

opposite directions flow through the coils, the magnetic field in the centerline of the coil arrangement symmetry is $H(0,0,0) = 0$. According to [16], the magnetic field along the z -axis is as follows:

$$(1) \quad H(0,0,z) = \frac{J_0}{2} \left[\begin{array}{l} \left((a-z-Z_0) \ln \left[\frac{(R_0+b) + ((R_0+b)^2 + (a-z-Z_0)^2)^{1/2}}{(R_0-b) + ((R_0-b)^2 + (a-z-Z_0)^2)^{1/2}} \right] \right)^+ \\ \left((a+z+Z_0) \ln \left[\frac{(R_0+b) + ((R_0+b)^2 + (a+z+Z_0)^2)^{1/2}}{(R_0-b) + ((R_0-b)^2 + (a+z+Z_0)^2)^{1/2}} \right] \right)^- \\ \left((a-z+Z_0) \ln \left[\frac{(R_0+b) + ((R_0+b)^2 + (a-z+Z_0)^2)^{1/2}}{(R_0-b) + ((R_0-b)^2 + (a-z+Z_0)^2)^{1/2}} \right] \right)^+ \\ \left((a+z-Z_0) \ln \left[\frac{(R_0+b) + ((R_0+b)^2 + (a+z-Z_0)^2)^{1/2}}{(R_0-b) + ((R_0-b)^2 + (a+z-Z_0)^2)^{1/2}} \right] \right)^- \end{array} \right]$$

For the purpose of the investigation, we assumed that $J_0 = 250 \text{ A/m}^2$, $Z_a=0.7\text{m}$, $D=0.6 \text{ m}$, $d=0.25\text{m}$, $q=0.4\text{m}$, $L=1\text{m}$ and additionally introduced geometrical constraints as follows: $R_0+b \leq D$, $R_0-b \geq d$, $Z_0-a \geq q$, $Z_0 \leq L$, and $a \cdot b \cdot R_0 \leq 0.006\text{m}^3$. For these assumptions, the a , b , R_0 , Z_0 parameters are determined in such a manner to obtain the largest possible gradient of the magnetic field in an active area Z_a and to maintain simultaneously the maximal possible linearity of this gradient. The fitness function is defined as follows:

$$(2) \quad F = \frac{1000 \left[\left(\frac{4}{3} |H(z=0.75Z_a)| - |H(z=Z_a)| \right)^2 + 2 \left(2 |H(z=Z_a/4)| - |H(z=Z_a/2)| \right)^2 \right]^{1/2}}{|H(z=Z_a/2)|^k |H(z=Z_a)|}$$

The factor $k=0.15$ determines the priority of the field gradient magnitude with reference to its linearity.

Phase Angle Particle Swarm Optimization algorithm

In this method, a number of significant modifications and extensions in comparison to the original version was introduced. The operation mechanism and updating equations were rebuilt. The introduced changes concern both the way of the particle move and the way of the information collecting by the swarm during searching for the optimal solution. This helps the algorithm to explore the search space more efficiently. The PAPSO algorithm operation proceeds as follows:

1. Initialization, which relies on random attribution of the φ angle to each particles.
2. Establishing the particle position according to the formula [10]:

$$(3) \quad x_j^i = f(\varphi_j^i) = \frac{x_{max} - x_{min}}{2} \cdot \sin(\varphi_j^i) + \frac{x_{max} + x_{min}}{2}$$

3. Evaluation of the particle position by means of the fitness function
4. Reduction of the swarm cardinality by comparing the successive particles and the selection of better fitted particle (between two neighbored ones).
5. Establishing the best phase angle φ_{pb} , for each particle, by which the particle has managed to achieve its best position.
6. The choice of the particle of the φ_{gb} angle, by which the particle has achieved its best position within the whole swarm.
7. Updating the phase angle increment vector of each particle within the swarm according to the formula:

$$(4) \quad \begin{aligned} \Delta\varphi_{j+1}^i &= w\Delta\varphi_j^i + \sin(r_1)c_1(\varphi_{pb}^i - \varphi_j^i) \\ &+ \sin(r_2)c_2(\varphi_{gb} - \varphi_j^i) + c_3(\sqrt{3}(\varphi_{gb} + \varphi_{pb}^i)/4 - \varphi_j^i) \end{aligned}$$

8. Assigning two new phase angles to each particle, ($u \in [-1;1]$) according to the formulae:

$$(5) \quad \varphi_{j+1}^i = \varphi_j^i + \Delta\varphi_{j+1}^i$$

$$(6) \quad \varphi_{j+1}^{i2} = \varphi_j^i + u(-\Delta\varphi_{j+1}^i)$$

In this way, each particle with two new phase angles will obtain two proposals of its position.

9. Updating the particle location using (3) equation.
10. Evaluation of the particle position by means of the fitness function.
11. Comparison and selection of the new phase angle φ for each particle.
12. Repetition of 5-11 steps until the algorithm stopping criterion is met.

Where:

x_j^i -the position vector of the i -th particle in the j -th iteration,

$\Delta\varphi_{j+1}^i$ -the phase angle increment of the i -th particle,

φ_{pb}^i -phase angle of the personal best solution of the i -th particle,

φ_{gb} -the phase angle of the global best solution,

w -the inertial weight that determines the deviation of the particle original movement direction,

c_1, c_2, c_3 -the acceleration factor that determines how much the particle is influenced by its best phase angle and how much the particle is influenced by the whole swarm,

r_1, r_2 -the randomly generated angle in the range $(0^\circ, 90^\circ)$ for each iteration and for each particle.

The $\Delta\varphi_{j+1}^i$ value is limited by the $(\Delta\varphi_{min}, \Delta\varphi_{max})$ interval. As in case of $\Delta\varphi_{j+1}^i$ the value of the φ_j^i angle is limited and belongs to the $(\varphi_{min}, \varphi_{max})$ interval.

Results

Benchmark Test Functions

The first simulation tests of the new PAPSO algorithm were carried out on the benchmark function. The parameters of optimized functions are listed in Table 1. The obtained results were then compared with the achievements of the standard PSO and θ -PSO algorithms described in [10]. The computations were executed with inertia weight $w = 0,6$ and acceleration coefficients $c_1 = c_2 = 1,7$. The maximum number of iterations was established to 1000. The exemplary results of the tests performed for 40 particles of the swarm are depicted in Table 2. The presented values for PAPSO were averaged over 30 trials.

Table 2. PSO, θ -PSO and PAPSO algorithm performance

Function	Algorithm	Number of iterations		Number of achieved solution(%)
		Minimum	Average	
Camel	PAPSO	25	41	100
	θ -PSO	26	40	100
	PSO	32	60	100
Sphere	PAPSO	334	383	100
	θ -PSO	352	406	100
	PSO	577	716	100
Griewank	PAPSO	219	317	100
	θ -PSO	231	334	100
	PSO	375	461	100
Rosenbrock	PAPSO	180	279	100
	θ -PSO	194	283	100
	PSO	569	2268	55

The investigation of benchmark function confirmed the effectiveness of the proposed PAPSO algorithm. In almost all cases the new algorithm turned out to be more effective than the other algorithms used for the tests. For the tested functions (except Camel) the minimum and average

numbers of iterations were lower than those of θ -PSO and PSO. In order to find the optimum of this function with required accuracy fewer number of iterations was therefore needed. For Camel function the results of PAPSO and θ -PSO were almost the same but they were much better than the standard PSO.

Coil Arrangement

The study on an effectiveness of the proposed method used to determine an optimal geometry of the coil arrangement was undertaken by means of a computer program written in Mathematica. The computations were executed with acceleration coefficients $c_1=0.4$, $c_2=c_3=0.6$. The parameter w was used in the range 0.8 to 0.3 with a linearly decreasing, whereas the maximum number of iterations was fixed to 1000. The results were then compared with the achievements of the standard PSO and the θ -PSO algorithms.

The exemplary results of the tests performed for 10, 30, 50 and 80 particles in the initial population of the swarm are depicted in Fig. 2. All the values were averaged over 50 trials.

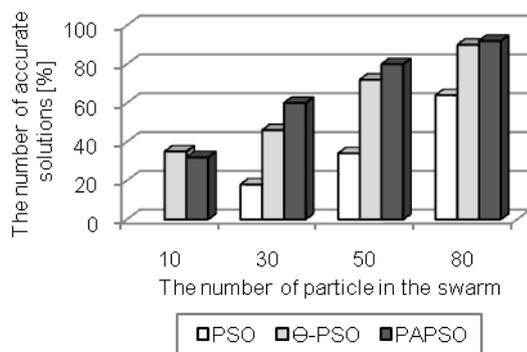


Fig.2. The number of accurate solutions versus the swarm cardinality for standard PSO, θ -PSO and PAPSO

The proposed algorithm turned out to be effective with respect to the number of accurate results and the number of iterations needed to achieve them. In comparison to the standard PSO (Table 3), the new algorithm was able to find more accurate solutions within a few times lower iteration number.

Table 3. The relationship between the swarm cardinality and the number of iterations to achieve the accurate solutions for the PSO, θ -PSO and PAPSO

Algorithm	The number of particles in the swarm			
	10	30	50	80
PAPSO	107	82	64	57
θ -PSO	131	112	76	64
PSO	-	642	567	531

Moreover, the algorithm also achieved to be efficient for small populations of particles. In case of PSO for 10 particles in the swarm, no accurate solution was managed to be obtained when the number of iterations was as high as 1000, whereas the PAPSO could find 32% optimal solutions within only 107 iterations (on average). In most cases, the results were also more accurate than for the θ -PSO algorithm. Only for the population size of 10 particles, the PAPSO algorithm found 3% fewer accurate solutions than θ -PSO, but for lower number of iterations of even 18%.

As far as the optimal solutions are concerned, the best results were obtained for large swarm. For the populations comprising 80 particles, PAPSO algorithm could find over 80% accurate solutions. It was also found, that the

proposed new method gave the solutions of the same accuracy, and faster exploration of the search space.

Summary

In the following study, the new PAPSO optimization algorithm with the phase angle vector was proposed. In the proposed method, determination of the velocity vector was omitted. Instead of the velocity vector, the phase angle vector was introduced. The equation of the phase angle increment update is responsible for the particle movement within the search space. The particle position is determined by the mapping of the corresponding angles.

The proposed algorithm was tested on benchmark test functions and was applied to determine the optimal geometry of the coil arrangement generating magnetic field with the specific parameters. The results were then compared with performances of the standard PSO and the θ -PSO algorithms. The calculations obtained in this study confirmed efficiency of the proposed algorithm.

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Author: dr inż. Bożena Borowska, Politechnika Łódzka, Instytut Informatyki, ul. Wólczajska 215, 93-005 Łódź, E-mail: bozena.borowska@p.lodz.pl