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Enhancing the ACS Heuristic: Parameter Tuning and Computational Experiments for Solving MIP Problems

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Abstract

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Chapter 1

Introduction

1.1 Mixed Integer Programming (MIP)

Mixed Integer Programming (MIP) is a powerful optimization technique used to model complex real-world problems where discrete decisions are essential—such as resource allocation, scheduling, and logistics. MIP models aim to find the best solution according to a given objective function, which is either maximized or minimized. A generic mixed-integer program (MIP) will be defined as

$$\begin{cases} \min & c^T x \\ \text{s.t.} & Ax = b \\ & x_i \in \mathbb{Z}, \forall i \in \mathcal{I} \\ & l \leq x \leq u \end{cases} \quad (1.1)$$

where $c \in \mathbb{R}^n$, $A \in \mathbb{R}^{m \times n}$ and $\mathcal{I} \subseteq \{1 \dots n\}$ is the subset of integer variables indices. The solution vector x is bounded by $l \in \bar{\mathbb{R}}^n$ and $u \in \bar{\mathbb{R}}^n$ where $\bar{\mathbb{R}} = \mathbb{R} \cup \{-\infty, \infty\}$. Here, some decision variables are constrained to take integer values, which increases the model's complexity.

1.1.1 Heuristic in MIP solving

The presence of integer variables significantly increases the computational complexity of MIP problems, especially as the number of variables and constraints grows. Solvers such as IBM ILOG CPLEX^[1] and GUROBI^[2] incorporate a variety of built-in strategies to improve performance. Among these strategies are heuristics, which are specialized algorithms designed to quickly find feasible (but not necessarily optimal) solutions. Although heuristics do not guarantee optimality, they are computationally efficient and can significantly accelerate the solution process by providing good initial solutions or guiding the search within the solution space.

1.2 Alternating Criteria Search (ACS)

Finding high-quality feasible solutions is a key aspect of the discrete optimization process, and constructing them has been one of the main focuses of research in the last few decades. Primal heuristics differ in whether they require a starting feasible solution or not. In the first case, the heuristics are called starting heuristics, and state-of-the-art techniques rely on Large Neighborhood Search (LNS)^[3] to find high-quality feasible solutions. The effectiveness of such strategies relies on the power of the MIP solver to solve the subproblems generated during the LNS heuristic. Improvement heuristics, in contrast, require a feasible starting solution and aim to improve its quality with respect to the objective. The simplest improvements are the 1-opt and 2-opt^[4] methods, but—as with starting heuristics—there are many improvement heuristics based on LNS ideas. The neighborhoods explored are defined based on branch-and-bound information—such as the best incumbent and the LP relaxation. This is essentially a limitation in terms of diversification: the strategies may reach good solutions, but it is difficult to find them in the early stages of the search. Therefore, to address this issue, the Alternating Criteria Search (ACS)^[5]—or Parallel ACS, its straightforward parallel implementation—can be utilized. The search neighborhoods are defined based on randomization instead of branch-and-bound information, allowing for a wider exploration of the search space. The aim of this strategy is to find high-quality feasible solutions at the beginning of the search, increasing the heuristic’s effectiveness.

1.2.1 Parallelization of ACS

Parallelism can be exploited in MIP optimization to accelerate computation and enhance solver performance, in terms of execution time and scalability. The exploration of the branch-and-bound tree in parallel is now incorporated in most state-of-the-art MIP solvers—such as CPLEX or GUROBI—and it essentially consists of solving multiple subproblems of the original MIP simultaneously. However, this strategy may not scale well to a large number of cores^[6]. To address this limitation, the proposed heuristic—Parallel ACS (PACS)—is a parallel algorithm that integrates elements of both starting solution and improvement heuristics, offering the capability of generating starting solutions and improving them with respect to the original objective. To leverage parallelism, Parallel ACS performs a large number of LNS runs simultaneously over a diversified set of neighborhoods, with the aim of increasing the chances of finding higher-quality solutions. After each LNS exploration, the strategy consolidates the local improvements through a recombination phase. PACS combines parallelism and diversified LNS in order to address large instances of MIPs arising from various application domains.

1.2.2 Limitations of the PACS Approach

Although the experiments about the PACS strategy bring evidence about the effectiveness on hard MIP instances, there are several considerations that must be addressed. First of all, the PACS algorithm was executed on an 8-node computing cluster, each equipped with two Intel Xeon X5650 6-core processors and 24 GB of memory, totaling 96 cores and 192 GB of RAM. Such computing resources are not typically available in consumer-grade or general-purpose environments. Furthermore, as discussed in the original study, the PACS strategy could assist the MIP solver in accelerating convergence. This suggests that integrating PACS heuristics within a MIP solver could provide a meaningful contribution to solver performance. Second, before the PACS execution, an initial calibration phase is required to tune the parameters prior to execution. This undermines the objective of designing a self-contained, efficient heuristic. With these considerations in mind, the goal of this thesis is to refine the PACS strategy to make it more compatible with MIP solvers. In particular, the aim is to establish a fixed, robust parameter configuration that remains effective regardless of the input instance.

Chapter 2

The ACS Framework

The Alternating Criteria Search (ACS) heuristic is designed to pursue a twofold objective: to identify a feasible solution and to subsequently enhance its quality with respect to the objective function $c^T x$. To this end, ACS employs a Large Neighborhood Search (LNS) strategy, which iteratively solves two auxiliary mixed-integer programming (MIP) subproblems in order to address both objectives. The heuristic requires an initial vector, which is not mandated to be a feasible solution for the original MIP. This vector is progressively refined by solving sub-MIPs in which a subset of variables is fixed to the values of the initial input.

2.1 Feasibility-MIP (FMIP)

In linear programming theory, it is well established that the feasibility problem can be addressed using the two-phase Simplex method, in which an auxiliary optimization problem is solved to obtain a feasible starting basis, and hence a feasible solution. In a similar manner, the following auxiliary MIP problem, denoted as the Feasibility-MIP (FMIP), is formulated to identify a feasible starting solution:

$$\left\{ \begin{array}{ll} \min & \sum_{i=0}^m \Delta_i^+ + \Delta_i^- \\ \text{s.t.} & Ax + I_m \Delta^+ - I_m \Delta^- = b \\ & x_i = \hat{x}_i, \forall i \in F \\ & l \leq x \leq u \\ & x_i \in \mathbb{Z}, \forall i \in \mathcal{I} \\ & \Delta^+ \geq 0, \Delta^- \geq 0 \end{array} \right. \quad (2.1)$$

Here, I_m denotes an $m \times m$ identity matrix, while Δ^+ and Δ^- are vectors in R^m corresponding to the m constraints. These are introduced as slack variables, and the objective is to minimize their sum. Analogous to the two-phase Simplex method, a vector x is feasible for the original MIP if and only if it can be extended to a solution of value 0 for

the associated FMIP. Since solving an FMIP is as computationally demanding as solving the original problem, the neighborhoods are restricted by fixing a given subset F of the integer variables to the values of an input vector $[\hat{x}, \Delta^+, \Delta^-]$. Because of the introduction of slack variables, the vector \hat{x} is not required to be a feasible solution itself, but it must be integral and within the variable bounds to preserve the feasibility of the model. In this way, the FMIP guarantees that feasibility is preserved under any arbitrary variable-fixing scheme. Once a feasible solution is obtained, any LNS-based improving heuristic—such as RINS^[7], DINS^[8], or local branching^[9]—can then be applied to refine the solution.

2.2 Optimality-MIP (OMIP)

Rather than executing the FMIP until convergence to a feasible solution vector, the following auxiliary MIP problem, denoted as Optimality-MIP (OMIP), is designed to improve a partially feasible solution $[\hat{x}, \hat{\Delta}^+, \hat{\Delta}^-]$, which satisfies $\sum_{i=1}^m (\hat{\Delta}_i^+ + \hat{\Delta}_i^-) \neq 0$, with respect to the original objective $c^T x$:

$$\left\{ \begin{array}{ll} \min & c^T x \\ \text{s.t.} & Ax + I_m \Delta^+ - I_m \Delta^- = b \\ & \sum_{i=0}^m \Delta_i^+ + \Delta_i^- \leq \sum_{i=0}^m \hat{\Delta}_i^+ + \hat{\Delta}_i^- \\ & x_i = \hat{x}_i, \forall i \in F \\ & l \leq x \leq u \\ & x_i \in \mathbb{Z} \quad \forall i \in \mathcal{I} \\ & \Delta^+ \geq 0, \Delta^- \geq 0 \end{array} \right. \quad (2.2)$$

Analogous to the FMIP, the OMIP represents a reformulation of the original MIP model in which auxiliary slack variables are introduced for each constraint. This formulation enables the OMIP to enhance a solution even if it is not feasible for the original MIP. Moreover, to ensure that the optimal solution of the OMIP does not exceed the infeasibility of the input solution \hat{x} , an additional budget constraint is imposed, limiting the total slack to $\sum_{i=0}^m \hat{\Delta}_i^+ + \hat{\Delta}_i^-$.

By iteratively solving subproblems of both auxiliary MIPs, the ACS heuristic is designed to converge—although convergence is not formally guaranteed—to a high-quality feasible solution. By construction, infeasibility decreases monotonically after each iteration. However, the quality of the solution with respect to the original objective function may fluctuate.

2.3 Parallelization of ACS

The parallelization of ACS exploits parallelism by generating a diversified set of large neighborhood searches, which are solved simultaneously. Exploring multiple search neighborhoods in parallel is expected to increase the likelihood of identifying high-quality solutions. Following this parallelization step, the improvements obtained in parallel must be combined efficiently. To this end, an additional search subproblem is generated in which variables with identical values across different solutions are fixed. Consequently, the recombination phase constitutes a crucial step in PACS: by merging the improvements achieved during the parallel phase, the subsequent phase can explore a newly diversified set of large neighborhood searches based on the recombined solution, thereby enhancing the probability of further improvements. The pseudocode[1] illustrates the overall workflow. Each processor generates a set of randomized variable fixings and solves the associated

Algorithm 1 Parallel Alternating Criteria Search

```

function FMIP_LNS( $\mathcal{F}, \hat{x}$ )
    return  $\min\{\sum_i \Delta_i^+ + \Delta_i^- \mid Ax + I_m \Delta^+ - I_m \Delta^- = b, x_j = \hat{x}_j \forall j \in F, x_j \in \mathbb{Z} \forall j \in \mathcal{I}\}$ 
end function

function OMIP_LNS( $\mathcal{F}, \hat{x}, \hat{\Delta}$ )
    return  $\min\{c^t x \mid Ax + I_m \Delta^+ - I_m \Delta^- = b, \sum_i \Delta_i^+ + \Delta_i^- \leq \hat{\Delta}, x_j = \hat{x}_j \forall j \in F, x_j \in \mathbb{Z} \forall j \in \mathcal{I}\}$ 
end function

Ensure: Feasible solution  $\hat{x}$  if found
1: Initialize  $[\hat{x}, \Delta^+, \Delta^-]$  as an integer solution
2:  $T := numThreads()$ 
3: while  $timeElapsed() \leq timeLimit$  do
4:   if  $\sum_i \Delta_i^+ + \Delta_i^- > 0$  then
5:     for all threads  $t_i \in \{0, T-1\}$  in parallel do
6:        $F_{t_i} :=$  randomized variable index subset,  $F_{t_i} \subseteq \mathcal{I}$ 
7:        $[x^{t_i}, \Delta^{+t_i}, \Delta^{-t_i}] := \text{FMIP\_LNS}(F_{t_i}, \hat{x})$ 
8:     end for
9:      $U := \{j \in \mathcal{I} \mid x_j^{t_i} = x_j^{t_k}, 0 \leq i < k < T\}$ 
10:     $[\hat{x}, \Delta^+, \Delta^-] := \text{FMIP\_LNS}(U, x^{t_0})$ 
11:   end if
12:    $\Delta^{UB} := \sum_i \Delta_i^+ + \Delta_i^-$ 
13:   for all threads  $t_i \in \{0, T-1\}$  in parallel do
14:      $F_{t_i} :=$  randomized variable index subset,  $F_{t_i} \subseteq \mathcal{I}$ 
15:      $[x^{t_i}, \Delta^{+t_i}, \Delta^{-t_i}] := \text{OMIP\_LNS}(F_{t_i}, \hat{x}, \Delta^{UB})$ 
16:   end for
17:    $U := \{j \in \mathcal{I} \mid x_j^{t_i} = x_j^{t_k}, 0 \leq i < k < T\}$ 
18:    $[\hat{x}, \Delta^+, \Delta^-] := \text{OMIP\_LNS}(U, x^{t_0}, \Delta^{UB})$ 
19: end while
20: return  $[\hat{x}, \Delta^+, \Delta^-]$ 

```

sub-MIP—either FMIP or OMIP—until the allotted time limit is reached. Subsequently,

the solutions are exchanged, and the set U , containing the indices of variables common across solutions, is constructed. The recombination MIP then consists of a subproblem, again associated with either FMIP or OMIP, in which the variables in U are fixed. The best solution obtained will be either the most feasible or the most optimal, depending on whether a recombination FMIP or OMIP is employed. Moreover, every solution used as input can be incorporated as a MIP start, as it remains feasible under any variable-fixing strategy. Processor synchronization and memory communication are handled via the Message Passing Interface (MPI)^[10], owing to its efficient all-to-all collective communication primitives in distributed large-scale architectures.

2.4 Initialization of ACS

As introduced earlier in this chapter, ACS only requires a starting vector that is integer feasible and within the variable bounds. However, a stronger starting point is a solution that is as feasible as possible with respect to the objective function of the FMIP. The proposed algorithm provides a lightweight heuristic that seeks to minimize the infeasibility of the initial solution. This algorithm can be described by the pseudocode[2]. The algorithm first sorts the list of integer variables in order of increasing bound range. It

Algorithm 2 Starting vector heuristic

Require: Percentage of variables to fix θ , $0 < \theta \leq 100$, Fixed bound constant c_b

Ensure: Starting integer-feasible vector \hat{x}

```

1:  $V :=$  list of integer variables sorted by increasing bound range  $u - l$ 
2:  $F := \emptyset$ 
3: while  $\hat{x}$  is not integer feasible AND  $F \neq \mathcal{I}$  do
4:    $\mathcal{K} :=$  top  $\theta\%$  of unfixed variables from  $V$ 
5:   for  $k \in \mathcal{K}$  do
6:      $\hat{x}_k :=$  random integer value between  $[\max(l_k, -c_b), \min(u_k, c_b)]$ 
7:   end for
8:    $F := F \cup \mathcal{K}$ 
9:    $[x, \Delta^+, \Delta^-] := \min\{\sum_i \Delta_i^+ + \Delta_i^- \mid Ax + I_m \Delta^+ - I_m \Delta^- = b, x_j = \hat{x}_j \ \forall j \in F\}$ 
10:   $Q :=$  index set of integer variables of  $x$  with integer value
11:   $\hat{x}_q = x_q, \ \forall q \in Q$ 
12:   $F := F \cup Q$ 
13: end while
14: return  $\hat{x}$ 

```

then fixes the top $\theta\%$ of variables to random integer values within their respective bounds. The input parameter θ controls the trade-off between the difficulty of the LP relaxation and the quality of the resulting starting solution. In cases where the bounds are infinite, a constant value of 10^6 is used to clamp the bounds. The rationale behind the sorting step is to prioritize binary variables first, followed by the remaining integer variables. Until all integer variables are fixed, the LP relaxation of the FMIP is solved to optimize the unfixed variables. Any variables that attain integer values in this process are then fixed.

Since at least $\theta\%$ of the variables are fixed at each iteration, the algorithm is guaranteed to terminate after at most $\lceil 100/\theta \rceil$ iterations.

2.5 Variable Fixing Strategy

Selecting an appropriate variable fixing scheme is a challenging task: an overly restrictive strategy may fail to yield improvements, whereas an excessively loose strategy can lead to a search space that is too large to explore efficiently within a reasonable timespan. The proposed algorithm constitutes a simple yet intuitive variable fixing method. It incorporates randomness to promote diversification and allows for controlling the number of variables to be fixed through an adjustable parameter. The algorithm is described in the pseudocode[3]. The fixings are determined by selecting a random integer variable

Algorithm 3 Variable Fixing Selection Algorithm

Require: Fraction of variables to fix ρ , $0 < \rho < 1$

Ensure: Set of integer indices F

1: **function** RANDOMFIXINGS(ρ)

2: $i :=$ random element in \mathcal{I}

3: $F :=$ first $\rho \cdot |\mathcal{I}|$ consecutive integer variable indices starting from i in a circular fashion

4: **return** F

5: **end function**

x' and fixing a consecutive sequence of integer variables starting from x' up to a cap determined by ρ , an input parameter that specifies the number of variables to be fixed. The fixing is performed in a circular fashion: if the end of the set \mathcal{I} is reached before the required number of variables are fixed, the algorithm continues from the beginning of \mathcal{I} . The effectiveness of this strategy relies on the fact that, for many problems—such as network flow and routing—the variables are arranged consecutively, often defining a cohesive substructure within the problem.

Chapter 3

The PACS Tuning

3.1 PACS with Generalized Fixing

The baseline PACS algorithm enforces both the starting vector construction and the fixing scheme to operate exclusively on the set \mathcal{I} of integer variables. While this restriction may appear efficient and straightforward, it can, in fact, be limiting. In particular, if the MIP instance contains only a small fraction of integer variables relative to the total, focusing solely on them may reduce diversification and cause the algorithm to converge prematurely to a local minimum, which is undesirable in an optimization process. To overcome this issue, Algorithm 2 and Algorithm 3 are generalized into Algorithm 4 and Algorithm 5, respectively, thereby allowing both integer and continuous variables to be considered.

Algorithm 4 Generalized Starting vector heuristic

Require: Percentage of variables to fix θ , $0 < \theta \leq 100$, Fixed bound constant c_b

Ensure: Starting integer-feasible vector \hat{x}

```
1:  $V :=$  list of integer variables sorted by increasing bound range  $u - l$ 
2:  $F := \emptyset$ 
3: while  $\hat{x}$  is not integer feasible AND  $F \neq V$  do
4:    $\mathcal{K} :=$  top  $\theta\%$  of unfixed variables from  $V$ 
5:   for  $k \in \mathcal{K}$  do
6:      $\hat{x}_k :=$  random integer value between  $[\max(l_k, -c_b), \min(u_k, c_b)]$ 
7:   end for
8:    $F := F \cup \mathcal{K}$ 
9:    $[x, \Delta^+, \Delta^-] := \min\{\sum_i \Delta_i^+ + \Delta_i^- \mid Ax + I_m \Delta^+ - I_m \Delta^- = b, x_j = \hat{x}_j \ \forall j \in F\}$ 
10:   $Q :=$  index set of integer variables of  $x$  with integer value
11:   $\hat{x}_q = x_q, \ \forall q \in Q$ 
12:   $F := F \cup Q$ 
13: end while
14: return  $\hat{x}$ 
```

Algorithm 5 Generalized Variable Fixing Selection Algorithm

Require: Fraction of variables to fix ρ , $0 < \rho < 1$

Ensure: Set of integer indices F

```
1: function RANDOMFIXINGS( $\rho$ )  
2:    $i :=$  random element in  $\{1 \dots n\}$      $\triangleright n$ : number of variables in the original MIP  
3:    $F :=$  first  $\rho \cdot n$  consecutive integer variable indices starting from  $i$  in a circular  
   fashion  
4:   return  $F$   
5: end function
```

Since these algorithms are executed on the auxiliary MIP problems -FMIP or OMIP-, the variables subject to fixing correspond exactly to those defined in the original MIP formulation. By generalizing the fixing strategy to include both integer and continuous variables, diversification is enhanced, thereby reducing the likelihood of stagnation in local minima and potentially improving the exploration of the solution space.

3.2 Architecture-Agnostic Parallelization

In the original study, the Message Passing Interface (MPI) was employed to synchronize processors at each recombination phase. While this approach is well suited to high-performance computing environments, it may be unnecessarily complex in general-purpose scenarios, where a simpler multi-threading implementation is often preferable. In this thesis, communication is instead managed through a set of logical threads, which may differ from the number of available hardware threads. This abstraction ensures that, even on machines with fewer physical cores, the algorithm can reproduce the same behavior across different architectures, provided sufficient computational time is allowed. More specifically, during the coordination phase, either the most feasible or the most optimal solution—depending on whether a recombination FMIP or OMIP is performed—is shared among the logical processors. Each processor then continues working independently on its own copy, with updates to the incumbent solution handled exclusively through a thread-safe update function, formally described in Algorithm 6.

Algorithm 6 Parallel ACS Incumbent Update Procedure

Require: Candidate solution x with slack sum $S(x)$ and objective value $C(x)$; Incumbent \tilde{x} ; Zero-tolerance ϵ

Ensure: Updated incumbent \tilde{x} in a thread-safe manner

```
1: function UPDATEINCUMBENT( $x$ )  
2:   acquire lock  
3:   if ( $|S(x)| < |S(\tilde{x})| \vee (|S(x)| < \epsilon \wedge C(x) < C(\tilde{x}))$ ) then  
4:      $\tilde{x} \leftarrow x$   
5:   end if  
6:   release lock  
7: end function
```

This mechanism is crucial to ensure that the algorithm consistently improves and converges within the given time limit. The incumbent solution is updated whenever a better solution is identified: either one with a smaller slack sum, indicating improved feasibility, or one with a lower objective value $c^T x$ provided that the slack sum is less than the tolerance parameter ϵ , the zero-feasibility threshold.

3.3 Eliminating Calibration in PACS Parameter Selection

After adapting the PACS algorithm to a more general-purpose environment, another important challenge arises: parameter selection. Since PACS must be applicable to a wide range of hard MIP instances, it is necessary to identify parameter settings that generally perform well, both in terms of computational efficiency and heuristic solution quality. The goal is to remove the need for an explicit calibration phase, while still ensuring reliable performance. The tuning process specifically concerns the following parameters:

1. The time span assigned to each sub-MIP, either FMIP_LNS or OMIP_LNS
2. The parameter ρ governing the fixing strategy
3. The parameter θ governing the starting vector

3.3.1 Adaptive Determination of Sub-MIP Time Span

To guarantee determinism in the implementation, each sub-MIP is assigned both a time limit equal to the remaining computation time and a deterministic time limit. The latter is defined as the maximum number of instructions that the solver may execute before termination. The deterministic time limit is computed according to the following formula:

$$TL_{DET} = \max \left(x, \min \left(\frac{nz}{y}, X \right) \right) \quad (3.1)$$

which provides a dynamic mechanism for adapting the computational effort of each sub-MIP. In this formulation, x and X denote the minimum and maximum allowable deterministic time limits, respectively, nz represents the number of nonzeros in the constraint matrix A of the MIP problem, and y acts as a scaling factor. The chosen parameter values are:

1. Minimum deterministic time limit: $x = 10^3$
2. Scaling factor: $y = 10^2$
3. Maximum deterministic time limit: $X = 10^7$

This dynamic adjustment removes the need to explicitly set a fixed time limit for each subproblem, thereby eliminating the necessity of parameter tuning in this respect.

3.3.2 Adaptive Variable Fixing through ρ Adjustment

Fixed ρ Initialization

As discussed in Section 2.5, selecting an appropriate variable fixing scheme is a non-trivial task. In particular, the random fixings scheme presents additional challenges: if the fraction ρ of variables to be fixed is set too high, the procedure may fail to yield meaningful improvements; conversely, if ρ is set too low, the resulting search space may become excessively large, making it computationally intractable within a reasonable time frame.

For this reason, in the first instance, the parameter ρ is selected from a set of predetermined candidate values. Anticipating the results presented in Section ??, it can be observed that certain values of ρ are more effective in terms of solution quality, as they grant the solver greater flexibility during the optimization process.

Dynamic Adjustment of Parameter ρ

Since the LNS heuristics in the PACS algorithm restrict the search space by fixing a number of variables according to the value of ρ , it is natural to design a mechanism for dynamically adapting this parameter in order to increase the likelihood of discovering high-quality solutions.

Algorithm 7 Parallel ACS Rho Update (Parallel Phases)

Require: Status code MIP_{code} returned by the solver; Adjustment step Δ_ρ for ρ ; The variable fixing parameter ρ ; Number of parallel sub-MIPs num_{MIP}

Ensure: Updated value of ρ , synchronized across parallel sub-MIPs

```
1: function DYNRHOUPDATE( $MIP_{code}$ ,  $\Delta_\rho$ ,  $\rho$ ,  $num_{MIP}$ )
2:    $\hat{\Delta}_\rho \leftarrow \frac{\Delta_\rho}{num_{MIP}}$ 
3:   acquire lock
4:   if  $MIP_{code} = OPT \vee MIP_{code} = OPT_{TOL}$  then
5:      $\rho \leftarrow \rho - \hat{\Delta}_\rho$ 
6:   end if
7:   if  $MIP_{code} = FEAS_{TL} \vee MIP_{code} = FEAS_{DET.TL}$  then
8:      $\rho \leftarrow \rho + \hat{\Delta}_\rho$ 
9:   end if
10:  Clap  $\rho$  within  $[0.01, 0.99]$ 
11:  if *Tie Case detected* then
12:     $\rho \leftarrow \rho - \Delta_\rho$ 
13:  end if
14:  release lock
15: end function
```

The procedure, described in Algorithm 7, is applied after each sub-MIP optimization phase in the parallel step. Based on the status returned by the solver, the value of ρ is updated as follows:

- If the solver hits the time limit—either the deterministic limit or the global remaining time—while still producing a feasible solution, ρ is increased by $\frac{\Delta_\rho}{num_{MIP}}$. This adjustment suggests fixing more variables in subsequent phases, thereby simplifying the subproblem to be solved.
- Conversely, if the solver converges to an optimal solution within the tolerance, this indicates that the corresponding region of the search space has already been sufficiently explored. In this case, ρ is decreased by $\frac{\Delta_\rho}{num_{MIP}}$, enlarging the search space and granting the solver greater freedom in the following optimization steps.

Since each sub-MIP independently attempts to modify the value of ρ , synchronization through locking is required to prevent inconsistencies. The final update is determined as the average of the adjustments proposed by the parallel optimization phases. In case of a tie, a deterministic rule is applied: ρ is decreased by Δ_ρ .

Algorithm 8 Parallel ACS Rho Update (Recombination Phases)

Require: Status code MIP_{code} returned by the solver; Adjustment step Δ_ρ for ρ ; The variable fixing parameter ρ ; Number of parallel sub-MIPs num_{MIP}

Ensure: Updated value of ρ after recombination adjustment

```
1: function DYNRHOUPDATE( $MIP_{code}$ ,  $\Delta_\rho$ ,  $\rho$ ,  $num_{MIP}$ )
2:    $\hat{\Delta}_\rho \leftarrow \frac{2\Delta_\rho}{num_{MIP}}$ 
3:   if  $MIP_{code} = OPT \vee MIP_{code} = OPT_{TOL}$  then
4:      $\rho \leftarrow \rho - \hat{\Delta}_\rho$ 
5:   end if
6:   if  $MIP_{code} = FEAS_{TL} \vee MIP_{code} = FEAS_{DET.TL}$  then
7:      $\rho \leftarrow \rho + \hat{\Delta}_\rho$ 
8:   end if
9:   Clap  $\rho$  within  $[0.01, 0.99]$ 
10: end function
```

For the recombination phase, the procedure is slightly modified into Algorithm 8, where the adjustment step is doubled, i.e. $2\Delta_\rho/num_{MIP}$, in order to resolve potential tie cases.

Finally, although an initial value of ρ must be provided to start the fixing process, experimental results in Section ?? show that performance is only marginally affected by this initialization. Consequently, the effectiveness of the method does not critically depend on the initial choice of ρ .

3.3.3 Parameter-Free Strategy for Initial Solution Construction

Another non-trivial challenge concerns the selection of the initial solution vector. The parameter θ , which regulates the trade-off between solution quality and execution time, must be tuned to suit the most general cases. Similar to the fixing strategy, the analysis considers a set of predetermined values for θ , while also introducing a novel initialization strategy for the starting solution vector. The motivation is that using a small value of θ —which corresponds to generating a high-quality initial solution—may waste computation time, since the true strength of PACS lies in its diversification and parallelization capabilities.

Therefore, Algorithm 9 is designed with the following goal: to provide a heuristic starting solution that is deterministic and computationally lightweight, while remaining reasonably feasible for the original MIP problem.

Algorithm 9 Heuristic Initialization of Starting Solution

Require: *MIP* instance with n variables, objective coefficients c_i , bounds $[l_i, u_i]$; zero-tolerance ϵ

Ensure: Starting solution vector $x \in \bar{\mathbb{R}}^n$

```
1: function STARTSOLMAXFEAS(MIP)
2:    $x \leftarrow (+\infty, \dots, +\infty) \in \bar{\mathbb{R}}^n$ 
3:   for  $i = 1 \rightarrow n$  do
4:      $(l_i, u_i) \leftarrow$  variable bounds of variable  $i$ 
5:     if  $l_i = -\infty \wedge u_i = +\infty$  then
6:        $x_i \leftarrow 0$ 
7:     else if  $l_i > -\infty \wedge u_i = +\infty$  then
8:        $x_i \leftarrow l_i$ 
9:     else if  $l_i = -\infty \wedge u_i < +\infty$  then
10:       $x_i \leftarrow u_i$ 
11:     else
12:       if  $c_i \leq -\epsilon$  then
13:          $x_i \leftarrow l_i$ 
14:       else if  $c_i \geq \epsilon$  then
15:          $x_i \leftarrow u_i$ 
16:       else
17:          $x_i \leftarrow \text{RANDOMINTEGER}(l_i, u_i)$ 
18:       end if
19:     end if
20:   end for
21:   return  $x$ 
22: end function
```

The initialization procedure in Algorithm 9 operates as follows:

- If a variable has no finite bounds, it is set to 0.
- If a variable has only one finite bound, it is set to that bound.
- If both bounds are finite:
 - set the variable to the lower bound if $c_i \leq -\epsilon$, i.e. the coefficient is negative,
 - set it to the upper bound if $c_i \geq \epsilon$, i.e. the coefficient is positive,
 - choose a random integer within $[l_i, u_i]$ if $-\epsilon \leq c_i \leq \epsilon$, i.e. the coefficient is close to zero.

In this way, the solution initialization is almost entirely deterministic and computationally lightweight, although feasibility with respect to the original MIP problem may be partially

compromised. Furthermore, the proposed strategy is parameter-free, which makes it particularly well-suited for the purposes of this thesis.

3.4 Additional PACS Optimizations

While the previous modifications focused on tuning or slightly adjusting specific components of the PACS workflow, the following procedures—analyzed in Section ??—aim to further enhance the quality of the final solution.

3.4.1 Slack Upper Bound Enforcement and Budget Constraint Removal

The first idea is to exploit the values of the slack variables Δ^+ and Δ^- obtained in a previous iteration to enforce these values as upper bounds. In the PACS algorithm, this mechanism is already embedded in the budget constraint $\sum_{i=0}^m \Delta_i^+ + \Delta_i^- \leq \sum_{i=0}^m \hat{\Delta}_i^+ + \hat{\Delta}_i^-$, which provides only a global upper bound. The proposed strategy, instead, aims at fixing the upper bounds of individual slack variables, thereby assisting subsequent FMIP optimizations and improving feasibility. The procedure is implemented as shown in Algorithm 10.

Algorithm 10 Fixing Slack Variables to Upper Bound

Require: *MIP* model with n variables; Solution x of length $n + 2m$; Zero-tolerance ϵ

Ensure: Updated model with adjusted upper bounds for slack variables

```

1: function SLACKUPPERBOUNDFIXING(MIP,  $x$ )
2:    $S_{UB} \leftarrow \emptyset$  ▷ Set of variables-upper bound to be updated
3:   for  $i = n \rightarrow n + 2m$  do
4:      $u_i \leftarrow$  upper bound of variable  $i$ 
5:     if  $u_i - x_i > \epsilon$  then
6:        $S_{UB} \leftarrow S_{UB} \cup (i, x_i)$ 
7:     end if
8:   end for
9:   return  $S_{UB}$ 
10: end function

```

Here, the solution vector x corresponds to the current PACS incumbent—valid in both the parallel and recombination phases—and each slack variable is compared with its respective upper bound. If the variable value is smaller than the existing upper bound, the variable index together with its solution value is added to S_{UB} as a pair (i, x_i) . The resulting set S_{UB} is subsequently employed to update the upper bounds of the slack variables, thereby modifying the auxiliary FMIP and OMIP formulations as follows.

FMIP Reformulation

$$\left\{ \begin{array}{ll} \min & \sum_{i=0}^m \Delta_i^+ + \Delta_i^- \\ \text{s.t.} & Ax + I_m \Delta^+ - I_m \Delta^- = b \\ & x_i = \hat{x}_i, \forall i \in F \\ & \Delta_i^+ \leq \hat{u}_i, \forall (i, \hat{u}_i) \in S_{UB}^+ \\ & \Delta_i^- \leq \hat{u}_i, \forall (i, \hat{u}_i) \in S_{UB}^- \\ & l \leq x \leq u \\ & x_i \in \mathbb{Z}, \forall i \in \mathcal{I} \\ & \Delta^+ \geq 0, \Delta^- \geq 0 \end{array} \right. \quad (3.2)$$

OMIP Reformulation

$$\left\{ \begin{array}{ll} \min & c^T x \\ \text{s.t.} & Ax + I_m \Delta^+ - I_m \Delta^- = b \\ & \cancel{\sum_{i=0}^m \Delta_i^+ + \Delta_i^- \leq \sum_{i=0}^m \hat{\Delta}_i^+ + \hat{\Delta}_i^-} \\ & x_i = \hat{x}_i, \forall i \in F \\ & \Delta_i^+ \leq \hat{u}_i, \forall (i, \hat{u}_i) \in S_{UB}^+ \\ & \Delta_i^- \leq \hat{u}_i, \forall (i, \hat{u}_i) \in S_{UB}^- \\ & l \leq x \leq u \\ & x_i \in \mathbb{Z}, \forall i \in \mathcal{I} \\ & \Delta^+ \geq 0, \Delta^- \geq 0 \end{array} \right. \quad (3.3)$$

In these formulations, the set S_{UB} is partitioned into S_{UB}^+ and S_{UB}^- to remain consistent with the notation used for FMIP and OMIP. It is worth noting that the budget constraint in OMIP can now be removed, since the slack upper bound fixing provides a stronger restriction, as will be further confirmed by the results in Section ??.

3.4.2 WalkMIP: Violated Constraints Based Variable Fixing

It's worth to investigate a more sophisticated and hopefully efficient way to fix the variables values, enriching the diversification and consequently increasing the chances of find high-quality solution in a relative tiny time span. For this purpose, the WalkMIP strategy takes the strategy from the WalkSAT^[11] strategy and implement it to directly target the variables that cause the infeasibility of a partially feasible solution. The Algorithm 11 is the suiting of WalkMIP algorithm to cover constraint satisfiability problem.

Algorithm 11 Walk-based Repair Heuristic for MIP

Require: *MIP* model with $n + 2m$ variables (FMIP or OMIP); Initial solution x ; Fixing parameter ρ ; Random walk probability $p \in (0, 1)$; Solution kick probability p_{HUGE_KICK}

Ensure: Updated model with improved feasible solution

```
1: function WALKMIP(MIP,  $x$ ,  $\rho$ ,  $p$ ,  $p_{HUGE\_KICK}$ )
2:    $violConstr \leftarrow$  Set of indices of violated constraints, under  $x$ 
3:   if *First Execution* then  $numInitialConstr \leftarrow |violConstr|$ 
4:   end if
5:   if  $violConstr = \emptyset$  then return
6:   end if
7:   if with probability  $p_{HUGE\_KICK}$  then RANDOMFIXINGS( $\rho$ )
8:   end if
9:    $k \leftarrow \text{NUMVARTOFIX}(numInitialConstr, |violConstr|)$ 
10:  repeat
11:     $c \leftarrow$  Random violated constraint,  $c \in violConstr$ 
12:     $(i_{best}, val_{best}, minDMG) \leftarrow \text{MINDAMAGEMOVE}(c, x)$ 
13:    if  $minDMG \leq -\epsilon$  then
14:       $x \leftarrow \text{UPDATESOL}(x, i_{best}, val_{best})$   $\triangleright$  Apply best move to  $x$ 
15:    else if with probability  $1 - p \wedge i_{best} \neq -1$  then
16:       $x \leftarrow \text{UPDATESOL}(x, i_{best}, val_{best})$   $\triangleright$  Apply min-damage move to  $x$ 
17:    else
18:       $x \leftarrow$  Fix a random var in  $x$ 
19:    end if
20:     $violConstr \leftarrow$  Set of indices of violated constraints, under  $x$ 
21:    if  $violConstr = \emptyset$  then break
22:    end if
23:  until  $k$  times
24:  if  $k < \rho \cdot n$  then
25:    Fix remaining variables (up to  $\rho \cdot n_{mip}$ ) from original  $x$ 
26:  end if
27: end function
28: function MINDAMAGEMOVE( $c, x$ )
29:    $minDMG \leftarrow +\infty$ ,  $i_{best} \leftarrow -1$ ,  $val_{best} \leftarrow +\infty$ 
30:   for each variable  $i$  in  $c$  do
31:     if  $i.type = \text{BIN}$  then  $\hat{x}_i \leftarrow \neg x_i$ 
32:     else if  $i.type = \text{INT}$  then  $\hat{x}_i \leftarrow x_i + \text{RANDOMINTEGER}(-1, 1)$ 
33:     else if  $i.type = \text{DOUBLE}$  then  $\hat{x}_i \leftarrow x_i + \text{RANDOMDOUBLE}(-0.5, 0.5)$ 
34:     end if
35:     Clamp  $\hat{x}_i$  within its bounds  $[l_i, u_i]$ 
36:      $\delta \leftarrow \text{COMPUTEVIOLATION}(x, \hat{x}_i)$ 
37:     if  $\delta < minDMG$  then
38:        $i_{best} \leftarrow i$ ,  $minDMG \leftarrow \delta$ ,  $val_{best}^{18} \leftarrow \hat{x}_i$ 
39:     end if
```


It's important to notice that this procedure itself could only be used in a context in which the slack variable are different from zero and therefore, the original random fixing strategy, presented in the Section 3.3.2, must be still utilized in the OMIP optimization phases in which the slack sum is equal to zero.

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