



Enhanced decomposition-based hybrid evolutionary and gradient-based algorithm for many-objective optimization

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Abstract

This paper presents a novel decomposition-based hybrid many-objective optimization method using particle swarm optimization (PSO) and sequential quadratic programming (SQP) algorithms. In the proposed method, the objective function space decomposed into optimization sub-problems. In this respect, the origin of the coordinate systems of the scaled objective function space (i.e., ideal point) and uniformly distributed points in the scaled objective function space are used to generate reference directions. The scaling of the objective function space in each generation leads to the improvement of the accuracy of the ideal point, which is important because of its role in the generation of the reference directions. In the proposed method, optimization of all the reference directions performed simultaneously to increase the accuracy as well as the computational efficiency. In addition, the proposed method is based on the hybridization of multi-objective particle swarm optimization (MOPSO) using the reference direction and local fast and accurate search capabilities of the SQP to provide a computationally more efficient and accurate algorithm for solving many-objective optimization problems. In the proposed method, a new mutation operator was introduced into the MOPSO to enhance the computational performance, uniformity of the solution and to prevent the MOPSO from pre-convergence to the local Pareto front. The proposed method was compared with other many-objective optimization algorithms using several challenging DTLZ benchmark problems and real-world many-objective optimization problems involving 3 to 15 objective functions. The results show that the proposed method has a high diversity and very accurate Pareto solutions. In this respect, the proposed method outperforms other recently developed many-objective optimization algorithms used in this study. Based on the obtained results, the proposed method provides an effective way of solving many-objective optimization problems.

Keywords Many-objective optimization · Multi-objective particle swarm optimization (MOPSO) · Sequential quadratic programming (SQP) · Decomposition · Reference point · Multi-objective evolutionary algorithms (MOEAs)

1 Introduction

Most real-world engineering design optimization problems (e.g., aerospace [1, 2], automotive [3], and energy systems [4]) involve several and often conflicting design criteria. For example, the design optimization of aircraft structures involves several conflicting objectives such as weight, buckling, and reliability, and hence optimization must be conducted simultaneously to achieve a balanced optimum design among the various conflicting objectives. In the past decades, a significant number of research works

have focused on the multi-objective evolutionary algorithms (MOEAs) [5–8], which have been used successfully for solving multi-objective optimization problems [9–11]. However, the increased dimensionality of objective function search space and high computational costs limit the application of MOEAs (e.g., NSGA-II and MOPSO) for solving many-objective optimization problems [12–14].

There are several research works focused on solving many-objective optimization problems using evolutionary algorithms, which can be divided into the following five different categories: (i) preference-based relations [15], such as the ϵ -preferred [16], and the expansion dominance [17], that require the use of problem-specific parameters. (ii) indicator-based [18–20], which assign fitness to solutions using a scalar performance metric (indicator) [21]. (iii) dimensionality reduction [22], which are based on transforming

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many-objective optimization problem into a reduced number of objective functions [23–25], and depend on the characteristics of the problem [26], and the selection of parameters to find sub-problems [27]. (iv) aggregation-based [28], such as weighted sum [29], and weighted Tchebycheff [30], which transform a many-objective optimization problem into a multi-objective optimization problem [31]. (v) decomposition-based algorithms [32–34], which divides the problem into a number of sub-problems [35]. The decomposition-based approach has successfully solved many-objective optimization problems [36, 37]. In general, the decomposition-based approach can be divided into; (a) aggregation functions and (ii) reference points [38]. Recently, MOEA/D-M2M [39] and the NSGA-III [40], based on the decomposition-based methods, utilize reference points [41].

In addition, several research works focused on using many-objective evolutionary algorithms, such as Pareto-based MOEAs, in this approach, the original Pareto-based dominance relation has been modified to enhance the convergence pressure for solving many-objective optimization problems, such as fuzzy-dominance [42], corner sorting [43], reference point-based dominance [44], and generalized Pareto-optimality [45]. Several research works focused on enhancing diversity management in Pareto-based MOEAs, such as the shift-based density estimation in SPEA2-SDE [46] and the use of associated reference points in NSGA-III [40]. In decomposition based MOEAs, two adaptive generation approaches were developed using weight vectors [47]. For example, the Pareto-based dominance was integrated with the decomposition-based approach in reference [48], using two external archives to ensure convergence and diversity. Despite the above mentioned advantages for solving many-objective optimization problems, these methods suffer from some drawbacks [49]. For instance, these algorithms do not guarantee convergence to the true Pareto-optimal front [50]. To overcome these issues, several research works focused on hybrid based approaches [51]. For example, Zhang et al. [52], proposed a hybrid algorithm by combining NSGA-III and knowledge-based optimization [53], into a many-objective optimization algorithm. In addition, angle-based decomposition [54], and encouragement-based boundary intersection decomposition [55], have been integrated to provide a balance between convergence and diversity [56].

As described above, many-objective optimization problem introduces new challenges to the existing optimization methods. This paper presents a novel decomposition-based hybrid many-objective optimization algorithm to address the challenges associated with the many-objective optimization problems. The main advantages of the proposed algorithm are as follows; (1) Integration of global search capabilities of MOPSO based on reference direction, and local fast and accurate search capabilities of gradient-based,

sequential quadratic programming (SQP), to provide a computationally efficient and accurate algorithm for solving many-objective optimization problems. (2) Implementation of decomposition-based techniques to ensure the uniformity of the solution. (3) Implementation of a novel mutation operator for preventing the PSO from pre-convergence to the local optima as well as enhancing the computational performance of the MOPSO. In addition, the MOPSO is based on the reference direction, in which the information between reference points is utilized to enhance the efficiency of the MOPSO. Based on the obtained results, the proposed algorithm outperforms all other well-known many-objective optimization algorithms used in this study in terms of both diversity and convergence.

2 Methodology

The proposed methodology is based on the decomposition-based approach. In a decomposition-based approach, a set of reference directions or reference points are generated, in the objective function space, to guarantee uniformity of the results, which form the Pareto front. In reference point-based approaches, an objective function depends on the distance from the reference point.

The proposed methodology also utilizes the normal-boundary intersection (NBI) [57, 58], method to generate uniformly distributed reference points, which defines a set of reference points, as represented by (W), covering a hyper-plane, as shown in Fig. 1.

In the proposed methodology, a composite objective function is used to define the objective function. In this approach, the distance from a search point in the objective function space to its corresponding reference point is to be minimized. Therefore, reference points must be located between the origin of the objective function space coordinate system and the true Pareto frontier.

2.1 Decomposition-based hybrid evolutionary and gradient-based many-objective optimization algorithm

The organization of the proposed decomposition-based hybrid evolutionary and gradient-based many-objective optimization algorithm is shown in Fig. 2 and described as follows:

In Fig. 2, d_t is the search direction and a_t is the step size in iteration t . The parameter W represents the number of reference directions and z is used as the number of iterations, for checking that all the reference directions are utilized (i.e., $W = z$) in the SQP.

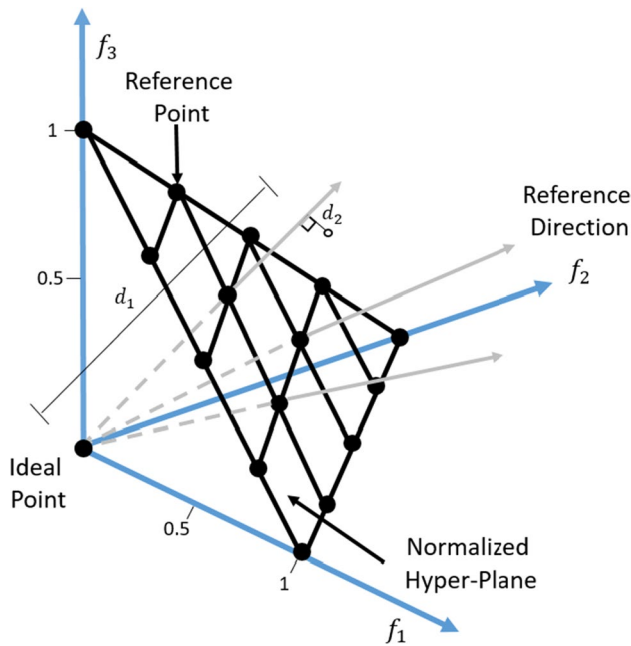


Fig. 1 A set of neighbouring particles and reference directions in a normalized hyper-plane for $M=3$ (M is the number of the objective functions). There are five reference points in the f_1 - f_2 plane

A general many-objective particle swarm optimization (MOPSO) can be stated as below [59]:

$$\begin{aligned} \min : f(x) &= [f_1(x), f_2(x), f_3(x), \dots, f_M(x)], x \in \Omega \\ \text{s.t.} \quad &g_q(x) \leq 0, \quad q = 1, 2, \dots, q \\ &h_s(x) = 0, \quad s = 1, 2, \dots, s \\ \text{side constraints : } &x_i^{lb} \leq x_i \leq x_i^{ub}, i = 1, 2, \dots, n \end{aligned} \quad (1)$$

where M is number of objective functions, Ω is the feasible search space, $x = (x_1, x_2, \dots, x_n)$, is the vector of n -dimensional decision variable. The mapping function, $f(x) : \Omega \rightarrow R^M$, defines M objective functions, which are bounded in the objective space R^M . The inequality and equality constraints, which are represented as $g_q : \Omega \rightarrow R^q$ and $h_s : \Omega \rightarrow R^s$, respectively. In addition, x_i^{lb} and x_i^{ub} , are the lower and upper bounds of i -th decision variables, respectively.

The proposed decomposition-based hybrid many-objective optimization algorithm for solving the above many-objective optimization problems utilizes several real reference points instead of a hypothetical point, as in the case of the Tchebycheff approach. In addition, the proposed methodology differs from the Tchebycheff approach as how the objective functions are defined. The main stages involved in the proposed methodology are shown in Fig. 2, as described below:

Stage 1: Decomposition of multi-objective optimization problems

In this stage, decomposition of multi-objective optimization problems into a single objective function, and initialization of the parameters of MOPSO (i.e., population size, the maximum number of generations, and stopping criteria), are defined. In the conventional MOPSO, the initial population size is provided as input to the MOPSO algorithm for the first reference point. In the proposed method, several evaluations are performed using the different versions of the MOPSO and Constriction Type 1. In order to update velocity and position of the i -th decision variable of the j -th particle [59], the following can be stated:

$$\begin{aligned} v_{ij}^{t+1} &= w(v_{ij}^t + \phi_1(Pbest_{ij}^t - x_{ij}^t) + \phi_2(Gbest_i^t - x_{ij}^t)) \\ x_{ij}^{t+1} &= v_{ij}^{t+1} \Delta t + x_{ij}^t \end{aligned} \quad (2)$$

where v_{ij} , and x_{ij} , are the current velocity and position of the i -th decision variable of the j -th particle. In the proposed method, $Pbest_{ij}^t$, is the i -th decision variable of the best position of the j -th particle in iteration t . $Gbest_i^t$ is the i -th decision variable of the best global position whole population ever achieved. ϕ_1 and ϕ_2 , are random number between $[0, \varphi/2]$. In this paper, φ is a constant and equals to 4.1. w is the inertia weight parameter, and the value of this parameter used in this paper is 0.7298 [60]. Δt is the time step for calculating the position of a particle from its velocity, which is conventionally equals to 1.

In the objective function space, reference points are based on the number of objective functions. The multi-objective optimization problem is decomposed into single-objective optimization problems by utilizing these reference points. Following this, the optimization process is performed for each reference point, as shown in Fig. 2. For all the reference points, the approximate range of the solution calculated using the MOPSO. A composite function defined as an objective function as described below in Stage 2.

Stage 2: Normal boundary intersection method for fitness assignment

In decomposition-based approaches, normal boundary intersection (NBI) [57], is a conventional way to obtain Pareto optimal. To obtain a point of Pareto front, the perpendicular distance to the reference line, d_2 , should be zero, as shown in Fig. 1. At the same time, the distance between the origin of the coordinate system of the objective function space and the projection of the reference direction to the f_1 and f_2 plane, and the reference direction, d_1 , to be minimized. In this respect, based on the preliminary studies and experimentations, d_2 is defined as a constraint. The definition of d_2 as a constraint is a problem dependent due to the different characteristics of the objective functions. The proposed algorithm utilizes a composite function as

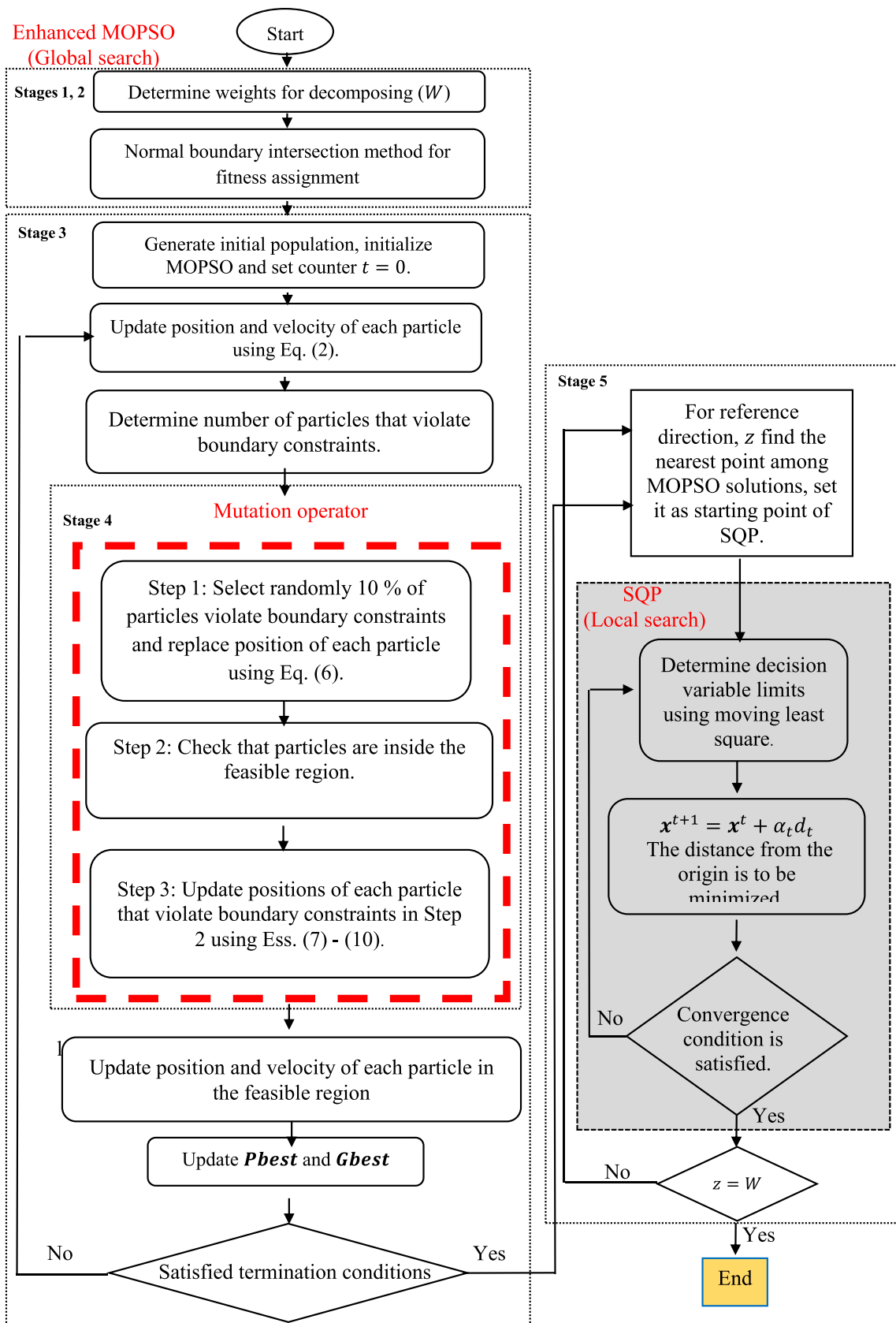


Fig. 2 The proposed decomposition-based hybrid many-objective optimization algorithm

the objective function for both MOPSO and SQP optimization process. In this respect, in the proposed algorithm, a recently developed composite function ($F(\mathbf{x})$) for many-objective optimization is utilized [61]:

$$F(\mathbf{x}) = F_1(\mathbf{x}) + F_2(\mathbf{x}) + F_3(\mathbf{x}) \quad (3)$$

where,

$$\begin{aligned} F_1(\mathbf{x}) &= \min \left(\frac{f_M(\mathbf{x}) - f_M^{\min}}{f_M^{\max} - f_M^{\min}} \right) \\ F_2(\mathbf{x}) &= \sum_{M=1}^M \left(\frac{f_M(\mathbf{x}) - f_M^{\min}}{f_M^{\max} - f_M^{\min}} \right) \\ F_3(\mathbf{x}) &= \left(\sum_{M=1}^M \left| \frac{f_M(\mathbf{x}) - f_M^0}{f_M^{\max} - f_M^0} \right|^p \right)^{1/p} \end{aligned} \quad (4)$$

where $f_M(\mathbf{x})$ is the value of the M -th objective function for \mathbf{x} , f_M^{\min} and f_M^{\max} are the minimum and maximum values of M -th objective function, f_M^0 is the M -th component of a reference point, and M represents number of the objective functions.

It should be mentioned that for the first iteration, the values obtained from the first function evaluation are used for determining f_M^{\max} and f_M^{\min} , as well as the $F_1(\mathbf{x})$, $F_2(\mathbf{x})$, and $F_3(\mathbf{x})$. In addition, as shown in Eq. (3), in the case of the $F_3(\mathbf{x})$, p is a parameter represented by L_p , which is a metric distance, and in this work, $p=2$ is used (Euclidian distance is considered).

The problem of the diversity (or the spread) of the solutions on the Pareto optimal frontier is associated with the different cases of distance functions, individually or in combination. The main problem is that the solutions can be sensitive to the choice of reference points. To overcome these problems, in this work, a new composite function (or fitness evaluation function) is utilized. This composite function involves the following three main features; (1) the best objective function value among the M objective functions, (2) the sum of all the objective functions, and (3) the metric distance function, which measures the distance between the objective functions, and their ideal values (aspiration levels, target values or the reference point). According to these definitions, $F_1(\mathbf{x})$ is the best normalized value through the objective functions, $F_2(\mathbf{x})$ is the proximity measure, the sum of the differences between the composite objective functions, and the best value obtained for the corresponding component of the composite function. In addition, $F_3(\mathbf{x})$ is defined as the summation of the normalized Euclidian distance of the obtained point and the intended reference point. It should be noted that in the proposed method, reference points must be located between the origin of the coordinate system of the objective function space and the real Pareto frontier. In this respect, prior to the start of the main stages of the proposed algorithm, the hyper-plane must be contracted to ensure

sufficient closeness of the reference points to the origin of the coordinate system of the objective function space.

Stage 3: Enhanced multi-objective particle swarm optimization (MOPSO):

As described above, the proposed hybrid decomposition-based many-objective optimization methodology integrates the global search capability of the MOPSO, with the local search capability of the SQP, as a hybrid many-objective optimization algorithm. In this respect, a new mutation operator (Stage 4) was introduced into the MOPSO to enhance the convergence of the proposed algorithm. In addition, the MOPSO is based on the reference direction, in which the information between reference points is utilized to enhance the efficiency of the MOPSO. The steps involved in the proposed method are described below:

- Step 1: If the reference point counter is greater than 1; do.
- Step 2: Among all the solutions, find the closest particle solution to the reference point (in the objective function space).
- Step 3: The final population values for the closest particle position are set as an initial population used in the MOPSO.

As stated above, in the first step of the proposed algorithm, global optimization is performed utilizing the enhanced MOPSO algorithm for each of the reference points based on the many-objective optimization formulation. The objective function using the penalty function stated as below:

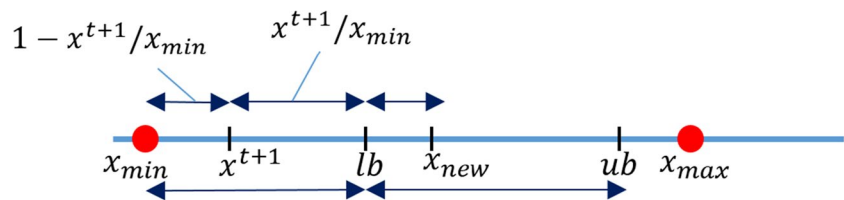
$$T(\mathbf{x}) = F(\mathbf{x}) + r_k \left\{ \sum_{i=1}^m [0, g_i(\mathbf{x})] + \sum_{j=1}^l [h_j(\mathbf{x})] \right\} \quad (5)$$

where $T(\mathbf{x})$ is the new objective function, r_k is the penalty coefficient and $\sum_{i=1}^m [0, g_i(\mathbf{x})] + \sum_{j=1}^l [h_j(\mathbf{x})]$, are the penalty terms that are added to the objective function, $F(\mathbf{x})$. Penalty terms are proportional to the constraint violation.

In order to avoid the effects of stochastic nature of the MOPSO algorithm on the result, a multi-start strategy is used [62, 63].

In the decomposition-based approaches, the ideal point is an important concept influencing the convergence of the algorithm. In several research works, the ideal point is used to normalize the objective function values, and subsequently to normalize the objective function space. In some algorithms such as MOEA/D and NSGA-III, this point is defined as; $\mathbf{z} = (f_1^{\min}, f_2^{\min}, \dots, f_i^{\min}, \dots, f_M^{\min})$, where in the optimization process, f_M^{\min} is the minimum of the M -th objective function, and M is the number of objective functions [57]. Based on this concept, the position of the ideal point is

Fig. 3 Replacing search points using the suggested mutation operator



improved during the search process, and the accuracy of the obtained ideal point gradually becomes close to the true ideal point. In this respect, the accuracy of the obtained ideal point is important for convergence.

In contrast, in some research works, the ideal point is defined as the origin of the coordinate system of the objective function space [58]. By considering the positive values for all the objective functions, the origin of the coordinate system of the objective function space is the absolute minimum for all the objective functions, because f_M^{\min} cannot be less than its corresponding value in the origin of the coordinate system of the objective function space, which is zero. Hence, based on this definition, the position of the ideal point is fixed.

Normalization effectively solves problems with objective functions in different orders of magnitude [61]. In some algorithms, reference directions are generated by connecting the ideal point to each of the reference points [58]. In the present work, the origin of the coordinate system of the objective function space is used as an ideal point. This makes the solution to be independent of the scales of objective functions space.

In the proposed method, the composite function, $F_1(\mathbf{x})$, $F_2(\mathbf{x})$, and $F_3(\mathbf{x})$, Eqs. (3) and (4) are used, and normalization is performed. Therefore, the accuracy of the obtained ideal point gradually becomes close to the true solution point. In addition, in some approaches [61], the ideal point is the base point of the decomposition of the Pareto front. In contrast, in the proposed method, the decomposition performed only at the start of the algorithm and does not depend on the ideal point. This satisfies the requirements of the uniformity of solutions of the Pareto front without the need for updating ideal point, whereas identification process of the ideal point requires additional computation costs [61]. In the proposed method, the search is performed for each weight independently.

Table 1 The variation of the coefficient δ , with respect to the reduced search space as the number of iterations increases

Iteration	Δ
Number of iterations is less than the maximum number of iterations divided by 3 ($\frac{iter_{max}}{3}$)	0.5
Number of iterations is greater than $\frac{iter_{max}}{3}$ and is less than $\frac{2 \times iter_{max}}{3}$	0.3
Number of iterations is greater than $\frac{2 \times iter_{max}}{3}$	0.1

In Stage 2 of the proposed method, a parallelization scheme adopted, and the calculations for all the reference directions are conducted simultaneously in each iteration. The projection of the objective function vector on the closest reference direction used as the objective function.

Stage 4: Enhanced mutation operator for MOPSO algorithm

MOPSO algorithm often converges to local optima and suffers from premature convergence [64]. To overcome this challenge, the mutation operator, which is a genetic algorithm operator, was introduced into the MOPSO algorithm. In this respect, several experiments were conducted on the main characteristics of many-objective optimization problems. Based on the experimentation results, an enhanced variant of the mutation operator was introduced into the MOPSO algorithm. The main purpose of this enhanced mutation operator is to prevent the algorithm from premature convergence to the local optima and to improve the global search capability of the MOPSO algorithm. The proposed enhanced mutation operator affects the constraints handling processes. The steps involved in the proposed enhanced mutation operator are described as follows:

Step 1: Select randomly 10% of particles that violate the boundary constraints and replace their positions in random, using the range of feasible distance as stated below:

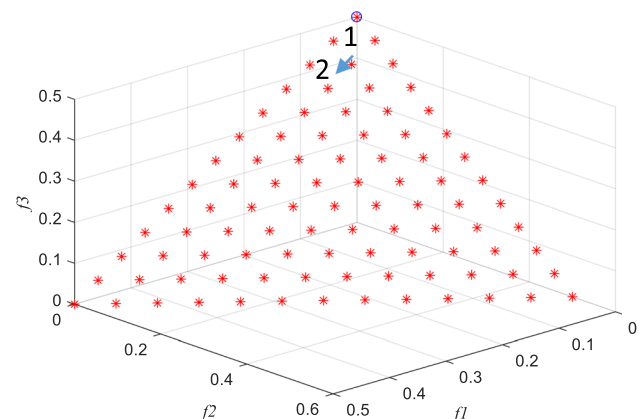


Fig. 4 Neighboring point concept (using first reference point solution as an initial point of reference point 2 in SQP search)

Table 2 DTLZ benchmark problems [65] (M = Number of objective functions)

Test problem	The term used as power (α)	Multiplying parameter (K)	M	Properties	Objective functions (that should be minimized)
DTLZ1	-	5	5, 8, 12, 16	Multimodal	$f_1(\mathbf{x}) = 0.5(1 + g) \prod_{i=1}^{M-1} x_i$ $f_{m=2:M-1}(\mathbf{x}) = 0.5(1 + g) \left(\prod_{i=1}^{M-m} x_i \right) (1 - x_{M-m+1})$ $f_M(\mathbf{x}) = 0.5(1 + g)(1 - x_1)$ <p>where $g = 100 \left(K + \sum_{i=M}^{M+K-1} (x_i - 0.5)^2 - \cos(20\pi(x_i - 0.5)) \right)$</p> <p>$\mathbf{x} = (x_1, x_2, \dots, x_{M+K-1})^T \in [0, 1]^{M+K-1}$</p>
DTLZ2	-	10	5, 8, 12, 16	Unimodal, concave	$f_1(\mathbf{x}) = (1 + g) \prod_{i=1}^{M-1} \cos(x_i \pi / 2)$ $f_{m=2:M-1}(\mathbf{x}) = (1 + g) \left(\prod_{i=1}^{M-m} \cos(x_i \pi / 2) \right) \sin(x_{M-m+1} \pi / 2)$ $f_M(\mathbf{x}) = (1 + g) \sin(x_1 \pi / 2)$ <p>where $g = \sum_{i=1}^{M+K-1} (x_i - 0.5)^2$, $\mathbf{x} = (x_1, x_2, \dots, x_{M+K-1})^T \in [0, 1]^{M+K-1}$</p>
DTLZ3	-	10	5, 8, 12, 16	Multimodal, concave	$f_1(\mathbf{x}) = (1 + g) \prod_{i=1}^{M-1} \cos(x_i \pi / 2)$ $f_{m=2:M-1}(\mathbf{x}) = (1 + g) \left(\prod_{i=1}^{M-m} \cos(x_i \pi / 2) \right) \sin(x_{M-m+1} \pi / 2)$ $f_M(\mathbf{x}) = (1 + g) \sin(x_1 \pi / 2)$ <p>where $g = 100 \left(K + \sum_{i=M}^{M+K-1} (x_i - 0.5)^2 - \cos(20\pi(x_i - 0.5)) \right)$</p> <p>$\mathbf{x} = (x_1, x_2, \dots, x_{M+K-1})^T \in [0, 1]^{M+K-1}$</p>
DTLZ4	100	10	5, 8, 12, 16	Concave	$f_1(\mathbf{x}) = (1 + g) \prod_{i=1}^{M-1} \cos(x_i \pi / 2)$ $f_{m=2:M-1}(\mathbf{x}) = (1 + g) \left(\prod_{i=1}^{M-m} \cos(x_i \pi / 2) \right) \sin(x_{M-m+1} \pi / 2)$ $f_M(\mathbf{x}) = (1 + g) \sin(x_1 \pi / 2)$ <p>where $g = 100 \left(K + \sum_{i=M}^{M+K-1} (x_i - 0.5)^2 - \cos(20\pi(x_i - 0.5)) \right)$</p> <p>$\mathbf{x} = (x_1, x_2, \dots, x_{M+K-1})^T \in [0, 1]^{M+K-1}$</p>
C1-DTLZ1	-	5	5, 8, 10, 15	Multimodal	<p>This is a C1-DTLZ1 test problem involving the following constraint:</p> $c(\mathbf{x}) = 1 - \frac{f_{M+1}(\mathbf{x})}{0.6} - \sum_{i=1}^{M-1} \frac{f_i(\mathbf{x})}{0.5} \geq 0$

Table 2 (continued)

Test problem	The term used as power (α)	Multiplying parameter (K)	M	Properties	Objective functions (that should be minimized)
C2-DTLZ2	1	10	5, 8, 10, 15	Unimodal, concave	This is a DTLZ2 test problem involving the following constraint: $c(\mathbf{x}) = \max \left\{ \max_{i=1, \dots, M} \left[(f_i(\mathbf{x}) - 1)^2 + \sum_{j=1, j \neq i}^M f_j^2 - r^2 \right], \left[\sum_{i=1}^M (f_i(\mathbf{x}) - 1/\sqrt{M})^2 - r^2 \right] \right\} \geq 0$
C3-DTLZ4	100	10	5, 8, 10, 15	Concave	The objective functions are the same as those were in the DTLZ2, with the following constraint: $c_j(\mathbf{x}) = \frac{f_j^2}{4} + \sum_{i=1, i \neq j}^M f_i^2 - 1^2 \geq 0 \forall j = 1, 2, \dots, M$

$$x_{ij} = r_{ij} \times (x_{ij}^{ub} - x_{ij}^{lb}) + x_{ij}^{lb}; \quad (6)$$

where r_{ij} is a random real number in the range of $[0, 1]$, x_{ij} is the current generation position of the i -th decision variable of the j -th particle, and x_{ij}^{ub} and x_{ij}^{lb} are the upper and lower bound of the i -th decision variable position of the j -th particle, respectively.

Step 2: Check that particles are inside the feasible region, which do not violate the boundary constraints.

Step 3: Calculate a ratio (r_{ij}) and multiply it by a random number to produce a new particle position. A schematic view of this concept shown in Fig. 3. In this figure, x_{ij}^{t+1} presents particles that violate the boundary constraints, $1 - x_{ij}^{t+1}/x_{ij(min)}$ and $x_{ij}^{t+1}/x_{ij(max)}$, which is equal to the coefficient r_{ij} , which guides the mutated particles to be located inside the feasible region that do not violate boundary constraints.

Indeed, suppose the position of a particle in iteration $t + 1$, which violates a boundary constraint. In this case, the position of the particle updated to a feasible point $x_{ij(new)}$, as described below.

To check, whether the position of the i -th decision variable of the j -th particle in iteration $t + 1$ violates the lower boundary constraint, the following statement can be used:

$$\text{If } x_{ij}^{t+1} < x_{ij}^{lb} \quad (7)$$

where, x_{ij} is the current generation position of the i -th decision variable of the j -th particle and t is the number of the iteration. x_{ij}^{ub} and x_{ij}^{lb} are the upper and lower bounds of the i -th decision variable position of the j -th particle, respectively.

If the lower boundary constraints violated, then the position of the particle can be updated using the following Eq. (8)

$$r_{ij} = 1 - x_{ij}^{t+1}/x_{ij(min)} \text{ and hence } x_{ij(new)} = r_{ij} \times (x_{ij}^{ub} - x_{ij}^{lb}) + x_{ij}^{lb} \quad (8)$$

Similarly, it can be examined whether the position of the i -th decision variable of the j -th particle in iteration $t + 1$ violates the upper boundary constraint using the following statement:

$$\text{If } x_{ij}^{t+1} > x_{ij}^{ub} \text{ then} \quad (9)$$

If the lower boundary constraints violated, then the position of the particle can be updated using the following Eq. (10) as below:

$$r_{ij} = 1 - x_{ij}^{t+1}/x_{ij(max)} \text{ and } x_{ij(new)} = 1 - r_{ij} \times (x_{ij}^{ub} - x_{ij}^{lb}) + x_{ij}^{lb} \quad (10)$$

where $x_{i,j(new)}$ is the new position of the i -th decision variable of the j -th particle, t is the number of iteration, $x_{i,j}^{lb}$ and $x_{i,j}^{ub}$ are the lower and upper bounds of i -th decision variable of the j -th particle, respectively. $x_{i,j(min)}$ and $x_{i,j(max)}$ are the minimum and maximum values of the corresponding variables, respectively.

Stage 5: Gradient-based optimization using SQP

In this stage, the results of MOPSO are transferred as input to the SQP. In this respect, as the maximum generation number in MOPSO is reached for the final reference point, the values of obtained decision variables are transferred to the SQP to attain an accurate solution. The optimization using SQP completes when the imposed stopping criteria on SQP are satisfied. The stopping criteria used in this study are satisfaction of the constraints and the maximum number of iterations (i.e., the difference between two consecutive iterations must be less than $1e-6$), which depends on the problem and required degree of accuracy, and this is defined by the user.

2.2 Sequential Quadratic Programming (SQP)

As described above, the proposed enhanced MOPSO algorithm is used to obtain global optimal solutions (corresponding to reference points), which are used to find a search direction using a linear approximation of constraints and quadratic approximation of objective functions in the SQP algorithm. The solution procedure is based on solving a quadratic programming sub-problem in each direction. Formulation of the SQP problem in iteration t can be stated as:

$$\begin{aligned} \min : & \quad \frac{1}{2}d^T H_t d + \nabla F(\mathbf{x})^T d \\ \text{constraints : } & \quad \nabla h_a(\mathbf{x})^T d + h_a(\mathbf{x}) = 0 \quad a = 1, \dots, a \\ & \quad \nabla g_b(\mathbf{x})^T d + g_b(\mathbf{x}) < 0 \quad b = 1, \dots, b \\ & \quad x^{lb} \leq \mathbf{x} \leq x^{ub} \end{aligned} \quad (11)$$

Search space : $[\mathbf{x}_0 - \delta \times \mathbf{x}_0, \mathbf{x}_0 + \delta \times \mathbf{x}_0]$

where d_t is the search direction, t and H_t is a positive definite approximation of the Hessian matrix of the Lagrangian function of the sub-problem. \mathbf{x} is the vector of n decision variables ($i = 1, 2, \dots, n$), (the optimal solution from the MOPSO that is considered as the initial search point (\mathbf{x}_0) to the SQP), x^{lb} and x^{ub} are the lower and upper bounds of the decision variables. $g_b(\mathbf{x})$ and $h_a(\mathbf{x})$ are the inequality and equality constraints, respectively. δ is a coefficient that defines the range of search space for the SQP and is set according to Table 1.

The composite function $F(\mathbf{x})$ used to transfer the many-objective optimization problem into a single objective optimization problem in each direction in the optimization process. This makes it possible to utilize the SQP algorithm (as a single

Table 3 Parameter settings of the population sizes and the number of function evaluations

Benchmark problems	Number of objective functions (M)	Population size (P)
DTLZ1-DTLZ4	5	210
	8	330
	12	364
	16	816
C1-DTLZ1	5	210
C2-DTLZ2	8	156
C3-DTLZ4	10	275
	15	135

objective optimization method). The Lagrangian function can be expressed as [62]:

$$L(\mathbf{x}, u, v) = F(\mathbf{x}) + \sum_{a=1}^a u_a h_a(\mathbf{x}) + \sum_{b=1}^b v_b g_b(\mathbf{x}) \quad (12)$$

where u_a and v_b are the Lagrangian multipliers.

The active set strategy used to solve the sub-problem as stated below:

$$\mathbf{x}^{t+1} = \mathbf{x}^t + a_t d_t \quad (13)$$

where a_t is the step size parameter in iteration t , which is determined using a line search method so that a sufficient decrease in an objective function is obtained [63].

It should be mentioned that the optimization using the SQP was conducted in MATLAB software. In this respect,

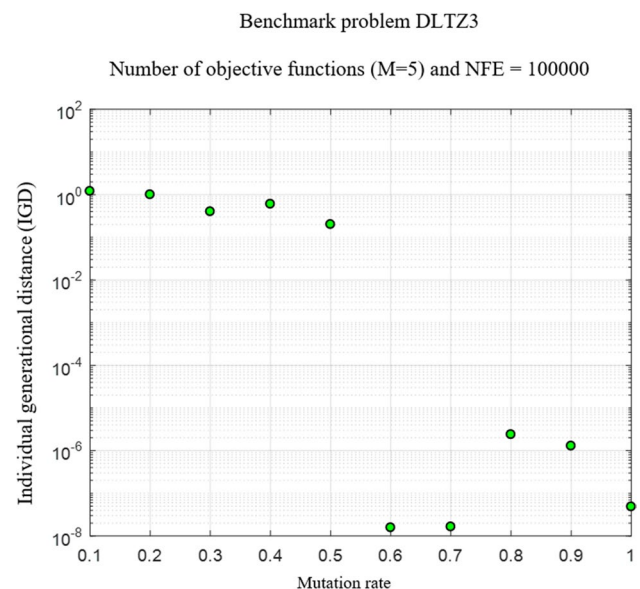


Fig. 5 Individual generational distance (IGD) values for different mutation rates

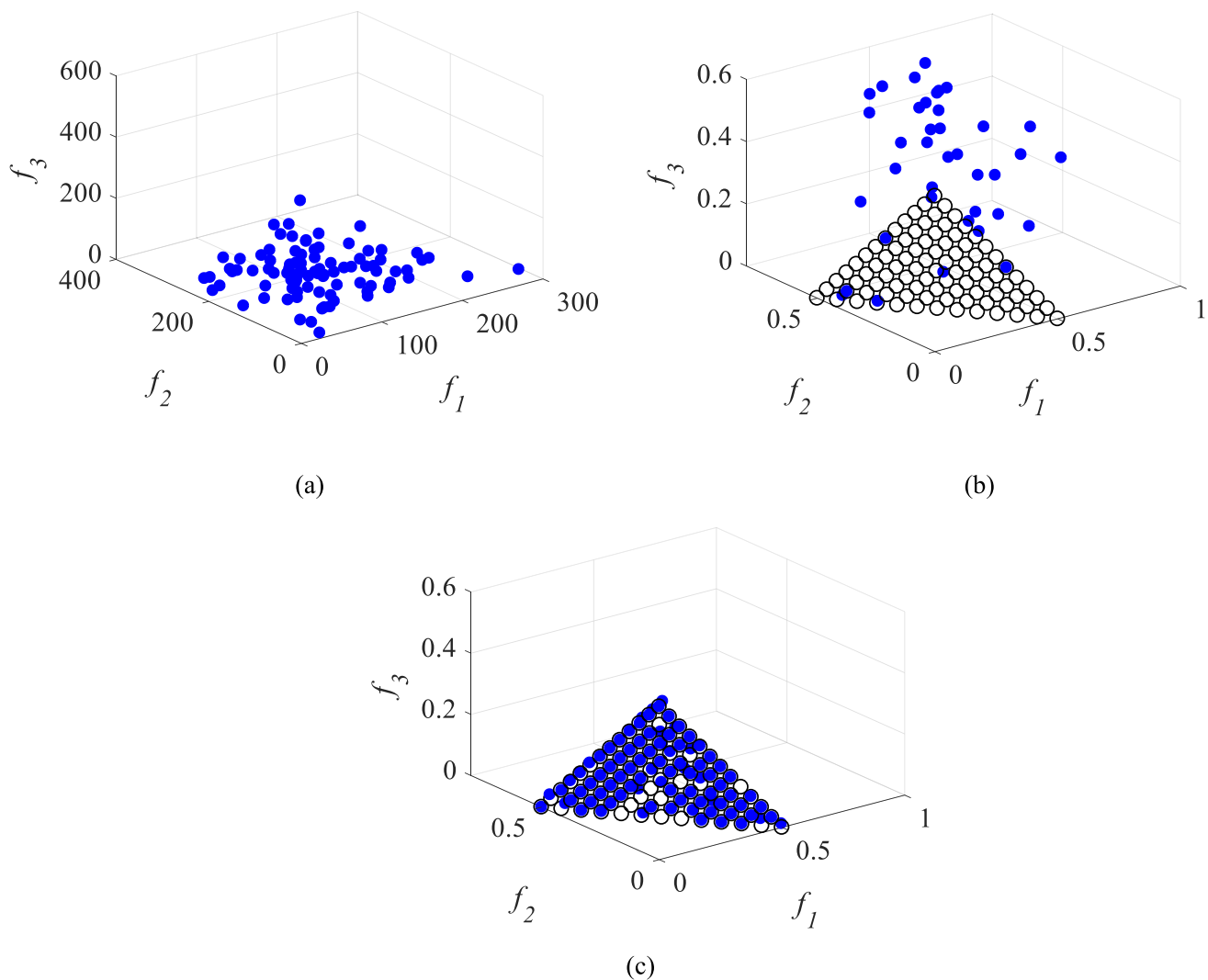


Fig. 6 DTLZ1 (3 objective functions): **a** Initial population, **b** MOPSO result (without mutation operator), **c** MOPSO result (with mutation operator)

both the gradients and the Hessians approximated numerically for the reference points.

In the proposed algorithm, the moving square bounds method utilized to decrease the range of the decision variables. Towards this, the search space for the SQP is based on the output of the MOPSO using coefficient δ , which decreases through the iterations of the SQP algorithm as defined in Table 1:

In Table 1, $iter_{max}$ is the maximum number of iterations, which is used in the SQP algorithm.

2.3 Neighboring point technique

In some problems, specifically those that involve smooth functions (e.g., problems with unimodal functions), only one initial point of the Pareto front can be sufficient for the local search. In this respect, in the first step, a global

search is performed using the proposed enhanced MOPSO algorithm, and an initial solution for the first reference point is obtained (point 1 in Fig. 4). In the second step, the result of the MOPSO algorithm used as an initial point for the SQP algorithm, which provides a more accurate solution for the reference point, which obtained in the first step. Following this, for reference point 2, a global search was not utilized, and only the local exploitation search using the SQP algorithm was conducted, in which the reference point 1 solution was considered as an initial point in the SQP algorithm. In this respect, for the next reference points, the closest reference point from the Pareto solutions is used as an initial point for the local search using the SQP algorithm without the need for a global search. In the optimization process, when the SQP algorithm converged, the solution added to the non-dominated solution set.

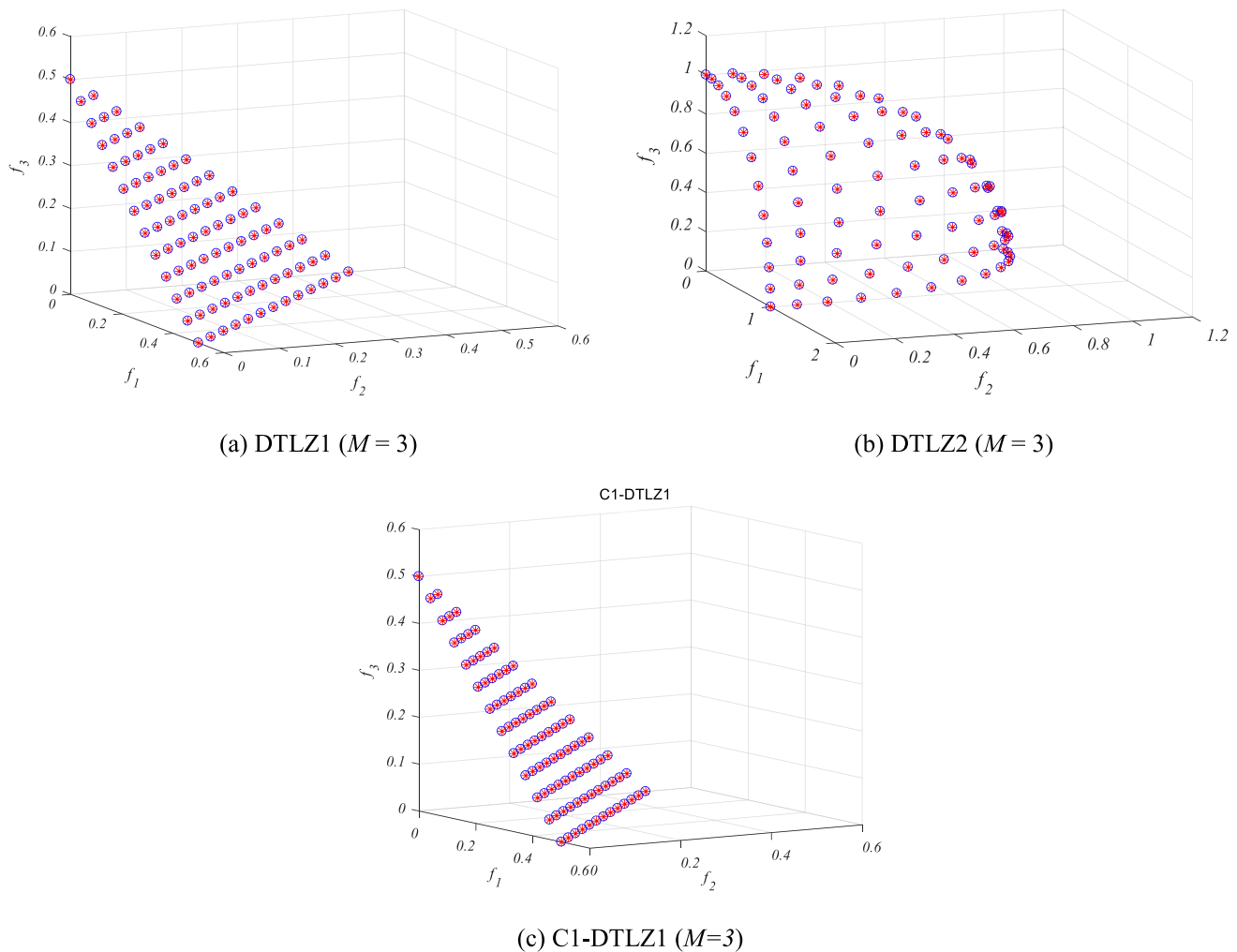


Fig. 7 Comparison between the proposed Pareto front and exact solution for **a** DTLZ1, **b** DTLZ2, **c** constrained DTLZ (C1-DTLZ1) with 3 objective functions

3 Experiments with benchmark problems

3.1 Experimental settings

Several experiments were conducted on well-defined optimization benchmark problems such as DTLZ1, DTLZ2, DTLZ3, and DTLZ4, with 3, 5, 8, 12, and 16 objective functions. In addition, C1-DTLZ1, C2-DTLZ2, and C3-DTLZ4, with 5, 8, 10, and 15 objective functions, to demonstrate the capability and computational efficiency of the proposed algorithm. These benchmark problems are mainly used to assess the computational accuracy and efficiency of many-objective optimization algorithms [38], as shown in Table 2.

In this comparative study of the performance of the proposed and other algorithms (e.g., RPD-NSGA-II, MOEA/DD, MOEA/D-DE, and NSGA-III, etc.), parameter setting was adopted as shown in Table 3. This involves, the population size, and the maximum number of function evaluations

for each number of objective functions M . In this experimentation, for the unconstrained benchmark problems (DTLZ1, DTLZ2, DTLZ3, and DTLZ4), the same population sizes as in [33], were used, and for constrained benchmark problems (C1-DTLZ1), the parameter setting is similar to [38]. In comparisons, the median of the performance criteria was obtained using 10 independent iterations.

To determine the number of function evaluations (NFE) of the proposed method (i.e., MOPSO and SQP). The total NFE of MOPSO and SQP are used. In general, for a smooth function, a gradient-based algorithm requires less iteration than an evolutionary algorithm to obtain the optimum point. In this respect and based on the results of several experiments using different benchmark problems, it has been concluded that for a fixed number of function evaluations, if 90% of the number of function calls are assigned to the evolutionary algorithm, the proposed algorithm reaches the most accurate solutions.

Fig. 8 Comparison of results of the proposed algorithm and other algorithms for the DTLZ1 benchmark problem

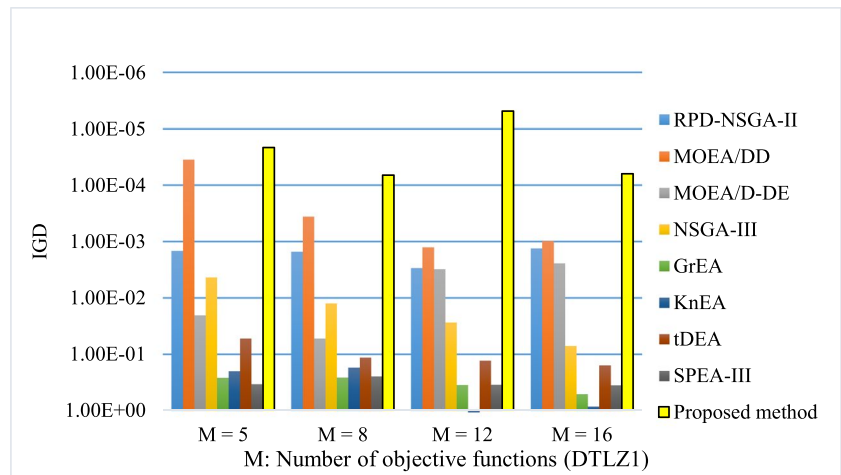


Table 4 The effect of hybridization of the MOPSO and SQP algorithms on the IGD. (M = Number of objective functions)

Benchmark problem	M	MOPSO		Proposed method		I-DBEA [57]	
		IGD	Simulation time (second)	IGD	Simulation time (second)	IGD	Simulation time (second)
DTLZ 1	3	1.27×10^{-1}	47.4	9.13×10^{-7}	162.3	1.75×10^{-3}	153.4
DTLZ 1	5	6.32×10^{-1}	30.6	2.14×10^{-5}	106.8	9.43×10^{-4}	156.5
DTLZ 2	3	5.47×10^{-2}	58.6	9.58×10^{-14}	53.8	6.37×10^{-4}	171.9
DTLZ 2	5	1.58×10^{-1}	52.7	8.70×10^{-14}	48.4	1.12×10^{-3}	231.2

Table 5 The effect of the proposed mutation operator in the MOPSO algorithm on the spacing and IGD performance metrics using the C1-DTLZ1 benchmark problem. (NFE = Number of function evaluation)

NFE	5,000			10,000			20,000		
	Criteria		Simulation time (second)	Criteria		Simulation time (second)	Criteria		Simulation time (second)
	Spacing	IGD		Spacing	IGD		Spacing	IGD	
GrEA	NaN	NaN	NaN	NaN	NaN	NaN	2.0759e-2	6.8772e-2	16.5
IBEA	NaN	NaN	NaN	NaN	NaN	NaN	2.2003e-1	1.7383e-2	17.9
KnEA	NaN	NaN	NaN	NaN	NaN	NaN	2.6951e-2	3.2562e-2	16.0
MOEAD	NaN	NaN	NaN	NaN	NaN	NaN	NaN	NaN	NaN
NSGAII	NaN	NaN	NaN	NaN	NaN	NaN	2.0115e-2	3.0906e-2	16.1
tDEA	NaN	NaN	NaN	NaN	NaN	NaN	NaN	6.0589e-2	16.3
EMOPSO	3.167e-2	-	51	2.1231e-2	-	103	1.9305e-2	-	15.0
MOPSO + the proposed mutation operator	3.989e-1	1.3594e+0	56	1.9400e-2	2.0600e-1	142	1.0100e-2	8.7000e-2	15.3

Values in boldface show the best performance of the algorithms as shown in the tables

In the process of the parameters setting of the proposed algorithm, consideration was given to establishing the best value of the mutation rate in the proposed enhanced MOPSO algorithm. In this respect, an experiment was carried out using the proposed MOPSO algorithm with ten different mutation rate values (i.e., between 0.1 and 1). The results of this experimentation are shown in Fig. 5. As can be seen from Fig. 5, for the mutation rate of 0.6 and 0.7, the accuracy

of results is better than other values of the mutation rate. Therefore, the parameter of the mutation rate was set to 0.6 for solving the benchmark problems.

Experiments with a benchmark problem with and without adopting the proposed mutation operator in the proposed MOPSO are performed to demonstrate the capability of the proposed mutation operator. The effect of the proposed mutation operator on the convergence of solutions to the

Table 6 Neighboring points affect the number of function evaluations (NFE) for DTLZ1. (M=Number of objective functions)

Problem	The proposed algorithm without the neighbouring point technique		The proposed algorithm with the neighbouring point technique	
	IGD	NFE	IGD	NFE
DTLZ1 (M=3)	8.14×10^{-5}	1,000,000	9.13×10^{-7}	100,000

Pareto optimal front is shown in Fig. 6, using the benchmark problem, DTLZ-1, with 3 objective functions and 91 reference points.

The results of this experimentation and association of the solutions corresponding to each reference point, with

10,000 fitness evaluations, are shown in Fig. 7. In this respect, Fig. 5(a) and (b), show that the solutions without the mutation operator limited particles are generated in the feasible region, which are not close to the Pareto front, whilst Fig. 7(c), with the mutation operator show that a large number of particles are generated in the feasible region. This experimentation, as described above, clearly illustrates the search capability of the proposed mutation operator in the EMOPSO algorithm. The proposed mutation operator provides fast convergence of the particles to the feasible region and hence generation of the Pareto front.

3.2 Performance metrics

In general, there are different performance metrics for evaluating multi-objective optimization algorithms, for

Table 7 Median IGD values for the DTLZ4 benchmark problem (bold items show the best performance). (M=Number of objective functions). (NFE=Number of function evaluations)

Problem	NFE	50,000		75,000		100,000	
		IGD	Simulation time (second)	IGD	Simulation time (second)	IGD	Simulation time (second)
DTLZ4 (M=12)	RPD-NSGA-II	3.39×10^{-3}	32.3	3.44×10^{-3}	47.8	3.13×10^{-3}	63.5
	MOEA/DD	5.90×10^{-4}	353.6	2.81×10^{-4}	527.7	2.60×10^{-4}	705.4
	MOEA/D-DE	6.26×10^{-3}	47.6	6.69×10^{-3}	71.2	6.79×10^{-3}	94.2
	NSGA-III	7.77×10^{-3}	17.9	7.07×10^{-3}	26.3	7.56×10^{-3}	41.4
	GrEA	4.5379×10^{-1}	18.8	4.5400×10^{-1}	28.7	4.5413×10^{-1}	36.3
	KnEA	4.9407×10^{-1}	19.4	4.9607×10^{-1}	28.5	4.8998×10^{-1}	39.9
	tDEA	5.00×10^{-1}	18.2	4.98×10^{-1}	27.6	4.98×10^{-1}	37.1
	SPEA-III	9.58×10^{-1}	17.8	6.75×10^{-1}	25.8	1.63×10^{-2}	35.3
	MOSA	1.87×10^0	18.5	7.643×10^{-2}	27.7	4.56×10^{-3}	37.1
	Proposed method	2.79×10^{-2}	17.5	2.03×10^{-10}	26.6	4.23×10^{-11}	46.3

Values in boldface show the best performance of the algorithms as shown in the tables

Table 8 Median IGD values with NFE=100,000 for the DTLZ1 benchmark problem (bold items show the best performance). (M=Number of objective functions)

Problem	DTLZ1 (M=5)		DTLZ1 (M=8)		DTLZ1 (M=12)		DTLZ1 (M=16)	
	IGD	Simulation time (second)	IGD	Simulation time (second)	IGD	Simulation time (second)	IGD	Simulation time (second)
RPD-NSGA-II	$1.46 \text{ e } -3$	104.4	$1.52 \text{ e } -3$	145.1	$2.95 \text{ e } -3$	200.0	$1.33 \text{ e } -3$	290.9
MOEA/DD	$3.54 \text{ e } -5$	102.2	$3.60 \text{ e } -4$	144.2	$1.27 \text{ e } -3$	194.3	$9.73 \text{ e } -4$	284.3
MOEA/D-DE	$2.03 \text{ e } -2$	107.1	$5.27 \text{ e } -2$	148.5	$3.10 \text{ e } -3$	204.5	$2.43 \text{ e } -3$	301.3
NSGA-III	$4.33 \text{ e } -3$	111.0	$1.26 \text{ e } -2$	156.9	$2.74 \text{ e } -2$	209.5	$7.18 \text{ e } -2$	309.1
GrEA	$2.5874 \text{ e } -1$	112.2	$2.5588 \text{ e } -1$	158.2	$3.4611 \text{ e } -1$	214.2	$5.0454 \text{ e } -1$	315.9
KnEA	$1.9834 \text{ e } -1$	116.1	$1.7076 \text{ e } -1$	162.2	$1.2822 \text{ e } 0$	221.2	$8.3971 \text{ e } -1$	325.0
tDEA	$5.2796 \text{ e } -2$	110.7	$1.1517 \text{ e } -1$	155.8	$1.3037 \text{ e } -1$	211.6	$1.5574 \text{ e } -1$	311.5
SPEA-III	<u>$3.36 \text{ e } -1$</u>	117.0	<u>$2.44 \text{ e } -1$</u>	163.1	<u>$3.46 \text{ e } -1$</u>	224.3	<u>$3.51 \text{ e } -1$</u>	329.3
MOSA	<u>$5.06 \text{ e } -3$</u>	117.9						
Proposed method	$2.14 \text{ e } -5$	106.8	$6.67 \text{ e } -5$	150.7	$4.81 \text{ e } -6$	204.1	$6.24 \text{ e } -5$	299.5

Values in boldface show the best performance of the algorithms as shown in the tables

Table 9 Median IGD values for the DTLZ2 benchmark problem (bold items show the best performance). (M=Number of objective functions)

Problem	DTLZ2 (M=5)		DTLZ2 (M=8)		DTLZ2 (M=12)		DTLZ2 (M=16)	
	IGD	Time (second)	IGD	Time (second)	IGD	Time (second)	IGD	Time (second)
RPD-NSGA-II	1.48 e -3	56.3	2.98 e -3	70.4	4.40 e -3	84.3	1.78 e -3	133.2
MOEA/DD	2.03 e -5	483.4	5.94 e -5	595.1	2.14 e -4	711.6	2.38 e -3	2317.1
MOEA/D-DE	1.61 e -3	87.3	4.23 e -2	88.9	4.43 e -3	76.3	2.44 e -3	108.9
NSGA-III	2.01 e -3	39.7	4.32 e -3	42.8	7.34 e -3	58.7	9.20 e -3	60.2
Proposed method	8.7 e -14	48.4	7.07 e -14	50.6	1.31 e -13	52.6	1.79 e -11	57.4

Values in boldface show the best performance of the algorithms as shown in the tables

Table 10 Median IGD values for the DTLZ3 benchmark problem (bold items show the best performance). (M=Number of objective functions)

Problem	DTLZ3 (M=5)		DTLZ3 (M=8)		DTLZ3 (M=12)		DTLZ3 (M=16)	
	IGD	Time (second)	IGD	Time (second)	IGD	Time (second)	IGD	Time (second)
RPD-NSGA-II	1.39 e -3	52.5	3.30 e -3	47.5	1.83 e -3	60.0	2.67 e -3	105.9
MOEA/DD	3.13 e -4	295.8	7.02 e -4	488.3	2.48 e -3	681.9	1.62 e -3	1752.6
MOEA/D-DE	1.17 e -2	59.9	4.05 e -2	63.3	1.92 e -3	67.1	2.80 e -3	85.4
NSGA-III	2.10 e -3	34.8	5.14 e -3	36.8	1.00 e -2	55.7	2.16 e -2	61.7
Proposed method	9.54 e -7	35.0	5.92 e -7	33.1	8.98 e -6	52.6	1.31 e -7	53.4

Values in boldface show the best performance of the algorithms as shown in the tables

Table 11 Median IGD values for the DTLZ4 benchmark problem (bold items show the best performance). (M=Number of objective functions)

Problem	DTLZ4 (M=5)		DTLZ4 (M=8)		DTLZ4 (M=12)		DTLZ4 (M=16)	
	IGD	Time (second)	IGD	Time (second)	IGD	Time (second)	IGD	Time (second)
RPD-NSGA-II	1.01 e -3	46.3	3.47 e -3	55.5	3.66 e -3	63.5	2.23 e -3	81.3
MOEA/DD	3.81 e -5	419.8	1.29 e -4	617.4	2.08 e -4	705.4	7.81 e -4	2483
MOEA/D-DE	4.17 e -3	89.1	7.31 e -2	92.6	6.98 e -3	94.2	3.66 e -3	116.0
NSGA-III	2.01 e -3	39.6	4.32 e -3	40.5	7.07 e -3	41.4	8.72 e -3	63.1
Proposed method	3.56 e -14	32.3	2.69 e -11	50.3	4.23 e -11	46.3	8.34 e -10	51.0

Values in boldface show the best performance of the algorithms as shown in the tables

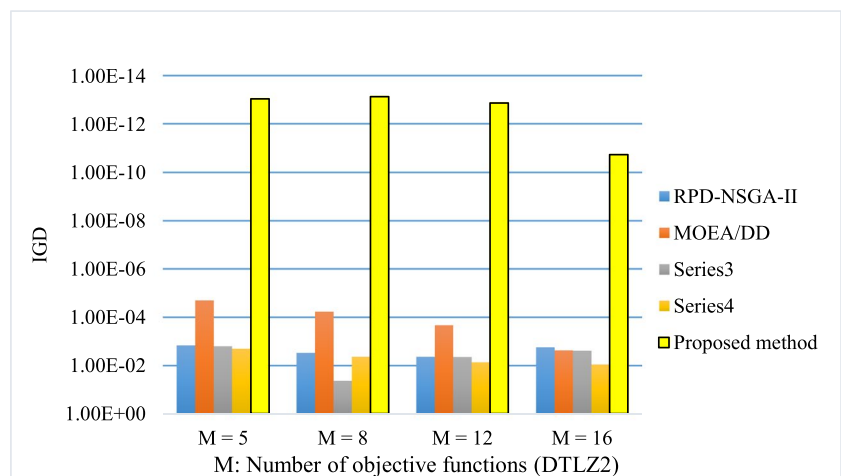
Fig. 9 Comparison of results of the proposed algorithm with other algorithms for DTLZ2 benchmark problem

Fig. 10 Comparison of results of the proposed algorithm and other algorithms for the DTLZ3 benchmark problem

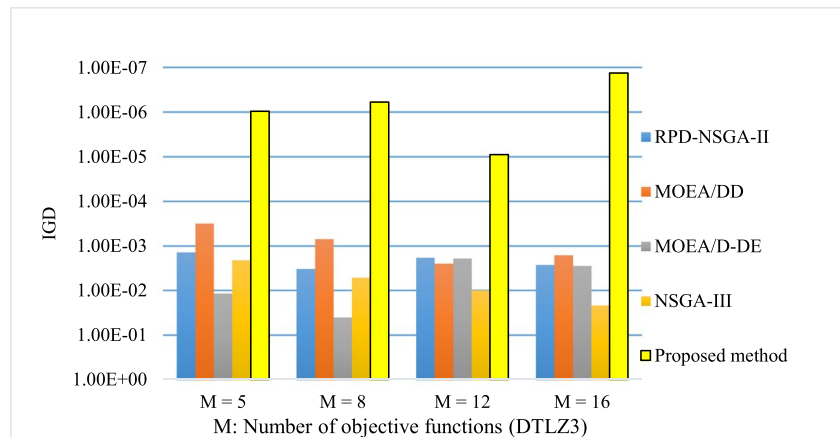


Fig. 11 Comparison of the results of the proposed algorithm and other algorithms for DTLZ4 benchmark problem

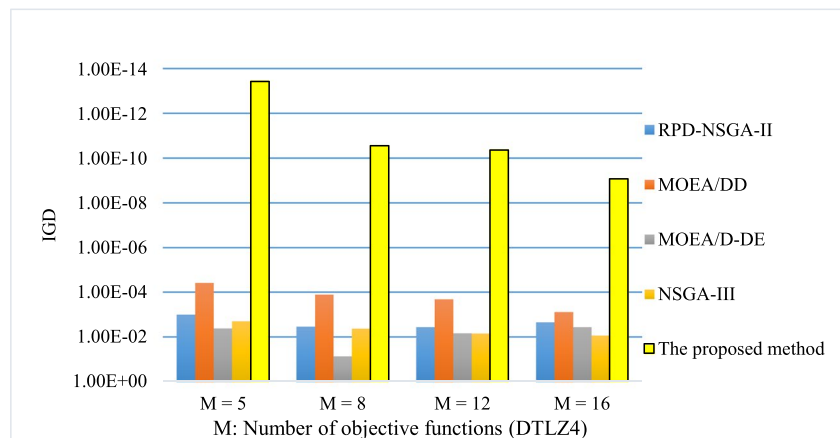
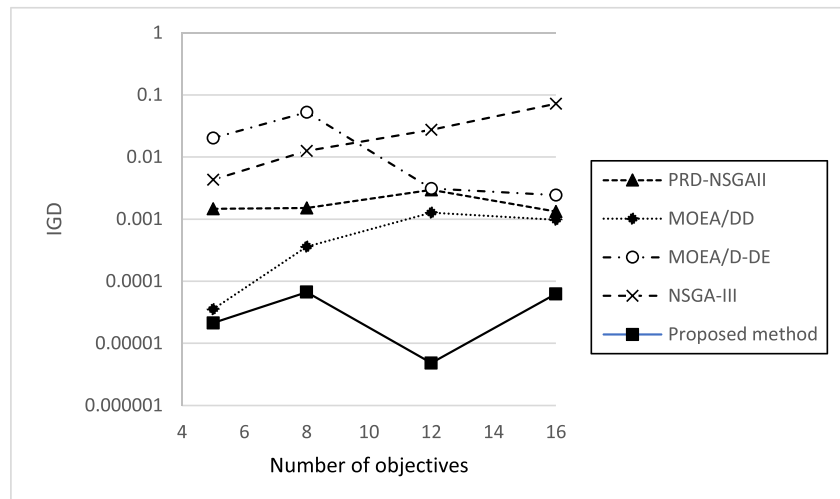


Fig. 12 Mean values for the IGD metric of DTLZ1 problem



example, spacing metric [66], pure diversity (PD) [67], generational distance (GD) [66], coverage of two sets [68] individual generational distance (IGD) [69], hyper volume (HV) [70]. In this paper, spacing, several function evaluations (NFE), GD, and IGD metric functions

are used to measure the performance of the proposed algorithm. IGD metric evaluates the convergence of the algorithm to the Pareto-optimal front. In addition, when the set of targeted Pareto-optimal points is defined with a suitable distribution, the IGD metric examines the

Fig. 13 Mean values for the IGD metric of the DTLZ3 problem

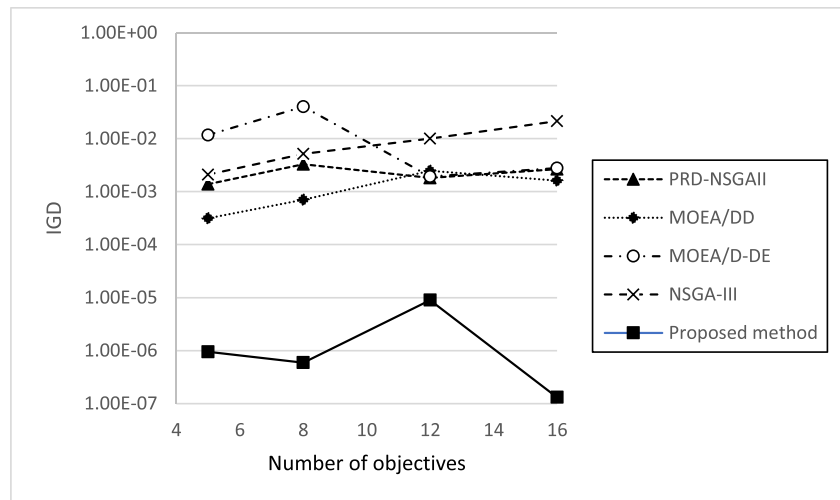
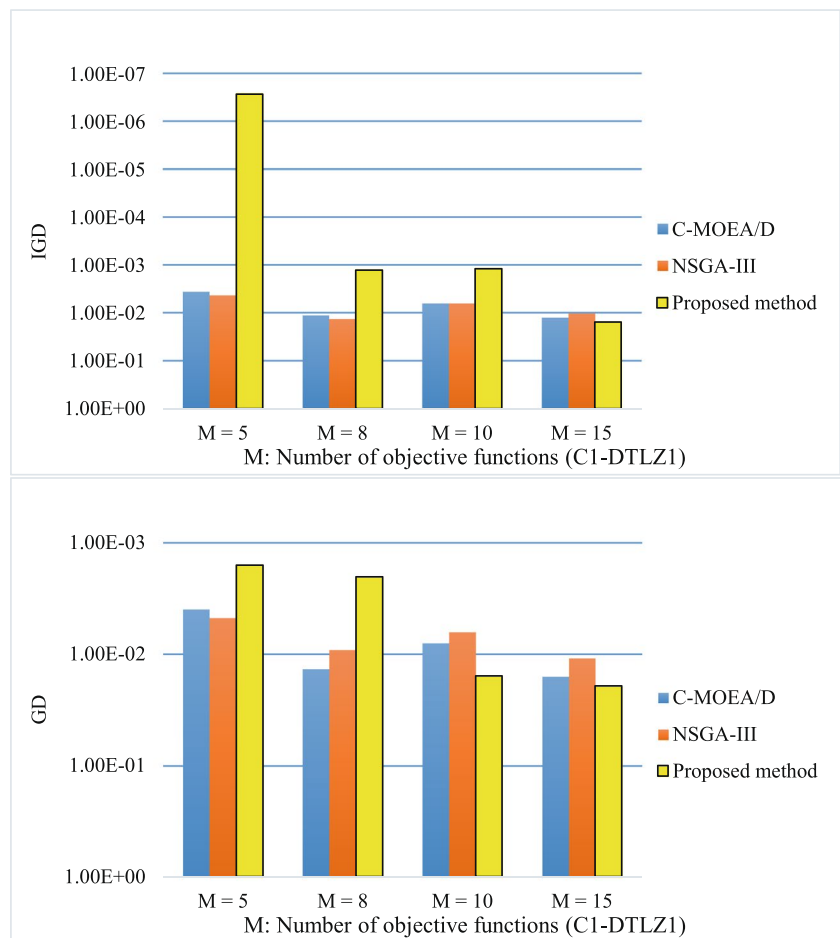


Fig. 14 Comparison of results of the proposed method and other algorithms for the C1-DTLZ1 benchmark problem



goodness of distribution of obtained solutions for the entire Pareto front. Another performance metric used to assess many-objective optimization algorithms is the GD metric, which measures how far the obtained solutions are away from the true Pareto front, which can reflect

the convergence of the algorithm. For unconstrained test problems, the solutions are close to the reference directions. Hence, the value of the GD is equal to the IGD. Therefore, GD values are presented for the constrained problems.

3.3 Validation

This section presents the validation and results of the proposed algorithm with other many-objective optimization algorithms using several challenging many-objective optimization benchmark problems. The obtained solutions for DTLZ1—DTLZ4 benchmark problems, with 3 objective functions, using the proposed algorithm and exact solutions, are shown in Fig. 7. The results of the proposed algorithm are shown with blue circles, and the exact solutions are characterized with red stars (Fig. 7). As can be seen from Fig. 8, the results of the proposed algorithm and those of exact solutions are in close agreement.

In addition, the IGD values that are calculated from the results of the MOPSO, and the proposed methods are compared with the recent many-objective optimization algorithms, as reported in I-DBEA [57]. The results of this comparative study using three and five-objective

functions, DTLZ1-DTLZ2 problems are shown in Table 4. It can be seen from Table 4, that MOPSO has performed a global search, and the obtained results are relatively close to the exact solution; however, the IGD performance metric obtained from the hybrid MOPSO-SQP (after applying SQP in the second step of the hybrid algorithm), show a considerable consistency of the obtained results with the exact solutions.

The effect of the proposed mutation operator in the MOPSO algorithm on the spacing and IGD performance metrics using a many-objective optimization benchmark problem C1-DTLZ1, compared to other many-objective optimization algorithms, is shown in Table 5. In Table 5, the best IGD and spacing performance metrics of the proposed and other algorithms shown in bold characters. As it can be seen from Table 5, the IBEA with 20,000 function evaluation has a slightly better IGD value, and EMOPSO with 5000 function evaluation spacing performance metric value,

Table 12 Median IGD and GD values for C1-DTLZ1 benchmark problem (bold items show the best performance). (M=Number of objective functions)

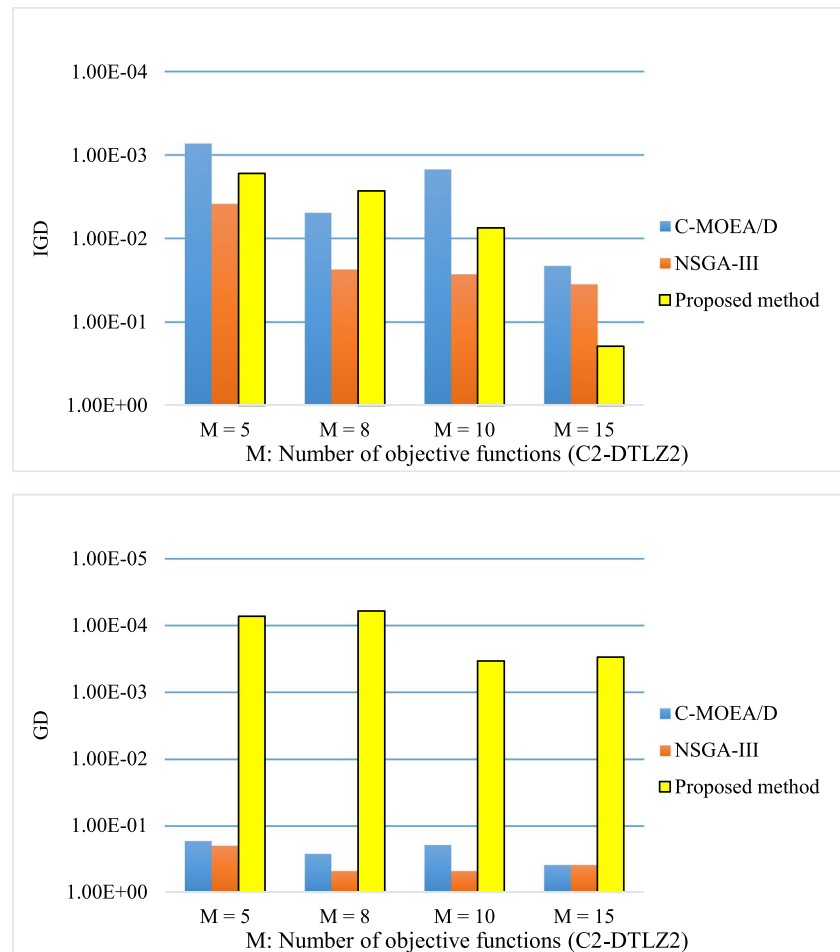
Problem	Max.Gen/ Pop.Size	Method	IGD	GD	Simulation Time (sec- ond)
C1-DTLZ1 (M=5)	600/210	C-MOEA/D	3.64 e -3	3.96 e -3	136.5
		NSGA-III	4.35 e -3	4.73 e -3	277.9
		Proposed method	2.62 e -7	1.6 e -3	61.2
C1-DTLZ1 (M=8)	800/156	C-MOEA/D	1.14 e -2	1.36 e -2	145.1
		NSGA-III	1.36 e -2	9.15 e -3	205.9
		Proposed method	1.26 e -3	2.04 e -3	44.7
C1-DTLZ1 (M=10)	1000/275	C-MOEA/D	6.41 e -3	7.96 e -3	343.8
		NSGA-III	6.36 e -3	6.36 e -3	226.1
		Proposed method	1.18 e -3	1.58 e -2	82.6
C1-DTLZ1 (M=15)	1500/135	C-MOEA/D	1.26 e -2	1.59 e -2	278.0
		NSGA-III	1.04 e -2	1.09 e -2	232.8
		Proposed method	1.54 e -2	1.94 e -2	131.2

Table 13 Median IGD and GD values for the C2-DTLZ2 benchmark problem (bold items show the best performance). (M=Number of objective functions)

Problem	Max.Gen/ Pop.Size	Method	IGD	GD	Simulation Time (second)
C2-DTLZ2 (M=5)	350/210	C-MOEA/D	7.30 e -4	1.68 e -1	92.9
		NSGA-III	3.87 e -3	1.98 e -1	29.4
		Proposed method	1.65 e -3	7.34 e -5	39.5
C2-DTLZ2 (M=8)	500/156	C-MOEA/D	4.97 e -3	2.60 e -1	105.4
		NSGA-III	2.35 e -2	4.73 e -1	40.8
		Proposed method	2.68 e -3	6.1 e -5	52.4
C2-DTLZ2 (M=10)	750/275	C-MOEA/D	1.49 e -3	1.91 e -1	287.9
		NSGA-III	2.69 e -2	4.72 e -1	85.4
		Proposed method	7.44 e -3	3.45 e -4	71.7
C2-DTLZ2 (M=15)	1000/135	C-MOEA/D	2.13 e -2	3.83 e -1	168.1
		NSGA-III	3.54 e -2	3.84 e -1	91.4
		Proposed method	1.96 e -1	3.02 e -4	110.7

Values in boldface show the best performance of the algorithms as shown in the tables

Fig. 15 Comparison of results of the proposed method and other algorithms for the C2-DTLZ2 benchmark problem



receptively, but in all other cases, the proposed method performs better than other algorithms used in this case study.

3.4 Effect of the neighboring point technique in the proposed algorithm

To illustrate the performance of the proposed neighboring point technique, the proposed algorithm with and without adopting the neighboring point technique is implemented on the DTLZ1 benchmark problem, and the results are shown in Table 6. As it can be seen from Table 6, the proposed method with the neighboring point technique performs better than without the neighboring point technique.

3.4.1 Examining the performance of the proposed method using unconstrained benchmark problems

This section presents a comparative study of the proposed method with other many-objective optimization algorithms using unconstrained many-objective optimization benchmark problems as described in Table 2. The results

of the proposed method are compared with the results of RPD-NSGA-II and MOEA/DD, MOEA/D-DE [71], and NSGA-III, as shown in Tables 7, 8, 9, 10 and 11 and Figs. 9, 10, 11, 12, 13, 14.

In this comparative study, both NFE and IGD performance metrics are used to evaluate the performance of the above algorithms, as shown in Tables 7–11 and Figs. 9–14. MOEA/DD (Table 7) performs slightly better than other algorithms when NFE is equal to 50,000, but the proposed method gives better accuracy for NFE of 75,000 and 1,000,000. For NFE = 75,000 and NFE = 100,000, the results of the proposed method have a high degree of accuracy than other algorithms such as; RPD-NSGA-II [33], MOEA/DD [72], MOEA/D-DE [73], NSGA-III [38], GrEA [74], KnEA [74], tDEA [75], SPEA-III [76], and MOSA [77].

As described above, among the different performance metrics for comparing the results of many-objective optimization algorithms, the IGD performance metric was selected because it reflects both convergence and divergence of results.

In addition, the proposed method was applied to the DTLZ1, DTLZ2, DTLZ3, and DTLZ4 many-objective optimization benchmark problems with a number of function

Fig. 16 Comparison of results of the proposed method and other algorithms for the C3-DTLZ4 benchmark problem

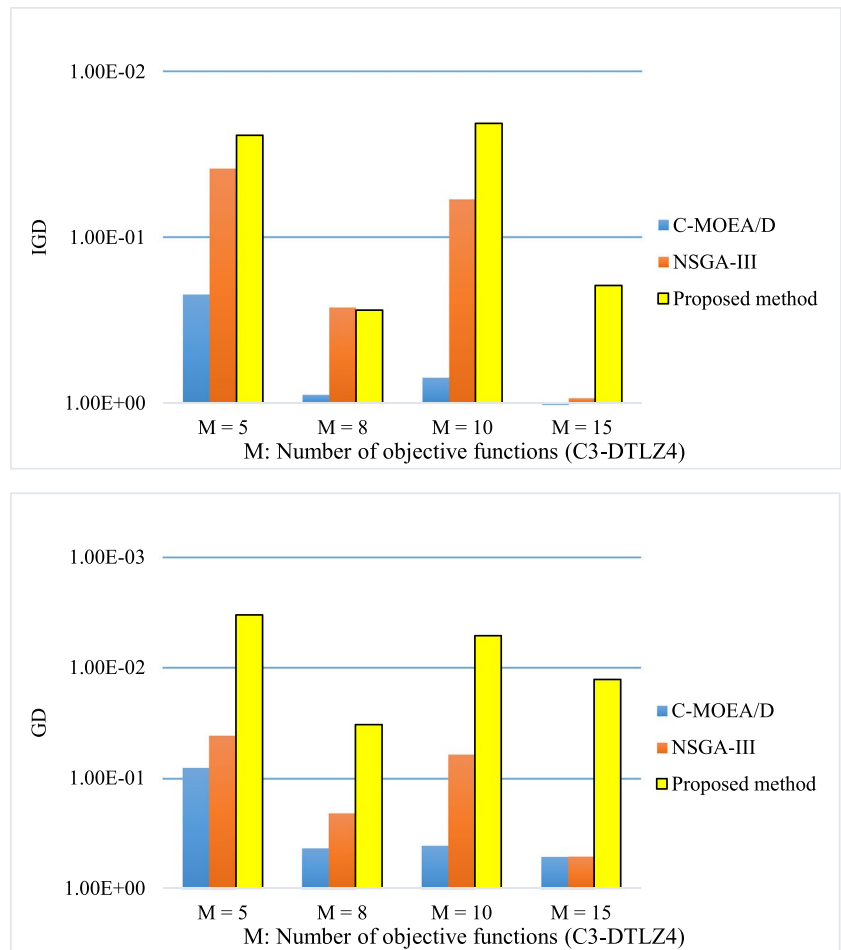


Table 14 Median IGD and GD values for the C3-DTLZ4 benchmark problem (bold items show the best performance). (M=Number of objective functions)

Problem	Max.Gen/ Pop.Size	Method	IGD	GD	Simulation Time (sec- ond)
C3-DTLZ4 (M=5)	1250/210	C-MOEA/D	2.21 e -1	8.01 e -2	364.0
		NSGA-III	3.86 e -2	4.09 e -2	98.4
		Proposed method	2.42 e -2	3.36 e -3	159.9
C3-DTLZ4 (M=8)	2000/156	C-MOEA/D	8.92 e -1	4.28 e -1	378.7
		NSGA-III	2.65 e -1	2.07 e -1	129.5
		Proposed method	2.75 e -1	3.30 e -2	250.0
C3-DTLZ4 (M=10)	3000/275	C-MOEA/D	7.01 e -1	4.06 e -1	1307.8
		NSGA-III	5.93 e -2	6.06 e -2	307.4
		Proposed method	2.05 e -2	5.18 e -3	302.8
C3-DTLZ4 (M=15)	4000/135	C-MOEA/D	1.45 e 0	5.12 e -1	698.6
		NSGA-III	9.32 e -1	5.08 e -1	339.9
		Proposed method	1.95 e -1	1.29 e -2	385.9

Values in boldface show the best performance of the algorithms as shown in the tables

evaluations of 100,000, and the results of the proposed method and other algorithms are shown in Table 8–11 and Figs. 9–14. Table 8 shows the results of the proposed and other algorithms using the DTLZ1 benchmark problem and

the IGD performance metric values. The IGD accuracy is in the order of 10^{-5} to 10^{-6} . In the DTLZ2 benchmark problem, $K=10$, and hence the number of decision variables for each objective function is more than that in the case of DTLZ1.

Fig. 17 Comparison of results with, and without parallelization of the proposed method

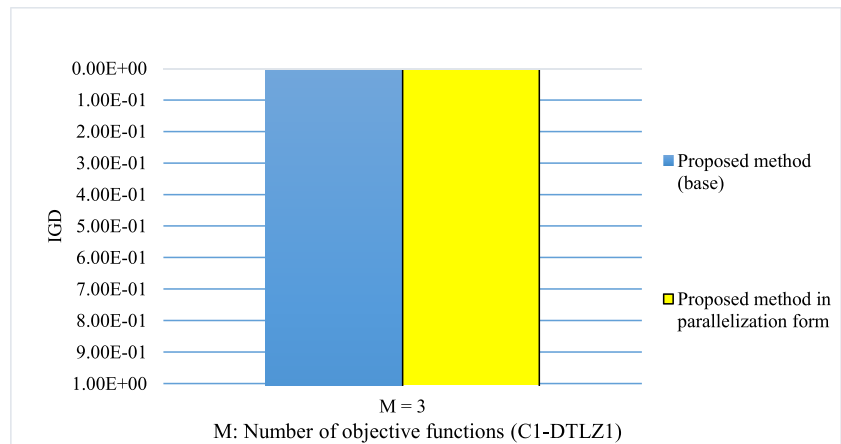


Table 15 Comparison of the proposed method with and without parallelization of the median IGD values for the C1-DTLZ1 benchmark problem (bold items show better performance). (M = Number of objective functions). (NFE = Number of function evaluations)

Problem	Method	IGD	NFE
C1-DTLZ1 ($M=3$)	The proposed method (without parallelization)	1.5156e0	477,360
	The proposed method with parallelization	1.1043e0	5000

Values in boldface show the best performance of the algorithms as shown in the tables

But it defines unimodal problems, and because of this characteristic, more accuracy is achievable (with the order of 10^{-11}).

The benchmark many-objective optimization problem DTLZ3 is more challenging because this benchmark problem

has more than 30^9 local Pareto optimal fronts [73], which is considered the most challenging many-objective optimization benchmark problem in the DTLZ family in terms of the difficulties associated with the convergence. This characteristic of the benchmark problem necessitates using a multi-start strategy in the global search stage of the proposed method.

Figures 9–14 show the IGD performance metrics for DTLZ, DTLZ2, DTLZ3, and DTLZ4 many-objective benchmark problems. These figures show that the proposed method outperforms all other algorithms in terms of convergence to the optimal Pareto frontier in all cases used in the above comparative study.

3.4.2 Examining the performance of the proposed method using constrained benchmark problems

This section presents the computational performance of the proposed method. In this respect, the proposed method was

Fig. 18 Comparison of results of the proposed method, FR-NSGA-II, and NSGA-II algorithms using CS_L -dominance metric

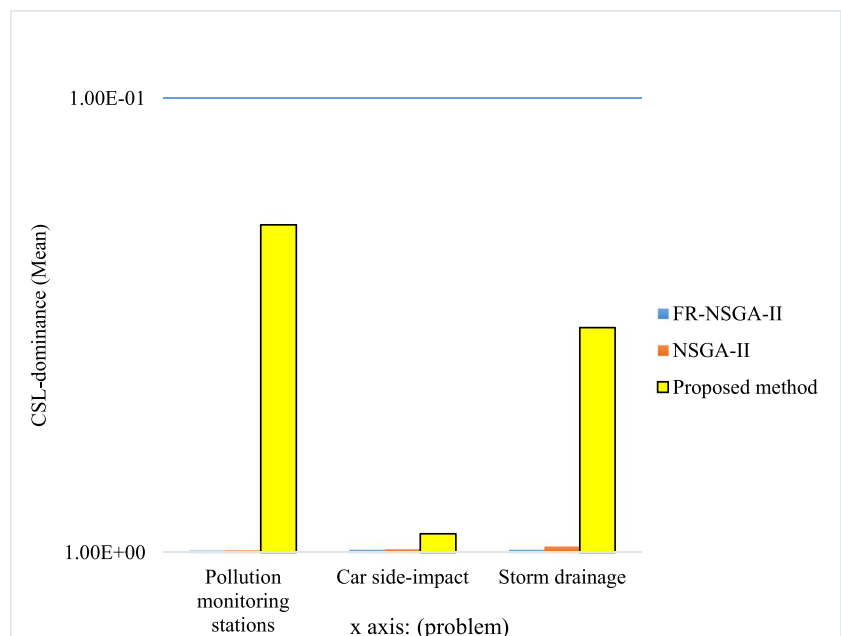


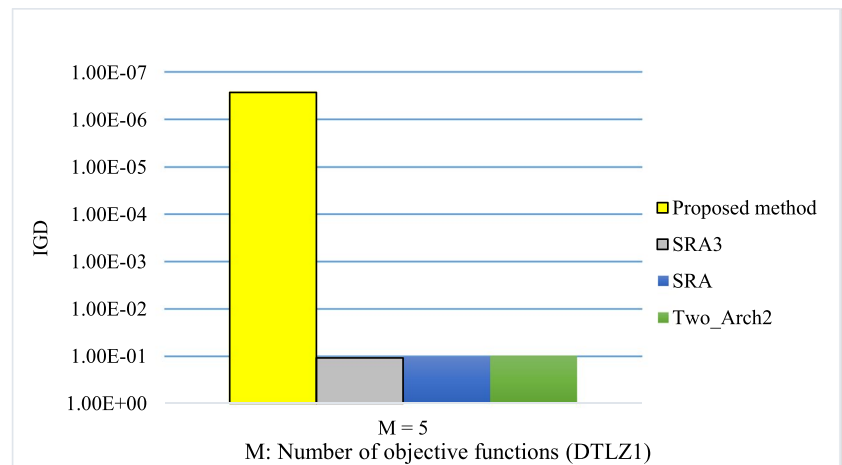
Table 16 Real-world test problems characteristics

Real-world test problem	Number of decision variables	Number of objective functions	Number of constraints
Location problem of pollution monitoring stations [78]	2	5	-
Car side-impact problem [79]	7	11	10
Storm drainage [80]	3	5	7

Table 17 Comparison of the proposed method, FR-NSGA-II, and NSGA-II algorithms using $CS_{L\text{-dominance}}$ metric (bold items show the best performance)

Problem	Algorithm	$CS_{L\text{-dominance}}$	
		Mean	Std
Pollution monitoring stations	Proposed method	0.9887	0.0048
	FR-NSGA-II	0.9871	0.0315
	NSGA-II	0.1900	0.3010
Car side-impact	Proposed method	0.9839	0.0031
	FR-NSGA-II	0.981	0.0094
	NSGA-II	0.9090	0.1908
Storm drainage	Proposed method	0.9846	0.0081
	FR-NSGA-II	0.9673	0.0153
	NSGA-II	0.32	0.1443

Values in boldface show the best performance of the algorithms as shown in the tables

Fig. 19 Comparison of results of the proposed and SEA3, SRA and Two_Arch2 method with five objective functions using DTLZ1 benchmark problem**Table 18** Median IGD values for the DTLZ1 benchmark problem (bold items show the best performance). (M=Number of objective functions)

Problem (Number of objective functions)	Number of function evaluations (NFE)	90,000
	Method	IGD
DTLZ1 (M=5)	SRA3	1.09×10^{-1}
	SRA	1.02×10^{-1}
	Two_Arch2	9.64×10^{-2}
	Proposed method	2.14×10^{-5}

applied to C1-DTLZ1 and C2-DTLZ2, constrained many-objective optimization benchmark problems. The results of the proposed method and other many-objective optimization algorithms for the above benchmark problems are shown in Tables 12 and 13. These results (Tables 12 and 13), show that for the benchmark problem C1-DTLZ1, with 15 objective functions, the NSGA-III has a better GD performance metric (Table 13 and Fig. 15). The C-MOEA/D algorithm for the benchmark problem C2-DTLZ2, with 15 and 10 objective functions has a better IGD performance metric. However, the proposed method in all other cases outperforms all other algorithms used in this study, both in terms of IGD, and GD performance metrics, as shown in Tables 12 and 13 and Figs. 15 and 16.

The results of the proposed methods and other algorithms for C3-DTLZ4, many-objective optimization benchmark

Table 19 Median IGD values for the DTLZ2 benchmark problem (bold items show the best performance). (M=Number of objective functions)

Problem (Number of objective functions)	Number of function evaluations (NFE)	90,000
	Method	IGD
DTLZ2 (M=5)	SRA3	1.80×10^{-1}
	SRA	1.76×10^{-1}
	Two_Arch2	1.52×10^{-1}
	Proposed method	8.7×10^{-4}

Values in boldface show the best performance of the algorithms as shown in the tables

Table 20 Median IGD values for the DTLZ3 benchmark problem (bold items show the best performance). (M=Number of objective functions)

Problem (Number of objective functions)	Number of function evaluations (NFE)	90,000
	Method	IGD
DTLZ3 (M=5)	SRA3	1.78×10^{-1}
	SRA	2.01×10^{-1}
	Two_Arch2	1.89×10^{-1}
	Proposed method	9.54×10^{-7}

Values in boldface show the best performance of the algorithms as shown in the tables

Table 21 Median IGD values for the DTLZ4 benchmark problem (bold items show the best performance). (M=Number of objective functions)

Problem (Number of objective functions)	Number of function evaluations (NFE)	90,000
	Method	IGD
DTLZ4 (M=5)	SRA3	1.79×10^{-1}
	SRA	1.78×10^{-1}
	Two_Arch2	1.53×10^{-1}
	Proposed method	3.56×10^{-14}

Values in boldface show the best performance of the algorithms as shown in the tables

problem is shown in Table 14 and Fig. 17. These results show that the proposed method significantly outperforms other many-objective optimization algorithms in all cases with 5, 8, 10 and 15, and objective functions in terms of both IGD and GD performance metrics used in this study (Table 14 and Fig. 17).

As described earlier in this paper, a parallelization scheme is adopted in the proposed method, in which optimization of all reference directions (weight vectors) is performed simultaneously. Table 15 shows a comparison between the

parallelization and without the parallelization of the proposed method on a benchmark problem C1-DTLZ1, with three objective functions. The obtained results (Table 15) show that the parallelization of the optimization has significantly enhanced the performance of the proposed method in terms of the NFE and, at the same time, has a comparable IGD performance metric concerning the case without parallelization of the proposed method as shown in Fig. 18.

3.4.3 Investigating the performance of the proposed method using real-world problems

This section presents applications of the proposed method to real-world many-objective optimization problems involving 5 and 11 objective functions. The computational performance of the proposed method compared with the other algorithms, in particular with one of the recent many-objective optimization algorithms (i.e., FR-NSGA-II [56]). The number of decision variables, objective functions, and constraints of the test problems is shown in Table 16. In the test problems of car side-impact and storm drainage problems, a penalty method was used for handling the constraints in MOPSO. In this respect, each test problem has been executed 10 times, and following this, average and standard deviations have been calculated to represent the computational performance of the algorithms used in this study.

Because true Pareto fronts are not known for real-world problems, the IGD metric is not suitable for comparing the performance of the algorithms. Therefore, the $CS_{L\text{-dominance}}$ metric, which is a modified version of two set coverage (CS) metrics, was used [56].

The results of the $CS_{L\text{-dominance}}$ metric for comparison of the proposed method and other algorithms using the above real-world test cases are given in Table 17 and Fig. 19. In all cases, the number of function evaluations is fixed to be 100,000. According to these results, the proposed method performs better than the FR-NSGA-II

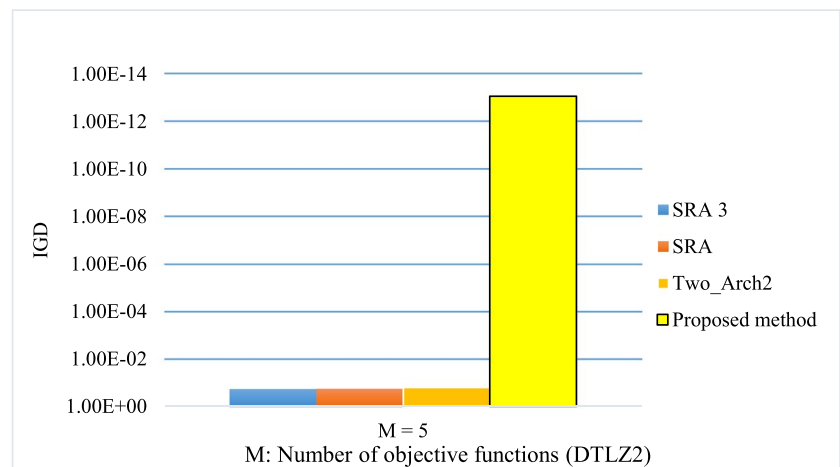
Fig. 20 Comparison of results of the proposed and SEA3, SRA and Two_Arch2 methods with five objective functions using DTLZ2 benchmark problem

Fig. 21 Comparison of results of the proposed and SEA3, SRA and Two_Arch2 methods with five objective functions using DTLZ3 benchmark problem

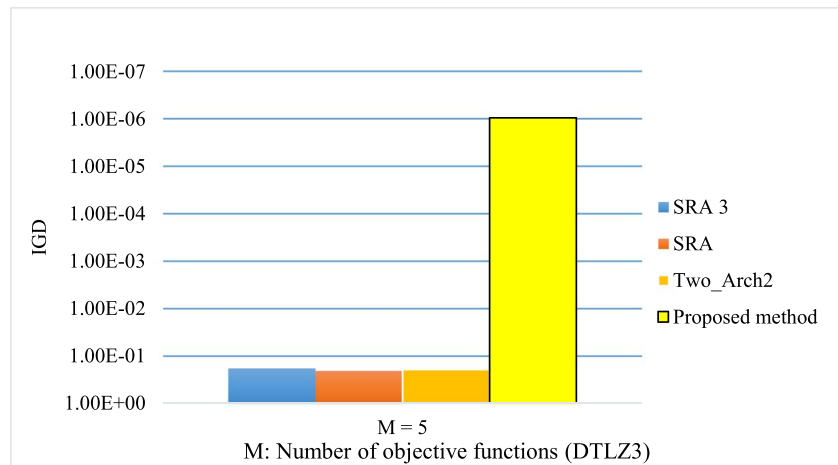
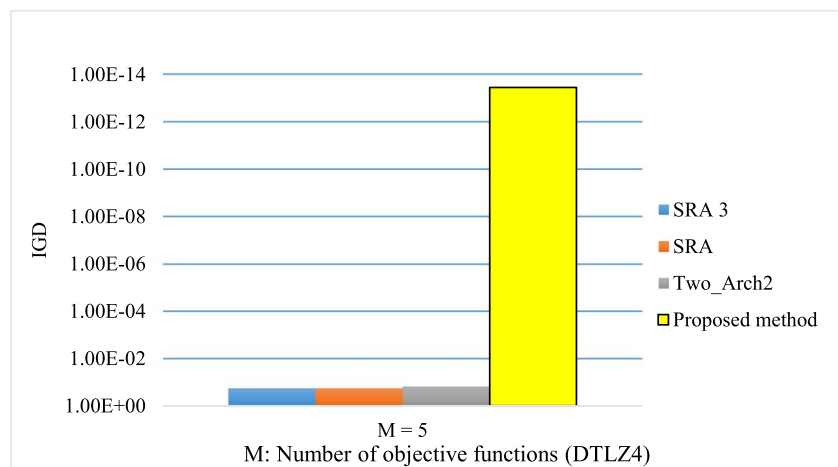


Fig. 22 Comparison of results of the proposed and SEA3, SRA and Two_Arch2 methods with five objective functions using DTLZ4 benchmark problem



and NSGA-II in terms of dominance for solving these real-world many-objective optimization problems. For calculating $CS_{L-dominance}$, the two last sets of Pareto solutions, which are obtained using the SQP algorithm of the proposed method, were compared based on the CS definition. As seen from Table 18, for the first test problem, $CS_{L-dominance}$, on average, is equal to 0.9887 for the proposed method, compared to the 0.9871 and 0.1900 values using FR-NSGA-II and NSGA-II, respectively. Similarly, with 11 objective car side-impact problems, the $CS_{L-dominance}$ obtained from the proposed method on average is 0.9839, compared to 0.981 for the FR-NSGA-II. Finally, for storm drainage application, the proposed algorithm reaches $CS_{L-dominance} = 0.9846$ on average, while this metric for FR-NSGA-II and NSGA-II when implemented are; 0.9673 and 0.32, respectively. The results of this experimentation of the proposed method's performance are shown in Fig. 19, in which the proposed method significantly outperforms the other algorithms used in this experimentation.

3.4.4 Comparative study of the performance of the proposed and other methods

In addition to comparative studies presented in the previous sections of this paper, in this section further performance of proposed many-objective optimization algorithms with SRA3, SRA, two_Arch2 [81], using DTLZ benchmark problems (Table 1), are compared. As described in the previous comparative studies, the results of this comparative study (Tables 18, 19, 20 and 21 and Fig. 19, 20, 21 and 22), also show that the proposed method outperforms SRA3, SRA and Two_Ach2.

4 Conclusions

This paper presented a novel decomposition-based hybrid many-objective optimization algorithm for solving many-objective optimization problems. Although several many-objective optimization algorithms have been developed; however, many-objective optimization algorithms present

new challenges in terms of the computational costs, convergence, and diversity of solutions. To address these research challenges, this paper focused on developing computationally efficient with a high degree of accuracy for solving many-objective optimization problems. The proposed method combines the global search capability of the PSO algorithm with the fast and accurate local search capability of the SQP for solving many-objective optimization problems. In addition, a new mutation operator introduced to enhance the performance of the MOPSO. The main feature of the proposed mutation operator is considering particles that violate the side constraint and returning them to the optimization process by replacing their position with a feasible location. In addition, a multi-start strategy utilized in the proposed method to overcome the stochastic nature of this evolutionary algorithm and to support the proposed method to reach more repeatable solutions.

The accuracy and computational performance of the proposed method were evaluated and compared with other many-objective optimization algorithms using several challenging many-objective optimization benchmarks and real-world test problems. The results of these experimental studies are shown in Tables 4–17 and Fig. 7–19. The results show that the proposed method outperforms other many-objective optimization algorithms in all cases used in this paper.

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Data availability Data will be made available on reasonable request.

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