

An improved semi-supervised prototype network for few-shot fault diagnosis

Zhenlian Lu¹, Kuosheng Jiang², Jie Wu³

^{1,2}School of Mechanical Engineering, Anhui University of Science and Technology, Anhui, 232001, China

^{1,2}Anhui Key Laboratory of Mine Intelligent Equipment and Technology, Anhui University of Science and Technology, Huainan, 232001, China

³School of Mechanical Engineering, Anyang Institute of Technology, Henan, 455000, China

³Corresponding author

E-mail: ¹2021200542@aust.edu.cn, ²jiangkuosheng333@stu.xjtu.edu.cn, ³jiewu06@163.com

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Abstract. The collection of labeled data for transient mechanical faults is limited in practical engineering scenarios. However, the completeness of sample determines quality for feature information, which is extracted by deep learning network. Therefore, to obtain more effective information with limited data, this paper proposes an improved semi-supervised prototype network (ISSPN) that can be used for fault diagnosis. Firstly, a meta-learning strategy is used to divide the sample data. Then, a standard Euclidean distance metric is used to improve the SSPN, which maps the samples to the feature space and generates prototypes. Furthermore, the original prototypes are refined with the help of unlabeled data to produce better prototypes. Finally, the classifier clusters the various faults. The effectiveness of the proposed method is verified through experiments. The experimental results show that the proposed method can do a better job of classifying different faults.

Keywords: intelligent diagnosis, semi-supervised learning, meta-learning, standard Euclidean distance, few-shot.

1. Introduction

The complete sample of monitoring data is the foundation of fault diagnosis for mechanical equipment [1], [2]. However, the collected data is usually limited and constitutes few-shot instances for mechanical condition monitoring, thus making it difficult to identify fault information. Therefore, intelligent diagnosis methods are studied to address this problem [3-6]. Specially, deep learning (DL) shows excellent performance in solving small sample problems [7-10].

The fault diagnosis model based on deep learning could effectively identify fault features from few-shot data, therefore solving the problem of fault diagnosis under small sample condition. Saufi et al. [11] proposed a deep learning model based on stacked sparse autoencoders (SSAE) to tackle the small sample size problem. The experimental results show that the diagnosis model has high diagnosis performance. Ren et al. [12] designed a capsule autoencoder model (CaAE) for intelligent fault diagnosis. The state capsule has undergone feature fusion, which makes it easier to achieve better diagnosis accuracy with smaller training samples. In addition, with the help of relevant datasets, feature adaptation based on transfer learning can also learn fault features from few-shot datasets to achieve accurate fault diagnosis. Xie and Duan used transfer component analysis (TCA) to extract transferable fault features from gearbox vibration signals. These two models can effectively extract and fuse cross-domain features [13], [14]. However, it is not ideal in terms of the actual accuracy for transfer learning method due to the complex high-dimensional data and negative transfer.

Meta-learning can more effectively complete the classification of faults by virtue of its unique training methods and strong generalization ability, addressing the problem of insufficient samples. Yu et al. [15] designed a meta-learning model based on time-frequency features and multi-label

convolutional neural networks. It can quickly identify new fault types by a few samples and a simple network update step. Wang et al. [16] proposed an improved metric meta-learning model. The approach makes full use of the attribute information of a single sample and the similarity information between sample groups. Chen et al. [17] have constructed an improved prototype network model with frequency domain data as input. The network uses Mahalanobis distance instead of Euclidean distance for prototype classification. Although the above meta-learning methods have achieved considerable results, they mainly focus on labeled samples and ignore the information existing in unlabeled samples.

In practical engineering, most of the collected data is unlabeled, leading to increased costs and time for labeling each sample individually. To address this issue, scholars have conducted further studies on semi-supervised learning due to its ability to generate stronger feature representations for each class [18-20]. Yu et al. [21] proposed an effective semi-supervised learning method based on the principle of consistent regularization, which could enhance the clustering between features through supervised and unsupervised loss terms. But this adds a large number of additional enhancement operations. Feng et al. [22] proposed a semi-supervised local Fisher discriminant network by adopting a pseudo-labeling strategy. It has an excellent performance for reducing the number of enhancement operations. Nevertheless, this network is more suitable for processing high-dimensional image data. This paper incorporates a semi-supervised learning strategy into the prototype network framework. The prototypes of all classes are refined multiple times by predicting the labels of unlabeled data to enhance the features of the individual prototypes [23-26].

In this paper, an improved semi-supervised prototype network (ISSPN) is designed to accomplish small sample fault classification with limited data. In this method, ProtoNet-based metric meta-learning network is used to extract fault features efficiently with limited data. Then, a semi-supervised learning strategy is employed, where unlabeled data is used to improve the quality of the prototypes for better fault clustering. Finally, the use of standard Euclidean distance instead of Euclidean distance in the prototype distance metric can eliminate the effect of differences in different feature scales. Meanwhile, the standardization of the data also reduces the impact of outliers on the distance calculation and improves the robustness of the model to outliers.

The rest of the paper is organized as follows: Section 2 outlines the relevant theory. Section 3 describes the proposed method and model in detail. Sections 4 and 5 are devoted to experimental analysis and conclusions, respectively.

2. Basic theory

2.1. Prototypical networks

Prototypical Networks (ProtoNet) is a typical metric-based meta-learning model. This network maps the data into different classes by the feature function, and then learns the similarity between each sample by the distance metric [27], as illustrated in Fig. 1. Specifically, since ProtoNet operates on a meta-learning framework, it adheres to a scenario training strategy. Firstly, the dataset is divided into multiple tasks using the “N-way K-shot” approach. This means that each task comprises N classes, with each class containing K samples. Moreover, each task consists of a support set S (training set) and a query set Q (test set). The task is divided into two parts during the training process: the meta-training set and the meta-test set. Simultaneously, the types of data in the two are different [28].

Next, the training process can be more accurately described as that the prototype network mainly learns a nonlinear feature mapping function $f_e(\cdot)$. It is parameterized as a neural network to map the prototype into a feature space. In this space, prototypes from the same class are closer, while prototypes from different classes are far away from each other. All parameters of prototype networks depend on the feature mapping function. To calculate the prototype P_d of each class d , the average value of the mapped feature vector represents the prototype. The calculation is as

follows:

$$P_d = \frac{1}{|\sum S_i|} \sum_{(x_i^s, y_i^s) \in |\sum S_i|} f_e(x_i^s), \quad (1)$$

where P_d is the prototype of class d , $|\sum S_i|$ represents the total size of samples in class d , x_i^s is the sample belonging to class d and y_i^s is the label corresponding to x_i^s .

Finally, the parameter e of the feature function is trained and updated by the query set samples under each meta-task. The classification of the sample is done by measuring the distance between the sample and all prototypes. The prediction probability of the sample belonging to the query set of class d is calculated as follows:

$$P(y = d | x_i \in Q) = \frac{\exp[-sE_{dist}\{f_e(x_i), P_d\}]}{\sum_{d'} \exp[-sE_{dist}\{f_e(x_i), P_{d'}\}]}, \quad d' \in \{1, 2, \dots, N\}. \quad (2)$$

For the prototype of all query sets, the loss function is defined to update the average negative logarithm of all query set samples in a given training set:

$$L_\theta(Q) = -\frac{1}{T_q} \sum_{(x_i, y_i) \in Q} \log p(y = y_i | x_i), \quad (3)$$

where T_q is the total number of samples in the query set Q . The training process updates the model parameters by minimizing L_θ .

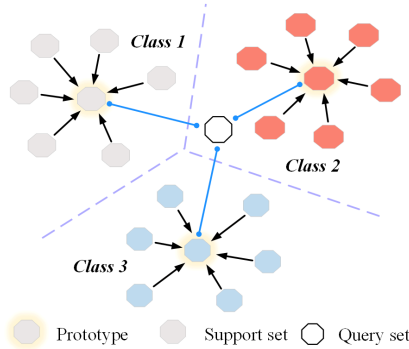


Fig. 1. Schematic representation of a metric-based meta-learning model

2.2. Prototype purification

Due to the small amount of sample data and the limited number of labeled samples, the initial prototype cannot correctly represent the cluster centroid. To get a better prototype, the unlabeled data is used to adjust the initial prototype. It can effectively solve the problem of position deviation caused by the limited labeled support data. The prototype can be guided slightly by the unlabeled samples, thereby better aggregating the information into the initial prototype. The prototype purification process is shown in Fig. 2.

First, the unlabeled data is assigned probability labels. And then, the probabilities of unlabeled data are aggregated into the existing prototypes to obtain more information. The initial prototype $P_d \in R^Z$ is constructed for each class with the labeled support set as shown in Eq. (1). Therefore, the probability of labeling the sample for each class can be given as follows:

$$P_{i,d} = p(y = d | x_i \in S_l) = \begin{cases} 1, & y = d, \\ 0, & y \neq d. \end{cases} \quad (4)$$

The S_l is a dataset that supports centralized labeling. The unlabeled data probability is shown in Eq. (5):

$$\tilde{P}_{i,d} = p(y = d | x_i \in S_u) = \frac{\exp[-sE_{dist}\{f_e(x_i), P_d\}]}{\sum_{d'} \exp[-sE_{dist}\{f_e(x_i), P_{d'}\}]} \quad (5)$$

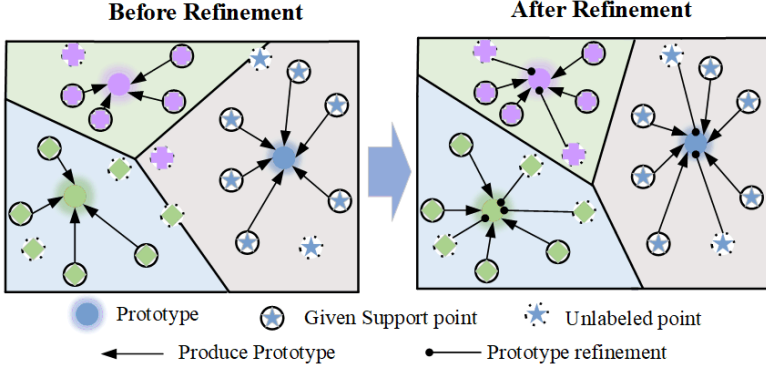


Fig. 2. Schematic diagram of prototype purification

The standard Euclidean distance function sE_{dist} is defined as:

$$sE_{dist}(o_1, o_2) = \|o_1 - o_2\| = \sqrt{\sum_{j=1}^Z \left(\frac{o_{1,j} - o_{2,j}}{s_j} \right)^2}, \quad (6)$$

where $o \in R^Z$; s_j is the standard deviation corresponding to the prototype. For the refined prototype, two probability weight coefficients are defined as:

$$q_{i,d} = \frac{P_{i,d}}{\sum_i P_{i,d} + \sum_j \tilde{P}_{j,d}}, \quad (7)$$

$$\tilde{q}_{i,d} = \frac{\tilde{P}_{i,d}}{\sum_i P_{i,d} + \sum_j \tilde{P}_{j,d}}. \quad (8)$$

Whereupon the original prototype is assigned by weight coefficient corresponding to each category:

$$\tilde{p}_d = \sum_{i=1}^{s_l} q_{i,d} f_e(x_i) + \sum_{i=s_l+1}^{s_u+s_l} \tilde{q}_{i,d} f_e(x_i). \quad (9)$$

The s_l and s_u are the number of labeled and unlabeled samples in the support set, respectively.

At last, the process of refining the prototype is completed by repeatedly executing the Eq. (9) and continuously assigning the value of \tilde{p}_d to P_d .

3. The proposed method

In this section, the improved semi-supervised prototype network (ISSPN) is described in three parts. The approach is based on an extension of metric meta-learning. In the first part, the learning style of the method is introduced. Next, the process of diagnosis is described in detail. Finally, the

optimizer is explained.

3.1. Learning style

The proposed method follows the way of semi-supervised few-shot learning in this paper. Foremost, the support set S is divided into two parts: labeled sample S_l and unlabeled sample S_u :

$$S = S_l \cup S_u = \{(x_i, y_i)\}_{i=1}^{s_l} \cup \{x_i\}_{i=s_l+1}^{s_l+s_u}. \quad (10)$$

After that, the labeled samples S_l are then randomly selected for “N-way K-shot” classification. The labeled samples after classification are input into the model to obtain the initial prototypes of n categories. Moreover, s_u samples in the unlabeled support set S_u are input into the convolution operation blocks (COBs) module for mapping into the feature space. The unlabeled samples are used to adjust the position of initial prototype through multiple iterations until full convergence. In addition, part of the remaining samples are extracted from the N classes as a query set Q , and then classify the samples in the query set q_n by prototype purification. The prediction formula for each sample in the query set is given by:

$$\tilde{y}_i = \operatorname{argmax}\{p_d(y = d | x_i \in Q)\}_{d=1}^{q_n}, \quad (11)$$

where, $p_d(y = d)$ indicates the probability of a sample in the query set belonging to class d . Eventually, the model parameters are updated by calculating the loss of the model, and then back-propagated.

3.2. The diagnosis process of the proposed method

The general framework of the proposed method in this paper is shown in Fig. 3. The training process of method can be divided into four steps.

Step 1. Sample division and data input: The vibration acceleration signals of the bearing under different health conditions are collected. The meta-training set and the meta-test set contain multiple sets of 2048 data points generated by the time series signal, i.e., $x_i \in R^{1 \times 2048}$. According to the learning method proposed in Section 3.1, the sample can be represented as $X = \{x_i\}_{i=1}^{s_l+s_u+q_n} \in R^{(s_l+s_u+q_n) \times 1 \times 2048}$. Where s_l is a labeled sample in the support set S_l , s_u is an unlabeled sample in the support set S_u , and q_n is the sample in the query set.

Step 2. COBs block: The data is mapped to a feature space and then merge all the channel features to obtain a mapped feature vector of $f_e(x_i^s)$. Therefore, A convolution operation block (COBs) consisting of multiple cobs of the same type is constructed, which include convolution block, batch normalization block, activation function ReLU and pooling layer. Consequently, a COBs block is designed to facilitate feature extraction. The module is capable of mapping input data to a feature space. The details about the convolutional operation block can be seen in Table 1. The flow chart of the structure is shown in the upper right block diagram in Fig. 3.

Specifically, a convolution operation F_{conv} is performed on the input data and the convolution kernel, thereby extracting the local input feature by sliding. The acceleration of the entire training process is achieved through the use of a batch normalization module. This approach also reduces the shift or disappearance of internal covariance, thus avoiding gradient explosion [29]. The input of k th BN layer can be set $x^k = \{x^{k(1)}, x^{k(2)}, \dots, x^{k(M)}\}$, and the minibatch size is M , so $x^{k(m)} = \{x_1^{k(m)}, x_2^{k(m)}, \dots, x_N^{k(m)}\}$. The BN operation can be shown as follows:

$$\bar{x}_n^{k(m)} = F_{BN}(x_n^{k(m)}) = \gamma_n^k \frac{x_n^{k(m)} - \mu_n}{\sqrt{\sigma_n^2 + \varepsilon}} + \beta_n^k, \quad (12)$$

$$\mu_n = \frac{1}{M} \sum_m x_n^{k(m)}, \quad (13)$$

$$\sigma_n^2 = \frac{1}{M} \sum_m (x_n^{k(m)} - \mu_n)^2, \quad (14)$$

where $\bar{x}_n^{k(m)}$ is BN layer output of the convolution result of the n th layer. μ_n and σ_n^2 are the mean and variance of x_n^k respectively. Usually, ε is a very small constant to prevent invalid calculations when the variance is zero, and the value is usually $\varepsilon = 1 \times 10^{-5}$. Here, the parameters γ_n^k and β_n^k are respectively the scale parameter and the offset parameter to be learned. They are initially set to 1 and 0, respectively.

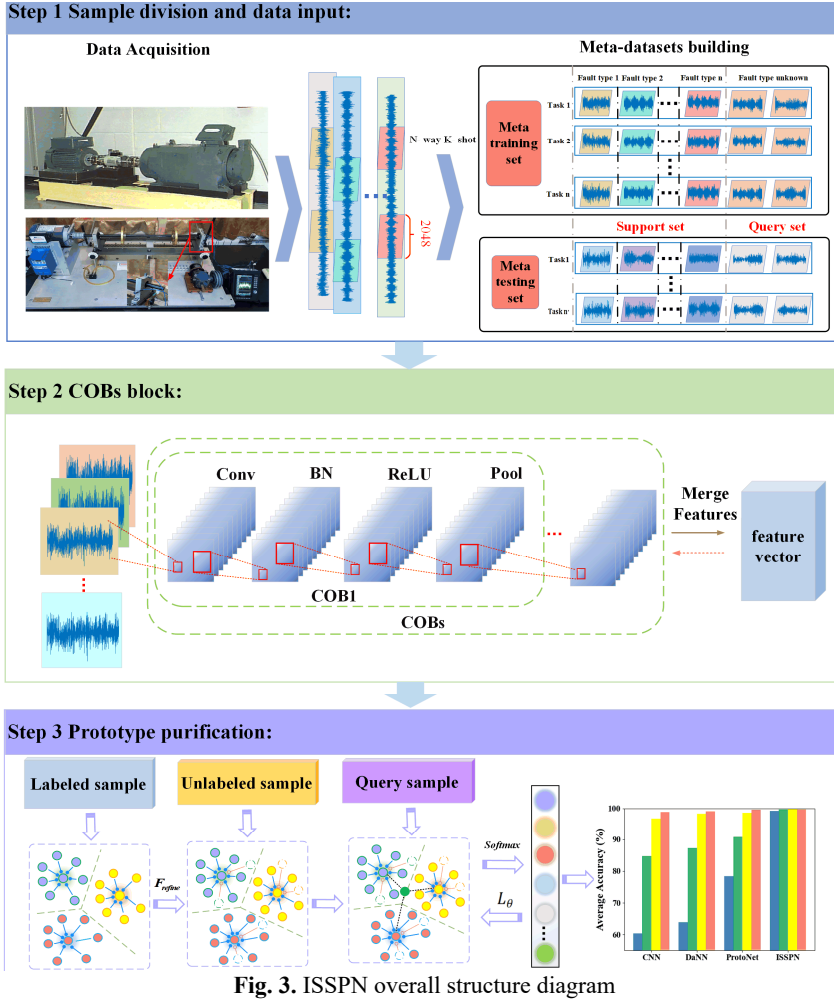


Fig. 3. ISSPN overall structure diagram

To make the features learned by the convolution layer easier to distinguish, the activation function becomes an indispensable part after convolution operation. The ReLU function is a frequently used to activation function in deep neural networks. It can enhance the sparsity of network. Its left saturation function alleviates the gradient vanishing problem of the neural network to a certain extent and accelerates the convergence speed of gradient descent. The ReLU is actually a ramp function, which is defined as follows:

$$\bar{x}_n^k = F_{ReLU}(\bar{x}_n^k) = \max(0, \bar{x}_n^k), \quad (15)$$

where \bar{x}_n^k is the output of upper layer (BN layer), and \bar{x}_n^k is the activation amount of \bar{x}_n^k after activation by the ReLU function.

Further, the pooling layer is used to select features and reduce the dimension of features. The overfitting phenomenon can be effectively avoided by reducing the number of parameters. The maximum pooling F_{MaxP} operation can better complete the current task, and it can complete the local maximum operation on the input feature map.

To sum up, set W and K to be the dimension and length of the channel, respectively. The output of sth convolution is $x_i^{(s)} \in R^{W \times K}$. Therefore, the above COBs can be described as:

$$x_i^{(s)} = F_{MaxP} \left\{ F_{ReLU} \left[F_{BN} \left(F_{conv} \left(x_i^{(s-1)} \right) \right) \right] \right\}. \quad (16)$$

Finally, all the channel features are combined. The input data is also mapped to the feature space. Therefore, it can be described as:

$$f_e(x_i) = [\hat{x}_i(1,1), \hat{x}_i(1, 2), \dots, \hat{x}_i(i, j), \dots, \hat{x}_i(W, K)], \quad (17)$$

where $f_e(x_i) \in R^Z$, $Z = W \times K$ is a Z -dimensional eigenvector. The input data also completes the feature mapping process by the COBs block, so the result is $f_e(x_i^s)$.

Step 3 Prototype Purification: The process of refining the prototype is completed by repeatedly executing the Eq. (9) and continuously assigning the value of \tilde{p}_d to P_d . To summarize, it can be generated the original prototype P_d from the output obtained in the previous step. Meanwhile, the unlabeled data s_u is assigned a predictable label \tilde{P}_d . Furthermore, it is aggregated into the original prototype to guide the original prototype with deviation for adjusting the position in multiple iterations until convergence, so as to purify the prototype.

Table 1. Details of the four methods

Block	Layer	Details
COB	1D-conv	64 @ 1×3
	BN	—
	Activation	ReLU
	1D-MaxPooling	1×2

Step 4. Update the model: The q_n sample in the query set is purified by the prototype to obtain its prediction probability and realize classification all at once. Furthermore, the model parameters are updated by calculating the loss of L_θ and propagating backward.

Although the stochastic gradient descent (SGD) optimizer can quickly reduce the loss in the early stage of training, it often struggles to converge to an optimal minimum value for the model loss. The ISSPN is trained more efficiently by using adaptive moment estimation, which computes bias-corrected first and second moment estimates to counteract bias. Moreover, it automatically adjusts the learning rate scale for different layers, eliminating the need for manual selection as in the case of SGD. Hence, it is easier to use [30].

4. Experimental analysis: The SQV bearing dataset

The performance of the proposed method is demonstrated by the collected bearing vibration dataset from a comprehensive mechanical failure test bed [31]. In order to fully prove the superiority of the method, three other methods are set up for comparison.

The first comparative experiment is set up on CNN whose feature extraction structure is consistent with the method in this paper. Similarly, DANN [32] can also be used as a good comparison object, which has the same feature extractor structure as ISSPN. DaNN is a

CNN-based domain-adaptive deep neural network. It adjusts the difference between two domains by minimizing the Maximum Mean Difference (MMD) in the source domain and the target domain. The last group of comparative experiments is set in ProtoNet, which is the basis of this paper. Table 2 gives the details of the four methods including the proposed one in this paper. In addition, the results are characterized by visualization of tables, graphs, confusion matrices, and t-distributed random neighborhood embedding (t-SNE) [33].

Table 2. Details of the four methods

Models	Learning style	Structure		Optimizer
		Block	Classifier	
CNN	Supervised	4*COB	Linear, Softmax	Adam
DANN	Semi-Supervised	4*COB	Linear, Softmax	Adam
ProtoNet	Supervised FSL	4*COB	Metric function	Adam
ISSPN	Semi-Supervised FSL	4*COB	Metric function	Adam

4.1. Description of the dataset

Due the fault signal characteristics of CWRU data set are very obvious. The validity of NSSPN method is further verified by collecting the bearing vibration signals of the comprehensive mechanical fault simulation test bench of the Spectral Mission (SQ) Company. Fig. 4 shows the integrated mechanical failure simulation test rig. The main part of the test bench is composed of driving motor, rotor system, motor control system, load and signal acquisition system. In the experiment, the vibration acceleration sensor is used to collect the bearing signal of the motor. The data acquisition instrument used is CoCo80, and the sampling frequency is 25.6 kHz. The experimental bearing is installed at the end of the drive motor, then a single point of failure is manually set on the experimental bearing. Specific information on the SQV bearing dataset is shown in Table 3. In addition, the experimental rotor system has four rotation frequencies (9 Hz, 19 Hz, 29 Hz, and 39 Hz). Fig. 5 shows a Time-Domain signal of the operating condition of the bearing at a rotation frequency of 39 Hz.

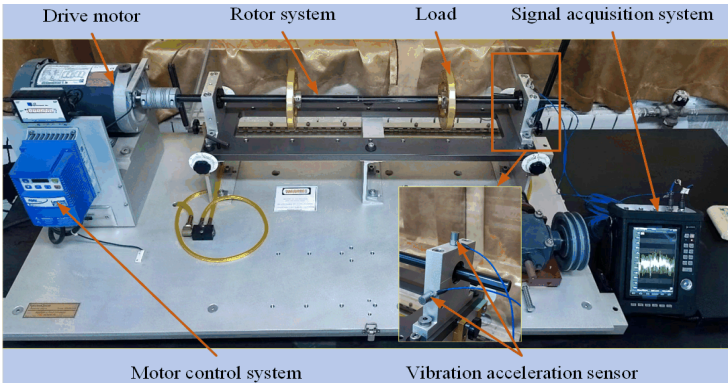


Fig. 4. SQV bearing dataset test bench

4.2. Experimental parameter setting

To further evaluate the performance of the proposed method, two experiments are conducted on the SQV bearing dataset. The first experiment is three fault identifications (NC, IF-3 and OF-3) at four different speeds. First of all, a training data set is constructed: 100 samples are generated for each of the three fault conditions at each speed. So a total of 100×4×3 samples are generated. One to four samples are selected for each fault at each speed, resulting in four to sixteen samples for each type of fault. The number of unlabeled samples for ISSPN follows the support set, i.e. {1, 5} for each class. Ultimately, a single experiment and five experiments are set up to

demonstrate the performance of the method, where each experiment is conducted for 30 iterations.

The second experiment is the identification of seven faults at one speed (39 Hz). Again, a data set needs to be constructed: 200 samples are generated for the 7 bearing States at 39 Hz, so a total of 1400 samples are generated. 5, 10, 20 and 30 samples are selected for each fault type. The training parameters are consistent with the above settings.

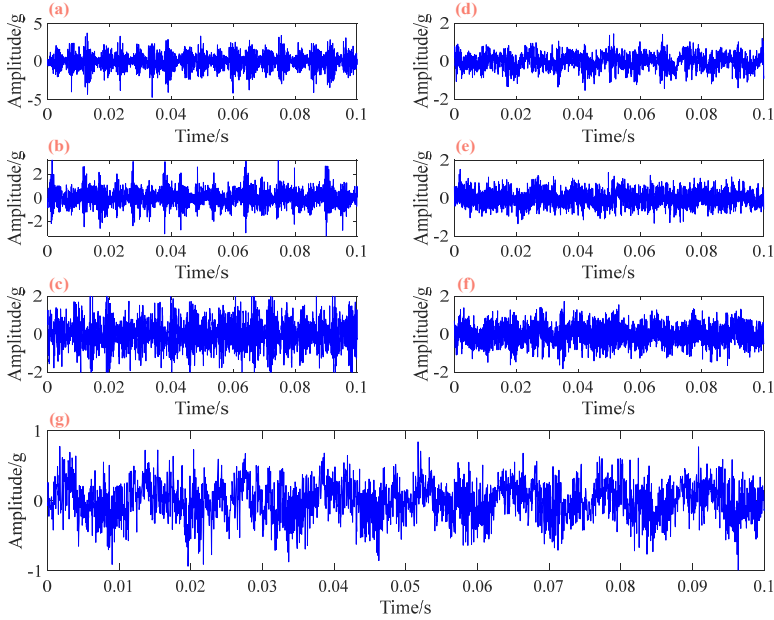


Fig. 5. Time-domain signal of seven kinds of bearing vibration signals: a) mild inner ring fault (IF-1), b) moderate inner ring fault (IF-2), c) severe inner ring fault (IF-3), d) mild outer ring fault (OF-1), e) moderate outer ring fault (OF-2), f) severe outer ring fault (OF-3), g) normal condition (NC)

Table 3. Details of SQV bearing dataset

Bearing condition	Fault damage degree	Damage scale (area/depth)
Normal condition (NC)	Mild fault (1)	4/0.5
Inner ring fault (IF)	Moderate fault (2)	8/4
Outer ring fault (OF)	Severe Fault (3)	12/2

4.3. Experimental results and analysis

Table 4 shows the results of the three types of fault identification accuracy, and Fig. 6. visually shows the identification accuracy of each method. From the above tables and figures, it can be seen that although the accuracy of CNN, DaNN and ProtoNet is not less than 85 %, the ISSPN method shows better performance. The fault recognition rate of ISSPN can still reach 98.64 % when there are only 4 samples in each class. To sum up, the ISSPN method can adapt to fault diagnosis at different speeds by comparing with the other three methods.

Table 4. Identification accuracy (%) of 3 types of faults on SQV dataset

Models	One experiment			Five experiments		
	4	10	16	4	10	16
CNN	85.94	91.67	99.38	90.75	99.05	100
DaNN	95.32	99.78	99.86	96.74	99.85	100
ProtoNet	91.27	97.69	99.79	97.19	100	100
ISSPN	98.64	100	100	100	100	100

Besides, the accuracy results for the seven types of faults are shown in Table 5 and Fig. 7. The ISSPN method also shows excellent performance in the identification of seven types of faults by comparing with the other three methods. Similarly, the fault identification of ISSPN is still excellent when there are the fewest fault samples of each type. So it can be seen that the ISSPN method can complete the multi-class fault diagnosis excellently.

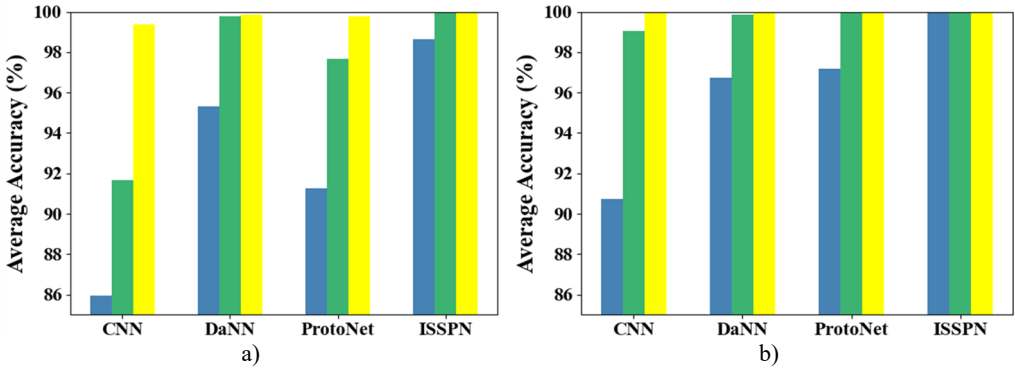


Fig. 6. Accuracy of identification of three types of faults on SQV dataset:
a) one experiment, b) five experiment

Table 5. Identification accuracy (%) of 7 types of faults on SQV dataset

Models	One experiment				Five experiments			
	5	10	20	30	5	10	20	30
CNN	60.38	84.76	96.78	98.89	63.46	87.94	98.04	99.43
DANN	63.92	87.26	98.34	99.08	65.52	89.5	98.54	99.63
ProtoNet	78.43	90.76	98.65	99.57	83.76	94.91	99.16	99.8
ISSPN	99.27	99.73	100	100	99.86	99.94	100	100

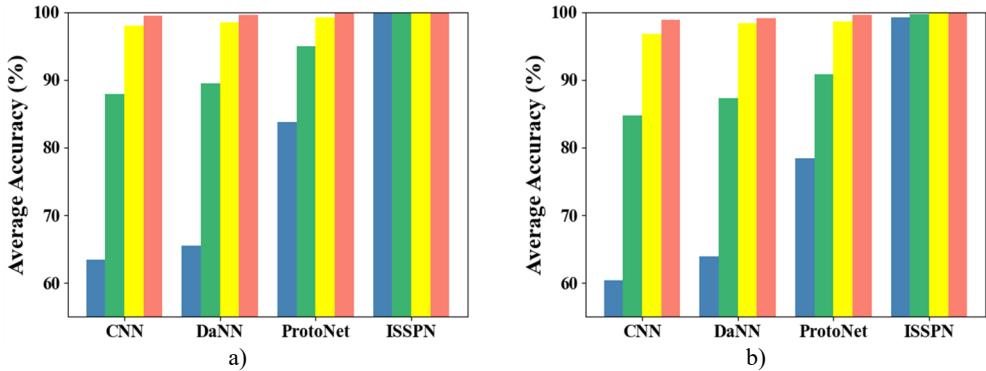


Fig. 7. Accuracy of identification of 7 types of faults on SQV dataset:
a) one experiment, b) five experiment

Taking the classification of seven types of faults as an example, Fig. 8 and Fig. 9 show the classification characterization results of ISSPN method in the form of confusion matrix and t-SNE visualization, respectively. It can be seen from the confusion matrix that the ISSPN method can accurately complete the classification of seven types of faults. The classification results of ISSPN method are also more obvious in the t-SNE visualization map. In summary, the ISSPN method has better feature extraction ability.

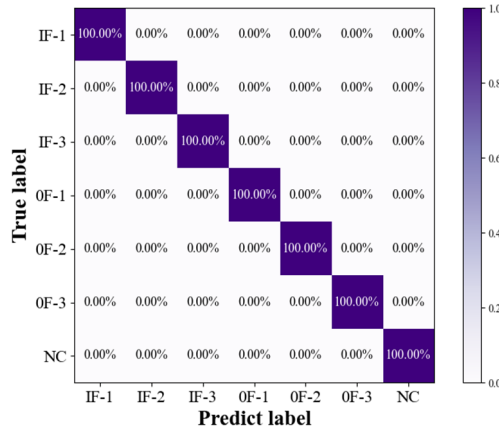


Fig. 8. Confusion matrix of ISSPN seven-class classification results on SQV dataset

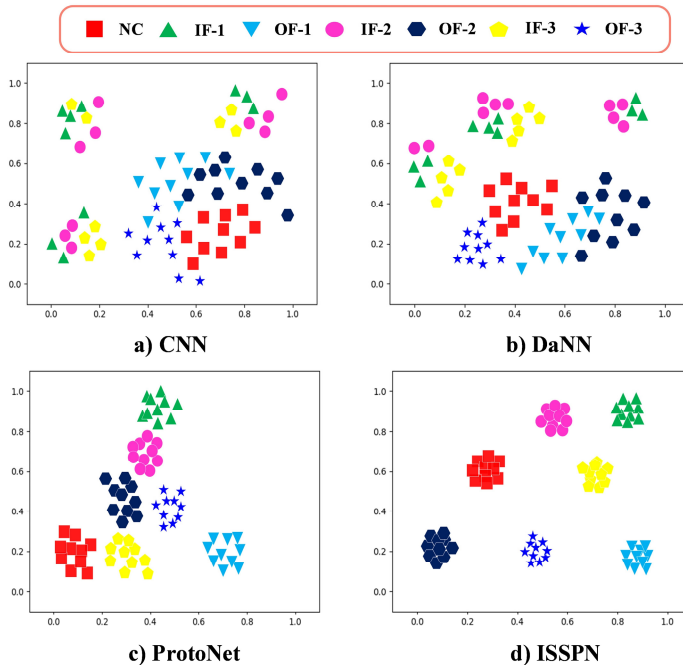


Fig. 9. Visualization of t-SNE for the four methods on SQV dataset

5. Conclusions

This paper presents an improved semi-supervised learning network (ISSPN) for few-shot fault diagnosis. It can obtain more effective features with limited data, thus effectively solving the shot less problem in machinery fault diagnosis. The method adopts a semi-supervised learning strategy to make full use of unlabeled data to obtain more accurate prototypes. Meanwhile, the variability due to different scales is eliminated by the standard Euclidean distance metric function, which improves the robustness of the model. The performance of the method is validated with one case, and results show that the method can accurately identify all kinds of faults under the condition of few-shot and have good generalization performance. In addition, the performance of ISSPN is better compared to other similar methods.

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Data availability

The datasets generated during and/or analyzed during the current study are available from the corresponding author on reasonable request.

Author contributions

Zhenlian Lu: writing-review and editing, writing-original draft, validation, software, methodology, conceptualization. Kuosheng Jiang: writing-review and editing, validation, investigation, resources, funding acquisition. Jie Wu: writing-review and editing, validation, supervision, investigation, formal analysis.

Conflict of interest

The authors declare that they have no conflict of interest.

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Zhenlian Lu received the B.S. degree in vehicle engineering from the Anhui University of Science and Technology, Huainan, Anhui, in 2021, where he is currently pursuing the M.Eng. degree. His main research interests include fault diagnosis and mechanical design.



Kuosheng Jiang received the Ph.D. degree from the State Key Laboratory for Manufacturing Systems Engineering, School of Mechanical Engineering, Xi'an Jiaotong University, Xi'an, China, in 2016. Since January 2020, he has been an Associate Professor with the School of Mechanical Engineering, Anhui University of Science and Technology, Huainan, China. His main research interests include mechanical equipment condition monitoring, fault diagnosis analysis, intelligent visual inspection, intelligent instrument, and virtual instrument design.



Jie Wu received the Ph.D. degree in mechanical engineering from Xi'an Jiaotong University, Xi'an, China, in 2019. He is currently an associate professor with the school of Mechanical Engineering, Anyang Institute of Technology. His research interests include intelligent diagnosis, complex equipment diagnoses and mechanical signal processing.