# Hyperspectral Image Restoration Based on Spatio-Spectral Structure Tensor Regularization

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Abstract—We propose a regularization function for hyperspectral image restoration based on a newly-designed structure tensor. We adopt a convex optimization approach with the use of the nuclear norm of a matrix, termed as spatio-spectral structure tensor. It consists of the gradient components of a hyperspectral image cube w.r.t. the spatio-spectral domain. The proposed approach allows to penalize variations in the spectral domain as well as the spatial domain to exploit the spatio-spectral correlations. Our experiments on denoising of hyperspectral images show that the proposed regularization leads to significant improvements in restoration performance over state-of-the-art methods.

#### I. INTRODUCTION

In recent years, image processing for hyperspectral images with a large number of spectral bands has become an important issue due to the development of remote sensing and imaging technologies [1], [2]. Since a hyperspectral sensor generally acquires spectral components covering narrow spectral bands, the illuminance captured by the sensor becomes low, so noise signals are amplified together with the luminance.

To extract signals buried in the noise, many denoising methods for the hyperspectral images have been proposed [3]–[8]. BM4D [3] is a highly accurate non-local method for images with three or more channels, which is an extended version of the state-of-the-art 2D image denoising method [9]. Another effective approach is regularization based on total variation-type functions exploiting variations in the spatial domain [5]–[7], which are stated as convex optimization problems and efficiently solved to estimate a noiseless latent image.

Low-rank-based regularization approaches have also been actively investigated [10]–[14]. In [13], the correlation of the spatial structure in a local region of a multi/hyperspectral image is exploited by using structure tensor total variation (STV), and high-performance restoration can be achieved without computationally inefficient non-local search. This is because STV is defined as the nuclear norm of the structure tensor, a matrix consisting of gradient components in a local region, and thus it can evaluate the semi-local spatial correlation of images. However, STV does not measure the correlation between bands because it deals with gradient

components for each band independently. To address this issue, its arranged version (ASTV) [14] has been proposed. ASTV simultaneously measures correlations between bands as well as spatial correlations, but it consists only of gradient components in spatial directions and does not explicitly exploit smoothness in the spectral direction.

In this paper, we propose a regularization function based on the structure tensor consisting of the gradient components both in the spectral domain in addition to the spatial domains. The regularization functions is then defined as the nuclear norm of the said structure tensor, so that our regularization can effectively promote the semi-local smoothness both in the spatial and spectral directions. Then, we formulate hyperspectral image restoration as a convex optimization problem involving our regularization function, which is efficiently solved by a primal-dual splitting method [15]–[17]. We apply our regularization to hyperspectral image denoising and demonstrate that it outperforms several state-of-the-art methods, namely, ASTV and BM4D.

### II. PROPOSED METHOD

### A. Spatio-spectral gradient filter

A hyperspectral image with M bands is denoted as  $\mathbf{u} = [\mathbf{u}'_1^\top, \dots, \mathbf{u}'_M^\top]^\top \in \mathbb{R}^{MN}$  (N is the number of pixels), where  $\mathbf{u}'_j \in \mathbb{R}^N$  ( $j = 1, \dots, M$ ) is the j-th band image. The gradient filters w.r.t. the vertical and horizontal directions in a single band  $\mathbf{u}'_j$  are defined as  $\mathbf{D}'_v$  and  $\mathbf{D}'_h \in \mathbb{R}^{N \times N}$ , respectively. The matrices are extended to apply  $\mathbf{D}'_v$  and  $\mathbf{D}'_h$  to  $\mathbf{u}$ , which are defined as  $\mathbf{D}_v = \operatorname{diag}(\mathbf{D}'_v, \dots, \mathbf{D}'_v) \in \mathbb{R}^{MN \times MN}$  and  $\mathbf{D}_h = \operatorname{diag}(\mathbf{D}'_h, \dots, \mathbf{D}'_h) \in \mathbb{R}^{MN \times MN}$ , respectively. The matrix  $\mathbf{D}_s \in \mathbb{R}^{MN \times MN}$  for derivation w.r.t. the spectral direction is defined as follows:

where  $\mathbf{I} \in \mathbb{R}^{N \times N}$  is an identity matrix and  $\mathbf{O} \in \mathbb{R}^{N \times N}$  is a zero matrix. Using  $\mathbf{D}_v$ ,  $\mathbf{D}_h$ , and  $\mathbf{D}_s$ , a *spatio-spectral* gradient filter matrix is defined as

$$\mathbf{D} = [\mathbf{D}_v^\top \ \mathbf{D}_h^\top \ \mathbf{D}_s^\top]^\top.$$

## B. Spatio-spectral structure tensor

For a given hyperspectral image  $\mathbf{u}$ , the gradient image is represented as  $\mathbf{D}\mathbf{u} \in \mathbb{R}^{3MN}$ , and constructed by two spatial gradient images,  $\mathbf{D}_v\mathbf{u}$  and  $\mathbf{D}_h\mathbf{u} \in \mathbb{R}^{MN}$ , and the spectral gradient image,  $\mathbf{D}_s\mathbf{u} \in \mathbb{R}^{MN}$  as follows:

$$\mathbf{D}\mathbf{u} = [(\mathbf{D}_{v}\mathbf{u})^{\top} (\mathbf{D}_{h}\mathbf{u})^{\top} (\mathbf{D}_{s}\mathbf{u})^{\top}]^{\top},$$

$$\mathbf{D}_{v}\mathbf{u} = [(\mathbf{D}_{v}'\mathbf{u}_{1}')^{\top} (\mathbf{D}_{v}'\mathbf{u}_{2}')^{\top} \dots (\mathbf{D}_{v}'\mathbf{u}_{M}')^{\top}]^{\top}, \quad (1)$$

$$\mathbf{D}_{h}\mathbf{u} = [(\mathbf{D}_{h}'\mathbf{u}_{1}')^{\top} (\mathbf{D}_{h}'\mathbf{u}_{2}')^{\top} \dots (\mathbf{D}_{h}'\mathbf{u}_{M}')^{\top}]^{\top},$$

$$\mathbf{D}_{s}\mathbf{u} = [(\mathbf{D}_{s}'\mathbf{u}_{1}')^{\top} (\mathbf{D}_{s}'\mathbf{u}_{2}')^{\top} \dots (\mathbf{D}_{s}'\mathbf{u}_{M}')^{\top}]^{\top}.$$

We divide the gradient images into 3D local blocks. The shape of each local block is a rectangular block with a square base in the spatial domain, and each block contains all the bands (e.g. the size of the block is  $10 \times 10 \times M$ ). The arrangement of the matrix is illustrated in Fig.1. We define a new structure tensor of the n-th local block by rearranging the matrices as follows:

$$\mathbf{L}_n = [\mathbf{D}\mathbf{u}_1^{(n)} \ \mathbf{D}\mathbf{u}_2^{(n)} \ \dots \mathbf{D}\mathbf{u}_M^{(n)}] \in \mathbb{R}^{N' \times 3M}, \quad (2)$$

where N' is the number of pixels in a single band of a local block.  $\mathbf{D}\mathbf{u}_j^{(n)} = [\mathbf{D}_v\mathbf{u}_j^{(n)} \ \mathbf{D}_h\mathbf{u}_j^{(n)} \ \mathbf{D}_s\mathbf{u}_j^{(n)}] \in \mathbb{R}^{N'\times 3}$   $(j=1,\ldots,M)$ , and  $\mathbf{D}_*\mathbf{u}_j^{(n)}$  consists of the j-th band in the n-th local block. We call  $\mathbf{L}_n$  in (2) a *spatio-spectral structure tensor*. In the next section, we construct a convex optimization problem with a regularization function based on the spatio-spectral structure tensor to restore a latent hyperspectral image.

# C. Regularization based on spatio-spectral structure tensor

In the paper, we focus on hyperspectral image denoising as an example of hyperspectral image restoration. We remark that our regularization can be applied to other problems such as non-blind image deblurring and compressive sensing.

We assume that the degradation process can be modeled as  $\mathbf{y} = \mathbf{u} + \mathbf{n}$ , where  $\mathbf{y} \in \mathbb{R}^{MN}$  is an observation image and  $\mathbf{n} \in \mathbb{R}^{MN}$  is the additive white Gaussian noise. Regularization based on the aforementioned spatio-spectral structure tensor can be defined as the sum of  $\mathrm{rank}(\mathbf{L}_n)$   $(n=1,2,\cdots,K)$  with two constraints. One is a constraint on the intensity range  $\mathcal{S} = [0,1]^{MN}$ . The other constraint represents data-fidelity, defined by the  $\mathbf{y}$ -centered L2-norm ball with the radius  $\epsilon > 0$  controlling the degree of fidelity to  $\mathbf{y}$ . Compared with

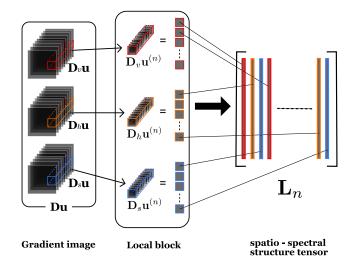


Fig. 1. Spatio-spectral structure tensor

the standard additive data-fidelity, such a constraint-type data-fidelity facilitates parameter setting because  $\epsilon$  has a clear physical meaning, so that it can be easily adjusted based on noise standard deviation, as addressed in [18]–[21]. Specifically we consider the following optimization problem:

$$\min_{\mathbf{x}} \sum_{n=1}^{K} \operatorname{rank}(\mathbf{L}_{n}) \quad s.t. \ \mathbf{x} \in \mathcal{S}, \ \|\mathbf{x} - \mathbf{y}\|_{2} \le \epsilon, \ (3)$$

where K is the number of extracted local blocks, and all pixels of  $\mathbf{x}$  are extracted at least once as a pixel of a local block. Here,  $\mathbf{L}_n$  is represented with an operator  $\mathbf{P}_n:\mathbb{R}^{3MN}\to\mathbb{R}^{3MN'}$  that extracts the n-th local block as

$$\mathbf{L}_n = \mathbf{P}_n \mathbf{D} \mathbf{x}. \tag{4}$$

Substituting this for Eq.(3) yields

$$\min_{\mathbf{x}} \sum_{n=1}^{K} \operatorname{rank}(\mathbf{P}_{n} \mathbf{D} \mathbf{x}) \quad s.t. \ \mathbf{x} \in \mathcal{S}, \ \|\mathbf{x} - \mathbf{y}\|_{2} \le \epsilon. \ (5)$$

Since  $\operatorname{rank}(\cdot)$  is the nonconvex function that counts the number of non-zero singular values, which makes it intractable to find the optimal solution, we introduce the nuclear norm  $\|\cdot\|_* = \sum_i \sigma_i$  ( $\sigma_i$  is the singular values) [22], which is a reasonable convex relaxation for  $\operatorname{rank}(\cdot)$ . Thus the problem (5) is converted to the following convex optimization problem:

$$\min_{\mathbf{x}} \sum_{n=1}^{K} \|\mathbf{P}_n \mathbf{D} \mathbf{x}\|_* \quad s.t. \ \mathbf{x} \in \mathcal{S}, \ \|\mathbf{x} - \mathbf{y}\|_2 \le \epsilon. \ (6)$$

We further reformulate (6) into the following form:

$$\min_{\mathbf{x}} \sum_{n=1}^{K} \|\mathbf{P}_{n} \mathbf{D} \mathbf{x}\|_{*} + \iota_{\mathcal{S}}(\mathbf{x}) + \iota_{\mathcal{B}_{\mathbf{y},\epsilon}}(\mathbf{x}), \tag{7}$$

where S is the closed convex set used in (3), and the indicator function is defined as

$$\iota_{\mathcal{S}}(\mathbf{x}) = \begin{cases} 0 & \mathbf{x} \in \mathcal{S}, \\ +\infty & \text{otherwise.} \end{cases}$$
 (8)

The set  $\mathcal{B}_{\mathbf{y},\epsilon}$  is defined as  $\mathcal{B}_{\mathbf{y},\epsilon} = \{\mathbf{x} \in \mathbb{R}^{MN} \mid \|\mathbf{x} - \mathbf{x}\| \}$  $\mathbf{y}|_2 \leq \epsilon$  using the prescribed error tolerance  $\epsilon$ , and  $\iota_{\mathcal{B}_{\mathbf{v},\epsilon}}(\mathbf{x})$  is defined in the same way as (8).

### D. Optimization

For solving the convex optimization problem in (7), we use a primal-dual splitting (PDS) method [15]–[17]. The PDS handles convex optimization problems of the form:

$$\min_{\mathbf{x}} F(\mathbf{x}) + G(\mathbf{x}) + H(\mathbf{A}\mathbf{x}), \tag{9}$$

where F is a differentiable convex function whose gradient is  $\beta$ -Lipschitz continuous, G and H are possibly nonsmooth convex functions whose proximity operators  $^{1}$  are computable, and **A** is a linear operator. The algorithm for (9) is given by

$$\mathbf{x}^{(k+1)} = \operatorname{prox}_{\gamma_1 G}(\mathbf{x}^{(k)} - \gamma_1(\nabla F(\mathbf{x}^{(k)}) + \mathbf{A}^{\top} \mathbf{z}^{(k)})),$$

$$\mathbf{z}^{(k+1)} = \operatorname{prox}_{\gamma_2 H^*}(\mathbf{z}^{(k)} + \gamma_2 \mathbf{A}(2\mathbf{x}^{(k+1)} - \mathbf{x}^{(k)})),$$
is given by the convex projection onto the  $l_2$  ball (10th

where  $\nabla F$  is the gradient of F,  $\mathbf{A}^{\top}$  is the transpose of A, x is the primal variable and z is the dual variable.  $H^*$  is the conjugate function of H, and its proximity operator can be computed using the proximity operator of H, given by

$$\operatorname{prox}_{\gamma_2 H^*}(\mathbf{v}) = \mathbf{v} - \gamma_2 \operatorname{prox}_{\gamma_2^{-1} H}(\gamma_2^{-1} \mathbf{v}). \tag{11}$$

Under appropriate conditions on  $\gamma_1$  and  $\gamma_2$ , the sequence generated by (10) converges to an optimal solution of (9). In order to apply PDS to our problem (7), we define F, G, H, and A as follows

$$F : \mathbb{R}^{MN} \to \mathbb{R}, \quad \mathbf{x} \mapsto 0,$$

$$G : \mathbb{R}^{MN} \to \mathbb{R} \cup \{\infty\}, \quad \mathbf{x} \mapsto \iota_{\mathcal{S}}(\mathbf{x}),$$

$$H : \mathbb{R}^{MN'+MN} \to \mathbb{R} \cup \{\infty\},$$

$$\mathbf{z} = [\mathbf{z}_{1}^{\top}, \mathbf{z}_{2}^{\top}]^{\top} \mapsto \sum_{n=1}^{K} \|\mathbf{z}_{1,n}\|_{*} + \iota_{\mathcal{B}_{\mathbf{y},\epsilon}}(\mathbf{z}_{2}),$$

$$\mathbf{A} : \mathbb{R}^{MN} \to \mathbb{R}^{3MN'+MN}, \quad \mathbf{x} \mapsto (\mathbf{P}_{n}\mathbf{D}\mathbf{x}, \mathbf{x}),$$

<sup>1</sup>The proximity operator (prox) [23], [24] is defined as  $\operatorname{prox}_{\sim H}(\mathbf{x}) =$  $\arg\min_{\mathbf{a}} H(\mathbf{a}) + \frac{1}{2\gamma} \|\mathbf{a} - \mathbf{x}\|_2^2$ 

**Algorithm 1** Algorithm for solving Prob. (7)

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1: input : \mathbf{x}^{(0)}, \mathbf{z}_{1,n}^{(0)}(n=1,...,K), \mathbf{z}_2^{(0)}
   2: set : \gamma_1, and \gamma_2 are given.
   3: while stopping criterion is satisfied. do
                     \mathbf{x}^{(k+1)} = P_{\mathcal{S}}(\mathbf{x}^{(k)} - \gamma_1(\sum_{n=1}^K \mathbf{D}^\top \mathbf{P}_n^\top \mathbf{z}_{1,n}^{(k)} + \mathbf{z}_2^{(k)})) for n = 1 to K do
                              \mathbf{r} = 1 \text{ to } K \text{ do} \\ \mathbf{v}_{1,n}^{(k)} = \mathbf{z}_{1,n}^{(k)} + \gamma_2 (\mathbf{P}_n \mathbf{D}(2\mathbf{x}^{(k+1)} - \mathbf{x}^{(k)})) \\ \mathbf{z}_{1,n}^{(k+1)} = \mathbf{v}_{1,n}^{(k)} - \gamma_2 \operatorname{prox}_{\frac{1}{\gamma_2} \|\cdot\|_*} (\frac{1}{\gamma_2} \mathbf{v}_{1,n}^{(k)})
   6:
   7:
                    \mathbf{v}_{2}^{(k)} = \mathbf{z}_{2}^{(k)} + \gamma_{2}(2\mathbf{x}^{(k+1)} - \mathbf{x}^{(k)})
\mathbf{z}_{2}^{(k+1)} = \mathbf{v}_{2}^{(k)} - \gamma_{2}\operatorname{prox}_{\frac{1}{\gamma_{2}}\iota_{\mathcal{B}_{\mathbf{y},\epsilon}}}(\frac{1}{\gamma_{2}}\mathbf{v}_{2}^{(k)})
12: end while
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where  $\mathbf{z}_{1,n}$  represents the *n*-th local block of  $\mathbf{z}_1$ . The steps of the algorithm for solving (7) are shown in Algorithm 1, where  $P_S$  in the 4th line is projection onto the convex set S, which transfers pixel values within the range. The proximity operator of the nuclear norm in the 7th line is obtained by applying singular value decomposition to  $L_n$  and then thresholding the singular values:

$$\operatorname{prox}_{\gamma \|\cdot\|_{*}}(\mathbf{L}_{n}) = \mathbf{U}\tilde{\Sigma}\mathbf{V}^{\top},$$

$$\tilde{\Sigma} = \operatorname{diag}(\max\{\sigma_{1} - \gamma, 0\}, \dots, \max\{\sigma_{3M} - \gamma, 0\}).$$
(12)

is given by the convex projection onto the  $l_2$  ball (10th

$$\operatorname{prox}_{\gamma_{\iota_{\mathcal{B}_{\mathbf{y},\epsilon}}}}(\mathbf{x}) = \begin{cases} \mathbf{x} & \mathbf{x} \in \mathcal{B}_{\mathbf{y},\epsilon}, \\ \mathbf{y} + \frac{\epsilon}{\|\mathbf{x} - \mathbf{y}\|_2} (\mathbf{x} - \mathbf{y}) & \text{otherwise.} \end{cases}$$
(13)

### III. EXPERIMENTAL RESULTS

In order to evaluate the performance of the proposed method, we compare it with ASTV [14] and BM4D [3] on denoising of hyperspectral images. We assume that input images are degraded by additive white Gaussian noise  $\mathbf{n} \in \mathbb{R}^{MN}$  with standard deviation of 0.1. The parameter  $\epsilon$  in (6) is appropriately changed depending on the noise intensity. All pixel values are normalized to within the range [0,1]. The size of the local blocks (in the block processing of the nuclear norm in both of the methods) was set to  $10 \times 10 \times M$ , and there is no overlap between them<sup>2</sup>. As for BM4D, we use the

<sup>&</sup>lt;sup>2</sup>On the basis of our experience, the block overlap significantly increases computational complexity, while we achieve it with only a little improvement in restoration performance.

PSNR[dB] / SSIM 19.99[dB] / 0.2703 32.42[dB] / 0.8241 33.49[dB] / 0.8438 **34.51[dB**] / **0.8923** 

PSNR[dB] / SSIM 19.99[dB] / 0.2863 32.11[dB] / 0.8273 33.44[dB] / 0.8502 **34.29[dB**] / **0.8965** 

PSNR[dB] / SSIM 19.99[dB] / 0.1805 34.22[dB] / 0.8338 35.01[dB] / 0.8617 **36.01[dB**] / **0.8933** 

Fig. 2. Results in PaviaC (102 bands) (upper), PaviaU (103 bands) (middle) and Frisco(148 bands) (lower): from left, Original, Noisy, ASTV [14], BM4D [3], and ours

MATLAB code published by the authors. Test images were taken from the images in [8], [25], [26]. In order to objectively evaluate denoising performance, two indexes: peak signal to noise ratio (PSNR) and SSIM [27] are used.

The quantitative evaluation of denoising performance by PSNR and SSIM is shown in Tables I and II, respectively. The experimental results for three hyperspectral images, PaviaC (102 bands), PaviaU (103 bands), and Frisco (148 bands) are shown in Fig.2, in which we show the 95th band in the PaviaC, the 98th band in the PaviaU and the 130th band in the Frisco. It can be seen

from Fig.2 that the proposed method can achieve satisfactory denoising performance while maintaining details compared to the conventional methods. In quantitative evaluations in Tables I and II, it can be confirmed that the quantitative performance of noise removal in the proposed method outperforms the conventional ones.

### IV. CONCLUSION

In this paper, we proposed a regularization method based on the spatio-spectral structure tensor aiming at restoration of hyperspectral images. The proposed method simultaneously exploits two properties. One is

TABLE I
COMPARISON WITH [14] AND [3] (IN PSNR[DB])

Image	(bands)	ASTV [14]	BM4D [3]	Ours
China	(7)	32.51	32.02	32.05
Ribeira	(33)	34.49	34.62	35.84
PaviaC	(102)	32.42	33.49	34.51
PaviaU	(103)	32.11	33.44	34.29
Frisco	(148)	34.22	35.01	36.01
Stanford	(148)	34.45	35.31	36.34
Salinas	(224)	35.03	36.76	36.50

TABLE II COMPARISON WITH [14] AND [3] (IN SSIM)

Image	(bands)	ASTV [14]	BM4D [3]	Ours
China	(7)	0.8729	0.8534	0.8521
Ribeira	(33)	0.8652	0.8740	0.9077
PaviaC	(102)	0.8241	0.8438	0.8923
PaviaU	(103)	0.8273	0.8502	0.8965
Frisco	(148)	0.8338	0.8617	0.8933
Stanford	(148)	0.8229	0.8649	0.8846
Salinas	(224)	0.8106	0.8720	0.8625

the low-rank property of gradient images w.r.t. the spatiospectral directions in the hyperspectral image, and the other is the correlation between spectral bands. By solving the convex optimization problem using some regularization, high performance method in denoising for hyperspectral images is achieved. The experimental results demonstrate that the proposed method outperforms the conventional methods.

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