# Benchmarking Permutation Flow Shop Problem: Adaptive and Enumerative Approaches Implementations via Novel Threading Techniques 

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# BENCHMARKING PERMUTATION FLOW SHOP PROBLEM: ADAPTIVE AND ENUMERATIVE APPROACHES IMPLEMENTATIONS VIA NOVEL THREADING TECHNIQUES 

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A thesis submitted in partial fulfillment of the requirements for the

Master of Science in Computer Science

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Benchmarking Permutation Flow Shop Problem: Adaptive and Enumerative Approaches Implementations via Novel Threading Techniques
is approved in partial fulfillment of the requirements for the degree of

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## Abstract

A large number of real-world planning problems are combinatorial optimization problems which are easy to state and have a finite but usually very large number of feasible solutions. The minimum spanning tree problem and the shortest path problem are some which are solvable through polynomial algorithms. Even though there are other problems such as crew scheduling, vehicle routing, production planning, and hotel room operations which have no properties such as to solve the problem with polynomial algorithms. All these problems are NP-hard. The permutation flow shop problem is also NP-hard problem and they require high computation. These problems are solvable as in the form of the optimal and near-optimal solution. Some approach to get optimal are exhaustive search and branch and bound whereas near optimal are achieved annealing, Genetic algorithm, and other various methods.

We here have used different approach exhaustive search, branch and bound and genetic algorithm. We optimize these algorithms to get performance in time as well as get the result closer to optimal. The exhaustive search and branch and bound gives all possible optimal solutions. We here have shown the comparative result of optimal calculation for 10 jobs with varying machine number up to 20 . The genetic algorithm scales up and gives results to the instances with a larger number of jobs and machines.

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## List of Acronyms

Cmax : Makespan
MPM : Multi Purpose Machine
NIQ : No Intermediate Queues
SDST : Sequence-Dependent Set-up Times
SMT : Surface Mount Technology
PCB : Printed Circuit Board
MILP : Mixed Integer Linear Program
Ex : Exhaustive Search
BB : Branch and Bound
GA : Genetic Algorithm
BFS : Breadth First Search
DFS : Depth First Search
BeFS : Best First Search
LB : Lower Bound
UB : Upper Bound
AR : Approximation Ratio

## Chapter 1

## Introduction

The world is developing in various fields with the aid of technologies. The sequencing is one of the common problem occurring frequently. The manufacturing sector, airlines sector, banking sector and various other day to day life activities involve sequencing to great extent ${ }^{[R W C M 03]}$. In economic and industrial field flow shop can be implied to great extent.This has lured many researchers to work on problem with diverse classical assumptions and different objective functions and by implementing various optimization techniques.There are two main elements of flow shop problem: (i) M machines for job processing and (ii) N jobs to be processed on those machines. The jobs and machines have their criteria. All jobs sequence execute in same machine order. The job is not executed multiple times on same machine. A job cannot be executed on multiple machines neither a machine can execute multiple jobs. At first Conway et al. (1967) have devised the notation for flow shop with makespan criterion as $n / m / F / c_{\max }{ }^{[R W C M 03]}$ later Graham et al. (1979) introduce new notation as $F \| c_{\max }{ }^{[R . L 79]}$. The makespan criterion is defined as the total time of completing all the jobs on all machines.This problem has $(n!)^{m}$ alternative sequence of jobs over the machines. The classical flow shop problem is assumed to have a buffer or queue to hold the jobs between the machines (Allahverdi et al. $1999{ }^{[A G A 99]}$ ). Later different variants were developed.Blocking flow shop problem was stated by Abadi, Hall and Sriskandarajah (1995) [AHSO0]. Aldowaisan and Allahverdi (1998) gave a concept on no-wait flow shop problem ${ }^{[A A 98]}$. The similar concept was earlier pitched by Piehler (1960), Reddi and Ramamoorthy (1972), Bonney and Gundry (1976), King and Spachis (1980), Gangadharan and Rajendran (1993) and Rock (1984) whereas no intermediate queues (NIQ) flow shop problem was mentioned by Stafford (1988), Stafford and Tseng (1990), and Wismer (1972) ${ }^{[505]}$. The survey of Hall and Sriskandarajah (1996) aided to deduce the computational complexity for a varieties of approach to problems describing also the different
application ${ }^{[H S 96]}$. Thus it make no-wait flow shop problem a topic of interest (e.g. Bertolissi $2000{ }^{[B e r 00]}$, Aldowaisan and Allahverdi $2004{ }^{[\text {AAA04] }}$, passim). In the beginning of 2000, hybrid flow shop problems came to eyes of researchers.

The flow shop environment in real life scenario is not as stated in classical flow shop problem where set-up times are assumed to be unaffected by the job's position in sequence. Hence set-up time can be added in processing time of job. In fact the real life flow shop problem has sequencedependent set-up times (SDST). The surface mount technology (SMT) and printed circuit board (PCB) manufacturing environments have SDST flow shop problem. A mixed integer linear program (MILP) model is suggested first by Srikar and Ghosh (1986) for SDST flow shop problem ${ }^{[S G 86]}$. Later, Stafford and Tseng (1990) ${ }^{[J T 90]}$ and Rios-Mercado and Bard (1999) ${ }^{[R M B 99]}$ did further work on the problem. Tseng and Stafford (2001 ${ }^{[\text {TJ01] }}, 2002{ }^{[\text {ST02] })}$ proposed two MILP models. The flow shop problem with SDST and makespan criterion was taken into consideration by Ruiz et al. (2004) ${ }^{[R M A 05]}$. The good performing metaheuristics of a regular flow shop problem as well as advanced genetic algorithms were proposed. The flow shop problem with setup times were divided into four categories by Allahverdi et al. (1999) [AGA99] as sequence independent non-batch set-ups, sequence-dependent non-batch set-ups, sequence independent batch set-ups, and sequencedependent batch set-ups ${ }^{[A G A 99]}$. The researchers put their effort based on these categories to solve the problem in real world. For PCB manufacturing environment Lee and Shaw (2000) have done a great deal of work through the minimizing the objective function concept ${ }^{[L S 00]}$. The travelling salesman problem is used to model the problem F/no-idle/C $\mathrm{C}_{\max }$ by Saadani et al. (2004) [SGM05]. All the previous mentioned are serial flow shop however the concept of a concurrency in flow shop is shown by Lee et al. (1993) [LCL93] and Potts et al. (1995) ${ }^{[P S S}{ }^{+95]}$.

Simultaneously, with the work on flow shop, the special case of the permutation flow shop was also studied by the researchers with great deal of interest since 1960. In the permutation flow shop problem the same job order is followed by all machines and is denoted as $F|p r m u| c_{m a x}{ }^{[R . L 79]}$. Dudek and Teuton (1964) ${ }^{[D T 64]}$ had explained the basic assumptions for this problem in detail. The optimal calculation approaches had been approached earlier by Szwarc (1997) ${ }^{[S z w 71]}$, Lageweg et $a l(1978){ }^{[L L R K 78]}$, Potts (1980) ${ }^{[P o t 80]}$, and Carlier and Reba(1996) ${ }^{[C R 96]}$. The permutation flow shop problem is known to be a NP-hard problem for three or more machines [GJS76]. Hence, heuristics approach had been a way of getting a near-optimal solutions. Yet, there is no perfect framework that had been the best. Many researchers had tried to classify problem and the heuristics for better results ${ }^{[\text {Lou96] }}$. Gupta(1979) ${ }^{[G u p 79]}$, King and Spachis(1980) ${ }^{[\text {KS80] }}$ and Parl et al.(1984) ${ }^{[P P E 84]}$ had
given the review on development of heuristics. All these heuristics happen to have different nature such as computation time, complexity order or memory requirement. Further, Widmer and Hertz (1989) [WH89], Moccellin (1995) ${ }^{[\mathrm{Moc95]}}$ and Nawaz et al.(1983) ${ }^{[\mathrm{NEH} 83]}$ have worked on new better heuristics. There were even earlier attempts to classify the flow shop heuristics as fixed functional heuristics, floating functional heuristics, and synthetic functional heuristics. Framimam et al.(2004) ${ }^{\text {[FGL04] }}$ presented a general framework for the development of the heuristics. There are approaches to handle the uncertainties that may occur during the job execution. Studies had been initiated by Gholami, Zandieh, and Alem-Tabriz(2009) [GZAT09] and Ouelhadj and Petrovic(2009) [OP08] which later grab more researchers towards the area. While the researchers are investing their effort on various nature of permutation flow shop problem, we here took a simple dig on the exhaustive search and branch and bound method for the optimal. The both methods are consuming the multithreading approach with some tweaks in the algorithm. In addition, for near-optimal solution, we took genetic algorithm and converted it into a concurrent approach with some tweaks in the algorithm. These methods requires extensive computation which we are trying to minimize and add the concurrent approach for faster execution. It is also known that the optimal solution though branch and bound techniques are most widely used to examine the performance for most studies involving heuristics ${ }^{[B H 91]}$. Hence, we here have expected these approach will aid in further research in the field of static permutation flow shop problem.

In the first chapter we started with the previous related works followed by short introduction to scheduling and shop problems. Later, flow shop and permutation flow shop problem are explained in detail with example. We concluded first chapter with the algorithm for makespan calculation. In chapter two, we briefly describe branch and bound algorithm along with the related history and explained the implementation of BB algorithm in permutation flow shop problem. At the end, we showed the multi-threaded approach of the BB algorithm. Genetic algorithm and its implementation in detail is described in chapter three. The multi-threaded approach to genetic algorithm for permutation flow shop problem is illustrated at the end of the chapter. The next chapter is the results comparison and analysis followed by chapter with conclusion and future work.

### 1.1 Scheduling

A schedule is the allocation of intervals for each job to be processed on machines. The jobs can be of $n$ different types and machines can be $m$. For the illustrative purpose of jobs and its relation to machines Gantt charts are good. There are two ways that are used in representing the Gantt號
chart, first, machine-oriented and next is job-oriented.


Figure 1.1: Machine oriented Gantt Chart

| $\mathrm{J}_{1}$ | $M_{1}$ | $\mathrm{M}_{2}$ |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| $\mathrm{J}_{2}$ | $M_{2}$ | M |  |  |  |  |  |
|  |  |  |  |  |  |  |  |
| $\mathrm{J}_{3}$ | M ${ }^{\text {a }}$ | $M_{1}$ |  |  |  |  |  |
|  |  |  |  |  |  |  |  |
| $\mathrm{J}_{4}$ |  |  |  |  |  |  |  |

Figure 1.2: Job oriented Gantt Chart

As seen in the figure above, machine-oriented has machines fixed and jobs are places as per intervals. Similarly, job oriented has jobs fixed and machines are places as per intervals.

### 1.1.1 Job Data

We follow Peter Brucker ${ }^{[\text {Bru07] }}$ notation stating for scheduling problems. $J_{i}$ is a job consisting of operations $O_{i 1}, O_{i 2}, \ldots, O_{i n_{i}}$ where $n_{i}$ is number of operations for the job. For each operation $O_{i j}$, there is certain processing time $p_{i j}$. Assume $J_{i}$ has $\mathrm{n}_{\mathrm{i}}=1$, its operation is now $O_{i 1}$ only then processing time is denoted by $p_{i} . r_{i}$ is a release date, on which the first operation of $J_{i}$ becomes available for processing may be specified. Associated with each operation $O_{i j}$ is a set of machines $\mu_{i j} \subseteq\left\{M_{1}, \ldots, M_{m}\right\} . O_{i j}$ may be processed on any of the machines in $\mu_{i j}$. Usually, all $\mu_{\mathrm{ij}}$ are one element sets or all $\mu_{i j}$ are equal to the set of all machines. Machine-oriented has dedicated machines while job oriented has machines as parallel. Multi-purpose machines (MPM) are adjusted such that they can process operation. These MPM are used in real-world problems like flexible manufacturing has various steps of operations performed by machines with different tools. While processing $O_{i j}$ may use all machines in the set $\mu_{i j}$, scheduling problems as such are called multiprocessor task scheduling problems. A cost function $f_{i}(t)$ is defined by a due date $d_{i}$ and a weight $w_{i}$ measures the cost of completing $J_{i}$ at time $t$. The symbols $p_{i}, p_{i j}, r_{i}, d_{i}$ and $w_{i}$ are all assumed to be integers. The scheduling problems are classified based on a three-fields $\alpha|\beta| \gamma$ classification where $\alpha$ specifies the machine environment, $\beta$ specifies the job characteristics and $\gamma$ denotes the optimality criterion. This classification scheme was introduced by Grahm et al. [R.L79].

### 1.1.2 Job Characteristics

There are at most six parameters to define the job characteristics usually $\beta$ is used as $\beta_{1}, \beta_{2}, \beta_{3}$, $\beta_{4}, \beta_{5}$, and $\beta_{6}$. Preemption is indicated by $\beta_{1}$ as $\beta_{1}=p m t n$ if allowed else $\beta_{1}$ is not present in $\beta$. Precedence relations is described through $\beta_{2}$. The values for $\beta_{2}$ varies as prec, intree (outtree), tree, chains and sp-graph depending on different nature of jobs. If precedence relation is not there $\beta_{2}$ is not included in $\beta$. Release dates specification for a job is handled through $\beta_{3}=r_{i} . \beta_{3}$ is not in $\beta$ when $r_{i}=0$ for all jobs. Processing time or operations behavior is stated as in $\beta_{4}$. Job deadline is specified through $\beta_{5}$. Batching nature is represented by $\beta_{6}$ as the values $p$-batch or $s$-batch if present else $\beta_{6}$ is not included in $\beta$.

### 1.1.3 Machine Environment

Similar to job characteristics, machine environment is addressed through $\alpha$ as $\alpha_{1}$ and $\alpha_{2}$. The values of $\alpha_{1}$ are $o, P, Q, R, P M P M, Q M P M, G, X, O, J$ and $F$ which has meaning as dedicated machines, identical parallel machines, uniform parallel machines, unrelated parallel machines, multi-purpose machines with identical speed, multi-purpose machines with uniform speed, general shop, mixed shop, open shop, job shop and flow shop respectively. $\alpha_{2}$ represents the number of machines if specified else machines count is arbitrary.

### 1.1.4 Optimality Criteria

$\gamma$ is used as optimality criteria definition. $C_{i}$ is denoted as finishing time for job $J_{i}$ and cost is $f_{i}\left(C_{i}\right)$. Bottleneck objectives and sum objectives are types of total cost functions denoted respectively below as equations.

$$
f_{\max }(C)=\max \left\{f_{\mathrm{i}}\left(C_{\mathrm{i}}\right) \mid i=1, \ldots, n\right\}
$$

and

$$
\sum f_{\mathrm{i}}(C)=\sum_{i=1}^{n} f_{\mathrm{i}}\left(C_{\mathrm{i}}\right)
$$

The schedule which minimizes these total cost functions is the aim of scheduling problem. Optimality Criteria $\gamma$ is set as $\gamma=f_{\text {max }}$ or $\gamma=\sum f_{\mathrm{i}}$ also there are special functions too. The makespan $C_{m a x}$, total flow time $\sum C_{\mathrm{i}}$ and weighted flow time $\sum w_{\mathrm{i}} C_{\mathrm{i}}$ are some common objective functions. In addition, objective functions with due date $d_{i}$ associated jobs $J_{i}$ have various definition as:
lateness $L_{\mathrm{i}}=C_{\mathrm{i}}-d_{\mathrm{i}}$
earliness $E_{\mathrm{i}}=\max \left\{0, d_{\mathrm{i}}-C_{\mathrm{i}}\right\}$
tardiness $T_{\mathrm{i}}=\max \left\{0, C_{\mathrm{i}}-d_{\mathrm{i}}\right\}$
absolute deviation $D_{\mathrm{i}}=\left|C_{\mathrm{i}}-d_{\mathrm{i}}\right|$
squared deviation $S_{\mathrm{i}}=\left(C_{\mathrm{i}}-d_{\mathrm{i}}\right)^{2}$
unit penalty $U_{\mathrm{i}}= \begin{cases}0 & \text { if } C_{\mathrm{i}} \leq d_{\mathrm{i}} \\ 1 & \text { otherwise }\end{cases}$
A regular objective function is nondecreasing with respect to all variables $C_{i} . E_{i}, D_{i}$ and $S_{i}$ are not regular and others are regular.

### 1.2 Flow Shop

It has been shown that the three or more machine permutation flow shop problems are NP-complete problems (Gonzalez and Sahni, 1978). Some of other shop scheduling problems are open shop problems, job shop problems, and mixed shop problems. These problems fall in general shop problems category. The general shop problem consists of n number of jobs $J_{1}, J_{2}, \ldots, J_{n}$. Each $J_{i}$ having operations $O_{1}, O_{2}, \ldots, O_{n_{i}}$ and each operations $O_{i j}$ have processing time $p_{i j}$. There are m Machines $M_{1}, M_{2}, \ldots, M_{m}$. Each operations $O_{i j}$ must be processed on a machine $\mu_{i j} \in\left\{M_{1}, M_{2}\right.$, ..., $\left.M_{m}\right\}$. The operations of jobs have a precedence relationship. A machine can process a job at a time and no two machines can process the same job at the same time. Job $J_{i}$ has some objective function of finishing time $C_{i}$ and our objective is to find a feasible schedule that minimizes the finishing times for all jobs $J_{1}, J_{2}, \ldots, J_{n}$. The assumption is that the objective function is regular.

### 1.2.1 Open Shop Problem

An open shop problem is defined such that there is no precedence relationship between operations and each job have exactly $m$ operations where operation $\mathbf{O}_{\mathbf{i j}}$ has to be processed in machine $\mathbf{M}_{\mathbf{j}}$.

### 1.2.2 Job Shop Problem

A generalized version of the flow shop problem is job shop problem. There are $n$ jobs and $m$ machines. Operations vary based on the job number yet it preserves the order. Each operation with their processing time is needed to be processed in a particular machine. Considering the
finishing time of the last operation, the solution is to find the best possible schedule which has a minimum value for the objective function.

### 1.2.3 Mixed Shop Problem

The mixed shop problem comprises the open shop problem and the job shop problem thus has open shop and job shop jobs. There are bound of a job on number count as $n_{O}$ for open shop and $n_{J}$ for job shop.

### 1.2.4 Flow Shop Problem

There are Jobs such as $J_{1}, J_{2}, \ldots, J_{n}$ and Machines such as $\mathrm{M}_{1}, \mathrm{M}_{2}, \mathrm{M}_{3}, \ldots, \mathrm{M}_{\mathrm{m}}$. These Jobs has different operations such as $\mathrm{O}_{1}, \mathrm{O}_{2}, \mathrm{O}_{3}, \ldots, \mathrm{O}_{\mathrm{m}}$ to be ran on m different machines each. For particular Job say $J_{i}$ running particular operation $\mathrm{O}_{\mathrm{j}}$ on machine $\mathrm{M}_{\mathrm{j}}$ it takes processing time as $\mathrm{p}_{\mathrm{ij}}$ units.

There is a relationship between the job 's operations and machines. The job $J_{i}$ 's any operations $\mathrm{O}_{\mathrm{j}}$ can only begin on machine $\mathrm{M}_{\mathrm{j}}$ when operations $\mathrm{O}_{1}, \mathrm{O}_{2}, \mathrm{O}_{3}, \ldots, \mathrm{O}_{\mathrm{j}-1}$ for the job $J_{i}$ are completed from machines $\mathrm{M}_{1}, \mathrm{M}_{2}, \mathrm{M}_{3}, \ldots, \mathrm{M}_{\mathrm{j}-1}$. Alternatively, no two operations for the same jobs are processed at the same time and all preceding operations have to be completed before beginning a new one. A machine can take only a job at one time. The job order may differ considering the operations performed on the particular machine. As being flow shop, the operations order hence machine order are followed by each job as $\mathrm{M}_{1}$ then $\mathrm{M}_{2}$ then $\mathrm{M}_{3}$ till $\mathrm{M}_{\mathrm{m}}$. Overall the aim of this problem is to reduce the makespan such that the completion times for all the jobs on all the machines is minimum. Hence the flow shop problem requires the effective job order for each machine.Computing for n jobs and m machines flow shop problem has ( $\mathrm{n}!)^{\mathrm{m}}$ different solutions in order to get the optimal one. This is highly unlikely a feasible way to get optimal results. Therefore there are other approaches as permutation flow shop where problem size reduces to n !.

## Example of Flow Shop Problem

The example shows three jobs and three machines problem with a solution.

The makespan for the example is found to be 20 .

| Jobs | Machines | M1 | M2 | M3 |
| :---: | :---: | :---: | :---: | :---: |
| J1 | 2 | 1 | 2 |  |
| J2 | 1 | 3 | 4 |  |
| J3 | 5 | 1 | 2 |  |

Table 1.1: Flow Shop Problem Example Input


Figure 1.3: Gantt chart for the flow shop example

### 1.2.5 Permutation Flow Shop

The huge size of the solution of the flow shop problem is somewhat reduced using permutation flow shop. Here instead of considering different job orders for each machine, just only one job order is chosen for all the machines. The problem size hence turned to n !.

## Example of Permutation Flow Shop Problem

The example shows three jobs and three machines problem with a solution for job order $\mathrm{J}_{1}, \mathrm{~J}_{2}$ and $\mathrm{J}_{3}$.

| Jobs | Machines | M1 | M2 |
| :---: | :---: | :---: | :---: |
| M3 |  |  |  |
| J1 | 3 | 1 | 2 |
| J2 | 1 | 3 | 6 |
| J3 | 5 | 3 | 2 |

Table 1.2: Permutation Flow Shop Problem Example Input

The makespan for the problem is found as 15 .


Figure 1.4: Gantt chart for the permutation flow shop example

### 1.2.6 Makespan Calculation

The job and machine along with their values are represented as matrix M. I have assumed machines as rows and jobs are columns in matrix M.The total number of machines are as totalMachine and the total number of jobs are as totalJob. The order or sequence of a job to be computed through the machines is represented as an array $\mathbf{O}$. The makespan calculated value is here represented as Cmax which is an integer value.

The algorithm 1 below shows steps for calculation of makespan.

```
Algorithm 1 Cmax Calculation Algorithm
    Input: (i) Job Machine Matrix M, (ii) Job Order \(O\)
    Output: Cmax Value
    Initialize variables
    Allocate space for cumulative job completion calculation matrix \(\mathbf{J}\)
    for each machine \(r\) in \(\mathbf{M}\) do
        Allocate memory for the machine \(r\) and assign to matrix \(\mathbf{J}\)
        if \(\mathrm{r}=0\) then
            for each job \(c\) in \(\mathbf{M}\) do
                if \(c \neq 0\) then
                        \(\mathbf{J}[r][c]=\mathbf{J}[r][c-1]+\mathbf{M}[r][\mathbf{O}[c]]\)
            else
                        \(\mathbf{J}[r][c]=\mathbf{M}[r][\mathbf{O}[c]]\)
        else
            for each job \(c\) in \(\mathbf{M}\) do
                if \(c=0\) then
                        \(\mathbf{J}[r][c]=\mathbf{J}[r-1][c]+\mathbf{M}[r][\mathbf{O}[c]]\)
                    else
                        \(\mathbf{J}[r][c]=(\mathbf{J}[r-1][c]>\mathbf{J}[r][c-1]) ? \mathbf{J}[r-1][c] \mathbf{J}[r][c-1]+\mathbf{M}[r][\mathbf{O}[c]]\)
    Cmax \(\leftarrow \mathbf{J}[\) totalMachine - 1][totalJob - 1]
    Release Memory from matrix \(\mathbf{J}\)
```


## Chapter 2

## Branch and Bound

### 2.1 History

Branch and Bound $(B \& B)$ is by far the most widely used tool for solving large scale NP-hard combinatorial optimization problems. B\&B is, however, an algorithm paradigm, which has to be filled out for each specific problem type, and numerous choices for each of the components exist. Even then, principles for the design of efficient B\&B algorithms have emerged over the years.

The branch and bound methods in flow shop scheduling have been widely used for finding optimal or near optimal solution methods. Ignall and Schrage (1965) [IS65] Lomnicki (1965) [Lom65], McMahon and Burton (1967) ${ }^{[\text {MB67] }}$, Ashour (1970) ${ }^{[\text {Ash70] }}$, Gupta (1971) ${ }^{[G u p 71]}$, Lageweg et al. (1978) ${ }^{[L L K 78]}$, and Bansal (1979) ${ }^{[B a n 79]}$ among others have developed different branch and bound methods for various measures of performance like makespan, mean flow time, mean tardiness and maximum tardiness. The difference and the efficiencies of a branch and bound algorithms are in the choice of the lower bound (LB) and elimination rules. The strong bounds and elimination rules eliminate relatively more nodes of the search tree which very often brings in more computation requirements as well. If such needs are excessively large, it may become advantageous to search through larger nodes using a weaker, but fast computable LB. However, the advantages of stronger bounds and elimination rules are more substantial in large scale problems (Baker, 1975) [Bak75]. In 1973 Salvador ${ }^{[S a 173]}$ suggested the permutation flow shop problem's solution through Branch and bound. Kochhar and Morris (1987) ${ }^{[K M 87]}$ report work on the development of the heuristics. The heuristics developed try to minimize the effect of setup times and blocking. Further work has been reported by Brah and Hunsucker in the development of mathematical formulation, primarily useful for small size problems ${ }^{[B H 91]}$.

### 2.2 Overview

The concept of the branch and bound provides a strong basis for constructing the algorithms to solve NP-hard discrete optimization problems. Starting with the whole solutions the B\&B algorithm searches for the best solution. However, the exponential number of solutions restricts explicit enumeration. In order to make solution space feasible for getting best solution bounds play a vital role. The bounds let search focus on only those space where there is a possibility of having the best solution.

The search begins with a pool containing all the possible solutions which are marked to be unexplored. The unexplored such subset is represented as nodes. For such node, B\&B algorithm processes a node at a time. The operations of B\&B algorithm consists of selection of a node, calculation of bound and branching. The order of these steps depends on the strategy of choosing the next node for processing. If the bound value of the subproblem is considered for selection as next subproblem, then branching is done after choosing the node. The branching means dividing the current node space into two or more subspaces base on certain criteria. The branching is done till the subspace has a single solution left which is compared to the current best solution and decision is made to keep or discard based on the result of the comparison. Otherwise, the bounding function for the subspace is calculated and compared to the current best solution. If the comparison results show the subspace cannot have an optimal solution it is pruned or discarded else it is kept for further iterations. Since the bounds are calculated first this is called eager strategy. Another way is calculating the bound of the selected node and then branch on the node if necessary. The nodes created are then stored together with the bound of the processed node. This strategy is called lazy strategy. This is good for the depth-first approach in the search tree. Ultimately, the search space is such reduced to have a solution and all live subspaces explored then the current best solution is the required solution.

### 2.3 Simple $B \& B$ algorithm

The simplest B\&B algorithm concept as an eager strategy is enumerated below:

1. Consider the whole unexplored solution which is a live node $\mathbf{S}$ in live.
2. Repeat until there is no node in live
(a) Get a node $\mathbf{P}$ from live
(b) Generate branches from $\mathbf{P}$
(c) for each branch of $\mathbf{P}$
i. Get bound value $\mathbf{X}$ for each branch of $\mathbf{P}$,
ii. If branch has only one solution then compare the current best value and update if better.
iii. If branch has solutions pool then compare the lower bound X obtained to global if better keep the branch in live else discard.
3. After completion the global results are required solution.

### 2.4 Figurative illustration of branch and bound

Let us understand the $\mathrm{B} \& \mathrm{~B}$ concept through the figure. Figure 2.1 has images shown having problem representation in form of eclipse area and tree. The figure 2.1a shows the initial state of any problem, on the left as eclipse whereas to the right is the root node. Based on certain properties the area is divided into smaller regions which also can be illustrated as in the form of a tree as shown in figure 2.1b. Meanwhile, there occurs bounding or pruning or fathoming which restricts the need to explore a certain portion of the solutions. Thus all the solutions falling to the region are not needed to be computed. In the figure 2.1c below, it is assumed S1 and S3 are pruned as their LB is no better than the current best solution where S2 and S4 are further iterated to get more deeper into the tree and hence towards the optimal. Let us assume, that S21, S22, S41 and S42 each have a solution only. Each solution is compared to the global value. If the LB of of the solution is better then new global value are set else it is discarded. Finally, the global value is the required optimal for the problem.


Assumed here S1 and S3 will yield no better result
(c)

Figure 2.1: Illustrative of Branch and Bound

### 2.5 B\&B algorithm in Permutation Flow Shop Problem

The B\&B requires the solutions to be searched for arranged in the form of a tree. The tree consists of nodes. In my case the nodes consist of node level, lower bound for the node, partial job order till that node, remaining job order, children count, children pointer, parent pointer, is fathomed and other navigation necessary properties. The strategy followed in my case is eager. The initial upper bound (UB) is obtained by iterating the random job order and computing makespan for the job order to a fixed number of times with a minimum makespan preserved from the iteration. We prune the node during traversal earlier by comparing UB to LB. This helps in the reduction of many branches and hence nodes search. The traversal in a node is the hybrid best first search (BeFS). The breadth-first search (BFS) is not suitable as it requires huge memory also depth-first search (DFS) may take a longer time to get an optimal solution.

### 2.5.1 Lower bound for root

Initially, the LB computation based on each machine $\left(\mathrm{LB}_{\mathrm{m}}\right)$ is achieved by computing the complete job processing on the machine. Here, for the base machine, it is assumed all the jobs are aggregated and in addition the minimum of a job completion on all machines before the base machine is done and similarly the minimum of a job completion on all machine after the base machine is added. The lower bound of the root ( $\mathrm{LB}_{\text {root }}$ ) is assigned by selecting the maximum value obtained from the lower bound computed on each machine base. The maximum value is selected because it is only the possible value. The mathematical representation is as:

$$
\begin{gather*}
L B_{\mathrm{m}}=\min _{i(0 \ldots N-1)}\left(\sum_{j=0}^{j=m-1} p_{\mathrm{ij}}\right)+\sum_{i=0}^{i=N-1} p_{\mathrm{im}}+\min _{i(0 \ldots N-1)}\left(\sum_{j=m+1}^{j=M-1} p_{\mathrm{ij}}\right)  \tag{2.1}\\
L B_{\mathrm{root}}=\max _{m(0 \ldots M-1)}\left(L B_{\mathrm{m}}\right) \tag{2.2}
\end{gather*}
$$

In equation 2.1, $m$ is the machine for which LB is to be calculated, $i$ is any job, $j$ is any machine, $N$ is total jobs, $M$ is total machines and $p_{i j}$ is processing time for any job $i$ for any machine $j$. It can be illustrated in the figure as below:


Figure 2.2: Lower bound for root calculation illustration figurative

### 2.5.2 Lower bound for node

Similar to LB for the root, lower bound for the node $\left(\mathrm{LB}_{\mathrm{m}}\right)$ also has three parts. The calculation is similar based on the machine. First part is partial Cmax based on partial order by considering only the machines up-to machine $m$. The second part is the summation of the processing time of jobs which are not in the partial order. The third part is the same as in case of LB for root. Taking only jobs that are not in partial order and machines that are after machine $m$ in machine order, the minimum sum of any single job is the third portion. Lower bound for the node ( $\mathrm{LB}_{\text {node }}$ ) is chosen from $\mathrm{LB}_{\mathrm{m}}$ whichever is largest. The mathematical representation is:

$$
\begin{gather*}
L B_{\mathrm{m}}=C m a x_{\mathrm{rm}}+\sum_{i \notin R} p_{\mathrm{im}}+\min _{i \notin R}\left(\sum_{j=m+1}^{j=M-1} p_{\mathrm{ij}}\right)  \tag{2.3}\\
L B_{\mathrm{node}}=\max _{m(0 \ldots M-1)}\left(L B_{\mathrm{m}}\right) \tag{2.4}
\end{gather*}
$$

In equation 2.3, $R$ is partial job order, $r$ is last job of partial job order $R$ and Cmax $_{\mathrm{rm}}$ is makespan value for order R and machines $(0, \ldots, \mathrm{~m})$. Figurative explanation is shown below.


Figure 2.3: Lower bound for node calculation illustration figurative

### 2.5.3 Node selection and branching

As stated earlier, the selection is the hybrid BeFS. BeFS is beneficial as it consumes less memory than BFS and also reaches optimal faster than DFS. We have fused BeFS with DFS. The fusion with DFS enabled in less consumption of memory. The node with LB is selected and branched as the eager strategy. On each iteration, the search goes deeper into the tree until it reaches the leaf. Then, it tries to complete search on all sister nodes before searching back on the parent level. As search return back to parent it then release the memory occupied by the children. The branching is always done to nodes which have better values than the current UB. The branching is not done further than N-1 level because the child after the node is just one. Instead, the LB of its child is calculated and placed earlier. The selection algorithm is as:

In algorithm 2, the first while loop makes sure there are children node to be selected. On line 13, the child fathomed helps in reduction of the tree earlier and adds in finding optimal earlier. Similarly, memory release on line 16 adds more efficiency of the algorithm.

### 2.5.4 The B\&B Algorithm

The B\&B Algorithm is composed using equation 2.2, equation 2.4 and algorithm 2.
The root and the first children are created separately then algorithm 3 further is used. The method GetPartialJoborders (A) generates the partial job order and remaining job order for the

```
Algorithm 2 Hybrid BeFS algorithm for node selection
    Input: Current Node \(C\)
    Output: Next Node \(C^{\prime}\)
    Declare variable tempNext \(C^{\prime}\)
    if \(C\) is leaf node then
        C.isfathomed \(=\) true
        \(C \leftarrow C\).parent
    while C.allchildfathomed \(=\) true do
        \(C \leftarrow C\).parent till root.
    while true do
        for child \(c\) in C.children do
                if \(c\).isfathomed \(=\) false then
                        if c.lowerbound \(<G L O B A L\).upperbound then
                        \(C^{\prime} \leftarrow c\) and break for loop
                    else
                            c.isfathomed \(=\) true
        if C.allchildfathomed \(=\) true then
                C.isfathomed \(=\) true
                Release memory occupied by all the children
                if C.parent \(=\) NULL then \(\quad \triangleright\) all nodes traversal completed.
                    Return NULL
                else
                    \(C \leftarrow C\).parent
    for child \(c\) in \(C^{\prime}\).parent.children do
        if c.lowerbound \(<C^{\prime}\).lowerbound and \(c\). isfathomed \(=\) false then
                \(C^{\prime} \leftarrow c\)
    Return \(C^{\prime}\)
```

child using the parent node. Based on the equation 2.4, GetLBforTheNode (A) method provide the LB for intermediate nodes. As stated earlier, on line 17 we have calculated the LB earlier in the parent node and reduced the further iteration shortening the tree. Also, the child is fathomed based on the calculated LB as they produce no better result further through the children. Eventually, reaching the last node of the tree, the result is saved if better else fathomed.

### 2.5.5 The B\&B algorithm multi-threading

The single threaded approach takes more time. In order to improve the performance and reach the optimal solutions fast, we have implemented simple threading. The root's children are enqueued to a queue from where each thread picks a child and applies algorithm 3. The threading approach is shown below in the flow chart:

```
Algorithm 3 Branch and Bound algorithm
    Declare variables myNode, child, children and childcount
    while !nodeSearchComplete do
        myNode \(\leftarrow\) GetMimimumChildNode()
        if myNode \(=\) NULL then
            nodeSearchComplete \(\leftarrow\) true
            Break while loop
        childcount \(\leftarrow\) totalJobs - myNode.level
        if childcount \(>1\) then \(\triangleright\) Create children for myNode
            Allocate memory for myNode.children
            for \(i \leftarrow(0, \ldots\), childcount-1) do
                child.level \(\leftarrow\) myNode.level +1
                child.parent \(\leftarrow\) myNode
                child.fathomed \(\leftarrow\) false
                child.childAllFathomed \(\leftarrow\) false
                    child.childCount \(\leftarrow 0\)
                GetPartialJoborders(child.partialOrder,child.remainingOrder,myNode,i)
                if childcount \(=2\) then
                    child.partialOrder[totalJobs-1] \(\leftarrow\) child.remaingingOrder[0]
                    child.lowerbound \(\leftarrow\) getCmax(child.partialOrder) \(\quad\) refer algorithm 1
                else \(\quad \triangleright\) refer equation 2.4
                    child.lowerbound \(\leftarrow\) GetLBforTheNode(child.partialOrder,child.remainingOrder,child.level)
                if child.lowerbound \(>\) GLOBAL.upperbound then
                    child.fathomed \(\leftarrow\) true
                    child.childAllFathomed \(\leftarrow\) true
                    myNode.children \([i] \leftarrow\) child
        if childcount \(=1\) then
            if child.lowerbound \(<=\) GLOBAL.upperbound then
                saveResult(myNode.partialOrder,myNode.lowerbound)
            child.fathomed \(\leftarrow\) true
            child.childAllFathomed \(\leftarrow\) true
```



Figure 2.4: Multi-Thread B\&B Approach illustration

## Chapter 3

## Genetic Algorithm

### 3.1 History

Earlier in the 1950s and the 1960s, independently several computer scientists studied evolutionary systems. They intend to use this idea for solving engineering problems. This started with the idea to evolve a population of candidate solutions to a given problem with the aid of operators mimicking natural selection and genetic variation. On that run, genetic algorithms (GA) were invented by John Holland and later Holland and his students and colleagues at the University of Michigan developed it. Holland, inspired by Darwinian theory, thought to implement similar in computer systems to design algorithms which follows adaptation phenomenon as in nature. A theoretical framework for adaptation under the GA emerge as Holland's book Adaptation in Natural and Artificial Systems was published in 1975. [Mit98]

### 3.2 Basic Concept

A new population is created going through a method described by GA. The method follows natural selection in creating the new population as well the essence of genetics are imprinted through operators crossover, mutation and inversion. The population consists of chromosomes basically strings of 0 's and 1 's. Each gene is composed of an allele ( 0 or 1 ) and such genes join together to form a chromosome. The new population is created as the selection operator chooses the chromosomes that better fits. Sub-parts of two chromosomes are exchanged forming a crossover,mutation occurs randomly altering the allele values of some genes in the chromosome and a contiguous section's order is reversed through inversion all these operations makes a new array of genes in the chromosome. These operators idea is that the newly formed chromosome may provide a better
solution to the problem.

## Simple GA Algorithm

Now, let us visualize a simple genetic algorithm concept in the digital world. It is explained as:

1. Initially create a population of size $n$ having chromosomes of length $l$-bits randomly.
2. Calculate the fitness function $f(x)$ over each chromosome x in the population.
3. Repeat the steps until $n$ offspring have been created
(a) From the current population, select a pair of parent chromosomes. The chromosomes are selected based on the fitness function's value higher or lower as preferred for type of problem. The selection is with replacement meaning same parent can be selected more than once.
(b) Now do crossover based on crossover probability $p_{c}$. If no crossover occurred then offspring are parent themselves. The point for crossover in chromosome is determined randomly. The crossover can be single point or multi-point. Ultimately, 2 offspring are formed.
(c) With a mutation probability $p_{m}$, mutate the offspring and put in new population.
4. Make the newly formed population as current population.
5. Based on required generations, go to step 2 and repeat.

Likewise, the biological generation, here a generation is marked as those process of an iteration. Depending upon the need number of generations varies. The entire set of generations forms a run. In overall operation, probabilities and randomness play a vital role hence at the end of the run we get different results.

### 3.2.1 Simple GA Example

Let's take an illustrative example. Assume chromosome has a string of length $l 7$ bit, fitness function $f(x)$ be the count of 1's in the string, population size $n$ be 4 , crossover probability $p_{c}$ be 0.8 and mutation probability $p_{m}$ be 0.005 .

Consider randomly generated initial population as

| Chromosome Name | Chromosome String | Fitness Value |
| :--- | :--- | :--- |
| P | 0011110 | 4 |
| Q | 1101101 | 5 |
| R | 1010101 | 4 |
| S | 0001000 | 1 |

Table 3.1: A Simple GA Example - Initial Data

Simply using "roulette-wheel sampling" as an implementation of fitness-proportionate selection in selection method we get parents. A circular roulette wheel is divided into arc areas based on individual fitness from the current population and each chromosome are assigned accordingly. The wheel is spun as to population size times to get that many parents here 4 times. Let from first 2 spins we get P and Q then on later we get Q and R . With $p_{c}=0.8$, the parents P and Q crossover at the second bit to form $\mathrm{T}=0001101$ and $\mathrm{U}=1111110$. While Q and R didn't crossover so they became offspring. Next step, the mutation occurred on offspring T on the second locus to form $T^{\prime}=0101101$ and R on the fourth locus to form $\mathrm{R}^{\prime}=1011101$. The final result then is:

| Chromosome Name | Chromosome String | Fitness Value |
| :--- | :--- | :--- |
| $\mathrm{T}^{\prime}$ | 0101101 | 4 |
| U | 1111110 | 6 |
| Q | 1101101 | 5 |
| $\mathrm{R}^{\prime}$ | 1011101 | 5 |

Table 3.2: A Simple GA Example - Result Data

Here we got a new better fit $\mathrm{U}(6)$, the average fitness also rose from 14/4 to 20/4. Eventually, with further iterations we will get a string with all ones.

### 3.3 GA in Permutation Flow Shop Problem

There are few adjustments in simple GA are made in order to advocate the issue of permutation flow shop problem. The chromosomes strings are job numbers exactly one time instead of 0's and 1's. We here represent chromosome string as job order. The crossover is two point and mutation swaps any two jobs in the job order. The crossover probability $p_{c}$ is assumed to be 1 . Based on the fitness function or objective function value from the current population, top certain (say $\delta \%$ ) job orders are taken directly to the new population. The remaining needed population are obtained through the same steps as in simple GA algorithm assuming $p_{c}=1$.

### 3.3.1 Initial Population

As the first step, the initial population is generated. A initial job order is set say $J_{1}, J_{2}, \ldots, J_{n}$ and with an advanced random number generator, we get any two position values in the job order to swap and create new job order. This newly created job order is used as the base job order to swap and generate another new job order. This process is repeated until the population size is reached.

Below are related algorithms : The next job order generator plays a vital role in providing randomly distributed candidate solutions.

```
Algorithm 4 GA Next Job Order Generator Algorithm
    Input: Job Order O
    Output: Job Order O
    Initialize the required parameters \(a, b\)
    flag \(\leftarrow\) true
    while flag do
        \(a=\operatorname{mynrand}(0\), totalJobs -1\() \quad \triangleright\) refer appendix A
        \(b=\) mynrand \((0\), totalJobs -1\()\)
        if \(a \neq b\) then flag \(\leftarrow\) false
    \(\operatorname{swap}(\mathbf{O}, a, b) \quad \triangleright\) refer appendix A
    return \(\mathbf{O}\)
```

In the algorithm 4, I have insured the random numbers generated are in the range and are unique. This makes sure that resultant job order $\mathbf{O}$ after swapping is different from that of input.

The random population generation algorithm is as follows:

```
Algorithm 5 Initial GA Population Generation
    Input: (i) Job Machine Matrix M, (ii) Job Order O
    Output: (i) Population \(\mathbf{P}\) (ii) Top \(\delta \%\) job orders \(\mathbf{T}\)
    Initialize the required parameters, \(\mathrm{i}=0, \mathrm{Cmax}, \mathbf{P}, \mathbf{T}\)
    while \(\mathrm{i}<\) population size do
        \(\mathbf{O} \leftarrow \operatorname{getNextJobOrder}(\mathbf{O}) \quad \triangleright\) refer algorithm 4
        Cmax \(\leftarrow\) jobSumulate \((\mathbf{M}, \mathbf{O}) \quad \triangleright\) refer algorithm 1
        Bind Cmax with O
        Insert \(\mathbf{O}\) into \(\mathbf{P}_{\mathbf{i}}\)
        for each element \(t\) in \(\mathbf{T}\) do \(\quad \triangleright\) To get ranked top \(\delta \%\) job orders.
            if \(t . C \max \leq C \max\) then
                    Replace in \(\mathbf{T}\) then terminate loop.
    Return \(\mathbf{P}, \mathbf{T}\)
```

The algorithm 5 uses algorithm 1 and algorithm 4 to yield the initial random population $\mathbf{P}$.The process of ranking the whole population is expensive in terms of performance to get the top $\delta \%$ job orders. We here avoid a load of ranking by comparing and inserting the better job orders in
a fixed sized array $\mathbf{T}$. This array $\mathbf{T}$ is transferred to new population directly. The remaining job orders for the new population are obtained through selection, crossover, and mutation.

### 3.3.2 Selection

The selection is done from the current population $\mathbf{P}$ with replacement. The job orders along with their Cmax values are used to create a range in which higher Cmax occupies more range. The range begins from 0 to sum of $C \max$ from the current population $\mathbf{P}$. I have designed an algorithm which helps in getting a better job order out of the current population for crossover. The algorithm is as:

```
Algorithm 6 Higher Probable Job Order Index Finder Algorithm
    Input: (i) Cumulative Cmax Array \(\mathbf{P}^{\mathbf{c}}\) (ii) Population Size \(\mathbf{S}\)
    Output: Higher Probable Job Order Index for Current Population Array \(\mathbf{P}\)
    Initialize the required parameters, Cumulative Cmax \(\mathbf{C}^{\mathbf{T}} \leftarrow \mathbf{P}^{\mathbf{c}}[\mathrm{S}]\)
    index \(\leftarrow \mathbf{S}\)
    random \(\leftarrow \operatorname{mynrand}\left(1, \mathbf{C}^{\mathbf{T}}\right) \quad \triangleright\) refer appendix A
    for \(i=\mathbf{S}-1\) till \(i=0\) do
        if random \(\leq \mathbf{P}^{\mathbf{c}}[i]\) then
            inde \(x \leftarrow\) index -1
    Return index
```

The Cumulative Cmax Array $\mathbf{P}^{\mathbf{c}}$ is created at the time of population generation so as to accommodate the selection process later. The index obtained from this algorithm 6 is then used as index in the current population $\mathbf{P}$ to get a parent for crossover. In addition, to make sure the two job order obtained are different, we have added following snippet before each index is fed for crossover.

$$
\begin{aligned}
& a=\text { higherProbJob }\left(\mathbf{P}^{\mathbf{c}}, \mathbf{S}\right) \\
& b=\text { higherProbJob }\left(\mathbf{P}^{\mathbf{c}}, \mathbf{S}\right) \\
& \text { while } a=b \text { do } \\
& \quad b=\text { higherProbJob }\left(\mathbf{P}^{\mathbf{c}}, \mathbf{S}\right)
\end{aligned}
$$

$$
\triangleright \text { refer algorithm } 6
$$

### 3.3.3 Crossover

The two unique job orders $\left(\mathbf{O}_{\mathbf{1}}\right.$ and $\left.\mathbf{O}_{\mathbf{2}}\right)$ are obtained through selection operation. The crossover we are implementing here is two point. Randomly the points are obtained such that first point
is always less than half of the job order and second is either equal to half or greater. This point selection can be of other ways too. The offspring generation has mainly three steps, copying the job order from the beginning until the start point from a parent (say $\mathbf{O}_{\mathbf{1}}$ ), copying the job order from the start point to the end point inclusive from another parent (say $\mathbf{O}_{\mathbf{2}}$ ) and finally copying the remaining job order from parent again(say $\mathbf{O}_{\mathbf{1}}$ ). While copying from the start point onward it is made sure that the job number has not occurred previously in the sequence. If, for a position in the new job order, the job number matches then the next job number is taken and if it reaches the end of the job order in parent then it again scans from the beginning to find unmatched job number and then copy it to the new job order. This occurs twice to create two offspring. Below is the algorithm illustrating it.

```
Algorithm 7 Crossover Algorithm
    Input: (i) Job Order \(\mathbf{O}_{\mathbf{1}}\) (ii) Job Order \(\mathbf{O}_{\mathbf{2}}\)
    Output: (i) New Job Order \(\mathbf{O}_{\mathbf{1}}{ }^{\prime}\) (ii) New Job Order \(\mathbf{O}_{\mathbf{2}}{ }^{\prime}\)
    Initialize the required parameters, start point \(s p\), end point \(e p\)
    \(s p \leftarrow \operatorname{mynrand}(1,(\) totalJob \(/ 2)-1) \quad \triangleright\) refer appendix A
    \(e p \leftarrow \operatorname{mynrand}(\) totalJob \(/ 2\),totalJob - 1)
    while \(i<s p\) do
        \(\mathbf{O}_{\mathbf{1}}{ }^{\prime}[i] \leftarrow \mathbf{O}_{\mathbf{1}}[i]\)
        \(\mathbf{O}_{\mathbf{2}}{ }^{\prime}[i] \leftarrow \mathbf{O}_{\mathbf{2}}[i]\)
    for \(i=s p\) till \(i=e p\) do \(\quad \triangleright\) for \(\mathbf{O}_{\mathbf{1}}{ }^{\prime}\)
        flag \(\leftarrow\) true
        while flag do
            if \(\left(\mathbf{O}_{\mathbf{1}}[i] \neq \mathbf{O}_{\mathbf{2}}[i]\right)\) and \(\left(\mathbf{O}_{\mathbf{2}}[i]\right.\) does not exists in any \(\left.\mathbf{O}_{\mathbf{1}}{ }^{\prime}[i]\right)\) then
                    \(\mathbf{O}_{\mathbf{1}}{ }^{\prime}[i] \leftarrow \mathbf{O}_{\mathbf{2}}[i]\)
                flag \(\leftarrow\) false
            else
                Select next job order from \(\mathbf{O}_{\mathbf{2}}\)
    Repeat similarly from step 7 to 14 for \(\mathbf{O}_{\mathbf{2}}{ }^{\prime}\)
    Return \(\mathbf{O}_{1}{ }^{\prime}, \mathbf{O}_{\mathbf{2}}{ }^{\prime}\)
```


### 3.3.4 Mutation

The obtained new job orders $\mathbf{O}^{\prime}$ after crossover are subjected to mutation. The mutation occurs based on the mutation probability $p_{m}$. Usually, $p_{m}$ is very low.

```
Algorithm 8 Mutation Algorithm
    Input: Job Order \(\mathbf{O}^{\prime}\)
    Output: Mutated Job Order \(\mathbf{O}^{\prime}\)
    Initialize the required parameters \(a, b\)
    flag \(\leftarrow\) true
    while flag do
        \(a=\operatorname{mynrand}(0\), totalJobs -1\() \quad \triangleright\) refer appendix A
        \(b=\) mynrand \((0\), totalJobs -1\()\)
        if \(a \neq b\) then flag \(\leftarrow\) false
    \(\operatorname{swap}\left(\mathbf{O}^{\prime}, a, b\right) \quad \triangleright\) refer appendix A
    return \(\mathbf{O}^{\prime}\)
```


### 3.4 Population Generation

The new population for each generation are generated using the operators selection, crossover and mutation. The steps for new population generation uses algorithm 6 , algorithm 7 and algorithm 8. There is chances of memory issues so proper cleanup are required after each generation. The algorithm for new population generation is illustrated as in algorithm 9 below:

### 3.5 GA multi-threading

Multi-threads are used in mainly two places in order to gain in performance. The initial population generation is done by multiple threads. Later the new population pool generation is also multithreaded. The threading approach is illustrated in the diagram below:

```
Algorithm 9 New GA Population Generation
    Input: (i) Job Machine Matrix \(\mathbf{M}\) (ii) Population \(\mathbf{P}\) (iii) Cumulative Cmax Array \(\mathbf{P}^{\mathbf{c}}\)
    (iv) Population Size \(\mathbf{S}\) (v) Top \(\delta \%\) job orders \(\mathbf{T}\)
    Output: (i) New Population \(\mathbf{P}^{\prime}\) (ii) Top \(\delta \%\) job orders \(\mathbf{T}\) (iii) New Cumulative Cmax Array
    \(\mathbf{P}^{\mathbf{\prime C}}\)
    Initialize the required parameters, \(\mathrm{i}=0, \mathbf{P}^{\prime}\)
    \(\mathrm{i} \leftarrow(\mathbf{S} * \delta \%)-1\)
    \(\mathbf{P}^{\prime} \stackrel{\text { insertall }}{\leftarrow} \mathbf{T} \quad \triangleright\) insert top job orders from current population to new.
    Update \(\mathbf{P}^{\mathbf{\prime c}}\) using \(\mathbf{T}\)
    while \(\mathrm{i}<\mathbf{S}\) do
        \(a=\) higherProbJob \(\left(\mathbf{P}^{\mathbf{c}} \mathbf{,} \mathbf{S}\right) \quad \triangleright\) refer algorithm 6
        \(b=\) higherProbJob \(\left(\mathbf{P}^{\mathbf{c}}, \mathbf{S}\right)\)
        while \(a=b\) do
                \(b=\) higherProbJob \(\left(\mathbf{P}^{\mathbf{c}}, \mathbf{S}\right)\)
        \(\mathrm{O}_{\mathbf{1}} \leftarrow \mathrm{P}[\mathrm{a}]\)
        \(\mathrm{O}_{2} \leftarrow \mathrm{P}[\mathrm{b}]\)
        \(\mathbf{O}^{\prime}{ }_{1}, \mathbf{O}^{\prime}{ }_{2} \leftarrow \operatorname{crossover}\left(\mathbf{O}_{\mathbf{1}}, \mathbf{O}_{\mathbf{2}}\right) \quad \triangleright\) refer algorithm 7
        \(\mathbf{O}^{\mathbf{m}} \mathbf{1} \leftarrow \operatorname{mutate}\left(\mathbf{O}^{\prime}{ }_{1}\right) \quad \triangleright\) refer algorithm 8
        \(\mathrm{O}^{\mathbf{m}}{ }_{\mathbf{2}} \leftarrow \operatorname{mutate}\left(\mathbf{O}^{\prime}{ }_{\mathbf{2}}\right)\)
        Insert \(\mathbf{O}^{\mathbf{m}}{ }_{1}\) and \(\mathbf{O}^{\mathbf{m}} \mathbf{2}^{\text {into }} \mathbf{P}^{\mathbf{\prime}}\) and update \(\mathbf{P}^{\mathbf{\prime} \mathbf{c}}\)
        for each element \(t\) in \(\mathbf{T}\) do \(\quad \triangleright\) To get ranked top \(\delta \%\) job orders.
            if \(t . C \max \leq \mathbf{O}^{\mathbf{m}}{ }_{1} . C \max\) then
                    Replace in \(\mathbf{T}\) with \(\mathbf{O}^{\mathbf{m}}{ }_{1}\) then terminate loop.
        for each element \(t\) in \(\mathbf{T}\) do
            if \(t . C \max \leq \mathbf{O}^{\mathbf{m}}{ }_{\mathbf{2}} . C \max\) then
            Replace in \(\mathbf{T}\) with \(\mathbf{O}^{\mathbf{m}} \mathbf{2}^{\text {then }}\) therminate loop.
    Release memory of \(\mathbf{P}, \mathbf{P}^{\mathbf{\prime}}\)
    Return \(\mathbf{P}^{\mathbf{\prime}}, \mathbf{P}^{\mathbf{c}}, \mathbf{T}\)
```



Figure 3.1: Multi-Thread GA Approach illustration

## Chapter 4

## Results

The algorithms were converted to program and results were obtained. The results were computed as a single thread as well as multi-thread. For the case of multiple threads, we have used eight threads. We have used C programming of version C99. The machine we have used is java.cs.unlv.edu and visual studio 2017 IDE is used. The problem files were obtained from New hard benchmark for flow shop scheduling problems minimising makespan ${ }^{[E v a 15]}$. The machine configuration we have used is:

```
OS
GNU/Linux
Kernel Version
#1 SMP Fri Feb 1 14:54:57 UTC 2019
Kernel Release
3.10.0-957.5.1.el7.x86_64
CPU
Architecture: x86_64
CPU op-mode(s): 32-bit, 64-bit
CPU(s): 8
Vendor ID: GenuineIntel
CPU family: 6
Model: 58
Model name: Intel(R) Xeon(R) CPU E3-1240 V2 @ 3.40GHz
CPU MHz: 1643.969
CPU max MHz: 3800.0000
CPU min MHz: 1600.0000
L1d cache: 32K
L1i cache: 32K
L2 cache: 256K
L3 cache: 8192K
MEMORY
Memory block size: 128M
Total online memory: 32G
Total offline memory: OB
```

Figure 4.1: Machine Configuration

### 4.1 Exhaustive Search Results

The results were generated only from problems of jobs 10 and machines 5, 10, 15 and 20 which are in small set problems from VFR benchmark problems. Each problem has 10 different sets. The number of solutions to calculate in order to get all the optimal is given by the factorial of the number of jobs (for 10 jobs, $10!=3628800$ solutions). In addition, we have done the calculation for all the optimal for the problem. The speedup is calculated by dividing the execution time of a single thread by the execution time of multi-thread. The comparative results for single thread and multi-thread along with speedup is shown below:

| S.N. | Problem | Count | Time (Th 1) | Time (Th 8) | Speedup |
| :--- | :--- | :--- | :--- | :--- | :--- |
| 1 | $10 \_5 \_1$ | 2228 | 1341.94202 | 932.01001 | 1.43984 |
| 2 | $10 \_5 \_2$ | 30 | 1310.16406 | 922.09302 | 1.42086 |
| 3 | $10 \_5 \_3$ | 36 | 1333.31397 | 930.43103 | 1.43301 |
| 4 | $10 \_5 \_4$ | 26 | 1327.45300 | 921.53998 | 1.44047 |
| 5 | $10 \_5 \_5$ | 12 | 1333.36597 | 939.31598 | 1.41951 |
| 6 | $10 \_5 \_6$ | 323 | 1312.76599 | 923.65302 | 1.42128 |
| 7 | $10 \_5 \_7$ | 66 | 1328.79004 | 930.16998 | 1.42855 |
| 8 | $10 \_5 \_8$ | 12 | 1329.18799 | 922.89801 | 1.44023 |
| 9 | $10 \_5 \_9$ | 18 | 1331.31006 | 923.76599 | 1.44118 |
| 10 | $10 \_5 \_10$ | 48 | 1325.68201 | 929.86597 | 1.42567 |
| 11 | $10 \_10 \_1$ | 2 | 2445.39307 | 1576.40100 | 1.55125 |
| 12 | $10 \_10 \_2$ | 548 | 2458.28809 | 1558.51807 | 1.57732 |
| 13 | $10 \_10 \_3$ | 24 | 2427.65991 | 1564.85107 | 1.55137 |
| 14 | $10 \_10 \_4$ | 4 | 2455.43604 | 1560.08594 | 1.57391 |
| 15 | $10 \_10 \_5$ | 5 | 2421.36206 | 1575.70496 | 1.53668 |
| 16 | $10 \_10 \_6$ | 317 | 2426.29590 | 1582.04297 | 1.53365 |
| 17 | $10 \_10 \_7$ | 1 | 2432.15088 | 1566.06494 | 1.55303 |
| 18 | $10 \_10 \_8$ | 48 | 2437.12695 | 1556.12598 | 1.56615 |
| 19 | $10 \_10 \_9$ | 6 | 2451.75708 | 1572.51502 | 1.55913 |
| 20 | $10 \_10 \_10$ | 15 | 2447.73193 | 1574.06604 | 1.55504 |
|  |  |  |  |  |  |

Table 4.1 continued from previous page

| S.N. | Problem | Count | Time (Th 1) | Time (Th 8) | Speedup |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 21 | 10_15_1 | 1 | 3527.47998 | 2177.64111 | 1.61986 |
| 22 | 10_15_2 | 1 | 3527.82007 | 2190.71191 | 1.61035 |
| 23 | 10_15_3 | 1 | 3509.75708 | 2196.44092 | 1.59793 |
| 24 | 10_15_4 | 5 | 3520.12891 | 2171.95313 | 1.62072 |
| 25 | 10_15_5 | 2 | 3503.35400 | 2196.25610 | 1.59515 |
| 26 | 10_15_6 | 1 | 3498.88794 | 2195.74512 | 1.59349 |
| 27 | 10_15_7 | 56 | 3539.15405 | 2186.43799 | 1.61868 |
| 28 | 10_15_8 | 9 | 3514.88208 | 2176.32690 | 1.61505 |
| 29 | 10_15_9 | 15 | 3531.58789 | 2202.57910 | 1.60339 |
| 30 | 10_15_10 | 1 | 3529.57495 | 2199.30005 | 1.60486 |
| 31 | 10_20_1 | 2 | 4639.98682 | 2852.52002 | 1.62663 |
| 32 | 10_20_2 | 1 | 4600.63721 | 2872.04102 | 1.60187 |
| 33 | 10_20_3 | 1 | 4594.14600 | 2889.40210 | 1.59000 |
| 34 | 10_20_4 | 4 | 4628.41992 | 2859.86401 | 1.61841 |
| 35 | 10_20_5 | 2 | 4594.84180 | 2841.83594 | 1.61686 |
| 36 | 10_20_6 | 218 | 4621.86523 | 2874.50195 | 1.60788 |
| 37 | 10_20_7 | 1 | 4625.17676 | 2847.19702 | 1.62447 |
| 38 | 10_20_8 | 2 | 4645.15820 | 2872.38989 | 1.61718 |
| 39 | 10_20_9 | 3 | 4601.27197 | 2869.10010 | 1.60373 |
| 40 | 10_20_10 | 2 | 4621.06104 | 2865.68604 | 1.61255 |

Table 4.1: Exhaustive search result single and multi threads with optimal results count.

The table 4.1 is summarized in terms of the number of machines and we get the following result.


Figure 4.2: Machine based Threaded and Non-Threaded comparison with speedup

The threading seems to give an average speedup of 1.5. It is expected to increase the speed of execution by 8 times as we use eight threads but only 1.5 is obtained. This happened because of Amdahl's law ${ }^{[G e n 67]}$, which states, there is only a certain percentage (i.e. less than $100 \%$ ) of execution time which can be subjected to the speedup. But as we go on increasing more jobs this approach wouldn't be enough to calculate optimal as the number of solution size increases exponentially.

### 4.2 Branch and Bound Algorithm Results

The next implementation Branch and Bound algorithm provided better results. The results for the same problem as of exhaustive search was taken. The comparative analysis of the algorithm based on execution time and nodes visited with and without threading along with speedup is tabulated below:

| S.N. | Problem <br> Size | BB |  | BB Threaded |  | BB <br> /BB Th | Ex Th <br> /BB Th |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  | Time | Node <br> Visited | Time | Node <br> Visited |  |  |
| 1 | 10_5_1 | 385.05801 | 1289946 | 250.62300 | 1124397 | 1.53640 | 3.71877 |
| 2 | 10_5_2 | 44.01800 | 172967 | 25.69200 | 144731 | 1.71330 | 35.89028 |
| 3 | 10_5_3 | 62.76900 | 250422 | 37.32700 | 217280 | 1.68160 | 24.92649 |
| 4 | 10_5_4 | 58.16800 | 234954 | 40.32100 | 214046 | 1.44262 | 22.85509 |
| 5 | 10_5_5 | 4.04000 | 12520 | 3.64300 | 16025 | 1.10898 | 257.84133 |
| 6 | 10_5_6 | 12.77600 | 47784 | 13.14300 | 51025 | 0.97208 | 70.27718 |
| 7 | 10_5_7 | 35.44700 | 140861 | 24.56200 | 132755 | 1.44316 | 37.87029 |
| 8 | 10_5_8 | 60.72302 | 244797 | 32.65100 | 174194 | 1.85976 | 28.26553 |
| 9 | 10_5_9 | 1.89900 | 4279 | 1.57200 | 4507 | 1.20802 | 587.63740 |
| 10 | 10_5_10 | 89.51600 | 361014 | 46.94700 | 280575 | 1.90675 | 19.80672 |
| 11 | 10_10_1 | 57.54300 | 151556 | 29.01200 | 143617 | 1.98342 | 54.33617 |
| 12 | 10_10_2 | 271.76502 | 790482 | 142.21503 | 731582 | 1.91094 | 10.95888 |
| 13 | 10_10_3 | 151.82001 | 245277 | 43.97800 | 229746 | 3.45218 | 35.58259 |
| 14 | 10_10_4 | 95.69500 | 224473 | 40.99800 | 200530 | 2.33414 | 38.05273 |
| 15 | 10_10_5 | 24.67600 | 57958 | 11.19000 | 55244 | 2.20518 | 140.81367 |
| 16 | 10_10_6 | 158.48300 | 453318 | 80.11900 | 422942 | 1.97810 | 19.74616 |
| 17 | 10_10_7 | 276.18103 | 803231 | 118.64200 | 632401 | 2.32785 | 13.19992 |
| 18 | 10_10_8 | 95.78300 | 130833 | 27.13900 | 124448 | 3.52935 | 57.33911 |
| 19 | 10_10_9 | 25.94100 | 41289 | 9.45500 | 38058 | 2.74363 | 166.31571 |
| 20 | 10_10_10 | 157.42000 | 384958 | 80.25500 | 407163 | 1.96150 | 19.61331 |
| 21 | 10_15_1 | 81.05500 | 133817 | 35.69400 | 130673 | 2.27083 | 61.00860 |
| 22 | 10_15_2 | 104.81200 | 189062 | 40.79900 | 159475 | 2.56898 | 53.69524 |
| 23 | 10_15_3 | 234.21800 | 435419 | 65.86400 | 285716 | 3.55609 | 33.34813 |
| 24 | 10_15_4 | 364.97803 | 725653 | 158.66400 | 671070 | 2.30032 | 13.68901 |
| 25 | 10_15_5 | 52.03400 | 70898 | 22.34500 | 78818 | 2.32866 | 98.28849 |
| 26 | 10_15_6 | 60.20400 | 97532 | 23.55500 | 89527 | 2.55589 | 93.21779 |
| 27 | 10_15_7 | 242.08200 | 475161 | 122.84802 | 458486 | 1.97058 | 17.79791 |

Table 4.2 continued from previous page

| S.N. | Problem <br> Size | BB |  | BB Threaded |  | BB <br> /BB Th | Ex Th <br> /BB Th |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- |
|  |  | Time | Node <br> Visited | Time | Node <br> Visited |  |  |
| 28 | $10 \_15 \_8$ | 164.58000 | 310466 | 57.57000 | 230850 | 2.85878 | 37.80314 |
| 29 | $10 \_15 \_9$ | 162.23900 | 293375 | 63.21900 | 276696 | 2.56630 | 34.84046 |
| 30 | $10 \_15 \_10$ | 71.40997 | 112719 | 31.05200 | 113137 | 2.29969 | 70.82636 |
| 31 | $10 \_20 \_1$ | 249.46500 | 295685 | 80.60500 | 267074 | 3.09491 | 35.38887 |
| 32 | $10 \_20 \_2$ | 166.57800 | 197686 | 61.16300 | 181222 | 2.72351 | 46.95717 |
| 33 | $10 \_20 \_3$ | 418.05899 | 594717 | 132.83200 | 475108 | 3.14728 | 21.75230 |
| 34 | $10 \_20 \_4$ | 79.84800 | 82861 | 35.50800 | 92656 | 2.24873 | 80.54140 |
| 35 | $10 \_20 \_5$ | 241.64301 | 299718 | 97.46400 | 271912 | 2.47931 | 29.15780 |
| 36 | $10 \_20 \_6$ | 144.31500 | 171803 | 58.46000 | 173378 | 2.46861 | 49.17041 |
| 37 | $10 \_20 \_7$ | 100.94400 | 109576 | 34.58300 | 93880 | 2.91889 | 82.32938 |
| 38 | $10 \_20 \_8$ | 82.06900 | 84133 | 35.27500 | 84345 | 2.32655 | 81.42848 |
| 39 | $10 \_20 \_9$ | 98.76000 | 109283 | 49.70100 | 144366 | 1.98708 | 57.72721 |
| 40 | $10 \_20 \_10$ | 161.92900 | 186145 | 76.32200 | 219059 | 2.12166 | 37.54731 |

Table 4.2: Execution time and node visited results for single and multi-threaded comparison with speedup

In table 4.2, it is clearly observed from the results that we gain performance when multithreading is used. The execution time of single and multi-threading are summarized in terms of average time, maximum time, minimum time and median time in order to understand the algorithm's performance from different prospect.

We found out the execution time was decreased for multi-threading which can be easily seen in figure 4.3. The data obtained is summarized and we get speedup summary as the maximum 3.556 , minimum 0.972 , average 2.252 and median 2.285 . Since the median is higher than average we can say the speedup is more skewed towards the left which is low. Yet, we can see there is good speedup factor. Similarly, we have calculated the execution time of multi-threaded exhaustive search to multi-threaded branch and bound on table 4.2. The summarized speedup results are way


Figure 4.3: Execution Time comparison for Single and Multi-Threaded
better with average 67.039 , maximum 587.637, minimum 3.718 and median 37.836 . The median is lower than average which means there are higher speedup values in comparison to lower ones.


Figure 4.4: Nodes Visited comparison for Single and Multi-Threaded

The next is, we put the nodes visited by the branch and bound method based on single and multi-
thread are summarized in terms of their average node visited, maximum nodes visited, minimum nodes visited and median of nodes visited. It is clearly visible from figure 4.4, in every aspect the multi-threaded approach outperforms the single thread. In order to get optimal, for the case of 10 jobs, the exhaustive search has to calculate 3628800 solutions which we can see has been drastically reduced in case of branch and bound. The number of nodes visited is summarized as maximum 1124397, minimum 4507, average 246068 and median 177708. This shows that we were able to reduce a lot of computation (only $30 \%$ get optimal even for maximum nodes visited).

The result is also compared to the result from $A$ branch-and-bound method to minimize the makespan in a permutation flow shop with blocking and setup times ${ }^{[T N 17] . ~ O n l y ~ t h e ~ r e l e v a n t ~}$ problems were taken and compared:

| Problem Size | BB(TN) |  | BB(here) |  | Speedup |
| :--- | :--- | :--- | :--- | :--- | :--- |
|  | Time | Nodes | Time | Nodes |  |
| $10 \_5$ | 54930 | 199743.05 | 47.64 | 235953.50 | 1153.02 |
| $10 \_10$ | 129950 | 270314.60 | 58.30 | 298573.10 | 2228.98 |

Table 4.3: BB results comparison with Takano and Nagano [TN17]

The table 4.3 shows our method visits more node in comparison to that of Takano and Nagano's results. However, there is a significant gain in terms of the execution time. In addition, the speedup is even better for a larger problem.

### 4.3 Genetic Algorithm Results

The genetic algorithm was used to calculate near-optimal solutions for all the small size problems from VFR benchmark problems with varied population size and generations. The results are generated varying the generation and the population size. Further, the set of 10 problems of each problem size are averaged in terms of the approximation ratio (AR) and the execution time. The AR here denotes how close to optimal we were able to get the makespan. In addition, the exact optimal obtained for that problem size is also noted.

We have results summarized for twenty-four different problem size into three different tables. Each table has records for two different values of generation and population. In the first table, we took 50 population size and generation of 50 and 100 . Similarly on the second table population is increased to 100 and generations to 100 and 1000. The third table has records for generation 2000 and population 100 as well for generations 1000 and population 1000 .

| S.N. | Problem <br> Size | Generation 50 Population 50 |  |  | Generation 100 Population 50 |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  | Ave AR | Ave Time | Count | Ave AR | Ave Time | Count |
| 1 | 10_05 | 1.00539 | 41.87093 | 4 | 1.00162 | 79.26700 | 7 |
| 2 | 10_10 | 1.00483 | 33.35743 | 4 | 1.00643 | 78.45760 | 3 |
| 3 | 10_15 | 1.00950 | 33.33077 | 0 | 1.00248 | 77.08480 | 2 |
| 4 | 10_20 | 1.00920 | 33.92823 | 1 | 1.00410 | 78.52070 | 2 |
| 5 | 20_05 | 1.04371 | 39.08390 | 0 | 1.03076 | 79.60833 | 0 |
| 6 | 20_10 | 1.07757 | 38.15840 | 0 | 1.06539 | 73.98053 | 0 |
| 7 | 20_15 | 1.07059 | 38.61563 | 0 | 1.06261 | 70.66593 | 0 |
| 8 | 20_20 | 1.06446 | 37.74750 | 0 | 1.05550 | 68.19177 | 0 |
| 9 | 30_05 | 1.02147 | 39.68787 | 0 | 1.01935 | 74.48783 | 0 |
| 10 | 30_10 | 1.10117 | 38.28437 | 0 | 1.09125 | 71.27473 | 0 |
| 11 | 30_15 | 1.11074 | 36.78903 | 0 | 1.09393 | 68.17823 | 0 |
| 12 | 30_20 | 1.09615 | 35.49190 | 0 | 1.08396 | 65.73140 | 0 |
| 13 | 40_05 | 1.02954 | 40.26257 | 0 | 1.02016 | 72.73687 | 1 |
| 14 | 40_10 | 1.09974 | 37.50417 | 0 | 1.08995 | 73.94120 | 0 |
| 15 | 40_15 | 1.11996 | 35.35737 | 0 | 1.10796 | 70.73783 | 0 |
| 16 | 40_20 | 1.12134 | 36.75620 | 0 | 1.10704 | 69.20780 | 0 |
| 17 | 50_05 | 1.02196 | 35.74777 | 0 | 1.01411 | 69.48460 | 0 |
| 18 | 50_10 | 1.09544 | 33.94417 | 0 | 1.08130 | 66.22093 | 0 |
| 19 | 50_15 | 1.13573 | 33.30110 | 0 | 1.11449 | 64.89777 | 0 |
| 20 | 50_20 | 1.13615 | 33.40837 | 0 | 1.12478 | 64.60513 | 0 |
| 21 | 60_05 | 1.03200 | 29.98170 | 0 | 1.01900 | 57.68743 | 0 |
| 22 | 60_10 | 1.10462 | 29.02387 | 0 | 1.08971 | 56.20353 | 0 |
| 23 | 60_15 | 1.13586 | 28.81720 | 0 | 1.12293 | 56.09087 | 0 |
| 24 | 60_20 | 1.14475 | 28.95177 | 0 | 1.13094 | 56.20207 | 0 |

Table 4.4: Genetic algorithm summary results 1

| S.N. | Problem <br> Size | Generation 100 Population 100 |  |  | Generation 1000 Population 100 |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  | Ave AR | Ave Time | Count | Ave AR | Ave Time | Count |
| 1 | 10_05 | 1.00184 | 70.42243 | 8 | 1.00014 | 637.15124 | 9 |
| 2 | 10_10 | 1.00304 | 67.30763 | 5 | 1.00055 | 563.00970 | 8 |
| 3 | 10_15 | 1.00451 | 66.25617 | 3 | 1.00145 | 554.71808 | 7 |
| 4 | 10_20 | 1.00441 | 67.27426 | 3 | 1.00000 | 549.31816 | 10 |
| 5 | 20_05 | 1.03209 | 68.53643 | 0 | 1.01390 | 558.81753 | 0 |
| 6 | 20_10 | 1.05253 | 63.50790 | 0 | 1.03567 | 548.49307 | 0 |
| 7 | 20_15 | 1.05213 | 55.56420 | 0 | 1.02947 | 540.93880 | 0 |
| 8 | 20_20 | 1.04463 | 55.47463 | 0 | 1.02646 | 537.31767 | 0 |
| 9 | 30_05 | 1.01604 | 56.46510 | 1 | 1.00770 | 551.06350 | 3 |
| 10 | 30_10 | 1.07758 | 55.59820 | 0 | 1.05031 | 541.44977 | 0 |
| 11 | 30_15 | 1.09166 | 55.69220 | 0 | 1.05935 | 1235.39293 | 0 |
| 12 | 30_20 | 1.07884 | 56.01157 | 0 | 1.04972 | 2297.76910 | 0 |
| 13 | 40_05 | 1.01835 | 56.01090 | 1 | 1.00590 | 998.44786 | 3 |
| 14 | 40_10 | 1.08696 | 55.90403 | 0 | 1.05410 | 630.12128 | 0 |
| 15 | 40_15 | 1.10128 | 55.95207 | 0 | 1.06491 | 544.53797 | 0 |
| 16 | 40_20 | 1.10191 | 56.79460 | 0 | 1.07217 | 552.93111 | 0 |
| 17 | 50_05 | 1.01267 | 56.25187 | 1 | 1.00408 | 630.04404 | 3 |
| 18 | 50_10 | 1.07916 | 56.33490 | 0 | 1.04923 | 547.22150 | 0 |
| 19 | 50_15 | 1.11664 | 56.96643 | 0 | 1.07751 | 556.66084 | 0 |
| 20 | 50_20 | 1.11920 | 58.15757 | 0 | 1.08277 | 565.84767 | 0 |
| 21 | 60_05 | 1.02101 | 56.49667 | 0 | 1.00436 | 1189.01226 | 2 |
| 22 | 60_10 | 1.08601 | 57.13873 | 0 | 1.05288 | 2486.81500 | 0 |
| 23 | 60_15 | 1.11879 | 58.59247 | 0 | 1.07877 | 635.25043 | 0 |
| 24 | 60_20 | 1.12831 | 59.89063 | 0 | 1.08899 | 580.39360 | 0 |

Table 4.5: Genetic algorithm summary results 2

| S.N. | Problem <br> Size | Generation 2000 Population 100 |  |  | Generation 1000 Population 1000 |  |  |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- |
|  |  | Ave AR | Ave Time | Count | Ave AR | Ave Time | Count |
| 1 | $10 \_05$ | 1 | 1231.2886 | 10 | 1 | 1912.20807 | 10 |
| 2 | $10 \_10$ | 1.00085 | 1127.04913 | 8 | 1 | 1937.68647 | 10 |
| 3 | $10 \_15$ | 1.00044 | 1109.87139 | 8 | 1 | 1944.28724 | 10 |
| 4 | $10 \_20$ | 1 | 1099.07697 | 10 | 1 | 1933.09803 | 10 |
| 5 | $20 \_05$ | 1.00956 | 1188.34078 | 0 | 1.00356 | 1902.75586 | 5 |
| 6 | $20 \_10$ | 1.03018 | 1097.1213 | 0 | 1.01747 | 1913.47974 | 0 |
| 7 | $20 \_15$ | 1.0255 | 1082.8509 | 0 | 1.01744 | 1913.43444 | 0 |
| 8 | $20 \_20$ | 1.01966 | 1075.11493 | 0 | 1.01378 | 2023.56199 | 0 |
| 9 | $30 \_05$ | 1.00662 | 1166.30976 | 3 | 1.00544 | 1945.76337 | 5 |
| 10 | $30 \_10$ | 1.04901 | 1083.87214 | 0 | 1.0433 | 1968.38396 | 0 |
| 11 | $30 \_15$ | 1.04903 | 1083.37943 | 0 | 1.04395 | 2069.26316 | 0 |
| 12 | $30 \_20$ | 1.04918 | 2744.781 | 0 | 1.03957 | 2277.33846 | 0 |
| 13 | $40 \_05$ | 1.00336 | 1148.62089 | 5 | 1.00301 | 1998.68504 | 4 |
| 14 | $40 \_10$ | 1.04807 | 1091.90557 | 0 | 1.04423 | 2140.90675 | 0 |
| 15 | $40 \_15$ | 1.06168 | 1238.30443 | 0 | 1.05737 | 2420.30684 | 0 |
| 16 | $40 \_20$ | 1.06874 | 4779.18795 | 0 | 1.06045 | 2785.26738 | 0 |
| 17 | $50 \_05$ | 1.00187 | 1171.33963 | 4 | 1.00287 | 2271.09966 | 3 |
| 18 | $50 \_10$ | 1.04451 | 1094.29702 | 0 | 1.04193 | 2452.44397 | 0 |
| 19 | $50 \_15$ | 1.07411 | 2802.18254 | 0 | 1.06927 | 4018.52802 | 0 |
| 20 | $50 \_20$ | 1.07756 | 4038.72514 | 0 | 1.07054 | 3988.47355 | 0 |
| 21 | $60 \_05$ | 1.00315 | 1172.21559 | 3 | 1.00462 | 2716.75994 | 2 |
| 22 | $60 \_10$ | 1.05169 | 1106.39646 | 0 | 1.04525 | 6880.58966 | 0 |
| 23 | $60 \_15$ | 1.06941 | 5149.50628 | 0 | 1.06913 | 3383.70755 | 0 |
| 24 | $60 \_20$ | 1.08586 | 1252.06204 | 0 | 1.0759 | 6904.88014 | 0 |
|  |  |  | 0 |  | 0 | 0 |  |

Table 4.6: Genetic algorithm summary results 3

We accumulated the results of table 4.4, table 4.5 and table 4.6 based on number of jobs and number of machines. The results are compared in terms of the execution time, the AR achieved
and the number of exact optimal reached. The accuracy calculation formula is:

$$
A R=\frac{G A \text { Value }}{\text { Optimal }}
$$

The AR $=1$ means the algorithm has generated the optimal perfectly. Any value higher than 1 means the result from GA is deviated from the optimal.

## Analysis per number of jobs

The figure 4.5 shows the time comparison. It is seen with an increase in the number of jobs the execution time increase. The execution time increase for change in configuration values which is also increasing.


Figure 4.5: Averaged execution time comparison summarized over number of jobs

For an increase in the number of jobs the AR indicate more deviation from the optimal as seen in the figure 4.6. But with a change in configuration GA provides better results and converge towards the optimal.

The third comparison in the figure 4.7 shows that more optimal is achieved as we increase the value of generation and population size. For a lower number of job we are able to get more optimal.


Figure 4.6: Averaged AR comparison summarized over number of jobs


Figure 4.7: Exact optimal count comparison summarized over number of jobs

## Analysis per number of machines

Based on the machine and the configuration values the GA tend to have an increase in the execution time. The respective execution time plot is shown in figure 4.8:


Figure 4.8: Averaged execution time comparison summarized over number of machines

The AR comparison is similar to the AR comparison based on the number of jobs as shown earlier. The result tends to be nearer to the optimal as the configuration is changed to higher values. Also, AR is better for problems with less number of machines. It is shown in the figure 4.9:


Figure 4.9: Averaged AR comparison summarized over number of machines

The trend is similar to that of job-based comparison for the number of optimal counts. More optimal is achieved as we change the configuration to higher values and count are higher for less number of machines. It is shown in the figure 4.10:


Figure 4.10: Exact optimal count comparison summarized over number of machines

### 4.4 Problems Faced

There are various technical problems that occurred during the thesis. The problem given by VFR was in a different format to read which was adjusted while reading from the file and creating the matrix. The multi-threading implementation initially raised the issue of race conditions as the original algorithm's approach didn't fit for the concurrent approach. With the use of multi-thread, the issue of memory leak arose which was handled by proper allocation and release of memory. The default library in C wasn't good enough for generating higher value random number. A different random number generator was used for the case. There was an issue of lack of memory for some case so many other problems weren't computed. For results comparison, there were no other results found in case of the VFR problems which are different from that of Taillard's.

## Chapter 5

## Conclusion and Future Work

### 5.1 Conclusion

We here have shown the improvement in algorithms for performance as well compared the obtained results. We achieved a maximum speedup of 1.6 in case of exhaustive search which can be further increased with an increase in the number of threads. We also have computed for all the optimal for the problem which can be reflected as a basis for near-optimal methods.

The next part, Branch and Bound performs way better in comparison to that of exhaustive search. A minimum speedup of 3 to 587 is obtained for the same problem set in compared to exhaustive search. Moreover, in terms of execution time, our branch and bound algorithm had speedup over 1153.02 compared to that of the branch and bound algorithm of Takano and Nagano. This suggests our approach of the branch and bound served well.

In the case of the genetic algorithm, we scale up our problem solving capacity than that of the other two approaches. Comparatively, we have observed the variation of results based on the configuration of generations and population size. It can be clearly seen that the higher the number of generation and population size the better are the results. But, there is the cost of higher execution time. Also, the population size seems to have a larger impact on the result than the generations.

Overall, we have computed a benchmark for the VFR permutation flow shop problems. These problems seem to have higher complexity with an increase in the number of jobs as well as an increase in the number of machines.

### 5.2 Future Work

The exhaustive search approach can be made as distributed with multi-threading approach which could give results for a larger problem. The branch and bound algorithm can be extended as distributed with multi-threading approach. In addition, the proper memory management may increase the algorithm's capacity of solving a larger problem. In genetic algorithm, a varying configuration can help in determining the best values of generations and population size based on the problem size. A proper approach to determine mutation percentage and best carry over chromosome percentage may yield better results. Also, properly random initial random population generation method may provide better result.

## Appendix A

## Selected Source Code

## Random Generator

```
/***Provides higher range random number ***/
unsigned long long llrand() {
    unsigned long long r = 0;
    for (int i = 0; i < 5; ++i) {
        r = (r << 15) | (rand() & 0x7FFF);
    }
    return r & 0xFFFFFFFFFFFFFFFFFULL;
}
/***Provides a random number between specified range***/
unsigned int mynrand(unsigned int lower, unsigned int upper)
{
    return (lower + llrand() % (upper - lower + 1));
}
```


## Swap

```
/***Swaps values at specified places***/
void swap(unsigned int *a, unsigned int i, unsigned int j)
{
    unsigned int temp = a[i];
    a[i] = a[j];
    a[j] = temp;
}
```

Lower Bound for Root
unsigned int GetInitialLBforRoot() \{

```
    unsigned int* sum1 = (unsigned int *)malloc(totalJobs * sizeof(unsigned int));
    unsigned int* sum3 = (unsigned int *)malloc(totalJobs * sizeof(unsigned int));
    unsigned int sum2 = 0;
    unsigned int minFirstJob = 0;
    unsigned int minLastJob = 0;
    for (unsigned int m = 0;m < totalMachines;m++) {
        tempLB = 0;
        sum2 = 0;
        minFirstJob = 0;
        minLastJob = 0;
        //find min sum for first jobs before the machine m.
        for (c = 0;c < totalJobs;c++) {
        sum1[c] = 0;
        for (a = 0;a < m;a++) {
        sum1[c] += originalMatrix[a] [c];
        }
        //sum all jobs in the machine m
        sum2 += originalMatrix[m][c];
        //find sum of last jobs on remaining machine after machine m
        sum3[c] = 0;
        for (a = m + 1;a < totalMachines;a++) {
            sum3[c] += originalMatrix[a][c];
        }
        if (c != 0) {
                if (sum1[minFirstJob] > sum1[c]) {
                minFirstJob = c;
                }
                if (sum3[minLastJob] > sum3[c]) {
                                    minLastJob = c;
            }
        }
    }
    tempLB = sum1[minFirstJob] + sum2 + sum3[minLastJob];
    if (tempLB > LB) {
        LB = tempLB;
    }
}
free(sum1);
free(sum3);
return LB;
```

\}

## Lower Bound for Node

unsigned int level) \{
register unsigned int m, c, a;
unsigned int tempLB, tempClr, sum2, minLastJobs, LB = 0;
unsigned int* sum3 = malloc(totalJobs * sizeof(unsigned int)); for ( $\mathrm{m}=0 ; \mathrm{m}$ < totalMachines; $\mathrm{m}++$ ) \{
tempLB $=$ sum2 $=$ minLastJobs $=0$;
tempClr = GetPartialCmax(jOrder, level, m);
for ( $c=0 ; c<t o t a l J o b s ~-~ l e v e l ; c++) ~\{~$
sum2 += originalMatrix[m][rOrder [c]];
sum3[c] = 0;
for ( $\mathrm{a}=\mathrm{m}+1 ; \mathrm{a}$ < totalMachines; $\mathrm{a}++$ ) \{
sum3[c] += originalMatrix[a][rOrder [c]];
\}
if (c ! = 0) \{
if (sum3[minLastJobs] > sum3[c]) \{
minLastJobs $=c$;
\} \}
\}
tempLB = tempClr + sum2 + sum3[minLastJobs];
if (tempLB > LB) \{ LB = tempLB;
\}
\}
free(sum3);
return LB;
\}

Generate Partial Order

```
void GetPartialJoborders(unsigned int *partialJobOrder, unsigned int *remJobOrder,
            unsigned int currLevel, unsigned int remJobNo) {
    register unsigned int r = 0, c = 0;
    //transferred specified job from remaining order to fixed order
    partialJobOrder[currLevel] = remJobOrder[remJobNo];
    //remove the job no from remJobDrder
    for (c;c < totalJobs - currLevel;c++) {
            if (c != remJobNo) {
                remJobOrder[r] = remJobOrder[c];
                r++;
            }
    }
```


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# Curriculum Vitae 

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