

REPORT DOCUMENTATION PAGE			Form Approved OMB NO. 0704-0188		
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1. REPORT DATE (DD-MM-YYYY)		2. REPORT TYPE		3. DATES COVERED (From - To)	
		New Reprint		-	
4. TITLE AND SUBTITLE Hyperspectral Image Classification via Kernel Sparse Representation			5a. CONTRACT NUMBER		
			W911NF-11-1-0245		
			5b. GRANT NUMBER		
6. AUTHORS Yi Chen, Nasser M. Nasrabadi, Trac D. Tran			5c. PROGRAM ELEMENT NUMBER		
			611102		
			5d. PROJECT NUMBER		
7. PERFORMING ORGANIZATION NAMES AND ADDRESSES Johns Hopkins University 3400 N Charles Street Maryland Hall Batlimore, MD 21218 -2608			5e. TASK NUMBER		
			5f. WORK UNIT NUMBER		
			8. PERFORMING ORGANIZATION REPORT NUMBER		
9. SPONSORING/MONITORING AGENCY NAME(S) AND ADDRESS(ES) U.S. Army Research Office P.O. Box 12211 Research Triangle Park, NC 27709-2211			10. SPONSOR/MONITOR'S ACRONYM(S) ARO		
			11. SPONSOR/MONITOR'S REPORT NUMBER(S) 60291-MA.10		
12. DISTRIBUTION AVAILABILITY STATEMENT Approved for public release; distribution is unlimited.					
13. SUPPLEMENTARY NOTES The views, opinions and/or findings contained in this report are those of the author(s) and should not be construed as an official Department of the Army position, policy or decision, unless so designated by other documentation.					
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15. SUBJECT TERMS Classification, hyperspectral imagery, joint sparsity model, kernel methods, sparse representation					
16. SECURITY CLASSIFICATION OF:		17. LIMITATION OF ABSTRACT		15. NUMBER OF PAGES	19a. NAME OF RESPONSIBLE PERSON
a. REPORT	b. ABSTRACT	c. THIS PAGE	UU		Trac Tran
UU	UU	UU			19b. TELEPHONE NUMBER
					410-516-7416

Report Title

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ARO Report Number 60291.10-MA
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Block 13: Supplementary Note

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Hyperspectral Image Classification via Kernel Sparse Representation

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Abstract—In this paper, a novel nonlinear technique for hyperspectral image (HSI) classification is proposed. Our approach relies on sparsely representing a test sample in terms of all of the training samples in a feature space induced by a kernel function. For each test pixel in the feature space, a sparse representation vector is obtained by decomposing the test pixel over a training dictionary, also in the same feature space, by using a kernel-based greedy pursuit algorithm. The recovered sparse representation vector is then used directly to determine the class label of the test pixel. Projecting the samples into a high-dimensional feature space and kernelizing the sparse representation improve the data separability between different classes, providing a higher classification accuracy compared to the more conventional linear sparsity-based classification algorithms. Moreover, the spatial coherency across neighboring pixels is also incorporated through a kernelized joint sparsity model, where all of the pixels within a small neighborhood are jointly represented in the feature space by selecting a few common training samples. Kernel greedy optimization algorithms are suggested in this paper to solve the kernel versions of the single-pixel and multi-pixel joint sparsity-based recovery problems. Experimental results on several HSIs show that the proposed technique outperforms the linear sparsity-based classification technique, as well as the classical support vector machines and sparse kernel logistic regression classifiers.

Index Terms—Classification, hyperspectral imagery, joint sparsity model, kernel methods, sparse representation.

I. INTRODUCTION

HYPERSPECTRAL imaging sensors capture images in hundreds of continuous narrow spectral bands, spanning the visible to infrared spectrum. Each pixel in a hyperspectral image (HSI) is represented by a vector whose entries correspond to various spectral-band responses. Different materials usually reflect electromagnetic energy differently at specific wavelengths. This enables discrimination of materials based on their spectral characteristics. One of the most important applications of HSI is image classification, where pixels are

labeled to one of the classes based on their spectral characteristics, given a small set of training data for each class. Various techniques have been developed for HSI classification. Among the previous approaches, the support vector machine (SVM) [1], [2] has proven to be a powerful tool to solve many supervised classification problems and has shown good performances in hyperspectral classification, as well [3]–[5]. Variations of SVM-based algorithms have also been proposed to improve the classification accuracy. These variations include transductive SVM, which exploits both labeled and unlabeled samples [6], and SVM with composite kernels (CKs), which incorporates spatial information directly in the SVM kernels [7]. Multinomial logistic regression [8] is another widely used classifier, which uses the logistic function to provide the posterior probability. A fast algorithm for sparse multinomial logistic regression has been developed in [9] and successfully adopted for HSI segmentation in [10], [11]. Some of the other recent HSI classification techniques can be found in [12]–[17]. In these recent methods, a feature extraction strategy is proposed in [12] for classification which generalizes the linear discriminative analysis and nonparametric discriminative analysis. In [13], the derivative information of the spectral signatures is exploited as features, and then decisions obtained from spectral reflectance and derivative information are fused for the final decisions. In [14], each image band is decomposed into intrinsic mode functions (IMFs) which are adaptive to local properties via empirical mode decomposition and then SVM is applied to the lower order IMFs for classification. In [15], the k -nearest-neighbor classifier is applied to the local manifolds to exploit the intrinsic nonlinear structure of HSIs. A semi-supervised classification algorithm is proposed in [16] in order to use a kernel machine which is iteratively updated by manifold regularization. In [17] the results from multiple classification/segmentation techniques are fused by postprocessing to generate the final spectral-spatial classification map. Most of the aforementioned HSI image classification techniques do not directly incorporate the spatial or the contextual information into the classifier.

Recently, sparse representation [18], [19] has also been proposed to solve many computer vision tasks [20]–[25], where the usage of sparsity as a prior often leads to state-of-the-art performance. Sparse representation has also been applied to HSI target detection and classification [26]–[28], relying on the observation that hyperspectral pixels belonging to the same class approximately lie in the same low-dimensional subspace. Thus, an unknown test pixel can be sparsely represented by a few training samples (atoms) from a given dictionary, and the corresponding sparse representation vector will implicitly

Manuscript received August 13, 2011; revised December 23, 2011 and March 22, 2012; accepted April 18, 2012. Date of publication July 10, 2012; date of current version December 19, 2012. This work was supported in part by the National Science Foundation under Grants CCF-1117545 and CCF-0728893, the Army Research Office under Grant 58110-MA-II and Grant 60219-MA, and the Office of Naval Research under Grant N102-183-0208.

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Color versions of one or more of the figures in this paper are available online at <http://ieeexplore.ieee.org>.

Digital Object Identifier 10.1109/TGRS.2012.2201730

encode the class information. The sparse representation-based classifier is different from the conventional sparse classifier SVM in the following aspects. SVM is a discriminative model, while the sparse representation method can be viewed as a generative model, where the signal (pixel) is expressed as a linear combination of atoms [19]. SVM is a binary classifier that finds the separating hyperplane between two classes (multi-class SVM requires a one-against-one or one-against-all strategy). The sparse representation-based classifier is from a reconstruction point of view. The sparse decomposition of the test pixel over the entire dictionary implicitly leads to a competition between the subspaces (classes), and thus the recovered sparse representation is discriminative. Moreover, in SVM, there is an explicit training stage. The SVM classifier is trained only once, and then this classifier with its fixed sparse support vectors is used to classify all of the test data. On the other hand, in our proposed approach, a new sparse representation vector is extracted for each test pixel and is thus adaptive, representing the sparsely selected atoms which are adapted to reconstruct the current test pixel.

HSIs are usually smooth in the sense that the pixels in a small neighborhood represent the same material and have similar spectral characteristics. Various techniques have been proposed recently to exploit the contextual correlation within HSI which have notably improved the classification and segmentation performance. Post-processing procedures are used in [29], [30] on the individually labeled samples based on certain decision rules to impose the smoothness. Markov random fields exploit the statistical dependency among neighboring pixels and are usually applied in Bayesian approaches [11]. The CK approach [7] is another way to incorporate the spatial information, which explicitly extracts spatial information for each spectral pixel and then combines the spectral and spatial information via kernel composition. Joint sparsity model (JSM) [31] is exploited in sparsity-based HSI target detection and classification [27], [28], where the neighboring pixels are simultaneously represented by a sparse linear combination of a few common training samples. Each pixel, although sharing the same common support, might have weighting coefficients taking on different values. In this way, the smoothness across neighboring spectral pixels is enforced directly in the classification stage, and no post-processing steps are performed. The details of CKs and the JSM will be further discussed in the following sections.

It is well-known that for the classical HSI image classification and target detection algorithms, the use of kernel methods yields a significant performance improvement [5], [32], because the kernel-based algorithms implicitly exploit the higher order structure of the given data which may not be captured by the linear models. Therefore, if the data set is not linearly separable, kernel methods [33]–[36] can be applied to project the data into a nonlinear feature space in which the data becomes more separable. In practical implementation, the kernel trick [37] is often used in order to avoid explicitly evaluating the data in the feature space.

In this paper, we propose a new HSI classification algorithm based on kernel sparse representation by assuming that a test pixel can be linearly represented by a few training samples in the feature space. The kernel sparse representation vector

is then obtained by decomposing the test pixel represented in a high-dimensional feature space over a structured dictionary consisting of training samples from all of the classes in the same feature space. The recovered sparse vector is used directly for classification. Although the proposed approach has a similar formulation as previous kernel regression approaches with a sparse prior such as kernel matching pursuit [33], kernel basis pursuit [34], and generalized LASSO [38], the underlying ideas are quite different. The objective of these previous approaches is to approximate a function as a linear combination of dictionary functions, which are the kernels centered at the training points, by minimizing certain loss function evaluated at these training points and subject to a sparsity prior. Therefore, the target vector for fitting consists of the observations of the function value at the training points, and the dictionary is then the dictionary functions evaluated at the training points which turns out to be the kernel matrix. In our proposed approach, the target vector is the test pixel itself in the feature space. It is not the similarity measure between the test sample and training samples and may not have an explicit expression. The dictionary also consists of the training samples in the feature space and cannot assume an explicit expression either. The recovered sparse representation vector can be viewed as a discriminative feature extracted from the test pixel and is used directly for classification.

The contextual correlation between pixels within a small spatial neighborhood can be incorporated into the kernel sparse representation through the JSM [31], where all neighboring pixels are simultaneously represented by a linear combination of a few common training samples in the feature space. Furthermore, the CK approach [7] can also be used with the proposed kernel sparse representation model in order to combine spectral and spatial information. Efficient kernel-based optimization algorithms are discussed in this paper for the recovery of the kernel sparse representations for both single-pixel and multi-pixel JSMs.

Notation-wise, vectors and matrices are denoted by lower and upper case bold letters, respectively. For a vector $\alpha \in \mathbb{R}^N$ and an index set $\Lambda \subseteq \{1, \dots, N\}$ with $|\Lambda| = t$, $\alpha_\Lambda \in \mathbb{R}^t$ is the portion of α indexed on Λ . For a matrix $S \in \mathbb{R}^{N_1 \times N_2}$, index sets $\Lambda_1 \subseteq \{1, \dots, N_1\}$ with $|\Lambda_1| = t_1$, and $\Lambda_2 \subseteq \{1, \dots, N_2\}$ with $|\Lambda_2| = t_2$, $S_{\Lambda_1, :} \in \mathbb{R}^{t_1 \times N_2}$ is a submatrix of S consisting of the t_1 rows in S indexed on Λ_1 , $S_{:, \Lambda_2} \in \mathbb{R}^{N_1 \times t_2}$ consists of the t_2 columns in S indexed on Λ_2 , and $S_{\Lambda_1, \Lambda_2} \in \mathbb{R}^{t_1 \times t_2}$ is formed by the rows and columns of S indexed on Λ_1 and Λ_2 , respectively.

The remainder of this paper is structured as follows. Section II briefly introduces the sparsity-based HSI classification technique. Section III defines the sparsity models in the feature space, then discusses how to incorporate spatial information, and describes the kernel sparse recovery algorithms. Experimental results are shown in Section IV, and conclusions are drawn in Section V.

II. SPARSITY-BASED HSI CLASSIFICATION

This section briefly introduces the sparsity-based algorithm for HSI classification, and more details can be found in

[26]–[28]. It is assumed that the spectral signatures of pixels belonging to the same class approximately lie in the same low-dimensional subspace. Thus, an unknown test sample $\mathbf{x} \in \mathbb{R}^B$, where B is the number of spectral bands, can be written as a sparse linear combination of all of the training pixels as

$$\mathbf{x} = \mathbf{A}\boldsymbol{\alpha} \quad (1)$$

where $\mathbf{A} = [\mathbf{a}_1 \ \mathbf{a}_2 \ \dots \ \mathbf{a}_N] \in \mathbb{R}^{B \times N}$ is a structured dictionary whose columns $\{\mathbf{a}_i\}_{i=1,2,\dots,N}$ are N training samples (referred to as atoms) from all classes, and $\boldsymbol{\alpha} \in \mathbb{R}^N$ is an unknown sparse vector. The index set on which $\boldsymbol{\alpha}$ have nonzero entries is the support of $\boldsymbol{\alpha}$. The number of nonzero entries in $\boldsymbol{\alpha}$ is called the sparsity level K of $\boldsymbol{\alpha}$ and denoted by $K = \|\boldsymbol{\alpha}\|_0$. Given the dictionary \mathbf{A} , the sparse coefficient vector $\boldsymbol{\alpha}$ is obtained by solving

$$\hat{\boldsymbol{\alpha}} = \arg \min \|\mathbf{x} - \mathbf{A}\boldsymbol{\alpha}\|_2 \quad \text{subject to} \quad \|\boldsymbol{\alpha}\|_0 \leq K_0 \quad (2)$$

where K_0 is a preset upper bound on the sparsity level. The problem in (2) is NP-hard, which can be approximately solved by greedy algorithms, such as orthogonal matching pursuit (OMP) [39] or subspace pursuit (SP) [40]. The class label of \mathbf{x} is determined by the minimal residual between \mathbf{x} and its approximation from each class subdictionary:

$$\text{Class}(\mathbf{x}) = \arg \min_{m=1,\dots,M} \|\mathbf{x} - \mathbf{A}_{:, \Omega_m} \hat{\boldsymbol{\alpha}}_{\Omega_m}\|_2 \quad (3)$$

where $\Omega_m \subset \{1, 2, \dots, N\}$ is the index set associated with the training samples belonging to the m th class. As pointed out in [25], the sparse representation-based classifier can be viewed as a generalization of the nearest neighbor classifier [41].

In HSI, pixels within a small neighborhood usually consist of similar materials and, thus, their spectral characteristics are highly correlated. The spatial correlation between neighboring pixels can be incorporated through a JSM [27], [31] by assuming the underlying sparse vectors associated with these pixels share a common sparsity pattern as follows. Let $\{\mathbf{x}_t\}_{t=1,\dots,T}$ be T pixels in a spatial neighborhood centered at \mathbf{x}_1 . These pixels can be compactly represented as

$$\begin{aligned} \mathbf{X} &= [\mathbf{x}_1 \ \mathbf{x}_2 \ \dots \ \mathbf{x}_T] = [\mathbf{A}\boldsymbol{\alpha}_1 \ \mathbf{A}\boldsymbol{\alpha}_2 \ \dots \ \mathbf{A}\boldsymbol{\alpha}_T] \\ &= \mathbf{A} \underbrace{[\boldsymbol{\alpha}_1 \ \boldsymbol{\alpha}_2 \ \dots \ \boldsymbol{\alpha}_T]}_{\mathbf{S}} = \mathbf{A}\mathbf{S}. \end{aligned} \quad (4)$$

In the JSM, the sparse vectors $\{\boldsymbol{\alpha}_t\}_{t=1,\dots,T}$ share the same support Λ , and, thus, \mathbf{S} is a sparse matrix with only $|\Lambda|$ nonzero rows. The row-sparse matrix \mathbf{S} can be recovered by solving the following optimization problem

$$\hat{\mathbf{S}} = \arg \min \|\mathbf{X} - \mathbf{A}\mathbf{S}\|_F \quad \text{subject to} \quad \|\mathbf{S}\|_{\text{row},0} \leq K_0 \quad (5)$$

where $\|\mathbf{S}\|_{\text{row},0}$ denotes the number of non-zero rows of \mathbf{S} , and $\|\cdot\|_F$ denotes the Frobenius norm. The problem in (5) can be approximately solved by the simultaneous versions of OMP (SOMP) [31] or SP (SSP) [28]. The label of the center pixel \mathbf{x}_1 is then determined by the minimal total residual

$$\text{Class}(\mathbf{x}_1) = \arg \min_{m=1,\dots,M} \left\| \mathbf{X} - \mathbf{A}_{:, \Omega_m} \hat{\mathbf{S}}_{\Omega_m} \right\|_F \quad (6)$$

where $\|\cdot\|_F$ denotes the Frobenius norm.

III. KERNEL SPARSE REPRESENTATION

If the classes in the dataset are not linearly separable, then the kernel methods can be used to project the data into a feature space, in which the classes become linearly separable [1]. The kernel function $\kappa : \mathbb{R}^B \times \mathbb{R}^B \mapsto \mathbb{R}$ is defined as the inner product

$$\kappa(\mathbf{x}_i, \mathbf{x}_j) = \langle \phi(\mathbf{x}_i), \phi(\mathbf{x}_j) \rangle. \quad (7)$$

Commonly used kernels include the radial basis function (RBF) kernel $\kappa(\mathbf{x}_i, \mathbf{x}_j) = \exp(-\gamma\|\mathbf{x}_i - \mathbf{x}_j\|^2)$ with $\gamma > 0$ controlling the width of the RBF, and order- d homogeneous and inhomogeneous polynomial kernels $\kappa(\mathbf{x}_i, \mathbf{x}_j) = (\mathbf{x}_i \cdot \mathbf{x}_j)^d$ and $\kappa(\mathbf{x}_i, \mathbf{x}_j) = (\mathbf{x}_i \cdot \mathbf{x}_j + 1)^d$, respectively. In this section, we describe how the sparsity models in Section II can be extended to a feature space induced by a kernel function.

A. Pixel-Wise Sparsity in Feature Space

Let $\mathbf{x} \in \mathbb{R}^B$ be the data point of interest and $\phi(\mathbf{x})$ be its representation in the feature space. The kernel sparse representation of a sample \mathbf{x} in terms of training atoms \mathbf{a}_i 's can be formulated as

$$\phi(\mathbf{x}) = \underbrace{[\phi(\mathbf{a}_1) \ \dots \ \phi(\mathbf{a}_N)]}_{\mathbf{A}_\phi} \underbrace{[\alpha'_1 \ \dots \ \alpha'_N]^T}_{\boldsymbol{\alpha}'} = \mathbf{A}_\phi \boldsymbol{\alpha}' \quad (8)$$

where the columns of \mathbf{A}_ϕ are the representations of training samples in the feature space and $\boldsymbol{\alpha}'$ is assumed to be a sparse vector.

Similar to the linear sparse recovery problem in (2), $\boldsymbol{\alpha}'$ can be recovered by solving

$$\hat{\boldsymbol{\alpha}'} = \arg \min \|\phi(\mathbf{x}) - \mathbf{A}_\phi \boldsymbol{\alpha}'\|_2 \quad \text{subject to} \quad \|\boldsymbol{\alpha}'\|_0 \leq K_0. \quad (9)$$

The problem in (9) can be approximately solved by kernelizing the OMP and SP algorithms (denoted by KOMP and KSP, respectively). Note that in the above problem formulation, we are solving for the sparse vector $\boldsymbol{\alpha}'$ directly in the feature space using the implicit feature vectors, but not evaluating the kernel functions at the training points.

In KOMP and KSP, essentially, each dot product operation in OMP/SP is replaced by the kernel trick in (7). Let $\mathbf{K}_A \in \mathbb{R}^{N \times N}$ be the kernel matrix whose (i, j) th entry is $\kappa(\mathbf{a}_i, \mathbf{a}_j)$, and $\mathbf{k}_{A, \mathbf{x}} \in \mathbb{R}^N$ be the vector whose i th entry is $\kappa(\mathbf{a}_i, \mathbf{x})$. Using the feature representations, the correlation (dot product) between a pixel $\phi(\mathbf{x})$ and a dictionary atom $\phi(\mathbf{a}_i)$ is then computed by

$$c_i = \langle \phi(\mathbf{x}), \phi(\mathbf{a}_i) \rangle = \kappa(\mathbf{x}, \mathbf{a}_i) = (\mathbf{k}_{A, \mathbf{x}})_i \quad (10)$$

the orthogonal projection coefficient of $\phi(\mathbf{x})$ onto a set of selected dictionary atoms $\{\phi(\mathbf{a}_n)\}_{n \in \Lambda}$ is given as

$$\mathbf{p}_\Lambda = \left((\mathbf{K}_A)_{\Lambda, \Lambda} \right)^{-1} (\mathbf{k}_{A, \mathbf{x}})_\Lambda \quad (11)$$

and the residual vector between $\phi(\mathbf{x})$ and its approximation using the selected atoms $\{\phi(\mathbf{a}_n)\}_{n \in \Lambda} = (\mathbf{A}_\phi)_{:, \Lambda}$ is then expressed as

$$\phi(\mathbf{r}) = \phi(\mathbf{x}) - (\mathbf{A}_\phi)_{:, \Lambda} \left((\mathbf{K}_A)_{\Lambda, \Lambda} \right)^{-1} (\mathbf{k}_{A, \mathbf{x}})_\Lambda. \quad (12)$$

Note that the feature representation of the residual vector $\phi(\mathbf{r})$ in (12) cannot be evaluated explicitly. However, the correlation between $\phi(\mathbf{r})$ and an atom $\phi(\mathbf{a}_i)$ can be computed by

$$\begin{aligned} c_i &= \langle \phi(\mathbf{r}), \phi(\mathbf{a}_i) \rangle \\ &= (\mathbf{k}_{A, \mathbf{x}})_i - (\mathbf{K}_A)_{i, \Lambda} \left((\mathbf{K}_A)_{\Lambda, \Lambda} \right)^{-1} (\mathbf{k}_{A, \mathbf{x}})_\Lambda. \end{aligned} \quad (13)$$

The KOMP and KSP greedy algorithms, similar to the linear OMP and SP algorithms, are used to locate the support Λ of the sparse vector $\hat{\alpha}'$. The KOMP algorithm augments the support set by only one index, which is given by $\lambda = \arg \max_{i=1, \dots, N} c_i$ with c_i being defined in (13) and $\phi(\mathbf{r})$ being the residual vector from the previous iteration, at each iteration until K_0 atoms are selected or the approximation error (i.e., norm of the residual vector in (12)) is within a preset threshold. The KSP algorithm maintains a set of K_0 indices with a backtracking mechanism. At each iteration, the index set is refined by adding K_0 new candidates, whose associated atoms have the K_0 highest correlation (13) to the residual vector from the previous iteration, to the current list and then discarding K_0 insignificant ones from the list of $2K_0$ candidates. This process repeats until certain stopping criterion is met. In both of the KOMP and KSP algorithms, after the support set Λ of $\hat{\alpha}'$ is determined, the entries of $\hat{\alpha}'$ indexed on Λ are computed by the orthogonal projection of the test pixel onto the selected dictionary atoms using (11). The KOMP/KSP algorithms can be viewed as special cases, with $T = 1$, of the kernelized SOMP/SSP (KSOMP/KSSP) algorithms (Algorithms 1 and 2) proposed in the next section, respectively. The details are thus omitted herein.

Once the sparse vector $\hat{\alpha}'$ is recovered, the residual between the test sample and the m th-class reconstruction in the high-dimensional feature space is then computed by

$$\begin{aligned} r_m(\mathbf{x}) &= \left\| \phi(\mathbf{x}) - (\mathbf{A}_\phi)_{:, \Omega_m} \hat{\alpha}'_{\Omega_m} \right\| \\ &= \left\langle \phi(\mathbf{x}) - (\mathbf{A}_\phi)_{:, \Omega_m} \hat{\alpha}'_{\Omega_m}, \phi(\mathbf{x}) - (\mathbf{A}_\phi)_{:, \Omega_m} \hat{\alpha}'_{\Omega_m} \right\rangle^{\frac{1}{2}} \\ &= \left(\kappa(\mathbf{x}, \mathbf{x}) - 2\hat{\alpha}'_{\Omega_m}{}^T (\mathbf{k}_{A, \mathbf{x}})_{\Omega_m} \right. \\ &\quad \left. + \hat{\alpha}'_{\Omega_m}{}^T (\mathbf{K}_A)_{\Omega_m, \Omega_m} \hat{\alpha}'_{\Omega_m} \right)^{\frac{1}{2}} \end{aligned} \quad (14)$$

where $\mathbf{k}_{A, \mathbf{x}}$ and \mathbf{K}_A are as defined above, and Ω_m is the index set associated with the m th class. The class label of \mathbf{x} is determined as

$$\text{Class}(\mathbf{x}) = \arg \min_{m=1, \dots, M} r_m(\mathbf{x}). \quad (15)$$

B. Joint Sparsity in Feature Space

The JSM in (4) can also be extended to the feature space as follows:

$$\begin{aligned} \mathbf{X}_\phi &= [\phi(\mathbf{x}_1) \dots \phi(\mathbf{x}_T)] = [\mathbf{A}_\phi \alpha'_1 \dots \mathbf{A}_\phi \alpha'_T] \\ &= \mathbf{A}_\phi \underbrace{[\alpha'_1 \dots \alpha'_T]}_{\mathbf{S}'} = \mathbf{A}_\phi \mathbf{S}' \end{aligned} \quad (16)$$

where the vectors $\{\alpha'_t\}_{t=1, \dots, T}$ share the same support. The row-sparse matrix \mathbf{S}' is recovered by solving

$$\hat{\mathbf{S}}' = \arg \min \|\mathbf{X}_\phi - \mathbf{A}_\phi \mathbf{S}'\|_F \quad \text{subject to} \quad \|\mathbf{S}'\|_{\text{row}, 0} \leq K_0. \quad (17)$$

In this paper, we propose the KSOMP and the KSSP algorithms in order to approximately solve the above joint sparse recovery problem in (17).

In KSOMP, at every iteration, the atom that simultaneously yields the best approximation to all the T pixels (or residuals after initialization) is selected. Specifically, let $\mathbf{C} \in \mathbb{R}^{N \times T}$ be the correlation matrix whose (i, j) th entry is the correlation between $\phi(\mathbf{a}_i)$ and $\phi(\mathbf{r}_j)$, where $\phi(\mathbf{r}_j)$ is the residual vector of $\phi(\mathbf{x}_j)$. The new atom is then selected as the one associated with the row of \mathbf{C} , which has the maximal ℓ_p -norm for some $p \geq 1$. The KSOMP algorithm is summarized in Algorithm 1. Note that when computing the projection in (11) and correlation in (13), a regularization term $\lambda \mathbf{I}$ is added in order to stabilize the matrix inversion, where λ is typically a small scalar (e.g., $\lambda = 10^{-5}$ is implemented in Section IV) and \mathbf{I} is an identity matrix whose dimensionality should be clear from the context. The parameter λ does not seriously affect the classification performance. It is, however, necessary for stable matrix inversion. Although the kernel matrix \mathbf{K}_A is usually invertible, $(\mathbf{K}_A)_{\Lambda, \Lambda}$ may be nearly singular for certain small index set Λ .

Algorithm 1: KSOMP

Input: $B \times N$ dictionary $\mathbf{A} = [\mathbf{a}_1 \dots \mathbf{a}_N]$, $B \times T$ data matrix $\mathbf{X} = [\mathbf{x}_1 \dots \mathbf{x}_T]$, kernel function κ , and a stopping criterion

Initialization: compute the kernel matrices $\mathbf{K}_A \in \mathbb{R}^{N \times N}$ whose (i, j) th entry is $\kappa(\mathbf{x}_i, \mathbf{x}_j)$ and $\mathbf{K}_{A, \mathbf{X}} \in \mathbb{R}^{N \times T}$ whose (i, j) th entry is $\kappa(\mathbf{a}_i, \mathbf{x}_j)$. Set index set $\Lambda_0 = \arg \max_{i=1, \dots, N} \|(\mathbf{K}_{A, \mathbf{X}})_{i, :}\|_p$ with some $p \geq 1$ and iteration counter $t = 1$.

while stopping criterion has not been met **do**

1) Compute the correlation matrix $\mathbf{C} \in \mathbb{R}^{N \times T}$

$$\begin{aligned} \mathbf{C} &= \mathbf{K}_{A, \mathbf{X}} - (\mathbf{K}_A)_{:, \Lambda_{t-1}} \left((\mathbf{K}_A)_{\Lambda_{t-1}, \Lambda_{t-1}} + \lambda \mathbf{I} \right)^{-1} \\ &\quad \times (\mathbf{K}_{A, \mathbf{X}})_{\Lambda_{t-1}, :} \end{aligned}$$

2) Select the new index as $\lambda_t = \arg \max_{i=1, \dots, N} \|\mathbf{C}_{i, :}\|_p$, $p \geq 1$

3) Update the index set $\Lambda_t = \Lambda_{t-1} \cup \{\lambda_t\}$

4) $t \leftarrow t + 1$

end while

Output: Index set $\Lambda = \Lambda_{t-1}$, the sparse representation $\hat{\mathbf{S}}'$ whose nonzero rows indexed by Λ are $\hat{\mathbf{S}}'_{\Lambda,:} = (\mathbf{K}_{\Lambda,\Lambda} + \lambda\mathbf{I})^{-1}(\mathbf{K}_{\Lambda,\mathbf{X}})_{\Lambda,:}$.

Similarly, KSSP is a simultaneous version of KSP where the K_0 atoms that best simultaneously approximate all of the T residuals in terms of the ℓ_p -norm are chosen. The KSSP algorithm is summarized in Algorithm 2. Note that the step for computing the residual vectors (12) is incorporated into the computation of the correlation vector in Step (1) of both KSOMP and KSSP.

Algorithm 2: KSSP

Input: $B \times N$ dictionary $\mathbf{A} = [\mathbf{a}_1 \dots \mathbf{a}_N]$, $B \times T$ data matrix $\mathbf{X} = [\mathbf{x}_1 \dots \mathbf{x}_T]$, kernel function κ , and a stopping criterion

Initialization: compute the kernel matrices $\mathbf{K}_{\mathbf{A}} \in \mathbb{R}^{N \times N}$ whose (i, j) th entry is $\kappa(\mathbf{x}_i, \mathbf{x}_j)$ and $\mathbf{K}_{\mathbf{A},\mathbf{X}} \in \mathbb{R}^{N \times T}$ whose (i, j) th entry is $\kappa(\mathbf{a}_i, \mathbf{x}_j)$. Set index set $\Lambda_0 = \{K_0 \text{ indices corresponding to the } K_0 \text{ largest numbers in } \|(\mathbf{K}_{\mathbf{A},\mathbf{X}})_{i,:}\|_p, p \geq 1, i = 1, \dots, N\}$, and set iteration counter $t = 1$.

while stopping criterion has not been met **do**

1) Compute the correlation matrix

$$\mathbf{C} = \mathbf{K}_{\mathbf{A},\mathbf{X}} - (\mathbf{K}_{\mathbf{A}})_{:, \Lambda_{t-1}} \left((\mathbf{K}_{\mathbf{A}})_{\Lambda_{t-1}, \Lambda_{t-1}} + \lambda\mathbf{I} \right)^{-1} \times (\mathbf{K}_{\mathbf{A},\mathbf{X}})_{\Lambda_{t-1}, :}, \in \mathbb{R}^{N \times T}$$

2) Find the index set $\mathcal{I} = \{K_0 \text{ indices corresponding to the } K_0 \text{ largest numbers in } \|\mathbf{C}_{i,:}\|_p, p \geq 1, i = 1, \dots, N\}$

3) Update the candidate index set $\tilde{\Lambda}_t = \Lambda_{t-1} \cup \mathcal{I}$

4) Compute the projection coefficients $\mathbf{P} = ((\mathbf{K}_{\mathbf{A}})_{\tilde{\Lambda}_t, \tilde{\Lambda}_t} + \lambda\mathbf{I})^{-1}(\mathbf{K}_{\mathbf{A},\mathbf{X}})_{\tilde{\Lambda}_t, :}, \in \mathbb{R}^{2K_0 \times T}$

5) Update the index set $\Lambda_t = \{K_0 \text{ indices in } \tilde{\Lambda}_t \text{ corresponding to the } K_0 \text{ largest numbers in } \|\mathbf{P}_{i,:}\|_p, p \geq 1, i = 1, \dots, N\}$

6) $t \leftarrow t + 1$

end while

Output: Index set $\Lambda = \Lambda_{t-1}$, the sparse representation $\hat{\mathbf{S}}'$ whose nonzero rows indexed by Λ are $\hat{\mathbf{S}}'_{\Lambda,:} = (\mathbf{K}_{\Lambda,\Lambda} + \lambda\mathbf{I})^{-1}(\mathbf{K}_{\Lambda,\mathbf{X}})_{\Lambda,:}$.

Once the matrix $\hat{\mathbf{S}}'$ is recovered, the total residual between the T neighboring pixels and their approximations from the m th-class training samples is computed by

$$r_m(\mathbf{x}_1) = \left(\sum_{i=1}^T \left(\kappa(\mathbf{x}_i, \mathbf{x}_i) - 2\hat{\mathbf{S}}_{\Omega_m, i}^{\prime T} (\mathbf{K}_{\mathbf{A},\mathbf{X}})_{\Omega_m, i} + \hat{\mathbf{S}}_{\Omega_m, i}^{\prime T} (\mathbf{K}_{\mathbf{A}})_{\Omega_m, \Omega_m} \hat{\mathbf{S}}'_{\Omega_m, i} \right) \right)^{\frac{1}{2}} \quad (18)$$

where $\mathbf{K}_{\mathbf{A},\mathbf{X}}$ and $\mathbf{K}_{\mathbf{A}}$ are as defined in Algorithms 1 and 2, and $\Omega_m \in \{1, 2, \dots, N\}$ is the index set associated with the m th class. The label for the center pixel \mathbf{x}_1 is then determined by the total residual

$$\text{Class}(\mathbf{x}_1) = \arg \min_{m=1, \dots, M} r_m(\mathbf{x}_1). \quad (19)$$

C. Kernel Sparse Representation With a Composite Kernel

Another way to address the contextual correlation within HSI is through a CK [7], which takes into account the spatial correlation between neighboring pixels by combining kernels dedicated to the spectral and spatial information. The CK approach has been shown to significantly outperform the spectral-only classifier in HSI classification [42]. This method, although originally proposed for SVM, can be readily incorporated into other classifiers which operate in the feature space, such as kernel logistic regression (KLR) and the kernel sparse representation-based classifier proposed in this paper. Specifically, let \mathbf{x}_i^w be the spectral pixel at location i in a HSI and \mathbf{x}_i^s be the spatial information extracted from a small neighborhood centered at location i , which is usually the mean and/or the standard deviation of the pixels within the neighborhood. The new pixel entity at this location can be redefined as $\mathbf{x}_i = \{\mathbf{x}_i^w, \mathbf{x}_i^s\}$. Note that in previous sections, \mathbf{x}_i contains only spectral information (i.e., $\mathbf{x}_i = \mathbf{x}_i^w$). The spectral and spatial information can then be combined in a variety of ways, including stacking, direct summation, weighted summation, and cross-information kernels [7]. In this paper, we consider the weighted summation kernel, which is shown to yield the best classification performance compared to other types of CKs [7]. The kernel function in this case is

$$\kappa(\mathbf{x}_i, \mathbf{x}_j) = \mu\kappa_s(\mathbf{x}_i^s, \mathbf{x}_j^s) + (1 - \mu)\kappa_w(\mathbf{x}_i^w, \mathbf{x}_j^w) \quad (20)$$

where $\mu \in (0, 1)$, and κ_s and κ_w are the kernel functions of the spatial and spectral features, respectively.

The CKs can be directly applied to the pixel-wise sparsity model in the feature space in (8). The sparse representation vector can be recovered using the KOMP or KSP algorithm, where the kernel matrix $\mathbf{K}_{\mathbf{A}}$ is now a weighted summation of the spectral and spatial kernel matrices of the training dictionary \mathbf{A} , and the vector $\mathbf{k}_{\mathbf{A},\mathbf{x}}$ also needs to be modified accordingly.

It is worth noting that the CK approach is different from the kernel JSM discussed in Section III-B in terms of incorporating localized spatial features. The JSM involves only the spatial information of the test pixels, and no prior knowledge about the neighbors of the training pixels is needed. On the other hand, for the CKs, the local spatial information for both training and test sets are necessary. Moreover, the JSM does not assume a sum or average of the same samples, but treats all pixels in a small neighborhood equally and finds the sparsity pattern that simultaneously represents these pixels. The CK approach, however, is not limited to local spatial features and can also be used to combine other types of features (e.g., centroid of the cluster as a global feature) and information sources.

TABLE I
THE 16 GROUND-TRUTH CLASSES IN THE AVIRIS INDIAN PINES IMAGE

No	Class	Samples	
		Train	Test
1	Alfalfa	6	48
2	Corn-notill	144	1290
3	Corn-min	84	750
4	Corn	24	210
5	Grass/Pasture	50	447
6	Grass/Trees	75	672
7	Grass/Pasture-mowed	3	23
8	Hay-windrowed	49	440
9	Oats	2	18
10	Soybeans-notill	97	871
11	Soybeans-min	247	2221
12	Soybean-clean	62	552
13	Wheat	22	190
14	Woods	130	1164
15	Building-Grass-Trees-Drives	38	342
16	Stone-steel Towers	10	85
Total		1043	9323

IV. EXPERIMENTAL RESULTS

In this section, we show the effectiveness of the proposed algorithms on classification of several hyperspectral data sets. For each image, we solve the sparse recovery problems in (2), (5), (9), and (17) for each test sample, and then determine the class by the minimal residual (the results are denoted by OMP/SP, KOMP/KSP, SOMP/SSP, and KSOMP/KSSP, respectively). The results of KOMP and KSP with CKs, as discussed in Section III-C, are denoted by KOMPCK and KSPCK, respectively. The classification results are then compared visually and quantitatively to those obtained by the classical SVM classifier and sparse multinomial KLR. For SVM and KLR classifiers, we use a spectral-only kernel (denoted by SVM/KLR), as well as a CK (denoted by SVMCK/KLRCK). In all classifiers with a CK, we use a weighted summation kernel, and the spatial information is the mean of pixels in a small neighborhood. The parameters for KLR, KLRCK, SVM, and SVMCK are obtained by cross-validation.

The first HSI in our experiments is the Airborne Visible/Infrared Imaging Spectrometer (AVIRIS) image Indian Pines [43]. The AVIRIS sensor generates 220 bands across the spectral range from 0.2 to 2.4 μm . In the experiments, the number of bands is reduced to 200 by removing 20 water absorption bands. This image has spatial resolution of 20 m per pixel and spatial dimension 145×145 . It contains 16 ground-truth classes. For each class, we randomly choose around 10% of the labeled samples for training and use the remaining 90% for testing, as shown in Table I and Fig. 1. RBF kernels are used in all kernel-based classifiers (i.e., SVM, SVMCK, KLR, KLRCK, KOMP, KSP, KSOMP, KSSP, KOMPCK, and KSPCK). Since this image consists of large homogenous regions, a large spatial window of size 9×9 ($T = 81$) is used in classifiers with a CK and the JSMs (4) and (16).

The classification performance for each of the 16 classes, overall accuracy (OA), average accuracy (AA), and the κ coefficient measure [44] on the test set are shown in Table II. The OA is the ratio between correctly classified test samples and the total number of test samples, the AA is the mean of the 16 class accuracies, and the κ coefficient is a robust measure

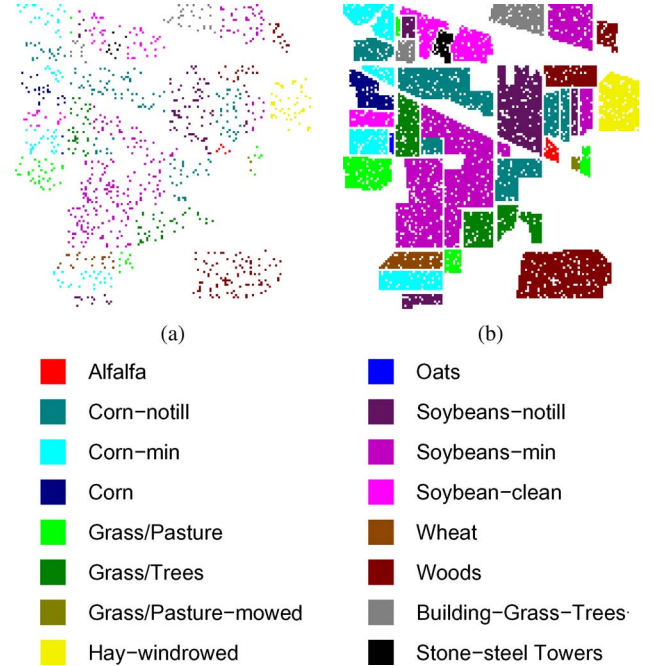


Fig. 1. (a) Training and (b) test sets for the Indian Pines image.

of the degree of agreement. The classification maps on labeled pixels are presented in Fig. 2, where the algorithm and OA are shown on top of each corresponding map. One can clearly see that incorporating the contextual correlation and operating in the feature space both have significantly improved the classification accuracy. The KOMPCK and KSPCK algorithms outperform all other classifiers—the OA for both of which are greater than 98%. The KSOMP and KSSP algorithms also yield superior performance, which have only about 1% lower OA than KOMPCK and KSPCK. Note that the kernel JSM for KSOMP and KSSP does not assume any prior knowledge of the neighbors of the training samples as the CK approach does.

The sparsity level K_0 and RBF parameter γ used in the above experiments are obtained from a small validation set. An n -fold cross validation would not be appropriate for finding the optimal sparsity level, unless n is large (e.g., leave-one-out cross validation). This is because the sparsity level K_0 is related to the size of dictionary, therefore, the optimal K_0 for part of the dictionary may not be optimal anymore for the entire dictionary. Now, we examine how these two parameters affect the classification performance on the Indian Pines image. We use randomly selected 10% of all labeled samples as the training set and the remaining samples as the test set, then vary K_0 from 5 to 80 and γ from 2^{-3} to 2^{12} in KOMP, KSP, KSOMP, and KSSP. The experiment for each γ , K_0 , and each of the four algorithm is repeated five times using different randomly chosen training sets to avoid any bias induced by random sampling. The window size is fixed at 9×9 for KSOMP and KSSP due to its smoothness. The OA on the test set, averaged over five independent realizations, are shown in Fig. 3. The bars indicate the maximal and minimal accuracies in five runs at each point, and we see that the fluctuation is usually within 2% and within 1% in a majority of cases. One can observe from Fig. 3(a) and (b) that for the pixel-wise kernel sparsity model,

TABLE II
CLASSIFICATION ACCURACY (%) FOR THE INDIAN PINES IMAGE USING 10% TRAINING SAMPLES AS SHOWN IN FIG. 1 AND TABLE I

Class	SVM	SVMCK	KLR	KLRCK	OMP	KOMP	SOMP	KSOMP	KOMPCK	SP	KSP	SSP	KSSP	KSPCK
1	81.25	95.83	64.58	75.00	68.75	72.92	85.42	97.92	97.92	68.75	72.92	81.25	91.67	95.83
2	86.28	96.67	89.46	96.43	65.97	86.36	94.88	97.21	99.22	74.65	87.91	95.74	97.98	99.15
3	72.80	90.93	70.67	95.47	60.67	77.47	94.93	96.67	96.93	63.20	78.53	92.80	97.73	96.93
4	58.10	85.71	67.14	86.19	38.57	62.86	91.43	93.33	95.24	40.00	62.86	82.38	96.67	97.14
5	92.39	93.74	90.60	96.42	89.49	90.38	89.49	95.75	98.43	89.04	90.60	93.29	94.85	98.21
6	96.88	97.32	98.07	98.66	95.24	97.17	98.51	99.55	99.11	95.98	96.88	98.81	98.96	99.11
7	43.48	69.57	17.39	82.61	21.74	21.74	91.30	60.87	100	21.74	21.74	82.61	17.39	100
8	98.86	98.41	98.86	97.95	97.05	98.18	99.55	100	100	99.09	98.64	99.77	100	99.97
9	50	55.56	16.67	50	33.33	55.56	0	0	88.89	61.11	55.56	0	0	100
10	71.53	93.80	74.97	93.80	68.20	77.61	89.44	94.60	98.05	70.72	79.33	91.27	94.37	97.70
11	84.38	94.37	84.87	95.54	75.96	85.68	97.34	99.28	97.43	77.94	86.90	97.43	98.33	98.20
12	85.51	93.66	81.16	91.85	54.53	77.90	88.22	95.65	98.73	61.23	78.44	89.13	97.46	98.73
13	100	99.47	100	100	100	100	100	100	100	100	100	99.47	100	100
14	93.30	99.14	95.02	96.56	92.87	95.70	99.14	99.83	99.40	95.62	95.96	99.05	99.91	99.48
15	64.91	87.43	61.70	88.01	41.23	55.85	99.12	91.81	97.95	48.25	55.56	97.95	97.08	97.37
16	88.24	100	57.65	88.24	94.12	92.94	96.47	91.76	97.65	92.94	94.12	92.94	94.12	95.29
OA	84.52	94.86	84.78	95.10	74.78	85.26	95.28	97.33	98.33	78.10	86.09	95.34	97.46	98.47
AA	79.24	90.73	73.05	89.55	68.61	78.02	88.45	88.39	97.81	72.52	78.50	87.12	86.03	98.31
κ	0.823	0.941	0.826	0.944	0.712	0.832	0.946	0.970	0.981	0.749	0.841	0.947	0.971	0.983

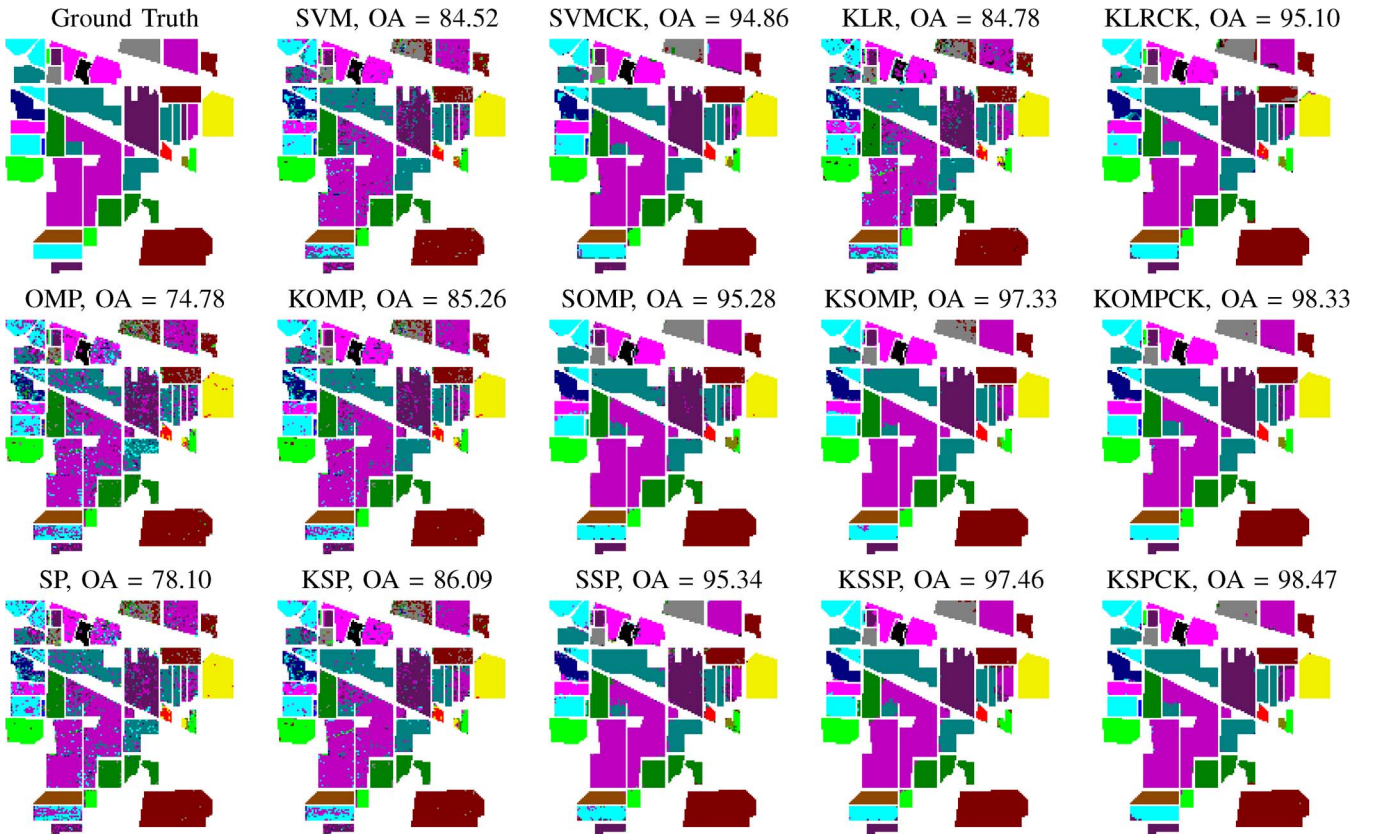


Fig. 2. Classification maps and overall classification accuracy (OA) for the Indian Pines image on labeled pixels with 10% training samples.

$\gamma = 512$ leads to the highest OA at all sparsity levels. For a fixed γ , the performance of KOMP and KSP generally improves as K_0 increases and tends to saturate as K_0 reaches 30–50. For KSOMP and KSSP, as shown in Fig. 3(c) and (d), the same tendency cannot be observed. However, the kernel JSM is more stable than the pixel-wise model, as for a large range of sparsity level K_0 and sufficiently large γ , the OA is always around 96% with a small variance. The stable performance suggests that we could also use empirical parameters K_0 and γ .

The next two HSIs used in our experiments, the University of Pavia and the Center of Pavia images, are urban images acquired by the Reflective Optics System Imaging Spectrometer (ROSIS). The ROSIS sensor generates 115 spectral bands ranging from 0.43 to 0.86 μm and has a spatial resolution of 1.3 m per pixel [42]. The University of Pavia image consists of 610×340 pixels, each having 103 bands, with the 12 most noisy bands removed. There are nine ground-truth classes of interests, as shown in Table III. For this image, we follow the

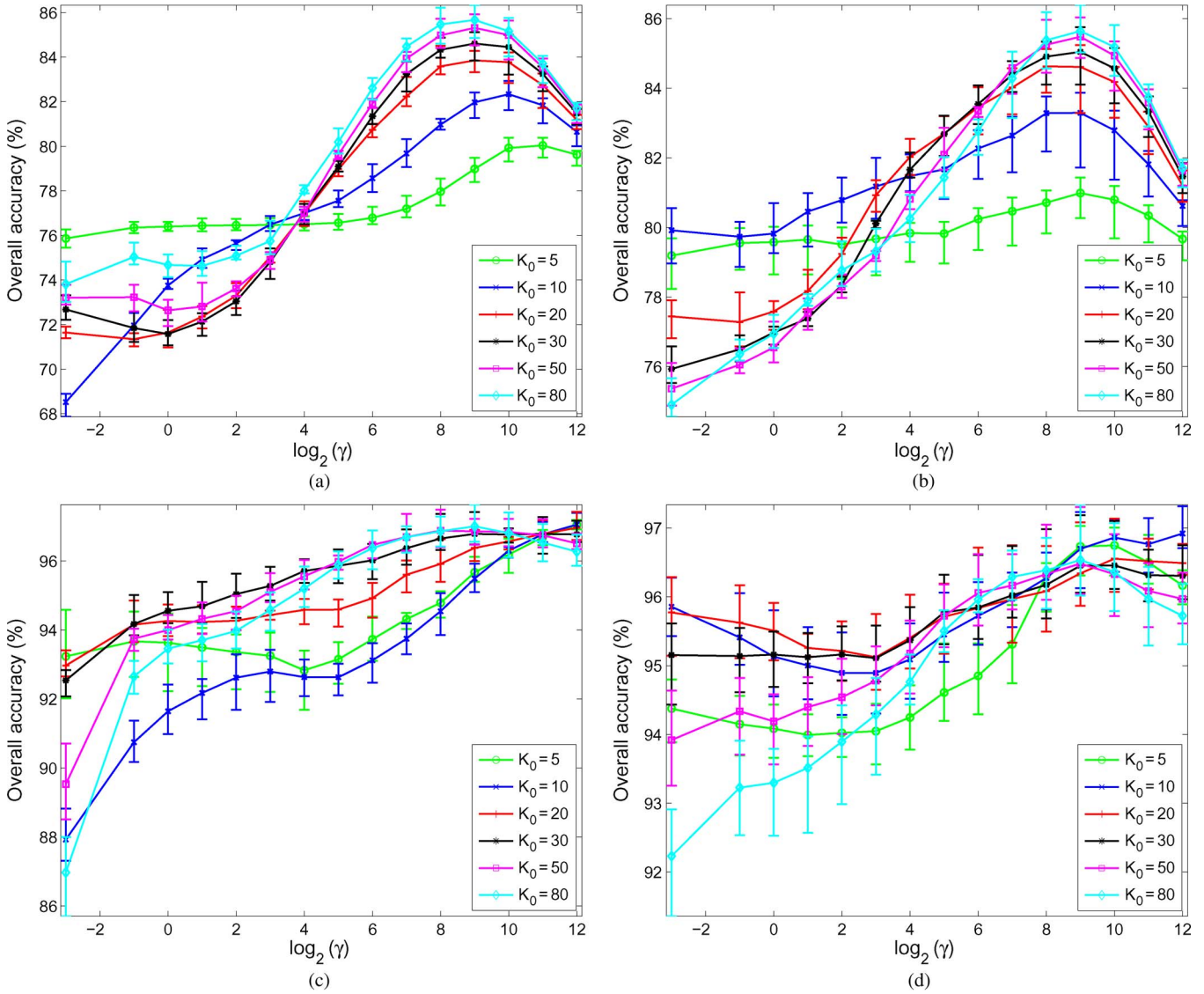


Fig. 3. Effect of sparsity level K_0 and RBF kernel parameter γ on the Indian Pines image using (a) KOMP, (b) KSP, (c) KSOMP, and (d) KSSP.

TABLE III
THE NINE GROUND-TRUTH CLASSES IN THE UNIVERSITY OF PAVIA IMAGE

No	Class Name	Samples	
		Train	Test
1	Asphalt	548	6304
2	Meadows	540	18146
3	Gravel	392	1815
4	Trees	524	2912
5	Metal sheets	265	1113
6	Bare soil	532	4572
7	Bitumen	375	981
8	Bricks	514	3364
9	Shadows	231	795
Total		3921	40002

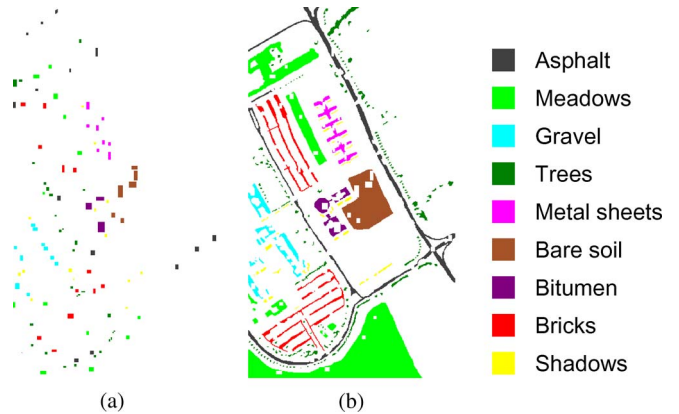


Fig. 4. (a) Training and (b) test sets for the University of Pavia image.

same experiment settings for the training and test sets as used in [30], [42], in which about 9% of labeled data are used as training, and the rest are used for testing, as shown in Table III and Fig. 4.

The classification accuracies and the κ coefficients on the test set using various techniques are shown in Table IV, and

the classification maps for all labeled pixels are presented in Fig. 5. Again, the RBF kernel is used for all kernel-based algorithms. This urban image lacks the large spatial homogeneity.

TABLE IV
CLASSIFICATION ACCURACY (%) FOR THE UNIVERSITY OF PAVIA IMAGE USING 3921
(AROUND 9%) TRAINING SAMPLES AS SHOWN IN FIG. 4 AND TABLE III

Class	SVM	SVMCK	KLR	KLRCK	OMP	KOMP	SOMP	KSOMP	KOMPCK	SP	KSP	SSP	KSSP	KSPCK
1	84.30	79.85	82.96	74.40	68.23	76.09	59.33	94.23	82.23	69.78	76.67	69.59	89.56	89.64
2	67.01	84.86	83.34	85.91	67.04	69.61	78.15	76.74	72.47	67.90	70.92	72.31	79.98	72.68
3	68.43	81.87	64.13	61.71	65.45	72.12	83.53	79.23	82.26	69.20	73.39	74.10	85.45	80.06
4	97.80	96.36	96.33	96.22	97.29	98.11	96.91	95.12	98.56	96.77	98.15	95.33	98.66	98.94
5	99.37	99.37	99.19	99.10	99.73	99.73	99.46	100	99.82	99.64	99.82	99.73	99.91	100
6	92.45	93.55	80.05	84.45	73.27	87.66	77.41	99.50	93.92	78.96	89.70	86.72	95.76	94.77
7	89.91	90.21	84.51	85.32	87.26	88.07	98.57	99.80	92.46	88.18	88.28	90.32	97.96	89.81
8	92.42	92.81	83.17	93.37	81.87	89.51	89.09	98.78	78.78	83.68	87.54	90.46	96.43	89.54
9	97.23	95.35	89.81	96.48	95.97	93.96	91.95	29.06	96.98	94.59	95.22	90.94	98.49	96.48
OA	79.15	87.18	83.56	84.77	73.30	78.33	79.00	85.67	81.07	74.86	79.18	78.39	87.65	83.19
AA	87.66	90.47	84.83	86.33	81.79	86.10	86.04	85.83	88.61	83.19	86.63	85.50	93.58	90.21
κ	0.737	0.833	0.784	0.799	0.661	0.725	0.728	0.815	0.758	0.681	0.735	0.724	0.840	0.785

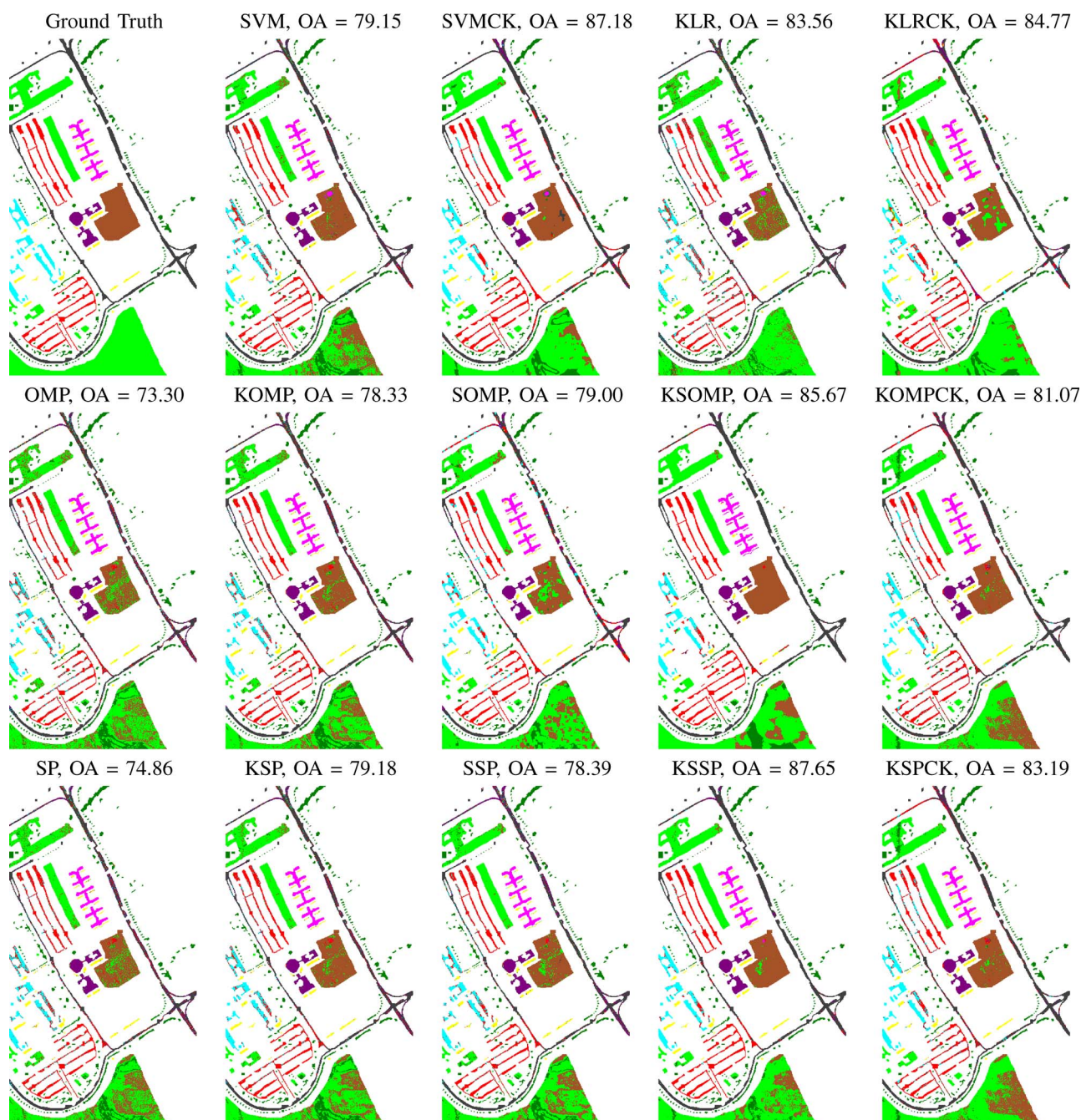
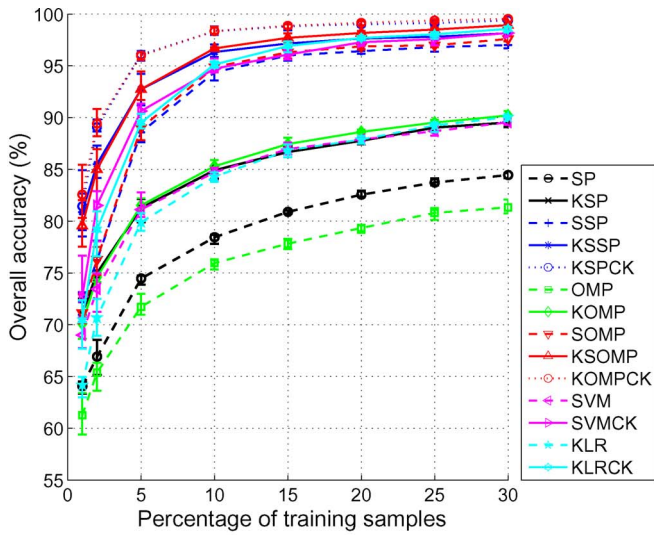
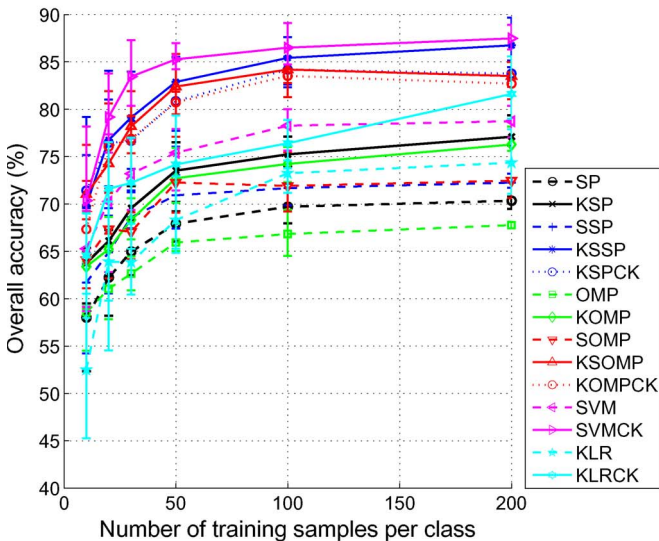


Fig. 5. Classification maps and overall classification accuracy (OA) for the University of Pavia image using around 9% labeled samples as training set.



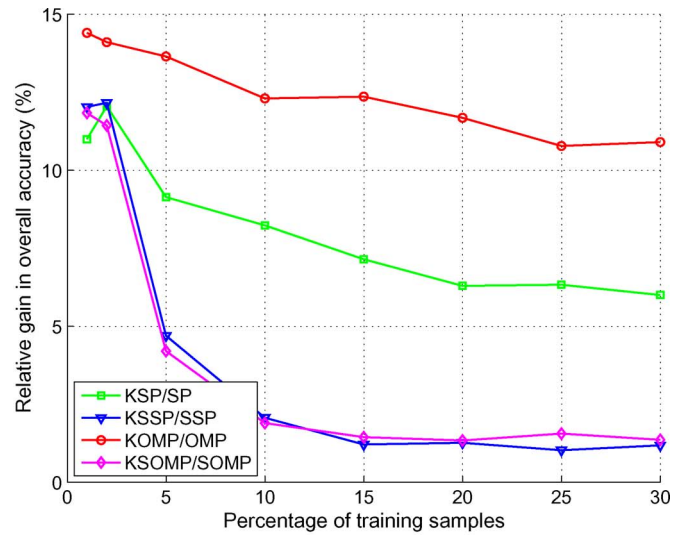
(a)



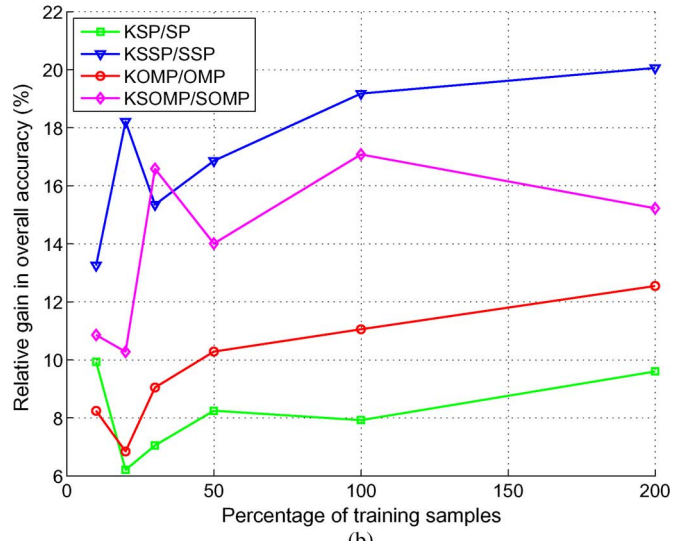
(b)

Fig. 6. Effect of dictionary size for (a) the Indian Pines image and (b) the University of Pavia image.

Therefore, a smaller neighborhood of size 5×5 is optimal for algorithms using a CK, and the linear and kernel JSMs. Similar to the Indian Pines image, the proposed KSOMP/KSSP algorithms achieve better or comparable performance when compared with the SVMCK classifier for most of the classes. KSOMP yields the best accuracy in five out of the total nine classes, and KSSP has the highest OA, AA, and κ coefficient. The overall performance of SVM, KOMP, and KSP, which are kernel methods for pixel-wise models, are comparable, and by incorporating the contextual information, the SVMCK, KSOMP, and KSSP techniques still have comparable performance. The sparsity-based algorithms generally do not handle the second class, representing Meadows, very well. For example, the accuracy for the second class for KSOMP and KSSP is 5%–9% lower than that for SVMCK and KLRCK, which affect the OA because this class contains more than 45% of the samples in the entire test set. This could be circumvented by selecting or learning a more representative training set which is sufficiently comprehensive to span the class subspace.



(a)



(b)

Fig. 7. Relative performance gain by kernelization for sparsity-based classifiers on (a) the Indian Pines image and (b) the University of Pavia image with different dictionary size.

In the sequel, we examine how the number of training samples affects the classification performance for various algorithms on the Indian Pines and the University of Pavia images. The algorithm parameters are fixed to be the same as those used to generate the results in Tables II and IV. For the Indian Pines image, in each test, we randomly choose 1% to 30% of the labeled data in each class as the training samples and the remaining samples as the test ones. The classification accuracy plots under various conditions are shown in Fig. 6(a) for the Indian Pines image, where the x -axis denotes the percentage of training samples from the total available labeled samples, and the y -axis is the OA on the test set. The accuracies are averaged over five runs for each classifier at each percentage level to avoid any bias induced by random sampling, and the bars indicate the maximal and minimal accuracies for each point in the five runs. The OA monotonically increases as the size of training set increases, and the variance is small (the difference between the maximum and minimum is within 1%) when at least 5%–10% training samples become available. The

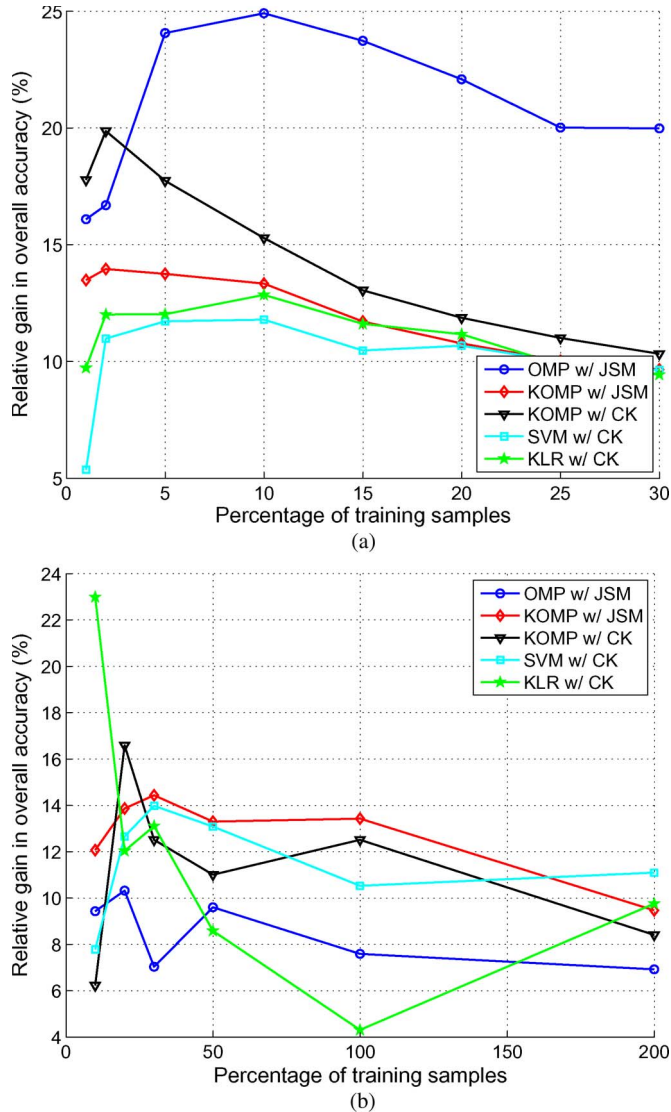


Fig. 8. Relative performance gain by contextualization for various classifiers on (a) the Indian Pines image and (b) the University of Pavia image with different dictionary size.

KOMPCK and KSPCK consistently yield higher OA than any other classifiers.

For the University of Pavia image, we create a balanced dictionary by randomly choosing $L = 10, 20, 30, 50, 100,$ and 200 training samples per class, and these training samples are a subset of the entire training set shown in Fig. 4(a). Since the dictionary is considerably small, the sparsity level K_0 is set to be no more than L . The classification accuracy plots are shown in Fig. 6(b), where the x -axis denotes the number of training samples per class, and the y -axis is the overall classification accuracy on the test set. Again, the accuracies are averaged over five runs for each classifier at each, L and the bars represent the maximum and minimum in the five runs. It is obvious that in most cases, the OA increases monotonically, and the variance decreases as the number of training samples increases. For the University of Pavia image, the performance at $L = 50$ is almost the same as that at $L = 100$ for all classifiers. The SVMCK classifier consistently outperforms all of the other classifiers when the number of training samples is small, but

TABLE V
THE NINE GROUND-TRUTH CLASSES IN THE CENTER OF PAVIA IMAGE AND THE TRAINING AND TEST SETS

Class		Samples	
No	Name	Train	Test
1	Water	745	64533
2	Trees	785	5722
3	Meadow	797	2094
4	Brick	485	1667
5	Soil	820	5729
6	Asphalt	678	6847
7	Bitumen	808	6479
8	Tile	223	2899
9	Shadow	195	1970
Total		5536	97940

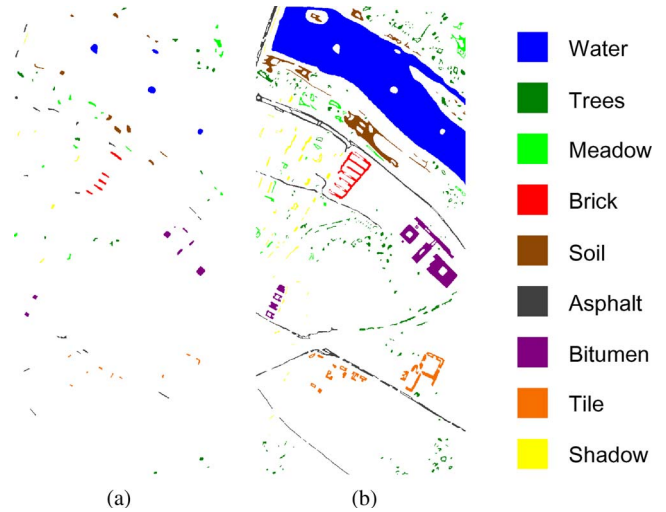


Fig. 9. (a) Training and (b) test sets for the Center of Pavia image.

the curves for SVMCK and KSSP tend to converge as more training samples are available (see Table IV for the performance comparison of SVMCK and KSSP with a large training set). It should also be pointed out again that during the training stage of algorithms using a CK, in order to extract the local spatial features for each training sample, one requires knowledge of the neighboring pixels or the location of the training sample, which may not be available in the training set. Moreover, the proposed sparsity-based algorithms rely on the approximation accuracy from each class subdictionary. Therefore, if the size of the subdictionary is too small, the training samples may not be sufficient to faithfully represent the subspace associated with each class, leading to a lower classification accuracy than the discriminative classifier SVM.

A closer inspection of the performance gain, as a function of the dictionary size, obtained by kernelization and contextualization is shown in Figs. 7 and 8, respectively. The y -axis represents the relative gain in percentage (averaged over five runs), which is the ratio between the improvement in accuracy and the OA of the algorithm before kernelization/contextualization. For example, in the case of contextualization of KOMP using the JSM, the relative gain is computed by

$$g = \frac{OA_{KSOMP} - OA_{KOMP}}{OA_{KOMP}} * 100\%$$

TABLE VI
CLASSIFICATION ACCURACY (%) FOR THE CENTER OF PAVIA IMAGE USING 5536 TRAINING SAMPLES
(AROUND 5% OF ALL LABELED SAMPLES AS SHOWN IN FIG. 9 AND TABLE V)

Class	SVM	SVMCK	KLR	KLRCK	OMP	KOMP	SOMP	KSOMP	KOMPCK	SP	KSP	SSP	KSSP	KSPCK
1	99.19	97.46	99.63	100	98.91	98.13	99.32	99.07	98.98	98.20	98.09	97.79	99.26	98.79
2	77.74	93.08	93.18	95.39	86.75	92.76	92.38	95.30	96.31	86.98	91.17	92.82	91.23	91.70
3	86.74	97.09	96.18	95.89	96.04	97.04	95.46	97.09	96.08	96.61	97.28	97.80	97.71	99.57
4	40.38	77.02	81.76	89.80	81.22	88.84	85.66	89.68	97.78	84.16	86.86	78.52	95.26	94.54
5	97.52	98.39	96.25	98.59	94.40	94.89	96.37	97.56	97.82	94.01	95.76	95.81	97.45	94.99
6	94.77	94.32	93.91	96.67	91.94	96.13	92.83	98.31	96.54	92.92	95.82	96.52	97.41	93.92
7	74.37	97.50	95.22	97.31	93.18	95.40	94.68	98.80	98.63	93.80	95.57	95.96	97.82	96.90
8	98.94	99.83	99.52	98.41	98.62	99.34	99.69	100	98.79	99.24	99.79	99.90	99.90	99.55
9	100	99.95	99.90	99.49	98.07	99.39	98.68	100	96.65	99.34	99.39	98.83	71.42	93.60
OA	94.63	96.81	97.99	98.92	96.68	97.19	97.66	98.53	98.46	96.40	97.08	96.93	97.82	97.55
AA	85.52	94.96	95.06	96.84	93.24	95.77	95.01	97.30	97.64	93.87	95.47	94.87	94.16	95.95
κ	0.899	0.943	0.963	0.980	0.940	0.949	0.958	0.973	0.972	0.935	0.947	0.945	0.960	0.956

where OA_{KSOMP} and OA_{KOMP} are the OA for the KSOMP and KOMP algorithms, respectively. The relative gain obtained by the kernelization of the SP, SSP, OMP, and SOMP algorithms is shown in Fig. 7(a) and (b) for the Indian Pines and University of Pavia images, respectively. One can observe that in most cases, kernelization consistently leads to a performance gain of 5% to 20%. The only exception exists in the KSSP and KSOMP algorithms for the Indian Pines image with a higher percentage of training samples, which is partly due to the fact that SSP and SOMP before kernelization already achieve an OA of at least 95%. In this case, an improvement of 2% to 3% means the error rate is reduced by half which could be considered significant.

The relative gain obtained by incorporation of spatial information in OMP, KOMP, SVM, and KLR is shown in Fig. 8(a) and (b) for the Indian Pines image and the University of Pavia image, respectively. The contextualization of SP and KSP has a similar effect to that of OMP and KOMP, and the results are not reported here. Note that with the kernel sparse representation model, the contextual correlation can be incorporated through either a JSM or a CK, and thus the relative gain of KSOMP (through JSM) and KOMPCK (through CK) over KOMP are both shown in Fig. 8. One can observe that for the India Pines image, the linear method OMP is the most sensitive to the spatial information, in which the relative gain is generally more than 20%. The other classifiers all work in the feature space, and the gain ranges from 10% to 15% in most cases, with a slight decrease as the number of training samples increases. For the University of Pavia image, the relative gain in classification accuracy is usually around 10% to 14%. Contrary to the case of the Indian Pines image, the improvement of the linear approach OMP is slightly less than the kernel methods. Moreover, the performance of KLR is not as consistent as the other methods.

The third image in our experiments, Center of Pavia, is the other urban image collected by the ROSIS sensor over the center of the Pavia city. This image consists of 1096×492 pixels, each having 102 spectral bands after 13 noisy bands are removed. The nine ground-truth classes and the number of training and test samples for each class are shown in Table V and illustrated in Fig. 9. For this image, about 5% of the labeled data are used as training samples. The classification results are summarized in Table VI, and the classification maps are shown in Fig. 10. KLRCK achieves a 100% accuracy on the

first class (water), which occupies 66% of the test set, and thus yields the best OA. The KSOMP and KSSP work very well on the other classes, except that KSSP fails for the ninth class (Shadow).

In general, one can observe from the experimental results on these three images that the incorporation of contextual information improves the classification performance (e.g., SP versus SSP, KSP versus KSSP, SVM versus SVMCK, etc). Moreover, operating in the kernel feature space also significantly improves the accuracy (e.g., SP versus KSP, SSP versus KSSP, etc).

V. CONCLUSION

In this paper, we propose a new HSI classification technique based on sparse representations in a nonlinear feature space induced by a kernel function. The spatial correlation between neighboring pixels is incorporated through a JSM. Experimental results on AVIRIS and ROSIS HSIs show that the kernelization of the sparsity-based algorithms improves the classification performance compared to the linear version. It is also shown that the proposed algorithm has a better or comparable performance to the recent spectral-spatial single-classifiers such as SVMCK.

The proposed sparsity-based classifier is different from the conventional sparse classifier SVM in many aspects. SVM is a discriminative model, which finds the separating hyperplane between two classes. A model with a fixed set of sparse support vectors is obtained by a training process over the whole training set, and then this SVM is used to classify all of the test data. However, our method can be considered as a generative model. The subspaces representing different classes implicitly compete with each other during the sparse recovery process, leading to a discriminative representation vector. This sparse representation vector is extracted for each test pixel and is thus adaptive. This will inevitably lead to an increase in the computational cost, but the kernel matrix K_A can be computed offline. Therefore, the most intensive part in the sparse recovery is the inversion of a matrix of at most size $K_0 \times K_0$ for the OMP-based algorithms and $(2K_0) \times (2K_0)$ for the SP-based algorithms. Moreover, in the OMP-based algorithms, since the support set is sequentially augmented by one index at a time, the inversion can be accelerated by Cholesky decomposition [45].

Our proposed dictionary-based classifier provides several advantages. First, new training samples can be easily added

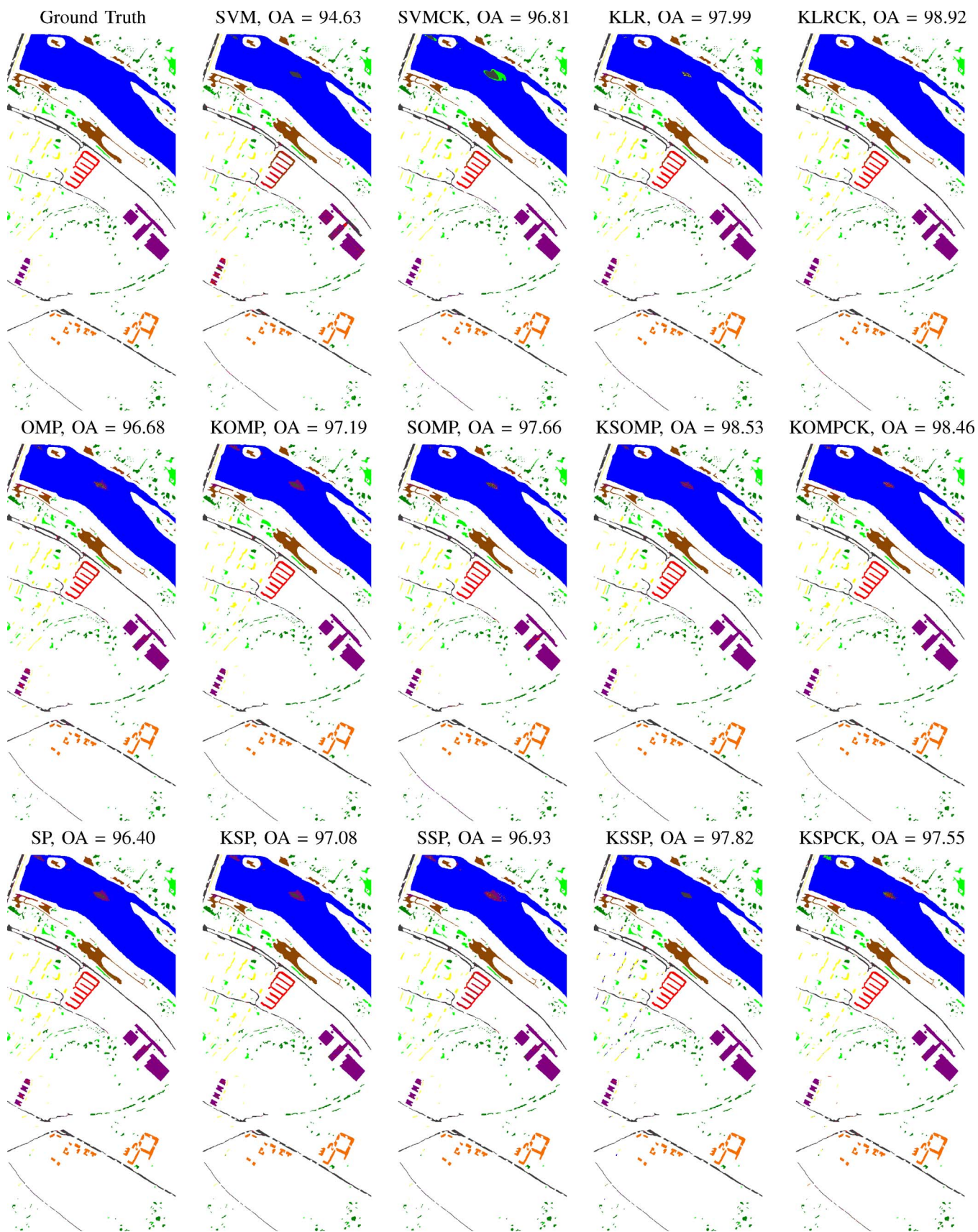


Fig. 10. Classification maps and overall classification accuracy (OA) for the Center of Pavia image using 5536 training samples (around 5% of all labeled samples) as shown in Fig. 10 and Table V.

to the dictionary without retraining the model, unlike the other classifiers (e.g., SVM and KLR) that need to retrain the model for the new training data. Also, our algorithm is particularly useful for creating a dictionary invariant to the environmental variations by adding synthetically generated spectral signatures that account for various illuminations and atmospheric conditions [46]. Moreover, the JSM in kernel space is still applicable when the training data is synthetically generated or from a spectra library rather than taken from the scene. On the other hand, in order to incorporate local spatial information, classifiers using composite-kernels require knowledge of the local spatial features of each training data which may not be available, and thus these classifiers may not be applicable in the case of unstructured synthetic or library training sets.

The classification accuracy can be further improved by a post-processing step, or combining the proposed technique with other state-of-the-art classifiers to generate a megaclassifier [17]. Another possible direction is the design/learning of a better dictionary such that the dictionary provides more accurate reconstruction, more discriminative power, and/or better adaptivity to the test data.

ACKNOWLEDGMENT

The authors would like to thank the University of Pavia and the HySenS project for kindly providing the ROSIS images of University of Pavia and Center of Pavia, and Prof. J. M. Bioucas-Dias for providing the sparse multinomial logistic regression code. Finally, we would like to thank all of the anonymous reviewers whose constructive criticisms have improved the paper significantly.

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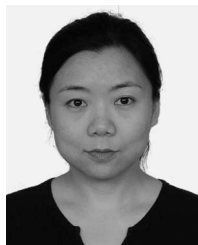
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