

Chapter 3: Independent Component Analysis^a

As in the chapter on factor analysis (Chapter 2) we assume that the observations $\{x_i^0\}_{i=1}^n$ are n realizations of an \mathbb{R}^p -valued random variable X^0 , and we let $X = X^0 - \mathbb{E}[X]$.

Then, the independent component analysis (ICA) **model** assumes that

$$X = \mathbf{B}S, \quad \mathbf{B} \in \mathbb{R}^{p \times p}, \quad \mathbb{E}[S] = 0, \quad \text{Var}(S) = \mathbf{I}_p \quad (3.1)$$

where the p components of the \mathbb{R}^p -valued random variable S are **independent random variables** and where \mathbf{B} is invertible.

Remark: The ICA model holds true if X is Gaussian and if $\text{Var}(X)$ is full rank, in which case \mathbf{B} and S can be obtained using population PCA^b.

Unlike the factor analysis model, where the factor F has typically no physical existence, in ICA the variable S is a real signal that we want to recover from X . Remark that this is the reason why the matrix \mathbf{B} is assumed to be invertible.

A typical (toy) application of ICA is to recover the $p = 2$ signals S_1 and S_2 emitted by two persons speaking simultaneously in a room from the signals X_1 and X_2 recorded by two microphones.

^aThe main reference for this chapter is [?].

^bTo see this recall that $X = \mathbf{\Gamma}Y$ where $\mathbf{\Gamma} \in O(p)$ and where $Y \sim \mathcal{N}_p(0, \mathbf{L})$, with \mathbf{L} a diagonal matrix having the eigenvalues of $\text{Var}(X)$ as non-zero entries (see Chapter 1, page 24). If $\text{Var}(X)$ is full rank the matrix \mathbf{L} is invertible, and thus $X = (\mathbf{\Gamma}\mathbf{L}^{1/2})(\mathbf{L}^{-1/2}Y)$. This shows that the ICA model holds with $\mathbf{B} = \mathbf{\Gamma}\mathbf{L}^{1/2}$ and with $S = \mathbf{L}^{-1/2}Y \sim \mathcal{N}_p(0, \mathbf{I}_p)$ (recall that $S \sim \mathcal{N}_p(0, \mathbf{I}_p)$ implies that $S_1, \dots, S_p \stackrel{\text{iid}}{\sim} \mathcal{N}_1(0, 1)$).

ICA model: Discussion of its assumptions and problem formulation

- In (3.1) there is no loss of generality to assume that $\text{Var}(S) = \mathbf{I}_p$. Indeed, if $X = \tilde{\mathbf{B}}\tilde{S}$ with $\Psi := \text{Var}(\tilde{S}) \neq \mathbf{I}_p$ then, letting $\mathbf{B} = \tilde{\mathbf{B}}\Psi^{1/2}$ and $S = \Psi^{-1/2}\tilde{S}$, we have

$$X = \tilde{\mathbf{B}}\tilde{S} = (\tilde{\mathbf{B}}\Psi^{1/2})(\Psi^{-1/2}\tilde{S}) = \mathbf{B}S, \quad \text{Var}(S) = \mathbf{I}_p \quad (3.2)$$

implying that (3.1) holds.

Remark: (3.2) shows that $\text{Var}(\tilde{S})$ cannot be recovered from X if we only know that $X = \tilde{\mathbf{B}}\tilde{S}$ for some matrix $\tilde{\mathbf{B}} \in \mathbb{R}^{p \times p}$.

- Recalling that S is a true signal that we want to recover, the assumption that the components of S are independent is necessary to make \mathbf{B} identifiable (the identifiability of \mathbf{B} will be discussed more precisely later). In particular, if we only assume that the components of S are uncorrelated then S can only be recovered up to an orthogonal transformation since, for all $\mathbf{G} \in O(p)$, $X = \mathbf{B}S = (\mathbf{B}\mathbf{G}^\top)(\mathbf{G}S)$ where $\text{Var}(\mathbf{G}S) = \text{Var}(S)$.

We assume for now that $\text{Var}(X) = \mathbf{I}_p$. In this case, the ICA model (3.1) implies that

$$\text{Var}(X) = \text{Var}(\mathbf{B}S) = \mathbf{B}\mathbf{B}^\top = \mathbf{I}_p$$

showing that if (3.1) holds then \mathbf{B} must be an orthogonal matrix^a

Therefore, under the assumption $\text{Var}(X) = \mathbf{I}_p$ we have $X = \mathbf{B}S$ if and only if $S = \mathbf{B}^\top X$, and thus under (3.1) and the assumption $\text{Var}(X) = \mathbf{I}_p$ the matrix $\mathbf{B} \in O(p)$ is such that the components of the random variable $\mathbf{B}^\top X$ are **independent**.

^aRecall that the ICA model assumes that \mathbf{B} is invertible

Defining B through an optimization problem

For a given random variable Z we denote by $g_Z(z)dz$ its probability distribution (w.r.t. some reference measure dz) and we let

$$H(Z) = - \int g_Z(z) \log(g_Z(z)) dz.$$

Remark: The quantity $H(Z)$ is called the entropy of Z .

We also recall that the Kullback-Leibler (KL) divergence between the distributions $p(z)dz$ and $q(z)dz$ is given by

$$\text{KL}(p||q) = \int \log\left(\frac{p(z)}{q(z)}\right)p(z)dz$$

and we also recall the following result:

Lemma 3.1 *For any distributions $p(z)dz$ and $q(z)dz$ we have $\text{KL}(p||q) \geq 0$, where the equality holds if and only if $p = q$.*

Easy computations show that if $Z = (Z_1, \dots, Z_p)$ is an \mathbb{R}^p -valued random variable then the KL divergence between the distributions $g_Z(z)dz$ and $\prod_{j=1}^p g_{Z_j}(z_j)dz$ can be written as follows:

$$\text{KL}\left(g_Z || \prod_{j=1}^p g_{Z_j}\right) = \sum_{j=1}^p H(Z_j) - H(Z) =: I(Z).$$

By Lemma 3.1, $I(Z) \geq 0$ and $I(Z) = 0$ if and only if all the components of Z are independent. Therefore, $I(Z)$ can be interpreted as a measure of independence between the components of Z . (The quantity $I(Z)$ is called the mutual information of Z .)

Under the assumption $\text{Var}(X) = \mathbf{I}_p$, it follows that the matrix \mathbf{B} in (3.2) verifies

$$\mathbf{B} \in \underset{\mathbf{G} \in O(p)}{\text{argmin}} I(\mathbf{G}^\top X). \quad (3.3)$$

Two key lemmas for finding an approximate solution to (3.3)

Lemma 3.2 *Let Y be a continuous and real-valued random variable with $\mathbb{E}[Y] = 0$ and $\text{Var}(Y) = 1$, and let $Z \sim \mathcal{N}_1(0, 1)$. Then, $H(Z) \geq H(Y)$, where the equality holds if and only if $Y \sim \mathcal{N}_1(0, 1)$.*

Proof: We have

$$\begin{aligned}
 \text{KL}(p_Y || p_Z) &= \int \log \left(\frac{p_Y(y)}{p_Z(y)} \right) p_Y(y) dy = -H(Y) - \int \log(p_Z(y)) p_Y(y) dy \\
 &= -H(Y) - \int \left(-\frac{1}{2} \log(2\pi) - \frac{y^2}{2} \right) p_Y(y) dy \\
 &= -H(Y) - \int \left(-\frac{1}{2} \log(2\pi) - \frac{z^2}{2} \right) p_Z(z) dz \\
 &= -H(Y) - \int \log(p_Z(z)) p_Z(z) dz \\
 &= -H(Y) + H(Z)
 \end{aligned}$$

where the third equality uses the fact that $\mathbb{E}[Y^2] = \mathbb{E}[Z^2]$. Then, the result follows from Lemma 3.1. \square

To proceed further for any \mathbb{R} -valued random variable Y such that $\mathbb{E}[Y] = 0$ and $\text{Var}(Y) = 1$ we let $J(Y) = H(Z) - H(Y)$, with $Z \sim \mathcal{N}_1(0, 1)$.

Remark: By Lemma 3.2, for any \mathbb{R} -valued random variable Y such that $\mathbb{E}[Y] = 0$ and $\text{Var}(Y) = 1$ we have $J(Y) \geq 0$, where the equality holds if and only if $Y \sim \mathcal{N}_1(0, 1)$. Hence, the quantity $J(Y)$, called the **negentropy** of Y , is a **measure of distance to normality**.

Lemma 3.3 *Assume $\text{Var}(X) = \mathbf{I}_p$. Then, the matrix \mathbf{B} in (3.2) is such that $\mathbf{B} \in \text{argmax}_{\mathbf{G} \in O(p)} \sum_{j=1}^p J(g_{(j)}^\top X)$.*

Proof: Let $p_X(\cdot)$ be the density of X and let $Y = \mathbf{G}^\top X$. Then, using the change of variables formula and the fact that $|\det(\mathbf{G})| = 1$, the density $p_Y(\cdot)$ of Y is defined by $p_Y(y) = p_X(\mathbf{G}y)$ and thus

$$\begin{aligned}
 H(\mathbf{G}^\top X) = H(Y) &= - \int p_Y(y) \log(p_Y(y)) dy = - \int p_X(\mathbf{G}y) \log(p_X(\mathbf{G}y)) dy = - \int |\det(\mathbf{G}^{-1})| p_X(x) \log(p_X(x)) dx \\
 &= - \int p_X(x) \log(p_X(x)) dx \\
 &= H(X)
 \end{aligned}$$

where the third equality uses the change of variables formula for integrals. Then the result follows from (3.3) and

from the definition of $I(\mathbf{G}^\top X)$. \square

Approximating the solution to (3.3)

Using Lemma 3.2 and Lemma 3.3, it follows if $\text{Var}(X) = \mathbf{I}_p$ then the matrix \mathbf{B} in (3.2) verifies

$$\begin{aligned} \mathbf{B} &\in \operatorname{argmax}_{\mathbf{G} \in O(p)} \sum_{j=1}^p J(g_{(j)}^\top X) \\ &= \operatorname{argmax}_{\mathbf{G} \in O(p)} \left\{ \sum_{j=1}^p \text{“departure from Gaussianity of } g_{(j)}^\top X \text{”} \right\}. \end{aligned} \quad (3.4)$$

The quantity $J(g_{(j)}^\top X)$ is usually intractable but its interpretation in term of distance to normality provides a simple way to compute an approximate solution to (3.4): replace $J(g_{(j)}^\top X)$ by another measure of distance to normality!

In ICA, the departure from Gaussianity of $g_{(j)}^\top X$ is often measured by

$$\left(\mathbb{E}[\varphi(g_{(j)}^\top X)] - \mathbb{E}[\varphi(Z)] \right)^2, \quad Z \sim \mathcal{N}_1(0, 1) \quad (3.5)$$

for some function $\varphi : \mathbb{R} \rightarrow \mathbb{R}$.

Remark: Using (3.5) with the function $\varphi(u) = \frac{1}{a} \log \cosh(au)$ for some $a \in [1, 2]$, or with the function $\varphi(u) = -e^{-u^2/2}$, often works well in practice and provide a reasonable approximation of the negentropy $J(g_{(j)}^\top X)$ [?].

Then, for a given choice of $\varphi : \mathbb{R} \rightarrow \mathbb{R}$ and assuming $\text{Var}(X) = \mathbf{I}_p$, the matrix \mathbf{B} in (3.2) can be approximated by the matrix $\tilde{\mathbf{B}}$ verifying

$$\tilde{\mathbf{B}} \in \operatorname{argmax}_{\mathbf{G} \in O(p)} \sum_{j=1}^p \left(\mathbb{E}[\varphi(g_{(j)}^\top X)] - \mathbb{E}[\varphi(Z)] \right)^2 \quad (3.6)$$

where $Z \sim \mathcal{N}_1(0, 1)$.

Estimation of the approximate solution $\tilde{\mathbf{B}}$

A possible approach for estimating the matrix $\tilde{\mathbf{B}}$ defined in (3.6) from the observations $\{x_i\}_{i=1}^n$ is as follows. For all $\mathbf{G} \in O(p)$ let

$$L(\mathbf{G}) = \sum_{j=1}^p \left(\mathbb{E}[\varphi(g_{(j)}^\top X)] - \mathbb{E}[\varphi(Z)] \right)^2$$

and

$$\hat{L}_n(\mathbf{G}, \{x_i\}_{i=1}^n) = \sum_{j=1}^p \left(\frac{1}{n} \sum_{i=1}^n \varphi(g_{(j)}^\top x_i) - \mathbb{E}[\varphi(Z)] \right)^2$$

so that $\hat{L}_n(\mathbf{G}, \{x_i\}_{i=1}^n) \approx L(\mathbf{G})^a$.

Then, noting that $\tilde{\mathbf{B}} \in \operatorname{argmax}_{\mathbf{G} \in O(p)} L(\mathbf{G})$, we can estimate $\tilde{\mathbf{B}}$ using the matrix $\tilde{\mathbf{B}}_n$ defined by

$$\tilde{\mathbf{B}}_n \in \operatorname{argmax}_{\mathbf{G} \in O(p)} \hat{L}_n(\mathbf{G}, \{x_i\}_{i=1}^n). \quad (3.7)$$

The definition of $\hat{L}_n(\mathbf{G}, \{x_i\}_{i=1}^n)$ requires to compute $\mathbb{E}[\varphi(Z)]$, a quantity which, depending on the choice of φ , may be intractable. However, being a one dimensional integral, we can easily (and efficiently) estimate $\mathbb{E}[\varphi(Z)]$ using numerical integration methods.

Letting $\hat{\varphi}$ be an estimate of $\mathbb{E}[\varphi(Z)]$, a computable estimate $\tilde{\mathbf{B}}'_n$ of $\tilde{\mathbf{B}}$ is therefore defined by

$$\tilde{\mathbf{B}}'_n \in \operatorname{argmax}_{\mathbf{G} \in O(p)} \sum_{j=1}^p \left(\frac{1}{n} \sum_{i=1}^n \varphi(g_{(j)}^\top x_i) - \hat{\varphi} \right)^2. \quad (3.8)$$

^aThe set $O(p)$ being compact it follows that if $\{X_i\}_{i=1}^n$ are i.i.d. copies of X then, under some conditions on φ , we have $\sup_{\mathbf{G} \in O(p)} |\hat{L}_n(\mathbf{G}, \{X_i\}_{i=1}^n) - L(\mathbf{G})| \rightarrow 0$ in probability. Note that, whatever the distribution of X is, this uniform convergence result holds if φ is continuous and bounded, as this is the case for the choice $\varphi(u) = -e^{-u^2/2}$ mentioned earlier.

Estimation of the approximate solution $\tilde{\mathbf{B}}$ (end)

A simple way to solve the optimization problem (3.8) is to use a **projected gradient descend** algorithm:

A projected gradient descend algorithm for solving (3.8)

Input: Matrix $\tilde{\mathbf{B}}^{(0)} \in O(p)$ and step-size $\gamma > 0$

for $s \geq 1$ **do**

(i) Let $\mathbf{C}_s = \tilde{\mathbf{B}}^{(s-1)} + \gamma \nabla_{\mathbf{B}} \sum_{j=1}^p \left(\frac{1}{n} \sum_{i=1}^n \varphi(x_i^\top \tilde{\mathbf{b}}_{(j)}^{(s-1)}) - \hat{\varphi} \right)^2$

(ii) Let $\tilde{\mathbf{B}}^{(s)} = (\mathbf{C}_s \mathbf{C}_s^\top)^{-1/2} \mathbf{C}_s \in O(p)$.

if Convergence=TRUE **then**

(iii) **return** $\tilde{\mathbf{B}}^{(s)}$.

(iv) **break**

end if

end for

Remark: In practice, more sophisticated (and more efficient) methods are used to estimate the matrix $\tilde{\mathbf{B}}$ defined in (3.6) [see ? , Section 6].

Remark: Estimating \mathbf{B} through the estimation of the matrix $\tilde{\mathbf{B}}$ is called the FastICA method.

Estimation of \mathbf{B} when $\text{Var}(X) \neq \mathbf{I}_d$

We now consider the general case where $\text{Var}(X) = \mathbf{\Sigma}$ for some $\mathbf{\Sigma} \neq \mathbf{I}_p$. We assume below that $\mathbf{\Sigma}$ is full rank and let \mathbf{L} be a diagonal matrix containing the eigenvalues of $\mathbf{\Sigma}$ and $\mathbf{\Gamma} \in O(p)$ be the corresponding matrix of orthonormal eigenvectors.

Let $\mathbf{K} = \mathbf{L}^{-1/2}\mathbf{\Gamma}^\top$, $\mathbf{B}_W = \mathbf{K}\mathbf{B}$ and $W = \mathbf{K}X$, and note that

- (i) Under the ICA model the random variable W is such that $W = (\mathbf{K}\mathbf{B})S = \mathbf{B}_W S$ and such that

$$\text{Var}(W) = \mathbf{L}^{-1/2}\mathbf{\Gamma}^\top \mathbf{\Sigma} \mathbf{\Gamma} \mathbf{L}^{-1/2} = \mathbf{L}^{-1/2}\mathbf{\Gamma}^\top (\mathbf{\Gamma} \mathbf{L} \mathbf{\Gamma}^\top) \mathbf{\Gamma} \mathbf{L}^{-1/2} = \mathbf{I}_p.$$

- (ii) Since $\mathbf{B}_W = \mathbf{K}\mathbf{B}$ we have $\mathbf{B} = \mathbf{K}^{-1}\mathbf{B}_W = \mathbf{\Gamma} \mathbf{L}^{1/2} \mathbf{B}_W$.

Let $\mathbf{\Lambda}$ be the diagonal matrix containing the eigenvalues of the matrix $\mathbf{S} := \frac{1}{n}\mathbf{X}^\top \mathbf{X}$, $\mathbf{A} \in O(p)$ be the corresponding matrix of orthonormal eigenvectors and $\mathbf{Y} = \mathbf{X}\mathbf{A}$ be the matrix of principal components.

Then, (i)-(ii) suggest the following three steps for estimating \mathbf{B} :

1. Compute $\mathbf{W} = \mathbf{Y}\mathbf{\Lambda}^{-1/2}$. Remark that $\frac{1}{n}\mathbf{W}^\top \mathbf{W} = \mathbf{I}_p$ and recall that (under some conditions) we can interpret $\{w_i\}_{i=1}^n$ as “approximate” realizations of W (see Chapter 1, pages 24–25). The transformation $\mathbf{X} \mapsto \mathbf{W}$ of the data is called **whitening**.
2. Use $\{w_i\}_{i=1}^n$ and the approach introduced in this chapter for estimating \mathbf{B} when $\text{Var}(X) = \mathbf{I}_p$ to compute an estimate $\tilde{\mathbf{B}}'_{n,W}$ of \mathbf{B}_W .
3. Estimate \mathbf{B} by

$$\hat{\mathbf{B}}_n := \mathbf{A}\mathbf{\Lambda}^{1/2} \tilde{\mathbf{B}}'_{n,W}. \quad (3.9)$$

Recovering the signals from the data

Recall that under the ICA model the observations $\{x_i\}_{i=1}^n$ are realizations of a random variable X such that $X = \mathbf{B}S$ with \mathbf{B} and S as in (3.1).

Under the ICA model we therefore have $x_i = \mathbf{B}s_i$ for all i and thus, denoting by \mathbf{S}_{sig} the $n \times p$ matrix with rows $\{s_i\}_{i=1}^n$,

$$\mathbf{X} = \mathbf{S}_{\text{sig}} \mathbf{B}^\top.$$

Given the estimate $\hat{\mathbf{B}}_n$ of \mathbf{B} defined in (3.9) we can estimate the matrix signals \mathbf{S}_{sig} using

$$\hat{\mathbf{S}}_{\text{sig}} := \mathbf{X} (\hat{\mathbf{B}}_n^\top)^{-1}.$$

Recalling that $\tilde{\mathbf{B}}'_{n,W} \in O(p)$, we have

$$\begin{aligned} \hat{\mathbf{S}}_{\text{sig}} &= \mathbf{X} (\hat{\mathbf{B}}_n^\top)^{-1} = \mathbf{X} \left((\tilde{\mathbf{B}}'_{n,W})^\top \mathbf{\Lambda}^{1/2} \mathbf{A}^\top \right)^{-1} \\ &= \mathbf{X} \mathbf{A} \mathbf{\Lambda}^{-1/2} \tilde{\mathbf{B}}'_{n,W} \\ &= \mathbf{W} \tilde{\mathbf{B}}'_{n,W} \end{aligned}$$

and therefore computing the estimate $\hat{\mathbf{S}}_{\text{sig}}$ does not require to compute the estimate $\hat{\mathbf{B}}_n$ of \mathbf{B} .

Remark: Estimating \mathbf{S}_{sig} by $\hat{\mathbf{S}}_{\text{sig}} = \mathbf{W} \tilde{\mathbf{B}}'_{n,W}$ is natural. Indeed, under the ICA model we have $W = \mathbf{B}_W S$ and thus, as we are interpreting the w_i 's as realizations of W , this implies that $w_i = \mathbf{B}_W s_i$ for all i , and thus that $\mathbf{W} = \mathbf{S}_{\text{sig}} \mathbf{B}_W^\top \Leftrightarrow \mathbf{S}_{\text{sig}} = \mathbf{W} \mathbf{B}_W$.

Identifiability of the ICA model

Assume that $\text{Var}(X) = \mathbf{I}_p$ and that the ICA model is correct, so that for some $\mathbf{B} \in O(p)$ we have $\mathbf{X} = \mathbf{S}_{\text{sig}} \mathbf{B}^\top$ where the rows of \mathbf{S}_{sig} are n realizations of S .

Then,

- If all the components of S are $\mathcal{N}_1(0, 1)$ random variables then $\mathbf{G}\mathbf{S} \stackrel{\text{dist.}}{=} S$ for any matrix $\mathbf{G} \in O(p)$ ^a. Therefore, in this case we have

$$X = \mathbf{B}\mathbf{S} \stackrel{\text{dist.}}{=} \mathbf{B}\mathbf{G}\mathbf{S} = (\mathbf{B}\mathbf{G})\mathbf{S}, \quad \mathbf{B}\mathbf{G} \in O(p), \quad \forall \mathbf{G} \in O(p)$$

showing that the matrix \mathbf{B} , and thus the signals \mathbf{S}_{sig} , can only be estimated up to an orthogonal transformation.

- The columns of \mathbf{B} , and thus the rows of \mathbf{S}_{sig} , can only be estimated up to a multiplicative sign. To see this let \mathbf{D} be a $p \times p$ diagonal matrix with such that $d_{jj} \in \{-1, 1\}$ for all j and let $\tilde{S} = \mathbf{D}\mathbf{S}$. Then,

$$X = \mathbf{B}\mathbf{S} = \mathbf{B}\mathbf{D}^{-1}\mathbf{D}\mathbf{S} = (\mathbf{B}\mathbf{D})\tilde{S}$$

where $\mathbf{B}\mathbf{D} \in O(p)$ and $\text{Var}(\tilde{S}) = \mathbf{I}_d$.

Remark: It can be shown that if at most one component of S is Gaussian and $\text{rank}(\mathbf{B}) = p$ then the columns of the matrix $\mathbf{B} \in O(p)$ are unique up to a multiplicative sign (see [?]).

^aUse the change of variable formula to show this.

Illustration of ICA

We let $p = 3$, $n = 1\,000$ and simulate the true signals matrix $\mathbf{S}_{\text{sig}} = [s_{ij}] \in \mathbb{R}^{n \times p}$ using

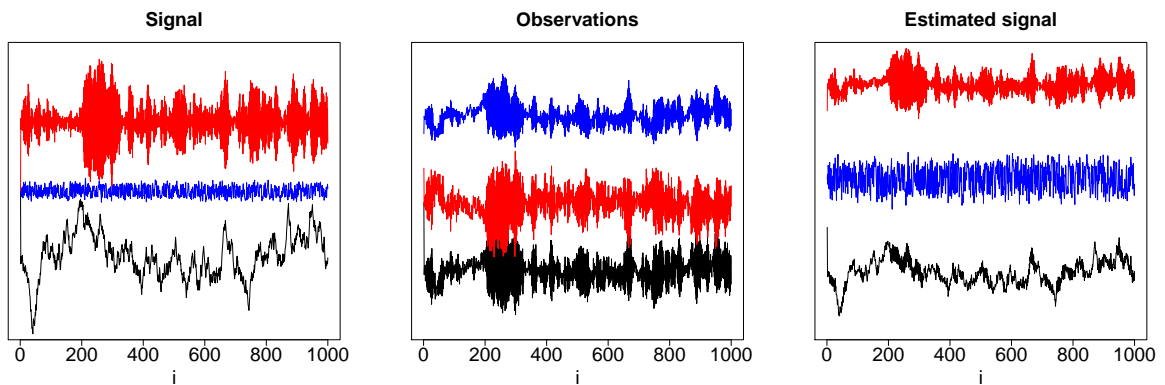
$$S_{ij} = \rho_j S_{(i-1)j} + \epsilon_{ij}, \quad \epsilon_{ij} \stackrel{\text{iid}}{\sim} \frac{1}{2} \mathcal{N}_1(-1, 0.25) + \frac{1}{2} \mathcal{N}_1(1, 0.25)$$

where $S_{(i-1)j} = 0$ for $i = 1$ and where $\rho_2 = -\rho_1 = 0.98$ and $\rho_3 = 0.2$. Remark that each row of the matrix \mathbf{S}_{sig} is a trajectory of a Markov chain. The three Markov chains (i.e. the three signals) are mixed using the matrix

$$\mathbf{B} = \begin{pmatrix} 1 & -1 & -3 \\ 1 & 1 & 2 \\ -1 & 3 & -3 \end{pmatrix}.$$

Remark: In this example the distribution of S_i is not exactly the same for all i .

The true signals \mathbf{S}_{sig} as well as the observations $\mathbf{X} = \mathbf{S}_{\text{sig}} \mathbf{B}^\top$ and the ICA estimate $\hat{\mathbf{S}}_{\text{sig}}$ of \mathbf{S}_{sig} (obtained with $\varphi(u) = \log \cosh(u)$) are shown in the following plots.



Remark: In the two plots the signals are shifted and coloured to facilitate the comparisons. In addition, the y -axis has been removed since, as shown above, we cannot estimate the variance of the components of \mathbf{S} .

References

- [1] Bishop, C. M. (2006). Pattern recognition. *Machine learning*, 128(9).
- [2] Mardia, K., Kent, J., and Bibby, J. (1979). *Multivariate analysis*. Probability and mathematical statistics. Academic Press Inc.