

# A Two-stage Genetic Algorithm for a Novel FJSP with Working Centers in a Real-world Industrial Application

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**Abstract:** Inspired by industrial issues and demands, we define a novel version of the Flexible Job Shop Scheduling Problem with Working Center. A working center is a group of machines performing the same type of operation. The job operations of different types follow a strict sequence across the working centers, while any order is allowed among operations of the same type. This paper illustrates a genetic algorithm with a two-stage chromosome representation, adapted genetic operators, local search, and social disaster technique to deal with a real-world industrial application. The algorithm has been tested on a classical benchmark to assess its adaptability and compare its performance with state-of-the-art techniques; then, we tested different variations of the proposed algorithm on a real-case test instance showing a consistent improvement when compared with the heuristic in use at the industrial company.

## 1 INTRODUCTION

Production scheduling is one of the most critical tasks in manufacturing systems and it has been extensively studied by the scientific community. The problem is of more than academic interest. The effective scheduling of operations processed by the shop floor reduces the working-process inventory and increases the throughput with a positive impact on the performances. The scheduling problem is concerned with allocating available production resources to tasks and determining the sequence of operations that can optimize business metrics. In literature, the job shop scheduling problem (JSP), which is NP-hard, is the standard formulation of the production scheduling problem. In the JSP, an operation can be performed by only one machine and a machine can perform a single operation at a time. However, to cope with a more complex real case scenarios, the JSP has been extended by allowing an operation executable on more than one machine (Flexible JSP), on all the machines (Total Flexible JSP), or on a subset of the machines (Partial Flexible JSP) (Xie et al., 2019).

A peculiar and practice-oriented specification of the FJSP is the one described by Behnke and Geiger. In their work, they illustrate a context where similar machines are pooled to working centers. In this setting, if an operation is assignable to a machine it is also assignable to any other machines belonging to the same working center. Another interesting property of the context is that the sequence of operations is fixed and the first and the last operations of each job must be processed by machines from the first and last working center (Behnke and Geiger, 2012).

The latter scenario comes close to that encountered during our research work, conducted with an Italian company. Here working centers aggregate machines performing a certain type of operation. Moreover, while the sequence of the operations from one working center to the next is fixed, the operations of the same working center can be performed in any possible order. The company had shown difficulties in finding an optimized scheduling strategy that would allow to comply with such constraints, not commonly studied in the FJSP field, and better distribute the working load. In order to answer this necessity, in this paper we introduce an extended variation of the FJSP with Working Center proposed by Behnke and Geiger and illustrate an adapted genetic algorithm with a two-stage chromosome representation, local search and social disaster technique.

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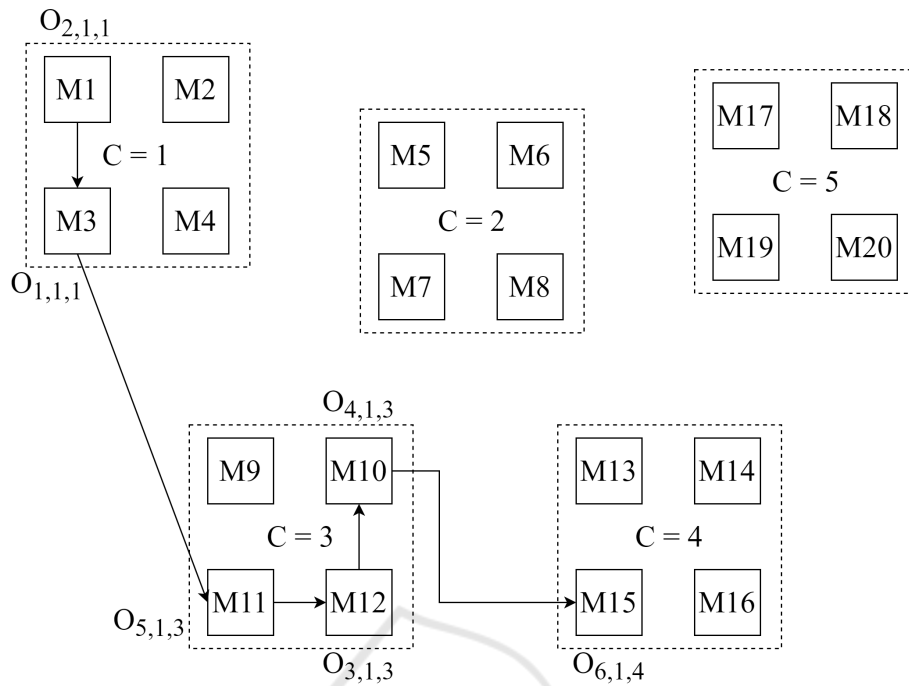


Figure 1: Possible operation sequence through working centers for a job  $j = \{o_{1,1,1}, o_{2,1,1}, o_{3,1,3}, o_{4,1,3}, o_{5,1,3}, o_{6,1,4}\}$ .

Our contribution consists of:

- the formalization of an industrial problem as a variation of the FJSP;
- the implementation of an adapted encoding/decoding system that makes a GA able to handle solutions complying with the presented constraints without performing corrections;
- the development of a scheduling strategy that improves human performances in an industrial context.

Population-based meta-heuristics have been indeed widely applied for solving the FJSP, with the genetic algorithm (GA) being the most popular among them. The algorithm integrates different strategies for generating the initial population, selecting the individuals for reproduction and reproducing new solutions (Xie et al., 2019). Genetic algorithms encode problems for which they aim at optimizing a cost function in chromosomes: the individuals of the population are vectors whose elements are called genes. When dealing with the FJSP and its relative subclasses, it is common practice to represent the problem using a two-sides chromosome. A chromosome can indeed be split in two parts that behave differently and encode different information: one encodes the machine assignment, while the other encodes the operation sequencing (Gao et al., 2008) (Yang et al., 2009) (Zhang et al., 2011) (Rahmati and Zandieh, 2012) (Defersha and Rooyani, 2020). In this work,

we extend the typical pipeline of the genetic algorithm with *local search* and *social disaster technique*. Local search can help avoid local minima by exploring the local space around a particular solution, possibly improving it. It can be used at different stages of the global search (Yun et al., 2013) or directly at the end of the global search process (Nouri et al., 2018). The second technique we use is called *social disaster technique*. The general idea is to diagnose the situation of loss of genetic diversity of the population, and in such case to apply a catastrophic operator to it. These operators have the purpose to return the population to an acceptable degree of genetic diversity, by replacing a number of selected individuals with others generated at random (Rocha and Neves, 1999). Both techniques are designed to help the model provide better solutions avoiding premature convergence to local optima, avoiding the waste of computational time around fruitless areas of the search space.

The structure of the paper is as follows: After detailing each constraint and assumption of our problem in Section 2, we illustrate chromosome representation, gantt building, the entire pipeline and each feature in Section 3. In Section 4 we discuss performances sustained on a test instance for the FJSP and on a test instance retrieved in our use case. Finally, in Section 5 we discuss obtained results, possible additional features and further improvements.

## 2 PROBLEM DEFINITION

Our problem can be described as a set of  $n$  jobs  $J = \{1, 2, \dots, n\}$  where each job  $j$  consists of a number of operations  $P_j = \{1, 2, \dots, p\}$ . Every operation is traceable to a class of operations called *type*  $k \in K$  with  $K = \{1, 2, \dots, l\}$ . Thus, we define the  $p$ -th operation of job  $j$  of type  $k$  as  $o_{p,j,k} \in O$ , where  $O$  is the set containing the operations of all jobs. The operations  $O$  are to be processed on a set of machines  $M$ , divided into subsets of machines called *Working Centers*  $c \in C = \{1, 2, \dots, w\}$ , where  $w = l$ . All machines grouped in a working center have the property to process only one type of operations  $k$ , making it possible to draw a one-to-one correspondence between the type of operations  $K$  and working centers  $C$ . Finally, a machine  $m_c$  belongs to one and only one working center  $c$  of the shop floor  $M$ . A machine  $m_c$  is described by a set of features that identify the capability of the machine to perform certain operations. Hence, there is no guarantee that a machine  $m_c$  can perform all operations of type  $k$ . The mapping between an operation and the subset of machines that are able to process it is defined by our real-world casuistry, where operations of the same type can widely differ from each other over several parameters and is not the object of this work. The sequence of operations for each job is not predetermined. The only constraint is imposed by the operation type. In fact, for each job, all the operations of type  $k$  have to be performed before any operation of type  $k + 1$ . Also, jobs are not necessarily characterized by the presence of operations belonging to all  $K$  types (*i.e.*, in Figure 1 there is no operation on machines where  $c = 2$ ), while the aforementioned sequence constraint is always valid. Figure 1 shows the assignment of the operations of job  $j = \{o_{1,1,1}, o_{2,1,1}, o_{3,1,3}, o_{4,1,3}, o_{5,1,3}, o_{6,1,4}\}$  to the working centers.

We aim at giving an ordered set of couplings as solution, *e.g.*,  $\langle (o_{2,1,1}, m1_1), (o_{1,1,1}, m3_1), (o_{5,1,3}, m11_3), (o_{3,1,3}, m12_3), (o_{4,1,3}, m10_3), (o_{6,1,4}, m15_4) \rangle$ , that associates each operation of each job to one of the available machines. A coupling  $(o_{p,j,k}, m_c)$  has a cost in terms of time  $t$ , which is the processing time of the operation on the machine. Different machines offer different performances for the same operations, thus resulting in different processing times. Newer machines, indeed, are expected to outperform older and consumed machines. Once retrieved where and when an operation is going to be processed, and having a forecast of the cost of the operation itself, we can build a gantt chart, as we shall detail in Section 3.1, to better visualize the solution and to eventually calculate the total cost of the solution itself.

Finally, we adopt *makespan* as the indicator to optimize while constructing suitable solutions for the described problem. Makespan is commonly defined as  $C_{max}$ : the completion time of the last operation of the last job in the system (Türkyılmaz et al., 2020).

To sum up, our problem can be summarized by the following constraints and assumptions:

- All jobs are available at release dates;
- Not all machines are necessarily available at time zero;
- Each operation can only be processed on one machine at a time;
- Each machine can only perform one operation at a time.
- Each operation cannot be interrupted during the processing process;
- Machines are grouped in Working Centers;
- Machines of each Working Center can only perform operations of the corresponding type  $k$ ;
- Not all machines belonging to a Working Center can perform all operations of the corresponding type  $k$ ;
- Operations can require a different processing time on different machines;
- There is no precedence constraint among jobs;
- For each job, the sequence of operations is constrained by operation types;
- There is no precedence constraint among operations of the same type.

## 3 PROPOSED APPROACH

In order to return suitable solutions in a real industrial context, according to constraints outlined in Section 2, we implement a dual-stage genetic algorithm with an adapted encoding-decoding chromosome representation and pipeline, local search and social distasters technique.

### 3.1 Chromosome Representation

The problem presented can be divided into two sub-problems: the machine assignment and the operation sequencing. For this reason, we utilize a two-sides chromosome to represent a possible solution to the problem: the left segment encodes the machine assignment while the right segment encodes the operation sequence (Figure 2).

For the sake of clarity, we start by detailing the right portion of the chromosome which represents the operation permutation. The right segment has  $n$  genes, each one representing one element of the operation sequence to schedule. The index  $i$  identifies a gene, while the value of the gene identifies an operation  $p$ . The operations are sorted by their respective type of operation, thus each operation of type  $k$  has an index  $i$  smaller than each operation of type  $k + 1$ . For instance, in Figure 2, operations with  $K = 1$  are given ids in the range [1, 2]. This way the grouping for operation types also determines  $s$  sub-portions of the chromosome, where  $s \leq l$ , where operations have no predefined sequence and guarantee the respect of sequence constraints. In Figure 2, for example, we have three operations of type  $K = 3$ ; these operations are assigned ids [3, 4, 5], hence are identified in the chromosome as the genes with values 3, 4 and 5. Moreover, these operations can occupy positions in the right vector at indexes 3, 4 and 5. It is to be noted how Figure 1 and Figure 2 only refer to a single job for convenience, while generally a chromosome represents a complete solution composed of multiple jobs.

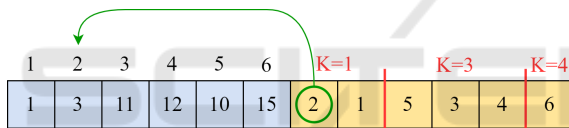


Figure 2: The chromosome representation of the solution described in Figure 1.

The left segment of the chromosome has  $n$  genes, each one indexed by operation  $p$ . The value of the  $p$ -th gene is the machine  $m_c$  assigned to perform the operation  $o_{p,j,k}$ . To sum up, the left segment of the chromosome provides the couplings  $(o_{p,j,k}, m_c)$ , i.e., the assignment of the operation to a machine, while the right segment encodes the order of the coupling in the solution.

We have seen how each id used as the value for genes of the right segment is eventually traceable to an operation  $o_{p,j,k}$ . In the left segment of the chromosome, each gene represents the id of the machine to which we assign the corresponding operation. This correspondence is mapped by the vector index of the gene, physiologically ordered from 1 to  $n$ . They indeed correspond to the ids we assigned to the operations in the right segment. Figure 2 shows graphically how a gene in the right segment of the chromosome identifies an operation for which the corresponding gene in the left segment of the chromosome reports the assigned machine. As for the example, the operation with id 2 is assigned to machine 8.

### 3.2 Gantt Building

To evaluate each generated solution we need to be able to decode and organize them under the form of a gantt chart. By iterating over the right segment of the chromosome we extract information about which operations must be inserted first into the gantt and on which machine, using the correspondences seen in Section 3.1. Then, for each operation, we set the earliest possible starting time just after the finish time of the latest operation of the same job already put in production.

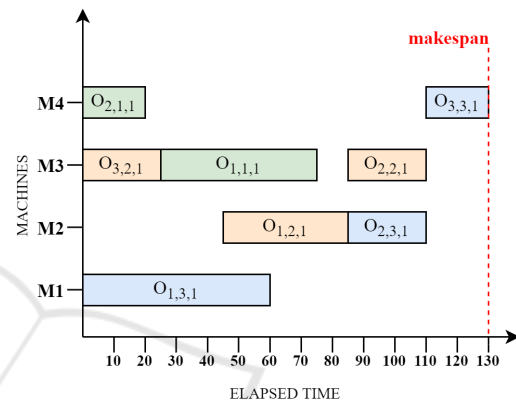


Figure 3: Example of a gantt chart utilized to evaluate the makespan.

In fact, for each operation we need to check when the last operation belonging to the same job terminated its processing and then check on the queue of the assigned machine; here we look for the earliest point where we can insert the operation without generating overlappings with other operations already placed on the machine (Demir and İşleyen, 2014). The procedure seen as a pseudo-code is:

```

FOR op in chromosome
  op_start = 0
  op_finish = 0

  prev_op = eventual latest operation
              of the same job

  IF prev_op not empty:
    op_start = prev_op_finish

  op_finish = op_start + op_exec_time

FOR i scheduled on op_machine:
  IF op_start >= i_finish:
    move forward in loop
  IF op_finish > i_start:
    op_start = i_finish
    op_finish = op_start +
                op_exec_time
    
```

Once the gantt chart is built we can use it to calculate  $C_{max}$ . In Figure 3 processing time is represented on the X-axis and is expressed in minutes; for three hypothetical jobs, we can see the positioning of their operations over time and space (machine assignment) and identify  $C_{max}$  of the schedule as 130 minutes.

### 3.3 Genetic Operators

Our two-stage representation of the problem deviates from the classical pipeline of the genetic algorithm to respect the dual nature of our chromosome representation and to introduce some variations.

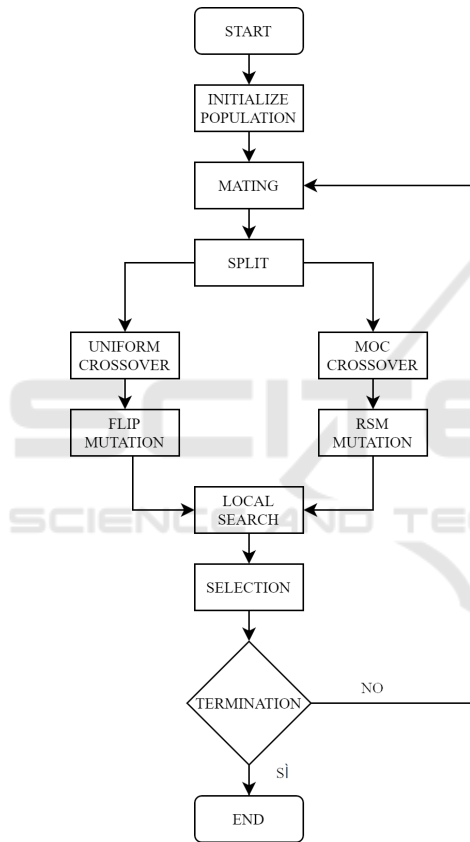


Figure 4: Our adjusted pipeline for the genetic algorithm.

We initialize the population by sampling  $N$  random suitable chromosomes, with  $N$  being a product of the length of the chromosome and a  $pop\_factor$  parameter:

$$N = 2 \cdot L \cdot pop\_factor \quad (1)$$

To emphasize the distance between starting individuals and cover wider portions of the search space without exploding the population size, we force the sampling process to generate distant chromosomes. To do so we introduce a dissimilarity distance  $Dist$  that can

measure the difference between two chromosomes. In the machine assignment segment, we verify whether the two chromosomes differ gene by gene. When they do,  $Dist$  is incremented by the number of alternative machines available to process the operation identified by the gene in question. In the operation sequencing segment, we increase  $Dist$  by 1 for each gene in which the two chromosomes differ.  $Dist_{max}$  is the maximum distance retrievable as defined by Nouri et al (Nouri et al., 2018). When sampling each chromosome we check its distance with all other individuals in the population. If  $Dist > Dist_{max} \cdot 0.5$  we add the chromosome to the initial population. This procedure ensures a good level of variability before starting the global search. We evaluate and sort each generated chromosome by  $C_{max}$ . Then we enter the crossover procedure by checking how many chromosomes are going to be mated, according to the  $Crossover Rate$ , and randomly select and mate chromosomes. At this point, our pipeline deviates and we split each chromosome into two segments, as seen in Section 3.1. Both segments of the chromosomes need to be treated separately and undergo different crossover and mutation procedures.

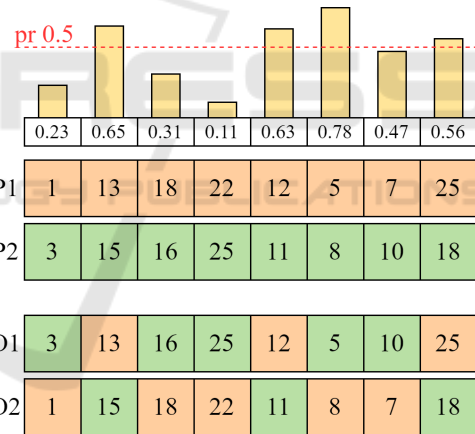


Figure 5: Uniform Crossover.

For the left segment of the chromosomes, parents undergo the crossover procedure with a *Uniform Crossover*. In Uniform Crossover two offspring are generated by combining genes of the two parents, following the direction impressed by a probability  $pr$ . Iterating in parallel over each gene of both parents, if  $pr > 0.5$  offspring 1 takes the gene from parent 1 while offspring 2 takes the gene from parent 2 and vice versa when  $pr < 0.5$  (Magalhaes-Mendes, 2013), as shown in Figure 5. Uniform Crossover always generates suitable chromosomes according to our constraints. The same can be said for the *Flip Mutation*, the chosen mutation operator. Offspring are mutated with a probability indicated as *Mutation Rate*, while



another parameter *Flip Rate* expresses the percentage of genes to flip. The flipping of each gene happens to respect the domain of the alternative machines available to process the operation identified by the gene in question, trying to adopt, when possible, a different value from the original one (Kala, 2016). For the right

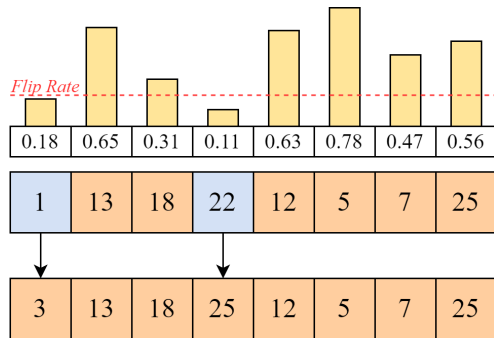


Figure 6: Flip Mutation.

portion of the chromosomes, parents undergo the crossover procedure with a *Modified Order Crossover (MOC)* operator. MOC Crossover constructs an offspring by choosing a number of genes from one parent and preserving the relative order of the elements of the other parent and a second offspring by repeating the procedure with reverted roles between the two parents (Umbarkar and Sheth, 2015). *MOC Rate* determines the number of genes to preserve as a percentage of the chromosome length. Given the nature of the chromosomes at hand, this crossover operation is to be performed separately for each of the *s* sub-portions of the chromosome defined in Section 3.1. Offspring are

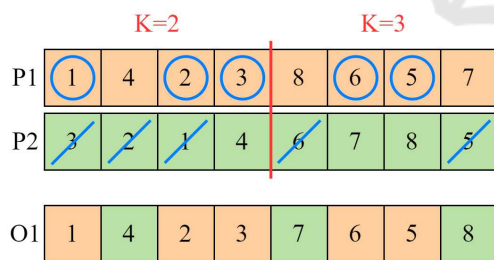


Figure 7: MOC Crossover.

mutated with the same probability determined by the Mutation Rate by a *Reverse Sequence Mutation* operator. This mutation operator reverses the order of a sub-vector of random length from the chromosome at hand (Abdoun et al., 2012). As for the crossover operation, this mutation operation is performed separately for each of the *s* sub-portions of the chromosome. Both the strategies adopted for the two respective segments are chosen to comply with the machine assignment and operation permutation constraints, in order to avoid the necessity to operate corrections on

generated offspring as well as overcoming and improving some solutions seen in literature (Moon et al., 2002) (Yun et al., 2013). Moreover, new solutions inserted into the population are always checked not to be already present in the population to avoid damaging repetitions.

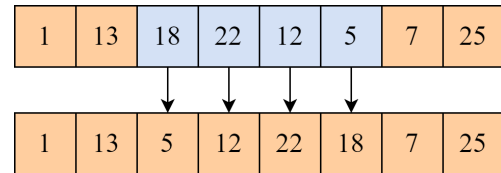


Figure 8: Reverse Sequence Mutation.

After crossovers and mutations are performed, the two segments of the offspring are recombined together. At this point, all offspring are evaluated and added to the initial population. Before moving forward to the following generation, the extended population is cut back to its original size by selecting the *N* best individuals, granting them a presence into the next generation (Chudasama et al., 2011).

### 3.4 Additional Features

In our model, we incorporate two additional features that are not part of the typical pipeline of the genetic algorithm: *local search* and *social disaster technique*. For the former, we utilize a Hill-climbing algorithm as seen in (Ceylan, 2006). At each generation we run the local search on two solutions: the best offspring and a randomly picked chromosome (Wan and Birch, 2013). The algorithm explores the local space around the interested solution by changing it one bit at a time. At each step, the local search algorithm changes the machine assignment segment by flipping a randomly picked gene and the operation permutation segment by swapping two randomly picked genes. If the produced solution improves the starting solution, the algorithm stops and the starting solution is updated into the population; otherwise, the algorithm performs another step until the maximum number of available steps *max step* is reached. For the latter, we use a parameter named *counter*, which is initialized at 1, and we introduce a hyper-parameter *patience*. We check after each generation whether we have improved the best solution. If it has not improved we increase the counter by 1. If it does improve we reset the counter to 0. If *counter == patience* we remove the worst *N/2* individuals from the population, randomly re-initialize an equivalent number of new individuals and reset the counter.

Table 1: Comparison of results on BCdata.

Instance	proposed GA		Rahmati et al., 2012		Nouri et al., 2018		Best	
	$C_{max}$	Avg $C_{max}$	$C_{max}$	Avg $C_{max}$	$C_{max}$	Avg $C_{max}$	$C_{max}$	Difference
<i>mt10c1</i>	930	942,4	946	947	<b>927</b>	930	927	0,32%
<i>mt10cc</i>	<b>915</b>	920	946	946	917	918.6	910	0,55%
<i>mt10x</i>	934	943,6	961	961	<b>923</b>	931.4	918	1,74%
<i>mt10xx</i>	940	945,2	945	945	<b>918</b>	924.4	918	2,40%
<i>mt10xxx</i>	934	941,4	954	954	<b>918</b>	921	918	1,74%
<i>mt10xy</i>	916	922,4	951	951	<b>908</b>	910	905	1,22%
<i>mt10xyz</i>	<b>855</b>	873,8	858	858	868	871.8	847	0,94%
<i>setb4c9</i>	970	983,8	959	959	<b>927</b>	936.6	914	6,13%
<i>setb4cc</i>	969	972	944	950	<b>938</b>	946.8	909	6,60%
<i>setb4x</i>	971	989,8	942	951	<b>944</b>	956.2	925	4,97%
<i>setb4xx</i>	967	984,8	967	967	<b>942</b>	953.6	925	4,54%
<i>setb4xxx</i>	987	992,4	991	991	<b>949</b>	958.6	925	6,70%
<i>setb4xy</i>	961	978,4	978	982	<b>931</b>	941.8	916	4,91%
<i>setb4xyz</i>	960	964,4	930	930.5	<b>926</b>	929.8	905	6,08%

#### 4 EXPERIMENTAL RESULTS AND DISCUSSION

We evaluated the proposed algorithm on two datasets: the benchmark dataset of (Chambers and Barnes, 1996) and a real-world scheduling dataset from an industrial company. The former dataset is well known and we compared our performances with other state-of-the-art approaches. For the latter dataset, we compared our approach with the scheduling heuristic currently in use at the company we collaborated with; finally, we also studied different combinations of genetic operators of the proposed algorithm.

Barnes and Chambers constructed a set of data (BCdata) from three of the most challenging classical job shop scheduling problems by replicating machines. The processing times for operations on replicated machines are assumed to be identical to the original (Chambers and Barnes, 1996). We tested the algorithm over 14 test instances by Barnes and Chambers, choosing the hyper-parameters experimentally as follows:

- *Pop factor*: 10;
- *Crossover Probability*: 0.6;
- *MOC Rate* = 0.7;
- *Mutation Rate* = 0.3;
- *Flip rate* = 0.4;

Due to the non-deterministic nature of the genetic algorithm, we ran our model 5 times on each instance over 200 generations. For each instance, we recorded the minimum and the average  $C_{max}$  found over the 5 runs. Table 1 compares our GA with

BBO algorithm of Rahmati and Zandieh (Rahmati and Zandieh, 2012), the hybrid metaheuristics-based multiagent model of Nouri (Nouri et al., 2018) and the best results obtained in literature also listing the percentage difference from our results.

Our algorithm can compete with the two proposed approaches to the point of obtaining better results on instances *mt10cc* and *mt10xyz*. However, it is to be noted that our algorithm does not converge toward the known global optimum, achieving better results for the *mt10* series, with margin ranging from 0.32% to 2.40%, than the *setb4* series, with margin ranging from 4.54% to 6.70%. This difference in the performances may be due to the unique setting of hyper-parameter which may require fine-tuning for each instance of the BCdata dataset.

For the second experimental setting, we are provided with a test instance composed of a total of 309 operations divided into 16 jobs with a mean of 19,3 operations each. The instance is anonymized so that jobs, operations and operation types are identified by numerical progressive ids. The operations are distributed among operation types as follows:

- *type 1*: 55 operations
- *type 2*: 29 operations
- *type 3*: 27 operations

The number of machines available to each operation is variable and ranges from 1 to 5. Machine features are not discussed in this work and information on alternative available machines for each operation is given in the test instance.

On this instance, we ran our algorithm 10 times over 500 generations. We varied our algorithm in four

different combinations in order to justify the choice made in terms of genetical operators and to use an additional feature such as the local search:

1. The proposed approach;
2. The proposed approach without the local search;
3. The proposed algorithm with a different set of genetic operators, namely:
  - *K-point Crossover* instead of Uniform Crossover (Umbarkar and Sheth, 2015);
  - *OX Crossover* instead of MOC Crossover (Magalhaes-Mendes, 2013);
  - *Twors Mutation* instead of Reverse Sequence Mutation (Abdoun et al., 2012);
4. The setting number 3 without the local search.

Hyperparameters are held the same for each approach and are chosen experimentally as follows:

- *Pop factor*: 1;
- *Crossover Probability*: 0.8;
- *MOC Rate* = 0.6;
- *Mutation Rate* = 1.;
- *Flip rate* = 0.2;

To resemble the schedule that would be sent to production in the original industrial environment, we used as a baseline a solution obtained with the scheduling heuristic used by the production manager, which is not different from the *min-min* heuristic as know in literature (Durasević and Jakobović, 2018). The heuristic can produce different solutions, so we ran it 1000 times to better represent the variability of human behavior. We then sorted the obtained results per  $C_{max}$ , selected the first and the fifth-best to show a range of possible results obtained by a human expert, and used them both as a baseline.

Figure 9 shows the obtained results in a line chart. The area around each line gives an idea of the variability of the results obtained, as it is delimited by the minimum and the maximum  $C_{max}$  value seen at a certain generation during the 10 runs. Both approach 1 and approach 3 perform better than the respective versions without the use of local search, namely approach 2 and 4, showing how important the effective use of a local search algorithm combined with the global search algorithm can be. The proposed approach performs better than the alternative one both in the short and in the long term: it is quicker in terms of generations to reach the fitness level obtained by the production manager’s heuristic; it is better at the 100 generations mark, being able to efficiently explore local portions of the search pace and improve rapidly; it is better after 500 generations, still being able to improve, managing to escape local minima with the use

of social disaster technique. When compared with the fifth baseline, the algorithm is able to obtain an improvement of 8.68% at the 100 generations mark and of 10% after 500 generations. When compared with the fifth baseline, the algorithm is able to obtain an improvement of 6.73% at the 100 generations mark and of 8% after 500 generations.

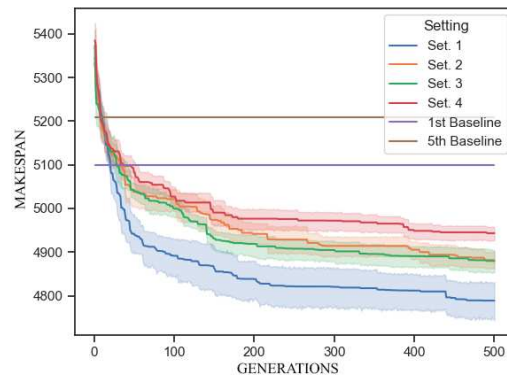


Figure 9: Comparison of results for the second experimental setting.

Eventually, our approach improves the baseline scheduling saving up to 412 minutes of working time when compared with the first baseline.

## 5 CONCLUSIONS

In this work, we presented a novel version of the Flexible Job Shop Scheduling Problem with Working Center, in which the operations have strict sequence constraints. In order to handle such constraints in an industrial environment, we developed a genetic algorithm with a two-stage chromosome representation with its encoding/decoding system, adapted genetic operators, local search and social disaster technique. To evaluate the proposed algorithm we first compared it with other approaches on a benchmark dataset obtaining good results. Finally, we tested our algorithm on a test instance provided by the company which inspired such formulation of the problem. We compared the results with the heuristic adopted by the production manager and with different settings of the proposed approach. We showed how in both analyzed settings the use of local search improves the performances and allows for quick convergence to good solutions. Also, we were able to show significant improvements against both baselines produced by a human heuristic by a margin of 8% to 10%. Further development of this work will be to dynamically handle the arrival of new operations during the scheduling process, considering attach-



ment/detachment setup time too, while accounting for computational performances, especially in comparison with other research works and with commercial products.

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