Relative Spectral Difference Occurrence Matrix: A Metrological Spectral-Spatial Feature for Hyperspectral Texture Analysis

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Abstract

We develop a spectral-spatial feature, Relative Spectral Difference Occurrence Matrix (RSDOM) for hyperspectral texture recognition. Inspired by Julesz's conjecture, the proposed feature is based on spectral difference approach and respects the metrological constraints. It simultaneously considers the distribution of spectra and their spatial arrangement in the hyperspectral image. The feature is generic and adapted for any number of spectral bands and range. We validate our proposition by applying a classification scheme on the HyTexiLa database. An accuracy comparable to local binary pattern approach is obtained, but at a much reduced cost due to the extremely small feature size.

Introduction

Hyperspectral imaging (HSI) has become increasingly popular with the advance of sensor technology and computational speed. Due to the ability of HSI in capturing the entire electromagnetic spectrum at hundreds of contiguous, narrow spectral bands, it allows non-destructive assessment of material with far superior discrimination. As such, hyperspectral texture analysis has received considerably attention and its application can be found in many fields including biomedical, industrial and remote sensing. For instances, hyperspectral texture analysis have been used in classification of biopsy samples [1, 2], food quality inspection [3, 4] and ground classification of aerial imagery [5, 6].

Many assessment of hyperspectral texture are the direct outgrowth of approaches originally adapted for monospectral (intensity) images [7, 8, 9]. Facing the complexity of hyperspectral data, the common strategy is band selection [10, 11, 12] or dimensionality reduction [13, 14, 15] using Principal Component Analysis (PCA), Non-Negative Matrix Factorization (NMF) and others. Then, the multivariate texture assessment is performed in a marginal way (band by band) or cross-channel processing [16]. Such approach is data dependent and consequently, results from different dataset are not incomparable. Besides, the act of feature transformation in dimensionality reduction could corrupt a part of the physical content of the dataset [17]. For metrological purposes, all physical content need to be preserved so that the processing accuracy, uncertainty, and bias are quantifiable [18].

In this work, we attempt to develop a spectral-spatial feature for hyperspectral texture analysis. We begin by considering the Julesz conjecture, which states that human discrimination is possible between textures differing in low order statistics [19, 20]. We next take inspiration from the pioneering work of Haralick and Ojala in texture analysis, namely Co-occurrence matrix [21] and Local Binary Pattern (LBP) [22]. We remark that the texture description using Co-occurrence matrix is complete, but its reduction into moments for similarity measure causes information losses. For LBP, the binarization provides a weak characterization of texture, but its similarity measure is extremely efficient. Following the work of [23] which relates Co-occurrence matrix

Figure 1: An illustration of RSDOM, combining spectral and spatial features

and histogram of differences, we combine the strength of both approaches and adapt it for HSI processing while respecting the metrological constraints. For enhanced discrimination, we incorporate the spectral feature in the texture assessment, thus inducing a spectral-spatial feature.

The rest of the article is organized in the following manner. We next describe our proposed spectral-spatial feature as well as its similarity measure. Then, we propose an experiment to test the feature with analysis and discussion, along with suggestions for future work. Finally, we present the conclusion for this work.

Proposed method

We are interested in a spectral-spatial feature for hyperspectral texture recognition that has the following properties: (I) computationally efficient; adapted for any spectral count and range, (II) metrological; conform to the science of measurement, and (III) compact; for concise representation. In the following, we describe the formulation of the feature in terms of its expected properties as well as its measure of similarity.

A spectral-spatial feature

We consider the description of hyperspectral texture from both spatial and spectral perspectives. The spatial feature examines the spatial arrangement in the image that gives rise to ideas such as smoothness, coarseness and pattern repeatability [24]. On the other hand, the spectral feature assesses the spectral information in the image that induces concepts such as intensity, saturation and purity. When combined, the spectral-spatial features provide a complete description of hyperspectral texture.

We begin by first considering the formulation of spatial feature M_{spa} . Following Julesz's conjecture on texture discrimination based on low order statistics [19, 20], we describe the spatial arrangement in a hyperspectral image H via the distribution of spectral difference between pixel pairs separated by a spatial vector $\vec{v}$. Suppose that $\vec{v}$ is specified by distance d and direction

θ , $M_{spa.}$ is expressed as a probability density function given by:

$$M_{spa.}^{(d,\theta)}(\Delta S) = \text{Prob}\left(d(S_i, S_j) = \Delta S\right), \forall i, j \in H, \quad (1)$$

$$\|\vec{ij}\| = d, \angle \vec{ij} = \theta$$

where $d(S_i, S_j) = \Delta S$ is the spectral difference between spectra S_i and S_j of pixels i and j respectively, separated by $\vec{v}$.

Next, we consider the formulation of spectral feature $M_{spe.}$. We characterize the spectral variation in H via spectral difference $d(S_i, S_r)$, where $i \in H$ and S_r is a chosen reference. Doing so, we are able to reduce the description from a P -dimensional spectral space to a p -dimensional difference space, where $p \ll P$. Expressed as a probability density function, $M_{spe.}$ is given by:

$$M_{spe.}(\Delta S_r) = \text{Prob}\left(d(S_i, S_r) = \Delta S_r\right), \forall i \in H \quad (2)$$

where $d(S_i, S_r) = \Delta S_r$ is the spectral difference between each pixel in the image and the reference. The choice of S_r is made such that it lies *just* outside the convex hull of the spectral distribution [25] of the entire image set. This is to maximize discrimination between the textured images.

Combining $M_{spa.}$ and $M_{spe.}$, we obtain the spectral-spatial feature M as joint probability density function given by:

$$M^{(d,\theta)}(\Delta S, \Delta S_r) = \text{Prob}\left(d(S_i, S_j) = \Delta S, d(S_i, S_r) = \Delta S_r\right), \quad (3)$$

$$\forall i, j \in H, \|\vec{ij}\| = d, \angle \vec{ij} = \theta$$

termed as *Relative Spectral Difference Occurrence Matrix* (RSDOM). An illustration of RSDOM is shown in Figure 1.

To explain further on the choice of S_r , consider the working example in Figure 2. Suppose there are two textured images (each with distinct spectral distribution $\mathcal{S}_1$ and $\mathcal{S}_2$) and three possible choices of S_r (S_{r1} , S_{r2} and S_{r3}). The task is to select an optimal S_r so that the discrimination between $\mathcal{S}_1$ and $\mathcal{S}_2$ is the maximum in the difference space. Referring to Figure 3 for expression in the difference space, S_{r1} is obviously a bad choice as it would have “perceived” $\mathcal{S}_1$ and $\mathcal{S}_2$ to be equidistant in the difference space. Both S_{r2} and S_{r3} lie outside the convex hull C of the spectral distribution, but S_{r2} would have been a better choice as S_{r3} may risk “perceiving” $\mathcal{S}_1$ and $\mathcal{S}_2$ to be equally far in the difference space. For maximized discrimination, it is thus evident that S_r has to be chosen such that it lies just outside the convex hull of the entire spectral distribution.

Metrological calculation of spectral difference

A spectrum is essentially a continuous function $S = f(\lambda)$ defined over a spectral range λ_{min} to λ_{max} that characterizes an electromagnetic radiation [26]. Due to the sampling operation during acquisition, a spectrum is represented as a sequence of measurements $S = \{s_j(\lambda), \forall \lambda \in [\lambda_{min}, \lambda_{max}]\}$ with spectral resolution depending on the sensor. Due to its discrete representation, a spectrum is often treated as if it were a set of independent measures, hence as vectors, probability density functions or sequences [27]. This leads to the usage of L2 norms in accessing spectral difference, for instances spectral angle mapper [28] and various divergence measures [29, 30, 31].

Figure 2: Representation in P -dimensional spectral space

Figure 3: Representation in p -dimensional difference space

However, the limits of such approach has been pointed out in [32]. While our formulation of RSDOM is not limited to any spectral difference formula, in this work we seek to express spectral difference based on the fact that spectrum is a *function*. As such, we use Kullback-Leibler pseudo-divergence (KLPD) [33] which is inspired from Kullback-Leibler divergence. Consider two spectra S_1 and S_2 , the KLPD is defined as the sum of spectral shape difference ΔG and intensity difference ΔW as:

$$d_{KLPD}(S_1, S_2) = \Delta G(S_1, S_2) + \Delta W(S_1, S_2) \quad (4)$$

where:

$$\Delta G(S_1, S_2) = k_1 \cdot KL(\bar{S}_1 \| \bar{S}_2) + k_2 \cdot KL(\bar{S}_2 \| \bar{S}_1) \quad (5)$$

$$\Delta W(S_1, S_2) = (k_1 - k_2) \log\left(\frac{k_1}{k_2}\right) \quad (6)$$

as the normalized spectra $\bar{S}$ is expressed as:

$$\bar{S}_j = \left\{ \bar{s}_j(\lambda) = \frac{s_j(\lambda)}{k}, \forall \lambda \in [\lambda_{min}, \lambda_{max}] \right\}, j \in \{1, 2\} \quad (7)$$

with the normalization constant given by:

$$k_j = \int_{\lambda_{min}}^{\lambda_{max}} s_j(\lambda) d\lambda, j \in \{1, 2\} \quad (8)$$

Employing KLPD, RSDOM is now a four dimensional joint probability density function given by:

$$M^{(d,\theta)}(\Delta G, \Delta W, \Delta G_r, \Delta W_r) =$$

$$\text{Prob}\left(d_{KLPD}(S_i, S_j) = (\Delta G, \Delta W), d_{KLPD}(S_i, S_r) = (\Delta G_r, \Delta W_r)\right),$$

$$\forall i, j \in H, \|\vec{ij}\| = d, \angle \vec{ij} = \theta \quad (9)$$

Representation using Gaussian Mixture Model

Considering RSDOM as probability density function, f-divergence is well adapted for to access the similarity between them [34]. Among all, Kullback-Leibler (KL) divergence has been identified as the most efficient similarity measure [35]. The KL divergence [36] $KL(P||Q)$ assess the quantity of information lost when Q is used to estimate P :

$$KL(P||Q) = \int_{-\infty}^{\infty} p(x) \ln \frac{p(x)}{q(x)} dx \quad (10)$$

where p and q are the respective densities of P and Q . In the particular case of multivariate normal distributions with mean μ and variance-covariance matrix Σ , KL divergence can be directly calculated by [37]:

$$KL(P||Q) = \frac{1}{2} \left[\log \frac{|\Sigma_Q|}{|\Sigma_P|} + \text{tr}(\Sigma_Q^{-1} \Sigma_P) - D + (\mu_Q - \mu_P)^T \Sigma_Q^{-1} (\mu_Q - \mu_P) \right] \quad (11)$$

where D is the dimension of data. As KL divergence is not symmetric, a similarity measure is defined as:

$$d_{KL}(P, Q) = KL(P||Q) + KL(Q||P) \quad (12)$$

Due to spectral variability and sensor noise, RSDOM is understandably huge and is approximately the size of the image. The direct calculation of KL divergence is therefore, computationally expensive, not to mention the huge size of RSDOM which induces problems in storage and transmission. We thus proceed to model RSDOM using Gaussian mixture model (GMM):

$$M = \sum_{i=1}^K \phi_i \mathcal{N}(\mu_i, \Sigma_i) \quad (13)$$

where ϕ is the mixture weight and each Gaussian component is modeled using $\mu \in \mathbb{R}^D$ and $\Sigma \in \mathbb{R}^{D \times D}$. Hence, RSDOM can be represented using just $K \times (D + D^2)$ scalar values with $D = 4$ which is extremely compact.

As there is no closed form solution for KL divergence between two GMMs, we resort to using variational approximation [38] KL_V , defined as:

$$KL_V(M||M') = \sum_i \phi_i \log \frac{\sum_{i'} \phi_{i'} e^{-KL(M_i||M_{i'})}}{\sum_j \phi'_j e^{-KL(M_i||M'_j)}} \quad (14)$$

as the two RSDOMs are given by $M = \sum_i \phi_i M_i$ and $M' = \sum_j \phi'_j M'_j$. Hence, the similarity measure is given by:

$$d_{KL}(M, M') = KL_V(M||M') + KL_V(M'||M) \quad (15)$$

At this point, it is worth mentioning that as KLPD is a distance measure that takes no account in direction. This causes the distribution of RSDOM to be one-sided in the ΔG - ΔW plane. This induces imperfection in the modeling as GMM assumes

Figure 4: Several images (rendered in RGB) taken from HyTexiLa dataset

double sided distribution. Therefore, we introduce a sign operator $\text{sign}(\tau)$ where τ is defined to be positive when S_i is further from S_r compared to S_j , or negative otherwise:

$$\tau = d_{KLPD}(S_i, S_r) - d_{KLPD}(S_j, S_r) \quad (16)$$

Hence, the final formulation of RSDOM is given by:

$$M^{(d, \theta)}(\Delta G, \Delta W, \Delta G_r, \Delta W_r) = \text{Prob} \left(\text{sign}(\tau) \cdot d_{KLPD}(S_i, S_j) = (\Delta G, \Delta W), d_{KLPD}(S_i, S_r) = (\Delta G_r, \Delta W_r) \right), \quad \forall i, j \in H, \|\vec{i} \vec{j}\| = d, \angle \vec{i} \vec{j} = \theta \quad (17)$$

Experiment and discussion

To access the performance of RSDOM in hyperspectral texture recognition, we apply a classification scheme on a hyperspectral texture dataset. As comparison, we benchmark our performance against local binary pattern based approach [16]. We analyze the result and lists down the suggestions for future work.

The HyTexiLa dataset

HyTexiLa (Hyperspectral Texture images acquired in Laboratory) [16] is a hyperspectral reflectance image dataset consisting of $N=112$ textured images from five categories: food (10 images), stone (4 images), textile (65 images), vegetation (15 images) and wood (18 images). The spectral count is 186, ranging from 405.37 nm to 995.83 nm at 3.19 nm interval, hence spanning both visible and near infrared region. Each image measures $1024 \times 1024 \times 186$ and is acquired using the HySpex VNIR-1800 hyperspectral camera. Several images from the dataset, rendered in RGB format, are presented in Figure 4.

Classification using nearest neighbor

We perform the classification using nearest neighbor (NN) approach. In NN, the query (test) image is compared against the training images of all classes and is assigned to the class which

Figure 5: Choosing the reference for HyTexiLa dataset

have the highest similarity with. In the context of our classification scheme, each image, regardless of its categories, is considered a class. Following the setting in [16], each image is split into $n=25$ patches, each measuring $204 \times 204 \times 186$, of which 12 is used for training and 13 for testing. We first perform the classification on the entire dataset (denoted as “All”) and later, intra-categorical classifications as some category is different from others in terms of texture directionality and spectral distinctness. If the query image is classified to its own class, the classification is marked correct or otherwise false. The classification accuracy is therefore defined as the ratio of correct classification to the total number of classification. To complement the classification result, we also report the F1-score, defined as the ratio of $2(\text{precision} \times \text{recall})$ to $(\text{precision} + \text{recall})$. To avoid any bias, we repeat all classification using 10 trials with random selection of training set and testing set in each trial. We report the average accuracy and F1-score along with standard error, with is standard deviation divided by $\sqrt{T}$ where T is the number of trials.

Choice of parameters

In [16], a local binary pattern (LBP) based approach is used with $d = 1$ on a neighborhood of 8 pixels. This is roughly equivalent to accessing spectral variation across three pixels. Therefore, we choose $d = 3$ in the calculation of RSDOM for equal comparison. Since RSDOM is a uni-directional texture descriptor, we arbitrary choose $\theta = 0$. For GMM modeling, we empirically determine that $K = 4$ is optimal for RSDOM representation.

We next consider the choice of reference. Let S_μ be the average spectrum from each patch of m pixels, defined as:

$$S_\mu = \left\{ s_\mu(\lambda) = \frac{1}{m} \sum_{i=1}^m s(\lambda), \forall \lambda \in [\lambda_{\min}, \lambda_{\max}] \right\} \quad (18)$$

and $S_{\mu, \text{global}}$ be the global average spectrum of the entire image set containing N images with n patches, defined as:

$$S_{\mu, \text{global}} = \frac{1}{N \times n} \sum_{i=1}^N \sum_{j=1}^n S_{\mu, i, j} \quad (19)$$

where $S_{\mu, i, j}$ is the average spectrum for the j^{th} patch from the i^{th} image. For simplicity, we consider $d(S_\mu, S_{\mu, \text{global}})$ to obtain a rough representation of the spectral distribution for the dataset.

To choose S_r , we recall that it has to be chosen such that it lies just outside the convex hull of the spectral distribution of the entire image set. However, we decide to adapt the rule considering

(a) coffee

(b) milkcoffee

(c) RSDOM, coffee

(d) RSDOM, milkcoffee

Figure 6: Representation of RSDOM for two textured images in the ΔW - ΔG plane

the large number of images (112) with varying texture complexity between the categories. In fact, we could divide the images into two groups. Group I, enclosed by convex hull C_1 , considers food, stone, vegetation and wood which are natural and complex textures. Group II, enclosed by convex hull C_2 , consists of textile which are man-made and relatively simple textures. For balanced performance, we choose to focus more on the classification of Group I. Consequently, we choose S_r such that it lies just outside C_1 as illustrated in Figure 5. By trial and error, we determine a choice of S_r that fulfills such requirement as:

$$S_r = 3 \left(0.5 + 0.5 \operatorname{erf} \left(\frac{(\lambda - \mu_\lambda)}{\sigma_\lambda} \right) \right) \quad (20)$$

where erf is the error function. μ_λ and σ_λ are given by:

$$\mu_\lambda = (\lambda_{\max} - \lambda_{\min})/2 \quad (21)$$

$$\sigma_\lambda = (\lambda_{\max} - \lambda_{\min})/3 \quad (22)$$

Result and analysis

Figure 6 illustrates RSDOM representation in the ΔG - ΔW plane for two textured images, “coffee” and “milkcoffee”. It can be seen that their spectral shape ΔG variation are similar, explainable from the fact that both “coffee” and “milkcoffee” are of the same food origin differing only on the processing level. On the spectral intensity ΔW , “coffee” exhibits higher variation than “milkcoffee” due to the fact that the former is coarser than the latter. This demonstrates that RSDOM is able to capture the textural properties in an easily understandable manner.

Table 1 presents the classification accuracy and F1-score using RSDOM. The classification result is excellent, with accuracy and F1-score above 90 % in all cases. For the entire dataset, the

Category	Accuracy (%)	F1-score (%)
All	95.6 ± 0.2	95.8 ± 0.2
Food	97.5 ± 0.4	97.5 ± 0.3
Stone	98.3 ± 0.5	98.3 ± 0.5
Vegetation	93.7 ± 0.5	94.1 ± 0.5
Wood	90.8 ± 0.7	91.2 ± 0.7
Textile	98.6 ± 0.1	98.7 ± 0.1

Table 1: Classification accuracy and F1-score on HyTexiLa database

classification accuracy reaches 95.6 %. To further analyze this result, we refer to the normalized confusion matrix as shown in Figure 7. As there are too many class (112 in total) which would rendered examination difficult, we display the confusion matrix in two parts. Figure 7(a) shows the confusion matrix with textile (T) classes clustered together, while Figure 7(b) shows the confusion matrix with food (F), stone (S), vegetation (V) and wood (W) classes clustered together. It is clear that the misclassification is largely due to two categories: wood and vegetation. This is also supported from the intra-categorical classification result in Table 1 which shows their lower accuracy compared to others.

We highlight three reasons to explain the misclassification. Firstly, the texture are mainly green for vegetation and brown for wood. Therefore, the makes discrimination within the categories more difficult. Secondly, our unitary choice of $d = 3$ prevents us from assessing texture with varying texton [39] size. This explains our higher score in food, stone and textile which consist of uniform and fine textons, but lower score in wood and vegetation which has varying texton size. Thirdly, our unitary choice of θ allows us to only evaluate texture in one direction. In fact, for $\theta = 0$ we are only able to assess texture varying in vertical direction but not in other directions. This is acceptable for isotropic texture, but not adapted for anisotropic texture such as those found in vegetation and wood categories.

In comparison to [16] which reports an accuracy of 98.76 % using Opponent Band Local Binary Pattern (OBLBP), our performance is lower although in the same efficiency range. The difference in performance can be attributed to the fact that OBLBP assesses texture in eight directions while RSDOM evaluates just one. However, it is worthwhile to mention that the feature size of RSDOM is extremely small. While OBLBP consists of 82944 scalar values, the size of RSDOM is only $K \times (D + D^2) = 80$ scalar values or 0.1 % of OBLBP's. Even if RSDOM is adapted to assess texture in eight directions as in the case of OBLBP, its feature size only increases to $8 \times K \times (D + D^2) = 640$ scalar values. This presents RSDOM as an extremely compact feature descriptor for rapid processing and storage.

Limitations and future work

The choice of $d = 3$ remain questionable, as it is possibly susceptible to image noise and it is expected that a larger d value could present better result. In fact, as texture varies in scale and direction, RSDOM could benefit from a multi-scale and multi-direction implementation. On the other hand, the choice of using variational approximation in this work to evaluate divergence between two RSDOM which are GMMs remains to be justified. Other methods ought to be studied in order to determine the best distance measure. ast but no least, other approaches of modeling RSDOM such as Generalized Gaussian Distribution (GGD) or shape descriptions (Fourier Descriptor, Zernike moments, Curvature Scale Space etc.) are also worth being studied.

(a) Textile classes clustered together

(b) Food, stone, vegetation and wood classes clustered together

Figure 7: Normalized confusion matrix for HyTexiLa classification (F=food, S=stone, V=vegetation, W=wood, T=textile)

Conclusion

We have proposed a spectral-spatial feature, Relative Spectral Difference Occurrence Matrix (RSDOM) for hyperspectral texture recognition. The proposed feature is in full accordance with psychophysical experiments on texture. It simultaneously considers the distribution of spectra and their spatial arrangement in the hyperspectral image. It is generic and adapted for any number of spectral bands and range. Thanks to its operation in the difference space, the curse of dimensionality is overcome without requiring band selection or dimensionality reduction techniques, therefore preserving the metrological properties. Its performance has been evaluated by applying a classification scheme on the HyTexiLa dataset with excellent accuracy (95.6%). The result is comparable to Local Binary Pattern (LBP) based approaches, but at a much smaller feature size (0.1 % of the LBP-based feature).

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