradient is found in [61] to be

$$\nabla \phi_{ML}[\boldsymbol{y}] = \mathbf{E} H_{\varphi}(\boldsymbol{y}) \tag{31}$$

where $H_{\varphi} \colon \mathbb{R}^n \mapsto \mathbb{R}^{n \times n}$ is

$$H_{\varphi}(\mathbf{y}) \stackrel{\Delta}{=} \varphi(\mathbf{y})\mathbf{y}^{\dagger} - I \tag{32}$$

with $\varphi \colon \mathbb{R}^n \mapsto \mathbb{R}^n$ the entry-wise nonlinear function

$$\varphi(\mathbf{y}) \stackrel{\Delta}{=} [\varphi_1(y_1), \cdots, \varphi_n(y_n)]^{\dagger}$$
 (33)

collecting the score functions related to each source. This is a remarkably simple result; this relative gradient merely is the expected value of a function H_{φ} of \boldsymbol{y} .

3) Interpretation: The ML contrast function ϕ_{ML} is minimum at points where its (relative) gradient cancels (i.e., by (31)) at those points which are solutions of the matrix equation $EH_{\varphi}(y) = 0$. This is interpreted by examining the (i,j)th entry of this matrix equation. For i=j, we find $E\varphi_i(y_i)y_i = 1$ which depends only on y_i and determines the scale of the *i*th source estimate. For $i \neq j$, the (i,j)th entry of $EH_{\varphi}(\boldsymbol{y})=0$ reads $E\varphi_i(y_i)y_j=0$ meaning that the jth output y_i should be uncorrelated to a nonlinear version $\varphi_i(y_i)$ of the *i*th output. Because φ_i and φ_i are nonlinear functions, the conditions for the pairs (i,j) and the pair (j,i) are (in general) not equivalent. Note that if the source signals are modeled as zero-mean unit-variance normal variables, then $\varphi_i(y_i) = y_i$ for all i and $H_{\varphi}(y) = yy^{\dagger} - I = H_{w}(y)$ [recalling (5)]. Then $\phi_{ML}[y]$ is minimum at points where $Eyy^{\dagger} = I$; we only obtain the whiteness condition. Again, this is not sufficient to determine a separating solution; score functions must be nonlinear (the source model must be non-Gaussian).

The idea of using nonlinear functions to obtain a sufficient set of independence conditions can be traced back to the seminal paper of Hérault and Jutten [43] (see [45] for an English translation), but the choice of the nonlinear functions was somewhat *ad hoc*. Féty [38] gave an interpretation of the nonlinear functions as amplifiers for the signals of interest; Bar-Ness also produced early work using nonlinear functions [7]. However, the ML principle makes it clear that the nonlinear functions are related via (30) a non-Gaussian model of the source distributions.

B. Estimating Functions

An estimating function for the BSS problem is a function $H: \mathbb{R}^n \mapsto \mathbb{R}^{n \times n}$. It is associated with an estimating equation

$$\frac{1}{T} \sum_{t=1}^{T} H(\mathbf{y}(t)) = 0$$
 (34)

thus called because H being matrix-valued, (34) specifies a priori as many constraints as unknown parameters in the BSS problem. Many BSS estimates can be characterized via an estimating function [17], [2].

A simple instance of an estimating function is $H_w(\boldsymbol{y})$, used in (5) to express the decorrelation between the entries of \boldsymbol{y} . Equation (34) with $H(\boldsymbol{y}) = H_w(\boldsymbol{y})$ is equivalent to $(1/T) \sum_t \boldsymbol{y}(t)\boldsymbol{y}(t)^\dagger = I$; it expresses the empirical whiteness of a batch of T samples of \boldsymbol{y} as opposed to the actual whiteness, i.e., $EH_w(\boldsymbol{y}) = 0$. The estimating function $H_w(\boldsymbol{y})$, however, is not appropriate for BSS, since whitening (or decorrelation) is not sufficient to determine a separating matrix.

The simplest example of an estimating function for BSS is obtained in the ML approach. The gradient of the likelihood (7) may be shown [61] to cancel at points A_{ML} , which are characterized exactly by (34) with $\mathbf{y} = A_{ML}^{-1} \mathbf{x}$ and $H = H_{\varphi}$, as defined in (32). In other words, maximum likelihood estimates correspond exactly to the solution of an estimating equation. This equation is nothing but the sample counterpart of $EH_{\varphi}(\mathbf{y}) = 0$ which characterizes the stationary points of $\phi_{ML}[\mathbf{y}]$. Recall that the latter is obtained [at (11) and (12)] as a limit of the log likelihood.

Because the value of an estimating function is a square matrix, it can be decomposed into a symmetric part (equal to its transpose) and a skew symmetric part (opposite of its transpose). This decomposition is simply

$$H(\mathbf{y}) = \frac{H(\mathbf{y}) + H(\mathbf{y})^{\dagger}}{2} + \frac{H(\mathbf{y}) - H(\mathbf{y})^{\dagger}}{2}.$$
 (35)

If the optimization of some regular contrast function corresponds to an estimating function $H(\boldsymbol{y})$, it is found that the optimization of the same contrast under the whiteness constraint corresponds to an estimating function $H^{\circ}(\boldsymbol{y})$ given by

$$H^{\circ}(\boldsymbol{y}) = H_w(\boldsymbol{y}) + \frac{1}{2} (H(\boldsymbol{y}) - H(\boldsymbol{y})^{\dagger}). \tag{36}$$

Thus, the symmetric part of H(y) is replaced by $H_w(y) = yy^{\dagger} - I$ [introduced at (5)], whose effect is to enforce the whiteness constraint. In particular, maximum likelihood estimates under the whiteness constraint are (again) solutions of (34) with the estimating function $H = H_{\varphi}^{\circ}$

$$H_{\varphi}^{\circ}(\mathbf{y}) \stackrel{\Delta}{=} \mathbf{y} \mathbf{y}^{\dagger} - I + \varphi(\mathbf{y}) \mathbf{y}^{\dagger} - \mathbf{y} \varphi(\mathbf{y})^{\dagger}.$$
 (37)

Other orthogonal contrast functions are associated to similar estimating functions. For instance, the simple fourth-order contrasts $\phi_4^{\circ}[y]$ and $\phi_m^{\circ}[y]$ [(23) and (26), respectively] yield estimating equations in the form (37) with nonlinear functions given respectively by

$$\varphi_i(y_i) = -k_i y_i^3$$
 and $\varphi_i(y_i) = y_i^3$. (38)

Recall that using the contrast function (26) assumes sources with negative kurtosis k_i . Thus, the two functions in (38) agree on the sign to be given to a cubic distortion (as was expected).

Some contrast functions, like ϕ_{ICA}° and ϕ_{JADE}° , when estimated from T samples are minimized at points which

cannot be represented exactly as the solution of (34) for a fixed estimating function. However, one can often find, as in [36], an asymptotic estimating function in the sense that the solution of the associated estimating equation is very close to the minimizer of the estimated contrast. For instance, the contrast $\phi_{\rm ICA}^{\circ}$ and $\phi_{\rm JADE}^{\circ}$ are asymptotically associated to the same estimating function as ϕ_4° . This implies that minimizing $\phi_{\rm ICA}^{\circ}$, $\phi_{\rm JADE}^{\circ}$ or ϕ_4° with cumulants estimated from T samples yields estimates which are equivalent (i.e., they differ by a term which is smaller than the estimation error) for large enough T.

Which functions are appropriate as estimating functions? One could think of using any H such that $\mathrm{E}H(s)=0$ as an estimating function because the estimating (34) would just be the sample counterpart of $\mathrm{E}H(y)=0$ and would a priori provide as many scalar equations as unknown parameters. However, the ML principle suggests the specific forms (32) and (37), with the nonlinear functions in $\varphi(y)$ being (approximations of) the score functions for the probability densities of the signals to be separated.

V. ADAPTIVE ALGORITHMS

A simple generic technique for optimizing an objective function is gradient descent. In most optimization problems, its simplicity is at the expense of performance; more sophisticated techniques—such as "Newton-like" algorithms using second derivatives in addition to the gradient—can often significantly speed up convergence. For the BSS problem, however, it turns out that a simple gradient descent offers "Newton-like" performance (see below). This surprising and fortunate result is obtained by descending along the relative gradient defined in Section IV-A.

A. Relative Gradient Techniques

1) Relative Gradient Descent: We first describe a generic relative gradient descent. Generally, the steepest descent technique of minimization consists in moving by a small step in a direction opposite that of the gradient of the objective function. The relative gradient of a contrast $\phi[y]$ is defined (Section IV-A) with respect to a relative variation of y, by which y is changed into $(I + \mathcal{E})y$. The resulting variation of $\phi[y]$ is (at first order) the scalar product $\langle \nabla \phi[y] | \mathcal{E} \rangle$ between the relative variation \mathcal{E} and the relative gradient $\nabla \phi[y]$, as in (27) or (28). Aligning the direction of change in the direction opposite to the gradient is to take $\mathcal{E} = -\mu \nabla \phi[y]$ for a small positive step μ . Thus, one step of a relative gradient descent can be formally described as

$$\mathbf{y} \leftarrow (I - \mu \nabla \phi[\mathbf{y}])\mathbf{y} = \mathbf{y} - \mu \nabla \phi[\mathbf{y}]\mathbf{y}.$$
 (39)

According to (28), the resulting variation of $\phi[\boldsymbol{y}]$ is $\delta \phi \approx \langle \nabla \phi[\boldsymbol{y}] | \mathcal{E} \rangle = \langle \nabla \phi[\boldsymbol{y}] | - \mu \nabla \phi[\boldsymbol{y}] \rangle = -\mu ||\nabla \phi[\boldsymbol{y}]||^2$, which is negative for positive μ .

The formal description (39) can be turned into off-line and on-line algorithms, as described next.

2) Off-Line Relative Gradient Descent: Consider the separation of a batch $x(1), \dots, x(T)$ of T samples based on the minimization of a contrast function $\phi[y]$ with relative

gradient $\nabla \phi[\boldsymbol{y}] = \mathrm{E} H(\boldsymbol{y})$. One looks for a linear transform of the data satisfying the corresponding estimating equation (1/T) $\Sigma_{t=1}^T$ $H(\boldsymbol{y}(t)) = 0$. The relative gradient descent to solve it is as follows. Set $\boldsymbol{y}^{(0)}(t) = \boldsymbol{x}(t)$ for $1 \leq t \leq T$ and iterate through the following two steps

$$\hat{\mathcal{H}} \leftarrow \frac{1}{T} \sum_{t=1}^{T} H(\mathbf{y}(t)) \tag{40}$$

$$\mathbf{y}(t) \leftarrow \mathbf{y}(t) - \mu \hat{\mathcal{H}} \mathbf{y}(t) \qquad (1 \le t \le T).$$
 (41)

The first step computes an estimate $\hat{\mathcal{H}}$ of the relative gradient for the current values of the data; the second step updates the data in the (relative) direction opposite to the relative gradient, as in (39). The algorithm stops when $(1/T) \sum_{t=1}^T H(\boldsymbol{y}(t)) = 0$, i.e., when the estimating equation is solved. It is amusing to note that this implementation does not need to maintain a separating matrix, rather, it directly operates on the data set itself with the source signals emerging during the iterations.

3) On-Line Relative Gradient Descent: On-line algorithms update a separating matrix B_t upon reception of a new sample $\boldsymbol{x}(t)$. The (relative) linear transform $\boldsymbol{y} \leftarrow (I+\mathcal{E})\boldsymbol{y}$ corresponds to changing B into $(I+\mathcal{E})B=B+\mathcal{E}B$. In the on-line mode, one uses the stochastic gradient technique where the gradient $\nabla \phi[\boldsymbol{y}] = \mathrm{E} H[\boldsymbol{y}]$ is replaced by its instantaneous value $H(\boldsymbol{y}(t))$. Hence the stochastic relative gradient rule

$$B_{t+1} = B_t - \mu_t H(\mathbf{y}(t)) B_t \tag{42}$$

where μ_t is a sequence of positive learning steps.

4) Uniform Performance of Relative Gradient Descent: A striking feature of the BSS model is that the hardness (in a statistical sense, as discussed in Section VI-C) of separating mixed sources does not depend on the particular value of the mixing matrix A; the problem is uniformly hard in the mixture. Very significantly, the device of relative updating produces algorithms which also behave uniformly well in the mixture. Right-multiplying the updating rule (42) by matrix A and using $\mathbf{y} = B\mathbf{x} = BA\mathbf{s}$, one readily finds that the trajectory of the global system $C_t \stackrel{\triangle}{=} B_t A$ which combines mixing and unmixing matrices is governed by

$$C_{t+1} = C_t - \mu_t H(C\mathbf{s}(t))C_t. \tag{43}$$

This trajectory is expressed here as a sole function of the global system C_t . The only effect of the mixing matrix A itself is to determine (together with B_0) the initial value $C_0 = B_0 A$ of the global system. This is a very desirable property since it means that the algorithms can be studied and optimized without reference to the actual mixture to be inverted. This is true for any estimating function H(y). However, uniformly good performance can only be expected if the H(y) is adjusted correctly to the distribution of the source signals, e.g., by deriving it from a contrast function. Algorithms based on an estimating function in the form (32) are described in [5] for the on-line version and in [61] for an off-line version; those based on form (37) are studied in detail in [21]. The uniform performance property was also obtained in [24].

5) Regular Gradient Algorithms: It is interesting to compare the relative gradient algorithm to the algorithm obtained by a regular gradient, that is, by applying a gradient rule to the entries of B for the minimization of $f(B) \stackrel{\triangle}{=} \phi[Bx]$. This is

$$B_{t+1} = B_t - \mu_t H(y(t)) B_t^{-\dagger}. \tag{44}$$

Not only is this form more costly because it requires (in general) the inversion of B_t at each step, but it lacks the uniform performance property. The trajectory of the global system depends on the particular mixture A to be inverted.

B. Adapting to the Sources

The iterative and adaptive algorithms described above require the specification of an estimating function H, for which two forms H_{φ} and H_{φ}° [(32) and (37)] are suggested by the theory. These forms, in turn, depend on nonlinear functions $\varphi_1, \dots, \varphi_n$ which, ideally, should be the score functions associated to the distributions of the sources (Section IV-B). When the source distributions are unknown, one may try to estimate them from the data (for instance, using some parametric model, as in [56]) or to directly estimate good nonlinear functions.

The first idea is to use Edgeworth expansions (e.g., see [51]) which provide approximations to probability densities in the vicinity of a Gaussian density. The simplest nontrivial Edgeworth approximation of a symmetric pdf q in the vicinity of the standard normal distribution is

$$q(s) = \frac{1}{\sqrt{2\pi}} \exp\left(-\frac{s^2}{2}\right) \left(1 + \frac{k}{24} \left(s^4 - 6s^2 + 3\right) + \cdots\right)$$

where k is the kurtosis of q. The corresponding approximate score function is

$$\varphi(s) = s - \frac{k}{6} (s^3 - 3s) + \cdots$$
 (45)

Thus, the Edgeworth expansion suggests that in a linearcubic approximation of the score function, the coefficient of the cubic part should be $-k_i/6$ for the *i*th source. Asymptotic analysis shows that such a choice at least guarantees the local stability (Section VI-A). There are other possibilities for deriving score functions by a density expansion, e.g., see [68] for a different proposal involving odd and even terms in φ .

A more direct approach than pdf expansion is proposed by Pham [61], who considers approximating ψ by a linear combination

$$\varphi_{\alpha}(s) \stackrel{\Delta}{=} \sum_{l=1}^{L} \alpha_{l} f_{l}(s) \tag{46}$$

of a fixed set $\{f_1, \dots, f_L\}$ of arbitrary basis functions. Rather surprisingly, the set $\{\alpha_1, \dots, \alpha_L\}$ of coefficients minimizing the mean square error $\mathrm{E}(\varphi_\alpha(s) - \psi(s))^2$ between the true score ψ and its approximation can be found without knowing ψ . The best mean square approximation involves only the expectation operator. It is

$$\varphi_*(s) = (\mathbf{E}F'(s))^{\dagger} (\mathbf{E}F(s)F(s)^{\dagger})^{-1} F(s) \tag{47}$$

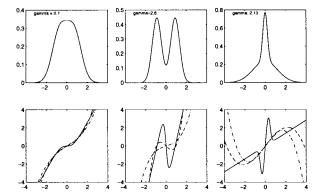


Fig. 12. Top row: three distributions and the values of γ_* as a measure of non-Gaussianity (see Section VI-B). Bottom row: the score function (solid) and its linear–cubic approximations, based on Edgeworth expansion (dashes-dots) and optimal (dashes).

where $F(s) \stackrel{\Delta}{=} [f_1(s), \cdots, f_L(s)]^\dagger$ is the $L \times 1$ column vector of basis functions and F'(s) is the column vector of their derivatives. This is a nice result because the expression of φ_* can be estimated simply by replacing in (47) expectations by sample averages and the values of s by the estimated source signals.

The two approaches of Edgeworth expansion and mean square fit, respectively leading to the approximations (45) and (47), are compared in Fig. 12.

Three pdf's are displayed in the top row; the bottom row shows the corresponding score function (solid line), the linear–cubic approximation by (45) (dash-dotted line), and the Pham approximation (dashed line) obtained from (47) with $F(s) = [s, s^3]$. Both approximations are similar in the first example when the pdf is close to Gaussian; in the second case, the optimal approximation fits the true score much better in the area of highest probability. None of the approximations seem good in the third example for the simple reason that the true score there cannot be well approximated by a linear–cubic function. However, the two approximations fit the score well enough to guarantee the stability of the gradient algorithms (see Section VI-A).

VI. PERFORMANCE ISSUES

This section is concerned with the performance of BSS algorithms and presents some asymptotic analysis results. It has been stressed repeatedly that it was not necessary to know the source distributions (or equivalently, the associated score functions) with great accuracy in order to obtain consistent BSS algorithms. There is, however, a limit to the misspecification of the source distributions, as illustrated by Fig. 7. This is elucidated at Section VI-A, which gives explicit stability limits. Even if a hypothesized distribution is good enough to preserve stability, one may expect a loss of estimation accuracy due to misspecification when a finite number of samples are available; this is quantified in Section VI-B, which also describes the ultimate achievable separation performance. Section VI-C discusses the general property of equivariance which governs the performance of BSS algorithms.

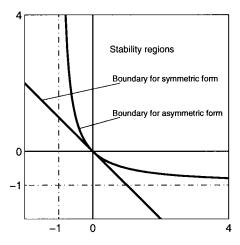


Fig. 13. Stability domains in the (κ_i, κ_j) plane.

A. Local Stability

A stationary (or equilibrium) point B of the learning rule (42) is characterized by $\mathrm{E}H(y)=\mathrm{E}H(Bx)=0$, i.e., the mean value of the update is zero. We have seen that separating matrices (with the proper scale) are equilibrium points; we are now interested in finding when they are locally stable, that is, when a small deviation from the equilibrium is pulled back to the separating point. In other words, we want the separating matrix to a (local) attractor for the learning rule (42). In the limit of small learning steps, there exists a simple criterion for testing local stability which depends on the derivative of $\mathrm{E}H(Bx)$ with respect to B. For both the symmetric form H_{φ}° and for the asymmetric form H_{φ} the stability condition can be worked out exactly. They are found to depend only the following nonlinear moments:

$$\kappa_i \stackrel{\Delta}{=} E\varphi_i'(s_i)Es_i^2 - E\varphi_i(s_i)s_i \tag{48}$$

where each s_i is rescaled according to EH(s) = 0, that is, $E\varphi_i(s_i)s_i = 1$ for $H = H_{\varphi}$, or $Es_i^2 = 1$ for $H = H_{\varphi}^{\circ}$.

Leaving aside the issue of stability with respect to scale, the stability conditions for the symmetric form (37) are [21]

$$(1 + \kappa_i)(1 + \kappa_j) > 1$$
 for $1 \le i < j \le n$ (49)

and for the asymmetric form (32), the conditions are [3] that $1 + \kappa_i > 0$ for $1 \le i \le n$ and that

$$\kappa_i + \kappa_j > 0 \quad \text{for} \quad 1 \le i < j \le n.$$
 (50)

Therefore, stability appears to depend on pairwise conditions. The stability domains for a given pair of sources are displayed in Fig. 13 in the (κ_i, κ_i) plane.

Note that the stability domain is larger for the symmetric form (37). This is a consequence of letting the second-order information (the whiteness constraint) do "half the job" (see Section II-B).

Some comments are in order. First, it appears that in both cases a sufficient stability condition is $\kappa_i > 0$ for all the sources. Thus, regarding stability, tuning the nonlinear functions φ_i 's to the source distributions should be understood as making the κ_i 's positive. Second, one can show

that if s_i is Gaussian, then $\kappa_i=0$ for any function φ_i . Therefore, the stability conditions can never be met if there is more than one Gaussian source, in agreement with the identifiability statements of Section II. Third, it can also be shown that if φ_i is taken to be the score function for the true density of s_i , then $\kappa_i \geq 0$ with equality only if s_i is Gaussian.

Section II-C illustrated the fact that the hypothesized source distributions should be close enough to the true distributions for the likelihood to still show a maximum around a separating point. The definition of κ_i provides a quantitative measure of how wrong the hypothesis can be—they should not allow κ_i to become negative.

We also note that it is not necessary for all the κ_i 's to be positive; if $\kappa_i < 0$ for one source at most, this can be compensated if the moments κ_j are large enough for all $j \neq i$. As seen from the stability domains (Fig. 13), one source at most can have an arbitrarily negative κ_i if the symmetric form is used, while the stability of the asymmetric form requests that $\kappa_i > -1$.

We have considered linear–cubic score functions in Sections IV and V. If $\varphi_i(s_i) = \alpha_i s_i + \beta_i s_i^3$ for two constants α_i and β_i , then $\kappa_i = \beta_i (3\mathrm{E}^2 s_i^2 - \mathrm{E} s_i^4) = -\beta_i k_i$ where, as above, k_i denotes the kurtosis. Note that the linear part of φ_i does not affect the stability and that stability is guaranteed if the coefficient β_i of the cubic part has a sign opposite to the sign of the kurtosis. Quite naturally, the functions in (38) and (45) come up naturally with the right sign. Therefore, if one wishes to use cubic nonlinearities, it is sufficient to know the sign of the kurtosis of each source to make separating matrices stable. For other than cubic scores, stability depends on the sign of κ_i , not on the sign of the kurtosis.

B. Accuracy of Estimating Equations

This section characterizes the accuracy of signal separation obtained by solving (34) with T independent realizations of \boldsymbol{x} .

If a matrix B is used for separation, the pth entry of $\mathbf{y} = B\mathbf{x} = BA\mathbf{s}$ contains the signal of interest s_p at power $(BA)_{pp}^2 \sigma_p^2$, and the qth interfering signal s_q at power $(BA)_{pq}^2 \sigma_q^2$. Therefore, for a given matrix B, the quantity

$$\rho_{pq}(B) \stackrel{\triangle}{=} \frac{(BA)_{pq}^2 E s_q^2}{(BA)_{pp}^2 E s_p^2} \qquad p \neq q$$
 (51)

measures the interference-to-signal ratio (ISR) provided by B in rejecting the qth source in the estimate of the pth source. Let \hat{B}_T be the separating matrix obtained via a particular algorithm using T samples. In general, the estimation error in regular statistical models decreases as $1/\sqrt{T}$ so that the limit

$$ISR_{pq} \stackrel{\triangle}{=} \lim_{T \to \infty} TE\rho_{pq} \Big(\hat{B}_T \Big)$$
 (52)

usually exists and provides an asymptotic measure of performance separation of a given off-line BSS technique. When $H=H_{\varphi}$ or $H=H_{\varphi}^{\circ}$ are used in the estimating

equation, the asymptotic ISR depends on the moments κ_i in (48) and

$$\gamma_i \stackrel{\Delta}{=} \mathrm{E}\varphi_i^2(s_i)\mathrm{E}s_i^2 - \mathrm{E}^2(\varphi_i(s_i)s_i) \ge 0.$$
 (53)

For simplicity, we consider identically distributed signals and identical nonlinear functions $\varphi_i(\cdot) = \varphi(\cdot)$, so that $\kappa_i = \kappa$ and $\gamma_i = \gamma$ for $1 \le i \le n$. With a symmetric estimating function H°_{φ} one finds

$$ISR_{pq}^{\circ} = ISR^{\circ} = \frac{1}{2} \left(\frac{\gamma}{\kappa^2} + \frac{1}{2} \right) p \neq q.$$
 (54)

Note that ISR° is lower bounded by 1/4 regardless of the value of γ . This is a general property of orthogonal BSS techniques [15] and is the price to pay for blindly trusting second-order statistics, i.e., for whitening. Thus rejection rates obtained under the whiteness constraint cannot be (asymptotically) better that 1/4T.

For an asymmetric estimating function H_{φ} , the ISR does not take such a simple form unless the common score φ is obtained by Pham's method (Section V-B). One then finds $ISR_{pq} = ISR$ and $ISR_{pq}^{\circ} = ISR$ ° as

$$ISR = \frac{1}{2} \left(\frac{1}{\gamma} + \frac{1}{\gamma + 2} \right) \quad ISR^{\circ} = \frac{1}{2} \left(\frac{1}{\gamma} + \frac{1}{2} \right) \quad (55)$$

where the last equation stems from (54) because Pham's method guarantees $\gamma = \kappa$. These expressions show that both ISR and ISR° are minimized by maximizing γ ; not surprisingly, γ can be shown to reach its maximum value γ_* precisely when $\varphi = \psi$, where ψ is the score function corresponding to the true density of the sources

$$\gamma_* \stackrel{\Delta}{=} E\psi^2(s)Es^2 - E^2[\psi(s)s]. \tag{56}$$

Note that the solution of (34) with $H=H_{\psi}$ then is the ML estimator based on the true model. It follows that the expression of ISR in (55) also is the asymptotic Cramér–Rao bound for source separation, i.e., the best achievable ISR rate with T independent samples (see [58], [69], and [61]).

Since the achievable performance depends on the magnitude of γ_* , this moment characterizes the hardness of the BSS problem with respect to source distribution. Not surprisingly, we can relate it to the non-Gaussianity of the sources as follows. As above, we denote ψ the score function for the (true) distribution of s, and we denote ψ_n the score function for the Gaussian distribution with the same variance as s (this is just $\psi_n(s) = s/\mathrm{E} s^2$). A "large" non-Gaussianity translates into a large difference between ψ and ψ_n . As we just saw, the measure of non-Gaussianity from the asymptotic point of view is measured by γ_* . Indeed, one finds

$$\gamma_* = \frac{\mathrm{E}(\psi(s) - \psi_n(s))^2}{\mathrm{E}(\psi_n(s))^2}.$$
 (57)

See Fig. 12 for the values of γ_* in three examples. For close-to-Gaussian sources, γ_* is (arbitrarily) small. In this case, according to (55), the best achievable rejection rates

are about $1/2\gamma_*T$ for both the symmetric and the asymmetric forms. This gives an idea of the minimum number of samples required to achieve a given separation. The other extreme is for sources which are far away from normality: the moment γ_* is not bounded above. In particular, it tends to ∞ when the source distributions tend to have a discrete or a bounded support. In the case of discrete sources, deterministic (error-free) blind identification is possible with a finite number of samples. In the case of sources with bounded support, the MSE of blind identification decreases at a much faster rate than the 1/T rate obtained for finite values of γ (see in particular [41]).

C. Equivariance and Uniform Performance

At first thought, the hardness of the BSS problem seems to depend on the distributions of the source signals and on the mixing matrix, with harder problems occurring when sources are nearly Gaussian and when the mixing matrix is poorly conditioned. This is not correct, however. The BSS problem is uniformly hard in the mixing matrix. Let us summarize the instances where this property appeared: the ultimate separation performance depends only on γ_* [(55)]; the asymptotic performance index in (54) and (55) depends only on some statistical moments; the stability of the adaptive algorithms (42) also depends only on the values of κ_i 's; even better, the trajectory (43) of the global system $C_t = B_t A$ does not depend on A, whose sole effect is to determine the initial point.

Therefore, not only does the problem appear to be uniformly hard in the mixing matrix, but there exist estimation techniques with a statistical behavior (regarding signal separation) which is independent of the particular value of the system to be inverted. This is a very desirable property: such algorithms can be studied and tuned independently of the particular mixture to be inverted and their performance can also be predicted independently of the mixture [16]. This is an instance of "equivariance," a property holding more generally in transformation models.

There is a simple prescription to design algorithms with uniform performance: adjust the separating matrix freely (i.e., without constraint) according to a rule expressed only in terms of the output y. To understand why the "output only" prescription ensures uniform performance, consider using a particular estimating function $H(\cdot)$ to separate a mixture of T samples $[s(1), \dots, s(T)]$. If the source signals are mixed by a given matrix A, then a solution of (34) is a matrix B such that $BA = \hat{C}$ where matrix \hat{C} is a solution of $T^{-1} \Sigma_{t=1}^T H(\hat{C}\mathbf{s}(t)) = 0$. Matrix \hat{C} does not depend on Aso that the global system $BA = \hat{C}$ is itself independent of A, and the estimated signals are $y(t) = \hat{C}s(t)$ regardless of A. In particular, the recovered signals are exactly identical to those that would be obtained with A = I, i.e., when there is no mixing at all. This argument, based on estimating equations, extends to the minimizers of contrast functions since the latter are defined as functions of the distribution of the output (the argument also applies to orthogonal contrast functions because the whiteness constraint is expressed only

in terms of y). The argument also justifies the specific definition of the relative gradient; a device was needed to express the first-order variations of a contrast function $\phi[y]$ in terms of a variation of y itself, i.e., without reference to B. Finally, it must be stressed that the argument does not involve asymptotics, rather, equivariance is observed exactly for any finite value of T.

Not all BSS algorithms are equivariant. For instance, Jutten and Hérault's original algorithm imposes constraints on the separating matrix resulting in a greatly complicated analysis (and behavior) (see [34], [41], and [48]). Other instances of nonequivariant techniques are to be found in most of the algebraic approaches (see Section VII) based on the structure of the cumulants of the observed vector \boldsymbol{x} . Precisely because the identification is based on \boldsymbol{x} and not on \boldsymbol{y} , such approaches are not generally equivariant unless they can be shown to be equivalent to the optimization of a contrast function of \boldsymbol{y} .

A word of caution is necessary before concluding: equivariance holds exactly in the noise-free model which we have considered so far. In practice, there is always some kind of noise which must be taken into account. Assume that a better model is $\mathbf{x} = A\mathbf{s} + \mathbf{n}$, where \mathbf{n} represents an additive noise. This can be rewritten as $\mathbf{x} = A(\mathbf{s} + A^{-1}\mathbf{n})$. As long as $A^{-1}\mathbf{n}$ can be neglected with respect to \mathbf{s} , this is a noise-free situation. This shows the limit of equivariance: a poorly conditioned matrix A has a large inverse which amplifies the effect of the noise. More precisely, we can expect equivariance in the high-SNR domain, i.e., when the covariance matrix of \mathbf{s} remains "larger" than the covariance matrix of $A^{-1}\mathbf{n}$.

VII. CONCLUSIONS

Due to limited space, the focus of this paper is on principles and many interesting issues have been left out, such as a discussion of the connections between BSS and blind deconvolution, convergence rates of adaptive algorithms, design of consistent estimators based on noisy observations, detection of the number of sources, etc. Before concluding, we briefly mention some other points.

A. Algebraic Approaches

The fourth-order cumulants of \boldsymbol{x} have a very regular structure in the BSS model

$$C_{ijkl}[\mathbf{x}] = \sum_{p=1}^{n} k_p A_{ip} A_{jp} A_{kp} A_{lp} \qquad 1 \le i, j, k, l \le n.$$
(58)

Given sample estimates of the cumulants, the equation set (58) (or some subset of it) can be solved in A in a least square sense. This is a cumulant matching approach [70], [42] which does not yield equivariant estimates. Optimal matching, though, can be shown to correspond to a contrast function [21]. However, the specific form of (58) also calls for algebraic approaches. Simple algorithms can be based on the eigen-structure of cumulant matrices built from cumulants [62], [64]. An exciting direction of

research is to investigate high-order decompositions that would generalize matrix factorizations like SVD or EVD to fourth-order cumulants [32], [29], [14], [26], [20].

B. Using Temporal Correlation

The approaches to BSS described above only exploit properties of the distribution of $\boldsymbol{x}(t)$. If the source signals are temporally correlated, time structures can also be exploited. It is possible to achieve separation if all the source signals have distinct spectra, even if each source signal is a Gaussian process [66]. Simple algebraic techniques can be devised (see [65] and [10]); the Whittle approximation of likelihood is investigated in [60]. Cyclostationary properties, when they exist, can also be exploited [12].

C. Deterministic Identification

As indicated in Section VI-B, sources with discrete support allow for deterministic identification (infinite Fisher information). Specific contrast functions can be devised [41] to take advantage of discreteness. There is a rich domain of application with digital communication signals coding information with discrete symbols by which deterministic identification is possible. See the review by Van der Veen [67] and the papers on CMA in this issue.

D. Open Problems and Perspectives

- 1) Learning Source Distributions: In the BSS problem, source distributions are a nuisance parameter. For a large enough sample size, it is possible to estimate the distributions and still obtain the same asymptotic performance as if the distributions were known in advance [2]; the design of practical algorithms achieving source adaptivity still is an open question.
- 2) Dealing with Noise: BSS techniques remaining consistent in the presence of additive noise has not been described here. For additive Gaussian noise, such techniques may resort to high-order cumulants or to noise modeling. It is not clear, however, that it is worth combating the noise. As a matter of fact, one may argue that taking noise effects into account is unnecessary at high-SNR and futile at low-SNR (because the BSS problem becomes too difficult anyway). Therefore, we believe it is still an open question to determine which application domains would really benefit from noise modeling.
- 3) Global Convergence: Some cumulant-based contrast functions can be proven to be free of spurious local minima in the two-source case (e.g., see [28]) or in a deflation approach (successive extractions of the source signals) [33], [34]. There is, however, a lack of general understanding of the global shape of contrast functions in the general case.
- 4) Multidimensional Independent Components: An interesting original variation of the basic independent component analysis (ICA) model would be to decompose a random vector in a sum of independent components with the requirement that the components are linearly independent but not necessarily one-dimensional. In the BSS model, this would be equivalent to grouping the source signals

in subsets with independence between the subsets but not within the subsets. This more general decomposition could be called multidimensional independent component analysis (MICA).

- 5) Convolutive Mixtures: The most challenging open problem in BSS probably is the extension to convolutive mixtures. This is a very active area of research, mainly motivated by applications in the audio-frequency domain where the BSS is often termed the "cocktail-party problem." The convolutive problem is significantly harder than the instantaneous problem: even input/output (i.e., nonblind) identification is very challenging because of the large number of parameters usually necessary to describe audio channels.
- 6) When the Model Does Not Hold: The introduction mentioned successful applications of BSS to biomedical signals. When examining these data it is very striking to realize that the extracted source signals seem to be very far from obeying the simple BSS model. The fact that BSS still yields apparently meaningful (to the experts) results is worth of consideration. A partial explanation stems from basing separation on contrast functions: even if the model does not hold (there are no independent source signals and no system A to be inverted), the algorithms still try to produce output which are as independent as possible. This does not tell the whole story, though, because for many data sets a stochastic description does not seem appropriate. We believe it will be a very interesting challenge to understand the behavior of BSS algorithms when applied "outside of the model."

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