

An Improved Identification Method for Multivariable System

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Abstract: Due to many classical identification methods cannot be directly used for closed-loop control system, an improved identification method is proposed to simultaneously identify model parameters and the structure. The improved identification method used the genetic algorithm to estimate the initial search scope for the PSO algorithm, and then used the search result as the initial value of the Rosenbrock algorithm. On the basis the genetic algorithm to estimate is introduced to provide the rough initial search scope for the presented algorithm to improve the validity and accuracy. Simulation results show that compare with the PSO algorithm, the inertia weight variation PSO algorithm and the PSO-SQP algorithm proposed by Qibing Jin et al, the presented algorithm improves the optimizing efficiency of the particle swarm.

Keywords: initial neighborhood, PSO-R, closed loop identification, multivariable identification

1 Introduction

To obtain object dynamic mathematical model is the basis of advanced control. In order to gain an accurate process model, many researchers proposed different identification methods. But these methods depend largely on the collection data, but if working under strong noise disturbance and insufficiency response condition, it may generate a large estimate bias.

In the open loop condition, it only needs persistent excitation, the system can always be identified. But in industrial control, in consideration of the factors of stability, security and economy, open loop identification always not be allowed, so closed loop identification is necessary needed, especially in the system is unstable in open loop or the system contains feedback mechanisms condition [1,2]. Conventional identification methods such as Least Squares estimate method and Maximum Likelihood method [3]. These methods can obtain accuracy results in open loop system. However, when these methods directly apply to closed-loop system may generate large estimation deviation, even lead to unidentified. For closed loop system, searching method is a very effective identification method. Pan proposed the nonlinear stochastic (NLJ) based on LJ [4]. NLJ method improved the convergence speed of the search. But the

identification results depend largely on the choice of initial parameters.

Particle swarm optimization (PSO) is an evolutionary computation technique developed by Dr. Eberhart and Dr. Kennedy in 1995 [5], inspired by social behavior of bird flocking. Compare with conventional identification methods, PSO method has many advantages such as simple computation, rapid convergence capability and without any requirement for the input and output data. PSO has been extended to many fields [6,7,8]. But, in practical application, PSO method has the limitations of converging to undesired local solution or premature convergence [9]. Later many improved PSO methods are proposed to solve these problems. Eberhart proposed a discrete binary PSO method which can limit dimensional position [10]. Shi proposed a linear decreasing weight PSO method which can improve the ability of local search by modifying weight value [11]. Clerc introduced constriction factor to improve the ability of local search [12]. These methods be restricted to change learning factors and weight values and cannot meet the requirements for identification accuracy in modern industrial control.

In this paper we present a improve method to estimate model parameter and the structure simultaneously. The method (Named PSO-R) combines the global search

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capability of PSO and exactly local optimization of Rosenbrock. To improve the algorithm performance, we introduce genetic algorithm to estimate the rough search scope of PSO, this step can avoid the local search trap or premature convergence. Compare with the basic particle swarm optimization algorithm, Inertia Weight Variation PSO algorithm and PSO-SQP [13] proposed by Qibing Jin, Through simulation can be seen that the PSO-R method is successfully applied to multivariable system identification and achieves better results in the experiment.

2 Preliminaries: the classic optimization algorithms - PSO and Rosenbrock

2.1 The main aspects of basic particle swarm optimization (PSO)

The mathematical description of Particle Swarm Optimization is: within a d dimension search place, each particle is looked as a node in place. Each particle depends on three vectors: a position vector $x_i = (x_{i,1}, x_{i,2}, \dots, x_{i,d})$ a velocity vector $v_i = (v_{i,1}, v_{i,2}, \dots, v_{i,d})$ and its experience vector $p_i = (p_{i,1}, p_{i,2}, \dots, p_{i,d})$. For each iteration, each particle according to the formulas below updated its speed and position:

$$v_{i,j}(t+1) = wv_{i,j}(t) + c_1r_1[p_{i,j} - x_{i,j}(t)] + c_2r_2[p_{g,j} - x_{i,j}(t)] \quad (1)$$

$$x_{i,j}(t+1) = x_{i,j}(t) + v_{i,j}(t+1) \quad (2)$$

where $v_{i,j}(t+1)$ and $x_{i,j}(t+1)$ are the current velocity and the current position, $v_{i,j}(t)$ and $x_{i,j}(t)$ are the previous velocity and the previous position. The learning factors c_1 and c_2 are set constant value, normally c_1 and c_2 are taken as 2, r_1 and r_2 are random numbers between $[0, 1]$, w is inertia weight which used to control the influence of previous velocity on the current velocity. The inertia weight w is a very important parameter in PSO algorithm and could be used to control algorithm exploration ability and exploitation ability. The larger inertia weight w is helpful to enhance the global search ability of algorithm and jump out the local optimum. In the later phase of the PSO algorithm, the smaller inertia weight is helpful to improve the local search ability and make the algorithm converge.

2.2 The basic principle of Rosenbrock algorithm

Rosenbrock algorithm is a method to solve unconstrained multiple optimization problem. The description of the

problem can use the minimum value of object function as follow:

$$\min f(x), x \in R^n \quad (3)$$

The basic principle of Rosenbrock algorithm is: construct n orthogonal vectors at the current position and then search in each direction, find the direction of the maximum function value decrease and move one step, reconstruct n orthogonal vectors at the new position and repetitive operation above.

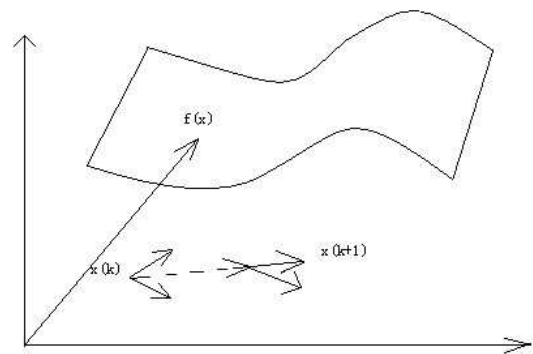


Fig. 1: The Rosenbrock method diagram

The algorithm is depicted as follow:

Step 1: Set initial estimated value $x^{(0)}$, $\{d^0, d^1, \dots, d^n\}$ are n initial standard orthogonal vectors, the initial step length is $\delta^0 = (\delta_1^0, \delta_2^0, \dots, \delta_n^0)^T > 0$, and set the parameters α, β and accuracy $\epsilon > 0$, set $k = 0$;

Step 2: Make $y = x^k$;

Step 3: Make y as the basic point, move axially parallel to $\{d^0, d^1, \dots, d^n\}$,

If $f(y + \delta_j^k d^j) \leq f(y)$, then

$$y = y + \delta_j^k d^j, \delta_j^k = \alpha \delta_j^k \quad (4)$$

If $y = y + \delta_j^k d^j, \delta_j^k = \alpha \delta_j^k$, then

$$y = y, \delta_j^k = -\beta \delta_j^k \quad (5)$$

Until failures appear in n directions;

Step 4: Set $x^{k+1} = y$, if $\|x^{k+1} - x^k\| \leq \epsilon$, then stop the iteration, the output is x^{k+1} , or go to the next step;

Step 5: Set $\tilde{d} = x^{k+1} - x^k$, build new orthogonal vectors $\{d_0, d_1, \dots, d_n\}_{k+1}$, set $\delta^{k+1} = \delta^0, k = k + 1$, turn to step 3.

According to the experience, $\alpha \in [2, 3]$, $\beta \approx 0.5$, in the initial stage orthogonal vectors can set as unit vectors.

3 PSO-R algorithm

3.1 Initial neighborhood

Most optimization algorithms exist this problem: search algorithm itself has the randomness and will cause blind inefficient in optimization process. In this paper, we use the genetic algorithm to estimate the rough range search space of PSO algorithm, improves the convergence and accuracy of search process, avoid randomness factor of the PSO algorithm cause the local search trap or premature convergence.

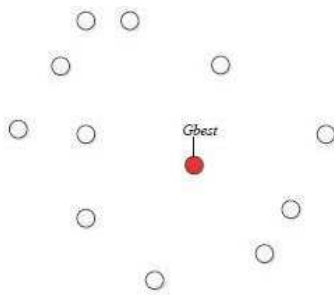


Fig. 2: Chart of Standard PSO algorithm searching process

Assume the red dot is the global optimum, the PSO algorithm random generate N nodes in the whole search scope, and then according the formulas of PSO algorithm to calculate the optimum.

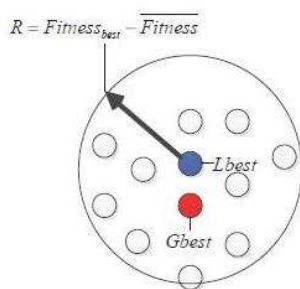


Fig. 3: Chart of Basing on the initial neighborhood of PSO algorithm optimization process

Assume the red dot is the global optimum, and the blue dot is the local optimum. The initial neighborhood in this paper, the center of initial value can be obtain by genetic algorithm. The radius R of initial neighborhood is the difference of the optimal value and average optimal value of GA algorithm, this constitutes the initial neighborhood of PSO algorithm.

3.2 PSO-R Algorithm

- Step 1: Using the GA algorithm to estimate the initial neighborhood of the PSO algorithm.
- Step 2: Initialized parameters in PSO algorithm, including initial search position and particle speed.
- Step 3: Comparison fitness value, if current fitness evaluation is better than the previous, let the current replace the previous.
- Step 4: Update the velocity of each particle and the coordinate position.
- Step 5: Judging the iteration number whether satisfy the set value, if satisfy according the formula 3 to update the position and speed of particle, and go to the next step, otherwise go to the step 2.
- Step 6: Using the result of PSO algorithm as initial point of Rosenbrock, the results of the parameters estimation are $\hat{\theta}_q$, precision is ϵ_q , q is the cycling time. According $J = \sum_{t=0}^N (y(t) - y_e(t))^2$ to update the initial value of PSO and the cycling time.

$$\hat{\theta}_{q+1}^0 = (\hat{\theta}_{q+1}^+ + \hat{\theta}_{q+1}^-) / 2 \quad (6)$$

where

$$\begin{cases} \hat{\theta}_{q+1}^+ = \hat{\theta}_q + abs(rand(\hat{\theta}_q - \hat{\theta}_{q-1})) \\ \hat{\theta}_{q+1}^- = \hat{\theta}_q - abs(rand(\hat{\theta}_q - \hat{\theta}_{q-1})) \end{cases}$$

$$\epsilon_{q+1} = rand \times \epsilon_q \quad (7)$$

- Step 7: Gain the global optimal value $\hat{\theta}_{k,l}$, and the corresponding optimal fitness value $f(\hat{\theta}_{k,l})$, where the main iteration time in PSO algorithm is k , l is the serial number of the particle in the group.
- Step 8: Through calculate obtain the optimal value and the corresponding fitness function value. In turn down as $\hat{\theta}_r$, $f(\hat{\theta}_r)$, if $f(\hat{\theta}) < f(\hat{\theta}_r)$ let $\hat{\theta} = \hat{\theta}_r$.
- Step 9: Along the $\hat{\theta}$ axis direction to search. Searching make fitness function minimum variable value, if $f(\hat{\theta} + v_1 d_j) \leq f(\hat{\theta})$, $\hat{\theta} = \hat{\theta} + v_1 d_j$.
- Step 10: Determine whether meet the requirement $||\theta_{k+1} - \theta_k|| \leq \epsilon_q$, if satisfy, end the computation and output the optimal result θ_{k+1} or go to step 11.
- Step 11: Update orthogonal vector group for n search directions, and use to set the initialization parameter forward σ^{k+1} step, go to step 9. Figure 5 is the flow chart of the PSO-R algorithm.

4 Simulation

4.1 First order plus dead time (FOPDT) model

$$Gs = \frac{k_p}{\tau_p s + 1} e^{-\theta s}$$

Based on closed loop step test in terms of a P-type controller ($kc = 0.5$), with a step change of $h = 0.05$ to

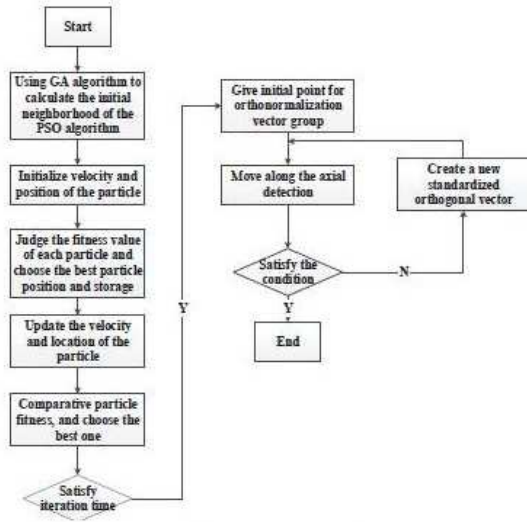


Fig. 4: Algorithm chart of PSO-R

the set point, we choose $Gs = \frac{12.8}{16.7s+1}e^{-s}$ as an example of FOPDT model. The basic PSO algorithm, the inertia weight variation PSO Algorithm, the PSO-SQP algorithm and the proposed algorithm are respectively used to estimate the selected model.

A Gaussian white noise respectively generating $NSR = 5\%$, 10% and 20% is added to the output to test the robustness.

Table 1: The results of various algorithm for FOPTD model in no noise

Actual Value	12.8	16.7	1
PSO	12.7989	16.6989	-1.0022
WPSO	12.7999	16.6982	-1.0091
PSO-SQP	12.8000	16.7000	-0.9989
PSO-R	12.8000	16.7000	-1.0000

Table 2: The results of various algorithm for FOPTD model in $NSR = 5\%$ noise

Actual Value	12.8	16.7	1
PSO	12.7841	16.6916	-0.9901
WPSO	12.8089	16.7110	-0.9598
PSO-SQP	12.7998	16.6970	-0.9972
PSO-R	12.8000	16.7007	-1.0002

Fig. 5 depicts the result of the PSO-R from step response test for FOPDT model. Table 1 to table 4 demonstrates that the search optimization method with frequency response to estimate the initials can obtain an excellent performance of FOPDT model identification in

Table 3: The results of various algorithm for FOPTD model in $NSR = 10\%$ noise

Actual Value	12.8	16.7	1
PSO	12.7928	16.7140	-0.8338
WPSO	12.8014	16.8495	-1.0572
PSO-SQP	12.8008	16.7017	-1.0012
PSO-R	12.8000	16.6987	-0.9982

Table 4: The results of various algorithm for FOPTD model in $NSR = 20\%$ noise

Actual Value	12.8	16.7	1
PSO	12.8006	16.7435	-0.8338
WPSO	12.8214	16.9926	-1.1901
PSO-SQP	12.8214	16.6928	-1.1901
PSO-R	12.8000	16.6999	-1.0060

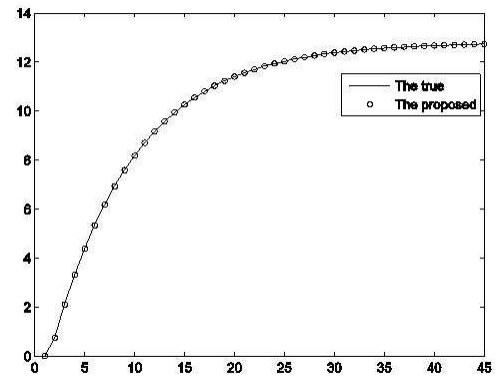


Fig. 5: Step responses test for FOPDT model estimated by PSO-R

spite of the existence of the noise and the complexity of the equivalent input; among the four algorithms, the proposed PSO-R achieves better results than PSO, WPSO and PSO-SQP in terms of accuracy and robustness.

4.2 Second order plus dead time (SOPDT)

$$\text{model } Gs = \frac{k_p}{t_1s^2+t_2s+1}e^{-\theta s}$$

A unit step is given to the SOPDT model of $Gs = \frac{1}{12s^2+8s+1}e^{-s}$ in terms of a P-type controller ($kc = 0.5$), with a step change of $h = 0.05$ to the set point. Experiments with the measurement noises of $NSR = 5\%$, 10% and 20% are respectively taken for PSO, WPSO, PSO-SQP and the proposed PSO-R with initial parameter estimate. The results are listed in Table 2.

The persistency of the excitation is a key issue for the identification of the system, whereas many methods need a persistent excitation during the whole process. In the

Table 5: The results of various algorithm for SOPTD model in no noise

Actual Value	1	12	8
PSO	1.0006	11.1810	7.9299
WPSO	0.9999	12.3526	8.0258
PSO-SQP	1.0000	12.1500	8.0049
PSO-R	1.0000	12.0080	8.0007

Table 6: The results of various algorithm for SOPTD model in NSR = 5% noise

Actual Value	1	12	8
PSO	1.0004	11.0869	7.9235
WPSO	1.0002	11.5059	7.9583
PSO-SQP	1.0000	11.7049	7.9933
PSO-R	1.0000	11.9868	7.9991

Table 7: The results of various algorithm for SOPTD model in NSR = 10% noise

Actual Value	1	12	8
PSO	1.0013	10.7953	7.9078
WPSO	1.0006	10.4549	7.8918
PSO-SQP	1.0004	11.0517	7.8997
PSO-R	1.0001	11.7738	7.9819

Table 8: The results of various algorithm for SOPTD model in NSR = 20% noise

Actual Value	1	12	8
PSO	1.0004	10.5602	7.8801
WPSO	1.0004	10.4046	7.8629
PSO-SQP	1.0002	11.0049	7.8777
PSO-R	1.0002	11.6753	7.9776

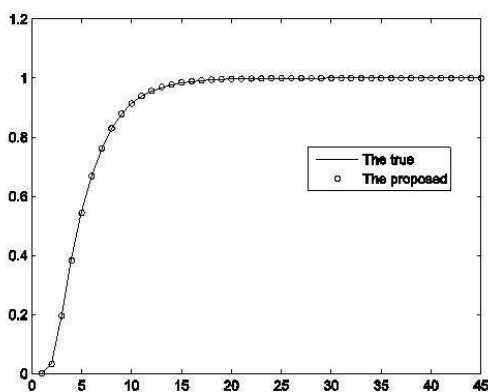


Fig. 6: Step responses test for SOPDT model estimated by PSO-R

paper, we respectively make the experiment with different excitation time and the results are as Fig. 7.

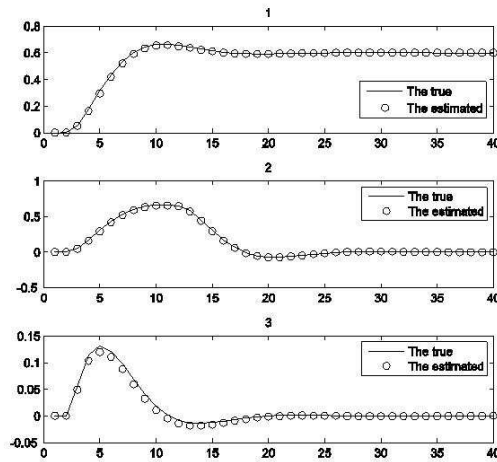


Fig. 7: The results of PSO-R with different excitation time

Subfigure 1 indicates a persistent excitation all time, Subfigure 2 depicts the excitation time $t = 10s$ and Subfigure 3 shows the excitation time $t = 1s$. The presented method is feasible for the non-sustainable excitation whereas the stability and accuracy of the results may be affected at a certain degree, hence a persistent excitation is preferable in the simulations of SISO system.

4.3 Woodberry model

Woodberry model is proposed by Wood and Berry in 1973 and widely adopted in the identification of multivariable system because of its characteristics of strong coupling and multiple time delays between the various loops. We get a good result with PSO-R in the experiment, have also made very good identification results.

Woodberry model

$$G(s) = \begin{bmatrix} \frac{12.8e^{-s}}{1+16.7s} & \frac{-18.9e^{-3s}}{1+21s} \\ \frac{6.6e^{-7s}}{1+10.9s} & \frac{-19.4e^{-3s}}{1+14.4s} \end{bmatrix}$$

For the decentralized closed loop system with $kc1 = 0.5$, $kc2 = -0.09$, the step tests are respectively taken by $r1 = 1$, $r2 = 2$ and $r1 = 4$, $r2 = 2$. Choose $G_{ci}(r_i - y_i)$ as the equivalent inputs in the simulation and results of PSO, WPSO, PSO-SQP and the proposed method are listed in Table 9 to table 12.

4.4 Multivariable model

The transfer function of system being identified is

$$y = 1 + x_1 - 1.5x_2 + 1.6e^{-x_3} + 2 \sin(1 + x_4) \quad (8)$$

Table 9: The results estimated channel $G11$ by various algorithms

Actual Value	12.8	16.7	-1
PSO	12.7929	15.6	-1.21
WPSO	12.7949	16.6193	-1.0044
PSO-SQP	12.7984	16.3140	-1.0027
PSO-R	12.8000	16.6942	-1.0014

Table 10: The results estimated channel $G12$ by various algorithms

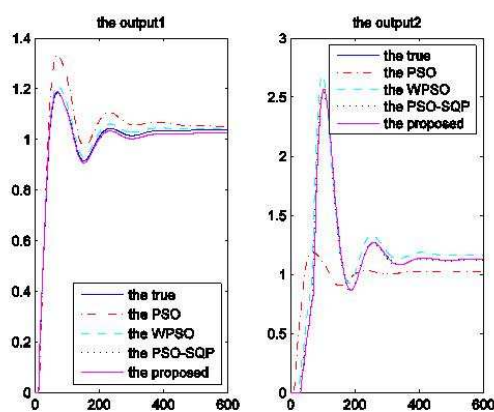
Actual Value	-18.9	21	-3
PSO	-18.7700	18.2000	-2.8500
WPSO	-18.8419	20.2100	-2.93
PSO-SQP	-18.7280	20.9420	-3.0017
PSO-R	-18.9000	21.0000	-3.0002

Table 11: The results estimated channel $G21$ by various algorithms

Actual Value	6.6	10.9	-7
PSO	6.5656	10.9218	-6.5500
WPSO	6.5766	10.9109	-6.7630
PSO-SQP	6.5856	10.9081	-6.9852
PSO-R	6.6000	10.9000	-6.9975

Table 12: The results estimated channel $G22$ by various algorithms

Actual Value	-19.4	14.4	-3
PSO	-19.8800	13.2700	-2.4000
WPSO	-19.3074	14.4339	-2.7300
PSO-SQP	-19.3942	14.4140	-3.0060
PSO-R	-19.4000	14.4001	-3.0007

**Fig. 8:** The results of the various algorithm for Woodberry model

According to formula 8 generate 30 groups sample model. Assumed we didn't know the structure of model in advanced, chosen all of the meta-model to identify the

system. Set the parameters of BPSO algorithm in identification process as follow: particle number $N = 20$, inertia weight $w = 1$, accelerate factors $C_1 = 2$, $C_2 = 1.65$, $X_{\max} = 2$, the max evolution number is 1200. Set inertia weight in WPSO algorithm as $X_{\max} = 0.9$, $X_{\min} = 0.4$. Other parameters are the same as BPSO algorithm.

After many times simulation, chosen a better group result as follow:

PSO algorithm is

$$y = 0.2584 + 1.4775x_1 - 1.9802x_2 + 1.9991e^{-1.3560x_3} + 2.1390 \sin(1.9882 + 1.0138x_4)$$

WPSO algorithm is

$$y = 0.3793 + 1.0683x_1 - 1.5760x_2 + 1.2469e^{-0.8776x_3} + 2.0376 \sin(0.9791 + 1.0013x_4)$$

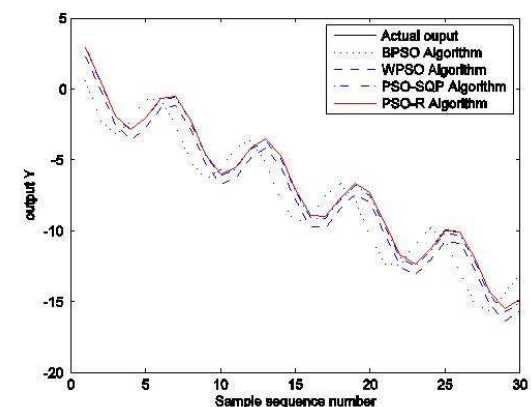
PSO-SQP algorithm is

$$y = 0.9970 + 0.9956x_1 - 1.5042x_2 + 1.5896e^{-0.9734x_3} + 2.0027 \sin(0.9988 + 1.0001x_4)$$

PSO-R algorithm is

$$y = 0.9984 + 0.9997x_1 - 1.5000x_2 + 1.5963e^{-0.9734x_3} + 2.0000 \sin(1.0000 + 1.0000x_4)$$

Identification parameters of each algorithm are shown in table 13.

**Fig. 9:** Without noise cases of various algorithm system output curve

From the contrasts above we could see, the hybrid algorithm combined particle swarm optimization algorithm and sequential quadratic programming algorithm identify system output curve is more close to the real value, precision obviously improved.

Table 13: Various algorithm system parameter identification results

Actual Value	1	1	-1.5	1.6	-1	2	1	1
BPSO	0.2584	1.4775	-1.9802	1.9991	-1.3560	2.1390	1.9882	1.0138
WPSO	0.3793	1.0683	-1.5760	1.2469	-0.8776	2.0376	0.9791	1.0013
PSO-SQP	0.9970	0.9956	-1.5042	1.5896	-0.9734	2.0027	0.9988	1.0001
PSO-R	0.9984	0.9997	-1.5000	1.5963	-0.9897	2.0000	1.0000	1.0000

5 Conclusion

In this paper, we present an improved identification method for multivariable system. The idea is to search the local optimization with Rosenbrock algorithm at the set iteration of PSO and get a better result as the best global location. With the rough search scope of PSO algorithm estimated by the genetic algorithm, the search optimization approaches are improved in convergence speed and robustness, experiment is utilized to estimate the FOPDT and SOPDT under the disturbances of different NSR use PSO, WPSO, PSO-SQP and PSO-R are respectively. The results of simulation prove that PSO-R with the advantage of the global search capability of particle swarm optimization (PSO) algorithm and exactly local optimization of Rosenbrock algorithm, is an approximate unbiased and effective identification method that can be successfully applied to the model of the closed loop identification with large noise, time delay. Finally, we also use PSO-R method to identify multivariable closed loop system, compared with other methods the method proposed by this article get a better result as well.

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