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Research and Education Center "Higher IT School" (HITs)

APPROVED BY

Head of the Main Educational Program

Professor, Doctor of Physical and Mathematical Sciences

car sciences

O.A. Zmeev

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BACHELOR'S THESIS

WIND SPEED PREDICTION PERFORMANCE BASED ON MODEL DECOMPOSITION METHODS AND OPTIMIZATION ALGORITHMS

Main Educational Program

02.03.02 – Fundamental Computer Science and Information Technology
Specialization "Software Engineering"

Hu Zhichao

Bachelor's Thesis Supervisor Associate Professor, Cand.of Science (Physics & Mathematics)

Zh. N. Zenkova

____ 2022

Consultant

Assistant Lecturer of Software Engineering Department of TSU

D. O. Zmeev

«<u>08</u>» <u>06</u> 2022

Written by

Student of Group

No.<u>971812</u>

Zhickoo HuZh. Hu

« 06 » 06 2022

The Ministry of Science and Higher Education of the Russian Federation NATIONAL RESEARCH TOMSK STATE UNIVERSITY (NR TSU)

REC "Higher IT School"

THE TASK

Hu Zhichao

algorithms

"Software Engineering"

алгоритмов оптимизации

a) to academic office -

3. Description:

б) to State Examination Board -

2. The deadlines for the completion of task:

needs for better management of wind farm.

The organization that is being researched:

Final qualification work supervisor

Final qualification work consultant

Assistant Lecturer of Software Engineering Department of TSU

The task is accepted for execution

18.02

(Physics & Mathematics)

Associate Professor, Cand. of Science

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APPROVED MAP Head hys.-Math.), professor O.A. Zmeev Regarding the bachelor's final qualification work implementation by a student (student's full name) In the field of study Fundamental Computer Science and Information Technology, specialization 1. The bachelor's final qualification work theme (in the English and in the Russian language) Wind speed prediction performance based on model decomposition methods and optimization Эффективность прогнозирования скорости ветра на основе методов декомпозиции модели и The purpose is to explore the wind speed prediction models with the best prediction accuracy, goals and objectives of the FQW, expected results Expected results: The error analysis of the selected model can be carried out according to the evaluation index, and the model with the smallest error and the optimal model can be obtained. Zhanna N. Zenkova (signature) (initials, surname) Denis O. Zmeev (signature) (initials, surname) Zhichao Hu

(signature)

(initials, surname)

ABSTRACT

With the problems such as climate change caused by fossil energy, various countries have started to pay attention to renewable energy, and wind energy has become a key research object due to the problem of non-pollution. Since the accuracy of wind energy prediction mainly depends on the accuracy of wind speed prediction. Therefore, exploring and seeking methods to improve the accuracy of wind speed prediction has become the most important issue at present. In this paper, propose to use Back-Propagation Neural Networks(BPNN) and Support Vector Regression (SVR) and Kernel Based Extreme Learning Machine (KELM) as three different sets of base models and optimize them by using different machine learning algorithms such as Genetic Algorithm (GA), Particle Swarm Optimization (PSO), Sparrow Search Algorithm (SSA), Sine Cosine Algorithm (SCA), and then combine them with different decomposition methods such as Empirical Mode Decomposition (EMD), Ensemble Empirical Mode Decomposition (EEMD), Variational Mode Decomposition (VMD), Wavelet Decomposition (Wavelet) to form the corresponding combined prediction models to explore which combined prediction model has higher accuracy. Based on the data from the National Center for Meteorological Sciences, the experiments show that the prediction accuracy of the combined prediction model formed by the VMD method is significantly higher than that of other combined prediction models in some models of BPNN and most of SVR, while the prediction accuracy of the combined prediction model formed by the Wavelet combination is higher in the models of KELM and part of models of SVR.

Keywords – Wind speed prediction, Model Decomposition Method, Back-Propagation Neural Networks, Support Vector Regression, Kernel Based Extreme Learning Machine, Genetic Algorithm, Particle Swarm Optimization, Sparrow Search Algorithm, Sine Cosine Algorithm.

Bachelor's thesis 70 pages, 22 figures, 12 code samples, 5 tables, 63 sources, 1 appendix.

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GLOSSARY

EMD¹⁹– Empirical Mode Decomposition

EEMD²⁰ – Ensemble Empirical Mode Decomposition

VMD²¹ – Variational Mode Decomposition

Wavelet²² – Wavelet Decomposition

IMF – Intrinsic Mode Functions

BPNN¹⁴ – Back Propagation Neural Network

SVR¹⁵ – Support Vector Regression

SVM – Support Vector Machine

KELM¹⁶ – Kernel Based Extreme Learning Machine

GA¹² – Genetic Algorithm

PSO¹³ – Particle Swarm Optimization

SSA¹⁷ – Sparrow Search Algorithm

SCA¹⁸– Sine Cosine Algorithm

MAPE – Mean Absolute Percentage Error

MSE – Mean Squared Error

MAE – Mean Absolute Error

RMSE – Root Mean Squared Error

Matlab¹ – It's a programming and numeric computing platform used by millions of engineers and scientists to analyze data, develop algorithms, and create models

INTRODUCTION

Since the 21st century, as the new round of technological revolution and industrial change continues to develop in depth, a new generation of technologies and applications represented by artificial intelligence has penetrated into every aspect of society, and people are enjoying the convenience and prosperity brought by these new technologies. However, climate change, caused by the use of fossil fuels, is threatening everything. According to the UN survey elevation, the temperature on earth has increased by 1.1°C in the last decade (2011-2020) compared to the late 19th century, and extreme droughts, water shortages, major fires, and sea level rise caused by climate change are threatening human life². As a result, renewable energy has become a key research object for countries around the world, and renewable energy targets and support policies have spread almost all over the world, and according to the report, at least 164 countries have set renewable energy targets and 145 countries have introduced support policies³. And wind energy, as a non-polluting renewable energy source, has become a key research target for countries around the world, and according to IRENA, solar and wind energy account for 91% of all new installed renewable energy capacity⁴. However, the stochastic and fluctuating nature of wind energy complicates the entry of a high percentage of wind energy into the grid system. Between the accuracy of wind energy prediction results depends mainly on the accuracy of wind speed prediction. Wind speed prediction can effectively reduce the risk associated with wind-related uncertainty⁵. How to use modern machine learning techniques to improve the accuracy of wind

speed prediction results and how to use machine learning algorithms to explore the optimal prediction model have become the fields that this paper wants to explore deeply.

According to the time scale of prediction, wind speed predictions can be classified into ultra-short-term predictions (mainly in minutes), short-term predictions (mainly in days), medium-term predictions (mainly in months, weeks) and long-term predictions (mainly in years)⁶. For wind speed prediction methods are mainly divided into physical and statistical methods. Physical methods use real-time meteorological conditions for prediction, however, due to the complexity of numerical meteorological models, they are not suitable for short-term and ultrashort-term wind speed prediction⁷. Statistical methods are used to establish a functional relationship between historical data and wind power output by means of mathematical statistics, such as time series method⁸, neural network method⁹, Kalman filter method¹⁰, support vector machine method¹¹, etc. However, many scholars have combined with other optimization algorithms and other techniques in order to build a better combined prediction model and thus to improve the prediction accuracy. For example, some scholars have used machine learning optimization algorithms such as genetic optimization algorithm¹² and Particle Swarm Optimization algorithm¹³ to optimize the parameters of a single prediction model in order to improve the wind speed prediction accuracy.

Based on this, this paper proposes Back-Propagation Neural Networks (BPNN)¹⁴ and Support Vector Regression(SVR)¹⁵ and Kernel Based Extreme

Learning Machine(KELM)¹⁶ as the three base models, and the base models are optimized by machine learning algorithms such as Genetic Algorithm (GA)¹², Particle Swarm Optimization (PSO)¹³, Sparrow Search Algorithm (SSA)¹⁷, Sine Cosine Algorithm (SCA)¹⁸, etc. to optimize the base model, and then with different decomposition methods such as Empirical Mode Decomposition (EMD)¹⁹, Ensemble Empirical Mode Decomposition (EEMD)²⁰, Variational Mode Decomposition (VMD)²¹, Wavelet Decomposition (Wavelet)²², etc., to form a combined prediction model to predict the wind speed, and to explore which combined prediction model has a higher The accuracy of the combined prediction model is higher.

The paper is organized as follows: the first part is a background introduction, the second part focuses on the theoretical knowledge of decomposition and prediction methods and the corresponding optimization algorithms, and the third part analyzes different combinatorial prediction models through experimental studies. The fourth part is introduction to the main program code. The fifth part presents the corresponding research conclusions based on the experimental data analysis. Part A of the Appendix A introduces the specific code realization of the optimization algorithm in Matlab.

1 Introduction to the methodologies

1.1 Introduction to the base model algorithms

1.1.1 Back-Propagation Neural Networks

Back-Propagation Neural Networks (BPNN) is a multilayer neural network trained according to the error back-propagation algorithm. The BPNN model has a powerful computational power and a very complex mapping capability. Based on this, the BPNN model can be trained adaptively for a large number of unstructured, non-exact laws²³.

The network structure of BPNN consists of: input layer, hidden layer and output layer. The structure is illustrated in Figure 1²⁴ below. The input layer is the first layer of the neural network, which takes the input information and passes the signal to the next layer. The input layer does not perform any processing on the data signal. Hidden layer: The other layers between the input layer and the output layer are called hidden layers, which usually do not receive signals directly from the outside world and do not send signals to the outside world. It is a general term for the different functional layers. Output layer: as the last layer of the neural network, it is mainly used to receive the input from the hidden layer and to output the resultant values predicted by the model, The code construction of the keras-based BPNN algorithm is shown in code sample 1²⁵, For example, the neural network shown in code sample 1 is built quickly by keras and contains an input layer with 7129 nodes,

and a hidden layer with 128 nodes, and an output layer, which is a common binary classification model.

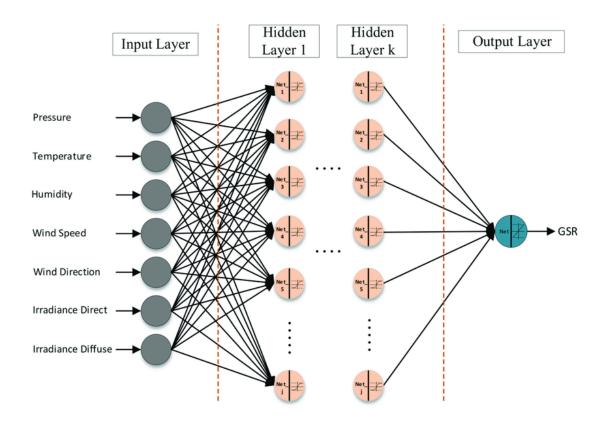


Figure 1 – Backward propagation neural network structure

Code sample 1 – Neural networks built on keras code

1.1.2 Support Vector Regression

Support Vector Regression are a new type of classifier based on the theoretical foundation of Support Vector Machine (SVM). The purpose of SVM²⁶ is to find a hyperplane so that the sample points are as far away from the hyperplane as possible. The SVR, on the other hand, wants the sample points to be as close to the hyperplane as possible. The model is optimized by minimizing the width between intervals²⁷. The structure is schematically shown in the following Figure 2 and Figure 3²⁸.

So, it can be considered statistically that the traditional regression method is only considered correct when the regression function f(x) is exactly equal to y, when and only when this condition occurs, all kinds of complex losses need to be calculated, which makes our computation much more. On the contrary, support vector regression considers that if the regression function f(x) deviates from y within an acceptable range, this acceptable range is called the threshold (N). As long as |f(x) - y| > N. Then the prediction can be considered correct. The equation for the specific SVR regression model can be expressed as formula $(1)^{29}$:

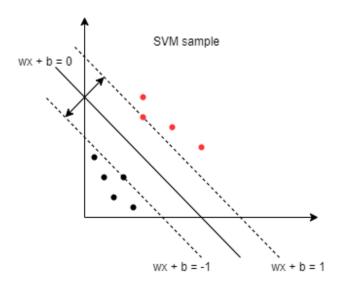


Figure 2 – Schematic diagram of SVM

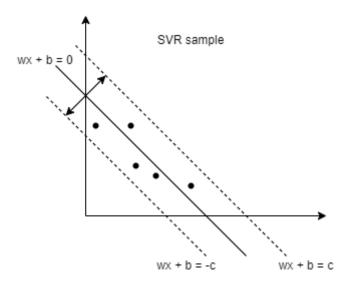


Figure 3 – Schematic diagram of SVR

$$f(x) = w^T x + b \tag{1}$$

In the formula (1), w and b are the model parameters to be determined.

1.1.3 Kernel Based Extreme Learning Machine

Kernel Based Extreme Learning Machine (KELM) is an improved learning algorithm based on the theoretical basis of Extreme Learning Machine (ELM)³⁰ and combined with kernel functions. By referring to the kernel function in ELM, the KELM algorithm increases the stability and robustness of the classification model on the basis of the original ELM algorithm³¹, which makes the KELM algorithm have better performance to deal with more classification problems³².

ELM is feed-forward neural network, whose objective function f(x) can be expressed as formula 2^{33} :

$$f(x) = a(x) \times b = A \times b = L \tag{2}$$

In the formular 2, x is the input vector, a(x), A is the output of the hidden layer node, b is the output weight, L is the desired output.

By introducing the regularization factor C and the unit matrix I on the basis of formula 2, the expression of the output weights is shown in formula 3^{33} :

$$b = A^T (AA^T + \frac{I}{c})^{-1}L \tag{3}$$

Then, the kernel function is introduced, and the matrix equation of the kernel function can be expressed as formula 4^{33} :

$$\Omega ELM = AA^{T} = \alpha(x_i)\alpha(x_i) = K(x_i, x_i)$$
(4)

In formula 4, x_i, x_i is the test input vector.

Finally, the expression can be expressed as formula 5^{33} :

$$f(x) = [K(x, x_1); \dots; K(x, x_n)](\frac{l}{c} + \Omega ELM)^{-1}L$$
 (5)

In formula 5, $(x_1, x_2, x_3, ..., x_n)$ is the given training sample, n indicates the sample size, K() is denoted as the kernel function.

1.2 Optimization algorithms

1.2.1 Genetic Algorithm (GA)

The main idea of Genetic Algorithm is derived from Darwin's theory of biological evolution of natural selection and the computational model of biological evolution of genetic mechanism, and it is a machine learning algorithm that searches for the optimal solution through the natural selection process of natural evolution of superiority³⁵.

The Genetic Algorithm is an efficient search for each running parameter code by randomization technique using all the individuals in the population as the object³⁶. The algorithm operation process is shown in Figure 4 below, which selects the optimal parameter code through three evolutionary operations: selection, crossover and mutation. Because of its own efficiency and class parallel processing and global optimality has a very obvious advantage in dealing with various nonlinear problems and solving multi-objective constrained problems³⁷.

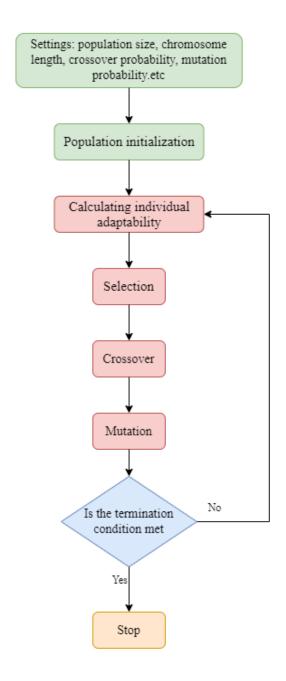


Figure 4 – Execution process of Genetic Algorithm

The three important operations of Genetic Algorithm are selection, crossover and mutation. For the code implementation, please refer to the Genetic Algorithm in Appendix A^{38} .

1.2.2 Particle Swarm Optimization (PSO)

The main idea of the Particle Swarm Optimization algorithm originates from the process of foraging for food by flocks of birds and the process of collaborating with each other to complete the search for food. It is simpler than the Genetic Algorithm because there is no "crossover" or "mutation" operation³⁹. The Particle Swarm Optimization algorithm is based on an iterative approach to find the global optimal solution by iteratively updating the optimal solution over and over again⁴⁰.

It has become a popular research area in optimization algorithms because of its simple principle, few adjustable parameters and easier implementation⁴¹. The algorithm operation process is shown in Figure 5 below:

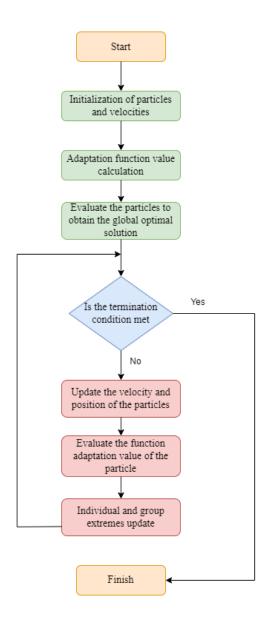


Figure 5 – Execution process of Particle Swarm Optimization

The PSO algorithm finds the optimal solution by iteration. During each iteration, the particle updates its velocity and position by individual and population extremes. The optimal solution found by the particle itself is called the individual extremum, which can be represented by pbest. The other optimal solution found in the population is called the population extreme, which can be represented by gbest. The formula for the particle to update its own velocity and position can be represented by the following formula 6 and formula 7⁴²:

Velocity:

$$v_{i+1} = w \times v_i + c_1 \times rand_1 \times (pbest_i - x_i) + c_2 \times rand_2 \times (gbest_i - x_i)$$
 (6)

Position:

$$x_i = x_i + v_{i+1} \tag{7}$$

In the above two formulas, w is the inertia factor, generally taken as 1. c_1 and c_2 are the learning factors, generally taken as 2. $rand_1$ and $rand_2$ are random numbers between (0, 1). v_i and x_i denote the velocity and position of the particle in the i-th dimension, respectively. $pbest_i$, $gbest_i$ denote the value of the i-th dimension of the best position of a particle, the value of the i-th dimension of the best population, respectively.

For code implementation of Particle Swarm Optimization algorithm, please refer to Particle Swarm Optimization in Appendix A.

1.2.3 Sparrow Search Algorithm (SSA)

The sparrow search optimization algorithm was proposed in 2020, and its main idea is derived from sparrow foraging behavior and anti-predation behavior, which is a new type of swarm intelligence optimization algorithm⁴⁴. The sparrow search optimization algorithm can be considered as an improvement of the artificial swarm algorithm, and it has almost the same structure as the artificial swarm algorithm, only with some differences in the search. Its algorithm flow is mainly shown in Figure 6 below⁴⁵.

In the Sparrow Search Algorithm, in which the discoverers, especially those possessing better adaptation worthy discoverers, are given priority in the search process to obtain food. At the same time, the discoverers can also provide foraging directions for the joiners in all populations, so that the discoverers can forage in a larger area. Similarly, the joiners will always monitor the discoverers during the foraging process. If the entire sparrow population is threatened by a predator, it will immediately engage in anti-predatory behavior⁴⁶.

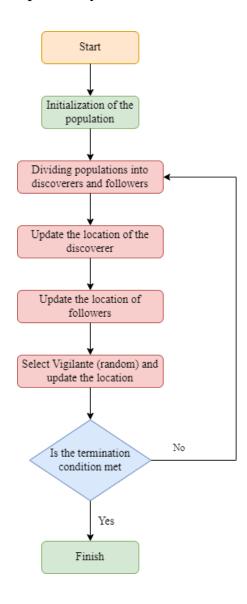


Figure 6 – Flow chart of the execution process of the Sparrow Search Algorithm

For code implementation of Sparrow Search Algorithm, please refer to Sparrow Search Algorithm in Appendix A.

1.2.4 Sine Cosine Algorithm (SCA)

The Sine Cosine Algorithm is a new stochastic optimization algorithm that is an intelligent optimization algorithm proposed by Seyedali Mirjalili, Australia⁴⁷.

The algorithm works by creating multiple initial stochastic candidate solutions and then fluctuates towards the optimal solution using mathematical models of sine and cosine. The stochastic solutions are evaluated iteratively by the objective function so that different regions in the space can be used to achieve the global optimum, very efficiently avoiding the local optimum problem⁴⁸.

The sine cosine optimization algorithm is widely used in various fields of optimization problems because of its flexibility, simplicity and ease of implementation. The operation process of the sine cosine optimization algorithm can be divided into the following two phases: the exploration phase and the development phase. In the exploration phase, the initial stochastic solution is combined with the optimization algorithm to quickly search for feasible spatial regions among all the stochastic solutions. During the development phase, the random solution changes, but the rate of change is always lower than that of the exploration phase. Its formulation can be expressed by the following equation 8⁴⁹.

$$X_i^{t+1} = \begin{cases} X_i^t + r_1 \times \sin(r_2) \times |r_3 P_i^t - X_i^t| & r_4 < 0.5 \\ X_i^t + r_1 \times \cos(r_2) \times |r_3 P_i^t - X_i^t| & r_4 > 0.5 \end{cases}$$
(8)

In this equation, X_i^t is the i-th dimension of the current individual and the position of the t-th generation. r_2 denotes a random number from 0 to 2π . r_3 denotes a random number between 0 to 2. r_4 denotes a random number from 0 to 1. P_i^t denotes the position of the *i-th* dimension of the optimal individual position variable at the *t-th* iteration.

The specific execution flow chart of the SCA is shown in Figure 7 below:

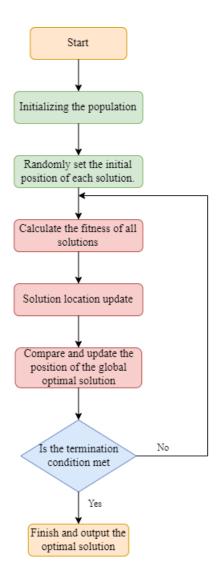


Figure 7 – Flow chart of the execution process of the Sine Cosine Algorithm

The code to initialize the random result set of the Sine Cosine Algorithm and the main loop is shown in Sine Cosine Algorithm in Appendix A.

1.3 Decomposition methods

1.3.1 Empirical Mode Decomposition (EMD)

The EMD algorithm was proposed by Norden e. Huang, an American scientist, and it can decompose a signal by virtue of the time-scale characteristics of the data itself without setting any basis function in advance⁵¹. EMD algorithm has a very obvious advantage in dealing with various nonlinear and non-stationary signals because of its data-driven adaptive nature⁵².

EMD obtains the first-order residual quantity $r_1(t)$ by differentiating the original signal X(t) and $c_1(t)$, and then replaces $r_1(t)$ with the original signal X(t) for the corresponding processing, and the n-th-order modal function $c_n(t)$ and the final residual quantity $r_n(t)$ can be obtained after repeating n times. The final decomposition expression is⁵:

$$X(t) = \sum_{1}^{n} c_n(t) + r_n(t)$$
 (9)

During the EMD method, it is very easy to generate modal mixing because of the presence of signals with similar scales in the IMF component. Also because the signal has no sample points to be considered after the beginning and the end, it can also lead to the polar packets diverging at the end points, thus leading to errors,

which is often referred to as the end effect. Therefore, in order to solve these problems, the EEMD method is proposed⁵³.

1.3.2 Ensemble Empirical Mode Decomposition (EEMD)

The EEMD algorithm is an improved algorithm proposed by Norden e. Huang and other scientists to suppress the modal aliasing phenomenon caused by the EMD algorithm. By superimposing a Gaussian white noise auxiliary signal processing (NADA), the modal aliasing phenomenon is effectively suppressed⁵⁵.

The specific steps of the decomposition of the EEMD algorithm are shown below⁵:

The white noise $M_i(t)$ with normal distribution is added to the original signal X(t) to obtain the new signal $X'(t)^5$:

$$X'(t) = X(t) + M_i(t)$$
 (10)

2) Then the obtained new signal X'(t) is decomposed by EMD, so that the IMF component and the other remaining components can be obtained⁵:

$$X'(t) = \sum_{j=1}^{n} c_j(t) + r_n(t)$$
 (11)

Then repeat the execution of step 1) and step 2) N times, and then the above corresponding IMF will be averaged⁵, and finally can obtain the final IMF $(C_n(t), \text{ after EEMD as follows:}$

$$C_n(t) = \frac{1}{n} \sum_{i=1}^{N} c_i(t)$$
 (12)

1.3.3 Variational Mode Decomposition (VMD)

The VMD algorithm is an adaptive signal processing method whose core idea is to construct and solve variational problems. Unlike the methods proposed by scientists such as Norden e. Huang, the VMD algorithm aims to decompose the original signal into subsignals of different frequencies⁵⁶. This is because the VMD algorithm considers the original signal as formed by the superposition of subsignals with different frequency dominance. Therefore, IMF is defined in the VMD algorithm as a bandwidth-constrained frequency modulation function, and the original signal is decomposed into a number of specified IMF components by constructing and solving a variational problem⁵⁷. It is also because of this feature that the VMD algorithm can effectively avoid the phenomenon of modal aliasing and has the advantage of better decomposition accuracy for processing complex data, and it is for this reason that the VMD algorithm is often used to deal with various non-smooth and non-linear signal problems⁵⁸.

1.3.4 Wavelet Decomposition (Wavelet)

The Wavelet Decomposition algorithm can be seen as an optimization of the Fourier transform, By breaking up a continuous periodic signal into a linear

combination of trigonometric signals of different frequencies, when the period of this periodic signal is infinite, this is the Fourier transform⁵⁹. The basic equation of Fourier transform is sin function and cos function, and the basic equation of wavelet transform is wavelet function. The change made by wavelet transform lies in turning the infinitely long trigonometric function basis into a finite length wavelet basis that will decay, and different wavelets have very big differences in waveform, and similar wavelets form a wavelet family⁶¹. The functional representation of the wavelet transform is shown below. Due to the sparse coding property of Wavelet Decomposition, wavelet packet decomposition is commonly used to deal with various signal filtering problems and to deal with various signal noise reduction problems and various data compression problems⁶².

$$\alpha = W^T f \tag{13}$$

where α denotes the wavelet coefficients obtained from the transform and W is the orthogonal matrix. f is the input signal.

2 Experimental research

2.1 Data preparation

The experimental data for this experiment were obtained from the National Center for Meteorological Sciences (NCMS), and the station located in Qingdao, Shandong Province, with a latitude and longitude position of (36.07°N, 120.33°E) and an elevation of 76 m. The specific experimental idea of this experiment is shown in Figure 8 below:

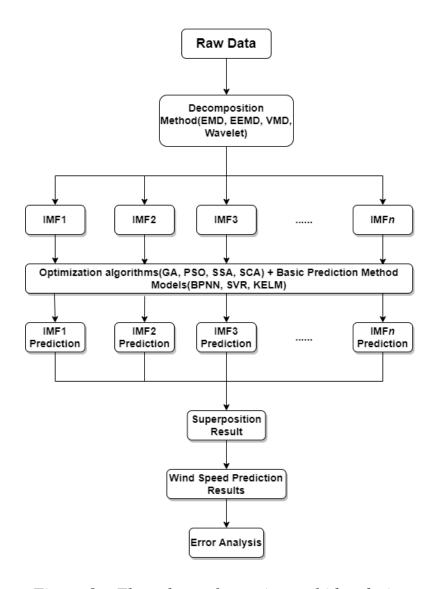


Figure 8 – Flow chart of experimental idea design

2.2 Data decomposition

The sample data for this experiment has a total of 300 and is divided into two different data sets, the training set and the test set, for training and prediction. The first 80% of the sample data is used as the training set for the whole model, and the remaining 20% of the sample data is used as the test set for the whole model.

Firstly, by using different decomposition methods (EMD, EEMD, VMD, Wavelet) to decompose the original wind speed data to obtain IMF1, IMF2, IMF3.....IMFn. The specific decomposition results are shown in the following Figure 9,10:

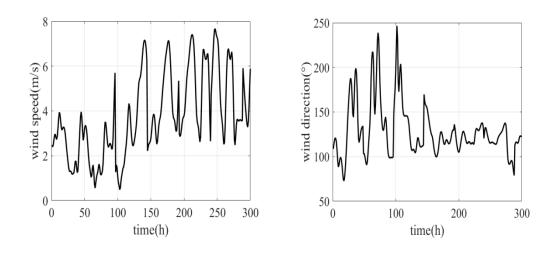


Figure 9, 10 – Raw data signal decomposition

The raw data decomposed by the EMD method is shown in the following Figure 11:

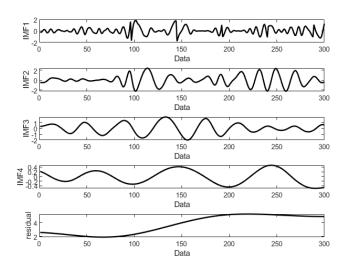


Figure 11 – Raw data signal decomposed by EMD

The raw data by the EEMD method through decomposition is shown in the following Figure 12, 13:

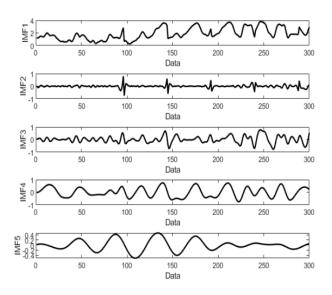


Figure 12 – Raw data signal decomposed by EEMD

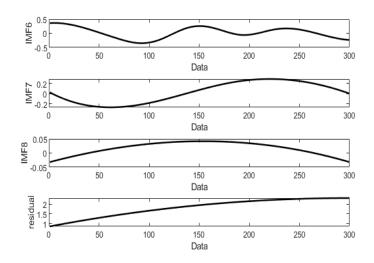


Figure 13 – Raw data signal decomposed by EEMD

The raw data decomposed by the VMD method is shown in the following Figure 14, 15:

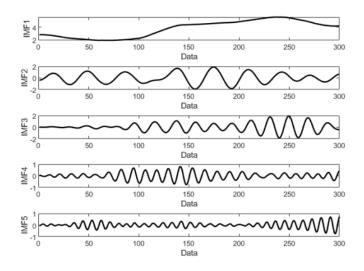


Figure 14 – Raw data signal decomposed by VMD

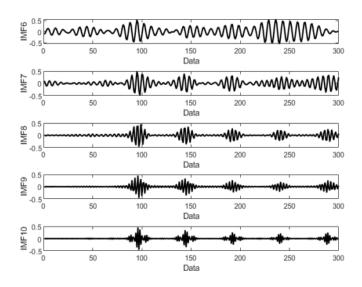


Figure 15 – Raw data signal decomposed by VMD

The raw data decomposed by the Wavelet Decomposition method is shown in the following Figure 16:

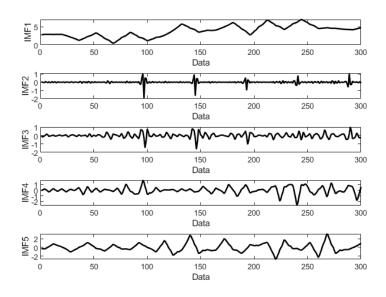


Figure 16 - Raw data signal decomposed by Wavelet

2.3 Predictive models

In this paper, BPNN, SVR and KELM are used as three groups of basic prediction models. In order to improve the accuracy of the model, four different optimization methods (GA, PSO, SSA and SCA) are used to optimize the three groups of basic models, and then they are combined with the decomposition method to obtain the following different groups of combined prediction models.

The prediction results of the combined prediction model formed by combining the EMD method with the prediction model are shown in the following Figure 17:

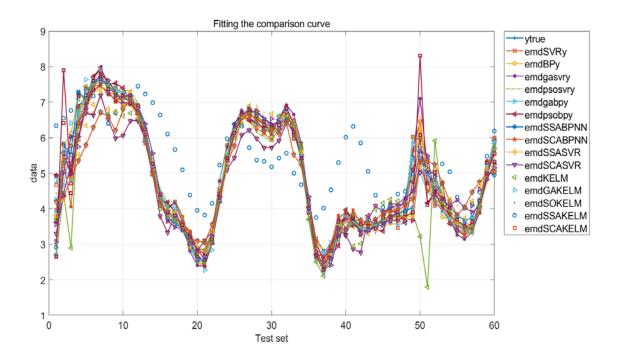


Figure 17 – Prediction diagram of EMD combined prediction model

The prediction results of the combined prediction model formed by combining the EEMD method with the prediction model are shown in the following Figure 18:

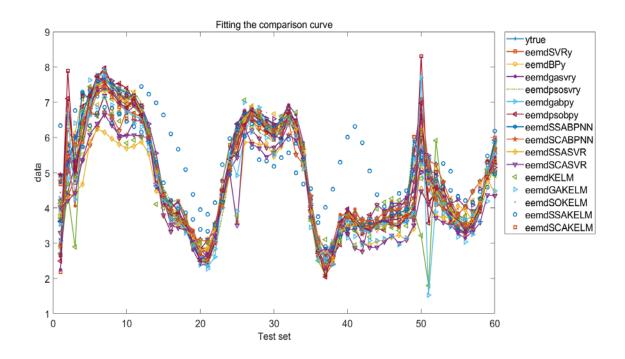


Figure 18 – Prediction diagram of EEMD combined prediction model

The prediction results of the combined prediction model formed by combining the VMD method with the prediction model are shown in the following Figure 19:

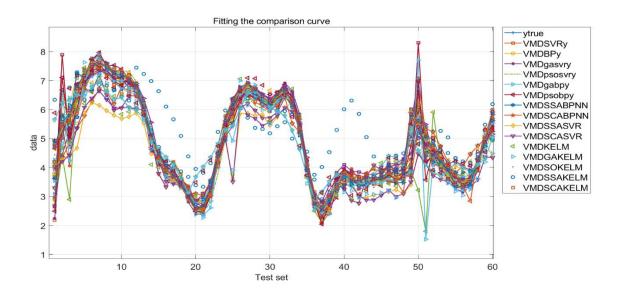


Figure 19 – Prediction diagram of VMD combined prediction model

The final prediction results of the prediction model formed by combining the Wavelet Decomposition method with the prediction model are shown in the following Figure 20:

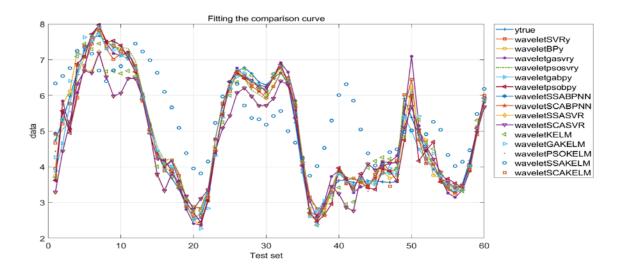


Figure 20 – Prediction diagram of Wavelet Decomposition combined prediction model

2.4 Evaluation indicators

After predicting the model, it is necessary to analyze the error of the model to measure the prediction accuracy of the model. In this paper, the prediction accuracy of the combined prediction model is analyzed by using four common evaluation indexes: mean absolute percentage error, mean square error, mean absolute error, and root mean square error. Through them, quantitative error analysis is carried out. The evaluation indexes range from $[0, +\infty)$, and when the error between the predicted value and the true value is infinitely closer to 0, it means that the prediction accuracy

of the model is closer to the perfect combination prediction model. Among them, the definitions and formulas of the four evaluation indicators are shown below:

$$MAPE = \frac{100\%}{n} \sum_{i=1}^{n} \left| \frac{\hat{y}_i - y_i}{y_i} \right|$$
 (14)

$$MSE = \frac{1}{n} \sum_{i=1}^{n} (\hat{y}_i - y_i)^2$$
 (15)

$$MAE = \sum_{i=1}^{n} |\hat{y}_i - y_i|$$
 (16)

$$RMSE = \sqrt{\frac{1}{n} \sum_{i=1}^{n} (\hat{y}_i - y_i)^2}$$
 (17)

In the formula: n represents the number of test samples, y_i and y_i represent the true and predicted values at moment i-th.

2.5 Error analysis

After the combination prediction model predicts the original data, it is necessary to analyze and compare the prediction results of the combination prediction model according to the evaluation index to obtain the optimal combination prediction model.

The error table of the combined prediction model formed by combining the EMD method with the basic prediction model is shown in the following Table 1:

Table 1 Error analysis of EMD combined prediction method

Model\Index	MAPE(%)	MSE(m/s)	MAE(m/s)	RMSE(m/s)
emdBPNN	5.44	0.1784	0.2469	0.4223
emdGABPNN	6.11	0.2244	0.2863	0.4738
emdPSOBPNN	9.20	0.4299	0.4364	0.6557
emdSSABPNN	3.20	0.032	0.1519	0.1519
emdSCABPNN	4.29	0.0429	0.2032	0.2032
emdSVR	6.83	0.4635	0.3345	0.6808
emdGASVR	5.44	0.2561	0.2631	0.5061
emdPSOSVR	6.42	0.2408	0.3000	0.4907
emdSSASVR	8.87	0.0887	0.5958	0.5958
emdSCASVR	8.87	0.0887	0.5961	0.5961
emdKELM	7.02	0.0702	0.5061	0.5061
emdGAKELM	5.36	0.0536	0.5004	0.5004
emdPSOKELM	7.06	0.0706	0.4474	0.4474
emdSSAKELM	5.07	0.0507	0.4377	0.4377
emdSCAKELM	6.87	0.0687	0.4680	0.4680

The error table of the combined prediction model formed by combining the EEMD method with the basic prediction model is shown in the following Table 2:

Table 2 Error analysis of EEMD combined prediction method

Model\Index	MAPE(%)	MSE(m/s)	MAE(m/s)	RMSE(m/s)
eemdBPNN	5.75	0.2335	0.2752	0.4832
eemdGABPNN	6.27	0.2964	0.2938	0.5445
eemdPSOBPNN	7.00	0.3480	0.3419	0.5899
eemdSSABPNN	2.83	0.0283	0.1343	0.1343
eemdSCABPNN	0.54	0.0054	0.0255	0.0255
eemdSVR	6.65	0.2903	0.3245	0.5388
eemdGASVR	6.25	0.2527	0.3104	0.5027
eemdPSOSVR	6.14	0.2386	0.2926	0.4456
eemdSSASVR	9.34	0.0934	0.7235	0.7235
eemdSCASVR	12.48	0.1248	0.8973	0.8973
eemdKELM	8.19	0.0819	0.626	0.626
eemdGAKELM	6.06	0.0606	0.4897	0.4897
eemdPSOKELM	5.54	0.0554	0.4013	0.4013

Model\Index	MAPE(%)	MSE(m/s)	MAE(m/s)	RMSE(m/s)
eemdSSAKELM	11.08	0.1108	0.6324	0.6324
eemdSCAKELM	5.83	0.0583	0.4065	0.4065

The error table of the combined prediction model formed by combining the VMD method with the basic prediction model is shown in the following Table 3:

Table 3 Error analysis of VMD combined prediction method

Model\Index	MAPE(%)	MSE(m/s)	MAE(m/s)	RMSE(m/s)
VMDBPNN	1.89	0.0127	0.0820	0.1126
VMDGABPNN	1.99	0.0110	0.0852	0.1048
VMDPSOBPNN	2.17	0.0158	0.0946	0.1256
VMDSSABPNN	2.54	0.0254	0.1204	0.1204
VMDSCABPNN	2.33	0.0233	0.1106	0.1106
VMDSVR	5.08	0.0597	0.2058	0.2443
VMDGASVR	4.95	0.0759	0.2291	0.2755
VMDPSOSVR	3.04	0.1986	0.1419	0.1737
VMDSSASVR	7.25	0.1225	0.642	0.642

Model\Index	MAPE(%)	MSE(m/s)	MAE(m/s)	RMSE(m/s)
VMDSCASVR	7.27	0.1227	0.6417	0.6417
VMDKELM	10.62	0.0702	0.5061	0.5061
VMDGAKELM	5.36	0.0536	0.5004	0.5004
VMDPSOKELM	7.06	0.0706	0.4474	0.4474
VMDSSAKELM	5.07	0.0507	0.4377	0.4377
VMDSCAKELM	6.87	0.0687	04.68	0.468

The error table for the combined prediction model formed by combining the Wavelet Decomposition method with the basic prediction model is shown in the following Table 4:

Table 4 Error analysis of Wavelet combined prediction method

Model\Index	MAPE(%)	MSE(m/s)	MAE(m/s)	RMSE(m/s)
waveletBPNN	3.57	0.0357	0.2555	0.2555
waveletGABPNN	3.47	0.0347	0.2148	0.2148
waveletPSOBPNN	5.08	0.0508	0.3403	0.3403
waveletSSABPNN	4.17	0.0417	0.1978	0.1978

Continuation of Table 4

Model\Index	MAPE(%)	MSE(m/s)	MAE(m/s)	RMSE(m/s)
waveletSCABPNN	3.61	0.0361	0.1711	0.1711
waveletSVR	5.57	0.057	0.3358	0.3358
waveletGASVR	5.85	0.0585	0.3979	0.3979
waveletPSOSVR	4.80	0.048	0.2651	0.2651
waveletSSASVR	8.67	0.0867	0.5454	0.5454
waveletSCASVR	8.79	0.0873	0.5463	0.5463
waveletKELM	6.65	0.0665	0.3883	0.3883
waveletGAKELM	4.16	0.0416	0.246	0.246
waveletPSOKELM	3.42	0.0342	0.1961	0.1961
waveletSSAKELM	5.08	0.0508	0.1993	0.1993
waveletSCAKELM	3.3	0.032	0.0906	0.0906

In order to make a more intuitive comparison and analysis of the prediction error results of these 60 different sets of combined prediction models, the Table 5 obtained by collation is shown below:

Table 5 Error analysis of all combined prediction method

EMD/EEMD/VMD/Wavelet	MAPE(%)	MSE(m/s)	MAE(m/s)	RMSE(m/s)
BPNN	5.44/5.75/1.89/3.57	0.1784/0.2335/0.0127/0.0357	0.2469/0.2752/0.0820/0.2555	0.4223/0.4832/0.1126/0.2555
GABPNN	6.11/6.27/1.99/3.47	0.2244/0.2964/0.0110/0.0347	0.2863/0.2938/0.0852/0.2148	0.4738/0.5445/0.1048/0.2148
PSOBPNN	9.20/7.00/2.17/5.08	0.4299/0.3480/0.0158/0.0508	0.4364/0.3419/0.0946/0.3403	0.6557/0.5899/ <mark>0.1256</mark> /0.3403
SSABPNN	3.20/2.83/2.54/4.17	0.032/0.0283/0.0254/0.0417	0.1519/0.1343/0.1204/0.1978	0.1519/0.1343/0.1204/0.1978
SCABPNN	4.29/0.54/2.33/3.61	0.0429/0.0054/0.0233/0.0361	0.2032/0.0255/0.1106/0.1711	0.2032/0.0255/0.1106/0.1711
SVR	6.83/6.65/5.08/5.57	0.4635/0.2903/0.0597/0.057	0.3345/0.3245/0.2058/0.3358	0.6808/0.5388/0.2443/0.3358
GASVR	5.44/6.25/4.95/5.85	0.2561/0.2527/0.0759/0.0585	0.2631/0.3104/0.2291/0.3979	0.5061/0.5027/ <mark>0.2755</mark> /0.3979
PSOSVR	6.42/6.14/3.04/4.80	0.2408/0.2386/0.1986/0.048	0.3000/0.2926/0.1419/0.2651	0.4907/0.4456/0.1737/0.2651
SSASVR	8.87/9.34/7.25/8.67	0.0887/0.0934/0.1225/0.0867	0.5958/0.7235/0.642/0.5454	0.5958/0.7235/0.642/ <mark>0.5454</mark>
SCASVR	8.87/12.48/7.27/8.79	0.0887/0.1248/0.1227/0.0873	0.5961/0.8973/0.6417/0.5463	0.5961/0.8973/0.6417/0.5463
KELM	7.02/8.19/10.62/6.65	0.0702/0.0819/0.0702/ <mark>0.0665</mark>	0.5061/0.626/0.5061/0.3883	0.5061/0.626/0.5061/ <mark>0.3883</mark>
GAKELM	5.36/6.06/5.36/4.16	0.0536/0.0606/0.0536/ <mark>0.0416</mark>	0.5004/0.4897/0.5004/0.246	0.5004/0.4897/0.5004/ <mark>0.246</mark>
PSOKELM	7.06/5.54/7.06/3.42	0.0706/0.0554/0.0706/0.0342	0.4474/0.4013/0.4474/0.1961	0.4474/0.4013/0.4474/0.1961
SSAKELM	5.07/11.08/5.07/5.08	0.0507/0.1108/0.0507/0.0508	0.4377/0.6324/0.4377/0.1993	0.4377/0.6324/0.4377/0.1993
SCAKELM	6.87/5.83/6.87/3.3	0.0687/0.0583/0.0687/0.032	0.468/0.4065/0.468/0.0906	0.468/0.4065/0.468/0.0906

Through the analysis of the above experimental data, the following conclusions can be obtained:

1) Among the combined prediction models formed by different decomposition methods, through the overall comparison, it can be found that the error of the combined prediction model formed by the combination of VMD methods is much smaller in BPNN and SVR models, indicating that in BPNN and most SVR (SVR, GASVR, PSOSVR) models, the combined prediction model formed by VMD has better prediction effect. The research also shows that the

combined prediction model formed by VMD overcomes the modal confusion caused by EMD method and solves the residual noise caused by EEMD.

- 2) In contrast, the combined prediction model formed by Wavelet presents a smaller experimental error in comparison with the KELM prediction model and part of SVR (SCASVR, SSASVR), indicating that the combined prediction model formed by Wavelet has a better prediction effect in the model of KELM. It also shows that Wavelet is also very good at overcoming the problems arising from the EMD method and the EEMD method.
- 3) Through the overall vertical comparison of the models, it can be found that although there are some deviations in different models, in general, the prediction model using SCA optimization algorithm shows more accurate prediction ability in BPNN prediction model. In the SVR prediction model, the prediction model using PSO optimization algorithm generally has more accurate prediction ability. Similarly, in the KELM model, the prediction model optimized by SCA optimization algorithm shows a more accurate prediction.
- 4) Through experimental analysis, it can be concluded that it is very undesirable to simply infer that a method must be better in theory, so as to ignore the actual situation.

INTRODUCTION OF CODE PROGRAM

In this section, the main program of the program as well as the drawing program will be introduced. The main programming tool used is Matlab.

In the main program, different decomposition models and different prediction models are selected for prediction, and then they are combined into a combined prediction model. Then the prediction accuracy is compared according to the evaluation index. The main program mainly includes three parts: parameter initialization, data decomposition and model selection.

The program code for the initialization of the parameters is shown below:

```
    %% Parameter initialization
    interval = 1;
    start_train = 1;
    lag = 4;
    data = wave;
    end_train = floor(0.8*length(data));
    [row col] = size(data);
    Y = [];
    mape = [];
    dstat = [];
    rmse = [];
```

Code Sample 8 – Parameter initialization in the main program

After initializing the parameters, a choice of different decomposition methods needs to be made. The procedure for data decomposition is shown below. The decomposition method model that can be chosen is indicated after %.

```
    %% Data decomposition
    decomp_type = 'VMD'; %emd,eemd,VMD, wavelet
    data_sample = data;
    [Comp] = tm_decomposition_method(decomp_type,data_sample);
```

Code Sample 9 – Data decomposition in the main program.

Then select different prediction models and combine them with decomposition methods to form a combined prediction model to predict and analyze data. The code is shown below. where the model's after % indicate the predictor models that can be selected.

```
    %% Selecting a predictive model
    Model = 'SVR';%SVR,BP,ga_svr,ga_bp,pso_svr,pso_bp;;;
    %KELM,PSO_KELM;SSA_KELM,SCA_KELM,GA_KELM, SSA_BPNN,SCA_BPNN,SSA_SVR,SCA_SVR
    [MAPE,RMSE,Dstat,MSE,MAE,Sum_t_all]

            emd_family_prediction(Comp,data_sample,Model);

    mz= sprintf([decomp_type '_' Model]);
    save(mz, 'Sum_t_all')
```

Code Sample 10 – Prediction model selection in the main program.

Then, by running the main program, the program will analyze the combined prediction model around the 4 set evaluation indicators. Then it will show the analysis result of the program on the right side. It will also load the data from the run into the taskbar on the left. This is shown in the Figure 21 below:

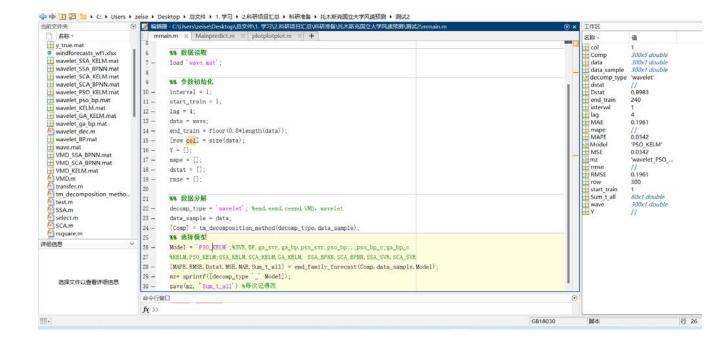


Figure 21 – The result of the main program

After all the combined prediction models are run, the analysis diagram of the combined prediction model can be drawn according to the running results. The drawing code is divided into two main parts. The first part loads the data for the run, and the second part is the plotting function. Taking VMD combination prediction method as an example.

The code to load the data is shown below:

```
1. clc
2. clear all
3. %%
4. load y_true.mat
5. y_true = s;
6. load VMD_SVR.mat
7. VMD_SVR = Sum_t_all;
8. load VMD_BP.mat
9. VMD_BP = Sum_t_all;
10. ...
```

Code Sample 11 - Loading data

The analysis diagram code for drawing the combined prediction model is shown below:

Code Sample 12 – Plotting of combined predictive models

Then, by running the plotting function, the following analysis diagram of the combined prediction model (Figure 22) is obtained.

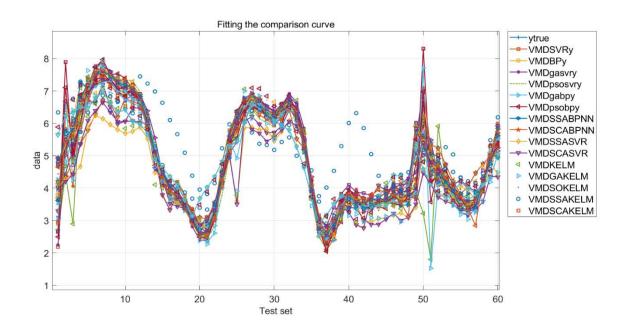


Figure 22 – Prediction diagram sample of VMD combined prediction model

CONCLUSION

In this paper, BPNN, SVR and KELM are used as three groups of basic prediction models. GA, PSO, SSA, SCA and other optimization algorithms are used to optimize the basic prediction model, and then four different decomposition methods are used to form different combined prediction models, and each combined prediction model is compared and analyzed.

The analysis of experimental data shows that

- The combined prediction model consisting of VMD method has higher prediction accuracy in the BPNN and most of SVR models compared with the combined prediction model formed by other decomposition methods. The accuracy of the combined prediction model formed by the Wavelet Decomposition method in the KELM model and part of SVR model is higher. It also shows that both the VMD method and the Wavelet Decomposition method can overcome the modal confounding problem generated by the EMD and the noise residual problem generated by the EEMD.
- Through the overall vertical comparison of the models, it can be found that although there are some deviations in different models, in general, the prediction model using SCA optimization algorithm shows more accurate prediction ability in BPNN prediction model. In the SVR prediction model, the prediction model using PSO optimization algorithm generally has more accurate prediction ability.

Similarly, in the KELM model, the prediction model optimized by SCA optimization algorithm shows a more accurate prediction.

- Through experimental analysis, it can be concluded that it is very undesirable to infer that a certain method is better only through theoretical analysis, while ignoring the actual situation. In the selection of prediction methods, more appropriate combinations should be selected according to different actual conditions to improve the prediction accuracy.
- Prospects and applications: This combined prediction model can be used not only for wind speed prediction, but also for other predictions, such as financial risk prediction models, etc.

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APPENDIX A

Part A Optimize algorithms for specific code implementation

1). Genetic Algorithm

Code Sample 2³⁸: Genetic Algorithm of selection: Selection of evolved individuals using the Roulette Wheel Selection method.

```
    for i=1:population_size

        r = rand * fitness_sum(population_size);
        first = 1;
3.
        last = population_size;
5.
        mid = round((last+first)/2);
6.
        idx = -1;
7.
        % Individual selection by the middle row method
9.
        while (first <= last) && (idx == -1)</pre>
             if r > fitness_sum(mid)
10.
11.
                 first = mid;
             elseif r < fitness_sum(mid)</pre>
12.
                 last = mid;
13.
             else
14.
                 idx = mid;
15.
                 break;
16.
17.
             end
             mid = round((last+first)/2);
18.
19.
             if (last - first) == 1
                 idx = last;
20.
                 break;
21.
22.
             end
23.
        end
24.
25.
       \% Generating a \textcolor{red}{\textbf{new}} generation of individuals
26.
       for j=1:chromosome_size
             population_new(i,j) = population(idx,j);
27.
28.
       end
29. end
30. for i=1:population_size
       for j=1:chromosome_size
32.
           population(i,j) = population_new(i,j);
```

```
33. end
34. end
```

Code sample 2³⁸ – matlab code implementation of selection in Genetic Algorithm

Code Sample 3³⁸: Genetic Algorithm of crossover.

```
1. for i=1:2:population_size
        % Crossover manipulation of chromosome strings of two individuals
3.
        if(rand < cross_rate)</pre>
            cross_position = round(rand * chromosome_size);
4.
5.
            if (cross_position == 0 || cross_position == 1)
6.
                continue;
7.
            end
8.
            % Swap binary strings for cross_position and beyond
9.
            for j=cross_position:chromosome_size
10.
                temp = population(i,j);
                population(i,j) = population(i+1,j);
11.
                population(i+1,j) = temp;
12.
13.
            end
14.
        end
15. end
```

Code sample 3³⁸ – Matlab code implementation of crossover in Genetic Algorithm

Code Sample 4³⁸: Genetic Algorithm of mutation.

```
1. for i=1:population_size
2.    if rand < mutate_rate
3.        mutate_position = round(rand*chromosome_size);
4.        if mutate_position == 0
5.             continue;
6.        end
7.        population(i,mutate_position) = 1 - population(i, mutate_position);
8.        end
9.    end</pre>
```

Code sample 4³⁸ – Matlab code implementation of Mutation in Genetic Algorithm

2). Particle Swarm Optimization

Code Sample of Particle Swarm Optimization algorithm ⁴³:

```
1.
      %% I. Emptying the environment
2. clc
3. clear all
4.
5. %% II. Plot the objective function curve
6. x = 1:0.01:2;
7. y = \sin(10*pi*x) ./ x;
8. figure
9. plot(x, y)
10. hold on
11.
12. %% III. Parameter initialization
13. c1 = 1.49445;
14. c2 = 1.49445;
15. maxgen = 50; % Number of evolutions
16. sizepop = 10;  % Population size
17. Vmax = 0.5; %Range of speed.
18. Vmin = -0.5;
19. popmax = 2; %Individual range of variation
20. popmin = 1;
21.
22.
23. %% IV. Generate initial particles and velocities
24. for i = 1:sizepop
        pop(i,:) = (rands(1) + 1) / 2 + 1;
26.
        V(i,:) = 0.5 * rands(1);
27.
        fitness(i) = fun(pop(i,:));
28. end
29.
31. %% V. Individual and group extremes
32. [bestfitness bestindex] = max(fitness);
33. zbest = pop(bestindex,:);
34. gbest = pop;
35. fitnessgbest = fitness;
36. fitnesszbest = bestfitness;
37.
38.
39. % VI. Iterative Optimization Search
```

```
40. for i = 1:maxgen
41. for j = 1:sizepop
            V(j,:) = V(j,:) + c1*rand*(gbest(j,:) - pop(j,:)) + c2*rand*(zbest - po
   p(j,:));
43.
            V(j,find(V(j,:)>Vmax)) = Vmax;
44.
            V(j,find(V(j,:)<Vmin)) = Vmin;</pre>
45.
46.
            pop(j,:) = pop(j,:) + V(j,:);
47.
            pop(j,find(pop(j,:)>popmax)) = popmax;
            pop(j,find(pop(j,:)<popmin)) = popmin;</pre>
48.
49.
50.
            fitness(j) = fun(pop(j,:));
51.
        end
52.
        for j = 1:sizepop
53.
54.
            if fitness(j) > fitnessgbest(j)
55.
                gbest(j,:) = pop(j,:);
                fitnessgbest(j) = fitness(j);
56.
            end
57.
58.
59.
            if fitness(j) > fitnesszbest
                zbest = pop(j,:);
60.
61.
                fitnesszbest = fitness(j);
62.
            end
        end
63.
64.
        yy(i) = fitnesszbest;
65. end
```

Code sample 5⁴³ – Matlab code implementation of PSO algorithm

3). Sparrow Search Algorithm

Code sample of Sparrow Search Algorithm⁴⁷:

```
    clear all;
    close all;
    clc;
    % Parameter Setting
    N=30;
    dim=2;
    N_discoverer=0.7*N;
    N_Followers=0.1*N;
    N_Vigilant=0.2*N;
```

```
11. Max_iter=100;
12. ST=0.6;
13.
14. %% Test Functions
15. f=@(x) sum(x.^2);
16. ub=10;
17. lb=-10;
18.
19. % Initialization
20. x=lb+rand(N,dim).*(ub-lb);
21. for i=1:N
22.
       fitness(i)=f(x(i,:));
23. end
24. [A,index]=sort(fitness);
25. x best=x(index(1),:);
26. x_worst=x(index(end),:);
27. best_fitness=A(1);
28. worst_fitness=A(end);
29. x best currently=x(index(1),:);
30. x_worst_currently=x(index(end),:);
31. best_fitness_currently=A(1);
32. worst_fitness_currently=A(end);
33. x_discoverer=x(index(1:N_discoverer),:);
34. x Followers=x(index(N discoverer+1:N discoverer+N Followers),:);
35. x_Vigilant=x(index(N_discoverer+N_Followers+1:N),:);
36. B=[-1,1];
37. F=best_fitness;
38. iter=1;
39.
40. %% Start of iterative updates
41. while iter<Max iter
       for i=1:dim
42.
           C(i)=B(round(rand)+1);
43.
44.
       end
45.
       A=C'*inv((C*C'));
46.
       R2=rand;
47.
       for i=1:N_discoverer
            for j=1:dim
48.
49.
                if R2<ST
                    x_discoverer(i,j)=x_discoverer(i,j)*exp(-i/rand*Max_iter);
50.
51.
                else
52.
                     x_discoverer(i,j)=x_discoverer(i,j)+randn;
53.
                end
54.
            end
```

```
55.
            ub_flag=x_discoverer(i,:)>ub;
            lb_flag=x_discoverer(i,:)<lb;</pre>
56.
57.
            x discoverer(i,:)=(x discoverer(i,:).*(~(ub flag+lb flag)))+ub.*ub flag+
   lb.*lb_flag;
58.
       end
59.
        for i=1:N_Followers
            for j=1:dim
60.
                if i>N/2
61.
62.
                    x_Followers(i,j)=rand*exp((x_worst_currently(j)-
   x Followers(i,j))/i^2);
63.
                else
64.
                    x_Followers(i,j)=x_discoverer(1,j)+abs(x_Followers(i,j)-
   x_discoverer(1,j))*A(j);
65.
                end
66.
            end
67.
            ub_flag=x_Followers(i,:)>ub;
68.
            lb_flag=x_Followers(i,:)<lb;</pre>
            x_Followers(i,:)=(x_Followers(i,:).*(~(ub_flag+lb_flag)))+ub.*ub_flag+lb
69.
    .*lb flag;
70.
       end
        for i=1:N_Vigilant
71.
72.
            for j=1:dim
73.
                if f(x_Vigilant(i,:))~=best_fitness_currently
74.
                    x Vigilant(i,j)=x best currently(j)+randn*abs(x Vigilant(i,j)-
   x_best_currently(j));
75.
                else
76.
                    x_Vigilant(i,j)=x_Vigilant(i,j)+B(round(rand)+1)*(abs(x_Vigilant
   (i,j)-x_worst_currently(j)))/abs(f(x_Vigilant(i,:))-
   worst_fitness_currently)+1;
77.
                end
78.
            end
79.
            ub_flag=x_Vigilant(i,:)>ub;
80.
            lb_flag=x_Vigilant(i,:)<lb;</pre>
            x_Vigilant(i,:)=(x_Vigilant(i,:).*(~(ub_flag+lb_flag)))+ub.*ub_flag+lb.*
81.
   lb_flag;
82.
       end
       x=[x_discoverer;x_Followers;x_Vigilant];
83.
        for i=1:N
84.
            fitness(i)=f(x(i,:));
85.
86.
        end
        [E,index]=sort(fitness);
87.
88.
        if f(x(index(1),:)) < best_fitness</pre>
89.
            best_fitness=f(x(index(1),:));
90.
            x_best=x(index(1),:);
```

```
91.
       end
92.
       if f(x(index(end),:))>worst_fitness
93.
            worst_fitness= f(x(index(end),:));
94.
            x_worst=x(index(end),:);
95.
       end
96.
       x_best_currently=x(index(1),:);
       x_worst_currently=x(index(end),:);
97.
       best_fitness_currently=E(1);
98.
       worst_fitness_currently=E(end);
99.
          x_discoverer=x(index(1:N_discoverer),:);
100.
101.
          x_Followers=x(index(N_discoverer+1:N_discoverer+N_Followers),:);
102.
          x_Vigilant=x(index(N_discoverer+N_Followers+1:N),:);
103.
          F=[F,best_fitness];
104.
          iter=iter+1;
```

Code sample 6⁴⁷ – Matlab code implementation of SSA algorithm

4). Sine Cosine Algorithm

Code sample of Sine Cosine Algorithm⁵⁰:

```
1.
      clear all
2.
  clc
3.
4. SearchAgents_no=30; % Number of search agents
5.
6. Function_name='F1'; % Name of the test function that can be from F1 to F23 (Tabl
   e 1,2,3 in the paper)
7.
8. Max_iteration=1000; % Maximum numbef of iterations
10. % Load details of the selected benchmark function
11. [lb,ub,dim,fobj]=Get_Functions_details(Function_name);

    [Best_score, Best_pos,cg_curve]=SCA(SearchAgents_no,Max_iteration,lb,ub,dim,fobj)

14.
15. figure('Position',[284
                                   660
                                         290])
                             214
16. %Draw search space
17. subplot(1,2,1);
18. func_plot(Function_name);
19. title('Test function')
20. xlabel('x_1');
```

```
21. ylabel('x_2');
22. zlabel([Function_name,'( x_1 , x_2 )'])
23. grid off
24.
25. %Draw objective space
26. subplot(1,2,2);
27. semilogy(cg_curve, 'Color', 'b')
28. title('Convergence curve')
29. xlabel('Iteration');
30. ylabel('Best flame (score) obtained so far');
31.
32. axis tight
33. grid off
34. box on
35. legend('SCA')
36.
37. display(['The best solution obtained by SCA is : ', num2str(Best_pos)]);
38. display(['The best optimal value of the objective funciton found by SCA is : ',
   num2str(Best score)]);
39.
40. function [Destination_fitness,Destination_position,Convergence_curve]=SCA(N,Max_
   iteration,lb,ub,dim,fobj)
41.
42. display('SCA is optimizing your problem');
44. %Initialize the set of random solutions
45. X=initialization(N,dim,ub,lb);
47. Destination position=zeros(1,dim);
48. Destination_fitness=inf;
49.
50. Convergence_curve=zeros(1,Max_iteration);
51. Objective_values = zeros(1,size(X,1));
53. % Calculate the fitness of the first set and find the best one
54. for i=1:size(X,1)
       Objective_values(1,i)=fobj(X(i,:));
55.
       if i==1
56.
            Destination_position=X(i,:);
57.
            Destination_fitness=Objective_values(1,i);
58.
       elseif Objective_values(1,i)<Destination_fitness</pre>
59.
60.
            Destination_position=X(i,:);
61.
           Destination_fitness=Objective_values(1,i);
62.
       end
```

```
63.
                         All_objective_values(1,i)=Objective_values(1,i);
64.
65. end
66.
67. %Main loop
68. t=2; % start from the second iteration since the first iteration was dedicated t
            o calculating the fitness
69. while t<=Max_iteration
70.
71.
                         % Eq. (3.4)
72.
                         a = 2;
73.
                         Max_iteration = Max_iteration;
74.
                          r1=a-t*((a)/Max_iteration); % r1 decreases linearly from a to 0
75.
                         % Update the position of solutions with respect to destination
76.
77.
                          for i=1:size(X,1) % in i-th solution
                                        for j=1:size(X,2) % in j-th dimension
78.
79.
80.
                                                     % Update r2, r3, and r4 for Eq. (3.3)
                                                      r2=(2*pi)*rand();
81.
82.
                                                      r3=2*rand;
83.
                                                      r4=rand();
84.
85.
                                                     % Eq. (3.3)
                                                      if r4<0.5
86.
87.
                                                                   % Eq. (3.1)
88.
                                                                   X(i,j)=X(i,j)+(r1*sin(r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position(j)-r2)*abs(r3*Destination_position_position(j)-r2)*abs(r3*Destination_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_position_positio
           X(i,j));
89.
                                                      else
90.
                                                                    % Eq. (3.2)
                                                                    X(i,j)=X(i,j)+(r1*cos(r2)*abs(r3*Destination_position(j)-
91.
            X(i,j)));
92.
93.
94.
                                        end
95.
                          end
```

Code sample 7⁵⁰ – Matlab code implementation of SCA algorithm's main loop and optimal solution



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