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# A FAST QUANTUM-BEHAVED PARTICLE SWARM OPTIMIZATION BASED ON ACCELERATING FACTOR

<sup>1</sup>ZHEHUANG HUANG

<sup>1</sup>School of Mathematics Sciences, Huaqiao University, Quanzhou, 362021, China Cognitive Science Department, Xiamen University, Xiamen, 361005, China E-mail: <u>hzh98@stu.xmu.edu.cn</u>

#### ABSTRACT

Quantum-behaved particle swarm optimization (QPSO) is a good optimization technique which has been successfully applied in many research and application areas. But traditional QPSO algorithm is easy to fall into local optimum and the rate of convergence is slow. To solve these problems, an improved algorithm based on dynamic adjustment of the acceleration factor is proposed. The experiments on high dimensional function optimization showed that the improved algorithm has more powerful global exploration ability and faster convergence speed.

Keywords: QPSO; Accelerating Factor; Function Optimization

#### 1. INTRODUCTION

Modern heuristic algorithms are considered as practical tools for nonlinear optimization problems. Particle swarm optimization (PSO) is a populationbased, self-adaptive search optimization technique first introduced by Kennedy and Eberhart [1, 2] in 1995. The method has been developed through a simulation of simplified social models.

PSO is a kind of algorithm based on social psychology. Like GA [3], PSO must also have a fitness evaluation function. Compared with GA, PSO has some attractive characteristics. It has memory, so knowledge of good solutions is retained by all particles [4]. But in GA, previous information is destroyed once the population changes.

The PSO method is becoming very popular due to its simplicity of implementation and ability to quickly converge to a reasonably good solution. In recent years, there has been increasing interest in developing the PSO algorithm. PSO has been applied successfully to all kinds of optimization problems such as nonlinear functions [5], neural networks [6-8], power and voltage control [9], and task assignment problem [10], etc.

This paper is organized as follows. In section 2 the background of PSO and QPSO is presented. The improved algorithm is presented in section 3. In Section 4 some experimental tests, results and conclusions are given. Section 5 concludes the paper.

#### 2. INTRODUCTION TO PSO

#### 2.1. Background

PSO is based on swarms such as fish schooling and bird flocking. According to the research results for bird flocking, birds find food by implementing an "information sharing" approach.

Each particle moves over the search space according to the historical behaviors of the particle and its companions. Suppose that the location of the ith particle is represented as  $X_i = (x_{i1}, x_{i2}, \dots x_{id})$ . The best previous position of the ith particle is represented as  $P_i = (p_{i1}, p_{i2}, \dots p_{id})$ , which is also called pbest. The location  $p_g$  is also called gbest. The velocity for the ith particle is represented as  $V_i = (v_{i1}, v_{i2}, \dots v_{id})$ .

If g is the index of the particle visited the best position in the swarm, and then  $P_g$  becomes the best solution found so far, and the velocity of the particle and its new position will be determined according to the following equations (1a) and (1b), respectively:

$$vid = w \times vid + c1 \times rand() \times (pid - xid)$$
$$+ c2 \times rand() \times (pgd - xid) \quad (1a)$$
$$xid = xid + vid \quad (1b)$$

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Where the function rand() can generate a random number between 0 and 1. w is called the inertia weight. c1 and c2 are two constant numbers, which are often called the cognitive confidence coefficients.

The procedure for implementing the PSO is given by the following steps:

Initialize the population  $x_i$ 

for each particle  $X_i$  do update the particle's best position if  $f(X_i) < f(pbi)$  then  $pbi = X_i$ end if update the global best position if f(pbi) < f(gb) then gb = pbiend if end for update particle's velocity and position for each particle  $X_i$  do for each dimension d in D do update particle's velocity by equation (1a) update particle's position by equation (1b) end for end for it = it + 1until iterations it < max

Algorithm 1. The PSO algorithm

## 2.2. Quantum Model of PSO

The development of quantum mechanics [11, 12] forced the scientists to rethink the applicability of classical mechanics and the traditional understanding of the nature.

In classical PSO algorithm, a particle is depicted by its position vector and velocity vector, which determine the trajectory of the particle. In quantum world, the term trajectory is meaningless, because position vector and velocity vector of a particle can not be determined simultaneously according to uncertainty principle. Some researchers consider a social organism is a system far more complex than that formulated by particle swarm optimization, and a linear evolvement equation is not sufficient to depict it at all. Quantum-behaved Particle Swarm Optimization (QPSO) was proposed [13, 14] in 2004. QPSO is inspired by quantum mechanics and fundamental theory of particle swarm. Wave function of position [15-17]depicts the state of the particle in quantized search space, not informing us of any certain information about the position of a particle that is vital to evaluate the fitness of a particle.

A global point denoted as mbest is introduced into PSO:

$$mbest = \frac{1}{M} \sum_{i=1}^{M} P_i$$
$$= \left(\frac{1}{M} \sum_{i=1}^{M} P_{i1}, \frac{1}{M} \sum_{i=1}^{M} P_{i2}, \cdots, \frac{1}{M} \sum_{i=1}^{M} P_{id}\right) (2)$$

The procedure of QPSO is given by the following steps:

Initialize the population  $X_i$ Do find out the mbest of the swarm by equation (2); for I=1 to population size M; if f( $p_i$ ) < f( $x_i$ ) then  $p_i = x_i$  $p_g = \min(p_i)$ for d=1 to dimension D  $f_{i1} = \operatorname{rand}(0,1), f_{i2} = \operatorname{rand}(0,1)$  $P=(f_{i1} * pid + f_{i2} * pgd)/(f_{i1} + f_{i2})$ L=beta\*abs(mbestd- $X_{id}$ ) u=rand(0,1);if rand(0,1) > 0.5;  $x_{id} = \text{p-abs(mbestd-} x_{id}) \times \log(1/u);$ else  $x_{id} = p + abs(mbestd - x_{id}) \times log(1/u);$ until the termination criterion is met.

Algorithm 2 The OPSO algorithm

## 3. THE IMPROVED ALGORITHM BASED ON ACCELERATING FACTOR (AQPSO)

From the QPSO algorithm, we can know local attractor q is determined by local optimum *pbest* and global optimum *gbest* :

$$q = \varphi \bullet pbest + (1 - \varphi) \bullet gbest ,$$
  
$$\varphi = \frac{c_1 r_1}{c_1 r_1 + c_2 r_2}, r_1 \sim U(0, 1), r_2 \sim U(0, 1)$$

From standard PSO algorithm, it is known that  $c_1$  and  $c_2$  as two accelerating factor, will not only affect the algorithm convergence speed, at the

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same time, also may lead to premature phenomenon occurs. In the early stage, if the local optimal has faster convergence speed and the global optimal has slower convergence speed, or in the late stage, the local optimal has slower convergence speed and the global optimal has faster convergence speed, can effectively avoid the happening of premature. In the early stages, the particles and the optimal location are far away. At later stage, the position of the particle and the optimal location are closer. We can use the error function represented the closer degree between particles and the global optimal position, so as to select the corresponding acceleration factor.

The particles closer to the global optimal position, the acceleration factor. Let error function as follows:

$$\Delta F = \frac{ABS(F_i - F_{gbest})}{MIN(ABS(F_i), ABS(F_{ebest}))}$$

Where  $F_i$  is the fitness of the ith particle,  $F_{gbest}$  is the fitness of global optimal,  $ABS(F_i)$  is absolute value of F, thus acceleration factor function is:

$$c1 = \\ \begin{cases} (Mc - mc) \times \frac{MAXITER - t}{MAXITER} + mc, & \Delta F > F_0 \\ (Mc - mc) \times \sin(\frac{MAXITER - t}{MAXITER}) + mc, & \Delta F < F_0 \end{cases}$$

$$\begin{cases} (Mc - mc) \times \sin(\frac{MAXITER - t}{MAXITER}) + mc, \\ (Mc - mc) \times \frac{MAXITER - t}{MAXITER} + mc, \end{cases} \quad \Delta F > F_0$$
(3) 
$$\Delta F < F_0$$

Where  $F_0 = 0.3$ .

c2 =

The procedure for implementing the AQPSO is given by the following steps:

Step 1 . Initialize all particles;

Step 2. Calculation of the value of the variable of  $\beta$ :

$$\beta = (1-0.5)*(MAXITER-t)/MAXITER+0.5$$

, where t is the number of iterations ;

Step3 . Calculating the fitness of each particle;

Step 4. Calculated error values according to the error function

$$\Delta F = \frac{f(x_i) - f(P_{gj})}{MIN(ABS(f(x_i)), ABS(f(P_{gi})))} ;$$

Step5. Calculated accelerating factor based on the acceleration factor function;

Step 6. Obtain the average optimal position:

mbest;

$$= \left(\frac{1}{N}\sum_{i=1}^{N}P_{i1}, \frac{1}{N}\sum_{i=1}^{N}P_{i2}, \frac{1}{N}\sum_{i=1}^{N}P_{i3}, \cdots, \frac{1}{N}\sum_{i=1}^{N}P_{ij}\right) ;$$

Step 7. Obtain the local attractor:

Update c1 and c2 by equation (3);

$$\varphi = \frac{c_1 r_1}{c_1 r_1 + c_2 r_2}, r_1 \sim U(0,1), r_2 \sim U(0,1)$$
$$p_{ij}(t) = \varphi_{ij}(t) \cdot P_{ij}(t) + (1 - \varphi_{ij}(t)) \cdot P_{gj}(t)$$

Step 8. update the particle's velocity and position

$$x_{ij}(t+1) = p_{ij}(t) \pm \beta \left| mbest_j(t) - x_{ij}(t) \right| \cdot \ln(1/\mu_{ij}(t))$$

Step 9. If iteration terminated, output optimal value; otherwise return 3 and continue.

#### 4. EXPERIMENTAL RESULTS

A set of unconstrained real-valued benchmark functions was used to investigate the effect of the improved algorithm. The functions are shown in Table 1.

The results are shown in Table 2. Each point is made from average values of over 30 repetitions.

For sphere function and Ackey function, AQPSO algorithm can effectively improve the accuracy such that the optimal value obtained is much closer to the theoretical one and the accuracy is improved compared with the standard QPSO algorithm. Shaffer function is a multimodal function, from the results of iteration, the convergence effect of this algorithm is not as good as the standard QPSO algorithm, but the difference is small and acceptable. On the convergence time, there is a significant improvement as expected. These experimental results show that improving algorithm can effectively improve the convergence speed with

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excellent convergence effect.

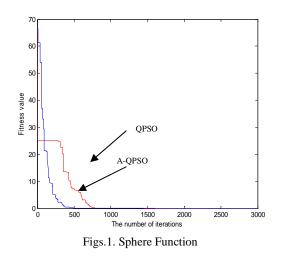
Table 1 Functions Used To Test The Effects Of AQPSO			
Function	Function expression		
Sphere function	$f_1(x) = \sum_{t=1}^n x_t^2$		
Rastrigrin function	$f_{3}(x) = \sum_{t=1}^{n} (x_{t}^{2} - 10\cos(2\pi x_{t}) + 10)$		
Griewank function	$f_2(x) = \frac{1}{4000} \sum_{t=1}^n (x_t - 100)^2 - \prod_{t=1}^n \cos(\frac{x_t - 100}{\sqrt{t}}) + 1$		
Ackey function	$f_{3}(x) = 20 + e - 20e^{-0.2\sqrt{\sum_{t=1}^{n} x_{t}^{2}}} - e^{\sum_{t=1}^{n} \cos(2\pi x_{t})}$		
Shaffer's function	$f_4(x) = 0.5 + \frac{(\sin\sqrt{x_1^2 + x_2^2})^2 - 0.5}{(1 + 0.001(x_1^2 + x_2^2))^2}$		

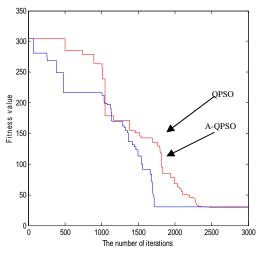
Table 2 The	Performances	Of AOPSO
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algorithm	ı OPSO		AQPSO	
	optimal	time(s)	optimal ti	me(s)
Sphere	5.29e-29	3.67	3.12e-30	3.1
Rastrigrin	23	4.08	23.8	3.27
Griewank	0	3.93	0	3.81
Ackey	2.79e-3	4.13	2.66e-3	3.84
shaffer	0	3.67	1.2e-9	3.02

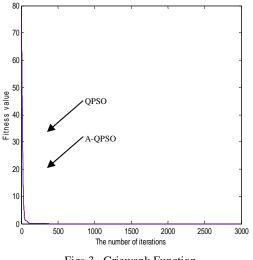
The comparison of two methods with convergent curves is shown in Figs.1. to 5.

The experiment results show the AQPSO algorithm has better result. Comparisons with QPSO, the AQPSO algorithm has both global search ability and fast convergence speed.





Figs.2. Rastrigrin Function

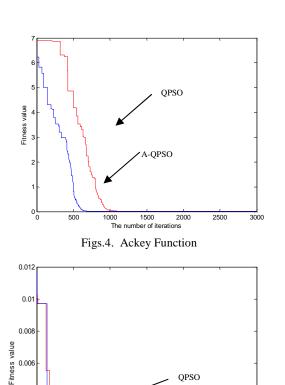


Figs.3. Griewank Function

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500

1000

0.004

0.002

0

From above, we can know the improved algorithm have more powerful global exploration ability and faster convergence speed and can be widely used in other optimization tasks.

1500

The number of iterations

Figs.5. Shaffer's Function

A-OPSO

2000

2500

3000

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