

# Advances in Hyperspectral Image and Signal Processing

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**Abstract**—Recent advances in airborne and spaceborne hyperspectral imaging technology have provided end users with rich spectral, spatial, and temporal information, which make a plethora of applications for the analysis of large areas of the Earth surface feasible. However, a huge number of factors, such as high dimensions and size of the hyperspectral data, the lack of training samples, mixed pixels, light scattering mechanisms in the acquisition process, and different atmospheric and geometric distortions, make such data inherently nonlinear and complex, which poses extreme challenges for existing methodologies to effectively process and analyze the data sets. Hence, rigorous and innovative methodologies are required for hyperspectral image and signal processing and have become a center of attention for researchers worldwide. This paper offers a comprehensive tutorial/overview focusing specifically on hyperspectral data analysis, which is categorized into seven broad topics: classification, spectral unmixing, dimensionality reduction, resolution enhancement, hyperspectral image denoising and restoration, change detection, and fast computing. For each topic, we provide a synopsis of the state-of-the-art approaches and numerical results for validation and evaluation of different methodologies, followed by a discussion of future challenges and research directions.

**Index Terms**—Hyperspectral image and signal processing, classification, spectral unmixing, dimensionality reduction, resolution enhancement, image restoration, denoising, change detection, fast computing, missions, tutorials.

## I. INTRODUCTION

REMOTE SENSING necessitates obtaining information from an object or a scene without having any physical contact with it. This is possible because different objects

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reflect, absorb, and emit electromagnetic radiation uniquely based on their molecular composition and texture. If the radiation arriving at a sensor is measured at a detailed wavelength range, the consequent spectral signature, also known as spectrum, can potentially be used to identify any given object of interest. To this end, the intent of hyperspectral imaging technology is to capture hundreds of spectral channels (i.e., to shape the spectra) from the immediate surface of the Earth, which can precisely characterize the chemical composition of different materials.

Hyperspectral sensors mostly sample the reflective portion of the electromagnetic spectrum, ranging from the visible region (0.4-0.7  $\mu\text{m}$ ) to the short-wave infrared region (almost 2.4  $\mu\text{m}$ ) in hundreds of narrow contiguous spectral channels 10nm wide. There are, however, other types of hyperspectral sensors that are able to characterize the emissive properties of objects by collecting data in the range of mid-wave and long-wave infrared.<sup>1</sup> Such detailed spectral sampling, making use of numerous small, commercial, high spatial and spectral instruments, has made hyperspectral images (HSIs) a valuable source of information for a wide variety of applications, including precision agriculture (e.g., monitoring the development and health of crops), the food industry (e.g., characterizing product quality), environmental monitoring, mineralogy, defense and security-based applications (e.g., identification of man-made materials), chemical imaging, astronomy, ecological sciences, and many others.

A better understanding of HSIs can be gained from Fig. 1. A three-dimensional *hyperspectral data cube* consists of  $n_1 \times n_2 \times d$  pixels in which  $n_1 \times n_2$  is the number of pixels in each spectral channel and  $d$  represents the number of spectral channels. An HSI can be defined using one of the following more detailed definitions:

- 1) *Spectral perspective (or spectral dimension)*: From this perspective, a hyperspectral data cube is composed of  $n_1 \times n_2$  pixels, where each pixel is a vector of  $d$  values. Each pixel corresponds to the reflected radiation of the specific region of the Earth and has multiple values in spectral bands. This detailed spectral information can be used to analyze different materials with precision. Fig.

<sup>1</sup>Please note that hyperspectral imaging covers a broad range of imaging systems, such as medical hyperspectral imaging, atmospheric sounding, close-range hyperspectral imaging and so on. Here, we focus solely on airborne or spaceborne remotely sensed hyperspectral images with a spectral coverage ranging from 0.4 to 2.5  $\mu\text{m}$ .

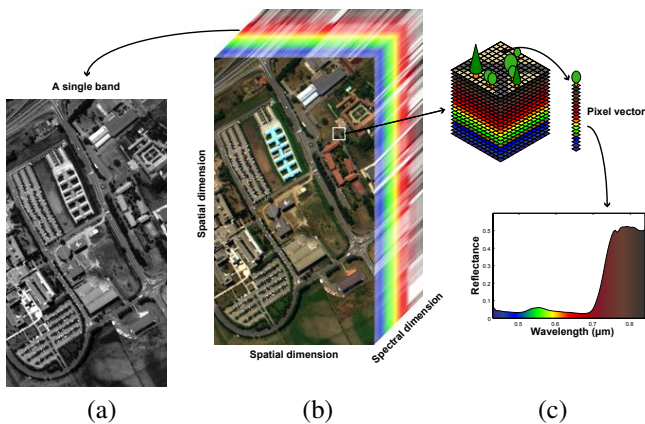


Fig. 1. An example of a hyperspectral data cube: (a) a gray scale image, (b) a hyperspectral data cube, (c) a pixel vector and its corresponding spectral signature.

1(c) shows a spectral profile of one pixel, with multiple values for each band in the spectral dimension.

- 2) *Spatial perspective (or spatial dimension)*: In this context, a hyperspectral data cube consists of  $d$  gray scale images with a size of  $n_1 \times n_2$ . The values of all pixels in one spectral band shape a gray scale image with two dimensions as shown in Fig. 1(a), which are spatial and spatial.

Although the greater dimensionality of HSI than multispectral images improves data information content considerably, it does introduce new challenges to the conventional image analysis techniques which have been specifically designed for multispectral data. Furthermore, it is almost impossible for humans to visualize spaces of higher dimensions than three (e.g., RGB images). A misunderstanding of high-dimensional spaces and conventional spaces sometimes leads to incorrect interpretations of HSI and the inappropriate choice of the data processing technique. Bearing this in mind, in the next subsection, we provide an overview of a few common challenges of HSIs and their possible solutions.

#### A. Main Challenges of HSI Analysis and Possible Solutions

Several factors make the analysis and processing of HSI a challenging task. Fig. 2 illustrates the main paths of HSI analysis, which have been developed primarily to address these factors. In this section, we take a closer look into each of the applications shown in Fig. 2. The common understanding of the HSI is that since such data contain a rich amount of spectral information, the whole dimensionality needs to be used to define precise boundaries in the feature space for a specific applications. The increasing spectral resolution of HSI benefits precision applications (e.g., Earth observation, precision agriculture, disease detection). However, it challenges conventional signal processing techniques and thus hampers its abilities in many real applications.

Taking classification as an example (as classification is one of the most popular applications for HSI), we found in [1] that when the number of training samples remains constant, after a few features, classification accuracy actually decreases

as the number of features increases. Two solutions have been widely exploited to address this problem.

- 1) **Dimension (Feature) Reduction**: As mentioned in several studies, such as [2–4], a high-dimensional space is almost empty and multivariate data can be represented in a lower dimensional space, where the undesirable effects of high-dimensional geometric characteristics and the “curse of dimensionality” are reduced. This fact has led to a chain of research on dimension (feature) reduction, which will be detailed in Section III.

- 2) **Robust Classifiers**: The imbalance between the number of bands and available training samples has a dramatic influence on supervised classifiers. In this context, HSIs often demand a huge number of training samples in order to estimate class parameters effectively. In order to benefit from the rich spectral information of HSI, one possible solution is based on the use of effective and efficient classification approaches that can handle high dimensionality even if a limited number of training samples is available. In addition, along with the detailed spectral information provided by HSIs, it is possible to take advantage of available spatial information (in particular for very high spatial resolution HSIs) to further improve the eventual classification map. Section IV will elaborate on advances in HSI classification.

Spectral mixing (including both linear and nonlinear models) is another bottleneck for analysis of HSI that occurs for a number of reasons, such as insufficient spatial resolution of the sensor and intimate mixing effect. When mixing takes place, it is not possible to distinguish the materials available in the pixels directly from the corresponding measured spectral vectors. However, detailed spectral information provided by HSI can be used to unmix hyperspectral pixels. Section V of this paper is dedicated to **spectral unmixing** to address these issues.

Spaceborne imaging spectrometers are usually designed to acquire HSIs with a moderate spatial resolution (e.g., a ground sampling distance of 30 m) because of inevitable trade-offs among spatial resolution, spectral resolution, temporal resolution, and signal to noise ratio (SNR). Spatial resolution enhancement of HSIs is a technology that is essential to expanding the range of applications for spaceborne hyperspectral missions. In Section VI, we will discuss techniques for the **resolution enhancement** of HSIs.

The degradation mechanisms associated with measurement process and atmospheric effects inject undesirable noise, which substantially downgrades the quality of hyperspectral data. SNR in HSIs is usually decreased during the imaging process depending on different noise sources. In remote sensing HSIs, there are often highly corrupted bands, which are usually removed before any further processing. Alternatively, HSI restoration can recover those corrupted bands and also improve the SNR of HSI, thereby improving the effectiveness of any further processing of the HSI. In this context, Section VII is dedicated to **HSI denoising and image restoration** techniques that address such effects.

Change detection (CD) is considered to be another emerging domain of research in the hyperspectral community. CD is the process of identifying and examining spectral-temporal

changes in signals. The detailed spectral sampling and representation in HSIs result in potential identification of more subtle spectral variations, which are usually not easily detected in traditional multispectral images. Accordingly, land-cover dynamic monitoring can be enhanced to a finer level. To this end, advanced CD techniques need to be designed to address CD issues in multitemporal HSIs, while at the same time overcoming the challenges caused by the hyperspectral data set. We will elaborate on different **change detection** methods in Section VIII.

Another vitally important aspect of HSI analysis that needs to be taken into account with precision is that hyperspectral remote sensors are now in the era of massive automatic data collection resulting from the improved spatial, spectral, and temporal resolutions provided by several hyperspectral instruments. As a result, **fast computing** (detailed in Section IX) is critical to accelerating the efficient exploitation and analysis of HSI.

### B. Missions and Statistics

Several hyperspectral imaging instruments are currently available for the purpose of remote sensing image and signal analysis, providing a large volume of images for various thematic applications. Airborne hyperspectral imaging sensors (e.g., AVIRIS, HYDICE, CASI, APEX, and HySpex) have been playing the main role in acquiring data sources for the hyperspectral scientific community. The European facility for airborne research (EUFAR)<sup>2</sup> has established standards and protocols in the field of airborne hyperspectral remote sensing, allowing transnational access to national infrastructures. In recent years, unmanned aerial vehicle (UAV) based real-time hyperspectral imaging has been becoming more common in various applications such as agricultural monitoring, while raising new challenges in image processing.

Table I presents the principal parameters of seven spaceborne imaging spectroscopy missions planned for the near future: CCRSS, DESIS [5], EnMAP [6], HISUI [7], PRISMA [8], Shalom [9], and HypIRI [10]. Many of those satellites are designed to have a ground sampling distance of 30 m, aiming at global coverage. Shalom's GSDs of 8 m and 10 m in the VNIR and SWIR ranges, respectively, are driven by both operational and commercial needs. DESIS and HISUI will be mounted on the international space station. The launch of these satellites will further accelerate research on hyperspectral image processing and its applications.

Fig. 3 shows statistics on papers related to hyperspectral image and signal processing published in IEEE journals during 2009–2012 and 2013–2016. All papers were searched via *IEEE Xplore* using “hyperspectral” as the main keyword and were categorized by broad topics by analyzing keywords in the titles. The size of each pie is proportional to the number of papers.

The totals returned by this search are an indicator of the hyperspectral community's recent growth. Seven topics under investigation represent 61.5% of all papers during 2013–2016. Classification is the most actively addressed topic in both

periods, while spectral unmixing is the second most common. Classification and unmixing related studies account for 41.6% of the total. The top two topics are followed by dimensionality reduction and image restoration. Image restoration shows a high growth rate, indicating that the improvement in the quality of HSIs is significant in subsequent processing. Resolution enhancement received particular attention during 2013–2016, as demonstrated by the highest growth rate. Although the number of papers related to CD increased steadily, the overall number is still small, probably due to limited data sets.

### C. Contribution

This paper introduces a detailed and organized overview of hyperspectral image and signal processing, which are categorized into seven different themes: dimensionality reduction, classification, spectral unmixing, resolution enhancement, image restoration, change detection, and fast computing. In each section, some numerical results, illustrations, a critical overview on state of the art, current challenges, and possible future works are provided. It is worth noting that the methodologies described or mentioned in this paper are mostly rooted in signal and image processing, statistical inference, and machine learning fields with a particular emphasis on methodologies developed since 2013, after the publication of a previous survey paper on hyperspectral remote sensing image analysis [11].

The rest of the paper is organized as follows. Section II describes three benchmark hyperspectral data sets referred to throughout the paper. Sections III, IV, V, VI, VII, VIII, and IX are dedicated to dimensionality reduction, classification, spectral unmixing, resolution enhancement, HSI denoising and image restoration, change detection, and fast computing, respectively. Finally, Section XI provides concluding remarks.

## II. DATA SETS

This section provides a description of three benchmark hyperspectral data sets (ROSI3-03 Pavia University, CASI Houston University, and Hyperion Umatilla County), which are referenced throughout the paper.

1) *Pavia University*: This data set was captured over the University of Pavia, Italy by the ROSIS-03 (Reflective Optics Spectrographic Imaging System) airborne instrument. The flight over the city of Pavia, Italy, was operated by the Deutsches Zentrum für Luft- und Raumfahrt (DLR, the German Aerospace Agency) within the context of the HySens project, managed and sponsored by the European Union. The ROSIS-03 sensor has 115 data channels with a spectral coverage ranging from 0.43 to 0.86  $\mu\text{m}$ . Twelve channels have been removed due to the existence of noise. The remaining 103 spectral channels are processed. The data have been corrected atmospherically, but not geometrically. The spatial resolution is 1.3 m per pixel. The data set covers the Engineering School at the University of Pavia and consists of different classes, including trees, asphalt, bitumen, gravel, metal sheet, shadow, bricks, meadow, and soil. The subset data set investigated in this review paper comprises  $640 \times 340$  pixels. Fig. 4 presents a false color image of ROSIS-03 Pavia University data and

<sup>2</sup><http://www.eufar.net/>

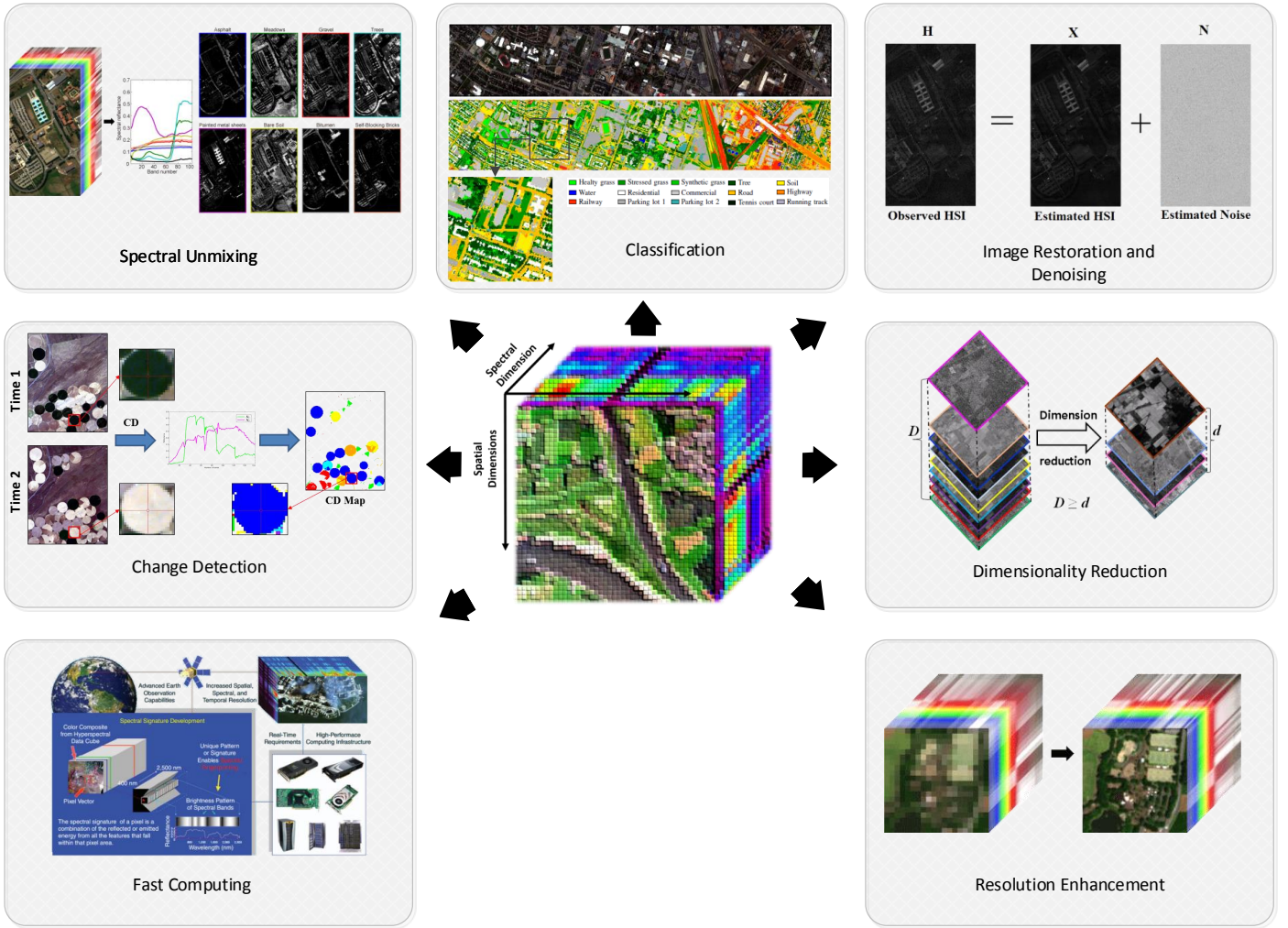


Fig. 2. Different paths of HSI analysis: spectral unmixing, classification, image restoration and denoising, change detection, dimensionality reduction, fast computing, resolution enhancement, and target and anomaly detection.

TABLE I  
PARAMETERS OF SEVEN SPACEBORNE IMAGING SPECTROSCOPY MISSIONS.

Parameter	CCRSS	DESIS	EnMAP	HISUI	PRISMA	Shalom	HypIRI
Altitude (km)	30	400	653	400	615	600	626
GSD (m)	30	30	30	30	30	10	30
Bandwidth (nm)	5–20	3.3	5.25–12.5	10–12.5	$\leq 12$	10	$\leq 10$
Spectral coverage ( $\mu\text{m}$ )	0.4–2.5	0.4–1.0	0.42–2.45	0.44–2.5	0.4–2.5	0.4–2.5	0.38–2.5
Number of bands	328	180	228	185	237	241	210
Swath width (km)	30	30.7	30	20	30–60	10	45
Other sensor	Pan	—	—	—	Pan	Pan	TIR

corresponding training and test samples that have already been separated.

2) *University of Houston*: This data set was captured by the Compact Airborne Spectrographic Imager (CASI) over the University of Houston campus and the neighboring urban area in June, 2012. The size of the data is  $349 \times 1905$  with spatial resolution of  $2.5\text{ m}$ . This data set is composed of 144 spectral bands ranging from  $0.38$  to  $1.05\ \mu\text{m}$ . These data consist of 15 classes, including grass healthy, grass stressed, grass synthetic, tree, soil, water, residential, commercial, road, highway, railway, parking lot 1, parking lot 2, tennis court, and running track. Parking lot 1 includes parking garages at the

ground level and also in elevated areas, while parking lot 2 corresponds to parked vehicles. Fig. 5 shows a three-band false color image and its corresponding already-separated training and test samples.

3) *Umatilla County*: A pair of real bitemporal Hyperion HSIs acquired on May 1, 2004 ( $\mathbf{X}_1$ ) and May 8, 2007 ( $\mathbf{X}_2$ ) were used to test several selected state-of-the-art CD approaches. This scene covers irrigated agricultural land in Umatilla County, Oregon, USA. The images under consideration have a size of  $180 \times 225$  pixels. The original Hyperion images contain 242 spectral bands, ranging from  $0.35$  to  $2.58\ \mu\text{m}$ , i.e., visible, near infrared, and short-wave infrared, with



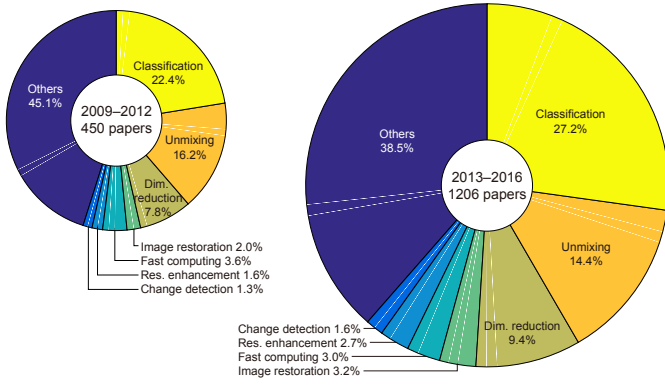


Fig. 3. Statistics on papers related to hyperspectral image and signal processing published in IEEE journals during 2009–2012 and 2013–2016. The size of each pie is proportional to the number of papers. The total number of papers for each time period is shown at the center of each pie chart.

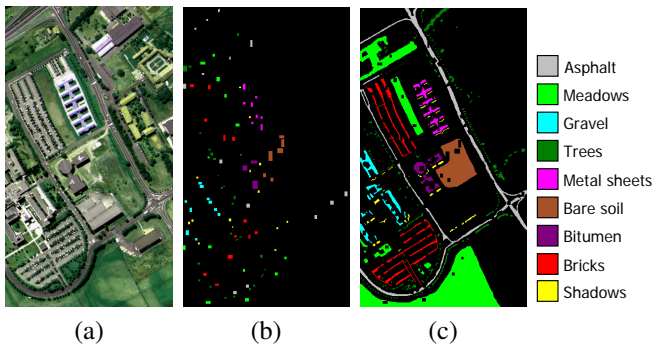


Fig. 4. ROSIS-03 Pavia University - (a) False color composite, (b) training samples, and (c) test samples.

a spectral resolution of  $0.01 \mu m$  and a spatial resolution of  $30 m$ . Preprocessing operations, such as the removal of uncalibrated and noisiest bands, bad stripes repair, atmospheric correction, and co-registration, have been carried out. Finally, 159 preprocessed bands (i.e., 8-57, 82-119, 131-164, 182-184 and 187-220) out of the original 242 bands were used in the CD experiment. Changes occurring in this scenario include the

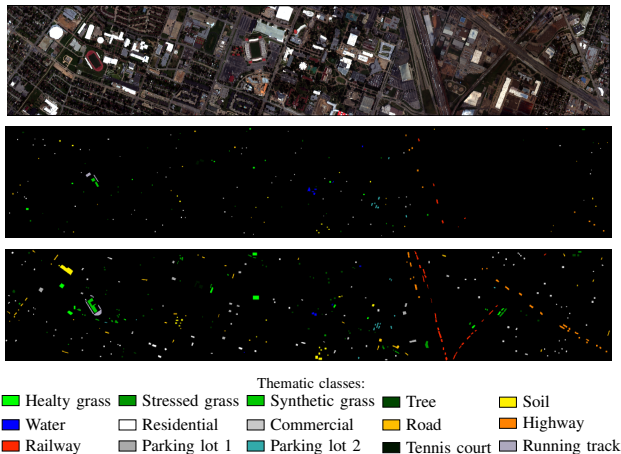


Fig. 5. CASI Houston - (From top to bottom) False color composite (R: band 70, G: band 50, B: band 20), training samples, and test samples.

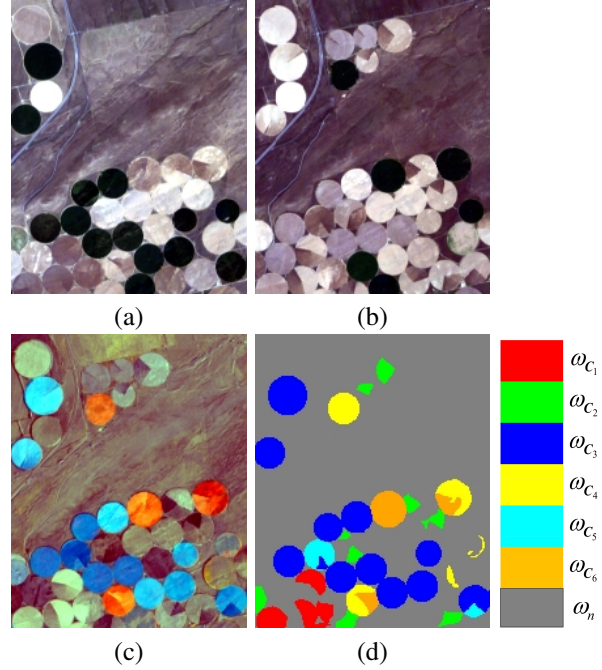


Fig. 6. Umatilla County - False color composite (R: 650.67 nm, G: 548.92 nm, B: 447.17 nm) of the bitemporal EO-1 Hyperion images acquired over an irrigated agricultural area in Umatilla County, OR (USA) in (a) 2004 ( $X_1$ ) and (b) 2007 ( $X_2$ ). (c) Composite of three SCV channels (R: 823.65 nm, G: 721.90 nm, B: 620.15 nm); (d) Multi-class change reference map, where six changes are in different colors, whereas the unchanged pixels are in gray.

land-cover class transitions between crops, bare soil, subtle variations in soil moisture, and water content of vegetation. More detailed descriptions of this data set can be found in [12]. Figs. 6 (a) and (b) show the false color composite of  $X_1$  and  $X_2$ , respectively. The false color composite of three spectral change vector (SCV) channels is shown in Fig. 6 (c); possible different changed pixels are illustrated in different colors, whereas the unchanged pixels are in gray. The multi-class change reference map is made based on careful image interpretation, as shown in Fig. 6 (d). Note that the possible subtle sub-pixel level changes (e.g., the one associated with the road surrounding the irrigated agricultural land [12]) was not considered in this paper in order to conduct the quantitative comparison with other pixel-level-based approaches fairly. Thus six pixel-level changes were considered, as shown in Fig. 6 (d).

### III. DIMENSIONALITY REDUCTION

The increasing spectral resolution of hyperspectral data benefits precision pattern recognition, but challenges both the memory capacity of ordinary personal computers and conventional signal processing techniques. For an HSI with spatial dimension of  $600 \times 400$  pixels at 16 bits-per-band-per-pixel, the data volume becomes 240 MB for 500 spectral bands. The data volume can be linearly increased when time series hyperspectral data are acquired to monitor environmental changes. The complexities of storing and processing the data will easily exceed the memory capacity of ordinary personal computers. Moreover, as discussed above, when the ratio between the spectral bands and the number of training

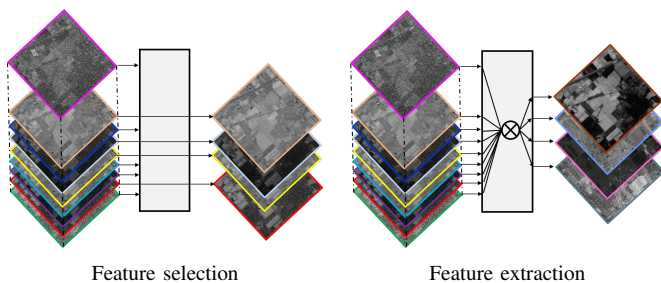


Fig. 7. Hyperspectral image dimensionality reduction.

samples is high, high-dimensional hyperspectral data suffer from the well-known issue of the curse of dimensionality. Dimensionality reduction (DR), aiming at identifying and eliminating statistical redundancies of hyperspectral data while keeping as much spectral information as possible, is widely used in hyperspectral data processing. Relatively few bands can represent most of the information in HSIs [13], making DR very useful for storage, transmission, classification, spectral unmixing, target detection [14], and visualization of remote sensing data [13, 15]. Recent work demonstrates the benefits of using dimensionality reduction when extracting relevant information from hyperspectral images for change detection [16], forest management [17], and urban planning [18]. The applications of dimensionality reduction are of interest well beyond hyperspectral data, for various applications of signal processing and computer vision [19], and wherever interpretation and analysis of high-dimensional data is of interest.

Hyperspectral DR consists of both feature selection (FS) and feature extraction (FE) [13]. FS tries to select a minimal subset of  $D$  features  $S = \{S_1, S_2, \dots, S_D\}$  from the original feature set  $F = \{F_1, F_2, \dots, F_d\}$  based on an adopted selection criterion, where  $D \leq d$  and  $S \subseteq F$ , while aiming to achieve improved performances for a specific application (e.g., classification, target detection, etc.). The objective of FE is to find a transformation function  $f : \mathbb{R}^d \rightarrow \mathbb{R}^D$  that can transform the high-dimensional data point  $\{\mathbf{x}_i \in \mathbb{R}^d\}_{i=1}^N$  to  $\mathbf{z}_i = f(\mathbf{x}_i)$ , where  $\{\mathbf{z}_i \in \mathbb{R}^D\}_{i=1}^N$  and  $D \leq d$ , such that most information of the high dimensional data is kept in a much lower dimensional subspace. The term  $f$  can be a linear or nonlinear transformation. Unlike FS, FE compresses the high-dimensional original data to generate a small number of new features, where each band often has a contribution to determining  $f$ , as shown in Fig. 7. DR methods can be categorized into unsupervised, supervised, and semisupervised approaches, depending on whether the class label information is being used. Each of these methods is discussed further below.

#### A. Unsupervised DR

Unsupervised DR methods deal with the cases where no labeled samples are available, aiming to find another representation of the data in the lower dimensional space by satisfying some given criterion. A variety of unsupervised DR methods has been introduced in the literature. The objective of these methods is not to optimize the accuracy for a given

classification task, since they do not consider class-specific information provided by labeled samples. For examples, principal component analysis (PCA) [20] reduces dimensionality by capturing the maximum variance in the data. Independent component analysis (ICA) [21] finds the project matrix by maximizing the statistical independence. Minimum noise fraction transform (MNF) [22] obtains the reduced features according to image quality measured by SNR, and local linear feature extraction (LLFE) [23–25] methods seek a projection direction in which neighborhood relationships are preserved in the feature spaces. The non-linear versions of these methods, such as kernel methods (e.g., kernel PCA, kernel ICA, kernel MNF [26]) and local methods (e.g., locally linear embedding [27], Laplacian eigenmap, and local tangent space alignment [19]) have been widely used to detect higher-order statistical redundancies. In the same manner, conventional unsupervised FS methods for DR select a subset of features from the original data according to a specific criterion, such as linear prediction error [28], entropy [29], or mutual information (by minimizing dependency) [30].

Recently, fusion-based methods and manifold learning methods have been widely explored for HSI unsupervised DR. Graph-based fusion methods couple data fusion and dimensionality reduction in a unified framework for classification [31, 32]. Borhani and Ghassemian present a kernel-based method to incorporate spectral and spatial information simultaneously for dimensionality reduction and classification of hyperspectral data [33], while Zhang et al. represent multiple features in a low-dimensional feature space where the complementary information of each feature is exploited by co-manifold learning and co-graph regularization [34]. In the approaches of [35], manifold learning has been exploited for feature extraction and salient band selection of HSIs. In [36], orthogonal total variation component analysis (OTVCA) is proposed, where a non-convex cost function is optimized to find the best representation for HSI in a low dimensional feature space while controlling the spatial smoothness of the features by using a total variation (TV) regularization. The TV penalty promotes piecewise smoothness (homogeneous spatial regions) on the extracted features, and thus helps to extract spatial (local neighborhood) information that is very useful for classification. It is shown that OTVCA is highly robust to noise because it exploits a penalized least squares minimization framework.

#### B. Supervised DR

Supervised methods rely on the existence of labeled samples to infer class separability. Several widely used supervised DR methods for HSIs are linear discriminant analysis (LDA) [37] and nonparametric weighted feature extraction (NWFE) [38], band selection based on Jeffries-Matsushita (J-M) distance [39], and mutual information (MI) [40]. Many extensions to these methods have been proposed in past decades, including modified Fishers LDA [41] and regularized LDA [42], modified NWFE using spatial and spectral information [43] and kernel NWFE [44], extended J-M to multiclass cases [40] and J-M distance for spatially invariant features [45], minimal-

redundancy-maximal-relevance based on mutual information [46], and normalized mutual information [47].

Recent supervised DR methods for hyperspectral data exploit the local neighborhood properties of data. Li et al. [48] employed local Fishers discriminant analysis [49] to reduce the dimensionality of the data while preserving the corresponding multimodal structure. In [50], local neighborhood information was exploited in both spectral and spatial domains to find a discriminative projection for DR of hyperspectral data. Cao et al. [51] proposed a supervised band selection, by introducing the local spatial smoothness of the HSI into the wrapper method. Dong et al. [52] presented an ensemble discriminative local metric learning method for DR, where local spatial information was incorporated into distance metric learning to learn a subspace, keeping the samples from the same class closer while pushing those from different classes further away. Sparse graph embedding (SGE) explores the sparsity structure of the data for hyperspectral DR. Ly et al. [53] proposed block sparse graph based discriminant analysis, which learns a block sparse graph for a supervised DR. Xue et al. [54] proposed a spatial and spectral regularized local discriminant embedding method for DR, where spatial information was integrated into the sparse graph learning process. In [55], a discriminative sparse multimodal learning is developed for multiple feature selection. However, the sparse coding used in SGE is helpful for learning on conditions where the coding is local [56], which means locality is more important than sparsity. Unfortunately, the converse is not true: sparsity does not always guarantee locality [56]. He et al. [57] proposed a weighted sparse graph to overcome the drawback of sparse coding in SGE, where both the locality and sparsity structure of the training pixels are integrated.

Other trends in supervised DR methods exploit various algorithms and learning techniques from soft computing, artificial intelligence, and machine learning. Genetic algorithms (GA) [58], particle swarm optimization (PSO) [59], and the combination of GA and PSO are used to optimize feature selection [60, 61]. Deep learning techniques, e.g., stacked autoencoder [62] and convolutional neural network [63], are used for spectral-spatial feature extraction for hyperspectral image classification [64, 65].

### C. Semisupervised DR

In real-world applications, labeled data are usually very limited and labeling a large amount of data may sometimes require considerable human resources or expertise. On the other hand, unlabeled data are available in large quantities at very low cost. For this reason, semisupervised methods [66–68], which aim at improved classification by utilizing both unlabeled and limited labeled data, have gained popularity in the machine learning community. Some of the representative semisupervised learning methods include Co-Training [66] and transductive SVM [67], and graph-based semisupervised learning methods [68]. Some semisupervised feature extraction methods add a regularization term to preserve certain potential properties of the data. For example, semisupervised discriminant analysis (SDA) [69] adds a regularizer into the objective

function of LDA. The resulting method makes use of a limited number of labeled samples to maximize class discrimination, and employs both labeled and unlabeled samples to preserve the local properties of the data. The approach of [70] proposed a general semisupervised dimensionality reduction framework based on pairwise constraints, and employs regularization with sparse representation. A semisupervised pairwise band selection method [71] was proposed for HSIs, in which an individual band selection process is performed only on each pair of classes. Other semisupervised feature extraction methods combine supervised methods with unsupervised ones using a trade-off parameter, such as semisupervised local Fisher discriminant analysis (SELF) [72].

However, it may not be easy to specify the optimal parameter values in these and similar semisupervised techniques, as mentioned in [70, 72]. Liao et al. [73] proposed a semisupervised local discriminant analysis (SELD) to overcome this problem by combining unsupervised methods (LLFE [23–25]) and a supervised method (LDA [37]) in a novel framework without any free parameters. They found an optimal projection matrix that preserves the local neighborhood information inferred from unlabeled samples, while simultaneously maximizing the class discrimination of the data inferred from the labeled samples. The approach of [74] improved SELD [73] by better modeling the differences and similarities between samples. Specifically, this method built a semisupervised graph, where labeled samples were connected according to their label information and unlabeled samples by their nearest neighborhood information. Graph embedding and manifold-based sparse representation were combined in a semisupervised framework for hyperspectral DR [75], where the sparse coefficients were exploited to construct the graph. Semisupervised manifold alignment [76] and semisupervised transfer component analysis [77] were proposed to find a transform matrix to project high-dimensional multimodal images into a lower-dimensional feature space, where the geometry of each modality can be preserved.

These semisupervised DR methods try to build a similar objective function, i.e., maximizing class discrimination while at the same time preserving the intrinsic geometric structure of the data. The optimal solutions are acquired by solving generalized eigenvalue problems in the same manner [78]. These methods can be further expanded to shape a manifold learning method by using the kernel trick, similarly to the approaches in [79].

### D. Experimental Results

Table II and Fig. 8 show the performances of some DR methods on the classification of the Pavia University HSI. We compare the performances when using raw hyperspectral data and seven DR methods (including unsupervised, supervised, and semisupervised DR methods) with three popular classifiers, whose parameter settings are the same as those in [74]. The training samples were randomly selected from the training set, with the sample size corresponding to different cases: 20, 40, and 80 samples per class, respectively. The results were averaged over 10 runs on different numbers of

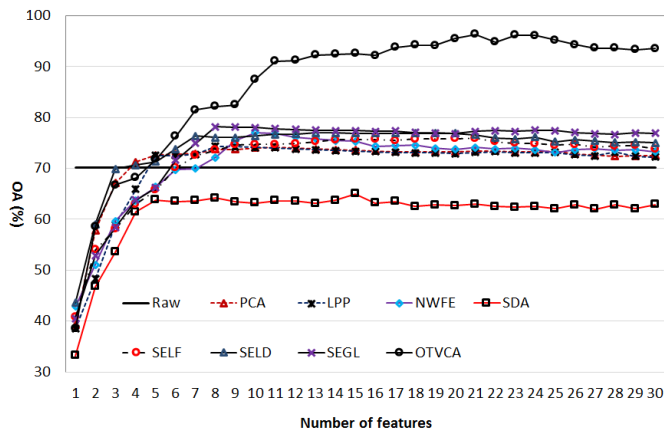


Fig. 8. The classification performance on Pavia University data as the number of features increases. Forty labeled training samples per class are randomly selected from the training set with an SVM classifier.

TABLE II  
DIMENSIONALITY REDUCTION FOR CLASSIFICATION OF PAVIA UNIVERSITY DATA, OA% (OPTIMAL NUMBER OF REDUCED FEATURES).

DR methods		Classifier	Number of labeled samples		
			20	40	80
Unsupervised	Raw	INN	63.14	68.36	69.88
		SVM	67.29	70.15	71.51
		RF	68.72	71.06	73.39
	PCA	INN	67.38(11)	69.46(12)	70.23(11)
		SVM	70.21(8)	74.77(8)	76.81(9)
		RF	73.46(11)	73.88(11)	76.13(11)
	LPP [24]	INN	66.97(18)	69.43(18)	70.17(12)
		SVM	71.80(9)	77.14(12)	77.95(18)
		RF	72.46(18)	74.20(19)	75.82(19)
	OTVCA [36]	INN	71.93(19)	75.81(20)	80.39(23)
		SVM	92.74(19)	96.32(21)	97.52(18)
		RF	<b>96.24(23)</b>	<b>97.75(21)</b>	<b>98.79(22)</b>
Supervised	NWFE [38]	INN	71.34(9)	73.34(9)	74.42(12)
		SVM	72.29(8)	76.62(8)	77.18(9)
		RF	75.84(9)	77.03(11)	78.74(8)
Semisupervised	SDA [69]	INN	52.67(7)	62.64(8)	70.76(8)
		SVM	51.62(8)	63.64(9)	70.96(10)
		RF	55.72(9)	66.97(9)	71.34(9)
	SELF [72]	INN	61.42(18)	68.75(18)	69.56(16)
		SVM	63.93(19)	75.85(19)	82.74(18)
		RF	67.98(18)	75.97(12)	80.56(12)
	SELD [73]	INN	77.27(17)	79.03(10)	81.95(11)
		SVM	75.71(11)	76.20(10)	82.03(11)
		RF	75.53(8)	77.50(9)	82.26(9)
	SEGL [74]	INN	79.40(9)	81.22(9)	83.57(10)
		SVM	77.57(8)	78.16(8)	82.26(8)
		RF	76.67(8)	79.14(8)	83.51(9)

extracted features from 1 to 30, and the averaged overall classification accuracy (OA) was recorded for each method. The results confirm that dimensionality reduction can improve classification performance on HSIs. As the size of the training sample increases, classification accuracy increases. Semisupervised DR methods (especially for [73, 74], designed for hyperspectral data) outperform both unsupervised and supervised methods for the INN classifier. DR methods that exploit spatial smoothness produce better results, even for unsupervised methods; for example, OTVCA [36] outperforms the other methods for both RF and SVM classifiers in terms of classification accuracy. We can also find from Fig. 8 that, as the number of features increases, the classification performances

do not always increase—in fact, some will decrease. In order to get the optimal classification performance, the number of reduced features needs to be optimized.

### E. Challenges in Dimensionality Reduction

Recent advances in sensor technologies and processing techniques strongly support the use of hyperspectral data. Moreover, global Earth observation missions (e.g., AVIRIS from NASA, PROBA series from European Space Agency, and the Gaofen series from China) make such data increasingly accessible. Furthermore, at lower altitudes, airplanes and Unmanned Aerial Vehicles (UAVs) can deliver very high resolution hyperspectral data from targeted locations. On the other hand, image processing techniques allow us to extract multiple level features from these ‘Big Hyperspectral Data.’

Two main challenges remain in hyperspectral dimensionality reduction: 1) mining complementary features (while reducing the dimension and redundancy) from multiple levels of ‘Big Hyperspectral Data’; 2) coupling dimensionality reduction and applications in a unified framework, ensuring that optimal features for applications are obtained. Most state-of-the-art research has separated dimensionality reduction and applications into two different steps. For example, morphological operators were employed in [80] to extract low-level features (such as size and shape of objects) from remote sensing images. In [81, 82], middle-level attribute features were extracted from HSIs for land-cover mapping. High-level features, such as object-based [83] and so-called deep learning features [84], have been used for change detection and classification. State-of-the-art DR methods typically deal with either lower-level or higher-level features, but not with a combination of both. The features extracted at each level have their own characteristics: high-level features are usually more powerful but less robust, while the low-level ones are less informative but more robust. On the other hand, classification is taken as one of the most popular applications to validate the performances of dimensionality reduction. Hyperspectral classification typically consists of two steps: (1) dimensionality reduction (via either feature extraction or feature selection) and (2) a training procedure for designing the classifier. However, it is difficult to ensure the best features from the first step to optimize the classification performance of the following one.

## IV. CLASSIFICATION

HSI classification is a fast-growing and highly active field of research in the hyperspectral community. A *classification algorithm* is utilized to distinguish between different land-covers by assigning unknown pixel vectors to one of the classes (or clusters). The individual classes are commonly differentiated based on the similarity to a certain class or by defining decision boundaries, which are constructed in the feature space. The initial set of features for classification usually encompasses spectral channels [4].

With reference to Fig. 1, two types of classification approaches can be broadly defined: spectral classifiers and spectral-spatial classifiers [4], where the former consider the HSI to be a list of spectral measurements with no spatial



organization, while the latter classify the input data by taking into account the spatial dependencies of adjacent pixels. Below, after discussing each of those two classification approaches, we take a glance at a few advanced classification approaches based on composite kernels, semisupervised and active learning, and sparse representation-based classifiers. Finally, a brief discussion of the main challenges in classifying HSI is provided.

### A. Spectral Classifiers

Based on the availability of *training samples* (also referred to as learning with a teacher) for the training stage, classification approaches can be grouped into three categories: supervised, unsupervised (also known as clustering), and semisupervised approaches.

Supervised approaches classify input data using *training samples*. These samples are usually collected in one of two ways: (1) the manual labeling of a small number of pixels in an image, or (2) based on some field measurements. In contrast, unsupervised classification does not consider training samples. The supervised type of approach classifies input data based only on an arbitrary number of initial “cluster centers” that may be either user-specified or selected quite arbitrarily. During the processing, each pixel is associated with one of the cluster centers, usually in an iterative way, based on a similarity criterion [85, 86]. In addition to unsupervised and supervised approaches, semisupervised techniques have been considered [87]. In such approaches, the training stage is based not only on labeled training samples, but also on unlabeled samples.

Since the consideration of training samples leads to higher classification accuracies than situations where there is no class-specific information, supervised approaches have gained more attention in the hyperspectral community than unsupervised ones. However, the curse of dimensionality is a bottleneck for supervised classification techniques. In theory, a large number of training samples is required to define precise class boundaries in the feature space. This problem intensifies when the number of bands (features) increases. However, in practice, there are not enough training samples to train supervised classifiers since the collection of such samples is time-consuming and/or costly. Therefore, classification approaches developed for HSIs need to be able to handle high dimensional data with only a limited number of training samples. The most widely used supervised spectral classifiers have been studied precisely and compared in [88]. Table III demonstrates classification accuracies—that is, overall accuracy (OA), average accuracy (AA), and kappa coefficient—obtained on the University of Houston data set by a number of widely-used supervised spectral classifiers in the hyperspectral community, including support vector machines (SVM) [89], random forest (RF) [90], rotation forest (RoF) [91], canonical correlation forest (CCF) [92, 93], back-propagation neural network (BP) [94], extreme learning machines (ELM) [95], kernel ELM (KELM) [96], 1D deep convolutional neural network (1D CNN) [84], and MLR [97]. For the algorithm setup, please see [88, 93].

Nowadays, a major contribution in the hyperspectral community is based on the use of deep learning for HSI classifica-

TABLE III  
SUPERVISED SPECTRAL CLASSIFIERS: CLASSIFICATION ACCURACIES  
OBTAINED ON THE UNIVERSITY OF HOUSTON DATA.

Class	SVM	RF	RoF	CCF	BP	ELM	KELM	1D-CNN	MLR
OA	80.1	72.9	79.1	83.3	80.9	79.5	80.6	78.2	80.6
AA	83.0	76.9	82.0	85.7	83.1	82.4	82.9	81.2	83.0
Kappa	0.786	0.709	0.775	0.820	0.793	0.778	0.790	0.784	0.790

tion. HSIs are highly influenced by various atmospheric scattering conditions, complicated light scattering mechanisms, inter-class similarity, and intra-class variability, which make the hyperspectral imaging procedure inherently nonlinear [65]. Compared to the so-called “shallow” models, deep learning approaches are expected to potentially extract high-level, hierarchical, and abstract features, which are, by nature, more robust when handling the nonlinearities of the input hyperspectral data. Although the use of deep learning in the hyperspectral community is in its early days, some contributions in the community have focused on the use of deep learning for HSI classification. A stacked auto-encoder (SAE) and auto-encoder with sparse constraint were proposed for HSI classification [98, 99], where hierarchical features were extracted from the input data. Another deep model, the deep belief network (DBN), was proposed for the classification of hyperspectral data by learning spectral-based features [100]. The critical comparison conducted in [88], specifically on supervised spectral classifiers, offered tantalizing hints about the logical selection of an appropriate classifier based on the application at hand. One of the main conclusions was that there is no classifier that can consistently provide the best performance in terms of classification accuracy when different data sets or different sets of training and test samples are considered. Instead, in addition to resulting classification accuracies, the consideration of an appropriate classifier should be based on the complexity of the analysis scenario (e.g., availability of training samples, processing requirements, tuning parameters, speed of the algorithm, etc.) and on the considered application domain.

### B. Spectral-Spatial Classifiers

Neighboring pixels in HSIs are highly related or correlated, since remote sensors acquire a significant amount of energy from adjacent pixels, and homogeneous structures in the image scene are generally larger than the size of a pixel. This is especially evident for images of high spatial resolution. Spatial and contextual information can provide useful information about the shape of different structures. In addition, such information reduces the labeling uncertainty that exists when only spectral information is taken into account, and helps to address the “salt and pepper” appearance of the resulting classification map. In general, spectral-spatial classification techniques are composed of three main stages:

- 1) Extracting spectral information (i.e., based on spectral classifiers discussed in Section IV-A);
- 2) Extracting spatial information (to be discussed later in this section);

- 3) Combining the spectral information extracted from (1) and spatial information extracted from (2).

In order to extract spatial information, two common strategies are available: the crisp neighborhood system and the adaptive neighborhood system. While the former considers spatial and contextual information in a predefined neighborhood system, the latter is more flexible and is not confined to a given neighborhood system. In the following two subsections, each neighborhood system will be briefly explained. It should be noted that these methods have been elaborated in detail in the book [4]. Table IV demonstrates several classification accuracies obtained on the Pavia University data set by different spectral-spatial classification approaches, which will be briefly discussed below.

1) **Crisp Neighborhood Systems:** Markov random fields (MRFs) are a family of probabilistic models that can be described as a 2-D stochastic process over discrete pixels lattices. MRFs have been widely used to integrate spatial context into image classification problems. In this family of approaches, it is assumed that for a predefined neighborhood of a given pixel, there is a high possibility that its closest neighbors belong to the same object. In [101], a classification framework was introduced by integrating SVM and MRF. The developed contextual generalization of SVMs was achieved by analytically relating the Markovian minimum-energy criterion to the application of an SVM. In [102], Ghamisi et al. proposed a spectral-spatial classification approach based on a generalization of the MRF called hidden MRF (HMRF). In that work, spectral and spatial information was extracted by using SVM and HMRF, respectively. Finally, the spectral and spatial information was combined via majority voting within each object. Xia *et al.* [103] integrated MRFs with the RoF classifier to further improve classification accuracy.

Another way of considering spatial information using the crisp neighborhood system is based on 2-D or 3-D deep CNNs [84]. CNNs consider local connections to deal with spatial dependencies using sharing weights, which can significantly reduce the number of parameters of the network compared to its 1-D fully connected version and extract spatial and contextual information using a predefined crisp neighborhood system [84]. In [104], an unsupervised approach was introduced to learn feature extracting frameworks from unlabeled hyperspectral imagery. That method extracts generalizable features by training on sufficiently large quantities of unlabeled data that are distinct from the target data set. The trained network is then able to extract features from smaller labeled target data sets and address the curse of dimensionality. In [65], a self-improving CNN-based approach was proposed for the classification of hyperspectral data. This approach solves the curse of dimensionality and the lack of available training samples by iteratively selecting the most informative bands suitable for the designed network. Table IV demonstrates the classification accuracies obtained by **SICNN** [65], and **2D CNN** [84] on the Pavia University data set. As can be seen, in all cases, the use of crisp neighborhood system-based spectral-spatial classification can improve the classification accuracy of spectral classifiers (e.g., **RF** and **SVM**). However, considering a set of crisp neighbors has some disadvantages:

- 1) The crisp neighborhood system may not contain enough samples, which downgrades the effectiveness of the classifier (particularly when the input data set is of high resolution and the neighboring pixels are highly correlated).
- 2) A larger neighborhood system may lead to intractable computational problems. Unfortunately, the closest fixed neighborhoods do not always accurately reflect information about spatial structures. For instance, they provoke assimilation of regions containing only a few pixels with their larger neighboring structures and do not provide accurate spatial information at the border of regions.
- 3) In general, the use of a crisp neighborhood system leads to acceptable results for big regions in the scene. Otherwise, it can make small structures in the scene disappear, merging them with bigger surrounding objects.
- 4) They may cause oversmoothing on the border of different classes. This problem, however, has been addressed using a gradient step in [102].

2) **Adaptive Neighborhood Systems:** In order to address the shortcomings of crisp neighborhood systems, an adaptive neighborhood system can be considered. One approach is to take advantage of different types of segmentation methods. Image segmentation is the process of partitioning a digital image into multiple non-overlapping regions or objects. In image segmentation a label is assigned to each pixel in the image such that pixels with the same label share certain visual characteristics [105]. These objects provide more information than individual pixels, since the interpretation of images based on objects is more meaningful than interpretations based on individual pixels.

For spectral-spatial classification of HSI using segmentation approaches, there are usually two methods to consider: (1) Segmentation and classification maps can be integrated using majority voting within each object by assigning the whole object to the most frequent classification label within that particular object [106],<sup>3</sup> and (2) segments can be considered to be input vectors for supervised classification [108]. In [109], however, a reverse view was employed where markers for spatial regions are automatically obtained from classification results and then used as seeds for region growing in the segmentation step. The classification accuracy of this segmentation method with an extra step (where the classification map is refined using the results of a pixelwise classification and a majority voting within the spatially connected regions) is shown in Table IV as **MSF**.

One common way to segment an image is based on histogram thresholding. A commonly used exhaustive search for optimal thresholds in terms of between-class distances is based on the Otsu criterion [110]. The approach is easy to implement, but it has the disadvantage of being computationally expensive. An exhaustive search for an  $n$ -level segmentation (i.e.,  $n - 1$  optimal thresholds) involves evaluations of fitness of  $n(L - n + 1)^{n-1}$  combinations of thresholds where  $L$  shows the number of intensity values. Therefore, this method is not suitable from a computational cost perspective. The task of determining  $n - 1$

<sup>3</sup>Majority voting within each object is described in [107].

optimal thresholds for  $n$ -level image thresholding could be formulated as a multidimensional optimization problem. In [106], a thresholding-based segmentation method was proposed, where an evolutionary-based optimization technique, named fractional order Darwinian particle swarm optimization (FODPSO), sought to find the best set of thresholds with the highest between-class distance. The classification accuracy obtained by this segmentation method is given in Table IV as **FODPSO**. This method is very fast, even for large data sets, since it works on the image histogram instead of the image space.

Morphological profiles (MPs) are another set of approaches based on adaptive neighborhood pixels. MPs comprise a number of features constructed by applying a set of openings and closings by reconstruction with a structuring element (SE) of an increasing size [111]. The result of the basic extension of the MP that is applicable to HSI is shown in Table IV as **EMP**, or extended MP. Although the MP is a powerful approach for the extraction of spatial information, the concept suffers from a few limitations:

- The shape of SEs is fixed, which imposes a constraint on model spatial structures within a scene.
- SEs are unable to characterize information about the gray-level characteristics of the regions, such as spectral homogeneity, contrast, and so on.

To address these shortcomings of the MPs, a morphological attribute profile (AP) was introduced, which provides a multilevel characterization of an image by using the sequential application of morphological attribute filters (AFs) [112]. A comprehensive survey on the use of APs for the classification of HSI can be found in [4, 113]. The classification accuracy obtained by the extension of AP on HSI, known as extended multi-AP (EMAP), is given in Table IV as **EMAP**. The main difficulties of using the AP, however, are not knowing (1) which attributes lead to a better discrimination ability for different classes, and 2) which threshold values should be considered to initialize each AP. To solve these issues, a few papers have tried to introduce automatic techniques for the use of APs, such as [114–116]. In [116] and [115], automatic spectral-spatial classification methods have been proposed based on the use of EMAP and supervised/unsupervised feature extraction approaches. The classification accuracy of (**APDAFE**) [116] and (**DBAPDA**) [115] are shown in Table IV.

MPs and APs produce extremely redundant features. To address this issue, a sparse classification using both spectral and spatial information was investigated in [117]. In [118], the performance of different feature extraction approaches, including linear, nonlinear, and manifold approaches, has been investigated to generate base images to construct EMAPs.

To further improve the conceptual capability of the AP and the corresponding classification accuracies, Ghamisi *et al.* proposed extinction profiles (EPs) in 2016 [119] by considering a set of connected idempotent filters and extinction filters. In contrast with the AP, the EP preserves the height of the extrema [119] and as a result shows better capability than the AP in terms of simplification for recognition. This advantage leads to higher classification accuracy for EPs than

for APs. In addition, EPs' parameters can be set automatically, independent of the kind of the attribute being used (e.g., area, volume, and so on). In other words, the main issue of conventional APs, the initialization of the threshold values, is solved by EPs [119]. In [120], the concept of EPs has been generalized to extract spatial and contextual information from HSIs, known as extended multi-EP (EMEP). The classification accuracy of EMEP using RF (**RF-EMEP**) and RoF (**RoF-EMEP**) are presented in Table IV.

### C. Composite Kernels

The main problem associated with the concept of spectral-spatial feature extraction approaches is that they usually increase the number of features, while the number of training samples remains the same. This can lead to the curse of dimensionality and high executable processing time. This problem has partially been addressed by combining different kernels for spectral and spatial information (i.e., *composite kernels*) [121] in the SVM classification process. However, classification using composite kernels and SVMs demands convex combination of kernels and a time-consuming optimization process. Therefore, the approach has been modified to deal with convex combinations of kernels through generalized composite kernels (results are shown as **GCK** in Table IV) [122] and multiple-kernel learning [123]. In [124], a classification framework was introduced that combines multiple features with linear and nonlinear class boundaries present in the data without requiring any regularization parameters to control the weights of considered features (results are shown as **MFL** in Table IV).

### D. Semisupervised and Active Learning

As discussed above, the number of training samples is usually limited, as the collection of such samples is either expensive or time-consuming. In such situations, the limited number of training samples available may not be representative of the statistical distribution of the data, which can downgrade the quality of the classification map obtained by supervised classifiers. To partially address this issue, active learning, which aims to find the most informative training set, has gained popularity in the hyperspectral community. Active learning starts an iterative process with a small and suboptimal initial training set and then selects a few additional samples from a large quantity of unlabeled samples. Active learning considers the result of the current model to rank the unlabeled samples according to a criterion that allows selection of the most informative samples to improve the model, thus minimizing the number of training samples while preserving discrimination capabilities as much as possible [125]. For a complete survey on the use of active learning for remote sensing image analysis, see [126] and [127].

Active learning and semisupervised learning share a similar conceptual background, since both types of learning try to address the issue of limited labeled samples. In this manner, both approaches start with a small set of labeled samples and a large set of unlabeled data. Active learning usually requires a labor-intensive human-labeling process, while semisupervised

TABLE IV  
SUPERVISED SPECTRAL-SPATIAL CLASSIFIERS: CLASSIFICATION ACCURACIES OBTAINED FROM PAVIA UNIVERSITY HYPERSPECTRAL DATA.

Class	RF	SVM	2D CNN	SICNN	FODPSO	MSF	EMP	EMAP	APDAFE	DBAPDA	RF-EMEP	RoF-EMEP	GCK	MFL
OA	71.3	78.8	78.8	83.4	88.1	91.1	77.7	90.7	97.0	98.0	96.1	96.3	98.0	97.5
AA	82.2	87.0	79.7	83.0	92.0	94.8	82.5	91.4	96.7	98.1	96.6	97.9	97.4	97.1
Kappa	0.648	0.735	0.734	0.778	0.848	0.880	0.710	0.877	0.960	0.974	0.949	0.952	0.974	0.967

learning, although avoiding human labeling by assigning pseudolabels to unlabeled data, may introduce incorrect pseudolabels and consequently downgrade classification performance [128]. Although active learning and semisupervised learning follow different work flows, they both aim to make the most of unlabeled data while reducing human-labeling efforts [125]. Therefore, it is common to use both of these strategies to make the most of these two paradigms for HSI classification. In [128], active learning and semisupervised learning were collaboratively integrated to form an approach called collaborative active and semisupervised learning that improves pseudo-labeling accuracy and thus facilitates semisupervised learning. This method was based on spectral information. In [125], active learning and hierarchical segmentation (HSeg) are combined for spectralspatial classification of HSI.

#### E. Sparse Representation Classification

Sparse representation classification (SRC)-based approaches with dictionary-based generative models [129, 130] have received considerable attention in the hyperspectral community. In this context, an input signal is represented by a sparse linear combination of samples (*atoms*) from a dictionary [129], where the training data is generally used as the dictionary. The main advantage of such approaches is that SRCs avoid the heavy training procedure that a supervised classifier generally conducts, and the classification is performed directly on the dictionary. Classification can be improved by incorporating contextual information from the neighboring pixels into the classifier. This can be done indirectly by exploiting the spatial correlation through a structured sparsity prior imposed in the optimization process. If an adequate number of training samples is available, discriminative as well as compact class dictionaries can also be developed to improve classification performance [131].

#### F. Challenges in Classification

The main challenges for the classification of HSI are not particularly related to methodology. They are, however, related to the lack of appropriate benchmark data sets and the corresponding training and test samples. As can be seen in Fig. 3, the highest numbers of contributions in the hyperspectral community is dedicated to HSI classification. The approaches are often capable of producing very accurate classification maps on the widely-used Indian Pines and Pavia data sets, which makes the real comparison of the approaches almost impossible. In other words, the existing data sets have already been saturated in terms of classification accuracies. Therefore, our community is in urgent need of more complex data sets

to share (e.g., highly nonlinear data sets with greater area coverage that are composed of many classes). In addition, a standard set of training and test samples should be defined for each particular data set, to make the proposed approaches fully comparable with each other.

### V. SPECTRAL UNMIXING

Spectral unmixing has been an alluring exploitation goal since the early days of hyperspectral image processing [132]. Mixed pixels are common in remotely sensed HSIs due to insufficient spatial resolution of the imaging spectrometer, or due to intimate mixing effects. However, the rich spectral resolution available in hyperspectral data cubes can be used to unmix hyperspectral pixels. In fact, mixed pixels can also be obtained with high spatial resolution data due to intimate mixtures. This means that increasing the spatial resolution often does not solve the problem. In other words, the mixture problem can be approached in a macroscopic fashion, which means that only a few macroscopic components and their associated abundances should be derived. However, intimate mixtures happen at microscopic scales, thus complicating the analysis with nonlinear mixing effects [133]. In addition to spectral mixing effects, there are many other interfering factors that can significantly affect the analysis of remotely sensed hyperspectral data. For instance, atmospheric interferers are a potential source of errors in spectral unmixing. Multiple scattering effects can also lead to model inaccuracies.

In linear spectral unmixing, the macroscopically pure components are assumed to be homogeneously distributed in separate patches within the field of view. In nonlinear spectral unmixing, the microscopically pure components are intimately mixed. A challenge is how to derive the nonlinear function, since nonlinear spectral unmixing requires detailed a priori knowledge about the materials. Resulting from this limitation, a vast majority of techniques have focused on linear spectral unmixing, where the goal is to find a set of macroscopically pure spectral components (called endmembers) that can be used to unmix all other pixels in the data. Unmixing then amounts to finding the fractional coverage (abundance) of each endmember in each pixel of the scene, which can be approached as a geometrical problem [134]. In the following section, we focus on the most relevant parts of the linear spectral unmixing chain. We also summarize the main efforts in nonlinear spectral unmixing.

#### A. Estimation of the Number of Endmembers

Determining the number of pure spectral endmembers in HSIs is a challenging problem. One of the most commonly



used approaches to this problem is the virtual dimensionality (VD) method [135], which follows the pigeon-hole principle. If we represent a signal source as a pigeon and a spectral band as a hole, we can use a spectral band to accommodate one source. Thus, if a signal source is present in our remotely sensed hyperspectral data set, we should be able to detect this particular source in the relevant spectral band. This can be done by calculating the eigenvalues of both the data-correlation and covariance matrices. A source is present if their difference is positive. Another popular approach has been hyperspectral signal identification with minimum error (HySime) [136]. The idea of HySime is to find the first  $k$  eigenvectors that contain the most data information, i.e., to find the  $k$  such that the mean square error (MSE) between the original data and its projection onto the eigenvector subspace is minimized. Subspace  $k$  is ranked in terms of data variance, but noise variance is not unitary in different directions and the contribution from signals may be smaller than from noise. HySime addresses this issue by using subspace projection techniques, thus contributing an additional feature with regard to VD: modeling of noise before the estimation. The eigenvalue likelihood maximization (ELM) method [137], in turn, implements a modification of the VD concept based on the following observations: 1) the eigenvalues corresponding to the noise are identical in the covariance and the correlation matrices, and 2) the eigenvalues corresponding to the signal (the endmembers) are larger in the correlation matrix than in the covariance matrix. The ELM takes advantage of this fact and provides a fully automatic method that does not need an input parameter (as does VD) nor estimation of the noise (as does HySime). Finally, the normal compositional model (NCM) [138] addresses the possibility that, in real images, there may not be any pure pixels. To address this issue, NCM assumes that the pixels of the HSI are linear combinations of an unknown number of random endmembers (the opposite of the deterministic approach). This model provides more flexibility with respect to the observed pixels and the endmembers, which are allowed to be a greater distance from the observed pixels.

### B. Endmember Extraction

The identification of endmembers is a challenging problem for which many different strategies have been proposed in the literature [134]. In order to categorize algorithms, we consider three different scenarios: 1) The data contains at least one pure pixel per endmember, i.e., there is at least one spectral vector in each vertex of the data simplex (pure pixel assumption); 2) The data do not contain pure pixels but contain enough spectral vectors on each facet. In this case, we may fit a minimum volume simplex to the data; 3) The data is highly mixed, with no spectral vectors near the facets. In this case, minimum volume algorithms fail and we need to resort to a statistical framework. We also consider algorithms that include spatial information in addition to spectral information for this purpose.

1) *Pure Pixel Assumption:* These methods assume a classic spectral unmixing chain with three stages: dimensional reduction, endmember selection, and abundance estimation. Here,

the endmembers are directly derived from the original hyperspectral scene. The pixel purity index (PPI) [139] is perhaps the most popular endmember extraction algorithm due to its availability in software packages. PPI has many parameters involved and is not an iterative algorithm. Manual intervention is required to select a final set of endmembers, which makes it unattractive for automatization purposes. An alternative is the N-FINDR [140], which assumes the presence of pure pixels in the original hyperspectral scene and further maximizes the volume that can be formed with pixel vectors in the data cube. Orthogonal subspace projection (OSP) [141], in turn, uses the concept of orthogonal projections. Vertex component analysis (VCA) [142] iteratively projects data onto a direction orthogonal to the subspace spanned by the endmembers, which have been already determined. In this regard, the algorithm is similar to OSP with: the main difference is that VCA applies a noise characterization process in order to reduce the sensitivity to noise. This is done by using singular value decomposition (SVD) to obtain the projection that best represents the data in the maximum-power sense. Another important concept in this category is the endmember bundle, explored by algorithms such as the multiple endmember spectral mixture analysis (MESMA) [143]. Although the shape of an endmember is fairly consistent, its amplitude generally varies due to illumination conditions, spectral variability, topographic modulation and others. MESMA addresses this issue using endmember bundles, which incorporate variability by representing each endmember by a set or bundle of spectra, each of which could reasonably be the reflectance of an instance of the endmember.

2) *Minimum Volume Algorithms:* If the data do not contain any pure signatures, we can fit a simplex of minimum volume in the case we have enough spectral vectors on the facets. This idea is the opposite of the concept of maximum volume adopted by N-FINDR: here the goal is to find the simplex with a minimum volume that encloses the data. From an optimization point of view, these algorithms are formulated by including a data term that minimizes the reconstruction error and a volume term that promotes mixing matrices of minimum volume. This is the case for the iterative constrained endmembers (ICE) [144] and the minimum volume constrained nonnegative matrix factorization (MVC-NMF) [145], whose main differences are related to the way they define the data volume term. The sparsity-promoting ICE (SPICE) [146] is an extension of the ICE algorithm that incorporates sparsity-promoting priors aimed at finding the number of endmembers. The minimum volume estimation (MVES) algorithm [147] integrates concepts of convex analysis and volume minimization to provide a solution similar to that of previously mentioned algorithms but using cyclic minimization with linear programming. Again, the assumption is that the enclosing simplex with minimum volume should coincide with the true endmember simplex (MVES uses hard positivity constraints). The minimum volume simplex analysis (MVSA) algorithm [148] follows a similar strategy, but allows violations of the positivity constraint. This is because, due to the presence of noise or perturbations, spectral vectors may lie outside the true simplex and this may introduce errors in the characterization. The MINVEST algorithm [149] also exploits this concept,

allowing a certain number of outliers when estimating the minimum volume that encompasses the HSI.

3) *Highly Mixed Data*: When the spectral mixtures are highly mixed, the geometrical-based methods yield poor results because there are not enough spectral vectors in the simplex facets. Statistical methods are a powerful alternative that usually comes with a price: higher computational complexity than with geometrical methods. Since, in most cases, the number of substances and their reflectances are not known, the problem can be approached as a blind source separation problem, with some statistical unmixing approaches proposing variations on the independent component analysis (ICA) [150]. However, ICA applicability is compromised by the statistical dependence existing among abundances. This has been addressed (among other strategies presented in the recent literature) by the dependent component analysis (DECA) algorithm [151]. Bayesian approaches have also been used. They have the ability to model statistical variability and impose priors to constrain solutions to physically meaningful ranges.

4) *Inclusion of Spatial Information*: Most available algorithms for endmember identification do not consider information about spatial-contextual information. In certain scenarios, it is important to include the spatial information in the analysis. The automatic morphological endmember extraction (AMEE) [152] uses extended morphological transformations to integrate spatial and spectral information. The spatial-spectral endmember extraction (SSEE) [153] uses a different approach. First, it processes the image using a local search window and applies singular value decomposition (SVD) to determine a set of eigenvectors that describe most of the spectral variance in the window. Then, it performs a projection of the entire image data onto eigenvectors to determine candidate endmember pixels. Finally, it uses spatial constraints to combine and average spectrally similar candidate endmember pixels (preserving similar but distinct endmembers that occupy unique image regions). In order to avoid modifying spectral-based algorithms for endmember extraction, spatial information can also be included as a preprocessing module, such as the spatial preprocessing (SPP) algorithm [154]. A region-based approach (RBSPP) [155] has also been developed to adaptively include spatial information. Finally, a spatial-spectral preprocessing (SSPP) approach [156] has been developed to derive a spatial homogeneity index that uses Gaussian filtering and is thus relatively insensitive to the noise present in the hyperspectral data.

### C. Abundance Estimation

Once the endmember signatures have been derived, different strategies can be used to estimate their fractional abundances different strategies [157]. The idea is to find the abundances that minimize the reconstruction error obtained after approximating the original hyperspectral scene using a linear mixture model. However, this is generally an unconstrained solution that does not satisfy the abundance non-negativity (ANC) and the abundance sum-to-one constraints (ASC). Whether abundance constraints should be imposed or not depends

on the practical application. It has been argued that, if the linear mixture model is accurate, the two constraints should be satisfied automatically. In any event, the ANC is more important than the ASC. Due to noise and spectral variability, reinforcing the ASC may be prone to introduce additional estimation error. When endmembers are unknown, endmember signatures should be extracted or estimated first. Some endmember extraction algorithms can provide abundance estimates simultaneously (e.g., algorithms without the pure pixel assumption) [134]. Another group of abundance estimation approaches based on blind source separation, which does not require endmember signatures to be known a priori, has been developed [134]. Widely used matrix-factorization-based blind source separation methods include independent component analysis (ICA) and non-negative matrix factorization (NMF), which have been mostly used in the context of unsupervised (soft) classification. If the linear assumption does not hold, nonlinear unmixing techniques should be used. In addition, as mentioned in previous section, if the spectral variability or endmember variability is being considered, the mixture model must be modified accordingly, which is traditionally accomplished by generating modified/extended linear mixture models [158–160].

### D. Sparse Unmixing

Spectral unmixing algorithms with the pure pixel assumption require the presence of pure pixels in the scene for endmember extraction. Due to spatial resolution and mixing phenomena, this assumption cannot be always guaranteed. Spectral unmixing algorithms without the pure pixel assumption generate endmember signatures that often do not relate to real physical signatures. A possible solution is to use ground spectral libraries to perform spectral unmixing, but libraries are very large, and hence the problem becomes sparse and difficult to solve. Another problem is the difference between the ground library and the image data. To address these issues, sparse unmixing [161] expresses pixel vectors as linear combinations of a few pure spectral signatures obtained from a potentially very large spectral library of ground materials. An advantage is that it sidesteps the endmember extraction step (including the estimation of the number of endmembers). In order to incorporate spatial information into the spectral unmixing formulation, a total variation (TV) regularizer has been developed to enforce spatial homogeneity by including this term in the original objective function [162]. It produces spatially smooth abundance fractions that improve sparse unmixing performance, even in very high noise conditions. Further, since it is generally observed that spectral libraries are organized in the form of groups with different variations of the same component (e.g., different mineral alterations), exploiting the inherent group structure present in spectral libraries can improve the results of sparse unmixing by selectively enforcing groups. For this purpose, a group-based formulation of sparse unmixing has been introduced [163]. A further development has also been recently introduced based on the concept of collaborativity, which promotes solutions with a minimum number of active endmembers (the number

of endmembers in a scene is generally low). This allows the number of endmembers participating in the final solution to be minimized [164], also partially circumventing the need to estimate the number of endmembers in the scene [165].

### E. Nonlinear Unmixing

Apart from the linear spectral unmixing algorithms, several alternatives can be found in the literature that deal with the mixture problem [166–168]. The bilinear mixing model (BMM) considers secondary illumination sources. This model represents a simplification of reality, as it only considers bilinear interactions (objects that can be illuminated by light reflected first by another object) [133]. BMMs can be generalized to deal with multiple endmembers, where they try to model bilinear interactions as new endmembers. The generalized bilinear model (GBM) provides more flexible solutions that can avoid some overfitting problems associated to bilinear models, but still assumes that there are not self-bilinear interactions [133]. BMM and GBM have inspired several posterior methodologies, some that use polynomial functions to model the nonlinearities provided by layer partitions and scattering properties in multilayered scenarios [169], and others that use different approaches to solve the proposed model [170–173]. In [174], the authors propose a modification of BMM where the transformation of the spectral information is based on a two-degree polynomial, thus defining the polynomial post-nonlinear model (PPNM), whose main advantage is that it is able to deal with self-interactions; PPNM has been recently extended to deal with all polynomial degrees [175]. In [169] a p-linear polynomial model is used to characterize nonlinearities in combination with a supervised artificial neural network (ANN) and polytope decomposition scheme. When the interactions are considered to exist at photon level, nonlinear unmixing methods try to model the optical characteristics of the intimate mixture from a theoretical analysis of the reflectance behavior attending at the specific geomorphical, chemical, and physical properties of the observed data. Several models have been proposed in this context, but the Hapke model [176] is still the most widely adopted approach. In addition, supervised techniques based on the use of methods such as kernel-based and ANNs have been proposed to perform unmixing on a microscopic scale [177–186]. The main advantage of this technique is that it can create a preliminary model of every nonlinear behavior; however, the need for reliable ground truth information about the training samples represents a major shortcoming.

### F. Challenges in Spectral Unmixing

Despite the availability of several consolidated techniques for linear spectral unmixing, and a suite of incipient techniques for nonlinear spectral unmixing, an important challenge remains that is related to the nonlinearity of mixing phenomena in HSIs. The inherently nonlinear nature of the process, and the dependence on the observed objects, creates the need to incorporate detailed information about the observed objects in order to properly model the multiple scattering phenomena occurring in the nonlinear case. The estimation of participating

endmembers in the mixture also remains a challenge, despite the availability of some techniques that can provide reasonable approximations. Another important hurdle for the automatic execution of sparse unmixing algorithms is the heterogeneity in available spectral libraries, although important efforts have been made toward the development of open-source libraries of spectral materials that can alleviate the need to estimate the number and the signatures of spectral endmembers in advance. In this regard, the availability of open libraries such as SPECCHIO<sup>4</sup> provides an important first step toward the general use of open spectral libraries for spectral unmixing purposes.

## VI. RESOLUTION ENHANCEMENT

The resolution enhancement of HSIs has attracted increasing attention in recent years, as shown in the statistical analysis of the trend described in the introduction. Resolution enhancement techniques for HSIs can be broadly categorized into four classes, as shown in Fig. 9: 1) hyperspectral super-resolution (multi-image/single-image), 2) subpixel mapping (or super-resolution mapping), 3) hyperspectral pan-sharpening, and 4) hyperspectral and multispectral data fusion (HS-MS fusion). The first, hyperspectral super-resolution, is an extension of ordinary super-resolution in computer vision to HSIs, where a high-resolution HSI is reconstructed from multiple low-resolution HSIs (or a single HSI) acquired by a single sensor. The second, subpixel mapping, is a resolution enhancement technique processed at a classification level using a single HSI as input. The remaining two classes, hyperspectral pan-sharpening and HS-MS fusion, are multisensor super-resolution techniques, in which an HSI is fused with an auxiliary higher-resolution data source (panchromatic or multispectral images) taken over the same area at the same time (or a similar period) for creating a high-resolution HSI. The following subsections provide an overview of recent advances in these four classes of techniques.

### A. Hyperspectral Super-Resolution

A variety of super-resolution image reconstruction techniques has been intensively developed over the past three decades in computer vision. Super-resolution techniques can be roughly divided into two types [187]: 1) classical multi-image super-resolution that obtains a high-resolution image (or sequence) from multiple low-resolution images for the same scene with different subpixel shifts, and 2) learning-based super-resolution that learns correspondence between low and high-resolution image patches from an external training database.

Multi-image super-resolution techniques have been extended to HSIs. Depending on the type of subpixel shifts in the low-resolution image, these techniques can be divided into two approaches: the first approach uses multiple HSIs acquired by the same sensor over the same scene, similar to the classical multi-image super-resolution; the second takes advantage of band-to-band misregistration (so-called *keystone*)

<sup>4</sup><http://www.specchio.ch/>

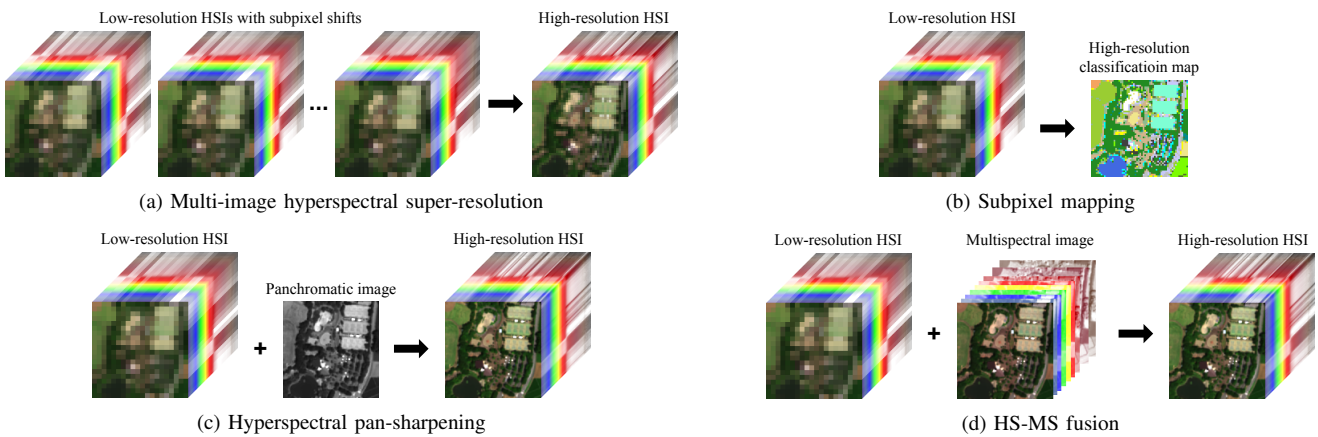


Fig. 9. Four classes of resolution enhancement techniques for HSIs: (a) multi-image hyperspectral super-resolution, (b) subpixel mapping, (c) hyperspectral pan-sharpening, and (d) HS-MS fusion.

in a single HSI. Akgum *et al.* first applied the multi-image super-resolution technique to multiple HSIs fully based on simulations [188].

In the remote sensing community, multi-image hyperspectral super-resolution has been mainly studied using multi-angular HSIs obtained by the compact high-resolution imaging spectrometer (CHRIS). In [189], conventional super-resolution techniques were applied to a set of multi-angular CHRIS images. Chan *et al.* developed a new super-resolution technique for multi-angular HSIs based on a thin-plate spline nonrigid transform model; this approach was intended to improve the image registration procedure and demonstrated its effectiveness with the experiments using multi-angular CHRIS images [190]. The impact of multi-angular super-resolution on classification and unmixing applications was further investigated in [191, 192].

Qian and Chen developed a technique for the second approach that enhances the spatial resolution of HSIs by exploiting the keystone characteristics [193]. Keystone is band-to-band misregistration in the cross-track direction caused by optical aberrations and misalignments in pushbroom systems. Different band images, including different subpixel misregistrations, can be used as input images in the multi-image super-resolution framework. The advantage of this method is that it requires only a single HSI, but the limitation is that the spatial resolution can be enhanced only in the cross-track direction.

Zhao *et al.* proposed sparse-representation-based algorithms for learning-based hyperspectral super-resolution [194, 195]. The high-resolution version of a given low-resolution patch is recovered by solving the sparse linear inverse problem with spectral regularization based on spectral unmixing, in which the patch dictionary is learned from a set of high-resolution panchromatic images or HSIs. Patel *et al.* developed a learning-based super-resolution method for HSIs based on an adaptive wavelet designed from training HSIs [196]. These learning-based super-resolution techniques do not require multiple images over the same scene, but require an external training database with target resolution. For the upcoming hyperspectral satellites, it is realistic to use images obtained

by operational multispectral satellites (e.g., Sentinel-2) for the training database.

### B. Subpixel Mapping

Subpixel mapping is a technique for enhancing the spatial resolution of spectral images by dividing a mixed pixel into subpixels and assigning land-cover classes to these subpixels [197]. Subpixel mapping techniques have been actively studied using multispectral images and extended to HSIs by exploiting rich spectral information. Subpixel mapping comprises two steps: estimating fractional abundances of classes (or endmembers) at a coarse resolution by soft classification or spectral unmixing; and determining the subpixel location of each class within a pixel, assuming spatial dependence. Many recent advances in subpixel mapping of HSIs aim at improving the accuracy of inverse problems involving these two steps.

Tong *et al.* proposed a method that exploits not only attraction but also repulsion between subpixels to better retrieve spatial dependence [198]. Zhang *et al.* integrated learning-based super-resolution into subpixel mapping, requiring an external training set [199]. The estimation of unknown spatial details of classes at a high resolution from a single low-resolution image is a typical ill-posed inverse problem. Different techniques for spatial regularization that have been recently studied in order to mitigate the ill-posedness, have assumed spatial prior models, such as Laplacian, total variation, bilateral total variation, and nonlocal total variation [200, 201].

Obviously, the accuracy of the abundance maps obtained in the first step greatly affects the final performance of subpixel mapping. To address this issue, Tong *et al.* proposed a genetic algorithm that can correct possible errors in the initial estimation of abundances using a mutation operator [202]. Xu *et al.* introduced a method that improves the accuracy of spectral unmixing in subpixel mapping by taking endmember variability into consideration, with the assumption that different representative spectra for each endmember are available [201].



### C. Hyperspectral Pan-Sharpening

Pan-sharpening is a technique that enhances the spatial resolution of multispectral imagery by fusing it with a higher-resolution panchromatic image. Hyperspectral pan-sharpening is an extension of conventional pan-sharpening to HSIs, and also a special case of HS-MS fusion. Naturally, there are two main approaches: 1) extensions of pan-sharpening methods, and 2) subspace-based methods (originally developed for HS-MS fusion). Hyperspectral pan-sharpening is motivated by spaceborne imaging spectroscopy missions that mount both hyperspectral and panchromatic imaging sensors, such as EO-1/Hyperion-ALI [203], PRISMA [8], and Shalom [9]. The main advantage of HSIs on multispectral images is rich spectral information, which enables discrimination and identification of spectrally similar materials. In other words, any spectral distortions will lead to inaccurate analysis results in the subsequent data processing. Therefore, the challenge of hyperspectral pan-sharpening is to enhance the spatial resolution of hyperspectral data, while minimizing spectral distortion.

Typical pan-sharpening techniques include component substitution (CS) [204–206], multiresolution analysis (MRA) [207, 208], and sparse representation (SR) [209, 210] algorithms. In the general CS-based pan-sharpening scheme proposed by Aiazzi *et al.* [206], a multispectral image is sharpened by adding spatial details obtained by multiplying the difference between a panchromatic image and a synthetic intensity component by a band-wise modulation coefficient. Gram-Schmidt Adaptive (GSA) is one of the benchmark CS-based pan-sharpening algorithms, in which the synthetic intensity component is computed via linear regression between a panchromatic image and lower-resolution bands. In the MRA-based pan-sharpening scheme, spatial details of each multispectral band are obtained by use of MRA, which calculates the difference between a panchromatic image and its low-pass version multiplied by a gain factor. The gain factor can be computed either locally or globally. Representative MRA pan-sharpening algorithms include the smoothing filtered-based intensity modulation (SFIM) method [207], the additive wavelet luminance proportional (AWLP) method [211], and the generalized Laplacian pyramid (GLP) method [208]. The SR-based pan-sharpening approach can be regarded as a special case of learning-based super-resolution, learning correspondence between low and high-resolution image patches from a panchromatic image [210] or an external database, including multiple high-resolution multispectral images [212]. In [213], Alparone *et al.* demonstrated the fusion of Hyperion and ALI panchromatic images using CS and MRA-based algorithms [214].

The subspace-based approach exploits the intrinsic spectral characteristics of the scene via a subspace spanned by a set of basis vectors, and enhances the spatial resolution of subspace coefficients. Since most of the subspace-based methods have been developed for HS-MS fusion, they will be discussed in the next subsection.

Eleven state-of-the-art algorithms from CS, MRA, and subspace-based approaches were applied to hyperspectral pan-

sharpening and compared in [215]. MRA and subspace-based algorithms demonstrated relatively high-quality and stable results. However, it was evident that large spectral distortion was inevitable for all algorithms under comparison, implying that there is room for further technology development [215].

### D. Hyperspectral and Multispectral Data Fusion

HS-MS fusion is a technique that fuses an HSI with a higher-resolution multispectral image to create high-resolution hyperspectral data. Unlike hyperspectral pan-sharpening, this technique yields higher spectral quality owing to spectral information from the high-resolution data source. Enormous efforts have made to develop algorithms in the last decade. Quite recently, a comparative review of HS-MS fusion was reported in [216]. Most of the HS-MS fusion algorithms can be categorized into at least one of the following six classes: 1) CS, 2) MRA, 3) SR, 4) unmixing, 5) Bayesian, and 6) hybrid approaches. The CS, MRA, and SR approaches are extensions of pan-sharpening techniques, whereas the unmixing and Bayesian approaches fall into the same broader category of subspace-based techniques.

CS-based methods can be adapted to HS-MS fusion by splitting the fusion problem into several pan-sharpening subproblems and applying CS-based pan-sharpening to these subproblems. A key procedure for the CS-based approach is to divide the HSI into several groups and assign a high-resolution image selected or synthesized from the MSI to each group. Chen *et al.* proposed a framework that solves the HS-MS fusion problem by dividing the spectrum of HS data into several regions and fusing hyperspectral and multispectral band images in each region by conventional pan-sharpening techniques [217]. Hyper-sharpening proposed by Selva *et al.* is a framework that effectively adapts MRA-based pan-sharpening methods to HS-MS fusion by synthesizing a high-resolution image for each HS band as a linear combination of MS band images via linear regression [218]. An SR-based pan-sharpening technique was adapted to HS-MS fusion with spectral grouping and joint sparse recovery in [219].

Subspace-based methods have been actively developed for HS-MS fusion problems. The HS-MS fusion task is formulated as the inverse problem of estimating the subspace basis and coefficients of the high-resolution HSI from the two input images [220]. The resolution-enhanced HSI can be reconstructed as the product of the basis matrix and the high-resolution coefficient matrix.

In recent years, a perspective of spectral unmixing has been attracting considerable attention in the context of subspace-based HS-MS fusion, owing to its straightforward interpretation of the fusion process: the basis matrix is defined as a set of spectral signatures of intrinsic materials (so-called endmembers) and the subspace coefficients correspond to the fractional abundances. Several unmixing-based methods have been proposed for HS-MS fusion with various optimization formulations [221–228]. In [216], it was noted that the high-resolution abundances estimation step has a significant influence on fusion performance: outstanding and stable performance can be achieved by minimizing the unmixing

reconstruction errors with respect to both HS-MS images rather than only the multispectral image, particularly when the overlap of SRFs between HS-MS sensors is small. Such algorithms include coupled nonnegative matrix factorization (CNMF) [225], coupled spectral unmixing with a projected gradient method [229], and HySure [230].

One important aspect of developing subspace-based methods has been determining how to mitigate the ill-posedness of inverse problems involving the estimation of high-resolution subspace coefficients. Regularization on subspace coefficients has been extensively explored. Sparsity-promoting regularization has been commonly adopted for subspace-based methods [223, 226, 227], assuming that there is a limited number of endmembers at each high-resolution pixel. HySure solves a convex optimization problem with vector-total-variation-based regularization of the spatial distribution of subspace coefficients, leading to implicit denoising effects in the fusion results [230]. Wei *et al.* developed a Bayesian HS-MS fusion methodology in which different types of regularization terms on subspace coefficients can be designed flexibly, based on information from the prior distribution in the observed scene [231]. A Sylvester equation-based explicit solution was further integrated into the Bayesian methodology to speed it up, leading to the fast fusion based on Sylvester equation (FUSE) [232]. Veganzones *et al.* [233] introduced local image processing into the unmixing-based approach, where well-posed inverse problems are solved for each small patch, assuming that the number of endmembers in each patch is smaller than the number of multispectral bands.

One interesting finding of the comparative study in [216] was that hyper-sharpening methods and unmixing-based methods (CNMF and HySure), which are entirely different approaches, showed high and comparable numerical performances, although the characteristics of resolution-enhanced HSIs are different. This finding implies that hybrid methods combining different approaches can be expected to further improve fusion performance.

Table V shows the overall quantitative assessment results of twelve HS-MS fusion algorithms for eight simulated HS-MS data sets used in [216], including those based on the Pavia University and University of Houston data sets. The quantitative assessment of fusion performance was carried out based on a version of Wald’s protocol presented in [216, 233].

The reconstruction performances of algorithms have been the greatest concern for researchers; however, their general versatility, computational costs, and impacts on applications are also important for users. Since each method has advantages and disadvantages, it is essential to choose a method based on the fusion and analysis scenarios. The impact of HS-MS fusion on applications has been recently investigated via classification and unmixing [216, 234]. Further research on real data is still necessary to verify the practicability of HS-MS fusion for the hyperspectral satellites of the near future.

### E. Challenges in Resolution Enhancement

The main challenges in the resolution enhancement of HSIs relate to practical issues. The development of multisensor

TABLE V  
AVERAGE QUALITY MEASURES FOR EIGHT DATA SETS FROM [216]

Category	Method	PSNR	SAM	ERGAS	Q2 <sup>n</sup>
CS	GSA [206]	39.221	2.063	1.885	0.8796
MRA	SFIM-HS [207, 218]	41.074	1.707	1.789	0.8867
	GLP-HS [208, 218]	41.322	1.664	1.733	0.8984
Unmixing	CNMF [225]	42.092	1.607	<b>1.681</b>	0.9060
	ECCV’14 [226]	38.101	2.658	4.561	0.8431
	ICCV’15 [228]	40.470	1.672	1.902	0.8847
	HySure [230]	<b>42.336</b>	<b>1.602</b>	1.766	<b>0.9089</b>
Bayesian	MAP-SMM [220]	40.008	1.822	1.997	0.8635
	FUSE [232]	40.568	1.881	1.908	0.8693
	FUSE-S [232]	41.177	1.803	1.804	0.8764
Ideal		$\infty$	0	0	1

super-resolution algorithms is recently active; however, very few publications in the literature discuss experiments on real data. Studies on the impact of the resolution enhancement of HSIs on applications are still lacking. It is necessary to clarify benefits and address practical issues of the resolution enhancement technology to promote its operational application to the hyperspectral satellite missions of the near future. For instance, temporal mismatches included in the input images raise challenges for resolution enhancement of HSIs. Furthermore, no-reference image quality assessment for HSIs needs to be developed to provide reliability information of resolution-enhanced hyperspectral data products for end-users. In the multisensor superresolution, there remains a problem that large spectral distortions are inevitable when the mismatch between the two imaging sensors is large in either the spatial or spectral domain (e.g., a large GSD ratio or hyperspectral pan-sharpening). To address this issue, a possible future direction for performance improvement lies in developing algorithms that exploit a spectral library or spatial information of a high-resolution image.

## VII. HSI DENOISING AND IMAGE RESTORATION

Image restoration generally refers to the reconstruction of the true image based on its corrupted version. The true image is unknown and therefore is estimated through the observed image, which has been degraded by different sources. Degradation sources depend on the imaging technology, system, environment, and other factors. When the observed signal is degraded by noise sources the estimation task is usually called denoising. Image restoration is generally used to refer to a broader set of methods that also include applications such as deconvolution, deblurring, and inpainting. However, in HSI analysis, the term restoration is often used to refer to denoising.

In hyperspectral imaging the received radiance at the sensor is degraded by different sources such as atmospheric haze and instrumental noise. The atmospheric effects are usually compensated by applying atmospheric corrections. Instrumental (sensor) noise includes thermal (Johnson) noise, quantization noise, and shot (photon) noise. Spectral bands are often corrupted to some degree. The presence of corrupted bands (also called junk bands) could degrade the efficiency of the image analysis technique and therefore they are usually removed from the data before any further processing. The information lost by removing those bands can be substantial; hence, an

alternative approach is to recover those bands and improve the SNR of the HSI. As a result, HSI restoration can be considered to be an important preprocessing step in HSI analysis.

An HSI can be modeled by

$$\mathbf{X} = \mathbf{H} + \mathbf{N}, \quad (1)$$

where  $\mathbf{X}$  is an  $n_{12} \times d$  matrix ( $n_{12} = n_1 \times n_2$ ) containing the vectorized observed image at band  $i$  in its  $i$ th columns,  $\mathbf{H}$  is the true unknown signal to be estimated and is represented as an  $n_{12} \times d$  matrix containing the unknown vectorized image at band  $i$  in its  $i$ th columns, and  $\mathbf{N}$  is an  $n_{12} \times d$  matrix containing the vectorized noise at band  $i$  in its  $i$ th columns. Note that all the aforementioned noises can be assumed to be additive noise. The restoration task is to estimate original (unknown) signal  $\mathbf{H}$ . Penalized (regularized) least squares is one of the most popular and common minimization framework used for estimation in HSI restoration. Penalized least squares usually is composed of a fidelity term and a penalty term. The penalty term is often chosen based on the prior knowledge of the signal and might be a combination of penalties. Penalized least squares might be also solved subject to some constraints.

#### A. Noise Source Assumptions in HSI

The presence of different noise sources in HSI makes its modeling and restoration very challenging. Therefore, HSI restoration often employs one of the following approaches or a mixture of them.

1) *Signal Independent Noise*: Thermal noise and quantization noise in HSI are modeled by signal independent Gaussian additive noise [235, 236]. Usually, noise is assumed to be uncorrelated spectrally, i.e., the noise covariance matrix is diagonal [236, 237]. The Gaussian assumption has been broadly used in hyperspectral analysis, since it simplifies the analysis considerably. In addition, noise parameter (variance) estimation is simpler under this assumption.

2) *Signal Dependent Noise*: Shot (photon) noise in HSI is modeled by a Poission distribution, since the noise variance is dependent on the level of signal. The noise parameter (variance) estimation is more challenging under this assumption compared to the signal independent case since it varies with respect to the signal level.

3) *Sparse Noise*: Impulsive noise, missing pixels, missing lines, and other outliers often exist in the acquired HSI, usually due to the malfunctioning of the sensor. In this review, we categorize them as sparse noise due to their sparse characteristics. Sparsity techniques or sparse and low-rank decomposition techniques have been used to remove sparse noise from the signal. In [238], impulsive noise was removed using an  $\ell_1$ -norm for both penalty and data fidelity terms in the minimization problem suggested.

4) *Striping Noise*: Hyperspectral imaging systems might also induce artifacts in HSIs usually referred to as pattern noise. For instance, in push-broom imaging systems, the target is scanned line by line and the image line is acquired in different wavelengths by an area-array detector (usually a charged coupled device or CCD). This line-by-line scanning causes an artifact called striping noise, which is often due to

calibration error and the sensitivity variation of the detector [239]. Striping noise reduction (also referred to as destriping in the literature) for push-broom scanning techniques has been widely studied in the remote sensing literature [240, 241], including work on HSI remote sensing [239, 242–244].

#### B. Evolution of HSI Restoration Approaches

Hyperspectral image restoration has been developed considerably during the past few years. Conventional restoration methods based on 2-D modeling and convex optimization techniques were not effective enough for HSI due to the lack of understanding of spectral information. The highly correlated spectral bands in HSI has been found very useful for HSI restoration. This is the main reason for the success of using multiple linear regression (MLR) as an estimation technique [245]. In [245], it is assumed that each band is a linear combination of the other bands, and therefore the  $i$ th band is estimated using least squares estimation. Note that this technique has been used for noise parameter estimation in several HSI restoration approaches [246, 247]. Many restoration approaches have been suggested in the literature to exploit the spectral information, and can be categorized into three main groups.

1) *Approaches that Use 3-D Models instead of 2-D Ones*: In [248], the discrete Fourier transform (DFT) was used to decorrelate the signal in the spectral domain and the 2-D discrete wavelet transform (2-D DWT) was investigated to denoise the signal in the spatial domain. In [249], HSI was treated as a 3-D datacube and an HSI restoration technique was proposed based on sparse analysis regularization and undecimated wavelet transform (UWT). The advantages of 3-D wavelets (orthogonal and undecimated) over 2-D ones for HSI restoration are also discussed in [249] and [237].

2) *Approaches that Propose New Penalties for Penalized Least Squares that Also Take into Account Spectral Information*: An algorithm given in [250] uses the 2-D DWT and a sparse restoration criterion based on penalized least squares having a group of  $\ell_2$  penalties on the wavelet coefficients. This method was improved in [251] for hyperspectral image restoration by taking into account the spectral noise variance in the minimization problem and solving it by using the alternating direction method of multipliers (ADMM). Subsequently, due to the redundancy and high correlation in the spectral bands in HSIs, penalized least squares using a first-order spectral roughness penalty (FOSRP) was proposed for hyperspectral image restoration [252]. The new cost function was formulated in the wavelet domain to exploit the MRA property of wavelets. The Stein unbiased risk estimator (SURE) was utilized to automatically select the tuning parameters. It was shown that FORSP outperforms sparsity penalties for HSI restoration. This method was improved using a combination of a spectral roughness penalty and a group lasso penalty [253]. Cubic total variation (CTV), given in [254], exploits the gradient in the spectral axis to improve the restoration results compared to TV denoising band by band. In [255], an adaptive version of CTV was applied for preserving both texture and edges simultaneously. In [256], a spatial-spectral

hyperspectral image denoising approach was developed. The spectral derivation was proposed to concentrate the noise in the low frequency. Then, noise was removed by applying the 2-D DWT on the spatial domain and the 1-D DWT on the spectral domain. A spatial-spectral prior for maximum a posteriori (MAP) was proposed in [257]. The prior was based on five derivatives, one along the spectral direction and the rest applied on the spatial domain for four neighborhood pixels.

3) *Approaches that Use Low-rank Models:* Due to the redundancy along the spectral direction, low-rank modeling has been widely used in HSI analyses and applications such as dimensionality reduction, feature extraction, unmixing, and compression. Principal component analysis (PCA) was used for hyperspectral restoration in [258] to decorrelate the noise and signal spectrally. A low-rank representation method is a three-mode factor analysis called Tucker3 decomposition [259], used for HSI restoration [260]. HSI is assumed to be a third order tensor and the "best" lower rank of the decomposition is chosen by minimizing a Frobenius norm. A similar idea was exploited for hyperspectral image restoration by applying more reduction spectrally [261]. A genetic algorithm (GA) was developed to choose the rank of the Tucker3 decomposition [262]. This work was followed by [263], in which the kernel function (Gaussian radial basis) is applied for each spectral band, with the idea of using multilinear algebra efficiently. Multidimensional Wiener filtering that exploits Tucker3 decomposition was used in [264], where the flattening of HSI was achieved by estimating the main direction that gives the smallest rank. Parallel factor analysis is also a low-rank modeling used in [265] for HSI denoising. A new 3-D linear model was proposed for HSI in [266], where 2-D wavelets were used for spatial projection and spectral-singular vectors of the observed HSI for spectral projection. A convex optimization was used for the restoration task based on the 3-D linear model and  $\ell_1$  penalty. Additionally, SURE was used for regularization parameter selection.

Low-rank modeling has also been used in synthesis and analysis penalized least squares [251, 267] and also TV regularization [251, 268]. A 3-D low-rank model in the form of model (2) was proposed in [247], where 2-D wavelets were used as the spatial basis while the spectral basis was assumed to be an unknown low-rank orthogonal matrix. Therefore, an orthogonality constraint was added to the optimization problem for the simultaneous estimation of the two unknown matrices in the minimization problem, which led to a non-convex optimization.

### C. HSI Model Selection

Further studies in HSI modeling and restoration [251] confirmed that capturing spectral redundancy by low-rank modeling is more appropriate than full-rank modeling for HSI modeling and restoration. In [251], a model selection criterion was given for a general model of the form

$$\mathbf{X} = \mathbf{A}\mathbf{W}_r\mathbf{M}_r^T + \mathbf{N}, \quad (2)$$

where  $\mathbf{A}$  ( $n_{12} \times n_{12}$  matrix) and  $\mathbf{M}_r$  ( $d \times r$  matrix,  $r \leq \min(n_{12}, d)$ ) could be two-dimensional and one-dimensional

orthogonal (known) bases, respectively, and  $\mathbf{W}_r$  ( $n_{12} \times r$  matrix) contains the corresponding coefficients for the unknown hyperspectral data,  $\mathbf{H}$ . The term  $\mathbf{W}_r$  is estimated by using  $\ell_1$  penalized least squares and the signal is restored by  $\hat{\mathbf{H}} = \mathbf{A}\hat{\mathbf{W}}_r\mathbf{M}_r^T$ . Ideally, the model that gives the lowest mean squared error (MSE) is the best choice among the candidates. However, the MSE is uncomputable in practice, since it depends on the true (uncorrupted) data. Therefore, SURE is suggested for use as an estimator of MSE for HSI. The results have confirmed that low-rank models give lower MSE (estimated by SURE) compared with full-rank ones.

### D. HSI Restoration with Mixed Noise Assumption

A mixed noise assumption has also been taken into consideration in HSI modeling and restoration. In a mixed noise assumption, HSI ( $\mathbf{X}$ ) in model (1) is assumed to be corrupted by a mixture of the noise sources described in Subsection VII-A. A mixture of the signal dependent noise ( $\mathbf{N}_{SD}$ ) and signal independent ( $\mathbf{N}_{SI}$ ) noise model has been taken into account in [246, 269, 270], as  $\mathbf{N} = \mathbf{N}_{SI} + \mathbf{N}_{SD}$ . Therefore, two parameters need to be estimated: the variances of  $\mathbf{N}_{SI}$  and  $\mathbf{N}_{SD}$ , which have Gaussian distribution and Poisson distribution, respectively. In [269], a 3-D (block-wise) non-local sparse restoration method is suggested for HSI. The minimization problem uses a group lasso penalty and a dictionary consisting of a 3-D discrete cosine transform (3D-DCT) and a 3-D discrete wavelet transform (3D-DWT), which is solved by using the accelerated proximal gradient method. In [270],  $\mathbf{N}_{SI}$  and  $\mathbf{N}_{SD}$  are removed sequentially. Maximum likelihood (ML) was used to estimate the two parameters of the noise model where MLR was investigated for an earlier estimation of the noise. In [242], a subspace-based approach was given to restore HSI corrupted by striping noise and  $\mathbf{N}_{ID}$ .

A widely used mixed noise assumption for HSI restoration is  $\mathbf{N} = \mathbf{N}_{SI} + \mathbf{N}_S$ , where  $\mathbf{N}_S$  represents the sparse noise defined in Subsection VII-A. This mixture assumption was utilized in [271], where low-rank and sparse matrix decomposition was taken into account to restore HSI. This method was improved by augmenting the total variation penalty to the restoration criterion [272]. Also, the noise-adjusted iterative low-rank matrix approximation given in [273] approximates HSI with a low-rank matrix while taking into account the changes of the noise variance through the spectral bands. In [274], using a weighted Schatten p-norm as a nonconvex low-rank regularizer was proposed for low-rank and sparse decomposition of HSI degraded by sparse and Gaussian noises.

### E. Experimental Results

In this subsection, we present some experimental results for HSI restoration. Fig. 10 shows the evolution of HSI restoration techniques based on SNR. The experiments were applied on a portion ( $128 \times 128 \times 96$ ) of the Pavia University data set where the variance of the Gaussian noise added varies along the



spectral axis ( $\sigma_i^2$ ) like a Gaussian shape centered at the middle band ( $d/2$ ), as

$$\sigma_i^2 = \sigma^2 \frac{e^{-\frac{(i-d/2)^2}{2\eta^2}}}{\sum_{j=1}^d e^{-\frac{(j-d/2)^2}{2\eta^2}}},$$

where the power of the noise is controlled by  $\sigma$ , and  $\eta$  behaves like the standard deviation for the Gaussian bell curve [245]. To evaluate the restoration results for the simulated data set, SNR in dB is used as

$$\text{SNR}_{\text{out}} = 10 \log_{10} \left( \frac{\|\mathbf{H}\|_F^2}{\|\mathbf{H} - \hat{\mathbf{H}}\|_F^2} \right),$$

where  $\|\cdot\|_F$  is the Frobenius norm and the noise input level for the whole cube is defined as

$$\text{SNR}_{\text{in}} = 10 \log_{10} \left( \frac{\|\mathbf{H}\|_F^2}{\|\mathbf{H} - \mathbf{X}\|_F^2} \right).$$

In Fig. 10, the results are shown when  $\text{SNR}_{\text{in}}$  varies from 5 to 40 dB in increments of 5dB. Note that the results shown are means over 10 experiments (adding random Gaussian noise) and the error bars show the standard deviations. In this experiment, six restoration methods are compared based on SNR:  $\ell_1$  penalized least squares using 2-D wavelet modeling (2-D Wavelet) [275] and 3-D wavelet modeling (3-D Wavelet) [249]; penalized least squares using first order spectral roughness penalty (FOSRPDN) [252];  $\ell_1$  penalized least squares using a wavelet-based low-rank model (SPAWMARS) [251]; low-rank matrix recovery (LRMR) [271]; and NAILRMA [273] are compared based on SNR. Note that 2-D Wavelet, 3-D Wavelet, FOSRPDN, and SPAWMARS all exploit SURE as a parameter selection technique and therefore are parameter free techniques. Also, note that SPAWMARS is a fully automatic version of SVDSRR [237] with rank selection. The Matlab codes for FORPDN and SPAWMARS are available online as [276] and [277], respectively.

The blue line in Fig. 10 indicates the noise levels, and therefore the effectivity of the HSI restoration techniques can be compared with respect to the noise levels. It can be seen that 3-D wavelet restoration considerably improves the conventional 2-D wavelet techniques. Also, FOSRPDN, which is based on a 2-D wavelet and a first order spectral roughness penalty, outperforms 3-D wavelet restoration, which confirms the importance of the spectral correlations. Finally, SPAWMARS, which utilizes a wavelet-based low-rank, outperforms the other techniques used in this experiment. Note that Wavelab Fast (a fast wavelet toolbox), which is provided in [278], was considered for the implementation of wavelet transforms. A Daubechies wavelet with 2 and 10 coefficients for spectral and spatial bases, respectively, in five decomposition levels is used in all the experiments. CPU processing time in seconds for different restoration approaches used in the experiment confirms that SPAWMARS is the most efficient method. All the experiments in this section are done in Matlab on a computer having an Intel®Core™ i7-4710HQ CPU at 2.5 GHz, 12GB of memory, and 64-bit operating system.

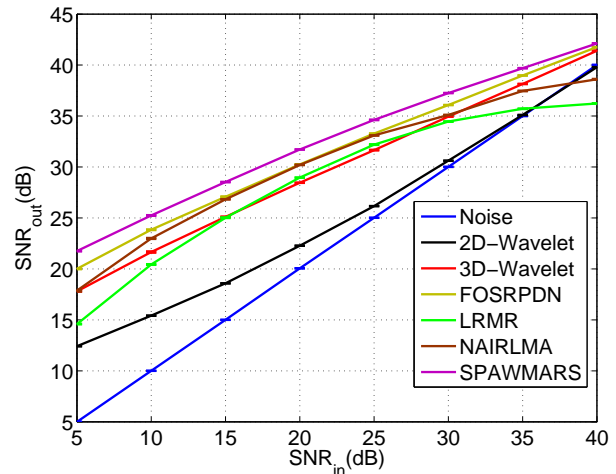


Fig. 10. Comparison of the performance of different HSI restoration methods for different levels of input Gaussian noise in dB.

TABLE VI  
CPU PROCESSING TIME IN SECONDS FOR DIFFERENT RESTORATION APPROACHES USED IN THE EXPERIMENT.

2-D Wavelet	3-D Wavelet	FOSRPDN	LRMR	NAILRMA	SPAWMARS
1.16	1.72	1.09	25.70	4.79	<b>0.62</b>

#### F. Challenges in HSI Denoising and Image Restoration

HSI restoration and modeling have several future challenges.

- 1) HSI model selection and noise parameter estimation need more attention. For instance, a model selection criterion that is not restricted to the Gaussian noise model would be very useful for HSI modeling and restoration. The main advantage of a model selection criterion is that it provides an instrument to compare the restoration techniques without using simulated (noisy) HSI but rather uses the observed HSI itself.
- 2) Investigating the contribution of the various HSI restoration approaches as a preprocessing step for further HSI analysis, such as change detection, unmixing, resolution enhancement, is of interest.
- 3) In mixed noise scenarios, investigating the dominant noise type within HSI should be considered.
- 4) HSI restoration approaches need to be computationally efficient if they are to be used as a preprocessing step in real-world applications. Fast computing techniques may be considered for the fast implementation of HSI restoration approaches in the future.

#### VIII. CHANGE DETECTION

In this section, we define a change detection problem in a pair of bitemporal HSIs (those in the multitemporal domain can be addressed straightforwardly, pair by pair). Let  $\mathbf{X}_1$  and  $\mathbf{X}_2$  be two HSIs acquired over the same geographical area at times  $t_1$  and  $t_2$ , respectively. The hyperdimensional difference image  $\mathbf{X}_D$ , i.e., the spectral change vectors (SCVs), can be computed by subtracting the bitemporal images pixel by pixel, i.e.,

$$\mathbf{X}_D = \mathbf{X}_2 - \mathbf{X}_1. \quad (3)$$

Let  $\Omega = \{\omega_n, \Omega_c\}$  be the set of all classes in  $\mathbf{X}_D$ , where  $\omega_n$  is the no-change class and  $\Omega_c = \{\omega_{C_1}, \omega_{C_2}, \dots, \omega_{C_k}\}$  is the set of  $K$  possible change classes. The considered binary CD problem can be formalized to separate the  $\omega_n$  and  $\Omega_c$  classes without distinguishing different classes in  $\Omega_c$ ; the objective of the multi-class CD task is to detect the changed pixels  $\Omega_c$  and to identify their classes in  $\{\omega_{C_1}, \omega_{C_2}, \dots, \omega_{C_k}\}$ .

Continuous satellite observation resulted in the acquisition of a large number of multitemporal remote sensing images. By analyzing these images, a better understanding of the changes and evolutions on the Earth's surface can be gained. Change detection (CD) is a technique that enables the land-cover changes occurring over a geographical area at different observation times to be identified [279]. In past decades, CD has played important roles in various multitemporal remote sensing applications, such as urban sprawl analysis, disaster loss evaluation, and forest and environmental monitoring. [280–282]. For optical remote sensing images, CD based on multispectral images has been intensively investigated due to the availability of multispectral sensors on board the last generation Earth Observation (EO) satellites. With the launch of new generation EO satellites carrying hyperspectral sensors, there are further opportunities to implement CD in multitemporal HSIs.

Unlike multispectral images, detailed spectral sampling in HSIs allows the potential detection of more spectral variations, especially those with spectrally insignificant subtle changes [283]. Compared with the abrupt changes that present in the coarse multispectral images, changes in HSIs are more sophisticated, implicit, and structurally complex. One emerging problem within this context is to detect small and subtle changes compared to the large unchanging background in multitemporal HSIs. This is related to the definition of an anomaly CD problem, which is not easily addressed by the classical CD techniques. The traditional CD identifies large land-cover changes characterized by significant spectral variations. An empirical definition of the “change” concept in multitemporal HSIs from the global and local spectrum discriminability can be found in [283]. For CD techniques in multispectral images, exhaustive investigations have been made in the past few decades. However, relatively few works covering CD in HSIs can be found. Recently, a book chapter [284] analyzed this challenging task and provided a literature review and comparison between the CD in multispectral and hyperspectral images. Problems and challenges in the existing methods were also discussed. This section reviews several recent CD techniques for HSIs, discussing and analyzing their properties. In addition, a quantitative analysis that compares the CD results obtained by some state-of-the-art CD techniques on the Umatilla County Hyperion data set is provided. A summary of some classical and recent techniques for multitemporal HSI CD techniques is provided in Table VII with some details, such as their categories and principal characteristics. Note that this summary is not exhaustive.

### A. Anomaly Change Detection

Anomaly change detection (ACD) in HSIs is intended to distinguish anomalous changes from the non-changes and pervasive changes in multitemporal HSIs [285]. The main idea of ACD is to design a robust detector that is able to maximize the difference between the anomaly changes and the unchanged scene background but minimize the possible class difference within the background. In this context, multivariate statistical techniques based on linear predictors, e.g., cross covariance (chronochrome, CC) [286] and covariance-equalization (CE) [287], are typical ACD and binary CD detectors. Recently, algorithms were proposed to detect anomalous changes in HSIs by modeling the data with elliptically contoured (EC) distributions [288]. A new ACD method was proposed in [285] that constructed the change residual image based on slow feature analysis (SFA) and detected anomaly changes by using the RX algorithm. A cluster kernel RX (CKRX) algorithm was developed in [289] that clustered the background pixels, then used the cluster centers in the anomaly detection. Other investigations have also focused on various specific changes, for example, work focusing on eliminating image parallax errors [290], vegetation and illumination variation [291], and diurnal and seasonal variations [292].

### B. Binary Change Detection

For binary CD, the aim is to separate the changed and unchanged pixels in the images. In this case, the classical techniques in multispectral images can still be considered, such as thresholding [293] or clustering [294] based on the computed magnitude of SCVs. Transform-based methods constitute an important class. Such techniques represent the original data in a feature space where the significant change information is concentrated in a few transformed components. This not only reduces data dimensionality and noise but also focuses on the specific changes of interest in specific components. The multivariate alteration detection (MAD) method originally proposed for multispectral images in [295] was applied to HSIs to detect vegetation change based on the canonical correlation analysis (CCA) [296]. It was extended to an iterative reweighted version (IR-MAD) [297] to better emphasize and detect changes. A temporal-principal component analysis (T-PCA) was proposed in [298], which exploits the temporal variances in the combined multitemporal HSIs. The no-change information is represented in the first principal component and the change information in the second, which is orthogonal to the first one. Recently, in [299], independent component analysis (ICA) was applied, combined with the uniform feature design strategy, by following a hierarchical framework to identify the specific vegetation changes. In [300], a subspace-based CD approach was designed by using the undesired land-cover class spectral signature as prior knowledge, where the subspace distance was computed to determine the anomalous pixels as change with respect to the background subspace. A semisupervised CD method was proposed in [301], where a Laplacian regularized metric framework was exploited to learn a distance metric for detecting change under noisy conditions. In [16], an unsupervised CD method was developed for slight

TABLE VII

SUMMARY OF SOME CLASSICAL AND RECENT TECHNIQUES FOR MULTITEMPORAL HYPERSPECTRAL IMAGES CHANGE DETECTION (NOT EXHAUSTIVE).

Application Purpose	Supervised/Semisupervised/Unsupervised	Techniques	Main Characteristics
Anomaly CD	Unsupervised	Chronochrome (CC) [285], covariance-equalization (CE) [286]	Anomaly change estimation based on linear predictors
	Unsupervised	Elliptically contoured (EC) distributions-based methods [287]	Anomalous change detectors based on elliptically contoured distributions
	Unsupervised	Slow feature analysis (SFA) - based RX algorithm [284]	SFA is used to construct the change residual image
	Unsupervised	Cluster kernel RX (CKRX) [288]	Background pixels are clustered and the cluster centers are used in ACD
	Unsupervised	Eliminating external factors (e.g., image parallax errors [289], vegetation and illumination variations [290], diurnal and seasonal variations [291])	Enhancing the CD performance
Binary CD	Unsupervised	Multivariate alteration detection (MAD) [294], iterative reweighted MAD (IR-MAD) [296], temporal-principal component analysis (T-PCA) [297], independent component analysis (ICA) [298]	Changes are represented and highlighted in transformed and reduced feature space
	Supervised	Subspace-based approach [299]	Changes are identified as anomalous pixels according to the subspace distance
	Semisupervised	Distance metric-based method [300]	Distance metric is learned in a Laplacian regularized metric framework
	Unsupervised	Slight change detection method [17]	Using block processing and locally linear embedding
	Unsupervised/Supervised	Unmixing-based methods [301-304]	Unmixing single-time or stacked multitemporal images to investigate the subpixel level change
Multiple CD	Supervised	Post-classification comparison (PCC) [305]	Classifying independently each single time image; detailed land-cover transitions (i.e., "from-to" information) can be obtained
	Supervised	3-D spectral modeling approach [306]	A spatial/spectral/temporal joint modeling and unmixing of multitemporal data
	Unsupervised	Hierarchical spectral change vector analysis (HSCVA) [282]	Hierarchical clustering and detecting spectral variations at different significant levels
	Semisupervised	Sequential spectral change vector analysis (S <sup>2</sup> CVA) [307]	Adaptive and sequential projection of SCVs to discover and detect multiple changes
	Unsupervised	Multitemporal spectral unmixing (MSU) [12]	Spectral-temporal unmixing to identify the unique multitemporal endmembers for representing multi-class changes
	Unsupervised/Supervised	Band selection (BS)-based method [309]	The most informative band subset on the difference image is selected to enhance CD performance

change extraction in a multitemporal hyperspectral image sequence. Feature space was built using block processing and locally linear embedding, and then a final CD map was generated by clustering the two change and no-change two classes. In [302, 303], CD in HSIs using unmixing was investigated on a single-time image and stacked multitemporal images, respectively. Also, in [304] and [305], sparse unmixing and a decision fusion based spectral mixture approach were exploited, respectively, in order to detect the subpixel-level change information.

### C. Multiple Change Detection

The aim of multiple CD is more complex and challenging than binary CD. Changed pixels are detected while distinguishing different classes of change. If the multitemporal ground reference samples are available, this task can be addressed according to a typical supervised post-classification comparison (PCC) method [306] by classifying independently two (or more) images at different times and then comparing the pixel class label to detect changes. The main advantages of PCC is that the detailed land-cover transitions are obtained (i.e., "from-to" information). However, the accuracy of the CD performance depends greatly on the accuracy of a single-time image classification result. In [307], a new approach for modeling the temporal variations of the reflectance response as a function of time period and wavelength was developed. A

library of known endmembers that depends mainly on the 3-D surface reconstruction quality and similarity measure is used to perform a classification task. Changes are detected, and it provides better modeling of seasonal variations. However, in practical applications, comprehensive multitemporal ground reference samples are usually not available. Accordingly, unsupervised techniques that do not rely on reference samples are more valuable. In [283], an unsupervised coarse-to-fine hierarchical spectral change vector analysis (HSCVA) approach was proposed. It was designed to cluster and detect changes that include spectral variations at different significant levels according to their discriminable spectral behaviors. A semisupervised sequential spectral change vector analysis (S<sup>2</sup>CVA) technique for discovering, identifying, and discriminating multiple changes in HSIs was proposed in [308]. Different kinds of changes are detected according to adaptive and sequential projection of SCVs at each level of the hierarchy. S<sup>2</sup>CVA successfully extends the compressed change vector analysis (C<sup>2</sup>VA) method [309], which is only suitable for dealing with multispectral CD cases with few spectral channels. In order to identify the sub-pixel level spectral changes so that a more accurate multiple CD result can be produced, an unsupervised multitemporal spectral unmixing (MSU) model was proposed in [12]. The MSU investigates in detail the spectral-temporal mixture properties in multitemporal HSIs, and the considered multiple CD problem is addressed by

TABLE VIII  
CHANGE DETECTION ACCURACIES AND ERRORS OBTAINED BY  
CONSIDERED STATE-OF-THE-ARTS TECHNIQUES ON THE UMATILLA  
COUNTY HYPERION DATA.

Methods	HSCVA[283]	S <sup>2</sup> CVA[308]	MSU[12]
OA(%)	95.26	95.36	95.37
$kappa$	0.8800	0.8822	0.8826
Omission (pixels)	1715	1878	1875
Commission (pixels)	1918	1676	1239
Total errors (pixels)	3633	3554	3114

analyzing the abundances of different distinct multitemporal endmembers at a subpixel level. In addition, both unsupervised and supervised band selection based CD approaches were designed to evaluate the potential detectability of the change detector in a reduced feature space [310]. Experimental results confirmed that the selected most informative band subset can help to improve the CD performance over the use of the original full dimensional HSIs. A multi-class CD experimental comparison was carried out on the Umatilla County Hyperion data set using the state-of-the-art CD techniques introduced above, including HSCVA [283], S<sup>2</sup>CVA [308], and MSU [12]. Numeric experimental results are provided in Table VIII, which includes the obtained overall accuracy (OA), kappa coefficient ( $kappa$ ), and omission and commission errors.

From Table VIII, we can see that the MSU resulted in the best performance in this data set, with respect to the highest OA (95.37%) and  $kappa$  (0.8826) values and lowest number of total errors (3114 pixels). This indicates the superior detection of spectral variations at a fine level in the MSU approach. The other two state-of-the-art methods also obtained a high level of OA and  $kappa$  values, which demonstrates the effectiveness of the considered methods. In practical applications, two pixel-level multiple CD techniques are sufficient to address the challenging multiple CD task, especially in an unsupervised/semisupervised fashion without using ground reference samples.

#### D. Challenges in Change Detection

For CD tasks in multitemporal HSIs, challenges come from the intrinsic properties of the hyperspectral data and design of sophisticated CD techniques to handle the complexities in CD process. From the hyperspectral data perspective, the high dimension inevitably leads to information redundancy and high computational cost. Moreover, changes are more implicitly to be represented and detected. In other words, from the spectral signature point of view, similar changes are more likely to be overlapped especially the subtle changes. Thus the difficulty to discriminate their classes in high dimensional feature space increases. From the CD methodology perspective, advanced approaches need to be designed, which can: 1) detect multiple complex changes (i.e., having different spectrally significance) more adaptively; 2) be implemented more automatically; and 3) be computationally effective. In particular, the development of unsupervised CD techniques for HSIs is more important for real applications. In fact, several sub-problems should be considered within a CD process, such as the identification of

the number of multiple changes, the separation of changed and unchanged pixels, and the discrimination of multi-class changes. Each of these sub-problems deserved to be investigated in details independently or simultaneously to generate more accurate CD results. In addition, the sub-pixel and super-pixel implementation of CD techniques at different detection scales are expected to enhance the CD performance. Learning from the limited prior change information transferred between the multitemporal HSIs could be another interesting research direction.

## IX. FAST COMPUTING

The high-dimensional nature of hyperspectral data sets, together with the complexity of the processing algorithms, call for advanced processing techniques for accelerating hyperspectral-related computations [311]. Traditionally, software for hyperspectral analysis has been written for serial computation, i.e., to be run on a single computer having a single central processing unit (CPU) in which a problem is broken into a discrete series of instructions. Instructions are then executed one after another, so that only one instruction may execute at any moment in time. In turn, parallel computing allows simultaneous use of multiple compute resources, i.e., to be run on multiple CPUs, a problem is broken into discrete parts so that each part can be executed simultaneously. Load balancing refers to the practice of distributing work among tasks so that all tasks are kept busy all of the time. This practice is considered to minimize task idle time.

Taking advantage of the aforementioned concepts, several techniques have been developed to accelerate hyperspectral imaging computation. In this section, we summarize some of the available strategies, which vary according to the adopted platform for fast computing and acceleration.

### A. Cluster Computers

Perhaps the most widely used type of high performance computing architecture employed to accelerate hyperspectral related computations is cluster computing. A cluster is a collection of commodity computers interconnected by a computer network. In order to efficiently execute a parallel problem in a cluster, partitioning strategies can be used so that the original problem is broken down into subtasks that are allocated to the different computers. In the case of hyperspectral imaging, two kinds of strategies have been used to partition the original hyperspectral data set into subsets for efficient cluster-based processing. In spectral-domain partitioning, a single pixel vector (spectral signature) may be stored in different processing units and communications would be required for individual pixel-based calculations, such as spectral angle computations [312]. In spatial-domain partitioning, every pixel vector (spectral signature) is stored in the same processing unit [313]. These concepts have not only been explored in clusters [314, 315], but also in heterogeneous networks of workstations [316].

## B. Hardware Accelerators

In addition to clusters, other kinds of high performance computing architectures have also been used. Specifically, in certain contexts onboard processing of HSIs is required. This is particularly the case in time-critical applications [317, 318]. For this purpose, other solutions have been explored. Particularly, reconfigurable computing provides higher performance (throughput and processing power) compared to other specialized hardware processors [319]. Reconfigurability is no longer a promise but a reality, and it is feasible to have several algorithms implemented on the same board, and to dynamically select one out of a pool of algorithms from a control station [320]. A field programmable gate array (FPGA) is a chip in which there is a matrix of blank cells called configurable logic blocks (CLBs). This device can be used to implement any circuit (provided there are a sufficient number of logic blocks). FPGAs have been widely used to accelerate hyperspectral imaging algorithms for onboard processing [321–323].

Another high performance computing architecture that has provided excellent performance when accelerating hyperspectral imaging computations has been the graphical processing unit (GPU) [324]. GPUs are now fully programmable using high-level languages such as NVIDIA CUDA.<sup>5</sup> The GPU specializes in computer-intensive, massive data parallel computation (which describes exactly graphics rendering). Therefore, more transistors can be devoted to data processing rather than data caching and flow control. The fast-growing video game industry exerts strong economic pressure for constant innovation. This has motivated the extended use of GPUs for accelerating many different hyperspectral imaging related tasks [317, 325–336].

## C. Cloud Computing

Recently, more sophisticated high performance architectures have been used for the processing of hyperspectral data. For instance, the use of cloud computing platforms for the processing of hyperspectral data in distributed architectures is becoming more widespread. In this sense, cloud computing offers advanced capabilities for service-oriented computing and high-performance computing [337]. Furthermore, the use of cloud computing for the analysis of large repositories of hyperspectral data can be considered a natural solution resulting from the evolution of techniques previously developed for other types of computing platforms [312]. In particular, the utilization of GPUs within distributed scenarios has been radically extended worldwide, thanks in part to the increasing development of deep learning based frameworks (e.g., Apache Spark, Caffe, Theano, Torch, and TensorFlow), which also have their application in the hyperspectral analysis community [64, 65, 84, 338, 339]. However, the recent literature still holds few examples of the use of cloud infrastructures for the implementation of hyperspectral analysis techniques in general and for the supervised classification of hyperspectral data in particular. This may be due to the lack of open repositories of HSIs available for public use, a situation that is expected to

change in the near future, since large distributed repositories of hyperspectral data for open use are expected to be available to the scientific community soon.

## D. Challenges in Fast Computing

The most important challenge related to the consolidation of fast computing techniques for the analysis of hyperspectral data (particularly in the context of real-time platforms) is still the high energy consumption required by high performance computing architectures, which reduces their applicability in real scenarios for onboard operation. Currently, the power consumption required by devices such as GPUs is too high for their incorporation into satellite platforms. Another shortcoming is the fact that these platforms are often subject to radiation tolerance issues. Future developments in hardware instruments for onboard operation are required for efficient real-time processing of HSIs, particularly in the context of satellite missions.

## X. CONCLUDING REMARKS

The role of hyperspectral image analysis cannot be underestimated for a plethora of applications, especially those related to change detection and scene classification. Without a doubt, the use of such valuable data has been well-established in the remote sensing community and the precise investigation of such data is increasing significantly. To this end, the considerable number of airborne and space-borne hyperspectral missions as well as the increasing number of scientific publications on this particular subject demonstrate that the area of hyperspectral image analysis is substantial, dynamic, and vibrant.

The field of hyperspectral imagery is extremely broad and it is impossible to investigate it comprehensively in one literature review. This paper focuses particularly on algorithmic approaches that have been developed, adapted, or proposed since 2013, covering a number of key research areas such as dimensionality reduction, classification, spectral unmixing, resolution enhancement, image restoration, change detection, and fast computing. It is indeed of high interest to summarize other survey papers from the application point of view where the usefulness of hyperspectral imagery can be demonstrated through different practical aspects such as mineralogy, environmental mapping and monitoring, geology, and so on.

In addition to the material that has been presented in this paper, hyperspectral data preprocessing plays a vital role in fostering application-oriented tasks. This important subject was out of the scope of this investigation, which mainly focused on the development of algorithms for hyperspectral image analysis and processing. To this end, we see the need for some tutorials and survey papers with a main focus on hyperspectral data preprocessing and preparation designed for atmospheric corrections, geometric and radiometric corrections, co-registration, and quality assessment.

As indicated several times throughout the paper, although the area of hyperspectral image analysis is well-established, there are still many doors left open for further investigation. We hope that this manuscript will raise new possibilities

<sup>5</sup><http://developer.nvidia.com>



for researchers to further investigate the remaining issues by developing fast, accurate, and automatic methodologies suitable for the application at hand.

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