Library-Based Linear Unmixing for Hyperspectral Imagery via Reversible Jump MCMC Sampling

Nicolas Dobigeon^{1,2} and Jean-Yves Tourneret²

¹ University of Michigan, Department of EECS, Ann Arbor, MI 48109-2122, USA
² University of Toulouse, IRIT/INP-ENSEEIHT, BP 7122, 31071 Toulouse cedex 7, FRANCE {nicolas.dobigeon, jean-yves.tourneret}@enseeiht.fr

Abstract—This paper studies a semi-supervised algorithm for linear hyperspectral unmixing. The proposed unmixing method assumes that the pure material spectra denoted as endmembers belong to a library that is a priori available. However, the number and the nature of endmembers appearing in the pixel are not known a priori, resulting in a model selection problem. This paper proposes to handle this model selection problem within a fully Bayesian framework. First, appropriate distributions are elected as prior distributions for the unknown parameters. Particularly, a distribution defined on a simplex is chosen as prior for an appropriate partial abundance vector to ensure the positivity and the sum-to-one constraints of the mixing coefficients. Due to the complexity of the posterior distribution, a reversible jump Markov chain Monte Carlo algorithm is proposed to estimate the number and the nature of the macroscopic materials, as well as their respective proportions in the pixel. The accuracy of the proposed method is illustrated by simulations on synthetic hyperspectral data.

1. INTRODUCTION

For few decades, hyperspectral imagery has been receiving increasing interest in the geoscience and image processing literatures [1]. Indeed, current spectro-imagers, such as AVIRIS, CASI or Hyperion, are able to acquire a same scene in several hundreds of contiguous bands, providing spectrally resolved images. Hyperspectral imagery has demonstrated its interest for numerous applications. These applications include geologic cartography [2], environmental [3] or military monitoring [4]. However, to benefit from the amount of information contained in these images, innovative strategies have to be developed to improve the accuracy of standard estimation, detection and classification algorithms [5]. One of the main steps in the analysis of such images consists of identifying the pure materials, or endmembers that are present in the observed scene. Then, an inversion step estimates their respective proportions in each pixel. This paper introduces a semi-supervised algorithm that can be used to perform this so-called spectral unmixing. In the literature, two models have been investigated to describe how endmember mixing is achieved. According to the first model, the observed pixel spectrum is assumed to be related to the unobserved material spectra via a linear mixture whose coefficients are the concentrations, or *abundances*, of the macroscopic components in the pixel [6]. The second model introduced in [7] refers to an intimate model that assumes a non-linear combination of the endmembers. However, as noted in [8], this intimate model can be easily linearized. Consequently, this paper concentrates on the Linear Mixing Model (LMM) that has received considerable attention in the literature [9], [10].

The proposed unmixing method assumes that the pure material spectra belong to a library that is a priori available. In our work, the number and the nature of endmembers appearing in the pixel are not known a priori. Estimating the number and the nature of the endmembers involved in the mixture is a classical model selection task. In this paper, this challenging problem will be handled within a fully Bayesian framework. Due to obvious physical reasons, the abundance vectors have to satisfy positivity and sum-to-one constraints. The Bayesian estimation procedure proposed here allows one to incorporate these constraints easily into the model. First, appropriate distributions are elected as prior distributions for the unknown parameters. Particularly, a distribution defined on a simplex is chosen as prior for an appropriate partial abundance vector to ensure the positivity and the additivity of the mixing coefficients, as in [11]. Because of the complexity of the posterior distribution, deriving standard Bayesian estimators remains difficult. In such cases, stochastic simulation methods can be used to overcome the difficulty. In this paper, a Markov chain Monte Carlo (MCMC) algorithm is proposed to generate samples asymptotically distributed according to the posterior distribution of the abundances and endmembers. Then, the generated samples will be used to approximate the Bayesian estimators. More precisely, as the number of endmembers is not a priori known, the unknown full parameter vector belongs to a space with unknown dimension. A reversible jump MCMC algorithm will allow ones to explore spaces with different dimensions [12].

This paper is organized as follows. The linear unmixing problem for hyperspectral images is formulated in Section 2. The hierarchical Bayesian model associated to this problem is described in Section 3. Section 4 details the reversible jump MCMC algorithm that generates samples distributed according to the abundance and endmember distribution. Simulation results obtained from synthetic hyperspectral data are presented in Section 5. Conclusions are reported in Section 6.

^{978-1-4244-2622-5/09/\$25.00 ©2009} IEEE

2. PROBLEM FORMULATION

In this section, we describe the observation model which is classically assumed to perform spectral unmixing of hyperspectral pixels. As noticed in the previous section, this paper focuses on the LMM that constitutes a good approximation in the reflective domain ranging from 0.4μ m to 2.5μ m (see [5] and [9]). The LMM assumes that the spectrum $\boldsymbol{y} = [y_1, \dots, y_L]^T$ of a mixed pixel is a linear combination of R endmember spectra \boldsymbol{m}_r corrupted by additive noise:

$$\boldsymbol{y} = \sum_{r=1}^{R} \boldsymbol{m}_r \boldsymbol{\alpha}_r + \boldsymbol{n}, \qquad (1)$$

where L is the number of available spectral bands of the pixel, $\boldsymbol{m}_r = [\boldsymbol{m}_{r,1}, \ldots, \boldsymbol{m}_{r,L}]^T$ is the spectrum of the rth material, α_r is the fraction of the rth material in the pixel and R is the number of pure materials (or *endmembers*) present in the observed pixel. Here we consider $\boldsymbol{n} = [n_1, \ldots, n_L]^T$ as an additive white noise sequence which is classically assumed to be an independent and identically distributed (i.i.d.) Gaussian sequence with variance σ^2 , denoted as $\boldsymbol{n} \sim \mathcal{N}(\boldsymbol{0}_L, \sigma^2 \mathbf{I}_L)$, where \mathbf{I}_L and $\boldsymbol{0}_L$ are the identity matrix of dimension $L \times L$ and the $L \times 1$ vector made of 0's, respectively. Note that more complicated noise models that take into account some correlations might be studied (see for instance [13]).

Due to physical reasons, the abundance vector $\boldsymbol{\alpha}(R) = [\alpha_r, \dots, \alpha_R]^T$ has to ensure the following non-negativity and sum-to-one constraints:

$$\begin{cases} \alpha_r \ge 0, \ \forall r = 1, \dots, R, \\ \sum_{r=1}^R \alpha_r = 1. \end{cases}$$
(2)

It is important to mention here that the number R of endmembers involved in the mixture is not known *a priori*. However, the endmember spectra m_r (r = 1, ..., R) belong to a spectral library

$$\mathcal{S} = \{\boldsymbol{s}_1, \dots, \boldsymbol{s}_{R_{\max}}\},\tag{3}$$

composed of R_{max} materials. These endmember spectra s_r $(r = 1, \ldots, R_{\text{max}})$ have been for instance identified by one of the numerous endmember extraction algorithms such as N-FINDR [14], MVT [15] or ICE [16].

The problem addressed in this paper consists of estimating the number of endmembers R, the spectra m_r (r = 1, ..., R), the fraction coefficient α_r (r = 1, ..., R) and the noise variance σ^2 under the constraints in (2) given the observed pixel y and the spectral library S.

3. HIERARCHICAL BAYESIAN MODEL

Likelihood

The LMM (1) and the statistical properties of the noise vector yield $\boldsymbol{y}|R, \mathbf{M}, \boldsymbol{\alpha}, \sigma^2 \sim \mathcal{N}(\mathbf{M}\boldsymbol{\alpha}, \sigma^2 \mathbf{I}_L)$, where $\mathbf{M} = [\boldsymbol{m}_1, \dots, \boldsymbol{m}_R]$. Consequently, the likelihood function of \boldsymbol{y} can be expressed as:

$$f(\boldsymbol{y}|R,\boldsymbol{\alpha},\mathbf{M},\sigma^2) = \left(\frac{1}{2\pi\sigma^2}\right)^{\frac{L}{2}} \exp\left[-\frac{\|\boldsymbol{y}-\mathbf{M}\boldsymbol{\alpha}\|^2}{2\sigma^2}\right],$$
(4)

where $\|\boldsymbol{x}\| = \sqrt{\boldsymbol{x}^T \boldsymbol{x}}$ is the standard ℓ_2 norm.

Parameter priors

Prior for the number of endmembers — A discrete uniform distribution on $[1, ..., R_{max}]$ is chosen for the prior associated to the number of mixture components R:

$$f(R) = \frac{1}{R_{\max}}, \quad R = 1, \dots, R_{\max}.$$
 (5)

Endmember matrix prior — Conditionally upon the number R of materials involved in the mixture, all combinations of R spectra belonging to the library S are assumed to be equiprobable:

$$f(\mathbf{M} \mid R) = {\binom{R_{\max}}{R}}^{-1} = \frac{\Gamma(R+1)\Gamma(R_{\max}-R+1)}{\Gamma(R_{\max}+1)},$$
(6)

where $\Gamma(\cdot)$ denotes the Gamma function.

Abundance prior — The abundance vector can be written as $\boldsymbol{\alpha} = \begin{bmatrix} \boldsymbol{c}^T, \alpha_R \end{bmatrix}^T$ with $\boldsymbol{c} = [\alpha_1, \dots, \alpha_{R-1}]^T$ and $\alpha_R = 1 - \sum_{r=1}^{R-1} \alpha_r$. The constraints (2) inherent to the LMM impose that \boldsymbol{c} belongs to the simplex \mathbb{S} defined by:

$$\mathbb{S}_{R} = \left\{ \boldsymbol{c} \middle| c_{r} \ge 0, \ \forall r = 1, \dots, R-1, \ \sum_{r=1}^{R-1} c_{r} \le 1 \right\}.$$
(7)

To reflect the absence of prior knowledge regarding the abundance vector, a uniform distribution on \mathbb{S} is chosen as prior distribution for *c*:

$$f(\boldsymbol{c}|R) = \frac{1}{\operatorname{vol}(\mathbb{S}_{\mathrm{R}})} \mathbf{1}_{\mathbb{S}_{R}}(\boldsymbol{c}), \qquad (8)$$

where $\mathbf{1}_{\mathbb{S}_R}(\cdot)$ stands for the indicator function defined on \mathbb{S}_R :

$$\mathbf{1}_{\mathbb{S}_R}(\boldsymbol{c}) = \begin{cases} 1, & \text{if } \boldsymbol{c} \in \mathbb{S}_R; \\ 0, & \text{otherwise.} \end{cases}$$
(9)

Noise variance prior — A conjugate inverse Gamma distribution is chosen as prior distribution for σ^2 :

$$\sigma^2 \mid \nu, \gamma \sim \mathcal{IG}\left(\frac{\nu}{2}, \frac{\gamma}{2}\right),\tag{10}$$

where ν will be fixed to $\nu = 2$ (as in [17]) and γ is an adjustable hyperparameter.

Hyperparameter prior distribution

The accuracy of the unmixing procedure may depend on the value of the unknown hyperparameter γ . As in [11], the hierarchical model proposed here uses a non-informative Jeffreys' prior for γ :

$$f(\gamma) \propto \frac{1}{\gamma} \mathbf{1}_{\mathbb{R}^+}(\gamma).$$
 (11)

Posterior distribution

The posterior distribution of the unknown parameter vector $\boldsymbol{\theta} = \{R, \mathbf{M}, \boldsymbol{\alpha}, \sigma^2\}$ can be computed from the following hierarchical structure:

$$f(\boldsymbol{\theta}|\boldsymbol{y}) \propto \int f(\boldsymbol{y}|\boldsymbol{\theta}) f(\boldsymbol{\theta}|\boldsymbol{\gamma}) f(\boldsymbol{\gamma}) d\boldsymbol{\gamma},$$
 (12)

where \propto means "proportional to", $f(\boldsymbol{y}|\boldsymbol{\theta})$ and $f(\gamma)$ have been defined in (4) and (11). Conditionally upon R, by assuming the prior independence between σ^2 , \boldsymbol{c} and \mathbf{M} i.e. $f(\boldsymbol{\theta}|\gamma) = f(R) f(\boldsymbol{c}|R) f(\mathbf{M}|R) f(\sigma^2|\nu,\gamma)$, the hyperparameter γ can be integrated out from the joint distribution $f(\boldsymbol{\theta},\gamma|\boldsymbol{y})$, yielding:

$$f\left(R, \boldsymbol{c}, \mathbf{M}, \sigma^{2} | \boldsymbol{y}\right) \propto \frac{1}{\sigma^{L+2}} \exp\left[-\frac{\|\boldsymbol{y} - \mathbf{M}\boldsymbol{\alpha}\|^{2}}{2\sigma^{2}}\right] \times \frac{1}{\operatorname{vol}\left(\mathbb{S}_{\mathrm{R}}\right)} f\left(R\right) f\left(\mathbf{M} | R\right) \mathbf{1}_{\mathbb{S}_{R}}(\boldsymbol{c}), \quad (13)$$

where the dimensions of \mathbf{M} and $\boldsymbol{\alpha}$ depend on the unknown parameter R. The next section shows that an appropriate hybrid Gibbs sampling strategy allows one to generate samples distributed according to the joint distribution $f(R, \boldsymbol{c}, \mathbf{M}, \sigma^2 | \boldsymbol{y})$.

4. REVERSIBLE JUMP MCMC ALGORITHM

This section studies an hybrid Metropolis-within-Gibbs algorithm that samples according to $f(R, c, \mathbf{M}, \sigma^2 | \mathbf{y})$. The vectors to be sampled belong to a space whose dimension depends on R, requiring to use a dimension matching strategy as in [12], [18], [19]. More precisely, the proposed algorithm consists of three different moves:

- 1. updating the endmember spectrum matrix M,
- 2. updating the endmember abundance vector c,
- 3. updating the noise variance σ^2 .

These three moves are scanned systematically as in [12]. They are detailed below.

Updating the endmember spectrum matrix M

The endmember spectra involved in the mixture (1) are updated by using three kinds of move. These moves are called *birth, death* and *switch* moves, as in [20, p. 53]. The first two moves consist of increasing or decreasing the number of pure components by 1. These moves allow the sampler to explore sets with different dimensions. Thus, they require the use of the reversible jump MCMC algorithm [12]. The third move does not affect the model dimension and requires a standard Metropolis-Hastings algorithm.

Assume that at iteration t of the sampler, the current model is defined by the vector $\boldsymbol{\theta}^{(t)} = \{R^{(t)}, \mathbf{M}^{(t)}, \boldsymbol{c}^{(t)}, \sigma^{2(t)}\}$. The three moves are defined as follows.

Birth — A birth move is proposed with the probability $b_{R^{(t)}}$. The current number of endmember $R^{(t)}$ is increased by 1

$$R^{\star} = R^{(t)} + 1.$$

A new spectrum m^* is randomly chosen with equal probability¹ among the available materials of the spectral library S. The new endmember matrix M^* is:

$$\mathbf{M}^{\star} = \left[\mathbf{M}^{(t)}, \boldsymbol{m}^{\star}
ight].$$

A new corresponding abundance has to be drawn ensuring the constraints in (2). This coefficient $\alpha^{(t)}$ is proposed according a rule inspired by [18]:

- draw a new abundance coefficient w^* from the Beta distribution $\mathcal{B}e(1, R^{(t)})$,
- re-scale the existing fractions so that all of them sum to 1:

$$\alpha_r^{\star} = \alpha_r^{(t)} \left(1 - w^{\star} \right), \ r = 1, \dots, R^{(t)}, \tag{14}$$

• set
$$\boldsymbol{\alpha}^{\star} = [\alpha_1^{\star}, \dots, \alpha_R^{\star}, w^{\star}]^T$$
.

Death — A death move is proposed with the probability $d_{R^{(t)}}$. The current number of endmember $R^{(t)}$ is decreased by 1:

$$R^{\star} = R^{(t)} - 1.$$

One of the spectra chosen with equal probability in the endmember matrix $\mathbf{M}^{(t)}$ is removed, as well as the corresponding abundance coefficient. The remaining abundances coefficients are re-scaled to sum to 1.

Switch — A switch move is proposed with the probability $u_{R^{(t)}}$. A spectrum randomly chosen in $\mathbf{M}^{(t)}$ is replaced by another spectrum randomly chosen in the library S.

At iteration t of the sampler, one of the moves *birth*, *death* and *switch* is randomly chosen with the probabilities $b_{R^{(t)}}$, $d_{R^{(t)}}$ and $u_{R^{(t)}}$, with $b_{R^{(t)}} + d_{R^{(t)}} + u_{R^{(t)}} = 1$. Of course, the death move is not allowed for $R^{(t)} = 1$ and the birth move is impossible for $R^{(t)} = R_{\max}$, i.e. $d_1 = b_{R_{\max}} = 0$. The different moves are assumed to be equiprobable such that $b_{R^{(t)}} = d_{R^{(t)}} = u_{R^{(t)}} = \frac{1}{3}$ for $R^{(t)} \in \{1, \dots, R_{\max} - 1\}$ and $b_1 = u_1 = d_{R_{\max}} = u_{R_{\max}} = \frac{1}{2}$. The acceptance rates of the *birth* and *death* moves are $\rho = \min\{1, A_b\}$ and $\rho =$

¹Of course, when prior knowledge regarding the probability of each spectrum to be in the image is available, this information can be used to randomly select a spectrum of the library.

 $\min\left\{1, A_{\rm b}^{-1}\right\}$, respectively, with (see [11] for more details):

$$A_{\rm b} = \exp\left[-\frac{\|\boldsymbol{y} - \mathbf{M}^{\star} \boldsymbol{\alpha}^{\star}\|^{2} - \|\boldsymbol{y} - \mathbf{M}^{(t)} \boldsymbol{\alpha}^{(t)}\|^{2}}{2}\right] \\ \times \frac{d_{R^{(t)}+1}}{b_{R^{(t)}}} \frac{1}{g_{1,R^{(t)}} (w^{\star})} (1 - w^{\star})^{R^{(t)}-1}$$
(15)
$$\times \frac{\Gamma\left(\delta R^{(t)} + \delta\right)}{\Gamma\left(\delta R^{(t)}\right) \Gamma\left(\delta\right)} w^{\star\delta-1} (1 - w^{\star})^{(\delta-1)R^{(t)}},$$

with $\delta = 1$.

The acceptance probability for the *switch* move is the standard Metropolis Hastings ratio $\rho = \min \{1, A_s\}$ with

$$A_{\rm s} = \exp\left[-\frac{\|\boldsymbol{y} - \mathbf{M}^{\star}\boldsymbol{\alpha}^{\star}\|^2 - \|\boldsymbol{y} - \mathbf{M}^{(t)}\boldsymbol{\alpha}^{(t)}\|^2}{2}\right].$$
 (16)

Updating the endmember abundance vector c

Updating the endmember abundance vector c is classically conducted by generating samples $c^{(t)}$ according to the posterior conditional distribution $f(c|R, \mathbf{M}, \sigma^2, y)$. By denoting \mathbf{M}_{-R} the matrix \mathbf{M} whose Rth column has been removed, i.e. $\mathbf{M}_{-R} = [m_1, \dots m_{R-1}]$, this conditional distribution is

$$f(\boldsymbol{c}|\boldsymbol{R}, \mathbf{M}, \sigma^{2}, \boldsymbol{y}) \propto \exp\left[-\frac{(\boldsymbol{c}-\boldsymbol{\mu})^{T} \boldsymbol{\Lambda}^{-1} (\boldsymbol{c}-\boldsymbol{\mu})}{2}\right] \mathbf{1}_{\mathbb{S}}(\boldsymbol{\alpha}), \quad (17)$$

where

$$\begin{cases} \mathbf{\Lambda} = \left[\frac{1}{\sigma^2} \left(\mathbf{M}_{-R} - \boldsymbol{m}_R \mathbf{u}^T\right)^T \left(\mathbf{M}_{-R} - \boldsymbol{m}_R \mathbf{u}^T\right)\right]^{-1}, \\ \boldsymbol{\mu} = \mathbf{\Lambda} \left[\frac{1}{\sigma^2} \left(\mathbf{M}_{-R} - \boldsymbol{m}_R \mathbf{u}^T\right)^T \left(\boldsymbol{y} - \boldsymbol{m}_R\right)\right], \end{cases}$$
(18)

with $\mathbf{u} = [1, \dots, 1]^T \in \mathbb{R}^{R-1}$. This distribution indicates that $\boldsymbol{c}|R, \mathbf{M}, \sigma^2, \boldsymbol{y}$ is distributed according to a truncated Gaussian distribution:

$$\boldsymbol{c}|\boldsymbol{R}, \mathbf{M}, \sigma^2, \boldsymbol{y} \sim \mathcal{N}_{\mathbb{S}}(\boldsymbol{\mu}, \boldsymbol{\Lambda}).$$
 (19)

As in [11], sampling according to a truncated Gaussian distribution can be achieved using a standard accept-reject procedure, when the number of endmembers is relatively small (as in the examples studied in this paper). More efficient simulation techniques based on Gibbs moves can be used for high dimension problems (see [21] or [22] for more details).

Updating the noise variance σ^2

Integrating out the hyperparameter γ from the joint posterior distribution $f(\theta, \gamma | y)$ leads to:

$$\sigma^2 | R, \boldsymbol{\alpha}, \mathbf{M}, \boldsymbol{y} \sim \mathcal{IG}\left(\frac{L}{2}, \frac{\|\boldsymbol{y} - \mathbf{M}\boldsymbol{\alpha}\|^2}{2}\right),$$
 (20)

from which it is straightforward to sample.

5. SIMULATION RESULTS

This section presents simulation results conducted on a synthetic mixed pixel. This pixel results from the combination of three endmembers (green grass, bare red brick, galvanized steel metal) with the abundance vector $[0.4, 0.2, 0.4]^T$. The observation is corrupted by an additive Gaussian noise with signal to noise ratio SNR = 20dB. A typical resulting mixed pixel is depicted in Fig. 1 (below).



Figure 1. Top: endmember spectra: green grass (solid line), bare red brick (dotted line), galvanized steel metal (dashed line). Bottom: resulting spectrum of the mixed pixel (SNR = 20dB).

This simulation considers a spectrum library composed of 6 endmembers that are provided by the ENVI software [23, p. 1035]. These spectra, representative of a urban or suburban environment, are depicted in Fig. 2.



Figure 2. Endmember spectra of the library.

The results presented here are obtained for $N_{\rm MC} = 20000$ iterations, including $N_{\rm bi} = 200$ burn-in iterations. The first step of the analysis consists of identifying the different combinations of endmembers in the matrices $\mathbf{M}^{(t)}$ generated by the sampler. Among the $N_r = N_{\rm MC} - N_{\rm bi}$ matrices $\mathbf{M}^{(t)}$ drawn by the sampler, 7 different combinations of spectra from the library S have been generated. The probabilities of each of these combination occurrence, depicted in Fig. 3, are clearly in favor of the endmember matrix $\mathbf{M} = [\mathbf{s}_2 \mathbf{s}_5 \mathbf{s}_6]$, which is in good agreement with the mixing scenario. Indeed, the maximum probability $P[\mathbf{M} = [\mathbf{s}_2, \mathbf{s}_5, \mathbf{s}_6] \mid \mathbf{y}] = 0.98$ corresponds to the actual spectra involved in the mixture.



Figure 3. Posterior probabilities of occurrence of the endmember combinations.

The posterior distributions of the corresponding abundances are finally estimated following the MMSE principle:

$$\hat{\boldsymbol{c}}_{\text{MMSE}} = \frac{1}{N_r} \sum_{t=1}^{N_r} \boldsymbol{c}^{(N_{\text{bi}}+t)}.$$
 (21)

Conditionally upon the maximum a posteriori estimate of the endmember matrix $\hat{\mathbf{M}}_{\text{MAP}} = [s_2, s_5, s_6]$, the posterior distributions of the corresponding abundance coefficients are depicted in Fig. 4. These posteriors are clearly in good agreement with the actual values of the abundances $\boldsymbol{\alpha} = [0.4, 0.2, 0.4]^T$.

The performance of the proposed algorithm is compared with a fully supervised Bayesian approach similar to [11]. First, 100 synthetic signals have been generated as described above: $\boldsymbol{\alpha} = [0.4, 0.2, 0.4]^T$, $\mathbf{M} = [\boldsymbol{s}_2, \boldsymbol{s}_5, \boldsymbol{s}_6]$, SNR = 20dB. The supervised algorithm assumes that the 6 endmember spectra of the library (depicted in Fig. 2) are involved in the mixture. Hence, with this approach, the vector to be estimated is $[0.0, 0.4, 0.0, 0.0, 0.2, 0.4]^T$. The mean squared errors (MSE) for the supervised algorithm and our library-based algorithm are reported in Table 1. These results show that the proposed semi-supervised approach outperforms the supervised algorithm.

6. CONCLUSIONS

This paper presented a hierarchical Bayesian model for hyperspectral linear unmixing. This model relied on appropriate prior distributions chosen for the unknown parameters.



Figure 4. Posterior distribution of the estimated abundances $\boldsymbol{\alpha} = [\alpha_1, \alpha_2, \alpha_3]^T$ conditioned upon $\mathbf{M} = [\boldsymbol{s}_2, \boldsymbol{s}_5, \boldsymbol{s}_6]$.

 Table 1. MSE for supervised and semi-supervised algorithms

	Supervised	Semi-supervised
MSE	$5.4 imes 10^{-2}$	$4.7 imes 10^{-2}$

It provided the posterior distribution of the abundance coefficients that can be used for their estimation. Assuming that the endmembers belong to a known spectral library, a reversible-jump MCMC method allowed one to identify the macroscopic materials actually involved in the mixture. In that sense, it constituted an original semi-supervised algorithm for linear unmixing of hyperspectral images.

7. REFERENCES

- C.-I Chang, Ed., *Hyperspectral data exploitation. The*ory and Applications. Hoboken (NJ): John Wiley & Sons, 2007.
- [2] Q. Jackson and D. A. Landgrebe, "An adaptive method for combined covariance estimation and classification," *IEEE Trans. Geosci. and Remote Sensing*, vol. 40, no. 5, pp. 1082–1087, May 2002.
- [3] M. Lewis, V. Jooste, and A. A. de Gasparis, "Discrimination of arid vegetation with airborne multispectral scanner hyperspectral imagery," *IEEE Trans. Geosci. and Remote Sensing*, vol. 39, no. 7, pp. 1471–1479, July 2001.
- [4] C.-I Chang, Hyperspectral imaging: techniques for spectral detection and classification. New York: Kluwer Academic, 2003.
- [5] N. Keshava and J. F. Mustard, "Spectral unmixing," *IEEE Signal Processing Magazine*, vol. 19, no. 1, pp. 44–57, Jan. 2002.
- [6] R. B. Singer and T. B. McCord, "Mars: Large scale mixing of bright and dark surface materials and implications for analysis of spectral reflectance," in *Proc. Lunar Planet. Sci. Conf.*, vol. 2, 1979, pp. 1835–1848.

- B. W. Hapke, "Bidirectional reflectance spectroscopy. I. Theory," *J. Geophys. Res.*, vol. 86, pp. 3039–3054, 1981.
- [8] P. E. Johnson, M. O. Smith, S. Taylor-George, and J. B. Adams, "A semiempirical method for analysis of the reflectance spectra of binary mineral mixtures," *J. Geophys. Res.*, vol. 88, pp. 3557–3561, 1983.
- [9] D. Manolakis, C. Siracusa, and G. Shaw, "Hyperspectral subpixel target detection using the linear mixing model," *IEEE Trans. Geosci. and Remote Sensing*, vol. 39, no. 7, pp. 1392–1409, July 2001.
- [10] D. C. Heinz and C.-I Chang, "Fully constrained least-squares linear spectral mixture analysis method for material quantification in hyperspectyral imagery," *IEEE Trans. Geosci. and Remote Sensing*, vol. 29, no. 3, pp. 529–545, March 2001.
- [11] N. Dobigeon, J.-Y. Tourneret, and C.-I Chang, "Semisupervised linear spectral unmixing using a hierarchical Bayesian model for hyperspectral imagery," *IEEE Trans. Signal Processing*, vol. 56, no. 7, pp. 2684–2695, July 2008.
- [12] P. J. Green, "Reversible jump MCMC computation and Bayesian model determination," *Biometrika*, vol. 82, no. 4, pp. 711–732, Dec. 1995.
- [13] N. Dobigeon, J.-Y. Tourneret, and Alfred O. Hero III, "Bayesian linear unmixing of hyperspectral images corrupted by colored gaussian noise with unknown covariance matrix," in *IEEE Int. Conf. Acoust., Speech, and Signal Processing (ICASSP)*, Las Vegas, USA, March 2008, pp. 3433–3436.
- [14] M. Winter, "Fast autonomous spectral end-member determination in hyperspectral data," in *Proc. 13th Int. Conf. on Applied Geologic Remote Sensing*, vol. 2, Vancouver, April 1999, pp. 337–344.
- [15] M. Craig, "Minimum volume transforms for remotely sensed data," *IEEE Trans. Geosci. and Remote Sensing*, pp. 542–552, 1994.
- [16] M. Berman, H. Kiiveri, R. Lagerstrom, A. Ernst, R. Dunne, and J. F. Huntington, "ICE: A statistical approach to identifying endmembers in hyperspectral images," *IEEE Trans. Geosci. and Remote Sensing*, vol. 42, no. 10, pp. 2085–2095, Oct. 2004.
- [17] E. Punskaya, C. Andrieu, A. Doucet, and W. Fitzgerald, "Bayesian curve fitting using MCMC with applications to signal segmentation," *IEEE Trans. Signal Processing*, vol. 50, no. 3, pp. 747–758, March 2002.
- [18] S. Richardson and P. J. Green, "On Bayesian analysis of mixtures with unknown number of components," *J. Roy. Stat. Soc. B*, vol. 59, no. 4, pp. 731–792, 1997.
- [19] —, "Corrigendum: On Bayesian analysis of mixtures with unknown number of components," *J. Roy. Stat. Soc. B*, vol. 60, no. 3, p. 661, 1998.
- [20] D. G. T. Denison, C. C. Holmes, B. K. Mallick, and A. F. M. Smith, *Bayesian methods for nonlinear classification and regression*. Chichester, England: Wiley, 2002.
- [21] N. Dobigeon and J.-Y. Tourneret, "Efficient sampling according to a multivariate Gaussian distribution

truncated on a simplex," IRIT/ENSEEIHT/TéSA, Tech. Rep., March 2007. [Online]. Available: http://www.enseeiht.fr/~dobigeon

- [22] C. P. Robert, "Simulation of truncated normal variables," *Statistics and Computing*, vol. 5, pp. 121–125, 1995.
- [23] RSI (Research Systems Inc.), ENVI User's guide Version 4.0, Boulder, CO 80301 USA, Sept. 2003.



Nicolas Dobigeon (S'05–M'08) was born in Angoulême, France, in 1981. He received the Eng. degree in Electrical Engineering from ENSEEIHT, Toulouse, France, and the M.Sc. degree in signal processing from the National Polytechnic Institute of Toulouse, both in June 2004. In October 2007, he ob-

tained the Ph.D. degree in Signal Processing from the National Polytechnic Institute of Toulouse.

From 2007 to 2008, he was a Research Associate at the Department of Electrical Engineering and Computer Science, University of Michigan. Since 2008, he has been an Assistant Professor with University of Toulouse (ENSEEIHT), France. He is also with the Signal and Communication group of the IRIT Laboratory. His research interests are centered around Bayesian estimation and Markov chain Monte Carlo (MCMC) methods for signal and image processing.



Jean-Yves Tourneret (M'94–SM'08) received the Ingénieur degree in Electrical Engineering from Ecole Nationale Supérieure d'Electronique, d'Electrotechnique, d'Informatique et d'Hydraulique in Toulouse (ENSEEIHT), France, in 1989 and the Ph.D. degree from the National Polytechnic Institute of

Toulouse, France, in 1992.

He is currently a professor in the University of Toulouse, France (ENSEEIHT) and a member of the IRIT laboratory (UMR 5505 of the CNRS). His research activities are centered around statistical signal processing with a particular interest to Markov Chain Monte Carlo methods.

Prof. Tourneret was the program chair of the European conference on signal processing (EUSIPCO), which was held in Toulouse (France) in 2002. He was also member of the organizing committee for the international conference ICASSP'06 which was held in Toulouse (France) in 2006. He has been a member of different technical committees including the Signal Processing Theory and Methods (SPTM) committee of the IEEE Signal Processing Society (2001–2007). He is currently serving as an associate editor for the IEEE TRANSACTIONS ON SIGNAL PROCESSING.