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Minimizing Makespan in Distributed Blocking Flowshops Using Hybrid Iterated Greedy Algorithms

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ABSTRACT This paper studies the distributed blocking flow shop scheduling problem (DBFSP) using metaheuristics. A mixed integer programming model for solving the problem is proposed, and then three versions of the hybrid iterated greedy algorithm (HIG**1**, HIG**2**, and HIG**3**) are developed, combining the advantages of an iterated greedy algorithm with the operators of the variable Tabu list, the constant Tabu list, and the cooling schedule. A benchmark problem set is used to assess empirically the performance of the HIG**1**, HIG**2**, and HIG**³** algorithms. Computational results show that all the three versions of the proposed algorithm can efficiently and effectively minimize the maximum completion time among all factories of the DBFSP, and $HIG₁$ is the most effective.

INDEX TERMS Flowshop scheduling, hybrid meta-heuristic, distributed blocking flowshop.

I. INTRODUCTION

In today's globalized economy, many companies have turned from traditional single-factory production to multi-factory production to reduce the production risk and the cost of transportation. Consequently, distributed scheduling problems that concern the assignment of jobs to various factories and their subsequent sequence have been increasingly attracting the attention of researchers during the last decade. With respect to the various distributed scheduling problems, Ruiz and Naderi [1] were the first to present a distributed permutation flowshop scheduling problem (DPFSP), which was a novel generalization of the most popular permutation flowshop scheduling problem (PFSP). Ruiz and Naderi [1] also extended the well-known NEH heuristic [2] and two variable neighborhood descent search algorithms, referred to as VND_a and VND_b , to minimize the global makespan, which is the maximal completion time of the last job to be processed in any factory. These methods provided much better computational results than typical heuristic approaches, and the VND^a heuristic yielded the best solutions, on average, among the above methods.

Following the presentation of their pioneering conference paper in 2009, Naderi and Ruiz [3] subsequently presented a full journal paper that provided several MIP models and heuristics for solving the same problem. Thereafter, some effective and efficient improvement heuristics were presented to solve the DPFSP. Among these, the NEH-based heuristic algorithm [4], the genetic algorithm (GA) [5], [6], the Tabu search algorithm (TS) [7], the estimation of distribution algorithm (EDA) [8], the modified iterated greedy algorithm (MIG) [9], and the bounded-search iterated greedy algorithm (BSIG) [10] were increasingly better approximate methods for solving the DPFSP. In particular, the BSIG algorithm of Fernandez-Viagas and Framinan [10] stands out as the state-of-the-art heuristic for solving the DPFSP.

This work studies the distributed blocking flowshop scheduling problem (DBFSP) using meta-heuristics. Distributed blocking permutation flowshops are common in manufacturing, especially in the chemical, metal, pharmaceutical, electronic, plastic, food-processing, and service industries, for example [11]. In a blocking flowshop, a finite number or possibly zero intermediate buffers exist between

each successive pair of machines. Consequently, a processing job that is completed on a machine will be blocked on that instrument until the next machine downstream becomes available. Reddi and Ramamoorthy [12] were the first to study the two-machine blocking flowshop scheduling problem (BFSP). Having reduced the problem to a special case of the travelling salesman problem, they applied the wellknown Gilmore-Gomory algorithm [13] to solve it in polynomial time. Since the BFSP is NP-hard in the strong sense for a shop that has more than two machines [14], finding the optimal solution within an acceptable computation time using exact methods (such as those of Suhami and Mah [15], Ronconi and Armentano [16], and Ronconi [17] is very difficult even for problems of moderate size. Most studies of this highly complex problem have concentrated on developing constructive heuristics and improvement heuristics that can find high-quality – but not necessarily optimal – solutions in a short computation time.

The most famous constructive heuristics for tackling the BFSP include the profile fitting (PF) heuristic [18], the minmax heuristic (MM) [19], the combined MM and NEH heuristic (MME) [19], and the combined PF and NEH heuristic (PFE) [19]. Recently, Companys and Ribas [20] proposed some constructive heuristics to minimize the maximum completion time among all factories on a distributed blocking flowshop. The experimental results in the aforementioned works reveal that these constructive heuristics can rapidly find feasible solutions, and they are more useful than exact methods for solving complex BFSPs. Noteworthy improvement heuristics for solving the BFSP include the GA [21], Ronconi's algorithm (RON) [16], the fast TS algorithm [22], the improved TS algorithm $(TS+M)$ [22], the hybrid discrete differential evolution algorithm (HDDE) [23], the iterated greedy algorithm (IG) [24], and the revised artificial immune system algorithm (RAIS) [25]. Researchers have confirmed that, among these heuristic algorithms, TS+M, RON, HDDE, IG, and RAIS are the best for solving the BFSP. A more detailed discussion of related methods and their applications to various BFSPs can be found in the comprehensive investigation of Hall and Sriskandarajah [14]. Although some studies have investigated different DPFSPs and BFSPs, to the best of the authors' knowledge, there is only one research [20] has been done on the DPFSP with the blocking constraint. Companys and Ribas [20] proposed some constructive heuristics to minimize the maximum completion time among the factories for the DPFSP with the blocking constraint. The computational results showed good performance of these constructive heuristics, which could be applied to obtain a fast solution or as the initial solution procedure in more sophisticated meta-heuristics for solving the DPFSP.

This work presents three versions of the hybrid iterated greedy (HIG) algorithm that combine the advantages of the IG algorithm with the operators of the variable Tabu list, the constant Tabu list, and the cooling schedule, to solve the DPFSP with the blocking constraint. Notably, this paper

is the first study to provide effective and efficient IG-based algorithms for solving this problem. This paper is organized as follows. Section 2 defines the DBFSP and formulates it using a mixed integer linear (MIP) model. Section 3 describes in detail the three versions of the HIG algorithm. Section 4 presents results of simulations and statistical evaluations of the proposed algorithm, applied to a benchmark problem set of instances. Finally, Section 5 draws conclusions and makes recommendations for future studies.

II. DESCRIPTION AND FORMULATION OF PROBLEM

The DBFSP that is considered herein is described as follows. A set of n jobs is to be assigned to, and processed in, one of *f* identical factories, each with a flowshop production system that comprises the same set of m machines in series. Every factory can process all jobs, and each job can be assigned to and processed in any one of these factories. All jobs must be sequentially processed through the m machines of the assigned factory in an identical sequence without preemption. The processing time for each job may vary among machines, but does not change from factory to factory. All jobs are ready for processing at the beginning of the planning horizon, and all machines are available over the scheduling period. No intermediate buffer, which could store jobs until the next operation is performed, is assumed to exist between any pair of consecutive machines in any factory, so no upstream machine can release a completed job to the succeeding machine if the latter is busy. In such a case, the completed job must be blocked on the upstream machine until the preceding job has been completed on the succeeding machine.

The objective of scheduling is simultaneously to assign jobs to various factories and to determine their sequences to be processed in each factory to minimize the maximum completion time among the factories (global makespan). Notably, a schedule that minimizes the global makespan for a DBFSP also minimizes the sum of the job waiting times and the sum of the machine idle times. The DBFSP can be designated by a triplet $DF_m |block|C_{\text{max}}$ using the well-known notation of Pinedo [26]. Since the $DF_m|block|C_{\text{max}}$ problem with only one factory reduces to a general BFSP, which is NP-hard in the strong sense [14] when the number of machines exceeds two, the $DF_m |block|C_{\text{max}}$ problem can be confidently concluded also to be NP-hard in the strong sense.

Based on the MIP model of the DPFSP that presented by Naderi and Ruiz [3], the $DF_m|block|C_{\text{max}}$ problem can be formulated as the following MIP mathematical model. First, the following notation is defined to simplify the formulation. *Parameters*:

- *n* : Number of jobs
- *m* : Number of machines
- f : Number of factories
i : Index of machines, *i*
- Index of machines, $i \in \{0, 1, 2, \ldots, m\}$, where $i = 0$ is a dummy machine
- *j* : Index of jobs, $j \in \{1, 2, ..., n\}$
- *k* : Index of job position in a given sequence, $k \in \{1, 2, \ldots, n\}$
- *l* : Index of factories, $l \in \{1, 2, \ldots, f\}$
- *pj*,*ⁱ* : Processing time of job*j*on machine *i*

Decision variables:

$$
X_{j,k,l} = \begin{cases} 1, & \text{if job } j \text{ occupies position } k \text{ in factory } l \\ 0, & \text{otherwise} \end{cases}
$$

- $d_{k,i,l}$ = Departure/Completion time of job in position *k* on machine *i* in factory *l*
- C_{max} = Maximal completion time (global makespan) of the last job to be processed in any factory

The objective function of the MIP formulation is

Minimize *C*max

and the constraints are

s.t.
$$
d_{1,0,l} = 0
$$
, $l = 1, ..., f$; $k = 1, ..., F$, (1)
 $d_{k,0,l} = d_{k-1,1,l}$, $k = 2, ..., n$; $l = 1, ..., f$, (2)

$$
d_{1,i,l} = d_{1,i-1,l} + \sum_{j=1}^{n} X_{j,1,l} \cdot p_{j,i},
$$

\n
$$
i = 1, \dots, m-1; l = 1, \dots, f,
$$
 (3)

$$
d_{k,i,l} \ge d_{k,i-1,l} + \sum_{j=1}^{n} X_{j,k,l} \cdot p_{j,i}, \quad k = 2, ..., n;
$$

$$
i = 1, ..., m-1; l = 1, f
$$
 (4)

$$
i = 1, ..., m - 1; l = 1, ..., f,
$$

\n
$$
d_{k,i,l} \ge d_{k-1,i+1,l}, \quad k = 2, ..., n;
$$
\n(4)

$$
i = 1, ..., m-1; l = 1, ..., f,
$$
 (5)

$$
d_{k,m,l} = d_{k,m-1,l} + \sum_{j=1}^{n} X_{j,k,l} \cdot p_{j,m},
$$

\n
$$
k-1 \qquad n: l-1 \qquad f
$$
 (6)

$$
k = 1, ..., n; l = 1, ..., f,
$$
\n(6)

$$
\sum_{k=1}^{n} \sum_{l=1}^{f} X_{j,k,l} = 1, \quad j = 1, \dots, n,
$$
 (7)

$$
\sum_{j=1}^{n} X_{j,k,l} \le 1, \quad k = 1, \dots, n, \ l = 1, \dots, f, \tag{8}
$$

$$
d_{k,i,l} \ge 0, \quad k = 1, \dots, n; \ i = 1, \dots, m; \ l = 1, \dots, f,
$$
\n(9)

$$
C_{\max} \ge d_{k,m,l}, \quad k = 1, \dots, n; \ l = 1, \dots, f, \qquad (10)
$$

$$
X_{j,k,l} \in \{0, 1\}, j = 1, \dots, n; k = 1, \dots, n; l = 1, \dots, f.
$$
\n(11)

The goal is to minimize the maximal completion time (global makespan) of the last job to be processed in any factory. Constraint sets (1) and (2) define the starting time of the job in position k on the first machine in factory *l*. Constraint set (3) defines the departure time of the job in the first position on machine *i* in factory *l*. Constraint sets (4), (5) and (6) specify the relationship between departure times of each assigned job on two successive machines in a factory. Constraint set (7) ensures that each job is dispatched to exactly one factory and to exactly one job position in

the assigned factory. Constraint set (8) ensures that only one job can be allocated to each job position at a factory. Constraint set (9) specifies the departure time of each job on each machine as non-negative. Constraint set (10) computes the maximum completion time among all factories. Finally, constraint set (11) defines all relevant binary variables.

FIGURE 1. An outline of the generic IG algorithm.

III. PROPOSED SELF-TUNING ITERATED GREEDY ALGORITHM

This work proposes three versions of the HIG algorithm, which combine the IG algorithm with the operators of the Tabu list and the cooling schedule. The IG algorithm (Fig. 1) is a simple but robust stochastic search algorithm that was developed by Jacobs and Brusco [27]. As can be seen in Fig. 1, after an initial solution ξ_0 is obtained, a generic IG algorithm improves the incumbent solution ξ and the best solution ξ*best* through the iterative execution of two main phases, *destruction* and *construction*, until a specified termination condition (e.g., a maximum number of iterations, or a maximum computation time) is satisfied. In the *destruction* phase, a fixed number (α) of elements of ξ is eliminated, yielding a partial solution ξ*p*. In the following *construction* phase, the eliminated elements are individually and sequentially reinserted into all possible positions of the current partial solution to yield a new solution ξ*new*. After ξ*new* has been yielded, some acceptance criteria are applied to determine whether it will replace ξ and ξ_{best} . Owing to their simplicity, flexibility, and high efficiency, IG-based algorithms have been successfully utilized to solve some classic scheduling problems, such as the single-machine [28], the parallelmachine [29]–[31], the permutation flowshop [32]–[34], the non-permutation flowshop [35], the multistage hybrid flowshop [36], and the distributed permutation flowshop [9] scheduling problems.

The three versions of the HIG algorithm that are proposed in this work $(HIG₁, HIG₂, HIG₃)$ combine the Tabu list and the cooling schedule with the IG algorithm, respectively. The main difference between the $HIG₁$, $HIG₂$, and $HIG₃$ algorithms is they use the variable Tabu list, the constant Tabu list, and without Tabu list, respectively.

A highly effective speed-up method is used in the construction phase of all three versions of the HIG algorithm to reduce the computational burden. The following subsections further describe the coding scheme of the solutions, the detailed procedures of the three versions of the HIG algorithm and the speed-up method.

A. CODING SCHEME OF SOLUTIONS

To signify the assignment of jobs to different factories and their production sequences in each factory, a solution is coded using a numerical sequence of *n* integers, separated into segments by $f - 1$ asterisks, where each segment corresponds to the sequences of the jobs in an assigned factory. Here, integers represent jobs and asterisks specify the partitioning of jobs in the factories. For instance, a solution that is encoded as {6, 5, 1, 9, 12, 7, ∗, 13, 10, 8, 11, 15, ∗, 3, 14, 2, 4, 16} is a solution with 16 jobs in three factories, with production sequences in the first, second and third factories of {6, 5, 1, 9, 12, 7},{13, 10, 8, 11, 15}, and {3, 14, 2, 4, 16}, respectively.

B. PROCEDURES OF PROPOSED HIG ALGORITHM

The main procedures of the three versions of the HIG algorithm for solving the $DF_m|block|C_{\text{max}}$ problem are as follows.

Step 1: Generation of initial solution

- *Step 1.1* Apply the PW heuristic ([37], see Section 3.3) to yield a job list $J = (J_{[1]}, J_{[2]}, \ldots, J_{[n]}).$
- *Step 1.2* Apply the NEH₂ heuristic [3] to insert a job, in order from the first job in the job list *J*, into the current partial solution until all jobs have been inserted and an initial solution Π has thus been obtained. That is, each job is inserted into all possible positions in the current partial solution, and the one with the lowest partial global makespan, is subsequently utilized to replace the current partial solution before the insertion of the next job.
- *Step 1.3* Set Π as both the incumbent solution (Π^*) and the best solution (Π_{best}) , and set the accessible Tabu list $TL := \phi$.

Step 2: Destruction phase

Step 2.1 Determine the accessible Tabu list (*TL*) for the removal of jobs based on the Tabu list tenure (*TLT*), which is obtained using the following formula;

$$
TLT = TLT^{low} + u(TLT^{high} - TLT^{low})
$$

where TLT^{high} and TLT^{low} are the maximal and the minimal Tabu list tenure, respectively *u* is set as a random number between 0 and 1 (meaning that a variable Tabu list is used) for the HIG₁ algorithm, u is set as 0 (meaning that a constant Tabu list is used) for the $HIG₂$ algorithm; TLT^{low} and TLT^{high} are set as 0 (meaning that no Tabu list is used) for the $HIG₃$ algorithm.

- *Step 2.2* Select one job that is not in *TL* from each factory that with the largest completion time. Move the α_1 selected jobs (which are assumed to be all of the selected jobs) from Π^* to Π_R and put them into TL , where Π_R is a list of the removed jobs, arranged in order of their selection. Concurrently, set as Π_{P1}^* the current partial sequence of Π^* with the α_1 removed jobs eliminated.
- *Step 2.3* Select one job that is not in *TL* from each factory that has the smallest completion time. Move the α_2 selected jobs (which are assumed to be all of the selected jobs) from $\prod_{p=1}^{\infty}$ to Π_R in the order in which they were selected, and add them to *TL*. Concurrently, set as $\prod_{P_2}^*$ the current partial sequence of Π_{Pl}^* with the α_2 removed jobs eliminated.
- *Step 2.4* Randomly select $(α α₁ α₂)$ distinct jobs that are not in *TL* from all of the factories, where $\alpha \in$ [α_{min} , α_{max}]. Move the ($\alpha - \alpha_1 - \alpha_2$) selected jobs from \prod_{P}^* to Π_R in the order in which they were selected and add them to *TL*.

Step 3: Construction phase

Sequentially reinsert the jobs in Π_R , from the first position to the last position, into \prod_{P}^{*} , until a new solution (Π_{new}^*) has been constructed. When reinserting a job, all possible positions in the current partial solution should be considered and the best one, and the best position, which is the one with the lowest partial global makespan, is subsequently utilized to replace the current partial solution before the insertion of the next job. To accelerate the insertion operation, the speed-up method that is described in Section 3.4 is used.

Step 4: Acceptance criterion

To improve the ability of the incumbent solution to escape from local minima, the following acceptance criterion and the cooling schedule are used in all three versions of the HIG algorithm to determine whether Π^* and Π_{best} will be updated by the newly obtained solution $\overline{\Pi}_{new}^*$.

IF
$$
C_{\text{max}}(\Pi_{new}^*) \leq C_{\text{max}}(\Pi_{best})
$$
 THEN
\nset $\Pi_{best} := \Pi_{new}^*$ and $\Pi^* := \Pi_{new}^*$;
\nELSE_I F $C_{\text{max}}(\Pi_{new}^*) \leq C_{\text{max}}(\Pi^*)$ THEN set
\n $\Pi^* := \Pi_{new}^*$;
\nELSE_I F $C_{\text{max}}(\Pi_{new}^*) > C_{\text{max}}(\Pi^*)$ TEHN
\nGenerator $r \sim U(0,1)$;
\nIF $r < e^{((C_{\text{max}}(\Pi^*) - C_{\text{max}}(\Pi_{new}^*))/T)}$ set
\n $\Pi^* := \Pi_{new}^*$

Otherwise, discard Π_{new}^* .

Here, $C_{\text{max}}(\cdot)$ represents the global makespan of solution (·); $r \in [0, 1]$ is a pseudo-random number that is sampled from the standard uniform distribution U (0,1); and *T* denotes the current temperature with an initial temperature $T_0 = T_{Value} \cdot \sum_{i=1}^{m}$ *i*=1 $\sum_{n=1}^n$ $\sum_{j=1}^{n} p_{j,i}$ and will be decreased from its preceding temperature, i.e., $T \leftarrow \lambda T \ (\lambda \in [0, 1])$, after running a preset number of iterations (*Iiter*) at a particular temperature.

Step 5: Stopping criterion

To test fairly the three versions of the HIG algorithm, iterate Steps 2–4 for each algorithm until the computation time reaches a specified threshold (T_{max}) .

The PW heuristic was proposed by Pan and Wang [37] for the single blocking flowshop problem, and The NEH₂ heuristic was proposed by Naderi and Ruiz [3] for the distributed permutation flowshop scheduling problem. In this study, we combined the PW heuristic and the NEH₂ heuristic to generate an initial solution for the $DF_m|block|C_{\text{max}}$ problem. To compare the solution quality on the same basis, the procedures for implementing the three versions of the HIG algorithm are the same, except for the use of different Tabu lists in Step 2.1: the $HIG₁$, $HIG₂$, and $HIG₃$ algorithms use a variable Tabu list, a constant Tabu list, and no Tabu list, respectively. Additionally, in Step 4, all the three versions of the HIG algorithm employ the decreasing temperature mechanism instead of the constant temperature mechanism to avoid searching processes that would be prone to stagnation. The process used is executed by generating a pseudo-random number $r \in [0, 1]$ and updating the incumbent solution Π^* by \prod_{new}^* whenever $r < e^{(\left[\mathcal{C}_{\text{max}}(\Pi^*) - \mathcal{C}_{\text{max}}(\Pi^*_{new})\right]/T)}$.

C. PW HEURISTIC

The PW heuristic is a simple construction heuristic that was proposed by Pan and Wang [37] for solving the blocking flowshop scheduling problem. The PW heuristic is applied as follows.

Step 1: Use the following equations to calculate the departure time, $d_{[k],i}$, of job *j* on machine *i* if it is at position *k* in the schedule.

$$
d_{[1],0} = 0,\t(12)
$$

 $d_{[k]0} = d_{[k-1]1}, k = 2, \ldots, n,$ (13)

$$
d_{[k],i} = \max\{d_{[k],i-1} + p_{[k],i}, d_{[k-1],i+1}\},\
$$

$$
k = 2..., n; i = 1..., m - 1, (14)
$$

$$
d_{[k],m} = d_{[k],m-1} + p_{[k],m}, k = 1...,n.
$$
 (15)

Step 2: Use the following equations to calculate the slope index $f_{j,k}$ ($\forall j, k$) of job *j* at position *k*.

$$
f_{j,k} = (n - k - 2)\delta_{j,k} + \chi_{j,k},
$$
 (16)

Where

$$
\delta_{j,0} = \sum_{i=1}^{m} \frac{m}{i + \frac{k(m-i)}{n-2}} (d_{[1],i} - p_{j,i}),
$$

\n
$$
\delta_{j,k} = \sum_{i=1}^{m} \frac{m}{i + \frac{k(m-i)}{n-2}}
$$

\n
$$
\times (d_{[k+1],i} - d_{[k],i} - p_{j,i}),
$$
 and
\n
$$
\chi_{j,k} = \sum_{i=1}^{m} \frac{m}{i + \frac{k(m-i)}{n-2}} (d_{[k+2],i} - d_{[k+1],i}) - \sum_{\substack{q \in U \\ q \neq j}} \frac{p_{q,i}}{(n-k-1)}.
$$

- **Step 3:** Set the job with the smallest $f_{j,0}$ value as the first job in current partial job list $J_P = (J_{[1]})$. In case of a tie, make the job has the smallest $\chi_{i,0}$ value as the first job. Let the unscheduled job set $U = J - \{J_{[1]}\}.$
- **Step 4:** Repeat the following procedure to select a job for adding to the next position in the current partial job list until $U = \phi$.
	- *Step 4.1* For each machine, use Eqs. 12-16 to calculate the departure time, $d_{[k],i}$ ($i = 1, \ldots, m$), of the job at the last position, say position *k*, in the current partial job list.
	- *Step 4.2* Use Eq. 17 to calculate the slope index $f_{j,k}$ of each job $j \in U$. Remove the job with the smallest $f_{i,k}$ value from the unscheduled job set, and add it at the next position in the current partial job list. In case of a tie, make the job has the smallest $\chi_{i,0}$ as the first job.

D. SPEED-UP METHOD

In the construction phase (Step 3) of the three versions of HIG algorithm, the jobs in Π_R are successively inserted at all possible positions in $\Pi_{p_2}^*$, and the best one, which is the one with the lowest partial global makespan, is chosen. In such a construction phase, substantial time is taken to calculate the makespan of each possible (partial) solution of the insertion. To accelerate the evaluation of the best insertion position, a speed-up method that is revised the scheme of Wang et al. [23] for solving the blocking flowshop problem is proposed. Based on the assumption that, in the current partial solution, *n^P* jobs have been assigned to a factory, the following speed-up method is applied to evaluate $n_p + 1$ sequences that are generated by inserting a job, J_j , at all possible positions in this current partial solution.

- **Step 1:** Use Eqs. 11-15 to calculate the departure times $d_{[k],i}$ ($k = 1, ..., np$; $i = 1, ..., m$) of the *n^P* jobs at the assigned factory in the current partial solution.
- **Step 2:** Use the following equations to calculate the tails *f*_{[*j*],*i*} (*j* = 1, ..., *n_P*; *i* = 1, ..., *m*) of the

n^P jobs in the assigned factory in the current partial solution.

$$
f_{[np],m+1} = 0, \t\t(17)
$$

$$
f_{[np], i} = f_{[np], i+1} + p_{[np], i}, i = m, \dots, 1,
$$
 (18)

$$
f_{[j],m+1} = f_{[j+1],m}, j = n_P - 1, \dots, 1,
$$
 (19)

$$
f_{[j],i} = \max\{f_{[j],i+1} + p_{[j],i}, f_{[j+1],i-1}\},\
$$

$$
j = np - 1, ..., 1; i = m, ..., 2,
$$
 (20)

$$
f_{[i],1} = f_{[i],2} + p_{[i],1}, j = np - 1, ..., 1.
$$
 (21)

- **Step 3:** Use Eqs. 12-16 to calculate the departure times $d_{[q],i}$ (*i* = 1, ..., *m*) of job $J_{j'}$ when it were to be inserted at position q in the assigned factory in the
- current partial solution. **Step 4:** Use the following equation to calculate the global makespan of the partial sequence $\prod_{P_q}^*$ when job J_j were to be inserted at position *q* in the assigned

factory in the current partial solution. $C_{\text{max}}(\Pi^*_{n})$

$$
C_{\max}(\Pi_{P_q}^*) = \max_{i=1,\dots,m} (d_{[q],i} + f_{[q],i}),
$$

$$
q = 1, \dots, np + 1.
$$
 (22)

Step 5: Choose the optimal insertion position, among the best ones across all factories, which minimizes the global makespan.

IV. COMPUTATIONAL RESULTS AND DISCUSSION

A. TEST PROBLEMS

To verify the effectiveness of the three versions of the HIG algorithm, the benchmark problem set that was presented by Naderi and Ruiz [3] for testing the DPFSP was used. The benchmark problem set was augmented using the 120 benchmark instances of Taillard [38], where the processing time $p_{i,i}$ ($j = 1, \ldots, n; i = 1, \ldots, m$) is an integer that is generated from the uniform distribution [1, 99]. Naderi and Ruiz [3] expanded these 120 test instances to 420 and 720 test instances in the small and large problem sets, respectively.

The instances in the small problem set featured the number of jobs $n = \{4, 6, 8, 10, 12, 14, 16\}$, the number of machines $m = \{2, 3, 4, 5\}$, and the number of factories $f = \{2, 3, 4\}$. The total number of combinations of distinct numbers of jobs, machines, and factories was 84, which was therefore the number of sub-problem sets. Five instances were generated for each sub-problem set, yielding a total of 420 (84 \times 5) instances in the small problem set. The instances in the large problem set were those in 72 sub-problem sets, featuring the number of jobs $n = \{20, 50, 100, 200, 500\}$, the number of machines $m = \{5, 10, 20\}$, and the number of factories $f = \{2, 3, 4, 5, 6, 7\}$. Ten instances were generated for each sub-problem set, yielding a total of 720 (72 \times 10) problem instances. The files of these test instances can be downloaded from http://soa.iti.es.

B. PARAMETER CALIBRATION

The proposed three versions of the HIG algorithm have seven parameters, which are $T_{Value}, I_{iter}, \lambda, TLT_R^{low}, TLT_R^{high}, \alpha_{min}$,

TABLE 1. Parameter values used in the two-parameter calibration experiments.

*: The best combination is bold face

TABLE 2. Performance comparisions on taillard's benchmark problems (Ave. RPD_{BKS}) for $t = 15$.

Problem size	HIG ₁	HIG ₂	HIG ₃	RAIS	HDDE	IG
2015	0.000	0.000	0.000	0.000	0.000	0.000
20 10	0.000	0.000	0.000	0.000	0.000	0.000
20 20	0.000	0.000	0.000	0.000	0.000	0.000
5015	0.059	0.016	0.211	0.027	1.000	0.219
50 10	0.035	-0.040	0.123	0.072	0.805	0.272
50 20	-0.145	-0.075	-0.053	0.084	0.439	0.088
10015	-0.093	0.010	-0.079	0.000	2.618	0.826
100 10	-0.215	-0.099	-0.163	0.000	1.912	0.890
100 20	-0.300	-0.171	-0.142	0.021	1.602	0.740
200 10	-0.419	-0.273	-0.315	0.000	2.833	0.331
200 20	-0.661	-0.463	-0.571	0.000	1.880	0.633
500 20	-1.161	-1.017	-1.038	0.076	1.631	0.091
Total average	-0.242	-0.176	-0.169	0.023	1.227	0.341

TABLE 3. Paried T-tests on AVE. RPDBKS for $t = 15$.

and α_{max} , where T_{Value} is used to determine the initial temperature $(T_0 = T_{Value} \cdot \sum_{i=1}^{m}$ *i*=1 $\sum_{n=1}^n$ $\sum_{j=1}^{n} p_{j,i}$; *I*_{iter} denotes the number of iterations in the search at a particular temperature; λ is the coefficient that directs the cooling schedule; TLT_R^{low} and TLT_R^{high} determine the minimal and maximal Tabu list tenures $(TLT^{low} = n \cdot TLT_R^{low}$ and $TLT^{high} = n \cdot TLT_R^{high}$, respectively; α_{min} and α_{max} are minimal and maximal number of jobs to be eliminated in the destruction phase.

Since the parameter values affect the computational results of the three versions of HIG algorithm, two sets of test instances are used to calibrate the parameters. In the two calibration experiments, the maximum computation time (T_{max}) to solve each selected instance was set to $10 \cdot n \cdot m$ (ms). The three versions of the HIG algorithm were executed in *C* language on a personal computer that had an Intel Core Quad CPU Q9400 @2.66 GHz processor and 20 GB of RAM. Each calibration experiment was conducted on 30 instances that are randomly produced using the same data generation

procedures as that of Naderi and Ruiz [3]. The test instances in the two sets featured the number of jobs $n = \{20, 50, 100\}$, the number of machines $m = \{5, 10, 20\}$, and the number of factories $f = \{2, 3, 4, 5, 6, 7\}$. For each combination p of parameter values, each test instance was solved five times, and the best known solution (C_m^p) lim_{min_i}) among the five replications was recorded for each test instance *i*. Then, the best known solution for each test instance $i(C_{\min_1}^{BKS})$ was obtained among all C_n^p $\lim_{m \to \infty}$. For choosing the best value for each parameter, the combination with the smallest average relative percentage rate $RPD_p = (C_{\min}^p - C_{\min}^{BKS})/C_{\min}^{BKS} \times 100\%$ was chosen.

In the first calibration experiment, several values of the parameters T_{Value} , I_{iter} , λ , TLT_R^{low} , TLT_R^{high} , α_{min} and α_{max} (see Table 1) were tested for the calibration and tuning of parameter combinations. In the second calibration experiment, the results of the first experiment were refined by adding additional values, which are presented in Table 1. As shown in Table 1, the following values were obtained using the optimal parameter combinations; in the first calibration experiment, $T_{Value} = 0.03$, $I_{iter} = 3000$, $\lambda = 0.9$, $TLT_R^{low} = 5\%, TLT_R^{high} = 10\%, \alpha_{\min} = 3 \text{ and } \alpha_{\max} = 6,$ while in the second calibration experiment (after refinement), $T_{Value} = 0.03, I_{iter} = 3500, \lambda = 0.915, TLT_R^{low} = 5\%,$ $TLT_R^{high} = 10.0\%$, $\alpha_{min} = 3$ and $\alpha_{max} = 6$. Three values of TLT^{low}_{R} , 5.0%, 7.5%, and 10.0%, are tested for use in HIG₂ algorithm. Because setting TLT_R^{low} to 7.5% yielded the best

TABLE 5. Ave. $\mathsf{RPD}_{\mathsf{LB}}$ values obtained using HIG₁, HIG₂, and HIG₃ algorithms grouped according to n and m for $t = 5$, 10, and 15 (large problem set).

result, TLT_R^{low} was fixed at this value in the HIG₂ herein. Using the above parameter settings, the maximal computational time was set to $t \cdot n \cdot m$ (ms) for various problem sets, where *t* is a scale factor that was set to (5, 10, 15, 20, 25, 30) and (5, 10, 15) for the single-factory and multi-factory problem, respectively. Each problem was executed five runs, and the best solution from the five replications was recorded. The following subsections present and discuss the computational results.

C. RESULTS OBTAINED USING PROBLEM SET OF SINGLE-FACTORY

To confirm the effectiveness of the three versions of the HIG algorithm, their performance was compared with

TABLE 6. Ave. $\mathit{RPD}_{\mathsf{LB}}$ values obtained using HIG₁, HIG₂, and HIG₃ algorithms grouped according to n and m for $t = 5$, 10, and 15 (large problem set).

	HIG.	HIG ₂	HIG ₃
2	1.102/0.280/0.118	1.244/1.244/0.224	1.163/0.279/0.117
3	0.941/0.293/0.143	0.974/0.974/0.179	0.934/0.314/0.150
$\overline{4}$	0.714/0.256/0.137	0.808/0.808/0.177	0.800/0.302/0.162
5	0.656/0.260/0.140	0.687/0.687/0.172	0.646/0.276/0.155
6	0.611/0.265/0.141	0.619/0.619/0.182	0.636/0.282/0.151
	0.660/0.320/0.196	0.499/0.499/0.102	0.688/0.363/0.241
Total Average	0.781/0.279/0.146	0.805/0.805/0.173	0.811/0.303/0.163

that of leading algorithms using the 120 single-factory benchmark instances of Taillard [38]. These leading algorithms are HDDE [23], IG [24], and RAIS [25], all of which were developed to solve the single-factory BFSP. Notably, the $HIG₃$ algorithm proposed in this study is an improved version of the IG algorithm, which is a state-ofthe-art algorithm, proposed by Lin et al. [9] for the distributed permutation flowshop scheduling problem. In the literature, HDDE is performed with ten replications, while IG and RAIS are performed with five replications. Therefore, the proposed $HIG₁$, $HIG₂$, and $HIG₃$ algorithms were conducted with five replications, and the relative percentage deviation (*RPDBKS*) of makespan from the best solution that was calculated using the following formula was used to compare the performance of these algorithms with that of HDDE, IG, and RAIS.

$$
RPD_{BKS} = (C_{\text{max}}^{alg} - C_{\text{max}}^{BKS}) / C_{\text{max}}^{BKS} \times 100\%
$$

where C_{max}^{alg} is the makespan value in the best solution that was obtained using the algorithm of interest, and C_{max}^{BKS} is the makespan value in the best solution that was obtained using HDDE, IG, and RAIS.

FIGURE 2. The average RPD_{BKS} values under different t vaules.

To reveal how the computational time affects quality of the solutions obtained using $HIG₁$, $HIG₂$, and $HIG₃$, various *t* values for these algorithms were tested. Figure 2 plots the total average *RPDBKS* values for the solutions that were obtained using the $HIG₁$, $HIG₂$, and $HIG₃$ algorithms with various *t* values. Evidently, as can be seen in Fig. 2, solution

TABLE 7. Paired t-tests on Min. RPD_{BKS}, Mean RPD_{BKS}, and Max. RPD_{BKS}.

$HIG1$ vs.	HIG ₂	HIG ₃
Test on Min. $RPD_{\scriptscriptstyle RKS}$		
Paired difference	$-0.024/-0.027/-0.027$	$-0.031/-0.023/-0.017$
t -value	$-2.102/-2.580/-2.638$	$-2.898/-2.592/-1.844$
Degree of freedom	719	719
P -value	0.018/0.005/0.004	0.002/0.005/0.033
Test on Mean $RPD_{\scriptscriptstyle RKS}$		
Paired difference	$-0.020/-0.021/-0.024$	$-0.025/-0.018/-0.019$
t -value	$-2.627/-2.976/-2.875$	$-3.528/-2.833/-3.015$
Degree of freedom	719	719
P -value	0.004/0.002/0.002	0.000/0.002/0.001
Test on Max. $RPD_{\scriptscriptstyle RKS}$		
Paired difference	$-0.017/-0.010/-0.017$	$-0.022/-0.012/-0.015$
t -value	$-1.618/-0.879/-1.618$	$-2.098/-1.292/-1.620$
Degree of freedom	719	719
P-value	0.053/0.190/0.053	0.018/0.098/0.053

quality increases with computation time. When *t* is equal to or larger than 15, the total average *RPD_{BKS}* values of the solutions that were obtained using of $HIG₁$, $HIG₂$, and $HIG₃$ algorithms are negative. Accordingly, $HIG₁$, $HIG₂$, and HIG_3 outperform HDDE, IG, and RAIS if t is equal to or greater than 15. Therefore, considering both of solution quality and computational efficiency, $t = 15$ is used in subsequent analysis of the 120 single-factory benchmark instances.

Table 2 lists the average *RPDBKS* (Ave. *RPDBKS*) value for the solutions of each problem size that are obtained using $HIG₁$, $HIG₂$, $HIG₃$, RAIS, HDDE, and IG. Each average *RPDBKS* value was taken over the 10 test instances for each problem size. Tables 8 and 9 present the best solutions that were obtained using these algorithms for each benchmark instance. As shown in Table 2, the total average *RPDBKS* value of the solutions that were obtained using the $HIG₁$ algorithm was -0.242% . For the HIG₂, HIG₃, RAIS, HDDE, and IG algorithms, the corresponding values were –0.176%, –0.169%, 0.023%, 1.227%, and 0.341%, respectively. Evidently, the three versions of the HIG algorithm outperform the three leading algorithms in solving the traditional BFSP, while $HIG₁$ is the best of them. Notably, as reported by Lin and Ying [25], RAIS is better than HDDE and IG, with its maximal computational time set to $100 \cdot n \cdot m$ (ms), while the maximal computational times for the $HIG₁$, $HIG₂$, and HIG₃ algorithms are set to $15 \cdot n \cdot m$ (ms). Therefore, the three versions of the HIG algorithm take significantly less time to compute better solutions to the traditional BFSP than taken by the RAIS algorithm.

To confirm further the effectiveness of the proposed $HIG₁$ algorithm, paired *t*-tests were performed on the average RPD_{BKS} values obtained using HIG_1 and those obtained using HIG₂, HIG₃, RAIS, HDDE, and IG. The results in Table 3 that the proposed HIG_1 algorithm significantly outperforms the HIG_2 , HIG_3 , RAIS, HDDE, and IG algorithms at a confidence level $\alpha = 0.05$.

n m	HIG ₁	HIG ₂	HIG ₃	RAIS	HDDE	IG	n m	HIG_1	HIG ₂	HIG ₃	RAIS	HDDE	IG
20 5	1374	1374	1374	1374	1374	1374	50 5	2993	2989	3001	2995	3033	3002
	1408	1408	1408	1408	1408	1408		3199	3182	3200	3191	3226	3201
	1280	1280	1285	1280	1280	1280		3001	3011	3013	3007	3039	3011
	1448	1448	1448	1448	1448	1448		3126	3118	3128	3125	3147	3128
	1341	1341	1341	1341	1341	1341		3150	3154	3165	3143	3192	3166
	1363	1363	1363	1363	1363	1363		3173	3173	3179	3169	3183	3169
	1381	1381	1381	1381	1381	1381		3028	3024	3028	3021	3054	3013
	1379	1379	1379	1379	1379	1379		3059	3053	3054	3058	3081	3073
	1373	1373	1373	1373	1373	1373		2900	2906	2913	2908	2929	2908
	1283	1283	1283	1283	1283	1283		3113	3118	3107	3114	3146	3120
20 10	1698	1698	1698	1698	1698	1698	50 10	3622	3641	3629	3633	3667	3638
	1833	1833	1833	1833	1833	1833		3489	3484	3497	3487	3523	3507
	1659	1659	1659	1659	1659	1659		3481	3475	3485	3482	3515	3488
	1535	1535	1535	1535	1535	1535		3662	3651	3660	3666	3685	3656
	1617	1617	1617	1617	1617	1617		3630	3627	3638	3634	3650	3629
	1590	1590	1590	1590	1590	1590		3590	3591	3593	3576	3622	3621
	1622	1622	1622	1622	1622	1622		3694	3674	3686	3683	3704	3696
	1731	1731	1731	1731	1731	1731		3562	3567	3567	3574	3590	3572
	1747	1747	1747	1747	1747	1747		3530	3517	3534	3541	3556	3532
	1782	1782	1782	1782	1782	1782		3619	3625	3621	3616	3642	3624
20 20	2436	2436	2436	2436	2436	2436	50 20	4502	4492	4499	4504	4516	4500
	2234	2234	2234	2234	2234	2234		4275	4283	4271	4291	4296	4276
	2479	2479	2479	2479	2479	2479		4261	4271	4269	4279	4290	4289
	2348	2348	2348	2348	2348	2348		4349	4354	4353	4368	4393	4377
	2435	2435	2435	2435	2435	2435		4262	4277	4263	4275	4284	4268
	2383	2383	2383	2383	2383	2383		4289	4276	4297	4283	4308	4280
	2390	2390	2390	2390	2390	2390		4302	4313	4310	4315	4325	4308
	2328	2328	2328	2328	2328	2328		4313	4314	4318	4319	4337	4326
	2363	2363	2363	2363	2363	2363		4313	4307	4312	4313	4332	4316
	2333	2333	2333	2333	2333	2333		4401	4410	4415	4419	4439	4428

TABLE 8. Results for the instances with $n = 20$, 50 and 100 for $t = 15$.

"Bold font means the best solution among various algorithms.

D. RESULTS OBTAINED USING SMALL PROBLEM SET OF MULTI-FACTORY

For the small problem set, the relative percentage deviation (*RPDBKS*) of makespan from the lower bound that is calculated using the following formula, was used to compare $HIG₁$, $HIG₂$ and $HIG₃$ algorithms with the proposed MIP mathematical model, in terms of solution quality.

$$
RPD_{LB} = (C_{\text{max}}^{alg} - C_{\text{max}}^{LB})/C_{\text{max}}^{LB} \times 100\%
$$

where C_{max}^{alg} is the makespan value of the best solution that is get using a given version of the HIG algorithm or by solving the proposed MIP mathematical model, and C_{max}^{LB} is the lower bound on the makespan value that is get by solving the proposed MIP mathematical model.

The MIP mathematical model was solved using a famous mathematical programming solver, Gurobi (Version 7.0), on a personal computer with an Intel Core Quad CPU Q9400 @ 2.66 GHz processor and 20 GB of RAM. The maximal computational time for each test instance was set to an elapsed CPU time of 3600 seconds. The final incumbent solution that was obtained by the Gurobi MIP solver was recorded as the feasible solution. The difference between the feasible solution and the lower bound is known as the gap; a gap of zero reveals that the solution is optimal.

Table 4 lists the statistical results concerning the average *RPDLB* (Ave. *RPDLB*) values that were obtained using the

small problem set using the MIP mathematical model, and the $HIG₁$, $HIG₂$, and $HIG₃$ algorithms. The three *t* values are separated by a slash $(t = 5/10/15)$, except in the result obtained using the MIP mathematical model, which includes only one Ave. *RPDLB* value. It should be noted that, because the number of jobs is smaller in these problems, the maximal number of jobs to be removed is set to *n*/2. As revealed in Table 4, the total average *RPDLB* values of the solutions that were obtained using $HIG₁$, $HIG₂$, and $HIG₃$ algorithms are smaller than those obtained using the MIP mathematical model. The MIP mathematical model found optimal solutions for all benchmark instances when the number of jobs did not exceed 10. When the number of jobs in benchmark instances was 12, 14, or 16, the MIP mathematical model obtained optimal solutions in 51, 18, and one out of 60 test instances, respectively. In total, the MIP mathematical model obtained optimal solutions in 310 out of 420 benchmark instances in the small problem set. Notably, all except one of the optimal solutions that were obtained using the MIP mathematical model were also obtained using $HIG₁$, $HIG₂$, and $HIG₃$ algorithms. Furthermore, the computational times required for $HIG₁$, $HIG₂$, and $HIG₃$ algorithms are much less than that of the MIP mathematical model. These analytical results confirm that the proposed $HIG₁$, $HIG₂$, and HIG³ algorithms exhibit excellent convergence to optimal solutions.

*Bold font means the best solution among various algorithms.

E. RESULTS OBTAINED USING LARGE PROBLEM SET OF MULTI-FACTORY

Because of the complexity of the $DF_m|block|C_{\text{max}}$ problem, a high-quality feasible solution to a large problem cannot be obtained using the proposed MIP mathematical model in a reasonable computation time. Therefore, for the test instances in the large problem set, the relative percentage deviation (*RPDBKS*) of makespan from the best solution obtained using the $HIG₁$, $HIG₂$, and $HIG₃$ algorithms was used to compare the solutions obtained using the $HIG₁$, $HIG₂$, and $HIG₃$ algorithms in terms of quality.

Tables 5 and 6 present the statistical results concerning the average *RPDBKS* values of 60 and 120 solutions, respectively, in the large problem set obtained using the three versions of the HIG algorithm; the results that were obtained using the three *t* values are separated by a slash $(t = 5/10/15)$. The statistical results reveal that more jobs yield a larger average *RPDBKS* (Table 5) and more factories are associated with a smaller average *RPD_{BKS}* (Table 6). Thus, the number of jobs and the number of factories in the benchmark instances affected the performance of the $HIG₁$, $HIG₂$, and $HIG₃$ algorithms, whereas the number of machines did not.

The best, mean, and worst makespan values of the solutions to each test problem, based on five trials, obtained using each version of the HIG algorithm, were used to compute *RPDBKS*

values, which were denoted as Min. *RPDBKS* , Mean *RPDBKS* , and Max. RPD_{BKS} . To determine whether the HIG_1 algorithm was better than the $HIG₂$ and $HIG₃$ algorithms, one-sided paired *t*-tests in terms of Min. *RPDBKS* , Mean *RPDBKS* , and Max. *RPDBKS* were performed for the various *t* values $(t = 5/10/15)$. The statistical results listed in Table 7 revealed that, at a confidence level of $\alpha = 0.05$, the proposed HIG₁ algorithm significantly outperformed the HIG_2 and HIG_3 algorithms in terms of Min. *RPDBKS* , Mean*RPDBKS* , and Max. *RPDBKS* , for most of the *t* values. These statistical results confirm that adopting the variable Tabu list operator of TS and the cooling schedule operator of SA significantly improves the performance of IG in solving the *DFm*|*block*|*C*max problem.

V. CONCLUSIONS AND RECOMMENDATIONS FOR FUTURE RESEARCH

The $DF_m|block|C_{\text{max}}$ problem is a variant of the DPFSP. Owing to the broad applications of distributed blocking flowshop systems, this work developed three versions of the HIG algorithm ($HIG₁$, $HIG₂$, and $HIG₃$) for solving the $DF_m|block|C_{\text{max}}$ problem to bridge the gap between theoretical progress in DBFSP and the industrial implication of distributed blocking flowshop systems. A comprehensive benchmark problem set is used to test the effectiveness and efficiency of the $HIG₁$, $HIG₂$, and $HIG₃$ algorithms.

In summary, the computational results that are presented in this work are very encouraging for the application of the HIG1 algorithm to the *DFm*|*block*|*C*max problem. The proposed $HIG₁$ algorithm exploits the PW and NEH2 heuristics to develop an initial schedule and combines the IG algorithm with the operators of the variable Tabu list and the cooling schedule. A highly effective speed-up method for evaluating of the best insertion position is used to reduce the computational burden. In view of the current lack of meta-heuristics to solve the $DF_m|block|C_{\text{max}}$ problem, this work provides an important basis to exploring this significant topic.

Many issues are worthy of further study in the area of this pioneering study. First, additional exact methods and meta-heuristic algorithms should be developed to solve effectively and efficiently the *DFm*|*block*|*C*max problem. Second, the proposed HIG algorithms could be modified to solve the distributed no-idle flowshop scheduling problem. Third, the distributed blocking flowshop scheduling problem that involves other sophisticated objectives is worthy of research. Fourth, the extension of the present consideration of the $DF_m|block|C_{\text{max}}$ problem to consider the multiobjective distributed blocking flowshop scheduling problem would increase the application of scheduling theory in industry. Finally, the novel theoretical research should be expanded from the deterministic *DFm*|*block*|*C*max problem to stochastic problems.

APPENDIX

See Tables 8 and 9.

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