
Advanced Mathematical Programming

Branch-and-Price for the Bin-Packing Problem

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Contents

1	Introduction	2
1.1	Bin-Packing Problem	2
1.2	Branch-and-Price approach	2
1.2.1	Set covering formulation	2
1.2.2	Restricted master problem	3
1.2.3	Subproblem	3
1.2.4	Branching	3
1.3	Resolution method	4
2	Branch-and-Price outline	5
3	Node processing	7
3.1	Ryan & Foster method	7
3.1.1	Branching rules	7
3.1.2	Subproblem resolution	7
3.2	Generic method	7
3.2.1	Branching rules	7
3.2.2	Subproblem resolution	9
3.3	Node processing algorithm	10
4	Heuristics	12
4.1	Root heuristics	12
4.2	Tree heuristics	12
4.2.1	MIRUP-based strategy	13
4.2.2	Rounding strategies	13
5	Numerical applications	14
5.1	Numerical results	14
5.2	Result analysis	14

1 Introduction

1.1 Bin-Packing Problem

In this report, we focus on the Bin-Packing Problem. It consists of putting objects of different size in bins with equal capacity while minimizing the number of bin used. Consider that set of items is given by $I = \{1, \dots, N\}$, their size is given by $S = \{s_1, \dots, s_n\}$ and that B bins with a capacity C are available. The number of bin is supposed large enough to store all the n objects. Thus, the Bin-Packing Problem can be written under the following form :

$$\left\{ \begin{array}{ll} \min & \sum_{b=1}^B y_b \\ \text{s.t.} & \sum_{i=1}^N s_i x_{ib} \leq C y_b \quad \forall b = 1, \dots, B \\ & \sum_{b=1}^B x_{ib} = 1 \quad \forall i = 1, \dots, N \\ & x_{ib} \in \{0, 1\} \quad \forall i = 1, \dots, N \quad \forall b = 1, \dots, B \\ & y_b \in \{0, 1\} \quad \forall b = 1, \dots, B \end{array} \right. \quad (\text{BPP})$$

Where $x_{ib} = 1$ if the item i is packed in the bin number b ($x_{ib} = 0$ otherwise) and $y_b = 1$ if the bin number b is non-empty ($y_b = 0$ otherwise). The first constraint ensure that the bin capacity is not exceeded and the second constraint ensure that each item is packed in a bin. This problem is known to be NP-hard.

1.2 Branch-and-Price approach

To address the NP-hardness of the problem, we can solve this problem with a Branch-and-Price algorithm.

1.2.1 Set covering formulation

We choose an other formulation of the (BPP). Let \mathcal{P} the set of all combination of items of I such that their cumulative size doesn't exceed the bin capacity C . Each element of \mathcal{P} is called a pattern. Let $P = |\mathcal{P}|$, we have

$$\mathcal{P} = \left\{ \mathbf{x} \in \{0, 1\}^{|N|}, \quad \sum_{i=1}^N x_i s_i \leq C \right\} = \left\{ \mathbf{x} = \sum_{p=1}^P \alpha^p \bar{x}^p, \quad \sum_{p=1}^P \alpha^p = 1, \alpha \in \{0, 1\}^{|P|} \right\}$$

Here, \bar{x}^p is a pattern and $\bar{x}_i^p = 1$ if the item i is used in the pattern p . Using this Dantzig reformulation of the set \mathcal{P} , we can write (BPP) under its set covering formulation :

$$\left\{ \begin{array}{ll} \min & \sum_{p=1}^P \alpha^p \\ \text{s.t.} & \sum_{p=1}^P x_i^p \alpha^p = 1 \quad \forall i = 1, \dots, N \\ & \alpha^p \in \{0, 1\} \quad \forall p = 1, \dots, P \end{array} \right. \quad (\text{SCBPP})$$

For this formulation, $x_i^p = 1$ if the item i is within the pattern p and $\alpha^p = 1$ if the pattern p is used in the solution. Thus, the solution is a set of patterns and each patter correspond to a bin. The constraint ensure that each item is packed in a pattern.

On thing to notice is that $|P|$ (the number of variables) is huge. At most, $|P|$ is the number of combination available with N items. Even for small instances, the number of variable is not tractable. Furthermore, it is needed to enumerate and store all the patterns possible in order to solve the problem which could involve a huge amount of memory.

1.2.2 Restricted master problem

To handle this problem, we introduce a new set $\mathcal{P}' \subset \mathcal{P}$ of cardinal P' containing only a fraction of the available patterns. We can solve (SCBPP) on this restricted number of pattern, which is more trackable. This leads to a sub-optimal bound. This new problem is called the restricted master problem :

$$\left\{ \begin{array}{ll} \min & \sum_{p=1}^{P'} \alpha^p \\ \text{s.t.} & \sum_{p=1}^{P'} x_i^p \alpha^p = 1 \quad \forall i = 1, \dots, N \\ & \sum_{p=1}^{P'} \alpha^p \leq B \\ & \alpha^p \in [0, 1] \quad \forall p = 1, \dots, P' \end{array} \right. \quad (\text{RMBPP})$$

If the set \mathcal{P}' is well chosen and contains the optimal patterns, then the optimal solution of (RMBPP) will be the same as the optimal solution of (SCBPP) (and also the solution of (BPP)). The problem is to choose the right patterns to include in \mathcal{P}' . As \mathcal{P}' will also have a big cardinal, we rather solve a relaxation of the master problem and that is why we have that $\alpha_p \in [0, 1] \quad \forall p = 1, \dots, P'$. We also add a constraint on the number of pattern allowed as we know that there are only B bins available. If we have some additional informations about the problem, it is possible to tighten this bound. Solving the relaxation will lead to an supra-optimal solution comparing to the integer form of (RMBPP) (with $\alpha^p \in \{0, 1\}$). We will see later how to obtain an integer solution in order to have the same solution as the optimal solution of (BPP).

1.2.3 Subproblem

We can write an other optimization problem called the subproblem (or pricing problem) which will generate "interesting" patterns to include in (RMBPP). If we note π and σ the optimal dual variables corresponding to the first and second constraint of (RMBPP), this pricing problem takes the following form :

$$\left\{ \begin{array}{ll} \min & 1 - \sum_{i=1}^N \pi_i y_i - \sigma \\ \text{s.t.} & \sum_{i=1}^N y_i s_i \leq C \\ & y_i \in \{0, 1\} \quad \forall i = 1, \dots, N \end{array} \right. \quad (\text{SPBPP})$$

The solution of the pricing problem is a feasible pattern (*i.e.* which doesn't exceed the bin capacity). The optimal cost is called the reduced cost. We can see that usually, the cost of a pattern is 1 (one pattern used correspond to one bin used) but here, the cost is penalized by a term containing π and σ . This second term will penalize patterns which can not improve the set \mathcal{P}' in the sense that if a pattern p has a high reduced cost, the solution of (RMBPP) with \mathcal{P}' or with $\mathcal{P}' \cup p$ will be the same. In fact, if z_{sp}^* denotes the optimal cost and \mathbf{y}^* the associated patter of (SPBPP) for a given π and σ , then \mathbf{y}^* can improve the solution of (RMBPP) if and only if $z_{sp}^* < 0$.

On thing very important to notice is that this pricing problem is equivalent to a knapsack problem with weights π , sizes s and capacity C . A knapsack problem is easy to solve. The key mechanism of the Branch-and-Price algorithm will be to solve sequentially (RMBPP) and (SPBPP) until a solution $z_{sp}^* \geq 0$ is obtained. At this stage, (RMBPP) will be solved at optimality. As (RMBPP) is a relaxation of (SCBPP), then we have to introduce a new mechanism in order to obtain feasible solutions.

1.2.4 Branching

Now, let's remember that (RMBPP) is a relaxation of (SCBPP) so it gives a supra-optimal solution. In order to have the optimal integer solution, we will introduce branching rules (new constraints) to break the fractional property of the variables. Let's consider a tree where each node correspond to a problem with its local branching rules. When a node is solved to optimality (*i.e.* (RMBPP) is solved sequentially until optimality), we will select fractional variables in the solution to branch on. Then, child nodes will be created under the node just solved with

new constraints in order to have a non-fractional variable for the rest of the branch. The tree will be explored node after node until all the nodes are explored. The key idea will be to cut non-improving branches during the explorations so as to explore only a part of the tree.

We can not branch only on the fractional α^p . Instead, we will select a pair (i, j) of items such that the variable

$$w_{ij} = \sum_{p \in \mathcal{P}' \mid x_i^p = x_j^p = 1} \alpha^p$$

is fractional. In one branch, we will impose that $w_{ij} \geq 1$ (items i and j always together) and in the other branch, we will impose that $w_{ij} \leq 0$ (items i and j always separated).

1.3 Resolution method

To solve the Bin-Packing Problem, we will use two different branching rules. The first one was proposed by Ryan & Foster in 1981 [RF81]. This branching rule allow to keep a single subproblem at each node but doesn't preserve the knapsack structure of the problem. The subproblems will rather be knapsack with conflicts problems, harder to solve than the usual knapsack problem. The second branching rule which will be used is a generic branching scheme introduced by Vance in 1994 [Van+94]. This branching rule preserves the knapsack structure of the subproblems but at each node, multiple subproblems will have to be solved.

For the Ryan & Foster branching rule, two methods will be proposed for the subproblem resolution. The first one is to use a classical solver so as to model and solve the subproblem with the new branching constraints. We will also see how to solve the subproblems with dynamic programming [Tot80]. For the Generic branching rule, we will also propose to solve the subproblem either with a solver or with a dynamic programming algorithm [SV13].

In addition, we will see how to construct an heuristic solution before the tree exploration in order to set better initial bounds. Three variations of a decreasing-size-order packing algorithm will be proposed [BHB09]. [TODO : Heuristics within the tree ?]

Finally, we will see two different way to add the nodes to the queue : FIFO and LIFO. We will see if the queueing methods affects the running time or the number of nodes explored during the BnP algorithm. [TODO : Diving strategies ?]

2 Branch-and-Price outline

The structure of the branch of price algorithm is the same whatever branching-rule, queueing method, heuristics and other parameters are chosen. It can be resumed by the following algorithm :

Algorithm 1: Branch and Price

Input: The items and their size s_1, \dots, s_N , the bin capacity C , an upper bound on the number of bins B , a precision ϵ

// Initialization

Initialize an empty tree

Initialize the queue with the root

Initialize the column pool with an artificial column

Initialize $UB \leftarrow B$

Initialize $LB \leftarrow \left\lceil \sum s_i / C \right\rceil$

Process a root heuristic to find a better UB (see 4.1)

// Tree exploration

while *the queue is non-empty* **do**

 Pop the first node in the queue

 Proceed the node according to the branching rule set (see 3.1, 3.2 and 3.3)

 Find the branching variable and add the two child nodes to the queue

 Process a tree heuristic to tighten UB (see 4.2)

 Update LB and UB with the solution found by the node or by the heuristic

 Cut branches that cannot improve the upper bound

if $|UB - LB| < \epsilon$ **then**

 | return the best solution associated to UB

end

end

In the following, the simplest steps of the BnP are explained. The heuristics, the branching method and the node processing will have a dedicated section to be explained.

Column management :

The column pool contains all the columns which are created by the subproblems. We need to add an artificial column which allow the master problem to always be feasible. This artificial column contains all the objects and have a very high cost. If the solution of a master problem contains this artificial column, we know that the master problem is infeasible because of the branching rules. Before each node processing, we will create a local node pool containing only the patterns satisfying the branching rules of the node in order to create a solution satisfying the current branching rules. In order to save speed and memory space, the column pool is global to all nodes and each node will have its proper node pool which is a filter on the global column pool. When a new pattern is created in a subproblem, it is added to the node pool and to the global column pool.

Initial bounds :

We could have set the initial bound at $UB = +\infty$ and $LB = -\infty$ but better initial bound can be found. Indeed, we can at least assume that the total number of bin used will be at most the number of items. Thus, we can set $B = N$. The integer relaxation of (BPP) gives also a lower bound for the integer (BPP) : we can set $LB = \left\lceil \sum s_i / C \right\rceil$ [SV13].

Queueing method :

In the while loop, we always choose the first node in the queue but the nodes won't always been stored in the same order in the queue. When adding the two child nodes to the queue, two methods can be used. The first one puts the new nodes at the beginning of the queue (LIFO) in order to proceed a Deep-First-Search in the tree. The second method is to put the new nodes at the end of the queue (FIFO) in order to proceed a Breadth-First-Search. The LIFO method will allow to find quickly good upper bounds but it will cut branches more deeply in the tree. The FIFO method will allow to cut branches at a high level in the tree but will be slower to get good bounds. The

best exploration method depends on the structure of the problem and in general, there is no one better than the other. An hybrid exploration of the tree can also be done. First, the LIFO method is set in order to find quickly a good upper-bound. Once this upper bound has been found, the queuing strategy switch to FIFO in order to explore the tree with a Breadth-First-Search. This method aims to find quickly an upper bound and then cut branches high in the tree. If a root heuristic is computed, then the Hybrid method is just like the LIFO method.

Feasible solution handling :

After the node processing, it is possible that the solution given is integer. In this case, we can see if this solution improve the current UB and if it is the case, we can update the value of UB. We also choose the best lower bound among all those of the nodes in order to update the global LB. This allow to see how close the algorithm is from the solution by bounding the optimal solution by LB and UB at each node processing.

Tree pruning :

Once the bounds are updated, we take a look at the nodes left in the queue. If a node has a lower bound larger than UB, then this node can't improve the current best solution. Thus, we can stop the exploration of the branch after this node. When $UB \simeq LB$, then the current best solution can not be improved and correspond to the optimal solution of (BPP). If the queue becomes empty while $UB \neq LB$, this means that either the problem has no solution or that ϵ is too small regarding to numerical errors. The method to branch, process a node, to run heuristics before and within the tree will be explained in the following sections. Then, we will evaluate the BnP with some performance criteria.

3 Node processing

Now that the core structure of the BnP has been presented, we can focus on the node processing where two methods can be used : the Ryan & Foster and the generic method.

3.1 Ryan & Foster method

The Ryan & Foster branching rule will allow to keep a single subproblem per node but it will not be as easy to solve as a classical knapsack problem.

3.1.1 Branching rules

When a node is solved to optimality and that two items i and j are found to create a branch, we have to create one child where $w_{ij} \leq 0$ (the down branch) and one child where $w_{ij} \geq 1$ (up branch). For the down branching rule, the Ryan & Foster branching rule simply add the constraint $x_i + x_j \leq 1$ in the subproblem in order to create a pattern with items i and j separated. For the up branching rule, the constraint $x_i = x_j$ will be added to the subproblem in order to create patterns containing i and j and patterns containing neither i nor j (as it is impossible to create a solution only with patterns with i and j in it). The subproblem will handle itself whether to generate patterns with (resp. without) i and j because the reduced cost will be too high if too many pattern with (resp. without) i and j have always be created.

While exploring the tree, the branching rules will be added sequentially to nodes. Thus, for nodes at the bottom of the tree, the subproblems will have several branching constraints. It is possible that some combination of constraints make the subproblem infeasible. In this case, the process of the node will be stopped and the branch will be cut as it will be impossible to create solutions for the rest of the branch.

Using such method for the node processing allow to keep a unique subproblem per node and it is simple to create the node pool as we only loop on the global column pool and keep only the columns satisfying the branching rules.

3.1.2 Subproblem resolution

The subproblem will always have to initial structure presented in 1.2.3 with the new node branching constraints. Thus, it is possible to solve this problem using a solver like Gurobi. [TODO : Dynamic prog]

3.2 Generic method

The generic method introduce a way more complex branching scheme but allow to keep subproblems with the knapsack structure. As the knapsack problem can be solve with dynamic programming, this will allow to have a faster resolution method for the subproblems.

3.2.1 Branching rules

When a node is solved to optimality and that two items i and j are found to create a branch, we have to create one child where $w_{ij} \leq 0$ (the down branch) and one child where $w_{ij} \geq 1$ (up branch). For the down branch, we will create a subproblem generating patterns with $x_i = 0$ and patterns with $x_i = 1, x_j = 0$. Thus, the solution constructed will be constituted of patterns with items i and j separated. But we only need one pattern with $x_i = 1, x_j = 0$ so we will also impose that

$$\begin{cases} \sum_{p \in \mathcal{P}' | x_i^p = 1, x_j^p = 0} \alpha^p = 1 \\ \sum_{p \in \mathcal{P}' | x_i^p = 0} \alpha^p = B - 1 \end{cases} \quad (3.2.1)$$

For the up branching rule, we will create one subproblem generating patterns with $x_i = x_j = 0$ and patterns with $x_i = x_j = 1$. Thus, the solution will be composed of patterns with i and j together (and patterns without i and j).

But we only need one pattern with $x_i = x_j = 1$ so we will also impose that

$$\begin{cases} \sum_{p \in \mathcal{P}' | x_i^p = x_j^p = 1} \alpha^p = 1 \\ \sum_{p \in \mathcal{P}' | x_i^p = x_j^p = 0} \alpha^p = B - 1 \end{cases} \quad (3.2.2)$$

For up and down branching rules, we can notice that the variables x_i and x_j are fixed so the subproblem keep its knapsack structure if a preprocess is made. After fixing the branching variable, a knapsack problem will remained to be solved.

The difficulty lies in the method to process several branches in a row in the tree. We have to introduce the notion of **branching set** which are composed of some **branching rules** (corresponding to the filter below the sum in (3.2.1) and (3.2.2)) and one **coefficient** (corresponding to the scalar 1 or $B - 1$ in (3.2.1) and (3.2.2)). Each branching rule has several variables set to zero and several variables set to one. For example,

$$\sum_{p \in \mathcal{P}' | \{x_1=0, x_2=1\} \cup \{x_3=0, x_4=1\}} \alpha^p = 1$$

is a branching set with branching rules $\{x_1 = 0, x_2 = 1\}$ and $\{x_3 = 0, x_4 = 1\}$ and with coefficient 1 and

$$\sum_{p \in \mathcal{P}' | \{x_1=x_2=0\}} \alpha^p = B - 1$$

is a branching set with branching rule $\{x_1 = x_2 = 0\}$ and with coefficient $B - 1$.

When a new node is created, we will loop over all the branching sets of the parent node and create new sets for the child node by adding the new branching rules. Four cases can happen :

- **The coefficient of the parent node set is $L > 1$:**

In this case, the branching set have a unique rule by construction of the child nodes.

- **We have to add a down branching rule with items i and j :** In this case, we will create two branching set for the child node. One with the same items set to one and zero as the parent node but we add the item i to the variable set to 0. This branching set have a coefficient $L - 1$. The second branching set will have the same items set to one and zero as the parent node but we add the item i to the variable set to 1 and the item j to the variable set to 0. This second branching set have a coefficient equal to 1.
- **We have to add an up branching rule with items i and j :** In this case, we will create two branching set for the child node. One with the same items set to one and zero as the parent node but we add the item i and j to the variable set to 0. This branching set have a coefficient $L - 1$. The second branching set will have the same items set to one and zero as the parent node but we add the item i and j to the variable set to 1. This second branching set have a coefficient equal to 1.

- **The coefficient of the parent node set is 1 :**

In this case, the branching set can have a multiple rule by construction of the child nodes so only one pattern satisfying one of the rule will be selected in the solution.

- **We have to add a down branching rule with items i and j :** Only one set is created in the child node. For each branching rule in the parent node branching set, we add two new rule to the child branching set containing the rule of the parent branching set and either $x_i = 0$ or $x_i = 1, x_j = 0$. This new branching set has a coefficient equal to 1.
- **We have to add an up branching rule with items i and j :** Only one set is created in the child node. For each branching rule in the parent node branching set, we add two new rule to the child branching set containing the rule of the parent branching set and either $x_i = x_j = 0$ or $x_i = x_j = 1$. This new branching set has a coefficient equal to 1.

The way new constraints are added to the child nodes is a bit complicated as we have to handle multiple subproblems and we have to ensure that the coefficient of each branching set add to B in order to satisfy the constraint $\sum_{p \in \mathcal{P}} \alpha^p = B$ of (RMBPP) for each node. Of course, some of the rules created may be infeasible if one item is both in the set of variables set to 0 and 1. In this case, the rule is not included in the branching set. If the parent branching set is split into one feasible and one infeasible branching set, then we only keep the feasible branching set and the coefficient of the new set is equal to the coefficient of the parent branching set.

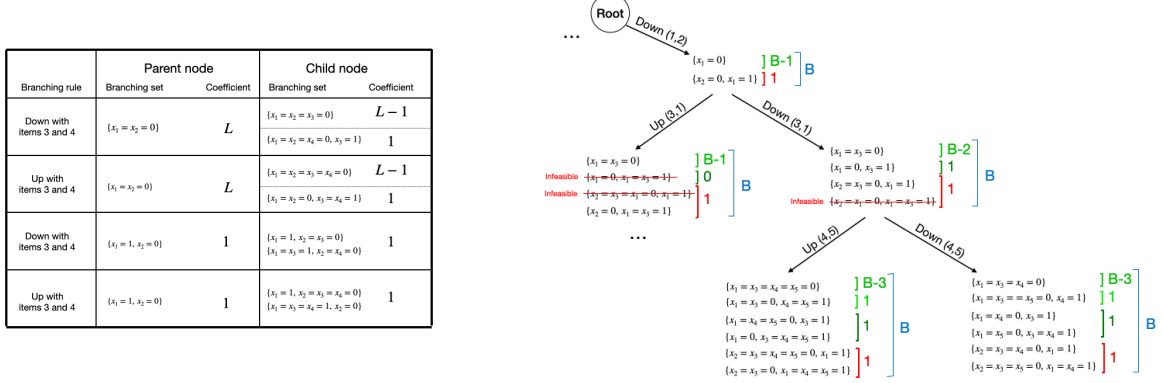


Figure 1: Left : The four cases presented above. Right : An example of how branching constraints are handled.

3.2.2 Subproblem resolution

In each node, there will be as many subproblems to solve as the total number of branching rules. As for the Ryan & Foster branching method, it is possible to use a solver like Gurobi to solve the subproblems. It will preprocess the problem by setting the variables to 0 or 1 according to the branching schemes before the solving process. But as the knapsack structure is preserved for each subproblem, it is possible to use dynamic programming instead of a classical solver.

The first step to do before the dynamic programming process is to preprocess the data and treat the variables already fixed by the branching rules. For each variable set to 0, we simply remove this variable from the problem. For each variable set to 1, we remove the variable from the problem, we also decrease the knapsack capacity by the size of the item and add its cost to a preprocess cost. At the end of the preprocess, if the capacity of the knapsack is still positive, then the dynamic programming process can start. If the capacity is negative, then the knapsack problem is infeasible and the resolution is stopped here. At the end of the dynamic programming process, the total cost is the cost outputted plus the preprocess cost and the items in the pattern are the items used in the knapsack solution plus the items set to 1 during the preprocess step.

To solve dynamically the knapsack problem considering a capacity $C^{knapsack}$, and items with sizes $s_i^{knapsack}$ and costs $p_i^{knapsack}$ for $i = 1, \dots, I^{knapsack}$, we have to introduce $t_{i,c}^{knapsack}$, the maximal cost of a knapsack problem with cost c and with i items among all the items. We can apply the following dynamic programming algorithm and backtracking to find the optimal cost of the knapsack and the items involved in this optimal cost :

Algorithm 2: Knapsack problem

```

for  $i = 1, \dots, I^{knapsack}$  do
  for  $c = 1, \dots, C^{knapsack}$  do
    if  $s_i^{knapsack} > c$  then
       $t_{i+1,c} \leftarrow t_{i,c}$ 
    else
       $t_{i+1,c} \leftarrow \max\{t_{i,c}; p_i^{knapsack} + t_{i,c-s_i^{knapsack}}\}$ 
    end
  end
end
end

```

Algorithm 3: Knapsack problem, backtracking

```

Pattern  $\leftarrow \emptyset$ 
 $c = C^{knapsack}$ 
for  $i = I^{knapsack}, \dots, 1$  do
  if  $t_{i,c} \neq t_{i-1,c}$  then
    Pattern  $\leftarrow$  Pattern  $\cup i$ 
     $c \leftarrow c - s_i^{knapsack}$ 
  end
end
end

```

The optimal cost is $t_{I^{knap}, C^{knap}}$ and the variable *Pattern* contains the items involved in the optimal cost.

We can observe that the complexity of the dynamic program is $o(I^{knap}C^{knap})$. Although it can be way better than solving the subproblem using a solver, some problems can happen when the capacity of the knapsack is large. The dynamic program is not only size-dependant but also data-dependant in the way that two problems with the same number won't be solved in the same amount of time. A problem with few items but with a huge capacity can be slower to solve than a problem with more items. In practice, it could be interesting to measure the solving time per number of items while using the solver and the solving time per number of items and capacity using dynamic programming in order to chose which solving method to chose. Furthermore, as the dynamic programming method is implemented "by hand" in this project while the solver have already been developed and optimized, the solver can still be faster in practice than the dynamic program even if the theoretical complexity is worst.

3.3 Node processing algorithm

To make things simpler to understand in the implementation, we only use on type of node for both branching method. A node always contains some branching sets but in the case of the Ryan & Foster method, it has only on branching set with no rule and a coefficient equal to B as we only have one subproblem per node. A node also has a local lower bound and a list of all up and down branching previously made in its branch in order to constructs the new constraints for the Ryan & Foster method (but also to debug the BnP when needed). In the following, π is the dual variable associated to the constraint of (RMBPP) and σ are the constraints associated to each branching subset (only one in the case of Ryan & Foster method). Whatever the method set to process the node, the algorithm is always the same :

Algorithm 4: Node process

```

// Initialization
nodePool ← filterColumnPool()
 $\pi, \sigma$ , solution, value ← solveMaster()
nodeLb ←  $\sum \pi_i$ , nodeUb ← value, minReducedCost ← 0
// Master problem and subproblems are sequentially solved
while true do
    if solution is integer then
        Update global UB if necessary
        Prune branches with a larger LB than the new UB
        break
    end
    for s in branching subsets do
        for r in branching rules of r do
            reducedCost, column ← solveSubproblem( $\pi, \sigma_s$ )
            if reducedCost <  $\infty$  then
                Update minReducedCost
                if reducedCost < 0 then
                    Add column to the node pool and the column pool
                    Update the master problem data with the new column
                end
                nodeLb = nodeLb + reducedCost +  $\sigma_s$ 
            else
                Node is pruned by infeasibility, break
            end
        end
    end
    if nodeLb  $\simeq$  nodeUb then
        Return the solution of the master problem (node is infeasible if it
        contains the artificial column), break
    end
     $\pi, \sigma$ , solution, value ← solveMaster()
end

```

In practice, a precision error is allowed when testing $\text{node}_l \simeq \text{node}_u$. for the Ryan & Foster method, the loop on s and r are only passed once as the node for the Ryan & Foster method has only one branching set with an empty branching rule.

4 Heuristics

Heuristics are proceeded during the algorithm in order to speed-up the resolution time by finding feasible solutions in order to cut branches faster.

4.1 Root heuristics

A first heuristic can be proceeded at root. This is a very important if we want to have a fast algorithm because the higher a branch is cut, the less node will be explored. If a very good bound is found at the root, it can prevent to explore a huge part of the tree.

The three heuristics which can be used at root are three variations of a more general heuristic called Decreasing Order Heuristic [Joh73]. The key idea is to sort the items in decreasing order and for each item, choose a bin to pack the object. The heuristics differs on the choice of the bin to pack the object. The general Decreasing Order Heuristic can be proceeded as following.

Algorithm 5: Decreasing Order Heuristic

```
while The are items that have not been packed do
    Pick the biggest item
    if The item can't be packed in any bin then
        | Add a new bin
    end
    Choose a bin able to pack the item
    Put the item in the bin
end
```

The three heuristics implemented are the First-Fit (FFD), the Best-Fit (BFD) and the Worst-Fit (WFD) algorithm and they select the bin to pack the item as following :

- **FFD** : Chooses the first bin in which the item can be packed
- **BFD** : Chooses the bin with the least amount of free space in which the item can be packed
- **WFD** : Chooses the bin with the most amount of free space in which the item can be packed

As all these heuristics have to sort the items (in $\mathcal{O}(N \log N)$) and to proceed a loop over each of the items (in $\mathcal{O}(N)$), their overall complexity is $\mathcal{O}(N \log N)$. Furthermore, if B^* is the optimal number of bins, it was shown that the result in the worst case for these heuristics are

- **FFD** : $\lfloor 1.7B^* \rfloor$ [DS13]
- **BFD** : $\lfloor 1.7B^* \rfloor$ [DS14]
- **WFD** : $2B^* - 2$ [Joh73]

We can see that these heuristics are very interesting both regarding to their complexity and their efficiency. It is very recommended to enable it when running the BnP algorithm. There exist many other heuristics which can be found in the [Wikipedia page of the BPP](#).

4.2 Tree heuristics

Dual bounds found through the tree are proven to be quite good [SV13]. Thus, an efficient primal heuristic allowing to get good primal bounds can lead to a very successful algorithm. Three different types of heuristics are implemented. The first type relies on the MIRUP property of the (BPP) [DIM16]. The second type of heuristics are described in [WG96] and are base on different rounding strategies, starting from the solution of the relaxed restricted master of a node. The third type of heuristics are diving heuristics described in [SV13] for the case of the Bin-Packing with Conflicts.

4.2.1 MIRUP-based strategy

The MIRUP property of the (BPP) is a conjecture which is still open. It has been conjectured [ST95] that given the solution $z_{\mathbb{C}\mathbb{R}}^*$ the solution of the continuous relaxation of (SCBPP), the following inequality holds :

$$z^* \leq \lceil z_{\mathbb{C}\mathbb{R}}^* \rceil + 1 \quad (4.2.1)$$

where z^* is the solution of (SCBPP) with the integrity constraint. $\lceil z_{\mathbb{C}\mathbb{R}}^* \rceil + 1$ is called the MIRUP bound. The idea is first to obtain the solution of (RMBPP) at each node and compute the MIRUP bound. Then, the restricted master problem is solved with integrity constraint and with a new constraint :

$$\sum_{p \in \mathcal{P}'} \alpha^p \leq \lceil z_{\mathbb{C}\mathbb{R}}^* \rceil + 1$$

If the solution is feasible and outperform the current best solution, then the best solution is updated. Adding this constraint allow to reduce drastically the search space for an integer solution and ensure that UB and LB will have only one unity of difference. However, the solving process can still be long as we solve an integer problem instead of a relaxed problem.

4.2.2 Rounding strategies

In 1995, Wäscher and Gau have presented several rounding strategies grouped in three different categories. This first group is called the Basic Pattern Approach and relies directly on the solution of the relaxed master problem. Method of this group have a name starting with **B**. The second group is called the Residual Pattern Approach and is based on the solution of the relaxed master problem where non-integer components have been rounded-down to zero. As this modification may lead to an infeasible solution, the methods of this groups aims to construct a feasible solution solving a Residual Problem (find new columns to add to the rounded-down solution to create a feasible solution). Method of this group have a name starting with **R**. Finally, the third group is called the Composite Approach and is a mix of the two first groups. Method of this group have a name starting with **C**.

Several method are proposed in [WG96] and methods **RSUC** and **CSTAOPT** are shown to be the most effective. Only the following methods has been implemented.

Procedure BRUSIM The simplest procedure consists of simultaneously rounding-up any non-integer component of the (RMBPP) to one. The advantages of this procedure are obvious : it is extremely fast and immediately result in a feasible solution. However, it produces very bad primal bounds.

Procedure BRURED Rounding-up the non-integer components simultaneously often creates an over-packaging of some items. Then it may be feasible to remove some patterns without violating the packaging constraint. Neumann and Morlock [Ban00] suggest to check whether a pattern can be eliminated without causing a violation of the packaging constraints.

Procedure BOPT This strategy simply solve (RMBPP) on the node pool with integrity constraints on the α^p . This can be done by any optimal solving procedure. However, it may be very long to run this procedure as the problem to solve has integrity constraints. This method provides quite good upper bounds.

Procedure BRUSUC Solving directly (RMBPP) with integrity constraints can be quite costly in time. The BRUSUC address to this problem by first fixing integer variable in the solution of (RMBPP). Then, (RMBPP) will be re-optimized (still with continuous constraints) until all the variables are integer. Before each re-optimization, the variable with the higher fractional value is fixed to one.

Procedure CSTAOPT A cutting pattern is a pattern containing an item which is packed more than one. The CSTAOPT procedure starts by finding a feasible solution by applying the BRUSUC procedure. Then, it removes iteratively cutting patterns until there are no cutting pattern left. It will remain a residual problem to be solved as removing cutting pattern can lead to an infeasible solution. The residual pattern is solved by an optimal method.

4.2.3 Diving heuristics

[TODO : Finish]

5 Numerical applications

5.1 Numerical results

5.2 Result analysis

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