# Clustering in Evolutionary Algorithms to Efficiently Compute Simultaneously Local and Global Minima

D.K. Tasoulis, V.P. Plagianakos, and M.N. Vrahatis

Department of Mathematics, University of Patras, GR-26110 Patras, Greece University of Patras Artificial Intelligence Research Center E-mail: {dtas,vpp,vrahatis}@math.upatras.gr

Abstract- In this paper a new clustering operator for Evolutionary Algorithms is proposed. The operator incorporates the unsupervised *k*-windows clustering algorithm, utilizing already computed pieces of information regarding the search space in an attempt to discover regions containing groups of individuals located close to different minimizers. Consequently, the search is confined inside these regions and a large number of global and local minima of the objective function can be efficiently computed. Extensive experiments shown that the proposed approach is effective and reliable, and greatly accelerates the convergence speed of the considered algorithms.

# **1** Introduction

Evolutionary Algorithms (EAs) refer to problem solving optimization algorithms which employ computational models of evolutionary processes. A variety of evolutionary algorithms have been proposed. The major ones include: Genetic Algorithms [11, 13], Evolutionary Programming [9, 10], Evolution Strategies [25, 27], Genetic Programming [16], and the Differential Evolution algorithm [31]. All these algorithms share the common conceptual base of simulating the evolution of the individuals that form the population using a predefined set of operators. Commonly two kinds of operators are used: the *selection* and the *search* operators. The most widely used search operators are *mutation* and *recombination*.

The selection operator mainly depends on the perceived measure of the fitness of each individual and enforces the natural selection and the survival of the fittest. The recombination and the mutation operators stochastically perturb the individuals providing efficient exploration of the search space. This perturbation is primarily controlled by the user defined recombination and mutation rates. Although simplistic from a biologist's point of view, these algorithms are sufficiently complex to provide robust and powerful search mechanisms and have shown their strength in solving hard optimization problems.

For the rest of the paper we consider the minimization problem of finding global minima of a continuous nonlinear, (possibly) nondifferentiable, multimodal objective function f. More specifically, our goal is to locate global minimizers  $x_t^*$  of the real-valued objective function  $f: \mathcal{E} \to \mathbb{R}$ :

$$f(x_t^*) \leqslant f(x), \quad \forall x \in \mathcal{E}$$

where t = 1, 2, ..., and the compact set  $\mathcal{E} \subseteq \mathbb{R}^n$  is a *n*-dimensional scaled translation of the unit hypercube.

In this paper, we propose a new clustering operator for EAs and we aim to study the above minimization problem focusing on the well-known Differential Evolution (DE) algorithm. DE utilizes mutations and recombinations as search mechanisms and selection to direct the search towards the most promising area of the solution space. It uses a non uniform mutation operator that can take child parameters from a particular parent more often than it does from others. DE has been applied to a large number of different optimization tasks. It has successfully solved many artificial benchmark problems [30], as well as hard real-world problems (see for example [6, 14, 26, 28]). In [15] DE has been applied to train neural networks and in [20, 21] we have proposed a method that made possible the efficient training of neural networks having arbitrary as well as constrained integer weights. The DE algorithm has also been implemented on parallel and distributed computers [7, 22].

Although DE is capable of optimizing difficult multimodal objective functions, the choice of the appropriate mutation operator depends on the problem at hand and in general is a nontrivial task that demands experimentation. In the literature numerous mutation operators have been proposed for DE [8, 31], having various effects on its exploration and exploitation dynamics. In this paper, we study the performance of some of the most commonly used mutation operators and propose a novel clustering operator that incorporates clustering techniques to accelerate the convergence of the algorithm. In general, clustering can be defined as the process of partitioning a set of patterns into disjoint and homogeneous meaningful groups, called clusters. In our cases the clusters discovered are regions of  $\mathcal{E}$  containing a minimum of the objective function f.

The key features of the proposed operator is that it utilizes already computed pieces of information regarding the search space and that it efficiently computes a large number of minima of the objective function in the process of finding the global ones. To achieve its design goals, the new operator utilizes the recently proposed unsupervised k-windows clustering algorithm [3, 4, 32, 36]. Historically, clustering techniques were originally conceived by Aristotle and Theophrastos in the fourth century B.C. and later during the 18th century were studied by Linnaeus [17]. However, their first comprehensive foundations were published in 1939 [41]. Clustering is considered fundamental in knowledge acquisition and has been applied in numerous fields including, statistical data analysis [1], compression and vector quantization [24], image analysis, etc.

In addition to the ongoing research aiming at the study of new clustering algorithms, a large amount of research has already been attributed to the development of clustering methods trying specifically to tackle global optimization problems. The starting point was Törn's work [39, 40]. Timmer in [37] considered several clustering methods and a book that tries to cover this field was published in 1989 [38]. Nevertheless, to the best of our knowledge, although novel efficient global optimization and clustering algorithms are being constantly proposed, not much work has been performed to identify the applicability and the performance of their synergy.

The rest of the paper is organized as follows. In Section 2 the Differential Evolution algorithm is outlined and a comparative study of different mutation operators is performed. In Section 3 the k-windows clustering algorithm is briefly described. In Section 4 the new DE operator is proposed and Section 5 is devoted to the presentation and the discussion of the experimental results. The paper ends with concluding remarks and some pointers for future work.

### 2 The Differential Evolutionary Algorithm

Differential Evolution [31] is a minimization method, capable of handling nondifferentiable, nonlinear and multimodal objective functions. To fulfill this requirement, DE has been designed as a stochastic parallel direct search method, which utilizes concepts borrowed from the broad class of evolutionary algorithms. The method typically requires few, easily chosen, control parameters. Experimental results have shown that DE has good convergence properties and outperforms other well known evolutionary algorithms [29, 31].

DE is a population-based stochastic algorithm that exploits a population of potential solutions, individuals, to effectively probe the search space. The population of the individuals is randomly initialized in the optimization domain with NP, n-dimensional vectors, following a uniform probability distribution and is evolved over time in order to explore the search space and locate the minima of the objective function. NP is fixed throughout the training process. At each iteration, called generation, new vectors are generated by the combination of vectors randomly chosen from the current population. This operation in our context can be referred to as *mutation*. The outcoming vectors are then mixed with another predetermined vector - the target vector - and this operation can be called as *recombination*. This operation yields the so-called trial vector. The trial vector is accepted for the next generation if and only if it yields a reduction in the value of the objective function f. Otherwise, target vector is retained in the next generation. This last operator can be referred to as selection.

#### 2.1 Search Operators

Here, we briefly describe the search operators that were studied in this paper. The search operators efficiently shuffle information among the individuals, enabling the search for an optimum to focus on the most promising regions of the solution space. The first operator considered is the mutation. Specifically, for each individual  $x_a^i$ , i = 1, ..., NP, where g denotes the current generation, a new individual  $v_{g+1}^i$  (mutant vector) is generated according to one of the following equations:

$$v_{g+1}^i = x_g^{\text{best}} + \mu(x_g^{r1} - x_g^{r2}), \tag{1}$$

$$x_{g+1}^i = x_g^{r_1} + \mu(x_g^{r_2} - x_g^{r_3}),$$
 (2)

$$v_{g+1}^i = x_g^i + \mu (x_g^{\text{best}} - x_g^i) + \mu (x_g^{r1} - x_g^{r2}),$$
 (3)

$$v_{g+1}^i = x_g^{\text{best}} + \mu(x_g^{r1} - x_g^{r2}) + \mu(x_g^{r3} - x_g^{r4}),$$
 (4)

$$v_{g+1}^{i} = x_{g}^{r1} + \mu(x_{g}^{r2} - x_{g}^{r3}) + \mu(x_{g}^{r4} - x_{g}^{r5}), \quad (5)$$

where  $x_g^{\text{best}}$  is the best member of the previous generation;  $\mu > 0$  is a real parameter, called *mutation constant*, which controls the amplification of the difference between two individuals so as to avoid the stagnation of the search process; and  $r_1, r_2, r_3, r_4, r_5 \in \{1, 2, \dots, i - 1, i + 1, \dots, NP\}$ , are random integers mutually different and not equal to the running index *i*.

Trying to rationalize the above equations, we observe that Equation (2) is similar to the crossover operator used by some Genetic Algorithms and Equation (1) derives from it, when the best member of the previous generation is employed. Equations (3), (4) and (5) are modifications obtained by the combination of Equations (1) and (2). It is clear that more such relations can be generated using the above ones as building blocks. For example, the recently proposed trigonometric mutation operator [8] performs mutations with probability  $\mu$  according to the following equation:

$$\begin{aligned} v_{g+1}^i &= (x_g^{r1} + x_g^{r2} + x_g^{r3})/3 + (p_2 - p_1)(x_g^{r1} - x_g^{r2}) + \\ &+ (p_3 - p_2)(x_g^{r2} - x_g^{r3}) + (p_1 - p_3)(x_g^{r3} - x_g^{r1}), \ (6) \end{aligned}$$

and with probability  $(1 - \mu)$  mutations according to Equation (2), where  $\mu$  is a user defined parameter. The values of  $p_m$ ,  $m = \{1, 2, 3\}$  and p' are obtained through the following equations:

$$\begin{aligned} p_1 &= \left| f(x_g^{r1}) \right| / p', \\ p_2 &= \left| f(x_g^{r2}) \right| / p', \\ p_3 &= \left| f(x_g^{r3}) \right| / p', \text{ and} \\ p' &= \left| f(x_g^{r1}) \right| + \left| f(x_g^{r2}) \right| + \left| f(x_g^{r3}) \right|. \end{aligned}$$

For the rest of the paper, we call  $DE_1$  the differential evolution algorithm that uses Equation (1) as the mutation operator,  $DE_2$  the algorithm that uses Equation (2), and so on.

The recombination operator is subsequently applied to further increase the diversity of the mutant individuals. To this end, the resulting individuals are combined with other predetermined individuals, called the target individuals. Specifically, for each component l (l = 1, 2, ..., n)of the mutant individual  $v_{g+1}^i$ , we randomly choose a real number r in the interval [0, 1]. Then, we compare this number with the *recombination constant*,  $\rho$ . If  $r \leq \rho$ , then we select, as the l-th component of the trial individual  $u_{g+1}^i$ , the l-th component of the target vector  $x_{g+1}^i$  becomes the l-th component of the target vector  $x_{g+1}^i$  becomes the l-th component of the trial vector. This operation yields the trial individual. Finally, the trial individual is accepted for the next generation only if it reduces the value of the objective function.

#### 2.2 Exploration vs. Exploitation

The main problem when applying Evolutionary Algorithms is to find a set of control parameters which optimally balances the exploration and the exploitation capabilities of the algorithm. There is always a trade off between the efficient exploration of the search space and its effective exploitation. For example, if the recombination and mutation rates are too high, much of the space will be explored, but there is a high probability of losing good solutions. In extreme cases the algorithm has difficulty to converge to the global minimum due to insufficient exploitation of the search space. Fortunately, the convergence properties of the DE typically do not heavily depend on its control parameters. However, since not all search operators have the same impact on the exploration of the search space, the choice of the optimal mutation operator can be troublesome. To illustrate this we utilize the following simple multimodal 2-dimensional function:

$$f(x_1, x_2) = \sin(x_1)^2 + \sin(x_2)^2,$$

where  $(x_1, x_2) \in \mathbb{R}^2$ . This function has an infinite number of global minima in  $\mathbb{R}^2$ , with function value equal to zero, at the points  $(\kappa \pi, \pi)$ , where  $\kappa, \in \mathbb{Z}$ . In the hypercube  $[-5, 5]^2$  the function f has 9 global minima. In Figure 1 a surface plot of the function f is exhibited.



Figure 1: 3–D Plot of the  $sin(x_1)^2 + sin(x_2)^2$  function

The six DE variants described above are applied to compute the global minima of the objective function f. Experimental results indicate that DE<sub>1</sub> exhibits very fast convergence to one of the global minima of f. On the contrary, DE<sub>2</sub> explores a large portion of the search space before converging to a solution. This fact is illustrated in Figures 2 and 3, where (for visualization purposes) a population consisting of 1000 individuals is plotted after 1, 5, 10, 20 generations of DE<sub>1</sub> and DE<sub>2</sub>, respectively.

A closer look at Equations (1) and (2) reveals that  $DE_1$  uses the best individual as a starting point for the computation of the mutant vector, thus constantly pushing the population closer to the location of the best computed point. On

the other hand, since  $DE_2$  utilizes three randomly chosen individuals for the computation of the mutant one, its exploration capability is greatly enhanced. However, it exhibits lower convergence speed.



Figure 2:  $DE_1$  population after 1, 5, 10, and 20 generations



Figure 3: DE<sub>2</sub> population after 1, 5, 10, and 20 generations

The performance of algorithms  $DE_3$  and  $DE_4$  resembles that of  $DE_1$ , due to the use of the best individual. However,  $DE_3$  and  $DE_4$  exhibited better exploration than  $DE_1$ , since they also incorporate randomly selected individuals. Algorithms  $DE_5$  and  $DE_6$  use only randomly selected individuals resulting in maximum exploration and the individuals of their populations are simultaneously attracted by more than one minimizers. In Figure 4 the population of the  $DE_6$ algorithm at different generations is plotted.

Figures 3 and 4 show that some mutation operators have the tendency to concentrate subsets of the population in the region of attraction of different minimizers of the objective function. This observation motivated the incorporation of a clustering algorithm to identify such subsets. Experiments indicate that the k-windows clustering algorithm discovers boxes capturing subpopulations of individuals located in the region of a minimizer. Consequently, the subpopulations are confined to search within each box. Thus, the optimization of the objective function proceeds without affecting the dynamics of the DE algorithm. This process gives all the minima existing in regions that the DE algorithm explored before the use of the clustering operator. It is obvious that DE



Figure 4: DE<sub>6</sub> population after 1, 5, 10, and 20 generations

algorithm must adequately explore the search space prior to the call of the clustering operator. In the next section, for completeness purposes, we give a brief description of the k-windows clustering algorithm.

# 3 The Unsupervised *k*-windows Clustering Algorithm

The recently proposed k-windows clustering algorithm uses a windowing technique to discover the clusters present in an n-dimensional dataset. More specifically, assuming that the dataset lies in n dimensions, the algorithm initializes a number of n-dimensional windows (boxes) over the dataset. Subsequently, it iteratively perturbs these windows using the movement and enlargement procedures, in order to capture within each window patterns that belong to a single cluster.

The movement and enlargement procedures are guided by the points that lie within each window. As soon as the movement and enlargement procedures do not significantly increase the number of points within each window they terminate. The final set of windows defines the clustering result of the algorithm.

A fundamental issue in cluster analysis, independent of the particular clustering technique applied, is the determination of the number of clusters present in a dataset. For instance well-known and widely used iterative techniques, such as the k-means algorithm [12] as well as the fuzzy cmeans algorithm [5], require from the user to specify the number of clusters present in the data prior to the execution of the algorithm. On the other hand, the unsupervised kwindows algorithm is capable to determine the number of clusters through a generalization of the original algorithm.

The unsupervised version of the k-windows algorithm was used in this paper, since the number of minima of an objective function is, in general, unknown. For a comprehensive description of the algorithm and an investigation of its capability to automatically identify the number of clusters present in a dataset, refer to [3, 4, 32, 36].

It must be noted that no objective function evaluations are necessary during the operation of the k-windows clustering algorithm. The computationally demanding step of the *k*-windows clustering algorithm is the determination of the points that lie in a specific window. This is the well studied *orthogonal range search* problem [23]. Numerous Computational Geometry techniques have been proposed [2, 23] to address this problem. All these techniques employ a preprocessing stage at which they construct a data structure that stores the patterns. This data structure allows them to answer range queries fast. For applications of very high dimensionality, data structures like the Multidimensional Binary Tree [23] are more suitable. On the other hand, for low dimensional data with a large number of points the approach of [2] seems more attractive. The *k*-windows clustering algorithm has been successfully applied in numerous applications including bioinformatics [33, 34], medical diagnosis [18, 35], and time series prediction [19].

# 4 The Proposed Clustering Operator

In this section, the clustering operator is described. The proposed operator utilizes the unsupervised k-windows algorithm and is called only once, after a user-defined number of generations. In practice, a small number of generations is sufficient for the DE algorithm to explore the search space.

Afterwards, the clusters of individuals are determined and subpopulations are confined within each region. Each subpopulation has NP/ individuals, where is the number of clusters found. If a region contains more individuals, the clustering operator selects the best NP/. On the other hand, if less individuals exist the clustering operator initializes new ones. The result of the algorithm is the location of many minima in a single run, including the global one. The modified DE algorithm using the clustering operator is outlined in Algorithm 1.

DIFF	FERENTIAL EVOLUTION ALGORITHM MODEL
0:	Initialize the population of NP individuals
1:	Evaluate the fitness of each individual
2:	Repeat
3:	For $i = 1$ to $NP$
4:	$Mutation(x_a^i) \rightarrow Mutant_a^i$
5:	Recombination(Mutant <sup>i</sup> <sub>a</sub> ) $\rightarrow$ Trial <sup>i</sup> <sub>a</sub>
6:	If $f(\text{Trial}_{a}^{i}) \leq f(x_{a}^{i}),$
7:	accept Trial <sup><math>i</math></sup> for the next generation
8:	EndFor
9:	Until the search space is sufficiently explored
10:	Find clusters using the clustering operator
11:	For $j = 1$ to
12:	Confine NP/ individuals within each cluster
13:	Use the DE algorithm to compute the minimum
14:	EndFor
15:	Return all the computed minima

To better utilize the proposed approach it is advisable to start the DE algorithm using a mutation operator that permits adequate exploration of the search space (for example DE<sub>2</sub> or DE<sub>6</sub>). Once the clusters around the minima have been determined, one can switch to a mutation operator that has faster convergence speed (for example  $DE_1$ ). Note that instead of using again the DE algorithm in Steps 12 and 13 of Algorithm 1, any other minimization algorithm (even not an evolutionary one) can be employed.

For very hard optimization problems, when the objective function is defined in many dimensions and possesses multitudes of local and global minima, the clustering operator could be called more than once. The same might be true for real–life optimization tasks, where the function value of the global minimum is unknown. Each consecutive call of the clustering operator will result in more promising subregions of the original search space and will save unneeded objective function evaluations, since the subpopulations will stay focused on regions containing desirable minima. For all the experiments conducted in this paper, one call of the clustering operator was sufficient for the algorithm to locate the global, as well as, many local minima.

To determine the applicability and the efficiency of the proposed clustering operator we incorporated it to the DE algorithm and applied the new method to the multimodal test function f, defined in Section 2.2, which posses 9 global minima in the hypercube  $[-5, 5]^2$ . The result of the application of the clustering operator on the function f is illustrated in Figure 5.



Figure 5: Clusters and global minimizers discovered in f

It must be noted that a number of independent experiments of the original DE algorithm gives no guarantee that all global minimizers will be detected, since the algorithm has no memory; no information concerning previously detected minimizers is kept. We performed 100 independent simulations, using each one of the six different mutation operators described above and Table 1 exhibits the average number of restarts needed for the DE algorithms to locate all the global minimizers of f. The modified algorithm that uses the clustering operator in most cases managed to find all the minimizers of f in a single execution.

# **5** Experimental Results

We implemented and tested the proposed clustering operator on a number of hard optimization tasks and it exhibited stable and robust performance. In this section we report the experimental results from 4 well–known minimization test functions. For each mutation operator we performed 100 independent experiments. A population consisting of 200

	Original DE Algorithm		DE with $k$ -win		-win	
	Min	Mean	Max	Min	Mean	Max
$DE_1$	54	110.2	203	1	17.5	66
$DE_2$	57	119.1	294	1	1.0	1
$DE_3$	60	133.0	239	1	1.1	7
$DE_4$	52	114.9	212	1	1.0	1
$DE_5$	49	106.7	245	1	1.0	1
$DE_6$	62	111.4	221	1	1.0	1

Table 1: Restarts needed to locate all the minimizers of f

individuals was used and the mutation and recombination constants had values  $\mu = 0.6$  and  $\rho = 0.8$ , respectively. The algorithm was terminated when the global minimum was located. The proposed clustering operator was called only once for the optimization of each of the four test function considered below, after 20, 20, 10, and 200 generations, respectively.

For each test function used, we present a table (Tables 2– 5) summarizing the average results for the 100 runs. The first column of the table indicates the name of the algorithm and the second column the average number of generations needed for the algorithm to locate the global minimum without the use of the clustering operator. The third and fourth columns give the average number of minima discovered (including the global one) and the corresponding average generations needed for the DE algorithm to locate the global minimum using the clustering operator.

#### 5.1 The Levy No. 5 Test Function

The first test function considered is the Levy No. 5:

$$f_1(\vec{x}) = {}_1 {}_2 + (x_1 + 1.42513)^2 + (x_2 + 0.80032)^2$$

where  $x_i \in [-10, 10], i = 1, 2$ , and  $_1$  and  $_2$  are given by:

$$_{1} = \sum_{i=1}^{5} i \cos[(i+1)x_{1}+i],$$
$$_{2} = \sum_{j=1}^{5} j \cos[(j+1)x_{2}+j].$$

There exist about 760 local minima and one global minimum with function value  $f_1^* = -176.1375$  located at  $\vec{x}^* = (-1.3068, -1.4248)$ . The large number of local optimizers makes it extremely difficult for any method to locate the global minimizer.

	w/o k-win operator	with $k$ -win operator		
	Generations	Minima located	Generations	
$DE_1$	33.21	5.97	34.36	
$DE_2$	70.66	20.52	54.26	
$\mathrm{DE}_3$	64.09	11.96	39.77	
$DE_4$	65.11	20.22	50.25	
$DE_5$	133.01	22.70	70.85	
$DE_6$	64.89	19.20	50.24	

Table 2: Average results for the Levy f	function
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The experimental results exhibited in Table 2 indicate that generally the use of the clustering operator enhances the performance of the DE algorithms. In detail, there is an average acceleration of the algorithm's convergence speed ranging from 30% to 80%. Additionally, as many as 20 minima (including the global one) were simultaneously computed. The only exception is DE<sub>1</sub> where a slight increase in the generations is observed (3%), but the modified algorithm locates the global as well as 5 local minima.

To better demonstrate the ability of the proposed approach to locate many minima at once, independent runs were conducted with the number of individuals in the population gradually increasing from 200 to 2000. In general, a larger population explores better the search space and more regions containing minima are located by the k-windows operator. In Figure 6 we exhibit the detailed results. The algorithm DE<sub>1</sub> locates on average 10 minima regardless of the size of its populations. On the contrary, the rest of the algorithms locate more minima as their population is increased. DE<sub>5</sub> exhibited the best performance finding simultaneously up to 85 minima.



Figure 6: The effect of the population size on the number of located minima for the Levy test function

#### 5.2 The Rastrigin Test Function

The Rastrigin test function is a typical example of a nonlinear multimodal function. This function was first proposed by Rastrigin as a 2-dimensional function [38] and is a fairly difficult problem due to its large number of local minima. It is given by the following equation:

$$f_2(\vec{x}) = \sum_{i=1}^{2} \left( x_i^2 - \cos(18x_i) \right)$$

where  $x_i \in [-1, 1], i = 1, 2$ . There exist 49 local minima.

Table 3 summarizes the average results from the 100 experiments. The results for the Rastrigin function are similar to the results for the Levy function. Faster convergence is exhibited for all the DE algorithms (except  $DE_1$ ). The average increase in convergence speed was from 10% to 130% and up to 20 minima were simultaneously located.

	w/o k-win operator	with k-win operator		
	Generations	Minima located	Generations	
$DE_1$	20.93	1.00	26.34	
$DE_2$	61.67	17.19	45.72	
$DE_3$	80.11	11.23	34.88	
$DE_4$	47.17	13.89	42.64	
$DE_5$	98.69	19.43	54.74	
$DE_6$	55.26	15.72	44.28	

Table 3: Average results for the Rastrigin function

### 5.3 The Shekel's Foxholes Test Function

This is a 2-dimensional test function given by the following equation:

$$f_3(\vec{x}) = \frac{1}{0.002 + \psi(\vec{x})}$$

where  $x_i \in [-65.536, 65.536], i = 1, 2$  and

$$\psi(\vec{x}) = \sum_{l=0}^{24} \frac{1}{1 + l + \sum_{i=1}^{2} (x_i - a_{li})^6}$$

The parameters for this function are:

$$a_{k1} = \{-32, -16, 0, 16, 32\}, \text{ where}$$

$$k = \{0, 1, 2, 3, 4\} \text{ and } a_{k1} = a_{k \mod 5, 1}$$

$$a_{k2} = \{-32, -16, 0, 16, 32\}, \text{ where}$$

$$k = \{0, 5, 10, 15, 20\} \text{ and}$$

$$a_{k2} = a_{k+m,2}, m = \{1, 2, 3, 4\}.$$

The global minimum of  $f_1(-32, -32) = 0.998004$ . This is a relatively easy test function and was included in the experiments in order to investigate the performance of the clustering operator when applied to problems for which the DE algorithm requires a relatively small number of generations to locate the global minimizer.

	w/o k-win operator	with k-win operator		
	Generations	Minima located	Generations	
$DE_1$	7.70	7.81	12.58	
$\mathrm{DE}_2$	12.95	12.58	13.69	
$DE_3$	2.04	9.34	12.43	
$DE_4$	10.75	8.13	12.24	
$DE_5$	11.86	9.87	16.85	
$\mathrm{DE}_6$	2.76	14.33	13.58	

Table 4: Average results for the Shekel's Foxholes function

In Table 4 the average performance of the algorithms is exhibited. It is clear that the use of the clustering operator results in the computation of many minima at once, but the problem is so easy that there is always an increase in the number of generations needed to locate the global minimum. This fact indicates that the proposed approach is better suited to difficult optimization tasks, when more than one minimizer is sought.

### 5.4 The 10-dimensional Griewangk Test Function

This test function is a generalization of the 2-dimensional Griewangk function, which is known to be relatively difficult to minimize and is given by the following equation:

$$f_4(\vec{x}) = \frac{1}{4000} \sum_{i=1}^{10} x_i^2 - \prod_{i=1}^{10} \cos \frac{x_i}{\sqrt{i}} + 1$$

where  $x_i \in [-10, 10], i = 1, 2$ . This test function is riddled with local minima. The global minimum of the function is  $f_4(0, \ldots, 0) = 0$ .

	w/o k-win operator	with k-win operator		
	Generations	Minima located	Generations	
$DE_1$	302.34	14.19	415.55	
$DE_2$	873.93	8.58	346.31	
$DE_3$	799.06	23.36	439.82	
$DE_4$	1280.42	4.40	356.52	
$DE_5$	1816.09	1.65	571.30	
$DE_6$	619.43	18.09	328.31	

Table 5: Average results for the Griewangk function

The experimental results illustrated in Table 5 show that the clustering operator greatly accelerates the DE algorithms ranging from 80% to 260%. Again, DE<sub>1</sub> is the exception; although on average 14 minima are located the proposed algorithm requires 27% additional generations.

# **6** Conclusion

In this work we propose a novel clustering operator and investigate its impact on the performance of the Differential Evolution optimization algorithm. This operator uses the recently proposed unsupervised k-windows clustering algorithm. The k-windows clustering algorithm uses previously computed pieces of information in an attempt to discover regions containing groups of individuals attracted by different minima.

Experiments show that the proposed approach greatly accelerates the convergence speed of the DE algorithms and that in addition to the global minimum is capable to locate simultaneously many local minima without extra function evaluations. To this end, the use of the proposed clustering operator is always suggested. In brief, the clustering operator has the following advantages:

- 1. locates local minima with relatively low function value, in addition to the global ones,
- 2. in general, fewer generations are required for the DE algorithm to converge,
- 3. there is no need for additional function evaluations,
- 4. utilizes the range search algorithm for fast and reliable execution,
- 5. its parallel implementation is straightforward [7],
- 6. is better suited to difficult high-dimensional multimodal objective functions.

In a future correspondence, we will present an adaptive scheme, based on the average movement of each individual, for the automatic call of the clustering operator. This scheme will ensure that the search space is sufficiently explored. We also intend to investigate the parallel implementation of the proposed approach, as well as its performance on difficult high-dimensional real-life problems encountered in bioinformatics, medical applications and neural network training.

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