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Pseudorandomness in Central Force Optimization

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Research Article

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Abstract

Central Force Optimization is a deterministic metaheuristic for an evolutionary algorithm that searches a decision space by flying probes whose trajectories are computed using a gravitational metaphor. CFO benefits from the inclusion of a pseudorandom component (a numerical sequence that is precisely known by specification or calculation but otherwise arbitrary). The essential requirement is that the sequence is uncorrelated with the decision space topology, so that its effect is to pseudorandomly distribute probes throughout the landscape. While this process may appear to be similar to the randomness in an inherently stochastic algorithm, it is in fact fundamentally different because CFO remains deterministic at every step. Three pseudorandom methods are discussed (initial probe distribution, repositioning factor, and decision space adaptation). A sample problem is presented in detail and summary data included for a 23-function benchmark suite. CFO's performance is quite good compared to other highly developed, state-of-the-art algorithms.

Keywords: Central force optimization; CFO; optimization; metaheuristic; evolutionary algorithm; pseudorandomness; decision space exploration and exploitation.

1 Introduction

This note examines the role of pseudorandomness in Central Force Optimization. CFO is a deterministic Nature-inspired search and optimization metaheuristic for an evolutionary algorithm (EA) based on gravitational kinematics [1-3]. CFO is similar to gradient-based optimization methods as discussed in detail in [4]. Proofs of convergence for CFO and an extended version have been developed [5,6], and the algorithm has been implemented on a GPU using various topologies [7-9]. The algorithm has been successfully applied to a variety of problems, among them: training neural networks [10]; power grid reliability assessment [11]; drinking water distribution networks [12]; solving nonlinear circuits [13]; array synthesis [14,15]; microstrip patch antenna design [16]; multiband slotted bowtie design [17]; rectangular microstrip patch design [18]; microwave broadband absorber design [19]; antenna optimization generally [20]; notched ultra wideband E-shape antenna design [21]; and increasing impedance bandwidth

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[22,23].

CFO analogizes Newton's mathematically precise laws of motion and gravity, so that its underlying equations are equally precise. The algorithm locates the global maxima of an objective function defined on a decision space Ω with unknown topology (landscape). CFO searches Ω by flying a group of probes whose trajectories are computed from two deterministic equations of motion at a series of discrete time steps (iterations). Details of the CFO metaheuristic are in the Appendix. CFO is fundamentally different from Nature-inspired EAs whose underlying equations are inherently stochastic. Particle Swarm Optimization (PSO) and Ant Colony Optimization (ACO) are examples. Their equations are formulated in terms of true random variables, and removing randomness causes these algorithms to fail completely. By contrast, CFO's equations are inherently deterministic. Every CFO run with the same setup returns precisely the same values step-by-step throughout the entire run. Nevertheless, effective implementations benefit from a pseudorandom component that enters the algorithm indirectly, not through its basic equations. Although pseudorandomness is not required in CFO, numerical experiments show that it is an important feature in effective implementations.

A pseudorandom variable is defined here as one whose value is precisely known but arbitrarily assigned. The value can be specified in advance (for example, an arbitrary sequence of numbers) or it can be calculated in a prescribed manner. The variable's randomness derives from the fact that its value is arbitrary and uncorrelated with Ω 's topology, not that it is uncertain in the sense of a true random variable. A true random variable's value is calculated from a probability distribution with successive calculations yielding different values that cannot be known in advance. This type of randomness is fundamentally different from CFO's pseudorandomness. Even when CFO includes pseudorandomness, it remains deterministic, always yielding the same result for runs with the same setup. Once a pseudorandom variable is specified, either explicitly or by calculation, its value is known with absolute precision, so that CFO's trajectory calculations are deterministic even in the presence of pseudorandomness. The original CFO implementation did not include a pseudorandom component [1], thereby limiting its ability to explore Ω ; but this limitation can be mitigated to some extent by introducing a measure of pseudorandomness. There are many ways this can be accomplished; three simple methods are described here.

2 CFO Methodology

This paper discusses a CFO implementation with pseudorandomness injected in the following ways: (1) the initial probe distribution (IPD); (2) the repositioning factor; and (3) changing the decision space boundaries. The algorithm is referred to as CFO-PR. Every CFO run starts with a user-specified IPD (total number of probes and their locations in Ω at the beginning of the run, step 0). An arbitrary, variable initial probe distribution is a convenient way to inject pseudorandomness, the effect of which is to provide better sampling of Ω 's topology than a static distribution. Each initial probe distribution in a set of distributions provides different information about $Ω$'s landscape. As the results below show, certain ones perform much better than others.

The second method of injecting pseudorandomness is the use of a step-by-step variable repositioning factor $\Delta F_{rep} \leq F_{rep} \leq 1$ where ΔF_{rep} is the step increment. Repositioning refers to the process of retrieving a probe that has flown outside the decision space (discussed in detail in the Appendix). A variable F_{rep} has the effect of pseudorandomly distributing probes throughout Ω , which provides better sampling of the decision space landscape as a run progresses. The third way pseudorandomness is injected is by shrinking the decision space around the best probe's location. This process coupled with variable F_{rep} redistributes probes in the smaller Ω in an arbitrary but precise, hence pseudorandom, manner. The effect is to speed CFO's convergence, but at the risk of premature convergence (on an empirical basis, this issue does not appear to be significant using the procedures described below). Pseudocode for CFO-PR appears in Fig. 1. The inner time step loop (*j* loop) is common to all CFO implementations, but the two outer loops inject initial probe pseudorandomness. The γ loop controls where initial probes are deployed, and the N_p/N_d loop determines their number. These two parameters define the initial probe distribution, and the two loops together create a variable, pseudorandom distribution (see Appendix for notation and equations). The variable F_{rep} procedure appears in step (f) of the pseudocode. And the pseudorandom decision space adaptation is in step (g). Ω 's boundaries shrink around the then best probe position vector every $20th$ step as discussed in detail below. A two-dimensional example is used to illustrate these techniques because it provides a concrete visualization of the different methods. In the actual CFO-PR implementation, of course, these techniques are generalized to the N_d -dimensional case.

2.1 Initial Probes

The manner in which initial probes are deployed using γ is shown schematically in Fig. 2, which provides a 2-dimensional (2D) schematic representation of a variable initial probe distribution comprising an orthogonal array of $\binom{1}{r}$ $\frac{N_p}{N_d}$ probes per axis deployed uniformly on probe lines parallel to the coordinate axes that intersect at a point along Ω 's principal diagonal. N_d is Ω 's dimensionality (here $N_d = 2$), and $x_i^{\min} \le x_i \le x_i^{\max}$, $1 \le i \le N_d$ define Ω 's domain (decision space).

For $N_p/N_d = 2$ to $(N_p/N_d)_{MAX}$ step 2: For $\gamma = \gamma_{start}$ to γ_{stop} by $\Delta \gamma$: (a.1) Compute initial probe distribution. (a.2) Compute initial fitness matrix. (a.3) Assign initial probe accelerations. (a.4) Set initial F_{rep} . For $j = 0$ to N_t (or earlier termination criterion): (b) Compute probe position vectors, \vec{R}_{i}^{p} , $1 \le p \le N_{p}$ [eq.(2), Appendix]. (c) Retrieve errant probes $(1 \le p \le N_p)$: If $\vec{R}_i^p \cdot \hat{e}_i < x_i^{\min}$: $\vec{R}^{\,p}_j \cdot \hat{e}^{\vphantom{\dagger}}_i = x^{\rm min}_i + F_{\it rep} (\vec{R}^{\,p}_{j\!-\!1} \cdot \hat{e}^{\vphantom{\dagger}}_i - x^{\rm min}_i)$ If $\vec{R}_i^p \cdot \hat{e}_i > x_i^{\text{max}}$: $\vec{R}_{i}^{p} \cdot \hat{e}_{i} = x_{i}^{\max} - F_{rep}(x_{i}^{\max} - \vec{R}_{i-1}^{p} \cdot \hat{e}_{i})$ (d) Compute fitness matrix for current probe distribution. (e) Compute accelerations using current probe distribution and fitnesses [eq. (1) , Appendix] (f) Increment F_{rep} by ΔF_{rep} : If $F_{rep} > 1$: $F_{rep} = \Delta F_{rep}$. (g) If $j \text{ MOD } 20 = 0$: Shrink Ω around \vec{R}_{host} . Next j Next γ Next N_p/N_d

For illustrative purposes, Fig. 2 shows nine probes on each probe line. The lines, which are parallel to the x_1 and x_2 axes, intersect at a point on Ω 's principal diagonal marked by position

vector
$$
\vec{D} = \vec{X}_{min} + \gamma (\vec{X}_{max} - \vec{X}_{min})
$$
, where $\vec{X}_{min} = \sum_{i=1}^{N_d} x_i^{min} \hat{e}_i$ and $\vec{X}_{max} = \sum_{i=1}^{N_d} x_i^{max} \hat{e}_i$ are

the diagonal's endpoint vectors. Parameter $0 \le \gamma \le 1$ specifies where along the diagonal the orthogonal probe array is placed by locating the probe lines' intersection point.

While Fig. 2 shows an equal number of probes on each line, a different number of probes per axis can be used instead. For example, if equal probe spacing were desired in a decision space with unequal boundaries, or if overlapping probes were to be excluded in a symmetrical space, then unequal numbers would be used. Unequal numbers also might be appropriate if *a priori* knowledge of Ω's landscape, however obtained, suggests denser sampling in one area. The initial

probe distribution in Fig. 2 with variable $\sqrt[p]{v}$ $\frac{N_p}{N_d}$ was used for the CFO-PR runs reported here,

but any number of other variable initial probe distributions could be used instead. The key idea is that the initial probe distribution must be pseudorandom, that is, arbitrary and therefore uncorrelated with the decision space landscape.

2.2 Repositioning Factor

A variable value for F_{ren} also adds pseudorandomness. F_{ren} starts at some arbitrary initial value that is incremented at each iteration by an arbitrary amount ΔF_{rep} so that $\Delta F_{rep} \leq F_{rep} \leq 1$. This errant probe retrieval scheme is an example of an arbitrary sequence of calculated numbers deterministically assigned to a CFO parameter. CFO's ability to search the decision space depends on where errant probes are reinserted in Ω , and this process is pseudorandom in nature because F_{ren} is deterministic but arbitrary and uncorrelated with Ω's topology. By placing errant probes pseudorandomly throughout the decision space, more information is developed about its topology as the run progresses. The scheme used here was empirically determined, and appears to work well across a wide range of objective functions. But, of course, there are many other procedures for setting F_{ren} 's value, some no doubt better than others.

Fig. 2. Variable 2-D initial probe distribution used for CFO runs reported here

2.3 Decision Space Adaptation

CFO-PR also includes adaptive reconfiguration of the decision space in order to improve convergence speed. This feature also is pseudorandom in nature because the way Ω 's boundaries are changed is arbitrary and uncorrelated with the landscape. Fig. 3 illustrates in 2D how Ω 's size is adaptively reduced in this case every $20th$ step around the probe's location with the then best Is adaptively reduced in this case every 20 step around the probe s location with the their best
fitness throughout the run up to the current iteration, \vec{R}_{best} . Ω's boundary coordinates are reduced by one-half the distance from the best probe's position to the each boundary on a coordinate-by coordinate basis, that is, $x_i^{\prime\prime\prime\prime\prime} = x_i^{\prime\prime\prime\prime\prime} + \frac{-b_{est} - y_i}{2}$ and $x_i^{\prime\prime\prime\prime\prime\prime} = x_i^{\prime\prime\prime\prime\prime\prime} - \frac{-b_{est} - y_i}{2}$ $\mathbf{x}_i^{\min} = x_i^{\min} + \frac{\vec{R}_{best} \cdot \hat{e}_i - x_i^{\min}}{2}$ and $x_i^{\max} = x_i^{\max} - \frac{x_i^{\max}}{2}$ $x_i^{\prime \min} = x_i^{\min} + \frac{R_{best} \cdot \hat{e}_i - x_i^{\min}}{2}$ and $x_i^{\prime \max} = x_i^{\max} - \frac{x_i^{\max} - R_{best} \cdot \hat{e}_i}{2}$ $x_i^{\min} = x_i^{\min} + \frac{R_{best} \cdot \hat{e}_i - x_i^{\min}}{2}$ and $x_i^{\max} = x_i^{\max} - \frac{x_i^{\max} - R_{best} \cdot \hat{e}_i}{2}$ but s position to the each boundary on and $x_i^{\text{max}} = x_i^{\text{max}} - \frac{y_i - y_{\text{best}} - y_i}{2}$, $\mathcal{L}^{\max}_i = x^{\max}_i - \frac{x^{\max}_i - \bar{R}_{best} \cdot \hat{\bm{e}}_i}{2},$ $x_i'^{\, \rm max} = x_i^{\, \rm max} - \frac{x_i^{\, \rm max} - R_{best} \cdot \hat{\pmb{e}}_i}{2} \, ,$ \vec{z} \vec{z} , where the primed coordinate is the new decision space boundary, and the dot denotes vector inner where the primed coordinate is the new decision space boundary, and the dot denotes vector limer product. For clarity, Fig. 3 shows \vec{R}_{best} as fixed, whereas generally it varies throughout a run. Changing Ω 's boundary every twenty steps instead of some other interval was chosen arbitrarily (another, probably better approach, might be a reactive adaptation based on performance measures such as convergence speed or fitness saturation).

3 Results and Discussion

3.1 A Sample Problem

The effectiveness of injecting pseudorandomness into CFO-PR will be illustrated with the 2D Goldstein-Price function (GP) plotted in Fig. 4. GP is defined as

$$
f(x_1, x_2) = -[1 + (x_1 + x_2 + 1)^2 \cdot (19 - 14x_1 + 3x_1^2 - 14x_2 + 6x_1x_2 + 3x_2^2)]
$$

× [30 + (2x₁ - 3x₂)² · (18 - 32x₁ + 12x₁² + 48x₂ - 36x₁x₂ + 27x₂²)]

where Ω : $-100 \le x_1, x_2 \le 100$ (note that in most published reports Ω is much smaller, *viz.*, $-2 \le x_1, x_2 \le 2$). GP's global maximum is -3 at $(0,-1)$. This function is multimodal with few local maxima, and it varies over nearly nineteen orders of magnitude as shown in Fig. 4.

Fig. 4. Goldstein-price (GP) function

The following parameter values were used for all runs reported in this note: $\alpha = 2$, $\beta = 2$, $G = 2$, $\Delta t = 1$, initial acceleration of zero, initial $F_{rep} = 0.5$, $\Delta F_{rep} = 0.05$, $\gamma_{start} = 0$, $\gamma_{stop} = 1$ with $\Delta \gamma = 0.1$ (eleven runs). For GP, $N_p/N_d = 4$ to 14 by 2, with different

,

ranges of this parameter for the other test functions as described below. In all cases, a run was terminated early if the average best fitness over 50 steps (including the current step) and the current best fitness differed by less than 10^{-6} . The pseudorandom IPD's computed using the procedure illustrated in Fig. 2 for GP are plotted in Fig. 5. A further illustration of the Probe Line IPD concept is shown in 3D in Fig. 6.

Table 1 shows a summary of the results for the GP function. A total of sixty-six optimization runs were made in six groups of eleven runs each (data for Run #0 are starting values). The column headings are for the most part self-explanatory. Each run began with $N_t = 500$, but, as the *#Steps* column shows, in no case were 500 iterations used because every run terminated early. N_{eval} is the number of function evaluations performed for the shortened run, and the total number of evaluations over all runs appears at the bottom of this column. The F_{ren} column lists F_{ren} 's value at the end of the run, and the V denotes that *Frep* was variable as discussed above. *Fitness* tabulates the best fitness returned during the run. The *Initial Probes* column shows the type of IPD, in this case probes uniformly spaced along probe lines parallel to Ω 's axes (notated I-AXIS) as described above and shown in Fig. 5.

The best fitness ranged from a low of $-236.6643...$ in run 12 to the global maximum of -3 at $(0,-1)$, which was returned in run #54 (best results highlighted in blue). Parameters for run #54 were $N_p/N_d = 12$, $N_p = 24$, and $\gamma = 0.9$. The total number of function evaluations over all runs was 180,472 while N_{eval} for the best run was 1,464 (60 iterations). Fig. 7 plots the evolution of GP's best fitness, which in only two steps increases from which in only two steps increases from $-2.992268247672x10^{-12}$ to GP's actual global maximum of -3 . This seems to be quite remarkable in view of the initial probe distribution for $\gamma = 0.9$ (Fig. 5) in which all probes are far removed from the maximum's location at $(0,-1)$. As the data in Table 1 clearly show, some sets of parameters are much better than others. Without pseudorandom initial probes, a single run would be made with, in this example, only 13.6% probability of locating the global maximum with a fractional accuracy of 0.03% (9 of 66 runs). This statistic highlights the importance of pseudorandomness in CFO.

Fig. 8 plots CFO's D_{avg} curve for GP. D_{avg} is the normalized average distance between the probe with the best fitness and all other probes at each time step, *viz.*, $\frac{1}{\cdot (N_p-1)}\sum_{p=1}^N \sqrt{\sum_{i=1}^n (x_i^{p,j}-x_i^{p^*,j})^2}$ where p^* is the number of the probe with the $=\frac{1}{(x_i-x_i)^2} \sum_{i=1}^n \left(\sum_{i=1}^n (x_i^{p,j} - x_i^{p^*,j})^2 \right)$ where p $\sum_{p=1}^{N_p} \sqrt{\sum_{i=1}^{N_d} (x_i^{p,j} - x_i^{p^*,j})^2}$ where p^* is the nu p **1)** $p=1$ **i** $i=1$ $\sum_{p \text{avg}} \frac{1}{L \cdot (N_p - 1)} \sum_{p=1}^{N} \sqrt{\sum_{i=1}^{N} (x_i^{p,j} - x_i^{p,j})^2}$ where p^* is the number of the $D_{\text{ave}} = \frac{1}{|X| \left(\frac{1}{2}X - 1\right)} \sum_{i=1}^{n} \left(\sum_{i=1}^{p} (x_i^{p,j} - x_i^{p^*,j})^2 \right)$ where p^* is the nu $i=1$ $i=1$ $(x_i^{p,j} - x_i^{p^*,j})^2$ where p^* is the number (N_p-1) $\sum_{p=1}^{\infty}$ $\sum_{i=1}^{\infty}$ $\sum_{i=1}^{\infty}$ $\frac{1}{\sigma^2} \sum_{i=1}^{N_p} \left(\sum_{i=1}^{N_d} (x_i^{p,j} - x_i^{p^*,j})^2 \right)$ where p^* is the number of the probe with the

best fitness, and $L = \sqrt{\sum_{i=1}^{N_d} (x_i^{\text{max}} - x_i^{\text{min}})^2}$ is the length of Ω 's principal diagonal (s $L = \sqrt{\sum_{i=1} (x_i^{\text{max}} - x_i^{\text{min}})^2}$ is the length of Ω's principal diagonal (see Appendix

for definitions). D_{avg} decreases monotonically through step 10 to 0.0496977, then increases very quickly to a peak of 0.4767301 at step 11, followed by another quasi-monotonic decrease through step 29 to 0.0274355. This cycle repeats through step 48 where D_{avg} is 0.0169657, followed by a jump to 0.1194779 at step 49. After a slight dip through step 53, *Davg* flattens out around a value of 0.11... The quasi-oscillatory behavior in D_{ave} usually correlates with local trapping, which in this case happens to be at the global maximum. Oscillation in D_{ave} may be a similar phenomenon to oscillation seen in ΔV curves for gravitationally trapped Near Earth Objects (NEOs), where ΔV is the velocity change needed to avoid earth impact, which suggests that NEO theory may hold the key to analytical mitigation or elimination of local trapping at local maxima and possibly another proof of convergence for CFO.

Because CFO-PR converges so quickly on GP's global maximum, the number of the probe with the best fitness is constant after step #1 as seen in the best probe plot in Fig. 9. The best probe number $(\#14)$ is the same for steps 0 and 1, but it switches to probe $\#2$ at step 2. Neither the number of the best probe nor the fitness change after step 2. Of course, for most functions the best probe number varies throughout a run, often quite erratically.

Figs. 10 and 11, respectively, plot the probe trajectories for the probes with the best ten fitnesses ordered by fitness and for the first sixteen individual probes ordered by probe number (number of trajectories plotted chosen as a matter of convenience). Both plots are very chaotic with no obvious sign of regularity in how probes gravitate to the global maximum. Nevertheless, there is some measure of regularity as reflected in the *Davg* curve because its appearance is not nearly as chaotic as the trajectory plots. In fact, in many cases D_{avg} exhibits a mathematically precise oscillation even when the probe trajectories look like Figs. 10 and 11 (see in particular [1]).

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Fig. 5. GP best run initial probe distributions

Fig. 6. Example of a 3D probe line IPD

3.2 A Benchmark Suite

CFO was tested against the same twenty-three function 2-30D benchmark suite used to evaluate Group Search Optimizer (GSO) [24]. GSO has gained some notoriety as an effective state-of-the art stochastic algorithm [25]. In [24] GSO is compared to two other algorithms, PSO and GA, using the benchmark suite described in detail in [26] (functions, decision spaces, and characteristics). PSO is a stochastic Particle Swarm Optimization algorithm implemented using PSOt, a MATLAB-based toolbox that includes standard and variant PSO algorithms. The standard PSO algorithm was used with recommended default parameters: population, 50; acceleration factors, 2.0; inertia weight decaying from 0.9 to 0.4. GA is a stochastic Genetic Algorithm implemented using the GAOT toolbox (Genetic Algorithm Optimization Toolbox). GA also employed recommended default parameter values with a fixed population size (50), uniform mutation, heuristic crossover, and normalized geometric ranking for selection. Thus, while the results reported here compare CFO and GSO directly, they also compare CFO to PSO and GA indirectly.

Table 1 summarizes results using the same function numbering as [24]. f_{max} is the known global *maximum* (note that the negative of each benchmark in [24] is used here because, unlike the other algorithms, CFO locates maxima, not minima). $\langle \cdot \rangle$ denotes average value. Because GSO, PSO and GA are inherently stochastic, their performance must be described statistically. Statistical data

in Table 2 for those algorithms are reproduced from [24], while Table 3 shows the number of function evaluations per GSO/GA/PSO run. The tabulated fitnesses are average values over 1,000 runs for the first thirteen benchmarks, and over 50 for the others. In marked contrast, CFO's results are repeatable over runs with the same parameters because the algorithm is completely deterministic, even when pseudorandomness is included. CFO therefore never requires a statistical description of its performance, which is a major advantage and a significant departure from the far more common stochastic approaches.

Test	N_d	$J_{\rm max}$	<best fitness="">/ other algorithm</best>	CFO-PR -					- - - - - - CFO - - - - - -
function				Best fitness	γ_{best}	Best N_p/N_d	N_{eval}		Best fitness with fixed
							Best Run	Total	DS boundary & fixed $N_p/N_d = 10$, $F_{rep} = 0.5$.
			Unimodal functions (other algorithms: average of 1000 runs)						
	30		$-3.6927x10^{-37}$ / PSO	$-4.8438x10^{-4}$	0.1		20,640	507,060	0
	30		$-2.9168x10^{-24}$ / PSO	$-4x10^{-8}$	0.5	2	5,040	716,400	
	30		$-1.1979x10^{-3}$ / PSO	$-6x10^{-8}$	0.5		10,260	1,534,260	
	30		$-0.1078 / GSO$	-4.2×10^{-7}	0.5		5,160	332,340	
	$30\,$		$-37.3582 / PSO$	-1.09289×10^{-3}	0.9		34,560	845,640	-29
	30		$-1.6000x10^{-2}$ / GSO		1.0		10,980	350,280	
	30		$-9.9024x10^{-3}$ / PSO	$-4.249x10^{-5}$	0.1		60,120	1,983,960	-0.002354
			Multimodal functions, many local maxima (other algorithms: avg 1000 runs)						
T ₈	30	12,569.5	12,569.4882 / GSO	12,569.4866	0.5		12,720	448,800	12,536.3016
	30		$-0.6509 / GA$	$-2.05x10^{-6}$	0.7		16,440	680,640	
f_{10}	$30\,$		$-2.6548x10^{-5}$ / GSO	$-1.5x10^{-7}$	0.5		5,100	904,980	
f_{11}	30		$-3.0792x10^{-2}$ / GSO	-9.97293×10^{-2}	0.1		42,660	489,060	-1.4141
f_{12}	30		$-2.7648x10^{-11}$ / GSO	-2.067×10^{-5}	0.5		3,660	341,400	-1.7671
f_{13}	30		-4.6948x10 ⁻⁵ / GSO	$-3.2853x10^{-3}$	0.6	6	16,920	679,620	-5.8
			Multimodal functions, few local maxima (other algorithms: avg 50 runs)						
f_{14}			$-0.9980 / GSO$	-0.9980	0.2	12	5,952	141,076	-1.4064
f_{15}		$-3.075x10^{-5}$	$-3.7713x10^{-4}$ / GSO	$-4.889x10^{-4}$	0	12	3,360	304,664	-6.4685×10^{-3}
f_{16}		1.0316285	1.031628 / GSO	1.031626	0.4	12	6,288	124,340	0.8535
f_{17}		-0.398	$-0.3979 / GSO$	-0.3979	0	8	1,872	108,340	-0.4001
f_{18}		-3	$-3/GSO$	-3	0.9	12	1,464	180,472	-3

Table 2. CFO-PR comparative results for 23 benchmark functions (GSO/GA/PSO data from [24])

 $(N_d =$ *Function Dimension;* f_{max} = *Known Global Maximum;* \lt > *Denotes Average Value)*

Function	GSO/GA/PSO	Function	GSO/GA/PSO
Ĵі	150,000	f_{I3}	150,000
f_2	150,000	f_{l4}	7,500
f_3	250,000	f_{l5}	250,000
J4	150,000	f_{l6}	1,250
f_5	150,000	f_{17}	5,000
f ₆	150,000	f_{l8}	10,000
J7	150,000	f_{I9}	4,000
J8	150,000	f_{20}	7,500
f_9	250,000	f_{2I}	10,000
f_{l0}	150,000	f_{22}	10,000
f_{II}	150,000	f_{23}	10,000
f_{12}	150,000		

Table 3. Number of GSO/GA/PSO Function Evaluations Per Run (from [24]).

Two columns of CFO-PR and CFO data are presented in Table 2. The CFO-PR data correspond to the best fitness returned by the single best run in the set of runs with variable N_p/N_d and variable γ . For functions $f_{14} - f_{23}$ eleven runs were made with $0 \leq \gamma \leq 1$ in increments of 0.1 and $4 \leq N_p/N_d \leq 14$ by 2 (66 runs total). For $f_1 - f_1$ the same procedure was used, but with $2 \leq N_p/N_d \leq 6$ by 2 (33 runs total) in order to avoid excessive runtimes. Table 2 shows the γ value corresponding to the best fitness, γ_{best} , and the corresponding best value of N_p/N_d . *N*_{eval} is the number of function evaluations, and it is tabulated for the single best run and for the group of runs used to determine γ_{best} and the best number of probes per axis. The CFO column contains results for runs in which pseudorandomness has been removed from the algorithm except for the probe line IPD, which was unchanged. Those runs employed fixed decision space boundaries, a fixed value of $F_{\text{ren}} = 0.5$, and a fixed value of $N_p/N_d = 10$.

Comparing CFO-PR to GSO/GA/PSO, in the first group of high dimensionality unimodal functions, f_1 - f_7 , CFO-PR returned the best fitness on five of the seven functions $(f_3 - f_7)$. PSO performed best on the first two. In the second set of high dimensionality multimodal functions with many local maxima, $f_8 - f_{13}$, CFO-PR performed best on two (f_9, f_{10}) and essentially the same as the best other algorithm (GSO) on *f8*. In last group of ten multimodal functions with few local maxima, f_{14} - f_{23} , CFO-PR returned the best fitness on four $(f_{20} - f_{23})$, equal fitnesses on three $(f_{14}$, *f17, f18*), and very slightly lower fitnesses on the remaining three. Even though it is in its infancy, CFO-PR performed very well against three other highly sophisticated algorithms. It returned the best, equal, essentially equal, or very slightly lower fitnesses on eighteen of the twenty three test functions. It is reasonable to conclude that, overall, CFO-PR performed as well or better than GSO, which in turn performed better than PSO or GA.

Comparing CFO-PR to CFO alone (most pseudorandomness taken out), CFO alone returned the best fitness overall on f_1 - f_4 , f_6 , f_9 , and on f_{10} . On f_{19} , GSO, CFO-PR and CFO all returned the same fitness, which happens to be the global maximum. On the other benchmarks, however, CFO with reduced pseudorandomness returned worse best fitnesses than CFO-PR. These results suggest that

CFO-PR performs better than CFO alone, but not by a wide margin.

Fig. 8. Evolution of GP function *Davg*

Fig. 9. GP Function best probe number

2D GP TRAJECTORIES OF PROBES WITH BEST FITNESSES (ORDERED BY FITNESS) 03-25-2013, 10:08:51

Fig. 10. GP Function trajectories of probes with best fitness

2D GP INDIVIDUAL PROBE TRAJECTORIES (ORDERED BY PROBE #) 03-25-2013, 10:08:51 100 "p1" "p9" "p10" "p2 **Yo11** $\sqrt{12}$ 50 $p13$ $x₂$ $\pmb{0}$ o 1d -50 -100 $\mathbf{0}$ -50 50 100 -100 $x1$

Fig. 11. GP function probe trajectories by probe number

4 Conclusion

This note suggests that pseudorandomness is an important, indeed perhaps essential, aspect of effective CFO implementations. A pseudorandom variable has an arbitrary but precisely known value that may be assigned or calculated. Its essential characteristic is that the value is uncorrelated with the decision space's topology, so that it has the effect of distributing probes pseudorandomly throughout the landscape. While in a general sense this process may appear to be similar to the randomness in an inherently stochastic algorithm, it is in fact fundamentally different. The equations underlying stochastic algorithms are formulated in terms of true random variables whose values are computed from probability distributions and consequently are unknowable until the calculation is made. Therefore successive calculations yield different values, and as a result every optimization run has a different outcome. By contrast, a pseudorandom variable in the context of CFO is known with absolute precision because of how its value is determined (assignment or deterministic calculation). This property allows CFO to compute probe trajectories precisely because it is inherently deterministic. Every CFO with the same setup, with or without a pseudorandom component, yields exactly the same results, step-by-step throughout the entire run. Importantly, CFO's reproducibility lends itself well to reactive implementations in which run parameters are tuned in response to performance metrics such as rate of convergence or fitness saturation as examples. Reactive stochastic algorithms, on the other hand, are very difficult to implement.

This paper provides examples of how pseudorandomness can improve CFO's performance. Three different approaches are used (initial probe distribution, repositioning factor, and decision space adaptation), and each was discussed in detail. A sample CFO-PR problem was presented in detail, and summary data included for a standard 23-function benchmark suite. CFO-PR's performance is quite good compared to other highly developed, state-of-the-art algorithms. In addition, data are presented for a CFO implementation in which all pseudorandomness except the IPD has been removed, and those results show that, as a general rule, injecting pseudorandomness improves CFO's performance. Hopefully these results will encourage further work on improved methodologies for injecting pseudorandomness into CFO, in particular where and how. Of course, any or all of CFO's run parameters can be pseudorandomized, not only the three considered here. But even with respect to those parameters, different approaches to how they are pseudorandomized may yield better results or faster runtimes. There are many fruitful areas of research on CFO, and it is the author's hope that this and the other CFO papers will provide the foundation and catalyst for that work.

Competing Interests

Author has declared that no competing interests exist.

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Appendix

CFO searches an N_d -dimensional decision space Ω for the global maxima of an objective function $f(x_1, x_2,...,x_{N_d})$ defined on $\Omega: x_i^{\min} \leq x_i \leq x_i^{\max}$, $1 \leq i \leq N_d$. The x_i are the decision variables, and *i* the coordinate number. The term fitness refers to the value of $f(\vec{x})$ at point \vec{x} in Ω . There is no *a priori* information about the objective function's maxima, that is, $f(\vec{x})$'s topology (landscape) is unknown [1].

CFO searches Ω by flying probes through the space at discrete time steps (iterations). Each probe's location is specified by its position vector computed from two equations of motion that analogize their real-world counterparts for material objects moving through physical space under the influence of gravity without energy dissipation.

Probe *p* 's position vector at step *j* is $R_j^p = \sum x_k^{p,j} \hat{e}_k$, where the $x_k^{p,j}$ are it N_d *k* p, \dot{f} \hat{a} *reparation* p, \dot{f} *are if* $k \in \mathcal{E}_k$, where the \mathcal{X}_k a $\vec{R}^p_j = \sum_{k=1}^{N_d} x_k^{p,j} \hat{e}_k$, where the $x_k^{p,j}$ are its coordinates and $\vec{R}_j^p = \sum_{k=1}^{N_d} x_k^{p,j} \hat{e}_k$, where the $x_k^{p,j}$ are if , where the $x_k^{p,j}$ are its coordinates and \hat{e}_k the unit vector along the x_k -axis. The indices p , $1 \leq p \leq N_p$, and j , $0 \leq j \leq N_t$, respectively, are the probe number and iteration number, where N_p and N_t are the corresponding total number of probes and total number of time steps.

Equations of Motion: In metaphorical CFO space each of the N_p probes experiences an acceleration created by the gravitational pull of masses in Ω . Probe p 's acceleration at step $j-1$ is given by

$$
\vec{a}_{j-1}^p = G \sum_{\substack{k=1 \ k \neq p}}^{N_p} U(M_{j-1}^k - M_{j-1}^p) \cdot (M_{j-1}^k - M_{j-1}^p)^\alpha \times \frac{(\vec{R}_{j-1}^k - \vec{R}_{j-1}^p)}{\left\| \vec{R}_{j-1}^k - \vec{R}_{j-1}^p \right\|^\beta},\tag{1}
$$

which is the first of CFO's two equations of motion. In equation (1), $(x_1^{p,j-1}, x_2^{p,j-1},..., x_N^{p,j-1})$ is the objective function's fith 2 $, \ldots, \Lambda_{N_d}$ *J* is the object $\mathcal{L}_{-1}^p = f(x_1^{p,j-1}, x_2^{p,j-1},..., x_{N_d}^{p,j-1})$ is the objective function's N_d \rightarrow 18 the objective function 8 fm $M_{j-1}^p = f(x_1^{p,j-1}, x_2^{p,j-1},...,x_{N_d}^{p,j-1})$ is the objective function's fitness at probe *p* 's location at time step $j-1$. Each of the other probes at that step (iteration) has associated with it fitness $M^k_{j-1}, k = 1,..., p-1, p+1,...,N_p$. *G* is CFO's gravitational constant, and $U(\cdot)$ is the Unit Step function, $U(z) = \begin{cases} 0, & otherwise \end{cases}$. $\left\{ \cdot \right\}$ $\mathbf{1}$ and $\mathbf{1}$ and $\mathbf{1}$ and $\mathbf{1}$ and $\mathbf{1}$ and $\mathbf{1}$ and $\mathbf{1}$ $\left\{\begin{array}{c} 0 & 0 & 0 \\ 0 & 0 & 0 \end{array}\right\}$ $\begin{bmatrix} 1, & z \ge 0 \end{bmatrix}$ $=\begin{cases} 1, & -1 \\ 0, & 1 \end{cases}$ *otherwise* $z \geq 0$ | $U(z) = \begin{cases} 1, & z = 0 \\ 0, & z = 1 \end{cases}$ 1, $z \ge 0$
0, otherwise $(z) = \begin{cases} z & -1 \\ 0 & 1 \end{cases}$.

The acceleration \vec{a}^p_{j-1} causes probe *p* to move from position \vec{R}^p_{j-1} at step $j-1$ to \vec{R}^p_j at step *j* according to the trajectory equation

$$
\vec{R}_{j}^{p} = \vec{R}_{j-1}^{p} + \frac{1}{2} \vec{a}_{j-1}^{p} \Delta t^{2}, \ j \ge 1 \tag{2}
$$

which is CFO's second equation of motion. Note that the original CFO paper [1] included a velocity term that was set equal to zero as a matter of convenience because it simply was a additive constant in the case of rectilinear motion. Upon further consideration, however, it became clear that this term should not be included in equation (2) because, in general, a probe's motion is not rectilinear. Instead it is curvilinear, in which case the acceleration and velocity vectors are in different directions. As an example, in the case of circular motion the velocity vector is tangent to the trajectory circle while the acceleration is inwardly directed along the radius, that is, perpendicular to the velocity. This limiting case illustrates why, in general, the velocity term appearing in real-world kinematic equations should not be included in metaphorical CFO-space because it changes the direction of each probe's acceleration.

The CFO equations of motion, (1) and (2), combine to compute a new probe distribution at each time step using masses discovered by the probe distribution at the previous step. Δt is the time interval between steps during which the acceleration is constant. Note that CFO's terminology has no significance beyond reflecting CFO's kinematic roots, as does the factor $\frac{1}{2}$ in eq. (2). The gravitational constant, G , and time increment, Δt , have direct analogues in Newton's equations of motion for real masses moving under real gravity through three-dimensional physical space. The CFO exponents α and β , by contrast, have no analogues in Nature. They provide added flexibility to the algorithm designer who, in metaphorical CFO space, is free to change how gravity varies with distance, or mass, or both, if doing so creates a more effective algorithm.

Mass: The concept of mass in CFO space is very important and quite different than it is in real space. Mass in the physical Universe is an inherent, immutable property of matter, whereas in CFO space it is a positive-definite user-defined function of the objective function's fitness, not necessarily the fitness itself. For example, in equation (1) mass is defined as $(M_{j-1}^k - M_{j-1}^p) \cdot (M_{j-1}^k - M_{j-1}^p)$ ^{*a*} [difference in fitness value] $p \rightarrow \alpha$ Lifferance in fitness $j-1$ ^{*j*} [unicicii $MASS_{CFO} = U(M_{j-1}^k - M_{j-1}^p) \cdot (M_{j-1}^k - M_{j-1}^p)^\alpha$ [difference in fitness values raised to the α power multiplied by the Unit Step]. A different function can be used if it results in a better performing CFO algorithm. In this specific implementation the Unit Step is a critical element because it prevents negative mass. Without the Unit Step CFO mass could be negative depending on which fitness is greater. But mass in the real Universe always is positive, and as a consequence the force of gravity always attractive. By contrast, mass can be positive or negative in metaphorical CFO space, depending on how it is defined, and undesirable effects may result from the wrong definition. Negative mass creates a *repulsive* gravitational force that flies probes away from maxima instead of toward them, thus defeating the very purpose of the algorithm. See [1,3] for graphical examples of the effect of repulsive gravitational force.

Errant Probes: At any iteration in a CFO run, it is possible that a probe's acceleration computed from eq. (2) may be too great to keep it inside Ω . If any coordinate $x_i < x_i^{\min}$ or $x_i > x_i^{\max}$, the probe enters a region of unfeasible solutions that are not valid for the problem at hand. The question is what to do with an errant probe, and it arises in many algorithms. There are many approaches. While many schemes are possible, a simple, empirically determined one is used here. On a coordinate-by-coordinate basis, probes flying out of the decision space are placed a fraction

 $\Delta F_{rep} \leq F_{rep} \leq 1$ of the distance between the probe's starting coordinate and the corresponding boundary coordinate. *Frep* is the variable repositioning factor (see, for example, [2,3] for a more detailed discussion). Its value, as well as those of all the CFO parameters, was determined empirically.

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