



Modified Water Evaporation Optimization based Design for Improving Efficiency of Three Phase Induction Motor

P. S. Prakash

Assistant Professor, Department of Electrical & Electronics Engineering, Annamalai University, Annamalainagar,
Tamilnadu, India

ABSTRACT: This paper presents a Modified Water Evaporation Optimization (MWEO) based design methodology for refining the efficiency of Induction Motor (IM). Modified Water Evaporation algorithm optimization, inspired by water evaporation from lake or river. MWEO is one of the biological and environmental science is computing models for solving multimodal optimization problems. Among the number of design variables of the IM, seven variables are identified as primary design variables and the MWEO based design methodology is tailored to optimize the chosen primary variables with a view to obtain the global best design. The developed methodology is applied in solving two IM design problems and the results are presented with a view of exhibiting the superiority of the developed algorithm.

KEYWORDS: Induction Motor Design, Water evaporation optimization

I. INTRODUCTION

Induction motors (IM) are the most widely used in domestic, commercial and various industrial applications. Especially, the squirrel cage IM is characterized by its simplicity, robustness and low cost, making it more attractive and hence captured a leading place in industrial and agricultural sectors. As millions of such motors are in use in various sectors, they consume a considerable percentage of overall produced electrical energy. The ever mounting pressure of oil crisis and the need for energy conservation necessitate designing the IMs with increased levels of efficiency through the selection of appropriate combination of the design parameters. The optimal design of IM (ODIM) is so complicated that it is still a combination of art and science. There are many geometrical parameters and their relationships connected with motor specifications, which are in general nonlinear (MehmetCunkas 2010).

Over the years, in addition to statistical methods (Han and Shapiro 1967) and the Monte Carlo technique (Anderson 1967), several mathematical programming techniques, which provides a means for finding the minimum or maximum of a function of several variables under a prescribed set of constraints, have been applied in solving the IM design problems. These techniques such as nonlinear programming, (Ramarathnam et al 1971), Lagrangian relaxation method (Gyeorye Lee et al 2013), direct and indirect search methods (Nagriyal et al 1979), Hooks and Jeeves method (Faiz et al 2001), Rosenbrock's method (Bharadwaj et al 1979-a), Powell's method (Ramarathnam et al 1973), finite element method (20/4:7 T. S. Parkin et al 1993) and sequential unconstrained minimization technique (Bharadwaj et al 1979-b) are most cumbersome and time consuming. Besides a few of them requires derivatives and exhibits poor convergence properties due to approximations in the derivative calculations.

Apart from the above methods, another class of numerical techniques called evolutionary search algorithms such as simulated annealing (Bhuvanewari et al 2005; Kannan et al 2010), genetic algorithm (GA) (Satyajit Samaddar et al 2013; Sivaraju et al 2011), evolutionary algorithm (Jan Pawel Wieczorek et al 1998), evolutionary strategy (Kim MK et al 1998), and particle swarm optimization (PSO) (Thanga Raj et al 2008; Sakthivel et al 2011) have been widely applied in solving the IM design problems. Having in common processes of natural evolution, these algorithms share many similarities; each maintains a population of solutions that are evolved through random alterations and selection. The differences between these procedures lie in the techniques they utilize to encode candidates, the type of alterations



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they use to create new solutions, and the mechanism they employ for selecting the new parents. These algorithms have yielded satisfactory results across a great variety of engineering optimization problems.

Recently, Water Evaporation Optimization (WEO) has been suggested for solving optimization problems. It is inspired by evaporation of water is very important in biological and environmental science, the decreasing wettability of surface can represent the decrease of objective function for a minimizing optimization problem. It has been applied to a variety of power system problems and found to yield satisfactory results.

The aim of this paper is to develop a MWEO based method for optimally designing IMs with a view of effectively exploring the solution space and obtaining the global best solution. The developed methodology has been applied in designing two IMs and the performances have been studied. The paper is divided into six sections. Section I provides the introduction, section II overviews WEO and elucidates the proposed method (PM), section III formulates the IM design problem, section IV discusses the results and section V concludes.

II. WATER EVAPORATION OPTIMIZATION ALGORITHM

The evaporation of water is very important in biological and environmental science. The water evaporation from bulk surface such as a lake or a river is different from evaporation of water restricted on the surface of solid materials. In this WEO algorithm water molecules are considered as algorithm individuals. Solid surface or substrate with variable wettability is reflected as the search space. Decreasing the surface wettability (substrate changed from hydrophilicity to hydrophobicity) reforms the water aggregation from a monolayer to a sessile droplet.

Such a behavior is consistent with how the layout of individuals changes to each other as the algorithm progresses. And the decreasing wettability of surface can represent the decrease of objective function for a minimizing optimization problem. Evaporation flux rate of the water molecules is considered as the most appropriate measure for updating individuals which its pattern of change is in good agreement with the local and global search ability of the algorithm and make this algorithm have well converged behavior and simple algorithmic structure. The details of the water evaporation optimization algorithm are well presented in (Kaveh and Bakhshpoori, 2016).

In the WEO algorithm, each cycle of the search consists of following three steps (i) Monolayer Evaporation Phase, this phase is considered as the global search ability of the algorithm (ii) Droplet Evaporation Phase, this phase can be considered as the local search ability of the algorithm and (iii) Updating Water Molecules, the updating mechanism of individuals.

(i) Monolayer Evaporation Phase

In the monolayer evaporation phase the objective function of the each individuals Fit_i^t is scaled to the interval [-3.5, -0.5] and represented by the corresponding $E_{sub}(i)^t$ inserted to each individual (substrate energy vector), via the following scaling function.

$$E_{sub}(i)^t = \frac{(E_{max} - E_{min}) \times (Fit_i^t - Min(Fit))}{(Max(Fit) - Min(Fit))} + E_{min} \quad (1)$$

Where E_{max} and E_{min} are the maximum and minimum values of E_{sub} respectively. After generating the substrate energy vector, the Monolayer Evaporation Matrix (MEP) is constructed by the following equation.

$$MEP_{ij}^t = \begin{cases} 1 & \text{if } rand_{ij} \leq \exp(E_{sub}(i)^t) \\ 0 & \text{if } rand_{ij} \geq \exp(E_{sub}(i)^t) \end{cases} \quad (2)$$

where MEP_{ij}^t is the updating probability for the j^{th} variable of the i^{th} individual or water molecule in the t^{th} iteration of the algorithm. In this way an individual with better objective function is more likely to remain unchanged in the search space.

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(ii) Droplet Evaporation Phase

In the droplet evaporation phase, the evaporation flux is calculated by the following equation.

$$J(\theta) = J_o P_o \left(\frac{2}{3} + \frac{\cos^3 \theta}{3} - \cos \theta \right) (1 - \cos \theta) \quad (3)$$

where J_o and P_o are constant values. The evaporation flux value is depends upon the contact angle Θ , whenever this angle is greater and as a result will have less evaporation. The contact angle vector is represented the following scaling function.

$$\theta(i)^t = \frac{(\theta_{\max} - \theta_{\min}) \times (Fit_i^t - Min(Fit))}{(Max(Fit) - Min(Fit))} + \theta_{\min} \quad (4)$$

Where the min and max are the minimum and maximum functions. The Θ_{\min} & Θ_{\max} values are chosen between $-50^\circ < \Theta < -20^\circ$ is quite suitable for WEO. After generating contact angle vector $\Theta(i)^t$ the Droplet Probability Matrix (DEP) is constructed by the following equation.

$$DEP_{ij}^t = \begin{cases} 1 & \text{if } rand_{ij} < J(\theta_i^{(t)}) \\ 0 & \text{if } rand_{ij} \geq J(\theta_i^{(t)}) \end{cases} \quad (5)$$

where DEP_{ij}^t is the updating probability for the j^{th} variable of the i^{th} individual or water molecule in the t^{th} iteration of the algorithm.

(iii) Updating Water Molecules

In the WEO algorithm the number of algorithm individuals or number of water molecules (nWM) is considered constant in all t^{th} iterations, where t is the number of current iterations. Considering a maximum value for algorithm iterations (t_{\max}) is essential for this algorithm to determine the evaporation phase and for stopping criterion. When a water molecule is evaporated it should be renewed. Updating or evaporation of the current water molecules is made with the aim of improving objective function. The best strategy for regenerating the evaporated water molecules is using the current set of water molecules ($WM^{(t)}$). In this way a random permutation based step size can be considered for possible modification of individual as:

$$S = rand \cdot (WM^{(t)} [permutel(i)(j)] - WM^{(t)} [permute2(i)(j)]) \quad (6)$$

where $rand$ is a random number in $[0,1]$ range, $permutel$ and $permute2$ are different rows of permutation functions. i is the number of water molecule, j is the number of dimensions of the problem. The next set of molecules ($WM^{(t+1)}$) is generated by adding this random permutation based step size multiplied by the corresponding updating probability (monolayer evaporation and droplet evaporation probability) and can be stated mathematically as:

$$WM^{(t+1)} = WM^{(t)} + S \times \begin{cases} MEP^{(t)} & t \leq t_{\max} / 2 \\ DEP^{(t)} & t > t_{\max} / 2 \end{cases} \quad (7)$$

Each water molecule is compared and replaced by the corresponding renewed molecule based on objective function. It should be noted that random permutation based step size can help in two aspects. In the first phase, water molecules are more far from each other than the second phase. In this way the generated permutation based step size will guarantee global and local capability in each phase.

III. IMPLEMENTATION OF WEO ALGORITHM TO SLOVE DESIGN PROBLEM

The detailed algorithmic steps for proposed WEO algorithm to solve design problem are presented below.

Step 1: Initialize total no of design variables, design constraints limits, number of water molecules, maximum number of algorithm iteration (t_{\max}),



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Step 2: Randomly initialize all water molecules.

Step 3: Obtain the design variable by applying priority list method and compute the objective function given by Eq. (1), Eq. (4) and Eq. (8) for all water molecules.

Step 4: Check whether t (current iteration) $\leq t_{\max}/2$.

Step 5: If step 4 is satisfied, then, water molecules are globally evaporated based on monolayer evaporation probability MEP using Eq. (2).

Step 6: For $t > (1+t_{\max}/2)^2$, Based on DEP (Eq. 5), evaporation occurs.

Step 7: Generate random permutation based step size matrix according to Eq. (6).

Step 8: Generate evaporated water molecules by adding the product of step size matrix and evaporation matrix to the current set of molecules MWM^(t) by using Eq. (7) and update the matrix of water molecules.

Step 9: Compare and update the water molecules.

Step 10: Return the best water molecule

Step 11: If the number of iteration of the algorithm (t) becomes larger than the maximum number of iterations (t_{\max}), the algorithm terminates. Otherwise go to step 3.

IV. PROPOSED METHOD

The proposed MWEEO based solution method for ODIM involves formulation of the problem, representation of water molecules through the chosen design variables and water molecules are accelerated to move as best solutions.

4.1 Problem Formulation

The ODIM problem involves large number of design variables. Many of these variables fortunately have a little influence either on the objective function or on the specified constraints. However, to ease the curse of high dimensionality, the following seven variables are identified as primary design variables.

$$X = [x_1, x_2, \dots, x_7] = \begin{bmatrix} \text{Core length to pole pitch} \\ \text{Average value of air gap flux density} \\ \text{Ampere conductor} \\ \text{Length of air gap} \\ \text{Stator current density} \\ \text{Rotor current density} \\ \text{Flux density in the core} \end{bmatrix}^T \quad (8)$$

The ODIM problem is formulated by defining an objective function and a set of constraints as

$$\text{Maximize } h(x) = \frac{KW}{KW + P_t} \quad (9)$$

Subject to

$$g(x) \leq 0 \Leftrightarrow \begin{cases} \text{maximum flux density of stator teeth} \leq 2 \\ \text{maximum flux density of rotor teeth} \leq 2.0 \\ \text{slip at full load} \leq 0.05 \\ \text{starting to full load torque ratio} \geq 1.5 \\ \text{stator temperature rise} \leq 70 \\ \text{per unit no load current} \leq 0.5 \end{cases} \quad (10)$$

$$x_i^{\min} \leq x_i \leq x_i^{\max} \quad i = 1, 2, \dots, nd \quad (11)$$

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Where $P_t = P_{nl} + P_{cus} + P_{cur}$ (12)

4.2 Representation of Design Variables

The water molecules s_i is represented to denote the chosen primary design variables, defined by Eq. (8), in vector form as:

$$s_i = [s_i^1, s_i^2, \dots, s_i^7] = [x_1, x_2, \dots, x_7] \quad (13)$$

4.3 Fitness Function

The algorithm searches for optimal solution by maximizing a fitness of Water molecules LI , which is formulated from the objective function of Eq. (9) and the penalty terms representing the limit violation of the explicit constraints of Eq. (10). The LI function is written as

$$\text{Maximize } LI = \frac{h(x)}{1 + w \sum_{i \in \eta} [g_i(x)]^2} \quad (14)$$

V. NUMERICAL RESULTS

The proposed MWEQ method (PM) is used to obtain the optimal design of two IMs. The first machine under study is rated for 7.5 kW, 400 V, 4 pole, 50 Hz and the second one for 30 kW, 400 V, 4 pole, 50 Hz. The effectiveness of the PM is demonstrated through comparing the performances with those of the ACO and WEO based design approaches. In this regard, the same set of primary design variables, fitness function and design equations, involved in the PM, are used to develop the ACO and WEO based design approaches. The software packages are developed in Matlab platform and executed in a 2.3 GHz Pentium-IV personal computer. There is no guarantee that different executions of the developed design programs converge to the same design due to the stochastic nature of the ACO and WEO, and hence the algorithms are run 20 times and the best ones are presented. The optimal design representing the values of the primary design variables for both the IMs and their efficiencies are presented in Table-1 and 2

Table 1 Comparison of Results for Motor-1

		ACO	WEO	PM
Primary Design Variables	x_1	1.31979	1.28631	1.29731
	x_2	0.42097	0.43608	0.42638
	x_3	23155.24	22965.20	22775.20
	x_4	0.46367	0.50578	0.48579
	x_5	3.64492	3.83688	3.73622
	x_6	2.00461	2.10381	2.01372
	x_7	1.10145	1.11679	1.11570
Constraints $g(x)$	$g_1 \leq 2$	1.621	1.720	1.711
	$g_2 \leq 2$	1.738	1.806	1.822
	$g_3 \leq 0.05$	0.021	0.022	0.020
	$g_4 \geq 1.5$	3.427	3.866	3.877
	$g_5 \leq 70$	45.728	48.768	48.826
	$g_6 \leq 0.5$	0.342	0.396	0.386
	$g_7 \geq 0.75$	0.854	0.799	0.811
Objective function $h(x)$	% Efficiency	86.727	86.757	86.899

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It is observed from these tables that the PM offers an efficiency of **86.899%** and **91.012%**, which are higher than those of ACO and WEO based approaches, for motor-1 and -2 respectively. These tables also include the values of the constraints of Eq. (10) along with their limits. It can also be observed from these tables that all the methods bring the constraints such as maximum flux density, slip at full load, starting to full load torque ratio, etc. to lie within the respective limit, as the constraints are added as penalty terms in the light intensity function of Eq. (14). It is obvious that the PM offers better %efficiency enhancement than those of the existing approaches for both the motors.

Table 2 Comparison of Results for Motor-2

		ACO	WEO	PM
Primary Design Variables x	x_1	1.19110	1.67666	1.68155
	x_2	0.44126	0.39535	0.39356
	x_3	28217.57	24855.40	24922.10
	x_4	0.89713	0.79880	0.79723
	x_5	2.69634	2.89081	2.88231
	x_6	2.01649	1.10505	1.1123
	x_7	1.10062	1.10540	1.11523
Constraints $g(x)$	$g_1 \leq 2$	1.487	1.882	1.892
	$g_2 \leq 2$	1.528	1.695	1.699
	$g_3 \leq 0.05$	0.016	0.001	0.011
	$g_4 \geq 1.5$	1.748	1.612	1.677
	$g_5 \leq 70$	46.449	34.275	34.783
	$g_6 \leq 0.5$	0.347	0.309	0.399
	$g_7 \geq 0.75$	0.774	0.719	0.701
Objective function $h(x)$	% Efficiency	90.582	90.989	91.012

VI. CONCLUSION

Indeed the MWEO is a powerful population based method for solving complex optimization problems. A new methodology involving MWEO for solving ODIM problem has been developed and applied on two IM design problems. It determines the optimal values for primary design variables. The ability of the PM to produce the global best design parameters that improves the efficiency of the motor has been projected. It has been chartered that the new approach fosters the continued use of MWEO and will go a long way in serving as a useful tool in design problems.

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