Single-agent Finite Impulse Response Optimizer vs Simulated Kalman Filter Optimizer

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ABSTRACT – Single-agent Finite Impulse Response Optimizer (SAFIRO) is a new estimation-based optimization algorithm which mimics the work procedure of the ultimate unbiased finite impulse response (UFIR) filter [1]. Introduced in 2018, the SAFIRO is the latest edition of estimation-based algorithms.

On the other hand, a metaheuristic algorithm called simulated Kalman filter (SKF), has been proposed in 2015 for numerical optimization problems [2-4]. It was intro-duced as population-based metaheuristics, where the search for optimal solution is conducted by a group of agents. The agents of SKF work like Kalman filters [5], where they go through prediction, measurement, and estimation process in every iteration. The measurement in SKF is a simulated measurement which is obtained using mathematical equation.

Many studies on SKF can be found in literature. For example, the SKF has been studied fundamentally [6-7]. The SKF also has been extended for binary optimization problems [8] and combinatorial optimization problems [9-11]. Hybridization of SKF with particle swarm optimization (PSO), gravitational search algorithm (GSA), and opposition-based learning [12-18] have also been proposed for better performance. Other variants called parameter-less SKF and randomized SKF algorithms were proposed in [19-20]. The SKF has also been applied for real world problems like the adaptive beamforming in wireless cellular communication [21-24], airport gate allocation problem [25-26], feature selection of EEG signal [27-28], system identification [29-30], image processing [31-32], controller tuning [33], and PCB drill path optimization [34-35].

Given the popularity of SKF algorithm and the amount of studies reported on the fundamental improvements and application of SKF, the performance of the SAFIRO is compared with that of SKF algorithm based on CEC2014 benchmark dataset. In the subsequent section, the SKF algorithm is firstly introduced. Then, the SAFIRO is explained. Next, experimental procedure is presented, and the result of the experiment is shown. Finally, the result of a statistical analysis is shown to conclude the findings of this study.

The Simulated Kalman Filter

The SKF algorithm follows the algorithm shown in Fig. 1. One iteration consists of fitness evaluation, update the best solution, predict, measure, and estimate.

Using n agents, a set of solution can be denoted as \(X(t) = \{X_1(t), X_2(t), ..., X_n(t)\}\), where \(t\) is the iteration number. The SKF starts with random initialization of solutions. In each iteration, the fitness of the agents’ are evaluated. Then, the agent with the best fitness value is identified as the best solution of the current population, \(X_{\text{best}}(t)\). Next, the best \(X_{\text{best}}(t)\) from the first iteration is selected as \(X_{\text{true}}\).
During the prediction phase, the current predicted state, $X_{i}(t|t+1)$, is assumed to be the estimated value:

$$X_{!}(t|t+1) = X_{!}(t)$$  \hspace{1cm} (1)

The error covariant is also updated as follows:

$$P(t|t) = P(t) + Q$$  \hspace{1cm} (2)

where $P(t)$ and $P(t|t+1)$ denote the current error covariant estimate and current transition error covariant estimate, respectively. Note that the error covariant estimate is influenced by the process noise, $Q$.

In SKF, measurements are simulated using an agent’s prediction and $X_{true}$. The dimensional wise calculation of measured value for each dimension of agent $i$th is calculated as follows:

$$Z_{i}(t) = X_{i}(t|t) + \sin(2\pi r_{i}(t)) \times |X_{i}(t|t) - X_{true}|$$  \hspace{1cm} (3)

where $r_{i}(t)$ is a random value within the range of [0,1]. The estimation phase follows the measurement phase and the estimated next value is updated using (4):

$$X_{i}(t+1) = X_{i}(t|t+1) + K(t) \times (Z_{i}(t) - X_{i}(t|t+1))$$  \hspace{1cm} (4)

where $K(t)$ is the Kalman gain, which is calculated as follows:

$$K(t) = P(t|t+1)/(P(t|t+1)+R)$$  \hspace{1cm} (5)

where $R$ is the measurement noise. Then, the current error covariant estimate is updated in estimation phase using (6):

$$P(t+1) = (1-K(t)) \times P(t|t+1)$$  \hspace{1cm} (6)

These steps continue until at the end of the iteration or at the end of the fitness evaluation.

### The Single-agent Finite Impulse Response Optimizer (SAFIRO)

The SAFIRO is a single-agent metaheuristic algorithm, recently proposed for single-objective numerical optimization problems, which mimics the framework of the estimation process in UFIR filter [1]. Like the UFIR filter, SAFIRO’s agent responsible to perform the measurement and estimation stage to estimate the solution. The solution in SAFIRO represented by the estimation of the agent’s position.

The flowchart of SAFIRO is depicted in Fig. 2. The horizon length, $N$ is the parameter that needs to be used in SAFIRO. As SAFIRO needs $N$ measurements to begin the optimization process, the value of $N$ is defined during the initialization stage. Then, by using a single agent, SAFIRO starts the process with the initialization of $N$ measurements, $Y(t) = rand(U[X_{min}, X_{max}])$. After initialization, these $N$ random initial measurements are evaluated by using the fitness function of the problem to determine the initial $X_{best \_ so \_ far}$. $X_{best \_ so \_ far}$ represents the best-so-far solution. For minimization problem, the initial measurement with the smallest fitness value is recorded as $X_{best \_ so \_ far}$, meanwhile, for the maximization problem, the initial measurement with the largest fitness value is recorded as $X_{best \_ so \_ far}$.

Next, the agent undergoes the measurement stage. In the UFIR filter operation, the measurement data can be obtained from the sensor. However, in SAFIRO, the measurement is simulated by random mutation of $X_{best \_ so \_ far}$ and shrinking local neighbourhood method. Each dimension of the problem to be optimized has a random value in the range of 0 to 1. The dimensions with a random value equal to or smaller than 0.5, keep the $X_{best \_ so \_ far}$ value as their measurement value, as in (7):
On the other hand, the dimensions with a random value larger than 0.5, are chosen to be mutated to generate a new candidate solution. This mutation process is conducted in a local neighbourhood of $X_{\text{best so far}}$. The measurement value for these dimensions can be calculated as in (8).

$$Y_d(t) = X_{\text{best so far}} + \text{rand}(U[-\delta, \delta])$$ (8)

This measurement process helps to encourage exploration through the mutation process, and at the same time creating a balance between the exploration and exploitation through the shrinking local neighbourhood. In the local search method, the search area is centred around $X_{\text{best so far}}$. The radius of the local neighbourhood, $\delta$, can be computed by using as follows:

$$\delta = \exp(-\beta t) \times 0.5(X_{\text{max}} - X_{\text{min}})$$ (9)

where $t$ is the number of the current iteration, $T$ is the number of maximum iteration, $X_{\text{max}}$ is the upper limit of search space, $X_{\text{min}}$ is the lower limit of search space, and $\beta$ is a coefficient value. This coefficient value controls the reduction speed of the neighbourhood’s size.

After the measurement stage, SAFIRO’s agent moves to the estimation stage. At this stage, the solution of SAFIRO is updated in a finite length according to the number of $N$. Each iteration, $t$, consists of sub-iteration, $k$. As depicted in Fig. 3, the first two points of the horizon are used for initial estimation, $\tilde{X}(k=2)$. Initial estimation is generated randomly between [lower limit, upper limit] of the first and the second point of the horizon. Then, the remaining points are used for iterative estimation update from $\tilde{X}(k=3)$ until $\tilde{X}(k=N)$. The solution of initial estimation is improved iteratively during this part which is depending on the value of $N$. The number of repetition for iterative estimation is equal to $N-2$ (as mentioned above, the first two points are used to generate the initial estimation). In this study, the effect of the number of $N$ towards SAFIRO’s performance is observed. Iterative estimation can be computed by using (10) and (11):

$$\tilde{X}(k) = \tilde{X}(k-1) + K(k)(Y(t-N+k) - \tilde{X}(k-1))$$ (10)

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**Figure 2.** The flowchart of the SAFIRO.
\( K(k) = (k)^1 \)  

where \( \tilde{X}(k) \) is the estimated solution for current point, meanwhile \( \tilde{X}(k - 1) \) is the previous sub-iteration point. The measurement value, \( Y(t - N + k) \) and the Kalman-like gain, \( K(k) \) influence the improvement of the solution. The value of \( K(k) \) can help to improve the estimation as the sub-iteration, \( k \) increases. The sub-iteration process end when \( k = N \) is reached. The final value of \( k \) is then stored as \( X(t) = \tilde{X}(k) \). The agent’s updated solution for that corresponding iteration is represented by the estimated value of \( X(t) \).

After obtaining the solution, \( X(t) \), the evaluation stage is done to evaluate its fitness. The fitness of \( X(t) \) is then compared to the fitness of \( X_{best \_so \_far} \). The \( X_{best \_so \_far} \) is updated when a better solution is found. For minimization problem, \( X_{best \_so \_far} \) is updated if \( fit(X(t)) < fit(X_{best \_so \_far}) \), meanwhile for maximization problem, \( X_{best \_so \_far} \) is updated if \( fit(X(t)) > fit(X_{best \_so \_far}) \). Measurement and estimation stages are repeated until the maximum iteration, \( T \) is met, and the \( X_{best \_so \_far} \) returns as the solution to the given optimization problem.

\section*{Experiment}

The CEC 2014 benchmark suite [36] is used for performance comparison. The experimental settings are as listed in Table 1. All parameter values listed in Table 1 follow the parameter setting in the CEC 2014 benchmark suite. The maximum number of function evaluation, maxFES = 10,000 iterations \( \times \) dimension, \( D \).

Regardless of the number of agents, a fair comparison of algorithms’ performance can be done by setting the same FES. To provide a fair evaluation for each comparison, the complexity of the problem is set as 50 dimensions, while the maxFES is set as 500,000. The stopping condition is set to be the maxFES. All experiments are run 51 times on each test function. Hence, the evaluations are based on the mean fitness value over 51 runs time with 500,000 maxFES of 50 problem dimensions. For all functions, the search space in the range of \([-100,100] \) is used for all dimensions.

For SKF, the value of the initial error covariance, \( P(0) \), the measurement noise, \( R \), and the process noise, \( Q \) are set as random \([0,1]\). For FIROs, the coefficient, \( \beta \) value is assigned as \( \beta = 10 \) to allow moderate transition from the exploration phase to the exploitation phase. The coefficient, \( \beta \), is a control parameter employed in the shrinking local neighborhood method of FIROs, which is computed during the measurement step. This parameter is used in the exponential decay equation to control the reduction of step-size.

\section*{Result and Discussion}

The experimental results are tabulated in Table 2, Table 3, Table 4, and Table 5. Even though the SAFIRO outperforms the SKF in majority case studies of hybrid and composition functions, mix results are observed for unimodal and simple multimodal functions. To obtain a concrete conclusion, Wilcoxon signed rank test [37] was used.
Table 1. Experimental setting.

| Parameter            | Value                       |
|----------------------|-----------------------------|
| Search space         | [-100,100]                  |
| No. of runs          | 51                          |
| No. of dimensions    | 50                          |
| No. of function evaluations | 10,000 iterations×50 dimensions = 500,000 maxFES |

Table 2. The mean fitness values of SAFIRO and SKF for unimodal functions.

| Function | Ideal Fitness | SAFIRO | SKF       |
|----------|---------------|--------|-----------|
| 1        | 100           | 7.98E+05 | 1.18E+07 |
| 2        | 200           | 7695.60 | 2.35E+08 |
| 3        | 300           | 300.0000 | 1.90E+04 |

Table 3. The mean fitness values of SAFIRO and SKF for simple multimodal functions.

| Function | Ideal Fitness | SAFIRO | SKF       |
|----------|---------------|--------|-----------|
| 4        | 400           | 488.72 | 569.32    |
| 5        | 500           | 520.0000 | 520.03 |
| 6        | 600           | 619.03 | 636.77    |
| 7        | 700           | 700.0100 | 702.87 |
| 8        | 800           | 994.29 | 818.99    |
| 9        | 900           | 1095.90 | 1086.43 |
| 10       | 1000          | 5785.20 | 1635.11  |
| 11       | 1100          | 6462.40 | 6810.65  |
| 12       | 1200          | 1200.10 | 1200.32  |
| 13       | 1300          | 1300.60 | 1300.58  |
| 14       | 1400          | 1400.60 | 1400.30  |
| 15       | 1500          | 1511.20 | 1627.16  |
| 16       | 1600          | 1620.60 | 1619.60  |

Table 4. The mean fitness values of SAFIRO and SKF for hybrid functions.

| Function | Ideal Fitness | SAFIRO | SKF       |
|----------|---------------|--------|-----------|
| 17       | 1700          | 4.60E+04 | 1.44E+06 |
| 18       | 1800          | 3980.80 | 6.33E+07 |
| 19       | 1900          | 1920.60 | 1960.45  |
| 20       | 2000          | 2476.10 | 3.84E+04 |
| 21       | 2100          | 6.11E+04 | 2.30E+06 |
| 22       | 2200          | 2920.90 | 3447.67  |
The Wilcoxon test usually is used when the population cannot be assumed to be normally distributed or it can be used to compare two related samples, matched samples, or repeated measurements on a single sample to assess whether their population mean ranks differ. Particularly, the null hypothesis for the test assumes that there is no significant difference between the mean error values of test algorithms while the alternative hypothesis tries to determine if there is a significant difference between test algorithms using 5% ($\alpha = 0.05$) significance level. Since the number of samples is 30, the critical value for the test is equal to 152. The sum of ranks where the SKF algorithm outperforms the SAFIRO is denoted as $R_-$ while the sum of ranks where the BH algorithm is outperformed by the SAFIRO is denoted as $R_+$. Hence, the SAFIRO is better than the SKF algorithm if $R_+ > R_-$ and the SAFIRO is significantly better than the SKF algorithm if $R_-$ value is less than the critical value. Based on the mean accuracy of SAFIRO and SKF, $R_+$ and $R_-$ obtained are 382 and 183, respectively. Hence, Wilcoxon test result shows that the SAFIRO algorithm is significantly better than the SKF algorithm.

Conclusions

The SKF is the most popular estimation-based optimization algorithm. On the other hand, the SAFIRO is the latest emerging estimation-based optimization algorithm. Since the original work of SAFIRO did not include the SKF in benchmarking, this paper brings these two algorithms to the same platform for a performance evaluation. It is found that for global optimization, the SAFIRO outperforms the SKF significantly.

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