

# Iterated Local Search and Particle Swarm Optimization for Floating-point Error Estimation

## Preliminary:

Both Iterated Local Search (ILS) and Particle Swarm Optimization (PSO) divide each value range into  $N_{gra}$  mutual exclusive divisions. ( $N_{gra}$  is a parameter that is decided by users.) Consequently, for an initial configuration  $C_{init}$ , ILS/PSO creates a (fixed) set of tighter configurations  $\mathbb{C}_{univ}$ .

$$C_{init} : \left\{ \begin{array}{l} I_1 \mapsto R_1 \\ \dots \\ I_n \mapsto R_n \end{array} \right\} \quad \mathbb{C}_{univ} : \left\{ \left\{ \begin{array}{l} I_1 \mapsto R_1^{p_1/N_{gra}} \\ \dots \\ I_n \mapsto R_n^{p_n/N_{gra}} \end{array} \right\} \right\}$$

$$0 \leq p_1, \dots, p_n < N_{gra}$$

Two configurations in  $\mathbb{C}_{univ}$ ,  $c_x$  and  $c_y$ , are “neighbors” if  $c_x = c_y$  or they satisfy all the following constraints: 1)  $c_x$  and  $c_y$  have the same domain, 2) there exists only one variable  $I_i$  such that  $c_x(I_i)$  and  $c_y(I_i)$  are adjacent ranges, and 3) for any variable  $I_j$  which is not equal to  $I_i$ ,  $c_x(I_j) = c_y(I_j)$ . A configuration,  $c_x$ , can “move” to another configuration,  $c_y$ , by one step if  $c_x$  and  $c_y$  are neighbors. A configuration,  $c_x$ , can “move”  $t$  steps to  $c_y^t$  if there exists a sequence of configurations,  $c_x, c_y^1, \dots, c_y^t$ , such that  $c_x$  and  $c_y^1$  are neighbors and  $c_y^i$  and  $c_y^{i+1}$  are neighbors ( $1 \leq i < t - 1$ ). ILS/PSO searches among  $\mathbb{C}_{univ}$  and tries to detect high floating-point errors.

## High-level Idea of ILS:

ILS starts from a single configuration in  $\mathbb{C}_{univ}$ , randomly moves to another configuration, evaluate the configuration with all its neighbors, and chooses the “best” configuration to start the next iteration.

## High-level Idea of PSO:

PSO holds a group of configurations called a “swarm.” It evaluates every configuration in the swarm and ranks them by their evaluation results (measured floating-point errors). To find a new swarm, PSO moves every configurations in the swarm according to their ranks: the higher measured floating-point error, the fewer step for the configuration to move.

## Helper Functions of ILS and PSO:

Algo. 1 shows the shared helper functions used by ILS and PSO. Function *RandConf* randomly returns a configuration in  $\mathbb{C}_{univ}$ . Function *NeighborConf*s takes a configuration and returns all its neighbor configurations. Function call *MoveConf*( $c, t$ ) moves a given configuration ( $c$ )  $t$  steps.

## ILS Algorithm

Before introducing ILS algorithm (in Algo. 2), we introduce an important ILS’s subroutine “*IFI*.” *IFI* takes a configuration, enumerates all its neighbors, and evaluates all the enumerated configurations. After evaluations, *IFI* returns the configuration that resulted in the highest floating-point error.

ILS algorithm (Algo. 2) is compromised by three phases:

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**Algorithm 1** Helper Functions for ILS and PSO

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1: global:  $C_{init}, N_{gra}$ 
2: procedure RANDCONF
3:   return: random  $C_{univ}$ 
4: end procedure
5:
6: procedure NEIGHBORCONFS( $conf$ )
7:    $conf_s \leftarrow \{conf\}$ 
8:   for each  $(I_i \mapsto R_i^{p/N_{gra}}) \in conf$  do
9:      $conf' \leftarrow conf \setminus \{(I_i \mapsto R_i^{p/N_{gra}})\}$ 
10:     $p^+ \leftarrow \max(p + 1, N_{gra} - 1)$ 
11:     $p^- \leftarrow \min(p - 1, 0)$ 
12:     $conf^+ \leftarrow conf' \cup \{(I_i \mapsto R_i^{p^+/N_{gra}})\}$ 
13:     $conf^- \leftarrow conf' \cup \{(I_i \mapsto R_i^{p^-/N_{gra}})\}$ 
14:     $conf_s \leftarrow conf_s \uplus conf^+ \uplus conf^-$ 
15:   end for
16:   return:  $conf_s$ 
17: end procedure
18:
19: procedure MOVECONF( $conf, n_{move}$ )
20:    $c \leftarrow conf$ 
21:   for  $i = 1$  to  $n_{move}$  do
22:      $c \leftarrow$  random  $NeighborConfs(c)$ 
23:   end for
24:   return:  $c$ 
25: end procedure
```

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- Phase 1 (line 22 to 29) randomly chooses  $N_{p1}$  configurations, evaluates them, and preserves the one ( $CurrConf$ ) which is evaluated to the highest error.
- Phase 2 (line 31) explores all  $CurrConf$ 's neighbors by invoking  $IFI$ .
- Phase 3 (line 33 to 40) randomly moves the  $CurrConf$  by  $N_{p2}$  steps, (by invoking  $MoveConf$ ) and evaluates all neighbors after moving.
- Repeat phase 3 until running out of "resource" (could be time or the total number of shadow value executions).
- Both  $N_{p1}$  and  $N_{p2}$  are user-decided parameters.
- Before moving the configuration in phase 3, we have an opportunity to choose a random starting point (line 37 to 39).

### Explanations of PSO (Algo. 3)

Before introducing PSO algorithm, we introduce some of PSO's terminologies first.  $\llbracket item_0 \dots item_n \rrbracket$  denotes an ordered list. Items,  $item_0 \dots item_n$ , are stored in order. For an ordered list  $l =$

$\llbracket item_0 \dots item_n \rrbracket$ , we use  $l[p]$  to denote the  $p$ 'th item of  $l$ :  $l[p] = item_p$ . A “swarm” is an ordered list of configurations. Function *RandSwarm* in Algo. 3 randomly generates a swarm.

PSO starts from a random swarm, and repeatedly evaluates the swarm until running out of resource. Evaluation of a swarm (*EvaSwarm* in Algo. 3) is composed by three phases:

- Phase 1 (line 15 to 20) evaluates all configurations in a swarm.
- Phase 2 (line 21) sorts all configurations by their. The configuration with the highest error will be given the highest rank.
- Phase 3 (line 22 to 24) moves configurations with different steps according to their ranks: the higher the rank, the fewer the steps to move.
- Note that *EvaSwarm* has side-effect on *WorstErr*.

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**Algorithm 2** Iterated Local Search for Floating-point Error Estimation

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1: Input:  $P, M_{init}, N_{p1}, N_{p2}, N_{gra}, N_{eva}$ 
2: Output: The Highest Floating-point Error.
3:  $CurrConf \leftarrow RandConf$ 
4: global  $WorstErr \leftarrow Eva(P, CurrConf, N_{eva})$ 
5:
6: procedure  $IFI(conf)$ 
7:    $LocalErr \leftarrow 0$ 
8:    $confs \leftarrow NeighborConfs(conf)$ 
9:   for all  $c \in confs$  do
10:     $err \leftarrow Eva(P, c, N_{eva})$ 
11:    if  $err > LocalErr$  then
12:       $LocalErr \leftarrow err$ 
13:       $LocalConf \leftarrow c$ 
14:    end if
15:  end for
16:  if  $LocalErr > WorstErr$  then
17:     $WorstErr \leftarrow LocalErr$ 
18:  end if
19:  return:  $LocalConf$ 
20: end procedure
21:
22: for  $i = 1$  to  $N_{p1}$  do
23:    $c \leftarrow RandConf$ 
24:    $err \leftarrow Eva(P, c, N_{eva})$ 
25:   if  $err > WorstErr$  then
26:      $WorstErr \leftarrow err$ 
27:      $CurrConf \leftarrow c$ 
28:   end if
29: end for
30:
31:  $CurrConf \leftarrow IFI(CurrConf)$ 
32:
33: while has resource do
34:    $c \leftarrow CurrConf$ 
35:    $c = MoveConf(c, N_{p2})$ 
36:    $CurrConf \leftarrow IFI(c)$ 
37:   if restart then
38:      $CurrConf \leftarrow RandConf$ 
39:   end if
40: end while
41: return:  $WorstErr$ 
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**Algorithm 3** Particle Swarm Optimization for Floating-point Error Estimation

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```
1: Input:  $P, C_{init}, N_{swarm}, N_{vel}, N_{gra}, N_{eva}$ 
2: Output: The Highest Floating-point Error.
3: global  $Swarm \leftarrow RandSwarm$ 
4: global  $WorstErr \leftarrow 0$ 
5:
6: procedure RANDSWARM
7:    $parts \leftarrow \square$ 
8:   for  $i = 1$  to  $N_{swarm}$  do
9:      $parts \leftarrow parts : (RandConf \mapsto 0)$ 
10:  end for
11:  return:  $parts$ 
12: end procedure
13:
14: procedure EVASWARM
15:  for all  $(c \mapsto e) \in Swarm$  do
16:     $err \leftarrow Eva(P, c, N_{eva})$ 
17:    if  $err > WorstErr$  then
18:       $WorstErr \leftarrow err$ 
19:    end if
20:  end for
21:   $Swarm \leftarrow \text{sort-by-err: } Swarm$ 
22:  for all  $Swarm[i]. 0 \leq i < N_{swarm}. Swarm[i] = c_i \mapsto err_i$  do
23:     $c_i = MoveConf(c_i, N_{vel} * i)$ 
24:  end for
25: end procedure
26:
27: while has resource do
28:    $EvaSwarm$ 
29: end while
30:
31: return:  $WorstErr$ 
```

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