# A new differential mutation base generator for differential evolution

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**Abstract** A new differential mutation base strategy for differential evolution (DE), namely best of random, is proposed. The best individual among several randomly chosen individuals is chosen as the differential mutation base while the other worse individuals are donors for vector differences. Hence both good diversity and fast evolution speed can be obtained in DE using the new differential mutation base. A comprehensive comparative study is carried out over a set of benchmark functions. Numerical results show that a better balance of reliability and efficiency can be obtained in differential evolution implementing the new generator of differential mutation base, especially in functions with high dimension.

Keywords Differential evolution · Differential mutation · Best of random · Diversity

# **1** Introduction

Differential evolution (DE) [1-6] is a simple, efficient and robust evolutionary algorithm (EA). It has found increasing applications in a number of scientific and engineering fields. Many researches have reported that DE outperformed some other EAs when optimizing benchmark functions or real-world problems [4-12].

In the community of differential evolution [2,13], differential evolution algorithms are marked as DE/x/y/z where x indicates how the differential mutation base is chosen,  $y \ge 1$  is the number of vector differences added to the differential mutation base vector, and z indicates which kind of crossover operation is used. It is well known that differential mutation

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is the crucial idea behind the success of differential evolution. Best and random are the two most often implemented differential mutation bases.

Through a comprehensive parametric study on differential evolution [5,14,15], it has been observed that differential evolution algorithms with best differential mutation base are in general more efficient while those with random differential mutation base are in general more reliable. Neither of them achieves harmony between reliability and efficiency, a challenge facing all evolutionary algorithms.

Generally speaking, differential mutation can be regarded as local search around differential mutation base. Guidance provided by the best differential mutation base leads to higher efficiency at the cost of diversity while blind search by using random differential mutation bases gains diversity and sacrifices efficiency. Hence, effectively balancing exploration and exploitation by using differential mutation base providing both guidance and diversity simultaneously might be a more promising solution. The new generator for differential mutation base, namely the best of random differential mutation base, is accordingly proposed.

The new generator for differential mutation base is very simple and introduces no additional intrinsic control parameter. Like random differential mutation, best of random differential mutation also randomly chooses several mutually different individuals. However, unlike random differential mutation, best of random differential mutation uses the best individual among these individuals as differential mutation base while the other worse individuals are donors for vector differences.

The new differential mutation base has been examined by comprehensive simulations over a set of benchmark functions. Numerical results show that a better balance of reliability and efficiency can be obtained in differential evolution implementing the new generator of differential mutation base, especially in functions with high dimension.

The rest of this paper is arranged as follows. Section 2 briefly describes differential evolution. Details of the new generator of differential mutation base are given in Sect. 3. A platform for evaluating differential evolution is constructed in Sect. 4. Section 5 presents numerical results of the conducted numerical evaluation. Conclusion is given in Sect. 6.

# 2 Differential evolution

## 2.1 Convention

Minimizing a single objective function  $f(\mathbf{x})$  with *N*-dimensional real optimization parameters  $\mathbf{x}$  without any constraint is considered in this paper. Natural real code is adopted.

Differential evolution operates on a population containing  $N_p$  individuals. It involves two stages, namely, initialization and evolution. Initialization generates initial population  $\mathbf{P}^0$ . Then  $\mathbf{P}^0$  evolves to  $\mathbf{P}^1$ ,  $\mathbf{P}^1$  evolves to  $\mathbf{P}^2$ , ..., and so on, until the termination conditions are fulfilled. While evolving from  $\mathbf{P}^n$  to  $\mathbf{P}^{n+1}$ , the three evolutionary operations, namely, differential mutation, crossover and selection are executed in sequence.

# 2.2 Classic differential evolution

The Fortran-style pseudo-code of classic differential evolution (CDE) is shown in Fig. 1 where the vector of optimization parameters  $\mathbf{x}_{i,n}$  is the *i*th individual in population  $\mathbf{P}^n$ ,  $\mathbf{v}_{i,n+1}$  is the mutant corresponding to  $\mathbf{x}_{i,n}$ ,  $\mathbf{b}_{i,n}$  is the differential mutation base,  $p_{1y}$  and  $p_{2y}$  are integer numbers, usually random,  $F_y$ , a constant usually in [0,1], is the mutation intensity for the *y*<sup>th</sup> vector difference  $\mathbf{x}_{p_{1y},n} - \mathbf{x}_{p_{2y},n}$ ,  $\mathbf{x}_{p_{1y},n}$  and  $\mathbf{x}_{p_{2y},n}$  are the donors for vector difference,

Initialization n = 0do  $i = 1, N_n$  $do_{j} = 1, N$  $x_{i,i,0} = b_i^L + \alpha_{i,i}(b_i^U - b_i^L)$ end do evaluate  $f(\mathbf{x}_{i,0})$ end do Evolution do while termination conditions are not satisfied do  $i = 1, N_p$ *differential mutation* to obtain mutant  $\mathbf{v}_{i, n+1}$ choose differential mutation base  $\mathbf{b}_{i,n}$ , donors  $\mathbf{x}_{p_{1,v},n}$  and  $\mathbf{x}_{p_{2,v},n}$  $\mathbf{v}_{i,n+1} = \mathbf{b}_{i,n} + \sum_{y>1} F_y(\mathbf{x}_{p_{1y},n} - \mathbf{x}_{p_{2y},n}), \ 1 \le i \ne p_{1y} \ne p_{2y} \le N_p$ regularize infeasible mutant  $\mathbf{v}_{i, n+1}$  $do_{j} = 1, N$ do while  $(v_{i, j, n+1} < b_i^L)$  $v_{i,i,n+1} = v_{i,i,n+1} + b_i^U - b_i^L$ end do do while  $(v_{i, j, n+1} > b_j^U)$  $v_{i,i,n+1} = v_{i,i,n+1} - b_i^U + b_i^L$ end do end do crossover mutant  $\mathbf{v}_{i, n+1}$  with target individual  $\mathbf{x}_{i, n}$  to deliver child  $\mathbf{c}_{i, n+1}$ evaluate child  $f(\mathbf{c}_{i, n+1})$ selection to get individual  $\mathbf{x}_{i, n+1}$ if  $f(\mathbf{c}_{i, n+1}) < f(\mathbf{x}_{i, n})$  $\mathbf{x}_{i, n+1} = \mathbf{c}_{i, n+1}$ else  $\mathbf{x}_{i,n+1} = \mathbf{x}_{i,n}$ end if end do n = n + 1end do

Fig. 1 Fortran-style pseudo-code of classic differential evolution

 $\mathbf{c}_{i,n+1}$  is the child or trial individual generated by crossover.  $\alpha_{i,j}$  is a real random number in [0,1],  $[b_i^L, b_i^U]$  is the search space of the *j*th optimization parameter,  $x_{i,j}$ .

# 2.2.1 Differential mutation base

There are various differential mutation base generators among which best and random differential mutation bases are the most often implemented.

A. Best differential mutation base

In this scheme, the best or the most dominant (in terms of objective function value) individual  $\mathbf{x}_{best,n}$  in  $\mathbf{P}^n$  is chosen as the differential mutation base.

B. Random differential mutation base

In this scheme, the differential mutation base is randomly chosen from  $\mathbf{P}^n$ .

#### 2.2.2 Crossover

Likewise, different crossover schemes are applicable, among which binomial crossover and exponential crossover are the most often implemented.

A. Binomial crossover

In this scheme,  $\mathbf{c}_{i,n+1}$  is generated as follows

$$c_{i,j,n+1} = \begin{cases} v_{i,j,n+1} & \beta_{i,j,n+1} \le C_r \\ x_{i,j,n} & \text{otherwise} \end{cases}$$
(1)

where  $\beta_{i,j,n+1}$  is a real random number uniform in the range [0,1] and  $C_r$ , a constant in [0,1], is the crossover probability.

There is an extreme case. The child duplicates  $\mathbf{v}_{i,n+1}$ . In this case, a randomly chosen parameter of  $\mathbf{x}_{i,n}$ ,  $x_{i,j,n}$ , will replace the corresponding parameter of the child  $\mathbf{c}_{i,n+1}$ ,  $c_{i,j,n+1}$ .

There is another extreme case.  $\mathbf{c}_{i,n+1}$  inherits no parameter from  $\mathbf{v}_{i,n+1}$  and hence no evolution happens. In this case, a randomly chosen parameter of the child  $\mathbf{c}_{i,n+1}$ ,  $c_{i,j,n+1}$ , will be replaced by the corresponding parameter of the mutant  $\mathbf{v}_{i,n+1}$ ,  $v_{i,j,n+1}$ .

B. Exponential crossover

The Fortran-style pseudo-code is given in Fig. 2.

```
do k = 1, N

C_{i,j,n+1} = x_{i,j,n}

end do

r = N * rand(0, 1) + 1

k = r

C_{i,k,n+1} = v_{i,k,n+1}

M = 1

\beta = rand(0, 1)

do while (\beta \le C_r and M < N - 1)

M = M + 1

\beta = rand(0, 1)

k = 1 + mod(k, N)

C_{i,k,n+1} = v_{i,k,n+1}

end do
```

Fig. 2 Fortran-style pseudo-code for exponential crossover for differential evolution

#### 2.2.3 Termination conditions

In this paper, the following two termination conditions are applied simultaneously. A. Objective met

This termination condition is only applicable to optimization problems whose optima are known. Such optimization problems include multi-dimensional benchmark functions and benchmark application problems for unconstrained single-objective optimization. Usually, they are involved in algorithm evaluation.

In this study, it can be mathematically expressed as

$$f\left(\mathbf{x}_{p,n}\right) - f\left(\mathbf{x}^{*}\right) \le \varepsilon_{f} \quad \exists 1 \le p \le N_{p} \tag{2}$$

where  $f(\mathbf{x}^*)$  is the known optima,  $\varepsilon_f$  is the predefined error tolerance. B. Limit of number of objective function evaluations

Unlimited time to search optimal solution is unaffordable in practice. The search process has to be terminated when the consumed time exceeds certain limit. Although limit on search time is straightforward, it does not capture the essence of optimization and is dependent on hardware. Instead, limit of number of objective function evaluations is a hardware-free marker for consumed time.

# 2.3 Dynamic differential evolution

Classic differential evolution is static in nature from the point of view of population updating and is therefore inherently inefficient. Inspired by the advantage of Gauss-Seidel method against Jacobi method for linear equations, the dynamic differential evolution (DDE) [16] was developed. The Fortran-style pseudo-code for dynamic differential evolution is shown in Fig. 3, where the generation indexes n and n + 1 do not show up any more. This indicates that all evolutionary operations are executed over individuals in the current population. There is only one population whose individuals are continuously updated by their competitive children. For more details of dynamic differential evolution, please refer to [16].

Extensive performance evaluation has been carried out [5]. As expected, in general, dynamic differential evolution outperforms classic differential evolution. But comparing with CDE, DDE has a disadvantage. Unlike CDE, it cannot be parallelized. For the problems where the function evaluation requires much computation time, parallelization is important hence CDE is more suitable than DDE to solve them.

# 3 Best of random differential mutation base

For most evolutionary algorithms including DE, how to balance two contradictory aspects: exploration and exploitation, is a crucial problem. In fact, this problem can also be understood from another point of view: the balance of diversity and evolution of the population. Diversity here means the dispersing of the individuals in the population, while evolution means constant improvement of the (average) performance of the population as they evolve from one generation to the next. For a good balance of them, in principle we should assign more search resource to the promising regions for fast evolution speed while at the same time keep the diversity of the population from decreasing rapidly for reliable convergence.

Generally speaking, differential mutation operation in DE could be regarded as a local search. Differential mutation base vector acts as the local search center and the scaled vector difference determines the search range or search step around the center. The diversity of the

Initialization
n = 0
do $i = 1, N_p$
do j = 1, N
$x_{i,j} = b_j^L + \alpha_{i,j} (b_j^U - b_j^L)$
end do
evaluate $f(\mathbf{x}_i)$
end do
Evolution
n = 1
i = 1
do while termination conditions are not satisfied
<i>differential mutation</i> to obtain mutant $\mathbf{v}_i$
choose differential mutation base $\mathbf{b}_i$ , donors $\mathbf{x}_{p_{1y}}$ and $\mathbf{x}_{p_{2y}}$
$\mathbf{v}_{i} = \mathbf{b}_{i} + \sum_{y \ge 1} F_{y}(\mathbf{x}_{p_{1y}} - \mathbf{x}_{p_{2y}}),  1 \le i \ne p_{1y} \ne p_{2y} \le N_{p}$
regularize infeasible mutant $\mathbf{v}_i$
do $j = 1, N$
do while $(v_{i,j} < b_j^L)$
$v_{i,j} = v_{i,j} + b_j^U - b_j^L$
end do
do while ( $v_{i,j} > b_j^U$ )
$v_{i,j} = v_{i,j} - b_j^U + b_j^L$
end do
end do
<i>crossover</i> mutant $\mathbf{v}_i$ with target individual $\mathbf{x}_i$ to deliver child $\mathbf{c}_i$
evaluate child $f(\mathbf{c}_i)$
selection to get individual $\mathbf{x}_i$
$\inf f(\mathbf{c}_i) < f(\mathbf{x}_i)$
$\mathbf{x}_i = \mathbf{c}_i$
additional selection
$\inf f(\mathbf{c}_i) < f(\mathbf{x}_{best}) \mathbf{x}_{best} = \mathbf{c}_i$
end if
$i = 1 + \operatorname{mod}(i, N_p)$
$if(i=1) \ n=n+1$
end do

Fig. 3 Fortran-style pseudo-code of dynamic differential evolution

population at the next generation depends on the diversity of the local search centers (base vectors). In DE/rand/\*/\*, the base vectors are randomly chosen from the population, hence a good diversity of base vectors (local search centers) is gained. But as the base vectors have no preference for promising regions or solutions, the evolution speed of DE/rand/\*/\* is relatively slow. On the other hand, all the individuals in DE/best/\*/\* use the same base vector: the best vector so far. This kind of concentrated search around the best individual often leads to fast evolution speed but is more inclined to stagnation due to the lack of enough diversity of base vectors.

From the above analysis, it is noticed that using base vectors with high quality in DE often brings fast evolution speed, and increasing the diversity of base vectors also promotes the diversity of the population. Differential mutation base providing both guidance and diversity should be a more promising choice. Based on this idea, a new generator for differential mutation base, namely best of random, is proposed. Best of random differential mutation base uses the best individual among several randomly chosen individuals as differential mutation  $p_{b} = N_{p} * \operatorname{rand}(0, 1) + 1$   $p_{1} = N_{p} * \operatorname{rand}(0, 1) + 1$ do while  $(p_{1} = p_{b}) p_{1} = N_{p} * \operatorname{rand}(0, 1) + 1$ if  $f(\mathbf{x}_{p_{1},n}) \leq f(\mathbf{x}_{p_{b},n})$  swap  $p_{1}$  and  $p_{b}$   $p_{2} = N_{p} * \operatorname{rand}(0, 1) + 1$ do while  $(p_{2} = p_{b} \text{ or } p_{2} = p_{1}) p_{2} = N_{p} * \operatorname{rand}(0, 1) + 1$ if  $f(\mathbf{x}_{p_{2},n}) \leq f(\mathbf{x}_{p_{0},n})$  swap  $p_{2}$  and  $p_{b}$ 

Fig. 4 Fortran-style pseudo-code for best of random differential mutation base

base and the other worse individuals as donors for vector differences. Thus both good quality and diversity of the base vectors can be obtained. The formula for best of random differential mutation strategy can be expressed as:

$$\mathbf{v}_{i,n+1} = \mathbf{x}_{p_b,n} + \sum_{y \ge 1} F_y(\mathbf{x}_{p_{1y},n} - \mathbf{x}_{p_{2y},n}), \ 1 \le i \ne p_{1y} \ne p_{2y} \le N_p$$
(3)

where, the individuals  $\mathbf{x}_{p_{1y},n}, \mathbf{x}_{p_{2y},n}$  and  $\mathbf{x}_{p_b,n}$  are randomly chosen from the population, and  $\mathbf{x}_{p_b,n}$  is the best one of them, i.e.  $f(\mathbf{x}_{p_b,n}) \leq \min(f(\mathbf{x}_{p_{1y},n}), f(\mathbf{x}_{p_{2y},n}))$ . The Fortran-style pseudo-code of best of random differential mutation base generator is shown in Fig. 4. DE implementing best of random differential mutation base is denoted as DE/BoR/\*/\*. In this paper, DE/BoR/1/\* is used. It should be pointed out that there is no need to modify any other evolutionary operations in differential evolution. What we need is to write a separate subroutine for the new differential mutation base and replace existing subroutines for differential mutation base, as highlighted by shadow in Figs. 1 and 3. It can be found easily that DE/BoR/1/\* is almost as simple as DE/rand/1/\* and introduce no extra intrinsic control parameters.

#### 4 Evaluation platform

# 4.1 Test bed

For a fair comparison of different DE algorithms, a test bed consisting of 13 benchmark functions with different representative mathematical features, such as modality, separability, differentiability, symmetry, is used. The definition and mathematical features of the benchmark functions are given in Appendix.

All the benchmark functions are scalable for the sake of investigating the effect of dimensionality. Two different dimensions, N = 8 and 50 representing low and high dimensional problems respectively, are used in the simulations. For dimension 50, as the simulation of some functions are computationally very expensive, only nine functions are simulated. Function Griewank of dimension 50 is only simulated with CDE schemes as it is also computationally very expensive and it is observed that the simulation results of multimodal functions of dimension 50 in CDE and DDE are similar.

## 4.2 Differential evolution algorithms

Twelve differential evolution algorithms are involved in this study

1. CDE/best/1/bin (CBB)

- 2. CDE/rand/1/bin (CRB)
- 3. CDE/BoR/1/bin (CBoB)
- 4. CDE/best/1/exp (CBE)
- 5. CDE/rand/1/exp (CRE)
- 6. CDE/BoR/1/exp (CBoE)
- 7. DDE/best/1/bin (DBB)
- 8. DDE/rand/1/bin (DRB)
- 9. DDE/BoR/1/bin (DBoB)
- 10. DDE/best/1/exp (DBE)
- 11. DDE/rand/1/exp (DRE)
- 12. DDE/BoR/1/exp (DBoE)

4.3 Intrinsic control parameters

Here, the finite set of intrinsic control parameters is  $S_{icp} = S_{Np} \cup S_F \cup S_{Cr}$ . All combinations of 3 intrinsic control parameters from their sets  $S_{Np}$ ,  $S_F$  and  $S_{Cr}$  are simulated in our experiments.

# 4.3.1 Population size

Two different sets of population sizes are used

$$\begin{split} \mathbf{S}_{Np} &= \{16, 24, 32, 40, 48, 56, 64, 72, 80, 120\}\\ \mathbf{S}_{Np} &= \{20, 30, 40, 50, 60, 70, 80\} \end{split}$$

The first set of population size is for simulating 8-dimensional benchmark functions while the second set for 50-dimensional benchmark functions. Initially, the sets of population sizes,  $\mathbf{S}_{Np} = \{8, 16, 24, 32, 40, 48, 56, 64, 72, 80, 120, 160, 200, 400\}$  and  $\mathbf{S}_{Np} = \{10, 20, 30, 40, 50, 60, 70, 80, 90, 100, 150, 200, 250, 500\}$ , are devised for 8 and 50-dimensional functions in our previous parametric study of DE/best/1/\* and DE/rand/1/\* [5]. It has been observed from the simulation results that DE does not benefit from increasing of population size if it has been above the optimal population size. Thus in this paper we choose smaller sets of population sizes, but if necessary, other population sizes will be added for some benchmark functions in order to include the optimal population size.

# 4.3.2 Mutation intensity

The set of tested mutation intensity takes the form as  $S_F = \{F | F = j \times \Delta_F, 1 \le j \le m/\Delta_F\}, m = 1$  unless specified otherwise.  $\Delta_F = 0.025$  is used in this paper.

# 4.3.3 Crossover probability

The set of tested crossover probability  $\mathbf{S}_{Cr} = \{C_r | C_r = j \times \Delta_{Cr}, 1 \le j \le 1/\Delta_{Cr}\}, \Delta_{Cr}$  is usually equal to  $\Delta_{F}$ .

# 4.4 Termination conditions

The error tolerance  $\varepsilon_f$  for each benchmark function is set to be  $1 \times 10^{-2}$ . The default limit of number of objective function evaluation is 2000 times the number of optimization parameters.

But it is increased to 64,000, 128,000 and 128,000 for 8-dimensional functions Griewank, Whitley and Weierstrass, respectively, and 200,000 for 50-dimensional Griewank function.

# 4.5 Performance indicators

Performances of evolutionary algorithms include reliability, efficiency, and robustness, in the order of importance, or priority. Reliability shows how confident the desired optima can be found by the evolutionary algorithm, efficiency shows how fast the evolutionary algorithm finds the desired optima, while robustness shows how sensitive the algorithm's efficiency is to other factors, such as randomness, intrinsic and non-intrinsic control parameters.

Evaluation of evolutionary algorithms can be done at two levels, i.e., parameter level and system level. Parameter level is the low level while system level is the high level.

# 4.5.1 Search and trial

A search is an execution of an evolutionary algorithm with fixed intrinsic control parameters. It is successful if it finds the desired optima. The number of objective function evaluations (NOFE) of a successful search is recorded.

To reduce randomness, each search with fixed intrinsic control parameters is repeated 100 times in this paper, such a repetition is a trial. The number of successful searches (NSS) in a trial is recorded. A trial with at least one successful search is a partially successful trial. In accordance, the average number of objective function evaluations (ANOFE) of all successful searches in a partially successful trial is computed. If all of the searches in a trial are successful, it is a successful trial.

#### 4.5.2 Parameter level evaluation

So far, attention has been focused on evaluation of evolutionary algorithms at parameter level. At this level, evolutionary algorithms are evaluated with fixed intrinsic and non-intrinsic control parameters. The number of successful searches among a trial indicates reliability, the average number of objective function evaluations of all successful searches among the trial indicates efficiency, while the distribution of number of objective function evaluations of all successful searches among the trial indicates robustness.

The interaction between evolutionary algorithm, intrinsic and non-intrinsic control parameters has not been touched at parameter level evaluation. Evolutionary crimes are often observable within such evaluations[5]. From the point of view of practical application, such evaluation is not enough.

## 4.5.3 System level evaluation

At system level, instead, evolutionary algorithms are evaluated with many different combinations of control parameters. Evaluation of evolutionary algorithms at system level regards evolutionary algorithm, its intrinsic and non-intrinsic control parameters, and even problem features as a system. The evolutionary algorithm is not isolated. On the contrary, the interaction between evolutionary algorithms, intrinsic and non-intrinsic control parameters is systematically addressed. The ultimate goal here is to find the optimal intrinsic control parameter values and the relationship between evolution algorithm, intrinsic and non-intrinsic control parameters, and problem features. It is exactly the demand of application engineers. However, by now, few people have paid attention to evaluation of evolutionary algorithms at system level.

At system level, many trials corresponding to different combinations of intrinsic control parameters, instead of one trial, are conducted. Attention is focused on successful trials. Under each population size,  $40 \times 40 = 1,600$  trials corresponding to different combinations of *F* and *C<sub>r</sub>* are conducted in this paper. The number of successful trials (NST) among the 1600 trials at each population size is recorded. The highest number of successful trials (HNST) among the NSTs of different population sizes is defined as the reliability indicator. By normalizing HNST to the total number of trials (TNT) under a fixed population size (TNT = 1,600 in this paper), the reliability indicator can also be expressed as the highest successful trial rate  $R_{HST}$  = HNST/TNT. The population size at which the HNST or  $R_{HST}$  is achieved is the optimal population size in terms of reliability. The minimal average number of objective function evaluations (MANOFE) among the ANOFEs of all successful trials at the optimal population size in terms of reliability is defined as the efficiency indicator. In this paper, no robustness indicator is defined since it is the least important performance for differential evolution at system level evaluation.

#### **5** Numerical results

# 5.1 8-Dimensional benchmark functions

The reliability and efficiency of differential evolution for solving 8-dimensional benchmark functions and the corresponding optimal intrinsic control parameters ( $N_p$ , F,  $C_r$ ,) are given in Table 1. The 12 participating DE algorithms are organized in 4 groups. Members in each group differ by differential mutation base. The best result in each group is highlighted in bold-face and the best result of all participating differential evolution algorithms is highlighted in bold italic.

With regard to the reliability, DE/BoR/1/\* is the most reliable among the three different differential mutation strategies in function f3. In function f6, DE/best/1/\* is the most reliable, DE/BoR/1/\* ranks second and DE/rand/1/\* is the least reliable. In functions f10 and f11, DE/rand/1/\* is a little bit more reliable than DE/BoR/1/\*, and DE/best/1/\* ranks the worst in terms of reliability. In the rest functions, the reliability of DE/BoR/1/\* is close to that of DE/rand/1/\*.

As for the efficiency, in functions  $f1 \sim f6$  and f8, DE/best/1/\* is the most efficient among the three different differential mutation strategies, and DE/BoR/1/\* ranks second. In functions f7, f9, f12 and f13, the efficiency of DE/BoR/1/\* is the best or close to the best among the three different differential mutation strategies. In function f10, CDE/rand/1/\* is a little bit more efficient than CDE/BoR/1/\* while DDE/rand/1/\* is almost as efficient as DDE/BoR/1/\*, and DE/best/1/\* is the least efficient. In function f11, DE/best/1/\* is the most efficient while the efficiency of DE/BoR/1/\* is close to that of DE/rand/1\*.

It is observed from Table 1 that the overall performance of DE/BoR/1/\* is never the worst in each function. In general, DE/BoR/1/\* is almost as reliable as but more efficient than DE/rand/1/\*. As a whole, a better balance of reliability and efficiency is obtained in DE/BoR/1/\* than in DE/best/1/\* and DE/rand/1/\*.

In order to find the relationship between DE algorithms and the mathematical features, such as modality, of the benchmark functions, the performance of DE algorithms with different differential mutation strategies in unimodal and multimodal functions are compared.

TUDY	•												
		CBB	CRB	CBoB	CBE	CRE	CBoE	DBB	DRB	DBoB	DBE	DRE	DBoE
Ų	$R_{ m HST}$ (%)	68.1	78.1	75.8	83.1	86.6	84.2	86.0	77.8	76.9	92.1	86.8	85.3
	MANOFE	1789	2765	1759	1624	3517	2255	<i>977</i>	2716	1506	707	2669	1343
	$N_p$	48	32	32	56	40	40	56	40	40	72	40	40
	F	0.500	0.250	0.275	0.350	0.250	0.325	0.250	0.275	0.250	0.175	0.325	0.300
	$C_r$	0.925	0.750	0.850	0.925	0.975	1.000	0.975	0.775	0.725	1.000	1.000	1.000
f2	$R_{ m HST}$ (%)	76.1	83.9	82.0	84.8	90.1	90.5	86.8	83.6	82.5	92.9	90.0	90.3
	MANOFE	1159	2902	1651	1296	3100	1854	575	1669	1104	586	2710	1384
	$N_p$	72	56	48	72	56	56	120	40	48	120	64	64
	F	0.375	0.225	0.275	0.450	0.025	0.250	0.050	0.225	0.325	0.100	0.250	0.275
	$C_r$	0.875	0.750	0.800	1.000	0.150	0.950	0.925	0.700	0.925	1.000	1.000	1.000
f3	$R_{ m HST}$ (%)	9.3	21.1	26.2	11.9	18.8	23.3	11.7	22.9	29.1	15.9	21.1	26.1
	MANOFE	4730	19195	13851	8716	15031	14634	1572	23493	11518	3728	14231	7746
	$N_{P}$	32	32	40	64	24	40	56	48	48	48	24	32
	F	0.575	0.650	0.575	0.550	0.625	0.600	0.450	0.600	0.525	0.525	0.675	0.550
	$C_r$	0.950	1.000	0.975	0.975	0.925	1.000	0.850	0.925	1.000	1.000	1.000	0.975
f4	$R_{ m HST}$ (%)	65.6	76.8	75.1	81.0	83.9	84.3	84.1	74.5	75.0	1.10	83.9	83.9
	MANOFE	1971	3159	1984	1791	4107	2552	936	2788	1380	818	3442	1978
	$N_{P}$	40	32	32	48	40	40	56	32	32	64	40	40
	F	0.475	0.225	0.325	0.375	0.225	0.350	0.275	0.350	0.350	0.175	0.275	0.400
	$C_r$	0.725	0.675	0.850	0.925	0.900	1.000	1.000	0.850	0.975	1.000	0.975	1.000
f5	$R_{ m HST}$ (%)	62.9	6.09	61.0	78.2	74.1	74.3	79.9	60.8	62.8	87.2	74.8	76.2
	MANOFE	1837	2050	1620	1622	2950	2188	959	1913	1278	842	2707	1667
	$N_p$	32	16	16	48	24	24	40	16	24	40	24	32
	F	0.500	0.175	0.450	0.350	0.100	0.325	0.375	0.300	0.375	0.325	0.075	0.350
	$C_r$	0.700	0.125	0.725	0.925	0.075	0.825	0.975	0.425	0.825	0.975	0.175	0.950

Table 1	continued												
		CBB	CRB	CBoB	CBE	CRE	CBoE	DBB	DRB	DBoB	DBE	DRE	DBoE
f6	$R_{ m HST}$ (%)	27.4	19.6	22.9	30.1	23.2	28.5	35.6	21.8	27.4	32.7	22.9	28.4
	MANOFE	2087	4074	3109	1994	4103	2304	1272	3825	2652	1867	2775	2464
	$N_p$	24	16	16	16	16	16	24	16	16	16	8	16
	F	0.525	0.575	0.550	0.650	0.550	0.375	0.425	0.550	0.600	0.650	0.700	0.475
	$C_r$	0.800	0.925	0.775	0.975	0.975	0.850	0.825	0.925	0.975	0.975	0.875	0.850
f7	$R_{ m HST}$ (%)	64.9	75.4	72.5	78.1	86.6	84.6	78.3	74.1	74.3	86.1	85.4	85.1
	MANOFE	2047	2363	1632	2796	3826	2822	1355	2109	1507	1749	2793	1997
	$N_{P}$	32	24	24	40	40	40	56	24	32	48	32	40
	F	0.300	0.200	0.175	0.325	0.125	0.200	0.300	0.225	0.200	0.350	0.200	0.225
	$C_r$	0.400	0.550	0.650	0.775	0.850	0.925	0.975	0.650	0.750	0.950	0.925	0.950
f8	$R_{ m HST}$ (%)	47.4	64.8	65.6	66.1	80.9	82.4	55.8	62.7	63.8	71.4	81.3	81.7
	MANOFE	3834	2154	2508	3002	3276	3265	1336	2045	1740	2798	3234	2887
	$N_{P}$	80	24	48	64	40	56	56	24	32	80	48	48
	F	0.350	0.100	0.050	0.100	0.100	0.075	0.200	0.200	0.125	0.300	0.025	0.100
	$C_r$	0.500	0.075	0.200	0.375	0.425	0.550	0.525	0.300	0.350	0.750	0.050	0.625
$f_{9}$	$R_{ m HST}$ (%)	18.4	34.2	29.4	40.5	63.3	61.3	17.4	33.9	33.4	39.6	62.2	63.2
	MANOFE	8609	4951	5165	5912	5024	4800	7116	4632	4342	6631	4311	4178
	$N_{p}$	56	32	40	56	32	40	64	32	40	64	32	40
	F	0.275	0.150	0.150	0.325	0.125	0.125	0.225	0.150	0.100	0.375	0.125	0.100
	$C_r$	0.175	0.125	0.025	0.450	0.350	0.300	0.050	0.125	0.100	0.475	0.125	0.300
f10	$R_{ m HST}$ (%)	6.8	44.4	37.9	20.0	74.8	68.1	6.8	41.8	41.9	22.4	75.7	71.8
	MANOFE	15992	6328	7213	25852	8062	9770	15744	7728	7882	19930	6620	6480
	$N_p$	72	40	56	140	56	72	72	48	64	120	48	64
	F	0.500	0.125	0.075	0.325	0.075	0.075	0.5009	0.100	0.150	0.500	0.075	0.225
	$C_r$	0.025	0.200	0.100	0.225	0.300	0.050	0.025	0.025	0.350	0.400	0.050	0.875

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		CBB	CRB	CBoB	CBE	CRE	CBoE	DBB	DRB	DBoB	DBE	DRE	DBoE
f11	$R_{ m HST}$ (%)	1.1	8.8	7.3	1.9	19.8	16.5	1.1	8.8	7.3	2.3	19.7	17.1
	MANOFE	19041	12689	12554	7766	10148	11231	10522	9850	11734	4262	11006	11700
	$N_p$	40	72	48	56	56	72	56	48	64	24	56	72
	F	0.775	0.250	0.275	0.275	0.325	0.225	0.400	0.300	0.300	0.450	0.325	0.250
	$C_r$	0.050	0.200	0.625	0.350	0.325	0.525	0.100	0.350	0.200	0.325	0.625	0.475
f12	$R_{ m HST}$ (%)	56.8	9.77	77.8	69.3	81.1	85.1	57.4	78.3	77.1	70.8	81.4	85.6
	MANOFE	2755	3250	2068	2499	3714	2175	2162	2663	1464	2129	3480	1690
	$N_{p}$	64	40	40	64	40	40	80	40	40	80	48	48
	F	0.425	0.250	0.250	0.525	0.250	0.300	0.425	0.225	0.300	0.475	0.325	0.250
	$C_r$	0.550	0.700	0.800	1.000	0.950	0.975	0.675	0.725	0.975	0.950	1.000	0.950
f 13	$R_{ m HST}$ (%)	37.9	40.2	40.8	76.7	83.6	82.3	38.6	40.0	40.5	78.6	83.4	82.8
	MANOFE	3365	2307	2274	3261	2329	2349	3160	2206	2231	3142	2287	2276
	$N_{p}$	40	24	24	40	24	24	40	24	24	40	24	24
	F	1.000	1.000	1.000	1.000	1.000	1.000	1.000	1.000	1.000	1.000	1.000	1.000
	$C_{r}$	0.100	0.050	0.050	0.200	0.025	0.075	0.150	0.075	0.025	0.200	0.075	0.050

In unimodal functions ( $f1 \sim f7$ ), DDE/best/1/\* is both most reliable and efficient in all the unimodal functions except f3 (Rosenbrock), in which DDE/best/1/\* has better efficiency but worse reliability compared with DDE/rand/1/\* and DDE/BoR/1/\*. CDE/best/1/\* also has best efficiency but not best reliability in most unimodal functions. As a whole, DE/best/1/\* performs best among 3 different differential mutation strategies. DE/BoR/1/\* ranks second. And DE/rand/1/\* performs worst.

In multimodal functions ( $f8 \sim f13$ ), DE/best/1/\* performs worst. It is much less reliable than DE/rand/1/\* and DE/BoR/1/\*, and has no better efficiency. DE/rand/1/\* is the most reliable while DE/BoR/1/\* is the most efficient in general.

From the results of the unimodal and multimodal functions, it is evident that the modality of the function greatly affects the performance of DE algorithms with different differential mutation bases. Generally, DE/best/1/\* performs better than DE/BoR/1/\* and DE/rand/1/\* in unimodal functions. While in multimodal functions better diversity is needed, hence DE/rand/1/\* and DE/BoR/1/\* perform better.

## 5.2 50-Dimensional benchmark functions

The reliability and efficiency of DE when solving 50-dimensional functions and the corresponding optimal intrinsic control parameters ( $N_p$ , F,  $C_r$ ) are given in Table 2.

It is observed from Table 2 that the reliability of DE/BoR/1/\* is close to that of the best one of the three different differential mutation strategies in all the functions except f5, in which DE/best/1/\* is the most reliable and DE/BoR/1/\* ranks second. In functions f2, f8, f9, f10 and f12, the reliability of DE/BoR/1/\* and DE/rand/1/\* is much better than that of DE/best/1/\*.

As for the efficiency, in functions f1, f4 and f5, DE/BoR/1/\* is almost as efficient as DE/best/1/\* and more efficient than DE/rand/1/\*. In functions f2, f7, f10 and f12, DE/BoR/1/\* is the most efficient among the three different differential mutation strategies. In function f8, the efficiency of DE/BoR/1/\* is close to that of DE/rand/1/\* and better than that of DE/best/1/\*. In functions f9, DE/rand/1/\* is the most efficient among the three different differential mutation strategies and DE/BoR/1/\* ranks second.

These simulation results show that DE algorithms with best of random differential mutation base generally perform better than those with other two differential mutation bases in 50-dimensional functions. The advantage of DE/BoR/1/\* becomes more obvious in high dimensional functions.

The effect of modality on the performance of DE algorithms with different mutation bases has been observed again. In the uni-modal functions, though DE/best/1/\* still performs better than DE/rand/1/\*, it is not better than DE/BoR/1/\*. In the multimodal functions, the performance of DE algorithms with 3 different differential mutation bases is similar to the case of dimension 8. DE/best/1/\* performs worst. DE/rand/1/\* is the most robust while DE/BoR/1/\* is the most efficient.

## 5.3 Effects of error tolerance

In order to study the effect of error tolerance on the performance of different DE algorithms, a smaller error tolerance, i.e. 1e-5, is used to re-simulate the 8-dimensional Sphere (f1), Rosenbrock (f3), Rastrigin (f9) functions and 50-dimensional Sphere, Rastrigin functions. The limit of number of objective function evaluations under the error tolerance 1e-5 is the same as that under the error tolerance 1e-2 in the simple function f1. While it is twice as

	•												
		CBB	CRB	CBoB	CBE	CRE	CBoE	DBB	DRB	DBoB	DBE	DRE	DBoE
ĥ	$R_{ m HST}$ (%)	35.1	33.8	34.6	72.8	75.1	75.8	43.1	34.0	39.3	74.4	75.3	76.3
	MANOFE	12851	15652	10845	21871	26497	21240	10934	15065	11221	19780	23549	19476
	$N_{p}$	30	30	20	40	40	40	30	30	30	40	40	40
	F	0.525	0.350	0.450	0.525	0.400	0.400	0.525	0.350	0.400	0.575	0.350	0.425
	$C_r$	0.375	0.325	0.275	0.900	0.950	0.925	0.450	0.300	0.450	0.950	0.925	0.950
f2	$R_{ m HST}$ (%)	18.2	35.1	33.7	67.6	79.5	78.7	17.1	34.1	32.9	65.5	80.3	78.6
	MANOFE	14443	15476	14163	22733	27878	18656	17151	12883	10832	26340	23763	15299
	$N_p$	50	60	80	70	70	60	60	50	60	70	70	50
	F	0.475	0.250	0.250	0.525	0.275	0.350	0.475	0.275	0.300	0.575	0.325	0.325
	$C_r$	0.200	0.500	0.450	0.875	0.925	0.925	0.175	0.400	0.450	0.900	0.950	0.900
f4	$R_{ m HST}$ (%)	29.2	23.4	27.6	62.5	58.4	60.6	35.3	24.0	31.9	62.9	58.2	60.5
	MANOFE	21353	26460	21530	31995	38118	31187	19092	26938	19285	31149	36620	30516
	$N_p$	20	20	20	20	20	20	20	20	20	20	20	20
	F	0.575	0.450	0.450	0.550	0.475	0.450	0.575	0.450	0.450	0.575	0.500	0.500
	$C_r$	0.325	0.325	0.325	0850	0.825	0.850	0.400	0.200	0.325	0.875	0.900	0.900
f5	$R_{ m HST}$ (%)	35.2	9.1	20.1	67.4	56.3	62.6	41.6	10.4	24.2	68.0	55.5	62.8
	MANOFE	18571	29673	20452	26239	25281	24048	15734	27534	17610	24608	26119	22834
	$N_p$	30	20	20	30	20	20	30	20	20	30	20	20
	F	0.550	0.350	0.450	0.525	0.200	0.425	0.550	0.300	0.475	0.575	0.225	0.475
	$C_r$	0.350	0.025	0.275	0.875	0.175	0.825	0.525	0.025	0.400	0.925	0.175	0.900
f7	$R_{ m HST}$ (%)	25.3	29.1	28.6	68.8	73.0	72.1	29.9	29.9	32.7	69.8	73.8	73.2
	MANOFE	19377	15255	16451	27415	26343	25066	20305	18239	14835	32556	31432	27900
	$N_p$	30	20	30	30	30	30	40	30	30	40	40	40
	F	0.500	0.375	0.350	0.400	0.300	0.325	0.475	0.300	0.350	0.450	0.275	0.300
	$C_r$	0.275	0.250	0.350	0.750	0.850	0.850	0.375	0.250	0.375	0.850	0.875	0.875

		CBB	CRB	CBoB	CBE	CRE	CBoE	DBB	DRB	DBoB	DBE	DRE	DBoE
f8	$R_{ m HST}$ (%)	5.6	10.4	10.3	46.6	67.8	63.0	5.5	10.9	10.3	46.3	9.99	63.6
	MANOFE	47775	29968	29388	46548	36793	40816	41674	29811	29150	46568	39576	38253
	$N_{p}$	70	40	40	70	60	70	09	40	40	70	70	70
	F	0.375	0.325	0.375	0.350	0.125	0.100	0.400	0.325	0.375	0.400	0.075	0.125
	$C_r$	0.025	0.050	0.050	0.425	0.400	0.100	0.025	0.075	0.050	0.550	0.150	0.525
$f_{9}$	$R_{ m HST}$ (%)	0.8	2.5	2.3	14.4	37.2	35.2	0.8	2.5	2.3	15.1	37.4	35.7
	MANOFE	69015	56199	65632	63926	54091	61770	68554	56066	56647	62946	53663	61564
	$N_{p}$	40	40	50	50	40	50	40	40	40	50	40	50
	F	0.950	0.350	0.325	0.475	0.325	0.250	0.950	0.350	0.475	0.325	0.325	0.250
	$C_r$	0.025	0.025	0.025	0.075	0.100	0.150	0.025	0.025	0.025	0.025	0.050	0.125
f  10	$R_{ m HST}$ (%)	14.5	41.9	39.6	47.6	81.8	81.1	I	I	I	I	I	I
	MANOFE	22996	21863	17913	68828	28372	27756	I	I	I	I	I	I
	$N_{p}$	60	70	80	120	60	80	I	I	I	I	I	I
	F	0.475	0.225	0.275	0.300	0.275	0.300	Ι	I	I	I	I	I
	$C_r$	0.175	0.450	0.525	0.500	006.0	0.925	I	I	I	I	I	I
f12	$R_{ m HST}$ (%)	20.9	32.6	32.1	65.5	72.8	73.7	20.4	31.8	33.9	64.7	73.8	73.4
	MANOFE	14253	16416	12607	25670	26333	21629	15586	17395	11454	24564	26610	17731
	$N_p$	30	40	40	50	50	50	40	50	40	50	50	40
	F	0.525	0.300	0.350	0.550	0.350	0.400	0.550	0.300	0.325	0.525	0.375	0.375
	$C_r$	0.200	0.350	0.450	0.900	0.950	0.950	0.300	0.525	0.400	0.875	0.950	0.925

Table 2 continued

many as that under the error tolerance 1e-2 in the more difficult functions f3 and f9. The reliability and efficiency of DE algorithms in these functions and the corresponding optimal intrinsic control parameters under the error tolerance 1e-5 are shown in Table 3.

It is observed from Table 3 that in 8-dimensional sphere function, DE/best/1/\* has best reliability and efficiency among the three different differential mutation strategies under the error tolerance 1e-5. The reliability of DE/BoR/1/\* is close to that of DE/rand/1/\* but it is more efficient than DE/rand/1/\*. In 50-dimensional sphere function, DE/best/1\* is a little bit more reliable and efficient than DE/BoR/1/\*, but the difference is small. DE/rand/1/\* performs worst in this function. In general, the comparison results among the three different differential mutation strategies are similar to those under the error tolerance 1e-2. In function f 1, the advantage of DE/best/1/\* becomes a little bit more obvious under the error tolerance 1e-5.

In 8-dimensional function f3, DE/BoR/1/\* performs best among the three differential mutation strategies under the error tolerance 1e-5. This result is similar to that under the error tolerance 1e-2. The performance of DE/best/1/\* in function f3 deteriorate much under the error tolerance 1e-5. It is the least reliable and efficient among the three differential mutation strategies.

In function f 9 with dimension 8 and 50, both the reliability and efficiency of DE/BoR/1/\* are similar to those of DE/rand/1/\*. DE/best/1/\* performs worst in this function. These comparison results are consistent with those under the error tolerance 1e-2.

From these simulation results, it is observed that using the error tolerance 1e-5 in these functions generally doesn't change the comparison results of the DE algorithms with different differential mutation strategies under the error tolerance 1e-2. It appears that 1e-2 is also sufficient as the error tolerance in these functions.

# 6 Conclusions

A new differential mutation base generator for DE, namely best of random, is presented in this paper. The proposed differential mutation base generator uses the best one among several randomly chosen individuals as the differential mutation base while the other worse individuals as donors for vector differences. Thus both good diversity and fast evolution speed can be obtained in DE using the new differential mutation base. An evaluation platform for evolutionary algorithms including DE is built. With full consideration of the demand from practical applications, differential evolution algorithms, intrinsic and non-intrinsic control parameters, and problem features are treated as a system. A comprehensive comparative study over a set of benchmark functions has been conducted at system level instead of parameter level. The results show that a better balance of reliability and efficiency can be obtained in DE implementing the new generator of differential mutation base, especially in high dimensional functions.

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		CBB	CRB	CBoB	CBE	CRE	CBoE	DBB	DRB	DBoB	DBE	DRE	DBoE
<i>f</i> 1 8D	$R_{ m HST}  (\%)$ MANOFE $N_p$ F	63.6 <b>2319</b> 40 0.450 0.675	<b>63.7</b> 4454 32 0.325 0.625	62.1 3032 32 0.375 0.700	74.8 <b>2431</b> 40 0.525 0.975	<b>75.5</b> 5608 40 0.250 0.800	73.4 3638 40 0.350 0.925	<b>80.6</b> 1110 48 0.300	62.5 3865 32 0.350 0.625	61.4 2383 32 0.425 0.825	<b>85.6</b> 1116 48 0.300 1.000	74.7 4725 40 0.375 0.975	74.0 2836 40 0.400 0.975
f1 50D	$R_{\rm HST}^{C}$ (%) MANOFE $N_p$ F	<b>31.4</b> <b>15111</b> 20 0.550	27.4 24337 30 0.400	27.7 27.7 19244 30 0.375 0.375	69.3 25580 30 0.550	67.5 67.5 30339 30 0.375 0.375	67.4 67.4 30660 40 0.450 0.950	<b>38.5</b> <b>15055</b> 30 0.550 0.475	27.5 18806 20 0.450	29.0 15265 20 0.475 0.275	70.1 24591 30 0.575 0.925	67.6 29539 30 0.425 0.900	67.8 2 <b>4581</b> 30 0.425
f 3 8D	$R_{\rm HST}^{C}$ (%) MANOFE $N_p$ F	0.9 72301 80 0.950	222.9 222.9 44634 64 0.575 1 000	<b>23.9</b> <b>34107</b> 72 0.575 1.000	2.6 2.6 107141 160 0.850 0.850	20.5 20.5 31042 40 0.625 1.000	<b>20.8</b> <b>26514</b> 48 0.600 0.975	0.9 135240 120 0.950	24.4 35228 56 0.550	<b>25.9</b> <b>24882</b> 64 0.550	2.7 2.7 105116 0.850 0.850	21.5 27864 40 0.625	<b>22.6</b> <b>22239</b> 48 0.625
f9 8D	$\substack{Cr}{R_{\rm HST}}$ (%) MANOFE $N_p$ $C_r$	16.1 16.1 12042 80 0.325 0.100	<b>34.5</b> <b>34.5</b> 10289 48 0.200 0.075	33.0 33.0 56 0.200 0.200	31.5 31.5 80 0.250 0.050	<b>62.5</b> 11691 56 0.200 0.025	61.9 61.9 56 0.200 0.150	15.9 11284 80 0.250 0.050	<b>35.3</b> 10243 48 0.225 0.050	32.8 32.8 56 0.200 0.025	30.8 30.8 80 0.275 0.050	62.9 9862 48 0.200 0.175	61.3 61.3 9864 56 0.175 0.100
f9 50D	$\begin{array}{c} R_{\rm HST} \left( \% \right) \\ {\rm MANOFE} \\ N_p \\ F \\ C_r \end{array}$	1.4 80297 50 0.475 0.025	<b>5.1</b> 81368 50 0.300 0.025	4.4 <b>78644</b> 50 0.300 0.025	27.5 116079 80 0.325 0.150	<b>57.3</b> 77756 50 0.250 0.050	55.9 86532 60 0.225 0.225	1.4 94612 60 0.475 0.025	<b>5.1</b> 80945 50 0.300 0.025	4.7 <b>79734</b> 50 0.325 0.025	27.0 115279 80 0.325 0.225	<i>57.9</i> 86865 60 0.200 0.200	56.4 <b>86606</b> 60 0.225 0.100

 Table 3
 Performance of differential evolution under the error tolerance 1e-5

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## Appendix

Benchmark functions in test bed

# f1 Sphere function

The sphere function is continuous, differentiable, separable, scalable, uni-modal, and symmetric. It is the benchmark for studying problem features.

The sphere function reads

$$f(\mathbf{x}) = \sum_{i=1}^{N} x_i^2 \tag{4}$$

Its minimum is  $f(\mathbf{x}^*) = 0$  at  $\mathbf{x}^* = \mathbf{0}$ . Standard search space is  $\mathbf{x} \in [-500, 500]^N$ 

# f 2 Step function 2

The step function 2 is discontinuous, non-differentiable, separable, scalable, uni-modal, and symmetric. It is designed for studying the effect of continuity.

The step function 2 reads

$$f(\mathbf{x}) = \sum_{i=1}^{N} (\lfloor x_i + 0.5 \rfloor)^2$$
(5)

Its minimum is  $f(\mathbf{x}^*) = 0$  at  $\mathbf{x}^* = \mathbf{0}$ . Standard search space is  $\mathbf{x} \in [-500, 500]^N$ 

#### f 3 Rosenbrock function

The Rosenbrock function is continuous, differentiable, non-separable, scalable, uni-modal, and asymmetric. It reads

$$f(\mathbf{x}) = \sum_{i=1}^{N-1} \left[ 100 \left( x_{i+1} - x_i^2 \right)^2 + (x_i - 1)^2 \right]$$
(6)

Its minimum is  $f(\mathbf{x}^*) = 0$  at  $\mathbf{x}^* = \mathbf{1}$ . Recommended search space is  $\mathbf{x} \in [-100, 100]^N$ 

#### f 4 Hyper-ellipsoid function

The hyper-ellipsoid function is continuous, differentiable, separable, scalable, uni-modal, and asymmetric. It is usually implemented to study the effect of symmetry.

The hyper-ellipsoid function reads

$$f(\mathbf{x}) = \sum_{i=1}^{N} 2^{i-1} x_i^2 \tag{7}$$

Its minimum is  $f(\mathbf{x}^*) = 0$  at  $\mathbf{x}^* = \mathbf{0}$ . Standard search space is  $\mathbf{x} \in [-500, 500]^N$ 

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#### f 5 Qing function

The Qing function is originally designed to study the effect of non-uniqueness. It is continuous, differentiable, separable, scalable, uni-modal, and asymmetric.

The Qing function reads

$$f(\mathbf{x}) = \sum_{i=1}^{N} (x_i^2 - i)^2$$
(8)

Its minima are  $f(\mathbf{x}^*) = 0$  at  $x_i^* = \pm \sqrt{i}$ . Standard search space is  $\mathbf{x} \in [-500, 500]^N$ 

#### f 6 Schwefel function 1.2

The Schwefel function 1.2 is a rotated sphere function. It is continuous, differentiable, nonseparable, scalable, uni-modal, and asymmetric. The effect of symmetry and rotation can be investigated through it.

The Schwefel function 1.2 reads

$$f(\mathbf{x}) = \sum_{i=1}^{N} \left( \sum_{j=1}^{i} x_j \right)^2 \tag{9}$$

Its minimum is  $f(\mathbf{x}^* = \mathbf{0}) = 0$ . Standard search space is  $\mathbf{x} \in [-500, 500]^N$ 

## f7 Schwefel function 2.22

The Schwefel function 2.22 is continuous, non-differentiable, non-separable, scalable, unimodal, and symmetric. It can be implemented to examine the effect of convexity because its landscape is non-convex.

The Schwefel function 2.22 reads

$$f(\mathbf{x}) = \sum_{i=1}^{N} |x_i| + \prod_{i=1}^{N} |x_i|$$
(10)

Its minimum is  $f(\mathbf{x}^*) = 0$  at  $\mathbf{x}^* = \mathbf{0}$ . Standard search space is  $\mathbf{x} \in [-500, 500]^N$ 

#### f 8 Schwefel function 2.26

The Schwefel function 2.26 is continuous, non-differentiable, separable, scalable, multimodal, and symmetric. The effect of differentiability and modality can be examined through it.

The Schwefel function 2.26 reads

$$f(\mathbf{x}) = -\frac{1}{N} \sum_{i=1}^{N} x_i \sin \sqrt{|x_i|}$$
(11)

Its minimum is  $f(\mathbf{x}^*) = -418.983$  at  $\mathbf{x}^* = 420.968746$ . Recommended search space is  $\mathbf{x} \in [-500, 500]^N$ 

#### f9 Rastrigin function

The Rastrigin function is continuous, differentiable, separable, scalable, multi-modal, and symmetric. It is very suitable for investigating the effect of modality.

The Rastrigin function reads

$$f(\mathbf{x}) = \sum_{i=1}^{N} \left[ x_i^2 - 10\cos\left(2\pi x_i\right) + 10 \right]$$
(12)

Its minimum is  $f(\mathbf{x}^*) = 0$  at  $\mathbf{x}^* = \mathbf{0}$ . Standard search space is  $\mathbf{x} \in [-100, 100]^N$ 

## f 10 Griewank function

The Griewank function is continuous, differentiable, non-separable, scalable, multi-modal, and asymmetric. It is also difficult to optimize. Accordingly, it is a good candidate for study the effect of modality, separability, and symmetry.

The Griewank function reads

$$f(\mathbf{x}) = \frac{1}{4000} \sum_{i=1}^{N} x_i^2 - \prod_{i=1}^{N} \cos\left(x_i/\sqrt{i}\right) + 1$$
(13)

Its minimum is  $f(\mathbf{x}^*) = 0$  at  $\mathbf{x}^* = \mathbf{0}$ . Recommended search space is  $\mathbf{x} \in [-500, 500]^N$ 

#### f 11 Whitley function

The Whitley function is a composite function. It is continuous, differentiable, non-separable, scalable, multi-modal, and symmetric. It is very difficult to optimize. It is also used to study the effect of modality and separability.

The Whitley function reads

$$f(\mathbf{x}) = \sum_{i=1}^{N} \sum_{j=1}^{N} \left[ \frac{y_{ji}^2}{4000} - \cos y_{ji} + 1 \right]$$
(14)

where  $y_{ji} = 100 \left(x_i - x_j^2\right)^2 + \left(x_j - 1\right)^2$ .

Its minimum is  $f(\mathbf{x}^*) = 0$  at  $\mathbf{x}^* = \mathbf{1}$ . Recommended search space is  $\mathbf{x} \in [-500, 500]^N$ 

# f 12 Ackley function

The Ackley function is continuous, differentiable, non-separable, scalable, multi-modal, and symmetric. It is therefore suitable for investigating the effect of modality and separability.

The Ackley function reads

$$f(\mathbf{x}) = 20 - 20e^{-0.2\sqrt{\frac{1}{N}\sum_{i=1}^{N}x_i^2}} + e - e^{\frac{1}{N}\sum_{i=1}^{N}\cos(2\pi x_i)}$$
(15)

Its minimum is  $f(\mathbf{x}^*) = 0$  at  $\mathbf{x}^* = \mathbf{0}$ , standard search space is  $\mathbf{x} \in [-30, 30]^N$ 

#### f 13 Weierstrass function

The Weierstrass function is continuous, differentiable, separable, scalable, multi-modal, and symmetric. It reads

$$f(\mathbf{x}) = \sum_{i=1}^{N} \sum_{k=0}^{K_{\text{max}}} a^k \cos\left[2\pi b^k \left(x_i + 0.5\right)\right] - N \sum_{k=0}^{K_{\text{max}}} a^k \cos\left(\pi b^k\right)$$
$$a = 1, b = 1, K_{\text{max}} = 10$$
(16)

Its minimum is  $f(\mathbf{x}^* = \mathbf{0}) = 0$ . Recommended search space is  $\mathbf{x} \in [-500, 500]^N$ 

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