An Improved Global Particle Swarm Optimization for Faster Optimization Process

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Abstract

An efficient Global Particle Swarm Optimization (GPSO) is proposed in order to overcome the concern of trapping in the local optimal point especially in high dimensional while using ordinary Particle Swarm Optimization (PSO). GPSO is able to bring all the particles to be closely clumped together faster than PSO. In this paper, an improved GPSO is proposed in order to get a closely clumped particles group faster than using GPSO. The original GPSO is improved by taking into account the global best fitness error and particle fitness clumping size of every iteration. The improved GPSO is simulated by using several two dimension mathematical function and benchmarked with the original GPSO. The improved GPSO is shown to be able to obtain closely clumped particles much more faster than the original GPSO up to 62%. The performances are also evaluated by comparing the standard deviation, average, best particle and worst particles obtained through a 50 independent runs. In term of the four factors mentioned, the improved GPSO performance is shown to be as good of the original GPSO.

Keywords: Particle swarm optimization; global particle swarm optimization; optimization

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1.0 INTRODUCTION

PSO is a popular optimization algorithm, which is proven to be robust and possess a well global exploration ability. It has been improved into many versions to be implemented in various kind of system such as Guided PSO for facial emotion detection [1,2], Gaussian PSO for bandwidth parameter determination [3] and genetic algorithm PSO for fault diagnosis [4]. PSO has tendency of getting trapped in local minima and slow convergence [5], most of the improved PSO are proposed due to these concerns.

In [6], an efficient GPSO algorithm is proposed in order to get an optimized point without getting trapped in the local optimal point as happened with ordinary PSO. A fast optimization method is essential in systems such as system scheduling [7,8] and human motion tracking [9,10]. The GPSO is said to be able to reach the optimal results in several iterations, which is faster than PSO. In [6], GPSO is compared to the ordinary PSO, iteration PSO and evolutionary PSO algorithms and has shown to have better performance.

In this paper, an improved GPSO is proposed in order to enhance the speed of optimization. The improved GPSO is taking into the account the global best fitness error and the particle fitness clumping size of every iteration. In Section 2, the modification done on the original PSO is discussed. In Section 3, the simulation results and discussions are presented. Lastly, Section 5 concludes this paper.

2.0 METHODOLOGY

2.1 An Improved Global Particle Swarm Optimization Algorithm

In ordinary PSO, the velocity algorithm is the most important part. Most of the modification of PSO revolves around this algorithm. The ordinary PSO velocity algorithm includes information such as local best position until k-th iteration for i-th particle (p_{best-i}^k) and global best position until k-th iteration (G_{best}^k) as shown below:

\[ v_{i}^{k+1} = w v_i^k + c_1 r_1 (p_{best-i}^k - x_i^k) + c_2 r_2 (G_{best}^k - x_i^k) \] (1)

where kis the current iteration, w is the inertia weight, v_i is the velocity of i-th particle, c_1 and c_2 are the acceleration coefficients for cognitive and social components, lastly r_1 and r_2 are random numbers between 0 and 1.

In GPSO, a fourth term is added into the velocity algorithm as below:

\[ v_{i}^{k+1} = w v_i^k + c_1 r_1 (p_{best-i}^k - x_i^k) + c_2 r_2 (G_{best}^k - x_i^k) + c_3 r_3 (G_{best}^k - x_i^k) \] (2)

where c_3 r_3 (G_{best}^k - x_i^k) is named as the Improvement Factor (IF), c_3 is acceleration coefficient and r_3 is another random
numbers from 0 to 1. \( E_{\text{best}} \) is the random value selected from \( G_{\text{best}} \) for each iteration.

In the proposed improved GPSO, the forth term, IF is modified with the addition of an exponential equation, which taking into consideration the particle fitness clumping size and global best fitness error until \( k \)-th iteration as below:

\[
y_{i}^{k+1} = \frac{wv_{i}^{k} + c_{1}r_{1}(p_{\text{best} - i}^{k} - x_{i}^{k}) + c_{2}r_{2}(G_{\text{best}}^{k} - x_{i}^{k})}{e^{s_{k} + \epsilon}}
\]

(3)

The added exponential equation \( e^{-s_{k} + \epsilon} \) is named correction acceleration factor (CAF) where \( s_{k} \) is the particle fitness clumping size of \( k \)-th iteration, which calculated using algorithm below:

\[
s_{k} = \text{abs}(F_{\text{max}}^{k} - F_{\text{min}}^{k})
\]

(4)

where \( F_{\text{max}}^{k} \) is the maximum value of current iteration particle fitness and \( F_{\text{min}}^{k} \) is the minimum value of current iteration particle fitness. Meanwhile \( e_{k} \) is the global best fitness error of \( k \)-th iteration determined with the difference between the current iteration global best fitness \( (G_{k}) \) and the expected fitness \((E)\). \( E \) is determined by the desire of the programmer.

\[
e_{k} = \text{abs}(G_{k} - E)
\]

(5)

The CAF is formed in order to control the acceleration of the particles depending on the difference between the local minimum and maximum fitness of current iteration and also the difference between global best fitness with the expected fitness that is to be determined.

During the start of optimization process, the value of Equations (4) and (5) are expected to be great, which means CAF value will also be great (not greater than 1). At this moment, the particles will accelerate following the determined IF value. As the iteration goes on, the value of Equations (4) and (5) will get smaller, causing the CAF value to be decrease, eventually to zero, render the IF to be null. At this time, particles start to slow down and eventually following the ordinary PSO velocity algorithm.

Within CAF, it is expected that the number of iteration required for GPSO particles to be closely clumped will decrease. This makes the improved GPSO to be the latest fast speed optimizer in PSO family. Figure 1 shows the flow chart of the improved GPSO algorithm.

3.0 RESULTS AND DISCUSSIONS

Two simulations had been done on six benchmark functions based on [6] as shown in Table 1 in order to evaluate the performance of the original GPSO and the improved (IGPSO). The first simulation is done in order to determine how many iterations are required for both GPSO and IGPSO to fulfill the requirement based on Equation (6), where \( \delta \) is the acceptable value determined by the desire of the programmer.

\[
s_{k} < \delta.
\]

(6)

The number of particles used are 200, maximum iteration is 1000 and \( c_{1} = c_{2} = c_{3} = 1 \). The simulation is done so that the simulation will stop only when both of GPSO and IGPSO fulfilled the requirement. Fitness average is calculated from the final iteration particles. Table 2 shows the first simulation result.

Table 1 Bench mark test functions

<table>
<thead>
<tr>
<th>Test Function ( n = 2 )</th>
</tr>
</thead>
<tbody>
<tr>
<td>( f_{1}(x) = \sum_{i=1}^{n} x_{i}^{2} )</td>
</tr>
<tr>
<td>( f_{2}(x) = \sum_{i=1}^{n} (10^i)^{(i-1)/(n-1)} x_{i}^{2} )</td>
</tr>
<tr>
<td>( f_{3}(x) = \sum_{i=1}^{n}</td>
</tr>
<tr>
<td>( f_{4}(x) = \sum_{i=1}^{n} (x_{i} + 0.5)^{2} )</td>
</tr>
<tr>
<td>( f_{5}(x) = \frac{1}{2} \sum_{i=1}^{n} (x_{i}^{4} - 6x_{i}^{2} + 5x_{i}) )</td>
</tr>
<tr>
<td>( f_{6}(x) = \sum_{i=1}^{n}</td>
</tr>
</tbody>
</table>

From Table 2, it is shown that IGPSO is able to fulfill the clumping size requirement more faster than GPSO at up to 62% rate. If IGPSO is to be simulated as long as GPSO is, the fitness average for final iteration for IGPSO is much more smaller than GPSO, which means the particles of IGPSO is much more closer when compared to GPSO. Even when requirement is not fulfilled in 1000 iterations, IGPSO is still producing smaller fitness average.
Table 2: Iteration and average fitness of single simulation results

<table>
<thead>
<tr>
<th>f</th>
<th>ε</th>
<th>Particle position limitation</th>
<th>Iteration required</th>
<th>Average Fitness after simulation stop</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td></td>
<td></td>
<td></td>
<td>GPSO</td>
</tr>
<tr>
<td>1</td>
<td>0.5</td>
<td>-100 to 100</td>
<td>392</td>
<td>0.0187</td>
</tr>
<tr>
<td>2</td>
<td>0.05</td>
<td>-10 to 10</td>
<td>212</td>
<td>7.03e-4</td>
</tr>
<tr>
<td>3</td>
<td>0.05</td>
<td>-10 to 10</td>
<td>158</td>
<td>4.20e-5</td>
</tr>
<tr>
<td>4</td>
<td>0.5</td>
<td>-100 to 100</td>
<td>269</td>
<td>0.0156</td>
</tr>
<tr>
<td>5</td>
<td>0.02</td>
<td>5 to 5</td>
<td>&gt;1000</td>
<td>-</td>
</tr>
<tr>
<td>6</td>
<td>0.05</td>
<td>-10 to 10</td>
<td>&gt;1000</td>
<td>0.2199</td>
</tr>
</tbody>
</table>

Figure 2 shows the sample of the CAF from Function 1. It can be seen that CAF never exceed the value of 1, so no over stepping of particles may happen. Over stepping of particles happens when particles moves to fast or too far. As the iteration increases, CAF decreases to zero, as expected. This figure shows that IF will only be applied at the beginning of the iteration, as the iteration goes on, the effective velocity value is obtained based on the original PSO algorithm.

Second simulation is done by executing 50 different runs of 50 iteration simulation. In each runs, the global best fitness particles is chosen. The number of particles used are 200, maximum iteration is 1000 and c1 = c2 = c3 = 1. At the end of the 50 runs, the worst fitness, best fitness, fitness average and standard deviation are determined from the 50 collected global best fitness particles. The result of the second simulation is shown in Table 3.

Table 3: Best fitness, worst fitness, average fitness and standard deviation of 50 different simulations

<table>
<thead>
<tr>
<th>f</th>
<th>Method</th>
<th>Worst</th>
<th>Best</th>
<th>Average</th>
<th>Std Dev</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>GPSO</td>
<td>0.0718</td>
<td>3.113e-5</td>
<td>0.0220</td>
<td>0.0219</td>
</tr>
<tr>
<td></td>
<td>IGPSO</td>
<td>0.0964</td>
<td>1.40e-4</td>
<td>0.0157</td>
<td>0.0188</td>
</tr>
<tr>
<td>2</td>
<td>GPSO</td>
<td>0.0029</td>
<td>5.085e-6</td>
<td>4.994e-4</td>
<td>5.722e-4</td>
</tr>
<tr>
<td>3</td>
<td>GPSO</td>
<td>0.0013</td>
<td>8.400e-7</td>
<td>1.482e-4</td>
<td>2.384e-4</td>
</tr>
<tr>
<td>4</td>
<td>GPSO</td>
<td>0.1122</td>
<td>5.093e-4</td>
<td>0.0221</td>
<td>0.0228</td>
</tr>
<tr>
<td>5</td>
<td>GPSO</td>
<td>-0.0047</td>
<td>-6.409e-6</td>
<td>-3.192e-4</td>
<td>0.0014</td>
</tr>
<tr>
<td>6</td>
<td>GPSO</td>
<td>0.1355</td>
<td>3.025e-4</td>
<td>0.0054</td>
<td>0.0192</td>
</tr>
<tr>
<td>7</td>
<td>IGPSO</td>
<td>0.0157</td>
<td>7.492e-5</td>
<td>0.0011</td>
<td>0.0023</td>
</tr>
</tbody>
</table>

4.0 CONCLUSIONS

From the simulation results and discussions, it can be concluded that the IGPSO is a faster version of GPSO, which able to reduce the optimization process up to 62% faster. The proposed correction acceleration factor (CAF) is proven to function as acceleration control mechanism for the improvement factor (IF). With the CAF as the global best fitness approaching the expected fitness, which in this case is 0. As the particles are getting close to each other, the particles will eventually slow down and starts to move based on original PSO velocity algorithm. By controlling the IF with CAF, overstepping of particles can be avoided. Overstepping of particles may cause more iterations are required in order to fulfill the clumping requirement. As a future work, IGPSO is to be implemented on high dimension or complex low dimension functions such as predicted mean vote (PMV) equations.

Acknowledgement

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References


