

# Dimension Reduction and Classification of Hyperspectral Images based on Neural Network Sensitivity Analysis and Multi-instance Learning

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**Abstract.** Hyperspectral remote image sensing is a rapidly developing integrated technology used widely in numerous areas. The rich spectral information from hyperspectral images aids in recognition and classification of many types of objects, but the high dimensionality of these images leads to information redundancy. In this paper, we used sensitivity analysis for dimension reduction. However, another challenge is that hyperspectral images identify objects as either a "different body with the same spectrum" or "same body with a different spectrum." Therefore, it is difficult to maintain the correct correspondence between ground objects and samples, which hinders classification of the images. This issue can be addressed using multi-instance learning for classification. In our proposed method, we combined neural network sensitivity analysis with a multi-instance learning algorithm based on a support vector machine to achieve dimension reduction and accurate classification for hyperspectral images. Experimental results demonstrated that our method provided strong overall classification effectiveness when compared with prior methods.

**Keywords:** Sensitivity Analysis, Artificial Neural Network, Ruck Sensitivity Analysis, Dimension Reduction, Classification, Hyperspectral Images, Multi-instance Learning, SVM

## 1. Introduction

Hyperspectral remote sensing technology is a rapidly developing field used in numerous areas of specialty, ranging from astronomy and geology to medicine and aerial surveillance, among others. Hyperspectral images provide rich spectral information that can aid in the recognition and classification of ground objects. However, the high dimensionality of these images leads to costly information redundancy, so the dimensions must be reduced. Sensitivity analysis using neural networks can be used for dimension reduction. However, another challenge remains. Hyperspectral images identify objects as either a "different body with the same spectrum" or "same body with a different spectrum." Therefore, it is difficult to maintain the correct correspondence

between ground objects and samples, which makes classification of the images difficult and computationally costly. This issue can be addressed using multi-instance learning for classification. In this research, we combine neural network sensitivity analysis with a multi-instance learning algorithm based on a support vector machine (SVM) to achieve dimension reduction for hyperspectral remote sensing images.

Hyperspectral imaging differs from general multispectral imaging insofar as hyperspectral imaging can display two-dimensional spatial information for a region of interest (e.g., the earth's surface), and it can add a dimension of spectral information. Multispectral images include only a few spectral bands, whereas hyperspectral images include hundreds of bands that are much narrower. Therefore, hyperspectral images can form an "image cube" [1]. In addition to multiple bands, hyperspectral images have distinct characteristics such as large amounts of data and high information redundancy that present difficulties for storing, transmitting, and processing the images. As a result, a band selection operation must be performed to reduce some of the unnecessary bands before processing hyperspectral images [2]. This operation helps to decrease the calculation requirements for hyperspectral image classification, and effectively avoids the Hughes phenomenon [3].

Hyperspectral image band selection can be considered an NP-hard combinatorial optimization problem [4]. Typically, search algorithms are used to search a band subset, allowing the evaluation standard to achieve its optimal value across all of the bands of the hyperspectral image. This queried band subset can then be treated as the optimal band combination. The common method of band selection and dimension reduction [5] for hyperspectral remote sensing images is to select several bands from the whole band to represent the whole band space. This approach requires that the band combination selected be able to provide effective improvement for classification accuracy in the subsequent classification. For this research, we select the band according to the band's contribution to classification. Bands that help to improve classification accuracy are selected first.

For classification of hyperspectral remote sensing images, neural network classifiers are commonly used because they work well for classifying images with high dimensionality and nonlinear structures. To provide a quantitative evaluation of the effects of one band on classification accuracy, a neural network sensitivity analysis can be employed that is based on a neural network classifier. Sensitivity analysis by a neural network [6] can provide a quantitative description of the influence of the input variables of a model on the output variables. The sensitivity coefficient of the model properties is sorted. Properties with larger sensitivity coefficients are chosen, and those with smaller ones are no longer included. In this way, the model is simplified, and the complexity of model processing is reduced. In this paper, we apply neural network sensitivity analysis to the band selection for hyperspectral remote sensing images, combined with a frequently used BP neural network classifier.

In addition to dimension reduction, the issue remains that hyperspectral images identify objects as either a "different body with the same spectrum" or "same body with different spectrum," which means they are unable to maintain the correct correspondence between target objects and samples. This challenge can be addressed using multi-instance learning for classification of hyperspectral images. In the mid- and late 1990s, T. G. Dietterich et al. proposed the concept of multi-instance learning in the study of drug activity prediction [7]. In this kind of learning, the training set consists of

several “bags” labeled with concept tags. Each bag contains several instances without concept notations. A bag is labeled positive if it contains at least one instance that is positive, and a bag is labeled negative if all of the instances in it are negative. By learning from the training bags, the learning system can predict the concept tag of a bag that is outside the training set as correctly as possible. Following on the work of T. G. Dietterich et al., many researchers began to devise practical multi-instance learning algorithms. Multi-instance learning stirred great interest in the machine learning field because it was a promising new learning framework in an area of machine learning previously unexplored. Multi-instance learning has unique properties and continues to offer good prospects for wide application.

A pixel-level classifier can help to divide the remotely sensed hyperspectral imagery, but if spectral background noise and clutter noise are present, classification accuracy will be decreased because of omissions or faults. Recently, some researchers proposed to adopt classification based on pattern spot images to solve the omissions or faults brought by spectrum changes. A pattern spot image refers to the single zone whose shape shares the same features with a spectrum. CHEN Jie et al. [8] put forward the rough set theory-based object-oriented classification of high resolution remotely sensed imagery. First, they abstracted the pattern spot image by means of watershed segmentation. Next, they analyzed the texture features of the abstracted images using Gabor wavelets, and divided the texture classification rules for abstracted images based on rough set theory. ZHANG Chuan et al. put forward object-oriented classification of high resolution remotely sensed imagery [9]. YANG Chang-bao et al. explored the object-oriented classification of remotely sensed imagery, and they determined the classification by segmenting the orthorectification SPOT image based on a distributed domain solver [10]. TAN Yu-min et al. put forward an object-oriented remote sensing image segmentation approach based on edge detection. [11]. Pattern spot image classification differs greatly from pixel-based classification because the former includes various regional texture space information, while the latter has only spectral features. Pixel-based classification considers only the features of a single pixel. Classification based on pattern spot images is less likely to be disturbed by spectral background noise and is more likely to retain regional integrity since it is abstracted by the regional spectra and space characteristics.

However, pattern spot image classification still has weaknesses in terms of its anti-noise properties that result in low classification accuracy with accompanying omissions or faults. Since the anti-noise property fails to provide the needed filtering, the classification will be less likely to be disturbed by noise with multi-instance learning spectra and space characteristics classified in the same zone, which is bigger than the object by making use of the regional relevance and object-oriented basis. DU Pei-jun et al. proposed that the cases in which different objects may have the same spectra characteristics or the same object may have different spectra characteristics, together with the noise in training instances, can be regarded as particular representatives of “ambiguity” in the training bag multi-instance learning. Therefore, when objects from the image segmentation are regarded as an instance in multi-instance learning, the object set of clutter is regarded as the bag, and multi-instance learning can be used in remote sensing classification [12].

The phenomena by which the same objects may have different spectra and different objects may have the same spectrum are the main source of land surface complexity for

remote sensing images. The complexity of land surface composition and the difficulty in selecting training samples cause the classification process to be highly dependent on human experience and prior knowledge. When using sensitivity analysis provided by an artificial neural network to realize dimension reduction for hyperspectral images, all of the bands are divided into several groups, as long as a lower correlation exists between adjacent bands. In addition, a differential evolution (DE) algorithm is used for optimizing the neural network structure. The bands that make only small contributions are given up based on the Ruck sensitivity analysis method.

Given the special advantages of multi-instance learning for solving ambiguous problems, and the advantages of neural network sensitivity analysis for dimension reduction, we suggest that integrating both multi-instance learning and sensitivity analysis for hyperspectral image classification can reduce the uncertainty of classification results. In view of this background and the benefits of applying new machine learning methods to remote sensing image classification, in this paper, we combine multiple-instance learning and an ensemble artificial neural network with embedded sensitivity analysis to improve the accuracy of hyperspectral image classification.

## 2. Related Work

### 2.1. Neural Network Sensitivity Analysis

Sensitivity analysis is an important research focus in the field of neural networks. In some practical applications, the availability of a huge amount of data can cause the trained neural network to become increasingly complex. The main task of sensitivity analysis is to determine how to analyze the parameters of the neural network effectively and simplify the scale of the network. Toward this end, we use the following procedure in this research.

Assume that the model is  $y = f(x_1, x_2, \dots, x_n)$ , where  $x_i$  is the  $i$ th property value of the model. It is necessary to ensure that any changes of each property are within the possible value range. The next step is to study and predict the influence of the change(s) of these properties on the output value of the model [13]. The degree of influence is called the sensitivity coefficient of property  $x_i$  on output value  $y$ : the higher the sensitivity coefficient, the greater the influence of the property on the output value. Thus, the sensitivity analysis can provide a quantitative description of the influence of the input variables on the output variable of a model. Furthermore, we can sort the sensitivity coefficients of the model's properties. We can choose the properties with larger sensitivities and give up the smaller ones according to the practical considerations of the problems. In this way, the model can be simplified and the computational complexity of processing the model reduced, which means dimension reduction is achieved.

Sensitivity analysis works based on the specific model. In most cases, however, when people face vast amounts of information, they are not clear about how the internal

mechanisms of the data work. In such cases, they cannot build the model expression  $y = f(x)$  directly, and therefore cannot conduct a sensitivity analysis either. However, researchers have shown that while a neural network does not need to model the physical concept of the research question, the neural network can provide more effective solutions for problems involving uncertainty or nonlinearity. The network provides a black box analysis model, and outputs reasonable results through the training and learning of input samples [14]. If the input data and output data are known, the neural network will use many simple neurons to simulate the non-linear relationships between the data. Neural network sensitivity analysis also uses the connection weights and the threshold between neurons to assess the influence of the input data on the output data [15].

Neural network sensitivity analysis can be divided into local sensitivity analysis and global sensitivity analysis. Some scholars have focused mainly on the study of local sensitivity analysis. There are four types of classical neural network sensitivity analysis. The first type is the sensitivity analysis method based on the connection weight, such as the Garson algorithm put forward in the early 1990s by Garson [16] and the Tchaban algorithm proposed by Tchaban [17]. The second type is sensitivity analysis based on the influence of the partial derivatives of output variables on input variables, such as Dimoponlos sensitivity analysis [18] and Ruck sensitivity analysis [19]. The third type is sensitivity analysis combined with statistics, such as methods based on random testing by Olden et al.[20]. The forth type is sensitivity analysis based on input variable disturbance, such as the method of adding white noise to the input data of a network and calculating the resulting change of output variables, an example of which was put forward by Scardi [21]

## 2.2. SVM Classification Methods and Multi-instance Learning

Training data are required to train the SVM model. However, these data cannot be separated without errors. The data points that are closest to the hyperplane are used to measure the margin, while the SVM attempts to identify the hyperplane that maximizes the margin and minimizes a quantitative proportion to the number of misclassification errors [22]. The SVM derives the optimal hyperplane as the solution of the following convex quadratic programming problem [23]:

$$\min_{w,b,\xi} \frac{1}{2} w^T w + C \sum_{i=1}^n \xi_i \quad s.t. \quad y_i (w^T x_i + b) \geq 1 - \xi_i, \quad \xi_i \geq 0, i = 1, 2, \dots, n, \quad (1)$$

where  $\{(x_1, y_1), \dots, (x_n, y_n)\}$  are the labeled training datasets with  $x_i \in R^d$  and  $y_i \in \{-1, 1\}$ ;  $w^*$  and  $b^*$  define a linear classifier in the feature space;  $C$  is the regularization parameter defined by the user; and  $\xi_i$  is a positive slack variable that handles permitted errors.

The optimization problem can be reformulated through a Lagrange function, where Lagrange multipliers can be found via dual optimization to generate a convex quadratic programming solution as follows [24] [25]:

$$\max Q(\alpha) = \sum_{i=1}^n \alpha_i - \frac{1}{2} \sum_{i=1}^n \sum_{j=1}^n \alpha_i \alpha_j x_i x_j K(x_i, x_j), \quad 0 \leq \alpha_i \leq C, i=1, 2, \dots, n \quad s.t. \quad \sum_{i=1}^n \alpha_i x_i = 0, \quad (2)$$

where  $\alpha = [\alpha_1, \alpha_2, \dots, \alpha_n]$  is the vector of the Lagrange multipliers, while  $K(\cdot, \cdot)$  is a kernel function. For a linearly non-separable case, a kernel function is introduced that satisfies the condition stated by Mercer's theorem and that corresponds to some types of inner product in the transformed (higher) dimensional feature space, as shown:

$$K(x_i, x_j) = \phi(x_i) \cdot \phi(x_j). \quad (3)$$

The final result is a discrimination function  $F(x)$  conveniently expressed as a function of the data in the original (lower) dimensional feature space:

$$F(x) = \text{sgn}[(w^*)^T \phi(x) + b^*] = \text{sgn} \left( \sum_{i=1}^n \alpha_i^* y_i K(x_i, x) + b^* \right). \quad (4)$$

Some popular kernel functions include the following:

a) Linear kernel

$$K(x, x_i) = (x \cdot x_i), \quad (5)$$

Polynomial kernel

$$K(x, x_i) = [(x^T x_i) + 1]^q, \quad (6)$$

where  $q$  is a constant.

b) Gaussian Radial Basis Function kernel

$$K(x, x_i) = \exp \left\{ -\frac{\|x - x_i\|^2}{\gamma^2} \right\}, \quad \gamma > 0, \quad (7)$$

Sigmoid kernel

$$k(x, x_i) = \tanh[v(x^T x_i) + c], \quad v > 0, c < 0. \quad (8)$$

Multi-instance learning was proposed as a new machine learning method to solve problems with relationships such as “1:N:1” using “object:description:label.” In multi-instance learning, labeled bags composed of several unlabeled instances are treated as training samples, and the goal of learning is to predict the labels of unknown new bags. In contrast to traditional supervised learning, the bag strategy employed by multi-instance learning offers special advantages in dealing with ambiguous problems. Multi-instance learning is viewed as the fourth machine learning framework, in parallel with reinforcement learning, supervised learning, and unsupervised learning (Bolton & Gader, 2011; Lozano-Perez, 1998; Bolton, et al., 2011; Zhang, et al., 2004; Zhou, et al., 2002; Andrews, et al., 2002; Zhang, et al., 2010). At the moment research on multi-instance learning has focused mainly on creating new multi-instance learning algorithms and on designing applications of multi-instance learning for various fields (Zhang, et al., 2004; Zhou, et al., 2002; Li, et al., 2004). In the image processing field, multi-instance learning has been used for image retrieval and scene classification, with good results obtained in some existing experiments (Li, et al., 2007; Li, et al., 2010; Wang, et al., 2010; Li, et al., 2008) [26].

Traditional supervised learning can be treated as a special case of multi-instance learning. The transformation of traditional supervised learning algorithms to make them capable of dealing with multiple instance problems is an important branch in multi-instance learning algorithm research. Considering the bag concept, multi-instance learning can be viewed as a generalization of traditional supervised learning. Combining multiple learners for the purpose of enhancing the performance of the base learner is an effective method in traditional supervised learning frameworks. According to the supervised nature and classification function of multi-instance learning, multiple instance ensemble learning is also a feasible approach, and some researchers have shown that performance is increased by using an ensemble. Research into the integration of ensemble learning and multi-instance learning is an active branch of machine learning, so advances have been seen in remote sensing image classification (Qi, et al., 2011; Zhou, et al., 2003; Auer & Ortner, 2004; Kittler, et al., 1998).

Multi-instance learning algorithms based on an SVM can be divided into two categories: multi-instance learning based on samples (mi-SVM), and multi-instance learning based on bags (MI-SVM) [27–28]. mi-SVM tries to identify a maximal margin hyperplane for the instances, subject to the constraint that at least one instance of each positive bag is located in the positive half-space while all instances of negative bags are in the negative half-space. MI-SVM tries to identify a maximal margin hyperplane for the bags by regarding the margin of the “most positive instance” in a bag as the margin of that bag.

$$\begin{aligned}
 \text{mi-SVM} : \min_{\{y_i\}} \min_{\{w\}} \frac{1}{2} \|w\|^2 + C \sum_i \xi_i, \quad \text{s.t. } \forall i : y_i (\langle w, x_i \rangle + b) \geq 1 - \xi_i, \xi_i \geq 0, y_i \in \{-1, 1\} \\
 \sum_{i \in I} \frac{y_i + 1}{2} \geq 1, \forall I \quad \text{s.t. } Y_I = 1, \text{ and } y_i = -1, \quad \forall I \quad \text{s.t. } Y_I = -1 \quad . \quad (9) \\
 \text{MI-SVM} : \min_{\{w, b, \xi\}} \frac{1}{2} \|w\|^2 + C \sum_I \xi_I, \quad \text{s.t. } \forall I : Y_I \max_{\{i \in I\}} (\langle w, x_i \rangle + b) \geq 1 - \xi_I, \xi_I \geq 0 \quad , \quad (10)
 \end{aligned}$$

where  $w$  and  $b$  are two parameters;  $\xi_i$  is a positive slack variable;  $x_i$  is the input value;  $y_i \in \{-1, 1\}$  is the output value; and  $C$  is the regularization parameter defined by the user.

### 3. Application of Neural Network Sensitivity Analysis and Multi-instance Learning to Band Selection

The back propagation (BP) neural network classifier is a common tool for the classification of hyperspectral remote sensing images. Combined with the neural network sensitivity analysis method introduced above, the band selection for BP can be carried out in the whole band space. The band combination that provides a large contribution to the classification is selected (as explained in Section 3.1 below), thereby realizing the purpose of dimension reduction. To get better analysis results, before conducting a sensitivity analysis the hyperspectral remote sensing image data should be pretreated. Some band combinations with weak correlation are preselected by subspace-partition. To avoid blindness in the selection of the initial weights and threshold of BP neural network [29], a differential evolution algorithm is used to optimize the BP neural

network. Last, the optimized BP neural network is used to conduct the sensitivity analysis. The sensitivity analysis results for all of the test samples are combined by using the comprehensive evaluation function, and finally the band with the biggest influence on classification results is selected. The specific process is explained in the following subsections of this paper.

### 3.1. Data Preprocessing and Band Selection

For the proposed method, data processing takes place using the following steps. Before employing the neural network sensitivity analysis, the original hyperspectral remote sensing image needs to be preprocessed to eliminate the bands that have interference from noise, water vapor, or other serious pollution. Select the object that has the largest number of samples as the pre-selected object that is good for classification. Typically, the original hyperspectral remote sensing image has a large number of bands, but there is a high correlation and high redundancy between bands. Therefore, to attain a good result from the sensitivity analysis, it is very important to choose as the input the band that also has a weak correlation.

To solve the above problem, the approach is to divide the whole band into several subspaces, and then select the band. The adaptive subspace decomposition (ASD) [30] method based on correlation filtering is used to divide the band set of the hyperspectral remote sensing image. First, we calculate the correlation coefficient denoted as  $R_{ij}$  between the two bands. Let  $u_i$  and  $u_j$  denote the number of  $i$  and  $j$  bands, respectively. As the value of the correlation coefficient grows larger, the correlation between bands becomes stronger, and as the correlation comes closer to 0, the correlation becomes weaker.  $R_{ij}$  is defined as:

$$R_{i,j} = \frac{E\{(x_i - \mu_i)(x_j - \mu_j)\}}{\sqrt{E\{(x_i - \mu_i)^2\}} \sqrt{E\{(x_j - \mu_j)^2\}}} \quad (11)$$

The value  $R_{ij}$  of the matrix  $R$  ranges between 0 and 1. As  $R_{ij}$  comes closer to 1, the correlation between the two bands becomes stronger.  $\mu_i$  and  $\mu_j$  are the mean values of  $x_i$  and  $x_j$ , respectively. They denote the gray mean values of the two bands.  $E[\bullet]$  is the mathematical value expected. When all  $R_{ij}$  values are identified, then the proper threshold  $T_b$  is set. The continuous bands of  $|R_{ij}| \geq T_b$  form a new subspace. We can control the number of subspaces and the number of bands in each subspace dynamically by changing the threshold  $T_b$ . Furthermore,  $R_s$  denotes the ratio of the number of bands in a subspace to the total number of bands in all subspaces. We select the bands in each subspace to create band combinations according to the ratio  $R_s$ . Then we reduce the correlation of the bands as much as possible. Combining the bands with the pre-object types and the object information of the original remote sensing image, we can determine the training sample  $P$  for input, expected value  $T$  for output, test sample  $P\_Test$  for

input, and expected value T\_Test for output, all of which are required for the training of a BP neural network. These steps make it convenient to determine the topological structure of the neural network.

### 3.2. Dimension Reduction and Classification based on Sensitivity Analysis and Multi-instance Learning

#### 3.2.1 Ruck sensitivity analysis

Ruck sensitivity analysis (1990) is based on the partial derivatives of output variables for input variables. This method is designed for feedback neural networks, such as BP neural networks and RBF neural networks. This approach, which is simple and fast, evaluates the partial derivatives with the activation function of the neural network, and calculates the influence of the input data on the output data. Therefore, Ruck sensitivity analysis is used to study band selection for a hyperspectral remote sensing image based on a BP neural network classifier.

The value  $R_{ij}$  of the matrix  $R$  ranges between 0 and 1. As  $R_{ij}$  comes closer to 1, the correlation between the two bands becomes stronger.  $\mu_i$  and  $\mu_j$  are the mean values of  $x_i$  and  $x_j$ , respectively. They denote the gray mean values of the two bands.  $E[\bullet]$  is the mathematical value expected. When all  $R_{ij}$  values are identified, then the proper threshold  $T_b$  is set. The continuous bands of  $|R_{ij}| \geq T_b$  form a new subspace. We can control the number of subspaces and the number of bands in each subspace dynamically by changing the threshold  $T_b$ . Furthermore,  $R_s$  denotes the ratio of the number of bands in a subspace to the total number of bands in all subspaces. We select the bands in each subspace to create band combinations according to the ratio  $R_s$ . Then we reduce the correlation of the bands as much as possible. Combining the bands with the pre-object types and the object information of the original remote sensing image, we can determine the training sample P for input, expected value T for output, test sample P\_Test for input, and expected value T\_Test for output, all of which are required for the training of a BP neural network. These steps make it convenient to determine the topological structure of the neural network.

The above Ruck sensitivity analysis is used only for the sensitivity analysis of the input value to the output value of a single sample test point. A comprehensive evaluation function is needed to synthesize the sensitivity analysis results of each single sample point. Here, we apply the MSA metrics proposed by Jacek.  $M$  [25] is chosen as the synthetic evaluation function. Assume  $S_{ik}$  denotes the sensitivity coefficient of the input variable  $i$  of all the samples to the output variable  $Y_k(k=1)$ , and  $s'_{ik}$  denotes the sensitivity coefficient of input variable  $i$  to the output variable  $Y_k(k=1)$  in the number  $t$  sample. The synthetic evaluation function can be expressed as (15):

$$S_{ik} = \sqrt{\frac{\sum_{t=1}^n (s_{ik}^t)^2}{n}}, \quad (15)$$

where  $n$  denotes the total number of samples, and  $S_{ik}$  is nonnegative.  $S_{ik}$  can be used to sort the sensitivity of the input bands, and then the influence of the input variable to the output results can be measured.

### 3.2.2 Band Classification Using Neural Network Sensitivity

First, according to the proportion  $R_s$  in each subspace, several bands are selected to form a band combination. Combined with the pre-selected object type, the training sample  $P$  and the test sample  $P\_Test$  are generated, which then serve as the input variables of the BP neural network. The number of band combinations is equal to the number of the neurons at the input end. Furthermore, the label value of the ground object type is used as the output of the BP neural network. Then, the training samples (with results  $T$ ) and test samples (with results  $T\_Test$ ) are generated. In addition, the weight and threshold of the BP neural network are optimized by a DE algorithm. All of the samples are classified by the optimized BP neural network, and the results of the sensitivity analysis are calculated. In addition, bands are sorted according to their sensitivity coefficients. Bands with smaller sensitivity coefficients will be given up, and bands with important effects on classification results will be identified. Last, the screened band combinations are classified using the neural network classifier to verify the effect of dimension reduction. The flow of the procedure is shown in Figure 1.

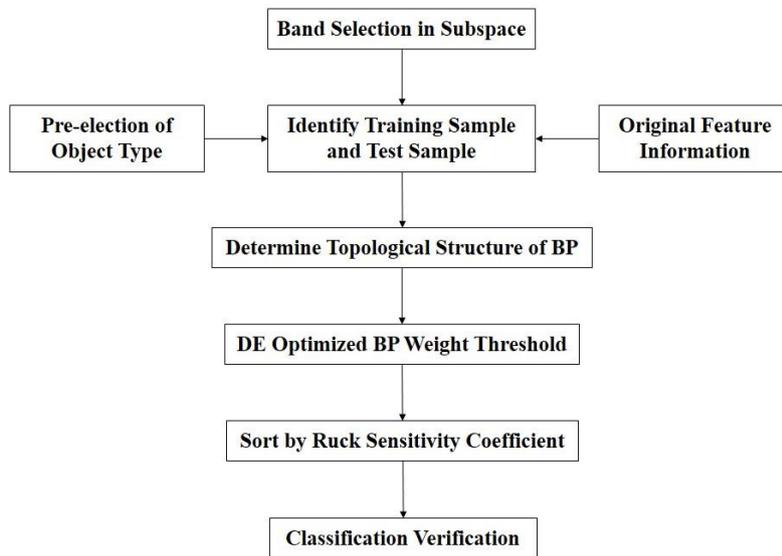
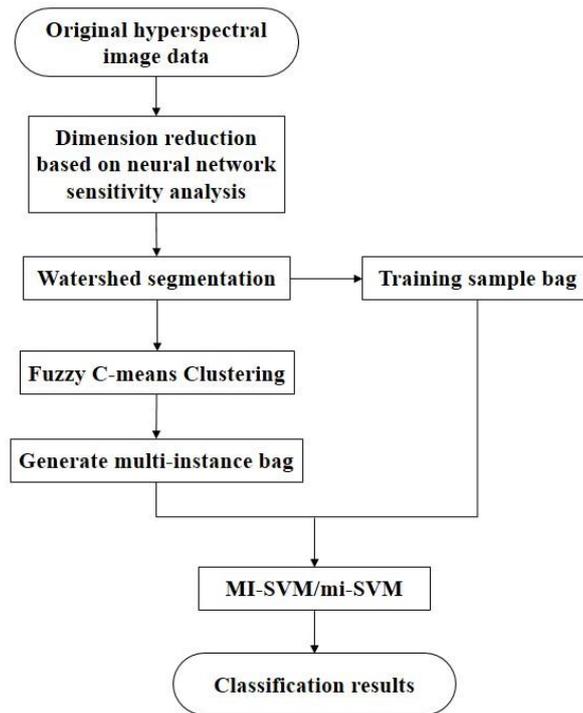


Fig. 1. The realization flow for band selection by sensitivity

### 3.2.3 Classification of Hyperspectral Remote Sensing Image based on Multi-instance Learning

As described above, in multi-instance learning [31–32], the training set is composed of labeled bags, and the bags are made up of non-labeled samples. The goal of multi-instance learning is to predict the label of previously unseen bags by learning the training set. Classification by bags as a unit includes many more features than classification by pixels or by the object as a unit, and both the anti-noise capability of the region and the accuracy of classification are improved. Each feature of the segmented object is taken as an instance; the object set generated by the clustering is taken as the packet. Using the packets, multi-instance learning based on an SVM is used for classification. In this paper, first, based on the dimension reduction of the band, the watershed transform algorithm is utilized to decompose the image into several objects that are used as instances for multi-instance learning. Then, the training samples are constructed with instances selected by artificial selection, and the unknown bags are obtained by the clustering algorithm. The new bag is marked with the nearest sample category that has maximum diversity. In other words, the number of generated bags is completely determined by the clustering algorithm. The steps of the algorithm are as follows.



**Fig. 2.** The realization flow of classification of hyperspectral remote sensing images based on multi-instance learning

- (1) Reduce the dimension of the hyperspectral band by using neural network sensitivity analysis.
- (2) Obtain the object by applying the common watershed algorithm.
- (3) For each band that is selected by the band selection method, calculate the property of the object, and construct the property space.
- (4) Multi-instance bags are generated by the fuzzy C-means clustering
- (5) Select samples for each kind of object to form a training sample bag, and use an SVM for classification.
- (6) Get the classification results.

The flow of the classification process is shown Figure 2.

## 4. Experiments and Results

### 4.1. Design of the Experiment

We designed some simulation experiments to demonstrate empirically the effectiveness of our proposed method for dimension reduction of hyperspectral remote sensing images using neural network sensitivity analysis and multi-instance learning. The experimental program was developed using MATLAB R2009b. The SVM classifier utilized the LIBSVM toolkit (<http://www.csie.ntu.edu.tw/~cjlin/libsvm/>) and adopted a radial basis function (RBF) to perform the experiments. The penalty factor of the SVM was 16. The BP neural network was realized by the built-in neural network toolbox of MATLAB. We designed the experiment using a standard hyperspectral image. The dataset was revised in MAT format, which can be obtained from the website [http://www.ehu.es/ccwintco/index.php?title=Hyperspectral\\_Remote\\_Sensing\\_Scens](http://www.ehu.es/ccwintco/index.php?title=Hyperspectral_Remote_Sensing_Scens).

The image was part of a hyperspectral image from a mixed agroforestry experimental zone in northwestern Indiana (USA). It was acquired by the AVIRIS sensors in June 1992. In this image, the range of wavelengths was 0.4~2.5  $\mu\text{m}$ , the size of image was 145\*145 pixels, and the spatial resolution of the image was 25m. In the pretreatment of the original image [24] (including image denoising and image deblurring), the bands that were heavily polluted by water vapor and noise (such as bands 1~4, 78, 80~86, 103~110, 149~165, and 217~224) were removed from the original bands. The remaining 179 bands were kept for experimental purposes. Figure 3 is an RGB false color image composite of the selected bands: 50, 27, and 17. Figure 4 is the real distribution image of the 7 objects chosen for the experiments.

There were 16 classes of ground objects in the image. Seven of these classes had larger sample sizes and were chosen for the classification experiments. The ratio between training samples and testing samples was selected evenly as 1:1. Table 1 displays the number of ground objects, as well as the names, and quantities of training and testing samples.



Fig. 3. The AVIRIS false color image

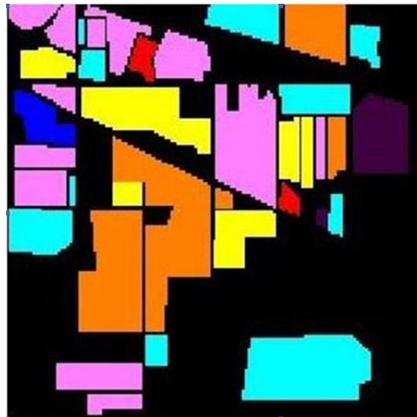


Fig. 4. The real distribution image

Table1. Training samples and test samples

Class number	Training samples	Training bags
Class 1	717	15
Class 2	417	9
Class 3	374	8
Class 4	484	10
Class 5	1234	25
Class 6	307	6
Class 7	647	13
Total	4180	86

## 4.2. Results and Analysis

A DE algorithm was used to optimize the BP neural network to get better results from the sensitivity analysis, which meant the band combination could improve the accuracy of classification effectively after the band with the smallest sensitivity coefficient was eliminated. Furthermore, to some extent, the multi-instance learning method based on the SVM could improve classification accuracy. We designed six experiments to demonstrate the superiority of the multi-instance learning method based on an SVM and optimizing the BP neural network with the DE algorithm. The band combinations used in the experiments were the same ones selected under the same subspace division. The detailed content of the experiments is shown in Table 2.

**Table 2.** Comparison experiments

Group	Content of Experiments
A	BP neural network classification (BP)
B	Sensitivity analysis and BP neural network classification (SA-BP)
C	Sensitivity analysis and Genetic algorithm optimizing BP neural network classification (SA-GABP)
D	Sensitivity analysis and Differential Evolution optimizing BP neural network classification (SA-DEBP)
E	Sensitivity analysis and MI-SVM (SA-MI-SVM)
F	Sensitivity analysis and mi-SVM (SA-mi-SVM)

In each experiment, six types of bands with different numbers were compared, and the values of  $R_s$  were selected relatively as 1/9, 1/6, 2/9, which meant selecting the band whose number was about 20, 30, or 40 from the divided subspace in accordance with the proportion of  $R_s$ . The topology structure of the BP neural network was set as the  $N$  number of neurons of the input layer equal to the number of bands in each group. The number of the types of primary features was 7, which was also the total number of the classification. The number of the neurons of output layer  $M$  was set as 7. The hidden layer was set to be single, and the number of its neurons  $L$  was set according to the expression  $L = \sqrt{N + M} + a$ . In this expression,  $a$  is the adjustment constant between 1 and 10.  $L$  changed with the training sample set, and ultimately it was determined that the error of the network was minimal when  $a = 5$ . The related parameters of the BP neural network training and DE algorithm were set as shown in Table 3 and Table 4.

**Table 3.** The setting of the parameters of BP

BP Parameters	Parameters setting
Frequency of training	1000
Minimum mean square error	0.01
Learning rate	0.1
Hidden layer activation function	Hyperbolic tangent function tansig

Output layer activation function	Linear function purelin
Training function	Levenberg-Marquadt Back propagation algorithm

**Table 4.** The setting of the parameters of DE

Parameters	Parameter values
Individual dimension D	$D = N*L+L*L*M+M$
Population size Nd	Nd = 20
Population size MAXGEN	MAXGEN = 50
Hybridization parameters CR	CR = 0.9
Differential evolution model	DE/best/1/bin

Except for group A, the other five groups of experiments needed to calculate the sensitivity coefficient. Considering the large number of sensitivity coefficients, Tables 5–7 show only the sensitivity coefficients of three kinds of bands. In the experiments with group D, these bands were analyzed using the Ruck method. In the table, the sensitivity coefficients are arranged in order from largest to smallest.

**Table 5.** Sensitivity coefficient of 20 Bands

No.	Band No.	Sensitivity coefficient	No.	Band No.	Sensitivity coefficient
1	117	1.0674	11	185	0.5610
2	43	0.9070	12	199	0.5477
3	140	0.8505	13	173	0.5102
4	57	0.7437	14	93	0.5099
5	5	0.7364	15	72	0.4385
6	102	0.6796	16	134	0.4076
7	68	0.6647	17	24	0.3984
8	212	0.6485	18	87	0.3313
9	37	0.6098	19	126	0.2332
10	15	0.5881	20	204	0.1486

**Table 6.** Sensitivity coefficient of 30 Bands

No.	Band No.	Sensitivity coefficient	No.	Band No.	Sensitivity coefficient
1	16	1.2224	16	141	0.3852
2	114	0.8933	17	73	0.3778
3	178	0.8743	18	37	0.3076
4	34	0.8259	19	87	0.3013
5	53	0.7306	20	117	0.2566

6	133	0.7176	21	148	0.2474
7	97	0.7028	22	191	0.1856
8	100	0.6639	23	200	0.1725
9	89	0.6369	24	170	0.1615
10	211	0.6300	25	8	0.1520
11	39	0.6259	26	42	0.1117
12	125	0.5951	27	11	0.0909
13	27	0.5242	28	57	0.0867
14	196	0.4252	29	68	0.0814
15	184	0.3990	30	166	0.0596

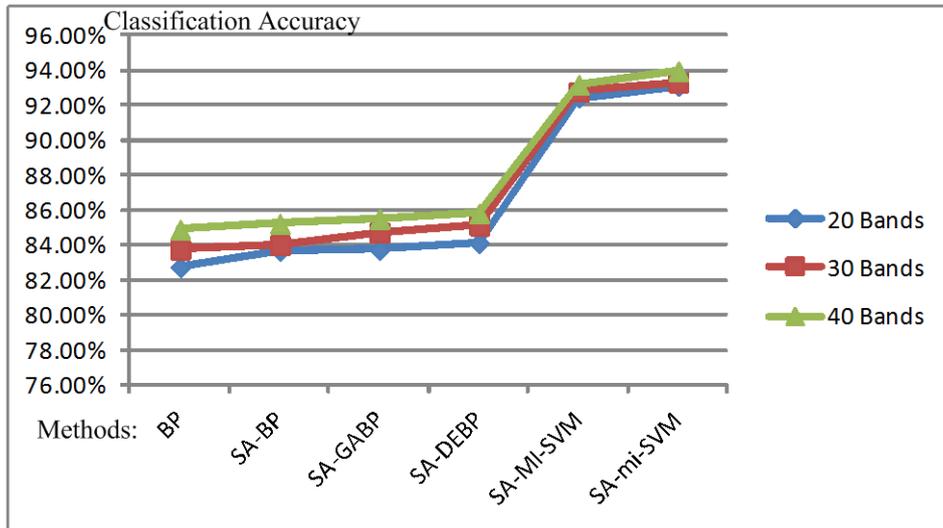
**Table 7.** Sensitivity coefficient of 40 Bands

No.	Band No.	Sensitivity coefficient	N o.	Band No.	Sensitivity coefficient
1	16	1.3678	21	114	0.4461
2	68	1.1337	22	141	0.4298
3	23	1.1020	23	170	0.3795
4	120	0.9650	24	125	0.3689
5	196	0.9390	25	39	0.3623
6	133	0.8981	26	180	0.3424
7	27	0.8731	27	49	0.3112
8	89	0.8263	28	73	0.3083
9	166	0.8159	29	211	0.2794
10	11	0.7527	30	8	0.2504
11	117	0.7440	31	87	0.2373
12	138	0.7327	32	191	0.2334
13	214	0.6805	33	200	0.2145
14	37	0.6289	34	42	0.1798
15	148	0.6065	35	178	0.1409
16	34	0.5631	36	57	0.1275
17	102	0.5511	37	184	0.1172
18	100	0.5422	38	78	0.0988
19	147	0.5087	39	97	0.0827
20	53	0.4860	40	14	0.0408

From Tables 5–7, we can see that the sensitivity coefficient of the final band was small enough, about the 1/20 of the first one. For several experiments regarding classification, first each band whose sensitivity coefficient was less than 0.3 was excluded. Then, combined with multi-instance learning, the band combinations after dimension reduction were classified again by the BP neural network, checking whether the dimension reduction with sensitivity analysis was effective in improving the accuracy of classification. Table 8 shows the classification accuracy of band combinations after dimension reduction in the experiments of groups A, B, C, D, E, F.

**Table 8.** Classification accuracy

Number of Bands	A	B	C	D	E	F
20	82.7	83.68%	83.7	84.13	92.4	93.0
	5%		5%	%	3%	9%
30	83.7	84.02%	84.7	85.14	92.7	93.3
	5%		3%	%	8%	2%
40	84.9	85.24%	85.5	85.83	93.1	93.9
	0%		2%	%	8%	6%

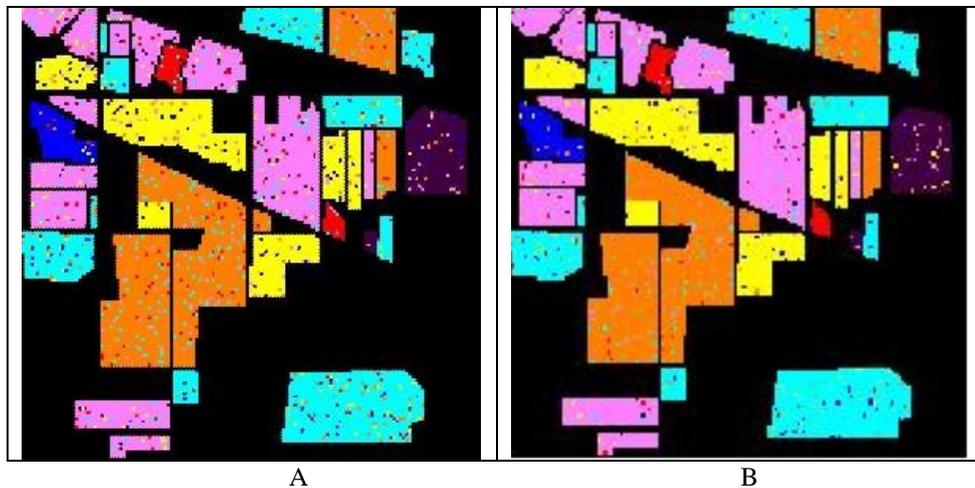


**Fig. 5.** Classification accuracy curves

From Table 8 and Figure 5, we note that with the same number of bands, the BP neural networks optimized with a genetic algorithm (GA) and DE algorithm in the experiments with groups C and D achieved better classification accuracy than groups A and B. This finding means that the sensitivity coefficients attained in the experiments with groups C and D provided more accurate reflection of the effects of each band on the classification results. Moreover, the BP neural network optimization with the DE algorithm in group D was better than for the GA algorithm adopted in group C. In the experiments with group B, the BP neural network was analyzed with the Ruck method after the division of the subspace without optimization, so the classification accuracy was lower than for groups C and D regardless of the number of bands. The BP neural network in experiments with group A was used in classification without dimension reduction, so its classification accuracy was the lowest. These results show that dimension reduction with sensitivity analysis was effective.

For classification, groups E and F used sensitivity analysis combined with multi-instance learning based on samples (mi-SVM) and multi-instance learning based on bags (MI-SVM), respectively. With the same number of bands, the classification accuracy of groups E and F was significantly higher than for groups B, C, and D, at more than 90%. These results demonstrate that in terms of classification, multi-instance learning performed better than the BP neural network. In addition, using multi-instance learning not only improved the classification accuracy, but also achieved an ideal classification effect. The classification accuracy using 30 bands was higher than when using 20 bands, and the classification accuracy using 40 bands was higher than when using 30 bands. This result indicates that with an increase in the number of bands, the classification accuracy of the same group improved. With the increased number of bands, the classification accuracy of groups E and F was more than 90%, which means that sensitivity analysis and multi-instance learning are very suitable for the classification of hyperspectral image. Figures 6–8 show the results of six groups of experiments using 20, 30, and 40 bands. It can be seen directly from these figures that groups A and B had the fewest incorrect classifications and higher classification accuracy regardless of the number of bands.

Finally, we compared the classification performance of the proposed method with some other hyperspectral image classification methods, as shown in Table 9. In this table, the accuracies outside the brackets were taken from corresponding references directly, and those accuracies in the brackets were obtained by our proposed method. All of the methods were tested under the same conditions with 7 classes and 5% training samples. Table 9 demonstrates that the classification performance of the proposed method outperformed all of the compared methods.



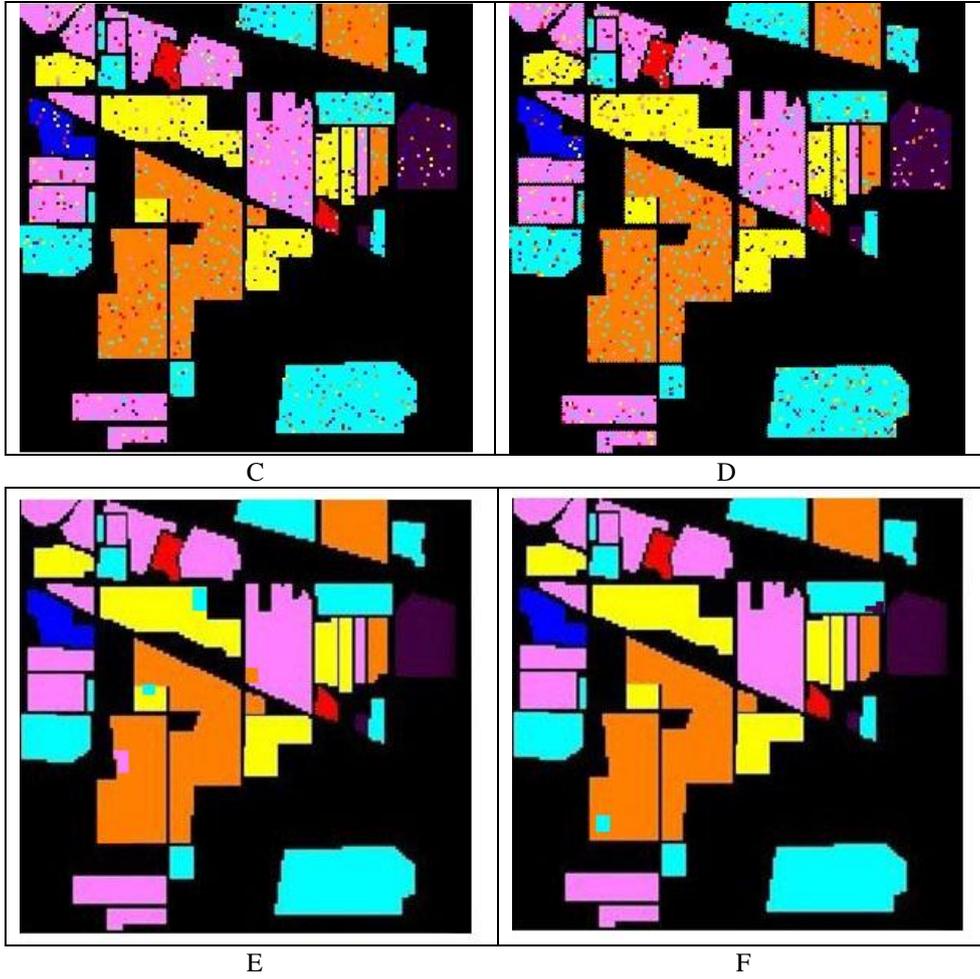


Fig. 6. Classification results of 20 bands

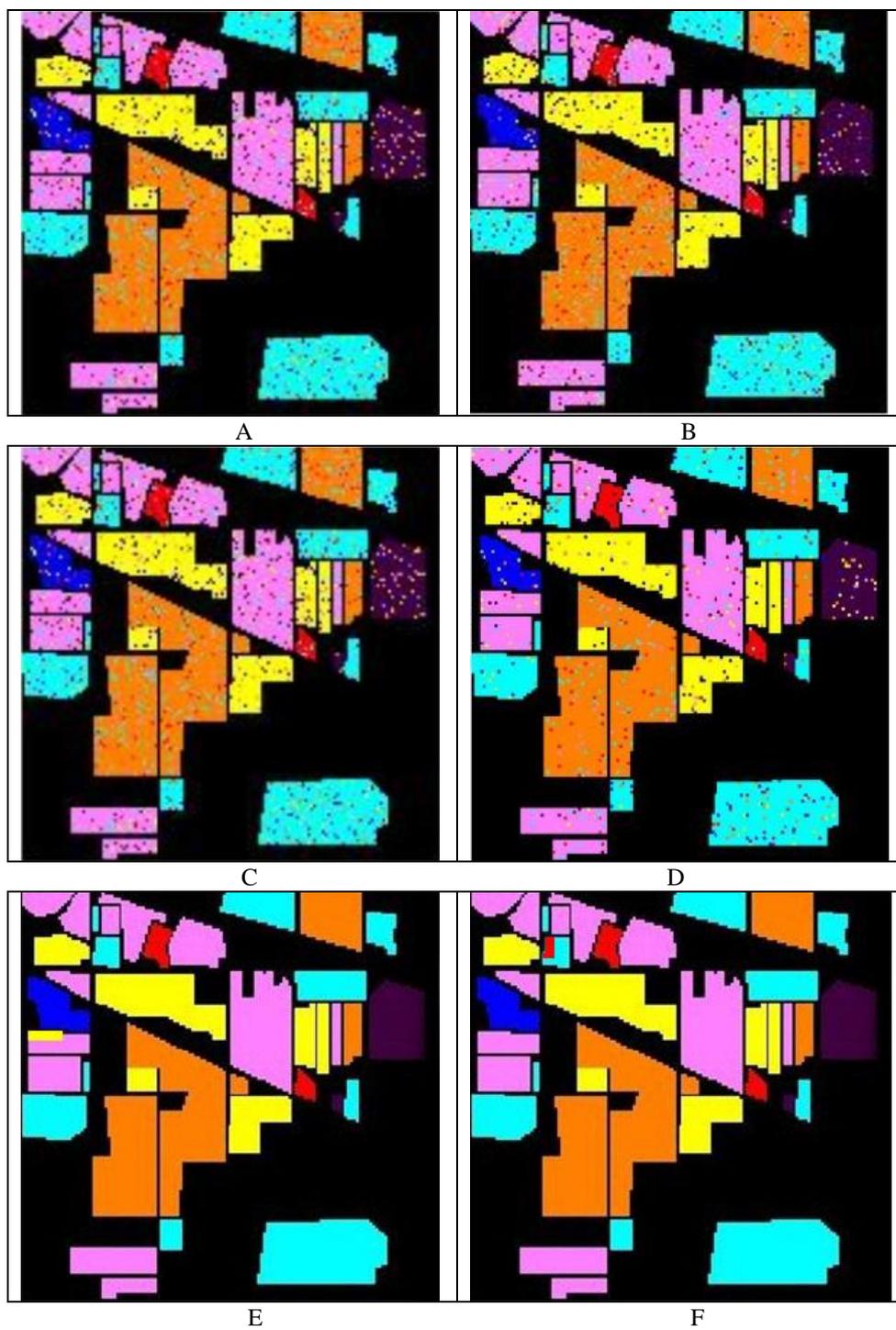


Fig. 7. Classification results of 30 bands

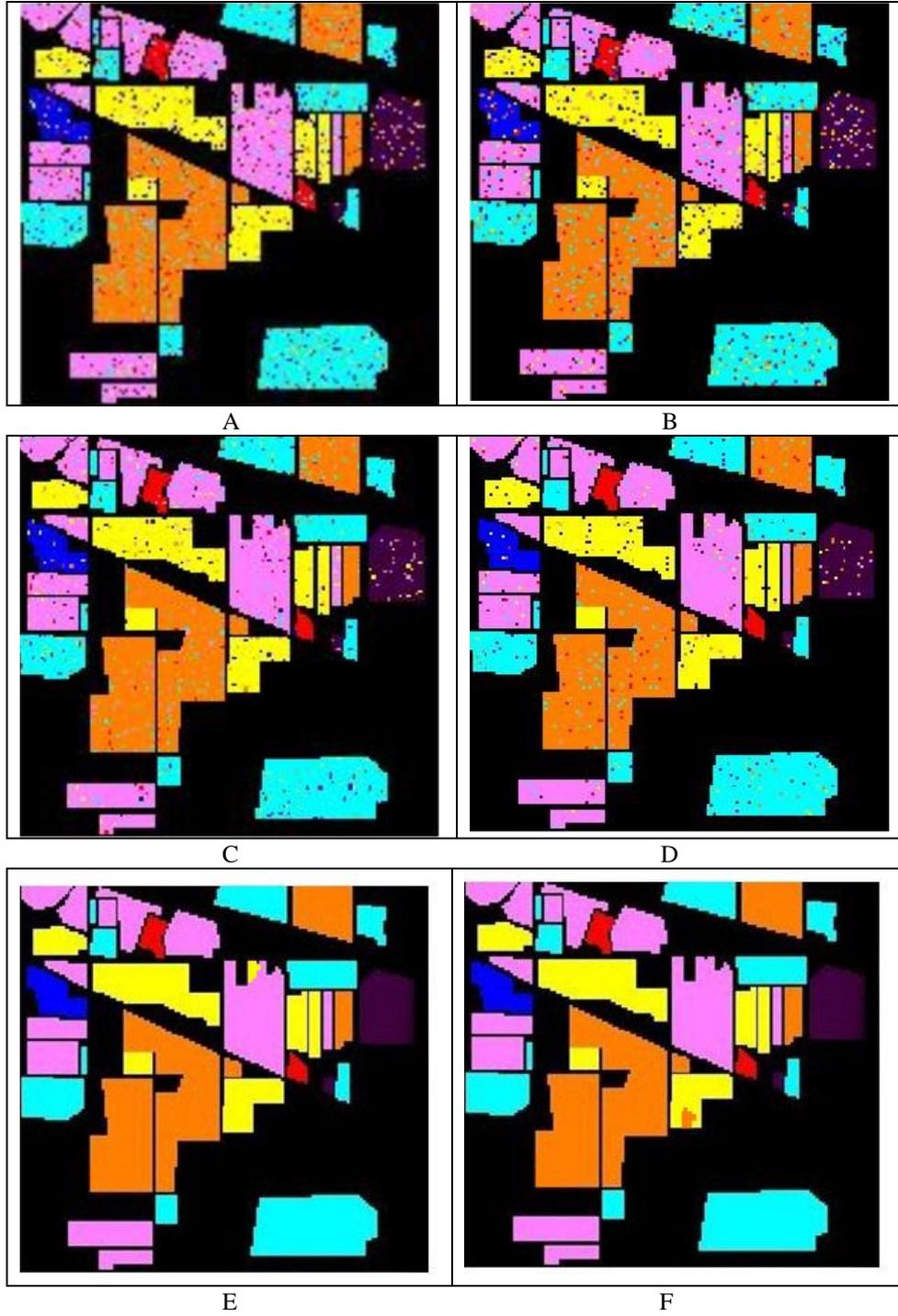


Fig. 8. Classification results of 40 bands

**Table 9.** Comparison with other methods

References	Baseline	Feature	Overall Accuracy
Sun et al. [33]	RWASL	Spectral	94.15%
Guo et al. [34]	FSAF	Spectral	84.36%
Li et al. [35]	KELM	Spectral	84.53%
Hu et al. [36]	RBF-SVM	Spectral	90.74%
Proposed method	SA-MI-SVM	Spectral	94.21%
Proposed method	SA-mi-SVM	Spectral	94.28%

## 5. Conclusions

Hyperspectral images take objects as the basic classification unit in a manner similar to the examples and packages used in multi-instance learning. Therefore, multi-instance learning can be used for the classification of remote sensing images. In this research, we combined neural network sensitivity analysis with a multi-instance learning algorithm based on an SVM to achieve dimension reduction and classification of hyperspectral remote sensing images. First, to reduce the correlation among the input properties, we used adaptive subspace division to select band combinations. In addition, a DE algorithm was adopted to optimize the BP neural network. To provide stable network connection weights and thresholds for sensitivity analysis by the neural network, we employed a sensitivity analysis method based on a partial derivative to calculate the sensitivity coefficient and remove bands that had very small coefficients. In this way, we achieved dimension reduction. To decrease the impact of the "different body with same spectrum" or "same body with different spectrum" phenomena on classification of hyperspectral images, a watershed transform algorithm was employed to decompose the image into several objects that were used as instances of multi-instance learning. The object set generated by clustering formed a bag, and the SVM was used for classification.

Our experimental results demonstrated that our method provided strong overall classification effectiveness when compared to prior methods. Based this study, we can draw the following conclusions.

1. Classification accuracy can be improved by means of the proposed neural network sensitivity analysis method because the contributions of bands for subsequent classification are sorted, and the bands with the largest contributions are selected.

2. A multi-instance learning method can learn robustly from highly noised training samples, but the bag label prediction rule for multi-instance learning also brings uncertainty to the classification results. The uncertainty problem can be controlled through the introduction of neural network sensitivity analysis.

3. The multi-instance learning algorithm based on an SVM can obtain higher classification accuracy under strong noise training conditions as long as the positive examples in the samples are selected properly.

4. For small sample training, the classification accuracy is higher based on the SVM and multi-instance learning algorithm.

Multi-instance learning has special advantages for resolving ambiguous problems compared with traditional supervised learning methods. Integrating multi-instance learning with neural network sensitivity analysis proved to have useful effects for controlling uncertainty. However, as the experiments showed, ensemble multi-instance learning needs a reasonable feature embedding scale factor. The main focus of future work should be on methods for implementing dynamically or self-adaptively chosen optimal scale factors.

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