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# Differential Evolution Via Exploiting Opposite Populations

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**Summary.** The concept of opposition can contribute to improve the performance of population-based algorithms. This chapter presents an overview of a novel opposition-based scheme to accelerate an evolutionary algorithm, differential evolution (DE). The proposed opposition-based DE (ODE) employs opposition-based computation (OBC) for population initialization and also for generation jumping. Opposite numbers, representing anti-chromosomes, have been utilized to improve the convergence rate of the classical DE. A test suite with 15 well-known benchmark functions is employed for experimental verification. Descriptions for the DE and ODE algorithms, and a comparison strategy are provided. Results are promising and confirm that the ODE outperforms its parent algorithm DE. This work can be regarded as an initial study to exploit oppositional concepts to expedite the optimization process for any population-based approach.

## 8.1 Introduction

Evolutionary algorithms (EAs) are well-established techniques to approach problems with mixed-type variables, many local optima, and with undifferentiable or non-analytical functions [1]. Among various kinds of evolutionary algorithms, differential evolution (DE) is well known for its effectiveness and robustness. Many comparative studies confirm that the DE outperforms many other optimizers [5]. Finding more accurate solution(s) in a shorter period of time for complex black-box problems is still a crucial target of research on evolutionary algorithms.

In this chapter, opposition-based schemes including opposition-based population initialization and generation jumping, will be described. The differential evolution (DE) is selected as a parent algorithm to verify the acceleration effect of the proposed schemes. A set of well-known complex benchmark functions is employed to experimentally compare and analyze the algorithms. Results confirm that Opposition-Based Differential Evolution (ODE) performs better than DE in terms of convergence speed and solution accuracy.

The main purpose of this and previous works has been to introduce a new notion into nonlinear continuous optimization via innovative metaheuristics, namely *the notion of opposition*. Although, all conducted experiments utilize DE as a parent algorithm, the proposed schemes are defined at the population level and, hence, have an inherent potential to be utilized for acceleration of other population-based algorithms.

The organization of this chapter is as follows: A short review of differential evolution is given in section 8.2. The main reasons to select DE as a parent algorithm are explained in section 8.3. Opposition-based differential evolution is described in section 8.4. Experimental verifications are elaborated in section 8.5. Finally, the chapter is concluded in section 8.6.

## 8.2 Differential Evolution (DE)

Differential evolution (DE) is a population-based optimization algorithm based on the idea of genetic annealing which was used to solve the Chebyshev polynomial fitting problem [1]. In order to solve the Chebyshev problem in continuous space, a modified genetic annealing algorithm from bit-string to floating-point encoding and a consequent switch from logical operators to arithmetic ones were proposed [2, 3, 4]. During these experiments, the differential mutation operator to perturb the population of vectors was discovered. Additionally, by using differential mutation, discrete recombination, and pair-wise selection, it was recognized that an annealing mechanism is not needed; it was removed completely and DE was born.

Let us assume that  $X_{i,G}(i = 1, 2, \dots, N_p)$  are candidate solution vectors in generation  $G$  ( $N_p$  : population size). Like other evolutionary algorithms, DE starts with an initial population, which is usually generated in a random manner. A typical vector of the initial population can be generated as follows [5]:

$$X_{i,j} = l_j + \text{RAND}_j(0, 1) \times (l_j - u_j) \quad \text{with } j = 1, 2, \dots, D, \quad (8.1)$$

where  $D$  is the problem dimensionality;  $l_j$  and  $u_j$  are the lower and the upper boundaries of the  $j^{\text{th}}$  variable, respectively, and  $\text{RAND}(0, 1)$  is a uniformly generated random number in  $[0, 1]$ .

Successive populations are generated by adding the weighted difference of two randomly selected vectors to a third randomly selected vector. For classical DE (see Algorithm 1), the mutation, crossover, and selection operators are straightforwardly defined.

### 8.2.1 Mutation

For each vector  $X_{i,G}$  in generation  $G$  a mutant vector  $V_{i,G}$  (see line 9 of Algorithm 1) is defined by

$$V_{i,G} = X_{a,G} + F(X_{c,G} - X_{b,G}), \quad (8.2)$$

where  $i = \{1, 2, \dots, N_p\}$ , and  $a$ ,  $b$ , and  $c$  are mutually different random integer indices selected from  $\{1, 2, \dots, N_p\}$ . Further, the variables  $i$ ,  $a$ ,  $b$ , and  $c$  are different so that  $N_p \geq 4$  is necessary. The factor  $F \in [0, 2]$  is a real constant which determines the amplification of the added differential variation of  $(X_{c,G} - X_{b,G})$  [5]. Larger values for  $F$  result in higher diversity in the generated population and lower values cause faster convergence.

### 8.2.2 Crossover

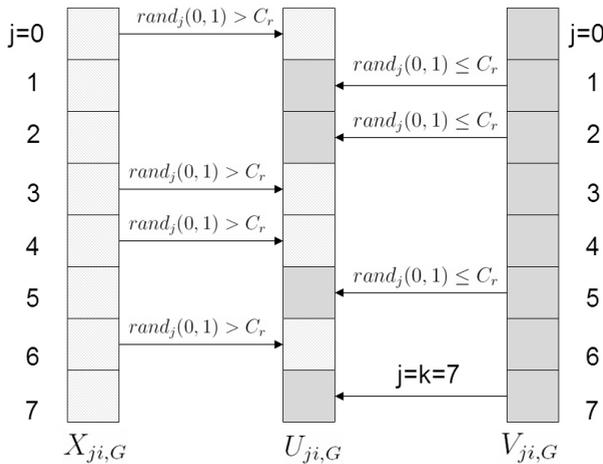
DE utilizes the crossover operation to generate new solutions by shuffling competing vectors and also to increase the population diversity. For the classical DE (lines 10 – 16 of Algorithm 1), the binary crossover is utilized. It defines the following trial vector:

$$U_{i,G} = (U_{1i,G}, U_{2i,G}, \dots, U_{Di,G}), \tag{8.3}$$

where

$$U_{ji,G} = \begin{cases} V_{ji,G} & \text{if } \text{RAND}_j(0, 1) \leq C_r \vee j = k, \\ X_{ji,G} & \text{otherwise.} \end{cases} \tag{8.4}$$

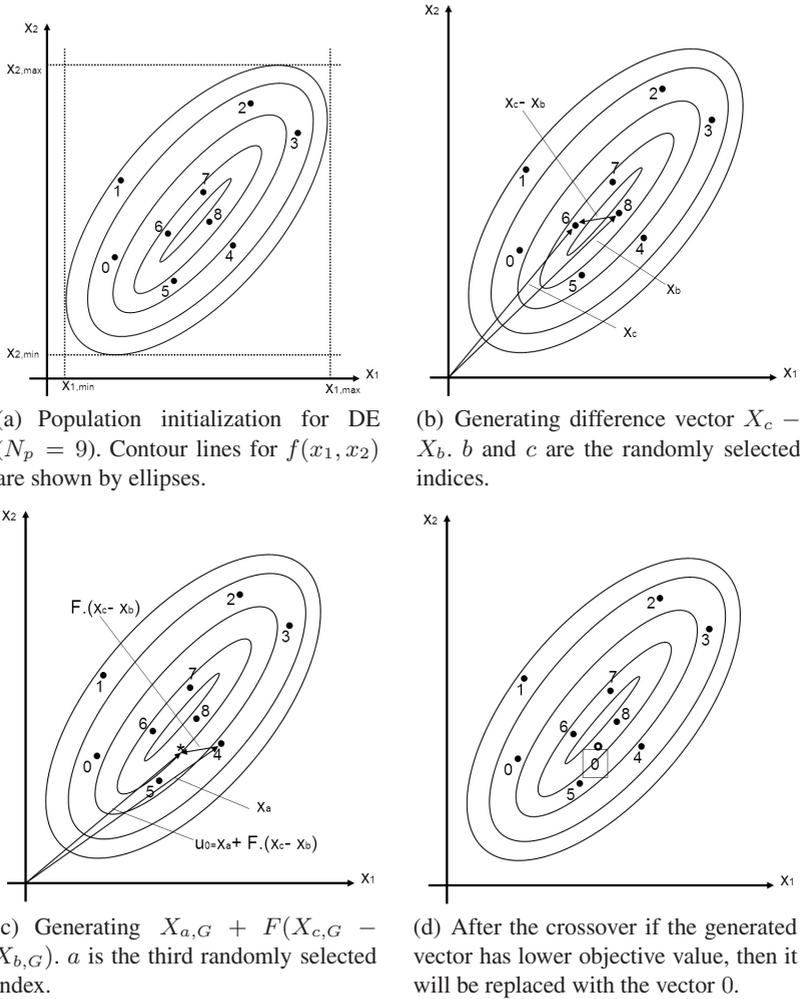
$C_r \in (0, 1)$  is the predefined crossover rate, and  $\text{RAND}_j(0, 1)$  is the  $j^{\text{th}}$  evaluation of a uniform random number generator. The parameter  $k \in \{1, 2, \dots, D\}$  is a random index chosen once for each  $i$  to make sure that at least one parameter is always selected from the mutated vector  $V_{ji,G}$ . The most common values for  $C_r$  are in the range of  $(0.4, 1)$  [17]. Figure 8.1 illustrates a pictorial example for the binary crossover.



**Fig. 8.1.** A pictorial example for the binary crossover in DE ( $k = 7$ ) [5]. When  $\text{RAND}_j(0, 1) \leq C_r \vee j = k$ , then the variable is copied from  $V_{ji,G}$ , otherwise copied from  $X_{ji,G}$  to  $U_{ji,G}$ .

### 8.2.3 Selection

Selection is a mechanism to decide which vector ( $U_{i,G}$  or  $X_{i,G}$ ) should be a member of next (new) generation  $G + 1$ . For a minimization problem, the vector with the lower objective function value is chosen (greedy selection). If  $f(U_i) \leq f(X_i)$ , then  $U_i$  is selected; otherwise  $X_i$  will be chosen (lines 17 – 23 of Algorithm 1).



**Fig. 8.2.** Illustration of one generate-and-test cycle for DE (starting from vector 0) [5]

This evolutionary cycle (i.e., mutation, crossover, and selection) is repeated  $N_p$  times to generate new populations. These successive generations are produced until the termination conditions are satisfied. One generate-and-test cycle for DE is presented in Figure 8.2.

The starting point for the mutation, crossover and selection is indicated by the comments in the algorithm. The algorithm terminates (line 5) when the best achieved fitness value (BFV) is smaller than the value-to-reach (VTR), or the number of function calls (NFC) exceeds the predefined maximum number of function calls ( $MAX_{NFC}$ ). The termination strategy can be defined differently based on the application or the purpose of the experiment. The number of generations, the execution time, or some population

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**Algorithm 1.** Differential Evolution (DE).  $P_0$ : Initial population,  $N_p$ : Population size,  $V$ : Noise vector,  $U$ : Trial vector,  $D$ : Problem dimension, BFV: Best achieved fitness value, VTR: Value-to-reach, NFC: Number of function calls,  $\text{MAX}_{\text{NFC}}$ : Maximum number of function calls,  $F$ : Mutation constant,  $\text{RAND}(0, 1)$ : Uniformly generated random number,  $C_r$ : Crossover,  $f(\cdot)$ : Objective function,  $P'$ : Population of the next generation.

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1: Generate uniformly distributed random population  $P_0$ 
2:  $\text{NFC} \leftarrow 0$ 
3: Evaluate individuals of  $P_0$ 
4:  $\text{NFC} \leftarrow \text{NFC} + N_p$ 
5: while (  $\text{BFV} > \text{VTR}$  and  $\text{NFC} < \text{MAX}_{\text{NFC}}$  ) do
6:   {Generate-and-Test-Loop}
7:   for  $i = 0$  to  $N_p$  do
8:     Select three parents  $X_a$ ,  $X_b$ , and  $X_c$  randomly from current population where  $i \neq a \neq b \neq c$ 
     {Mutation}
9:      $V_i \leftarrow X_a + F \times (X_c - X_b)$ 
     {Crossover}
10:    for  $j = 0$  to  $D$  do
11:      if  $\text{RAND}(0, 1) < C_r$  then
12:         $U_{i,j} \leftarrow V_{i,j}$ 
13:      else
14:         $U_{i,j} \leftarrow X_{i,j}$ 
15:      end if
16:    end for
     {Selection}
17:    Evaluate  $U_i$ 
18:     $\text{NFC} \leftarrow \text{NFC} + 1$ 
19:    if ( $f(U_i) \leq f(X_i)$ ) then
20:       $X'_i \leftarrow U_i$ 
21:    else
22:       $X'_i \leftarrow X_i$ 
23:    end if
24:  end for
25:   $X \leftarrow X'$ 
26: end while

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statistics (e.g., diversity or the improvement rate) are some commonly used termination criteria.

### 8.2.4 DE in Optimization Field

A summary classification of optimization methods can be seen in Figure 8.3. According to the proposed classification scheme for optimization methods, DE is a population-based, nonlinear, continuous and global optimization algorithm [1].

Studies have been conducted to enhance the performance of the classical DE algorithm by adaptive determination of DE control parameters. For instance, the fuzzy adaptive differential evolution algorithm (FADE) was introduced by Liu and Lampinen

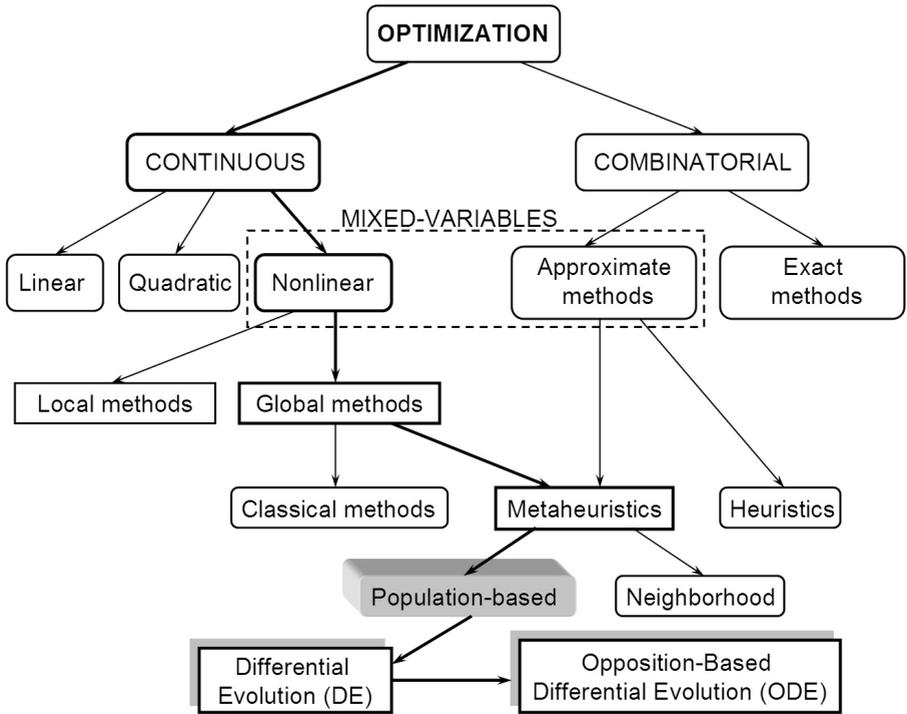


Fig. 8.3. A simple classification scheme of optimization methods [1]

[24]. They employed a fuzzy logic controller to set the mutation and crossover rates. In the same direction, Brest et al. [20] proposed self-adaptive DE. Teo [30] proposed a dynamic population sizing strategy based on self-adaptation, and Ali and Törn [25] introduced auxiliary population and automatic calculating of the amplification factor  $F$  for the difference vector.

Other researchers have experimented with multi-population ideas. Tasoulis et al. [31] proposed parallel DE where they assign each subpopulation to a different processor node. Shi et al. [32] partitioned high-dimensional search spaces into smaller spaces and used multiple cooperating subpopulations to find the solution. They called this method cooperative co-evolutionary differential evolution.

Hybridization with different algorithms is another direction for improvement of DE. Sun et al. [26] proposed a new hybrid algorithm based on a combination of DE and estimation of distribution algorithm. This technique uses a probability model to determine promising regions in order to focus the search process on those areas. Noman and Iba [35] incorporated local search into the classical DE. They employed fittest-individual refinement which is a crossover-based local search. Fan and Lampinen [33] introduced a new local search operation, trigonometric mutation, in order to obtain a better trade-off between convergence speed and robustness. Kaelo and Ali [34] employed reinforcement learning and some other schemes for generating fitter trial points.

All mentioned approaches are proposed to increase the convergence rate and/or the accuracy of DE. For more details about DE extensions, the reader is referred to literature [1, 5, 27].

### 8.3 Why Differential Evolution?

Differential evolution is a simple and compact metaheuristic which directly operates on continuous variables (arithmetic operators instead of logical operators). Unlike many binary versions of the genetic algorithms, DE works with the floating-point numbers. This removes encoding and decoding of the variables which is a source of inaccuracy. Consequently, this feature makes DE scalable for high-dimensional problems and also time and memory efficient.

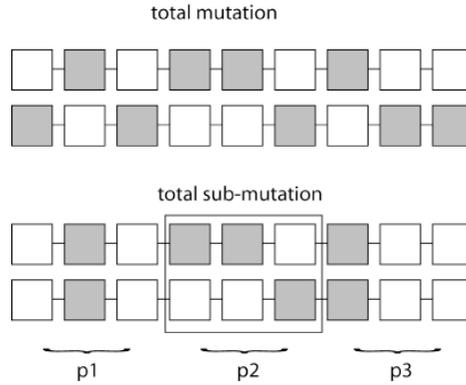
Another reason for choosing DE is that it does not need a probability density function to adapt the control parameters (unlike most evolutionary strategies) or any probability distribution pattern for the mutation (unlike genetic algorithms or evolutionary programming). DE's different mutation and crossover schemes distinguish it from other evolutionary algorithms [5].

Additionally, handling mixed integers, discrete and continuous variables makes DE more applicable for a wider range of real-world applications. The main advantage of DE while working with integer variables is that it internally works on a continuous space and only switches to the integer space during the evaluation of the objective function. This characteristic supports higher accuracy compared to some other well-known algorithms (e.g., GAs) which perform in the reverse manner [27]. Extensions of classical DE are capable of handling boundary constraints and also nonlinear function constraints which both are commonly required in the real-world problems [5].

Many comparative studies report higher robustness, convergence speed, and solution quality of the DE when compared to other evolutionary algorithms for both benchmark functions and real-world applications. A comprehensive performance study is provided in [5]. The authors first compare DE to 16 other optimizers against five well-known thirty-dimensional test functions (namely, Rosenbrock, Ackley, Griewangk, Rastrigin, and Schwefel). Consequently, they explored eight function-based comparative studies (e.g., unconstrained optimization, multi-constraints nonlinear optimization, and multi-objective optimization) and also eleven application-oriented performance comparison studies (e.g., multi-sensor fusion, earthquake relocation, image registration, and optimization of neural networks). Finally, they conclude [5] “[...] DE may not always be the fastest method, it is usually the one that produces the best results, although the number of cases in which it is also the faster is significant. DE also proves itself to be robust, both in how control parameters are chosen and in the regularity with which it finds the true optimum. [...] As these researchers have found, DE is a good first choice when approaching a new and difficult global optimization problem is defined with continuous and/or discrete parameters.”

### 8.4 Opposition-Based Differential Evolution

In his primary paper on opposition-based learning (OBL), Tizhoosh proposed to use *anti-chromosomes* for GAs [12]. For every selected chromosome a corresponding



**Fig. 8.4.** Generation of anti-chromosomes [12]

anti-chromosome can be generated. The initial chromosomes are generally generated randomly meaning that they can possess high or low fitness. However, in a complex problem it is usually very likely that the initial populations do not contain optimal solutions. Hence, in lack of any a-priori knowledge, it is reasonable to look at anti-chromosomes simultaneously. Considering the search direction and its opposite at the same time will bear more likelihood to reach the best population in a shorter time (for more motivation on OBL see also Chapters 1-2). Tizhoosh also suggested to use total or sub-total-mutation to generate opposite candidate solutions (Figure 8.4).

Similar to all population-based optimization algorithms, two main steps are distinguishable for DE, namely population initialization and producing new generations by evolutionary operations such as mutation, crossover, and selection. The opposition-based differential evolution (ODE) [6, 9, 13] will enhance these two steps by considering opposite solutions. For black-box optimization – which is a general assumption for optimization methods – there is no information about the shape of the problem landscape such that type II opposition can only be approximated via type I opposition (see Definitions 2 and 6 in Chapter 2). The pseudo-code of ODE is presented in Algorithm 2 [13].

### 8.4.1 Opposition-Based Population Initialization

In absence of domain knowledge, uniform random number generation is generally the only choice to create an initial population. But as mentioned before, by utilizing type-I opposition it is possible to obtain fitter starting candidates. Block (1) from Figure 8.5 shows the implementation of opposition-based population initialization (lines 5 – 12 of Algorithm 2). The following steps explain that procedure [6]:

- Step 1.* Initialize the population  $P(N_p)$  randomly,
- Step 2.* Create the opposite population OP by

$$OP_{i,j} = a_j + b_j - P_{i,j}, \quad \text{with } i = 1, 2, \dots, N_p; j = 1, 2, \dots, D, \quad (8.5)$$

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**Algorithm 2.** Pseudo-code of Opposition-Based Differential Evolution (ODE) in order to solve a minimization problem [Adopted from [13]].  $P_0$ : Initial population,  $OP_0$ : Opposite of initial population,  $P$ : Current population,  $OP$ : Opposite of current population,  $D$ : Problem dimension,  $[l_j, u_j]$ : Range of the  $j$ -th variable,  $J_r$ : Jumping rate,  $C_r$ : Crossover rate,  $\min_j^P/\max_j^P$ : Minimum/maximum value of the  $j$ -th variable in the current population. Lines **1-12** and **33-42** are implementations of opposition-based population initialization and opposition-based generation jumping, respectively.

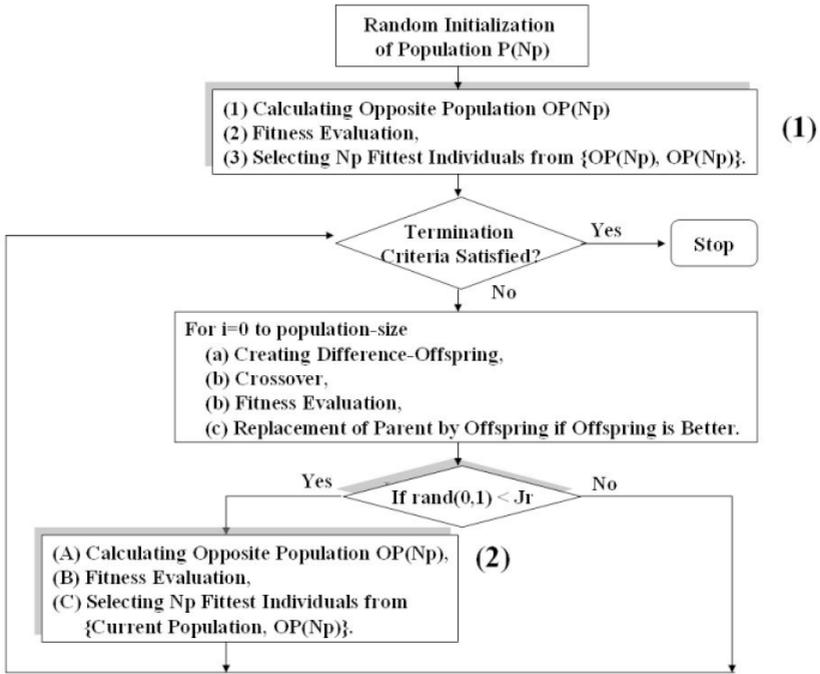
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1: Generate uniformly distributed random population  $P_0$ 
2:  $NFC \leftarrow 0$ 
3: Evaluate individuals of  $P_0$ 
4:  $NFC \leftarrow NFC + N_p$ 
   {**Begin of Opposition-Based Population Initialization**}
5: for  $i = 0$  to  $N_p$  do
6:   for  $j = 0$  to  $D$  do
7:      $OP_{0i,j} \leftarrow l_j + u_j - P_{0i,j}$ 
8:   end for
9: end for
10: Evaluate individuals of  $OP_0$ 
11:  $NFC \leftarrow NFC + N_p$ 
12: Select  $N_p$  fittest (best) individuals from  $P_0$  and  $OP_0$  as initial population  $P_0$ 
   {Begin of DE's Evolution Steps}
13: while ( $BFV > VTR$  and  $NFC < MAX_{NFC}$ ) do
14:   for  $i = 0$  to  $N_p$  do
15:     Select three parents  $P_{i_1}, P_{i_2},$  and  $P_{i_3}$  randomly from current population where
        $i \neq i_1 \neq i_2 \neq i_3$ 
16:      $V_i \leftarrow P_{i_1} + F \times (P_{i_2} - P_{i_3})$ 
17:     for  $j = 0$  to  $D$  do
18:       if  $RAND(0, 1) < C_r$  then
19:          $U_{i,j} \leftarrow V_{i,j}$ 
20:       else
21:          $U_{i,j} \leftarrow P_{i,j}$ 
22:       end if
23:     end for
24:     Evaluate  $U_i$ 
25:      $NFC \leftarrow NFC + 1$ 
26:     if ( $f(U_i) \leq f(P_i)$ ) then
27:        $P'_i \leftarrow U_i$ 
28:     else
29:        $P'_i \leftarrow P_i$ 
30:     end if
31:   end for
32:    $P \leftarrow P'$ 
   {**Begin of Opposition-Based Generation Jumping**}
33:   if  $RAND(0, 1) < J_r$  then
34:     for  $i = 0$  to  $N_p$  do
35:       for  $j = 0$  to  $D$  do
36:          $OP_{i,j} \leftarrow MIN_j^P + MAX_j^P - P_{i,j}$ 
37:       end for
38:     end for
39:     Evaluate individuals of  $OP_0$ 
40:      $NFC \leftarrow NFC + N_p$ 
41:     Select  $N_p$  fittest (best) individuals from  $P$  and  $OP$  as current population  $P$ 
42:   end if
43: end while

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**Fig. 8.5.** Gray boxes extend DE to ODE. Block (1): Opposition-based initialization, Block (2): Opposition-based generation jumping ( $J_r$ : jumping rate,  $RAND(0, 1)$ : uniformly generated random number,  $N_p$ : population size).

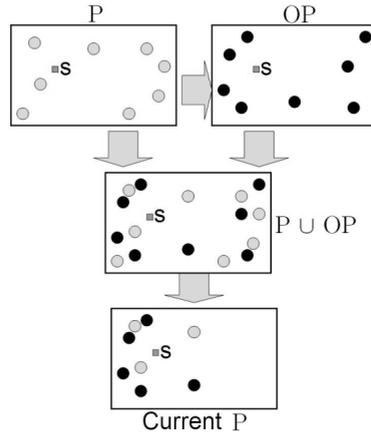
where  $P_i$  and  $OP_i$  denote the  $i^{th}$  individual of the current population and its corresponding opposite, respectively, and  $[l_j, u_j]$  is the range of the  $j^{th}$  variable.

*Step 3.* Select the  $N_p$  fittest (best) individuals from  $P \cup OP$  as the initial population.

According to the above procedure,  $2N_p$  function evaluations are required instead of  $N_p$  for the regular random population initialization. But, by the opposition-based initialization, the parent algorithm can start with the fitter initial individuals instead.

### 8.4.2 Opposition-Based Generation Jumping

By applying a similar approach mentioned in Sec. 8.4.1 to the current population, which means selecting  $N_p$  best individuals from the current and corresponding opposite populations, the evolutionary process can be forced to jump to a fitter generation (the generation with fitter individuals). After generating new populations, the opposite population is calculated and the  $N_p$  fittest (best) individuals are selected from the union of the current and opposite population based on a jumping rate  $J_r \in (0, 0.4)$  [13, 15]. In order to calculate the opposite population for generation jumping, the opposite of each variable is calculated dynamically; that is, the maximum and minimum values of each variable



**Fig. 8.6.** Example to visualize the opposition-based generation jumping in 2D space ( $N_p = 8$ )

in the current population ( $[\text{MIN}_j^p, \text{MAX}_j^p]$ ) are used to calculate opposite points instead of using variables' predefined interval boundaries ( $[l_j, u_j]$ ):

$$\text{OP}_{i,j} = \text{MIN}_j^p + \text{MAX}_j^p - P_{i,j}, i = 1, 2, \dots, N_p; j = 1, 2, \dots, D. \quad (8.6)$$

If the opposites are calculated within variables' static boundaries, it is possible to jump outside of the already shrunken search space and lose the knowledge of the already reduced space. Hence, we calculate opposite points by using variables' current interval in the population ( $[\text{MIN}_j^p, \text{MAX}_j^p]$ ) which is, as the search does progress, increasingly smaller than the corresponding initial range  $[l_j, u_j]$ . Block (2) from Figure 8.5 illustrates the implementation of opposition-based generation jumping (lines 33 – 42 of Algorithm 2).

A pictorial example for opposition-based generation jumping procedure in 2D space is illustrated in Figure 8.6. The letter 'S' indicates the location of the optimal solution. Dark and light circles represent the points and the opposite points, respectively.

In [14], we established mathematical proofs and experimental evidence to verify the advantage of opposite points compared to additional random points when dealing with high-dimensional problems (see also Chapter 2 for more discussions on the formalism of opposition). Both experimental and mathematical results confirmed that opposite points are more beneficial than additional independent random points. We can conclude that the opposition-based learning can be utilized to accelerate optimization methods since considering the pair  $x$  and its opposite  $\check{x}$  has apparently a higher fitness probability than pure randomness.

## 8.5 Experimental Verifications

In this section, the convergence metrics are defined and DE and ODE are compared experimentally over well-known benchmark functions (section 8.5.1). Also, the

contribution of opposite points to the achieved acceleration rate is investigated by replacing them with random points (section 8.5.2).

### 8.5.1 Comparison of DE and ODE

A set of 15 well-known benchmark functions [6, 13, 15], which contains 7 unimodal ( $f_1, f_2, f_3, f_6, f_{10}, f_{11}, f_{14}$ ) and 8 multimodal functions ( $f_4, f_5, f_7, f_8, f_9, f_{12}, f_{13}, f_{15}$ ), has been selected for performance verification of ODE. The definition of the benchmark functions is given in Table 8.1.

**Table 8.1.** List of employed benchmark functions (unimodal and multimodal)

Function	Search Space
$f_1(X) = \sum_{i=1}^D x_i^2$	$[-5.12, 5.12]^D$
$f_2(X) = \sum_{i=1}^D ix_i^2$	$[-5.12, 5.12]^D$
$f_3(X) = \sum_{i=1}^D \left( \sum_{j=1}^i x_j \right)^2$	$[-65, 65]^D$
$f_4(X) = 10D + \sum_{i=1}^D (x_i^2 - 10 \cos(2\pi x_i))$	$[-5.12, 5.12]^D$
$f_5(X) = \sum_{i=1}^D \frac{x_i^2}{4000} - \prod_{i=1}^D \cos\left(\frac{x_i}{\sqrt{i}}\right) + 1$	$[-600, 600]^D$
$f_6(X) = \sum_{i=1}^D  x_i ^{(i+1)}$	$[-1, 1]^D$
$f_7(X) = -20 \exp\left(-0.2 \sqrt{\frac{D \sum_{i=1}^D x_i^2}{D}}\right) - \exp\left(\frac{\sum_{i=1}^D \cos(2\pi x_i)}{D}\right) + 20 + e$	$[-32, 32]^D$
$f_8(X) = \sin^2(3\pi x_1) + \sum_{i=1}^{D-1} (x_i - 1)^2 (1 + \sin^2(3\pi x_{i+1})) + (x_D - 1)(1 + \sin^2(2\pi x_D))$	$[-10, 10]^D$
$f_9(X) = -\sum_{i=1}^D \sin(x_i) (\sin(ix_i^2/\pi))^{2m}, (m = 10)$	$[0, \pi]^D$
$f_{10}(X) = \sum_{i=1}^D x_i^2 + \left(\sum_{i=1}^D 0.5ix_i\right)^2 + \left(\sum_{i=1}^D 0.5ix_i\right)^4$	$[-5, 10]^D$
$f_{11}(X) = \sum_{i=1}^D  x_i  + \prod_{i=1}^D  x_i $	$[-10, 10]^D$
$f_{12}(X) = \sum_{i=1}^D (\lfloor  x_i  + 0.5 \rfloor)^2$	$[-100, 100]^D$
$f_{13}(X) = \sum_{i=1}^D  x_i \sin(x_i) + 0.1x_i $	$[-10, 10]^D$
$f_{14}(X) = \exp\left(-0.5 \sum_{i=1}^D x_i^2\right)$	$[-1, 1]^D$
$f_{15}(X) = 1 - \cos(2\pi \ x\ ) + 0.1 \ x\ $ , where $\ x\  = \sqrt{\sum_{i=1}^D x_i^2}$	$[-100, 100]^D$

We compare the convergence speed of DE and ODE by measuring the number of function calls (NFC) which is the most commonly used metric in literature [5, 7, 8, 9, 10, 11, 19]; a smaller NFC means higher convergence speed. The termination criterion is to find a value smaller than the value-to-reach (VTR) before reaching the maximum number of function calls  $\text{MAX}_{\text{NFC}}$ . In order to minimize the effect of the stochastic nature of the algorithms on the metric, the reported number of function calls (NFC) for

**Table 8.2.** Parameter settings for conducted experiments

<i>Parameter name</i>	<i>Setting</i>	<i>Reference</i>
population size ( $N_p$ )	100	[20, 21, 22]
differential amplification factor ( $F$ )	0.5	[11, 20, 23, 24, 25]
crossover probability constant ( $C_r$ )	0.9	[11, 20, 23, 24, 25]
jumping rate constant ( $J_r$ )	0.3	[9, 13, 18]
maximum number of function calls ( $\text{MAX}_{\text{NFC}}$ )	$10^6$	[9, 13, 18]
value to reach (VTR)	$10^{-8}$	[10]
mutation strategy	$DE/rand/1/bin$	[5, 20, 23, 26, 27]

**Table 8.3.** Comparison of DE, ODE, and RDE. The best result for each case is highlighted in boldface. Results for RDE have been discussed in section 8.5.2 (corresponding results for replacing the opposite points with random points).

<i>F</i>	<i>D</i>	<i>DE</i>			<i>ODE</i>			<i>RDE</i>		
		<i>NFC</i>	<i>SR</i>	<i>SP</i>	<i>NFC</i>	<i>SR</i>	<i>SP</i>	<i>NFC</i>	<i>SR</i>	<i>SP</i>
$f_1$	30	87748	1	87748	47716	1	<b>47716</b>	115096	1	115096
$f_2$	30	96488	1	96488	53304	1	<b>53304</b>	126780	1	126780
$f_3$	20	177880	1	177880	168680	1	<b>168680</b>	231152	1	231152
$f_4$	10	328844	1	328844	70389	0.76	<b>92617</b>	501875	0.96	522786
$f_5$	30	113428	1	113428	69342	0.96	<b>72231</b>	149744	1	149744
$f_6$	30	25140	1	25140	8328	1	<b>8328</b>	29096	1	29096
$f_7$	30	169152	1	169152	98296	1	<b>98296</b>	222784	1	222784
$f_8$	30	101460	1	101460	70408	1	<b>70408</b>	138308	1	138308
$f_9$	10	191340	0.76	<b>251763</b>	213330	0.56	380946	306900	0.60	511500
$f_{10}$	30	385192	1	385192	369104	1	<b>369104</b>	498200	1	498200
$f_{11}$	30	187300	1	187300	155636	1	<b>155636</b>	244396	1	244396
$f_{12}$	30	41588	1	41588	23124	1	<b>23124</b>	54316	1	54316
$f_{13}$	30	411164	1	411164	337532	1	<b>337532</b>	927230	0.24	3863458
$f_{14}$	10	19528	1	19528	15704	1	<b>15704</b>	23156	1	23156
$f_{15}$	10	37824	1	37824	24260	1	<b>24260</b>	46800	1	46800
$\text{SR}_{\text{ave}}$		0.98			0.95			0.92		

each function is the average over 50 different trials. The number of times, for which the algorithm successfully reaches the VTR for each test function is measured as the success rate SR:

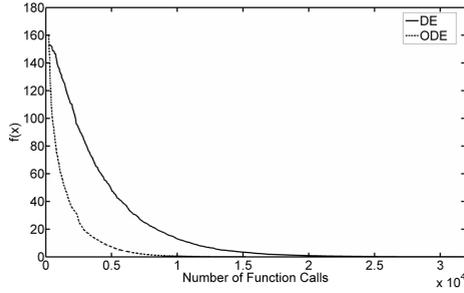
$$\text{SR} = \frac{\text{number of times reached VTR}}{\text{total number of trials}}. \quad (8.7)$$

In order to combine these two measures (NFC and SR), a new measure, called success performance has been introduced as follows [10]:

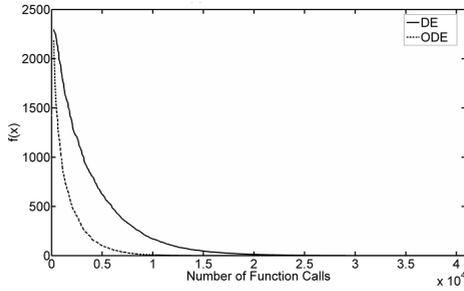
$$\text{SP} = \frac{\text{average of NFC over successful runs}}{\text{SR}}. \quad (8.8)$$

The parameter setting for all conducted experiments is summarized in Table 8.2.

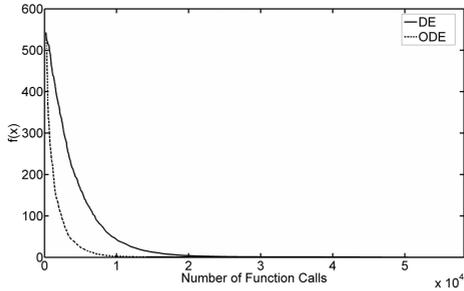
In order to maintain a reliable and fair comparison, the parameter settings are kept the same for all conducted experiments, unless we mention new settings. Besides, for all experiments, the reported values are the average of the results for 50 independent runs. In addition, and most importantly, extra fitness evaluations required for the



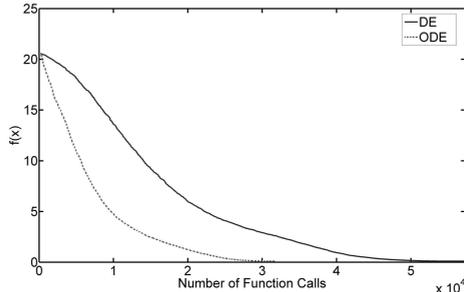
(a)  $f_1$ , ODE is 1.83 times faster



(b)  $f_2$ , ODE is 1.81 times faster



(c)  $f_5$ , ODE is 1.63 times faster



(d)  $f_7$ , ODE is 1.72 times faster

**Fig. 8.7.** Sample convergence graphs (best solution vs. number of function calls). As seen, ODE (dotted curve) shows better convergence speed than DE (solid curve) because it needs small amount of function calls to find the solution. is calculated by  $\frac{NFC_{DE}}{NFC_{ODE}}$ .

opposite points (both in population initialization and also generation jumping phases) are counted as well to accurately measure the benefit in spite of the additional overhead for computing the opposites.

The results for DE and ODE to solve the test problems are given in Table 8.3 (the results in the last column will be discussed in section 8.5.2). ODE outperforms DE on 14 benchmark functions with respect to the success performance. Some sample performance comparison graphs are presented in Figure 8.7. ODE (dotted curve) shows better convergence speed than DE (solid curve) because it needs smaller number of function calls to find the solution. With the same parameter settings for both algorithms and fixing the jumping rate for the ODE ( $J_r = 0.3$ ), their success rates are comparable while ODE shows better convergence speed than DE. DE has a better success rate (SR) than ODE on 3 functions ( $f_4, f_5$ , and  $f_9$ ). The jumping rate is an important control parameter which, if optimally set, can achieve even better results. Detailed discussions about this parameter can be found in [13].

On 7 multimodal functions (out of 8), ODE performs better than DE. This means that the opposition-based extension performs well even when the function contains many optima.

### 8.5.2 Contribution of Opposite Points

In this section, we verify whether the achieved acceleration rate for DE is really due to utilizing opposite points. For this purpose, all parts of the proposed algorithm remain unchanged and instead of using opposite points for the population initialization and the generation jumping, uniformly generated random points will be employed. In order to have a fair competition for this case, exactly like what we did for opposite points, the current interval (dynamic interval,  $[\text{MIN}_j^p, \text{MAX}_j^p]$ ) of the variables are used to generate new random points in the generation jumping phase. So, line 4 in Algorithm 2 should be changed to:

$$RP_{0i,j} \leftarrow l_j + (u_j - l_j) \times \text{RAND}(0, 1),$$

where  $\text{RAND}(0, 1)$  generates a uniformly distributed random number on the interval  $(0, 1)$ . In fact, instead of generating  $N_p$ ,  $2N_p$  random individuals are generated. In the same manner, line 30 in Algorithm 2 should be replaced with

$$RP_{i,j} \leftarrow \text{MIN}_j^p + (\text{MAX}_j^p - \text{MIN}_j^p) \times \text{RAND}(0, 1).$$

As mentioned before, the current boundaries of the variables ( $[\text{MAX}_j^p, \text{MIN}_j^p]$ ) are used to generate random numbers for generation jumping. And finally, in order to have the same selection method, lines 7 and 33 in Algorithm 2 are substituted with

*Select  $N_p$  fittest (best) individuals from  $P$  and  $RP$  as current population  $P$ ;*

After these modifications, the random version of ODE (called RDE) is established. Results for the current algorithm are presented in Table 8.3. As apparent, RDE can

not outperform DE or ODE on any of benchmark function with respect to the success performance. This clearly demonstrates that the achieved improvements are due to usage of opposite points, and that the same level of improvement cannot be achieved via additional random sampling [14, 15].

## 8.6 Conclusions and Future Work

In this chapter, we briefly reviewed how opposition-based optimization can be employed to accelerate convergence speed of differential evolution by embedding opposition-based population initialization and opposition-based generation jumping. The experimental results confirmed that ODE provides a higher performance than the classical DE. However, opposition-based optimization is still in its infancy and future research is required to fully investigate its benefits and drawbacks.

By replacing opposite points with uniformly generated random points in the same variables' range, the resulted algorithm (RDE) performs slower than the parent algorithm (DE). Therefore, the contribution of opposite points to the acceleration process was confirmed and was not reproducible by additional random sampling.

The benefits of opposition-based optimization is most likely not the same for different problems. This is because of using fixed settings for the parameters and/or the different characteristics of each problem (e.g., modality, dimension, surface features, separability of the variables and so on). Similar to all optimization approaches, ODE does not present a consistent behavior over different problems. However, over the employed benchmark test suite, ODE performed better than classical DE.

The proposed opposition-based schemes are general enough to be applied to other population-based algorithms. The opposition-based schemes work at the population level and leave the evolutionary part of the algorithms untouched. This generality gives higher flexibility to these schemes to be embedded inside other population-based algorithms.

Opposition-based optimization opens new perspectives to accelerate optimization processes. For most practical applications, we are faced with constrained functions and also with multi-objective problems. So far, there are many approaches for handling constraints in DE and also for multi-objective optimization using DE. All of these proposals can be borrowed and investigated to generalize ODE to solve multi-objective constrained problems.

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