# 3D Packing of Balls in Different Containers by VNS 

## A thesis submitted for the degree of <br> Doctor of Philosophy

by

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#### Abstract

In real world applications such as the transporting of goods products, packing is a major issue. Goods products need to be packed such that the smallest space is wasted to achieve the maximum transportation efficiency. Packing becomes more challenging and complex when the product is circular/spherical. This thesis focuses on the best way to pack three-dimensional unit spheres into the smallest spherical and cubical space. Unit spheres are considered in lieu of non-identical spheres because the search mechanisms are more difficult in the latter set up and any improvements will be due to the search mechanism not to the ordering of the spheres. The two-unit sphere packing problems are solved by approximately using a variable neighborhood search (VNS) hybrid heuristic. A general search framework belonging to the Artificial Intelligence domain, the VNS offers a diversification of the search space by changing neighborhood structures and intensification by thoroughly investigating each neighborhood. It is flexible, easy to implement, adaptable to both continuous and discrete optimization problems and has been use to solve a variety of problems including large-sized real-life problems. Its runtime is usually lower than other meta heuristic techniques. A tutorial on the VNS and its variants along with recent applications and areas of applicability of each variant. Subsequently, this thesis considers several variations of VNS heuristics for the two problems at hand, discusses their individual efficiencies and effectiveness, their convergence rates and studies their robustness. It highlights the importance of the hybridization which yields near global optima with high precision and accuracy, improving many bestknown solutions indicate matching some, and improving the precision and accuracy of others.

Keywords: variable neighborhood search, sphere packing, three-dimensional packing, meta heuristic, hybrid heuristics, multiple start heuristics.


## Declaration of Originality

I hereby certify that the work presented in this thesis is my original research and has not been presented for a higher degree at any other university or institute.

Signed: $\qquad$ Dated: $\qquad$
Abdulaziz Alkandari

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# Author's Publications 

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## Chapter 1

## Introduction

### 1.1 Background

The optimization of non-linear problem is a classical situation that is frequently encountered in nature. In most cases, finding the global optimum for these mathematical programs is difficult. This is due to the complexity of the topography of the search space and the exorbitant by high computational costs of the existing approaches. Despite the advancement of computational technologies, the computational costs remain excessively high. An alternative to these expensive methods are heuristic approaches which provide good quality solutions in reasonable computational time. There are several classes of heuristic methods. They can be grouped as local search and global search, or as nature-inspired and non-nature-inspired, or as single-start and multiple-start (or population based). The search can itself be a steepest descent/ascent or more elaborate, temporarily accepting, non-improving solutions; or of a prohibiting nature. A heuristic is successful if it balances the intensification and diversification of the search within the neighborhoods of the search
space.
A relatively new heuristic that has proven successful is the variable neighborhood search (VNS). The VNS searches for a (near-) global optimum starting from several initial solutions, and changes the size or structure of the neighborhood of the current local optimum whenever its search stagnates. In other words, it opts for an exploration phase every time its exploitation search fails in improving its current incumbent.

Another option in the search for a global optimizer is hybrid heuristics. These heuristics target overcoming limitations in terms of intensification and diversification through hybridization. For instance, genetic algorithms are known for their diversification whereas simulated annealing and tabu search are notorious for their intensification. Thus, their hybridization has resulted in the resolution of many complex combinatorial optimization problems.

In this thesis a particular non-linear program is addressed using hybrid heuristics inspired from the variable neighborhood search framework. Specifically, it considers the problem of packing three-dimensional unit spheres into three-dimensional containers, where the objective is to minimize the size of the container.

### 1.2 Problem description

This thesis considers packing $n$ identical spheres, of radius one, without overlap into the smallest containing sphere $S$. This problem, is a three-dimensional variant of the Open Dimension Problem: all small items (which are spheres) have to be packed into a larger containers (which should be a sphere or a cube) and the size of the container has to be minimized. The problem is equivalent to finding the coordinates
( $x_{i}, y_{i}, z_{i}$ ) of every sphere $i \in I=1, \ldots, n$, and the dimensions of the container such that every sphere $i \in I$ is completely contained within the object and no pair $(i, j)$ of spheres overlap.

### 1.3 Motivation

The sphere packing problem, which consists of packing spheres into the smallest sphere or cube, has many important real-life applications including materials science, radio surgical treatment, communication, and other vital fields. In materials science, random sphere packing is a model for the structure of liquids, proteins, and glassy materials. The model is used in the study of phenomena such as electrical conductivity, fluid flow, stress distribution and other mechanical properties of granular media, living cells, random media chemistry and physics. The model is also applied in the investigation of processes such as sedimentation, compaction and sintering. In radio surgical treatment planning, sphere packing is crucial to X-ray tomography. In digital communication and storage, it emerges in the packing of compact disks, cell phones, and internet cables. Other applications of sphere packing are encountered in powder metallurgy for three-dimensional laser cutting, in the arranging and loading of containers for transportation, in the cutting different natural by formed crystals, in the layout of computers, buildings, etc. Sphere packing is an optimization problem, but it is debatable whether it should be classified as continuous or discrete. The positions of the spheres are continuous whereas the structure of an optimal configuration is discrete. A successful solution technique should tackle these two aspects simultaneously.

### 1.4 Contribution

Packing unit spheres into three-dimensional shapes is a non-convex optimization problem. It is NP hard (Non-deterministic Polynomial-time hard), since it is an extension of packing unit circles into the smallest two-dimensional shape, which is, in turn, NP hard [32]. Thus, the search for an exact local extremum is time consuming without any guarantee of a sufficiently good convergence to an optimum. Indeed, the problem is challenging. As the number of unit spheres increases, identifying a reasonably good solution becomes extremely difficult. In addition, the problem has an infinite number of solutions with identical minimal radii. In fact, any solution may be rotated or reflected or may have free spheres which can be slightly moved without enlarging the radius of the container sphere. Finally, there is the issue of computational accuracy and numerical precision. Solving the problem via nonlinear programming solvers is generally not successful. Most solvers are not geared towards identifying the global optima. Subsequently, the problem should be solved by a mixture of search heuristics with local exhaustive (exact) searches of the local minima or their approximations. This thesis follows this line of research.

This thesis models the problems as non-linear programs and approximately solves them using a hybrid heuristic which couples a variable neighborhood search (VNS) with a local search (LS). VNS serves as the diversification mechanism whereas LS acts as the intensification one. VNS investigates the neighborhood of a feasible local minimum $(u)$ in search of the global minimum, where neighboring solutions are obtained by shaking one or more spheres of $(u)$ and the size of the neighborhood is varied by changing the number of shaken spheres, and the distance and the direction each sphere is moved. LS intensifies the search around a solution ( $u$ ) by subjecting its neighbors to a sequential quadratic algorithm with a non-monotone line search
(as the NLP solver).
The results and findings extracted from this research are beneficial both to academia and industry. The application of the proposed approach to other problems will allow solving larger instances of other difficult problems. Similarly, its application in an industrial setting will greatly reduce industrial waste when packing spherical objects; thus, reducing the costs, not only in its monetary aspects but also in its polluting aspect. VNS can be applied to other problems with high industrial relevance such as vehicle routing, facility location and allocation, and transportation. Thus, this research can contribute into the mainstream applications of economic and market-oriented packing strategies.

### 1.5 Outline

A tutorial and a detailed survey on VNS methods and applications, including the following is presented in chapter 2 of this thesis:

* a general framework for the principles of VNS.
* the principles of VNS.
* three earlier heuristic approaches.
* different VNS algorithms.
* four parallel implementations of VNS.

The pseudo code, strengths and limitations of each VNS heuristic, and some of its successful applications areas are also discussed in section 2. In Chapter 3 adopts a VNS algorithm is adopted to pack unit spheres into the smallest threedimensional sphere, and the most prominent literature on the subject is reviewed. The proposed hybrid approach proposed, and its relative and absolute performance
is detailed. The superiority of the proposed approach in terms of numerical precision is demonstrated, provide new upper bounds for 29 instances, and showing the utility of the local and variable neighborhood search. The effects of varying the VNS parameters is also presented. In Chapter 4 the problem of packing unit spheres into a cube is discussed, along with a detailed up-to-date literature review on the problem. A description of the solution, highlighting its symmetry reduction is presented along with the formulation augmenting techniques. The results presented compared are then to existing upper bounds. The results obtained matches 16 existing upper bounds and are accurate to $10^{-7}$ for the others. Finally, in Chapter 5 the thesis, is summarized future directions of research are recommended, and other applications of VNS' based hybrid heuristics are proposed.

## Chapter 2

## Variable Neighborhood Search

### 2.1 Introduction

The variable neighborhood search (VNS) is a meta-heuristic or a framework for building heuristics. The VNS has been widely applied during the past two decades because of its simplicity. Its essential concept consists in altering neighborhoods in the quest for a (near-) global optimal solution. In the VNS is investigated the search space are researched via a descent search technique, in which immediate neighborhoods are searched; then, deliberately or at irregular intervals, a more progressive search is adopted where in neighborhoods that are inaccessible from its current point are investigated. Regardless of the type of search, one or a few focal points within the current neighborhood serve as starting points for the neighborhood search. Thus, the search bounces from its current local solution to a new one, the search and only if it discovers a preferred solution, or undertakes a predefined number of successive searches without improvement. Hence, the VNS is not a trajectory emulating technique (as Simulated Annealing or Tabu Search) and does not define
prohibited moves as in Tabu Search.
An optimization problem can be defined as identifying the best value of a real-valued function $f$ over a feasible domain set $X$. The solution $x$ where $x \in X$ is feasible if it satisfies all the constraints for some particular problem. The feasible solution $x^{*}$ is optimal (or is a global optimum) if it yields the best value of the objective function $f$ among all $x \in X$. For instance, when the objective is to minimize $f$, the following holds:

$$
\begin{equation*}
f\left(x^{*}\right)=\min \{f(x): x \in X\} \tag{2.1}
\end{equation*}
$$

i.e., $x^{*} \in X$ and $f\left(x^{*}\right) \leq f(x), \forall x \in X$. The problem is combinatorial if the solution space $X$ is (partially) discrete. A neighborhood structure $N(x) \subseteq X$ of a solution $x \in X$ is a predefined subset of $X$. The solution $x^{\prime} \in N(x)$ is a neighbor of $x$. It is a local optimum of equation (2.1) with respect to (w.r.t.) the neighborhood structure $N(x)$ if $f\left(x^{\prime}\right) \leq f(x), \forall x \in N(x)$. Accordingly, any steepest descent method (i.e., technique that just moves to a best neighbor from the current solution) is trapped when it reaches a local minimum.

To escape from this local optimum, meta-heuristics, or general frameworks for constructing heuristics, adopt some jumping procedures which consist of changing the focal point (or incumbent solution) of the search, or accepting deteriorating moves, or accepting prohibiting moves, etc. The most widely known among these techniques are Genetic Algorithms (GA), Simulated Annealing (SA) and Tabu Search (TS) as detailed in Reeves [58] and Glover and Kochenberger [19].

A brief overview of the VNS and its different variants is presented in this chapter. The most attention will be paid to parallel VNS techniques. In section 2.2 background information is provided on generic VNS and its principles. Two
approaches that are predecessors of VNS are also presented. These approaches are the variable metric approach and iterated local search. In section 2.3 the different versions of VNS are detailed, along with the pseudo code, some applications, and evidence of their strengths for each version. In section 2.4 the parallelization of VNS four approaches are presented. Finally, In section 2.5 a summary is presented.

### 2.2 Preliminaries

The VNS adopts the strategy of escaping from local minima in a systematic way. It deliberately updates the incumbent solution in search of better local optima and in escaping from valleys $[44,45,25,27,28]$. VNS applies a sophisticated system to reach a local minimum; then investigates a sequence of diverse predefined neighborhoods to increase the chances that the local minimum is the global one. Specifically, it fully exploits the present neighborhood by applying a steepest local descent starting from different local minima. When its intensification stops at a local minimum, VNS jumps from the revamped local minimum into a different neighborhood whose structure is different from that of the present one. In fact, it diversifies its search in a pre-planned manner unlike in SA and TS, which permit non-improving moves within the same neighborhood or temper with the solution path. This systematic steepest descent within different neighborhood structures led to the VNS frequently outperforming other meta-heuristics while providing pertinent knowledge about the problem behavior and characteristics; thus, permitting the user to improve the design of the VNS heuristic both in terms of solution quality and runtime while preserving the simplicity of the implementation.

The simplicity and success of the VNS can be attributed to the following three observations:

Observation 1 A local minimum with respect to one neighborhood structure is not necessarily so for another neighborhood.

Observation 2 A global minimum is a local minimum with respect to all possible neighborhood structures.

Observation 3 For many problems, local minima with respect to one or several neighborhoods are relatively close to each other.

Differently stated, the VNS stipulates that a local optimum frequently provides some useful information on how to reach the global one. In many instances, the local and global optima share the same values of many variables.

However, it is hard to predict which ones these variables are. Subsequently, the VNS undertakes a composed investigation of the neighborhoods of this local optimum until an improving one is discovered.

The section 2.2 presents two fundamental variable neighborhood approaches proposed in a different context than the VNS. In section 2.2 .1 the variable metric procedure, originally intended for unconstrained continuous optimization problems, where using different metrics is synonym to different neighborhoods is discussed. In section 2.2.2 the iterated local search, also known as fixed neighborhood search, intended for discrete optimization problems is detailed.

### 2.2.1 The variable metric procedure

This procedure emanates from gradient-based approaches in unconstrained optimization problems. These approaches consist of taking the largest step in the best direction of descent of the objective function at the current point. Initially, the search space is an $n$-dimensional sphere. Adopting a variable metric modifies the
search space into an ellipsoid. The variable metric consists of a linear transformation. This procedure was applied to approximate the inverse of the Hessian of a positive definite matrix within $n$ iterations.

### 2.2.2 Iterated local search (LS)

The most basic form of VNS is the Fixed Neighborhood Search (FNS) [34], also known as the Iterated local search (ILS) [23]. In this strategy an initial solution $x \in X$ is chosen, and subjected to a local search Local $-\operatorname{Search}(x)$, and $x^{*}$ is declared as the best current solution. Then, iteratively undertaken the following three steps are: First, the current best solution $x^{*}$ is perturbed obtaining a solution $x^{\prime} \in X$ by using the procedure $\underline{\text { Shake }}\left(x^{*}\right)$. Second, the procedure Local - Search $\left(x^{\prime}\right)$ is applied to the perturbed solution $x^{\prime}$ to obtain the local minimum $x^{\prime \prime}$. Third and last, the current best solution $x^{*}$ is updated if $x^{\prime \prime}$ is a better local minimum. This three-step iterative approach is repeated until a stopping criterion is met. The stopping condition can be set as the maximal runtime of the algorithm or the maximal number of iterations or an optimality gap of the objective function if a good bound is available. However, this latter condition is rarely applied while the number of maximal iterations without improvement of the current solution is the most widespread criterion. Table 2.1 gives the pseudo code of FNS.

The perturbation of $x^{*}$, via procedure $\underline{\operatorname{Shake}}\left(x^{*}\right)$, is not neighborhood dependent; thus the adoption of the term "fixed". The size of the perturbation should not be too small or too large [23]. In the former case, the search will stagnate at $x^{*}$ since the perturbed solutions $x^{\prime} \in X$ will be very similar to the starting point $x^{*}$ (i.e., within the immediate neighborhood of $x^{*}$ ). In the latter case, the search will be assimilated to a random restart of the local search thus hindering the biased

Table 2.1: Pseudo Code of the FNS

1 Find an initial solution $x \in X$
$2 \quad x^{*} \leftarrow$ Local - Search $(x)$
3 Do While (Stopping condition is False):
$3.1 \quad x^{\prime} \leftarrow \underline{\operatorname{Shake}}\left(x^{*}\right)$
$3.2 \quad x^{\prime \prime} \leftarrow$ Local $-\operatorname{Search}\left(x^{\prime}\right)$
3.3 If $f\left(x^{\prime \prime}\right)<f\left(x^{*}\right)$ set $x^{*}=x^{\prime \prime}$
sampling of FNS where the sampled $x^{\prime}$ solutions are obtained from a fixed neighborhood whose focal point is $x^{*}$. The perturbation is to allow jumps from valleys and should not be easily undone. This is the case of the 4 -opt (known also as the double-bridge) for the traveling salesman problem where the local search initiated from $x^{\prime}$ yield good quality local minima even when $x^{*}$ is of very good quality [23]. The more different the sampled solution $x^{\prime}$ is from the starting point $x^{*}$, the stronger (and most likely the more effective) the perturbation is, but in such a case the cost of the local search is [23] more.

The procedure $\underline{\operatorname{Shake}}\left(x^{*}\right)$ may take into account the historical behavior of the process by prohibiting certain moves for a number of iterations (as in Tabu Search) or by taking into account some precedence constraints, or by fixing the values of certain variables, or by restricting the perturbation to only a subset of the variables.

The procedure Local $-\operatorname{Search}\left(x^{\prime}\right)$ is not necessarily a steepest descent type search. It is any optimization approach that does not reverse the perturbation of the Shake procedure while being reasonably fast and yielding good quality solutions, as is the case in Tabu Search, simulated annealing, a two-opt exhaustive search, etc.

In general, it is advantageous to have a local search that yields very good quality solutions.

Finally, the update of the biased sampling focal point (step 3.3 in Table 2.1) is not necessarily of a steepest-descent type. It may be stochastic; for example, the focal point of the search moves to a non-improving local minimum in search of a global optimum. This strategy resembles the acceptance criterion of simulated annealing. It may also be a combination of a steepest descent and of a stochastic search, which tolerates small deteriorating moves with a certain probability. The choice of the best updating approach should weigh the need and usefulness of the diversification versus intensification of the search [23].

Chiarandini and Stutzle [14] applied FNS (referring to it as ILS) to the graphcoloring problem, using the solution obtained by a well-known algorithm as the initial solution, Tabu Search as the local search, and two neighborhood structures with the first being a one-opt and the second, taking into consideration all pairs of conflicting vertices. They adopt a special array structure to store the effect of their moves and speed the search. Their ILS improves the results of many benchmark instances from the literature. Grosso et al.[9] employ FNS to identify max-min Latin hypercube designs. Their problem consists in scattering $n$ points in a $k$-dimensional discrete space such that the positions of the $n$ points are distinct. The objective is to maximize the distance between every pair of points. For their ILS, they generate the initial solution randomly but suggest that using a constructive heuristic may yield better results. They apply a local search that swaps the first component of two critical points, and a perturbation mechanism that extends the local search to larger neighborhoods. Their ILS improves some existing designs. It is as good as a multi-start search but faster. It is better than the latest implementation of simulated annealing to the problem and competes well with the periodic design approach.

Hurtgen and Maun [33] successfully applied FNS in positioning synchronized phasor-measuring devices in a power system network. They identified the bestknown solution for benchmark instances. When minimizing the total cost (expressed in terms of the number of phasor-mearument units to be installed) while ensuring full coverage of the network, they improved the best-known feasible solution by as much as $20 \%$. Burke et al. [18] compare the performance of ILS to that of six variants of hyper-heuristics a cross three problem domains. A variant combines one of the two heuristic selection mechanisms (uniformly random selection of a heuristic from a prefixed set of heuristics and reinforcement learning with Tabu Search) and one of three acceptance criteria (naive acceptance, adaptive acceptance, and great deluge). Even though it was not consistently the best approach for all tested instances of one-dimensional bin packing, permutation flow shop, and personnel scheduling problems, ILS outperforms, overall, the six variants of the hyper-heuristics. The authors stipulate that the robustness of ILS is due to its successful balancing of its intensification and diversification strategies, to its simple implementation which requires no parameter, and to the proximity of the local minima for the three classes of problems.

### 2.3 Elementary VNS algorithms

Because of its simplicity and successful implementations, VNS has gained wide applicability. This has resulted in several variants of VNS. A successful design takes into account the application area, the problem type, and nature of the variables, not to mention the runtime. Sections 2.3.1 to 2.3.6 present some mostly used VNS versions: the variable neighborhood descent (VND), the reduced variable neighborhood search (RVNS), the basic variable neighborhood search (BVNS), the general
variable neighborhood search (GVNS), the skewed variable neighborhood search (SVNS) and the variable neighborhood decomposition search (VNDS). In section 2.3.7 the distinctive characteristics of each of these variants are highlighted.

### 2.3.1 The variable neighborhood descent

VND is an objective form of VNS. It is founded on the first and third observations, which stipulate that a local optimum for a given neighborhood $N_{k}(x), k \in K=$ $\{1 \ldots, \bar{k}\}$ is not necessarily a local optimum for a second neighborhood $N_{k}(x), k^{\prime} \in$ $K$ and $k^{\prime} \neq k$ and that a global optimum is optimal over all $\bar{k}$ neighborhoods. The parameter $\bar{k}$ is the maximum number of neighborhoods. Therefore, it investigates a neighborhood $N_{k}(x), k \in K$ to obtain a local optimum and searches within other neighborhoods for an improving one. It returns to the initial neighborhood structure every time it identifies an improving solution while it moves to a more distant neighborhood every time it fails to improve the current solution. It stops when the current solution is optimum for all $\bar{k}$ neighborhoods. Table 2.2 gives a detailed algorithm of VND.

The VND is particularly useful for large-sized instances of combinatorial optimization problems where the application of local search meta-heuristics tend to use large CPU times (central processing unit). Nikolic et al. [48] implemented a VND for the covering design problem where neighboring solutions are built by adding and removing subsets to the current solution and resulted in the improving of 13 upper bounds.

Hansen et al.[50] consider berth allocation problem, i.e., the order in which ships will be allocated to berths. They minimize the total berthing costs where the total cost includes the costs of waiting and handling of ships as well as earliness

Table 2.2: Pseudo Code of the VND

1 Find an initial solution $x \in X$, set the best solution $x^{*}=x$, and $k=1$.
2 Do While $(k \leq \bar{k})$ :
2.1 Find the best neighbor $x^{\prime} \in N_{k}(x)$ of $x$
2.2 If $f\left(x^{\prime}\right)<f(x)$ then
set $x=x^{\prime}$ and $k=1$;
If $f\left(x^{\prime}\right)<f\left(x^{*}\right)$, set $x^{*}=x^{\prime}$
else
set $k=k+1$.
or tardiness. They propose a VNS, and compare it to a multi-start heuristic, a genetic algorithm, and a memetic algorithm. They show that the VNS is on average the better of the three approaches. Qu et al. [57] apply the VND to the delayconstrained least-cost multicast routing problem, which reduces to the NP-complete delay-constrained Steiner Tree. They investigate the effect of the initial solution by considering two initial solutions: one based on the shortest path and one on bounded-delay, and consider three types of neighborhoods. They show that the VND outperforms the best existing algorithms. Kratica et al. [35] use a VND within large shaking neighborhoods to solve a balanced location problem. This latter problem consists of choosing the location of $p$ facilities while balancing the number of clients assigned to each location. They show that their implementation is very fast and reaches the optimal solution for small instances. For large instances where the optimal solution value is unknown, the VNS outperforms a state-of-the-art modified genetic algorithm.

### 2.3.2 The reduced variable neighborhood descent

The RVNS is a slightly different provision of the VNS; yet, it retains its basic structure. It is based upon the third principle of the VNS which states that a global optimum is the best solution a cross all possible neighborhoods. It is a stochastic search. The outer loop constitutes the stopping. The inner loop conducts a search over a fixed set of $\bar{k}$ nested neighborhoods $N_{1}(x) \subset N_{2}(x) \subset \ldots, \subset N_{\bar{k}}(x)$ that are centered on a randomly generated focal point $x$. A solution $x^{\prime} \in X$ is generated using a stochastic procedure $\underline{\text { Shake }}(x, k)$ which slightly perturbs the solution $x$. For instance, in non-linear programming where the final solution of any optimization solver depends on the solution it is fed, procedure Shake $(x, k)$ would alter the value of one or more of the variables of $x$ by small amounts $\delta$ where the altered variable(s) and $\delta$ are randomly chosen. Subsequently, either the focal point of the search is changed to the current solution $x$ or an enlarged neighborhood search is to be undertaken. Specifically, when $f\left(x^{\prime}\right)<f(x)$, the inner loop centers its future search space on the most reduced neighborhood around $x^{\prime}$ (i.e., it sets $x=x^{\prime}$ and $k=1$ ). Otherwise, it enlarges its sampling space search by incrementing $k$ but maintains its current focal point $x$. Every iteration of the inner loop can be assimilated to injecting a cut to the minimization problem, if each new bound improves the existing one; thus, to fathoming a part of the search space.

Table 2.3 gives a detailed pseudo code of the RVNS. This approach is well suited for multi-start VNS-based strategies where RVNS can be replicated with different initial solutions $x \in X$ and the best solution over all replications is retained. It behaves like a Monte Carlo simulation except that its choices are systematic rather than random. It is a more general case of Billiard simulation; a technique that was applied to point scattering within a square.

Table 2.3: Pseudo Code of the RVNS

1 Find an initial solution $x \in X$; set the best solution $x^{*}=x$, and $k=1$.
2 Choose a stopping condition.
3 Do While (Stopping condition is False):
Do While ( $k \leq \bar{k}$ )
$3.1 x^{\prime} \leftarrow \underline{\text { Shake }}(x, k)$
3.2 If $f\left(x^{\prime}\right)<f(x)$, then
set $x=x^{\prime}$ and $k=1$;
If $f\left(x^{\prime}\right)<f\left(x^{*}\right)$ set $x^{*}=x^{\prime}$
else
set $k=k+1$.

RVNS is particularly useful when the instances are large or when the local search is expensive [53, 47]. Computational proof is given in section 2.3.4 for the high school timetabling problem, where the performance of the RVNS is been to better than be the GVNS, which applies simulated annealing as the local search [59]. Hansen et al. [53] further substantiated the claim regarding the speed of the RVNS. They compare the performance of the RVNS to that of the fast interchange heuristic for the $p$-median problem. They show that the speed ratio can be as large as 20 for comparable average solutions. For the same problem, Crainic et al. [72] provide additional performance analysis of a parallel implementation of the RVNS. Maric et al. [43] hybridize the RVNS with a standard VNS. They compare the performance of the resulting heuristic to a swarm optimization algorithm and a simulated annealing one for the bi-level facility location problem with clients' preference, but unlimited capacity for each facility. They show that the hybrid heuristic outperforms the other two in large-scale instances and is competitive in the case of smaller instances.

### 2.3.3 The basic variable neighborhood search

The BVNS is a descent, first-improvement method [53]. It is a hybridization of the VND and the RVNS. It evokes variable neighborhoods at irregular intervals but consistently applies a steepest descent to a (near-) global optimum. It is basically an RVNS where the inner loop applies to the solution $x^{\prime}$ obtained via $\underline{\operatorname{Shake}}(x, k)$ a steepest descent procedure Local $-\operatorname{Search}\left(x^{\prime}, k\right)$, which searches around the solution $x^{\prime}$ for the best first improving solution $x^{\prime \prime} \in N_{k}\left(x^{\prime}\right)$, where $k \in K$. In fact, it accelerates the search by opting for the first rather than the best improving solution. The BVNS adopts the same stopping criteria as the RVNS; that is, the inner loop stops if the investigation of $\bar{k}$ successive neighborhoods centered on $x$ yield no improvement, whereas the outer loop stops when a user-defined stopping criterion is satisfied. This condition is generally related to total runtime or the number of iterations without improvement of the best solution. Table 2.4 provides a detailed description of the BVNS.

The BVNS is well suited to multi-start VNS-based strategies, where a local search is applied to perturbed solutions. Toksari and Guner [73] provide computational proof that the local search of the BVNS is more efficient than the VND in the case of unconstrained optimization problems. They consider the case of a non-convex but differentiable function with many local minima but a single global minimum. Their VNS applies a standard descent heuristic with the directions of descent randomly generated. Their results are competitive with existing approaches when tested on existing benchmark instances. Sanchez-Oro and Duarte [62] show that the BVNS is superior to the RVNS and the VND where finding near-optima for both the min-max and min-sum variants of the vertex-cut minimization problems for short and long time horizons. M'Hallah and Alkandari [55] and M'Hallah et

Table 2.4: Pseudo Code of the BVNS

1 Find an initial solution $x \in X$; set the best solution $x^{*}=x$, and $k=1$.
2 Choose a stopping condition.
3 Do While (Stopping condition is False):
Do While ( $k \leq \bar{k}$ )
$3.1 x^{\prime} \leftarrow \underline{\operatorname{Shake}}(x, k)$.
$3.2 x^{\prime \prime} \leftarrow$ Local $-\operatorname{Search}\left(x^{\prime}, k\right)$.
3.3 If $f\left(x^{\prime \prime}\right)<f(x)$, then
set $x=x^{\prime \prime}$ and $k=1$;
If $f\left(x^{\prime \prime}\right)<f\left(x^{*}\right)$ set $x^{*}=x^{\prime \prime}$
else
set $k=k+1$.
al. [56] apply BVNS for packing unit spheres into the smallest cube and sphere, respectively, highlighting the utility of the local search.

### 2.3.4 The general variable neighborhood search

The GVNS is a low-level hybridization of the BVNS with the RVNS and the VND [51, 52]). A detailed description of the GVNS is given in Table 2.5. First, it applies the RVNS to obtain a feasible initial solution to the problem (step 2 in Table 2.5) in lieu of directly sampling a feasible solution $x \in X$ (as in step 1 in Table 2.4). This is particularly useful in instances where finding a feasible solution is in itself an NP-hard problem. Second, it replaces the local search (Step 3.2 in Table 2.4) of the BVNS with a VND (Step 5.2 in Table 2.5). In addition, it samples the points $x^{\prime}$ from the $k^{\text {th }}$ neighborhood $N_{k}(x)$ of the current solution $x$. In fact, it uses procedure

Shake $(x, k)$.

Table 2.5: Pseudo Code of the GVNS

1 Find an initial solution $x$.
$2 x^{\prime} \leftarrow R V N S(x)$ starting with $x$ to obtain a feasible solution $x^{\prime} \in X$.
3 Set the best solution $x^{*}=x^{\prime}$ the current solution $x=x^{\prime}$ and the neighborhood counter $k=1$.

4 Choose a stopping condition.
5 Do While (Stopping condition is False)
Do While $(k \leq \bar{k})$ :
$5.1 x^{\prime} \leftarrow \underline{\text { Shake }}(x, k)$.
$5.2 \quad x^{\prime \prime} \leftarrow V N D\left(x^{\prime}\right)$.
5.3 If $f\left(x^{\prime \prime}\right)<f(x)$, then
set $x=x^{\prime \prime}$ and $k=1$;
If $f\left(x^{\prime \prime}\right)<f\left(x^{*}\right)$ set $x^{*}=x^{\prime \prime}$
else
set $k=k+1$.

GVNS was applied to large-sized vehicle routing problems with time windows $[64,46]$ and to the traveling salesman problem [46]. The implementation of Mladenovic et al. [46] provides the best upper bounds in more than half of the existing benchmark instances. It was particularly useful because identifying initial feasible solutions and maintaining feasibility during the shaking procedure and the neighborhood investigation was a challenging task.

### 2.3.5 The skewed variable neighborhood search

The SVNS is a modified BVNS where a non-improving solution $x^{\prime \prime}$ may become the new focal point of the search, when this solution $x^{\prime \prime}$ is far enough from the current one $x$, but its value $f\left(x^{\prime \prime}\right)$ is not much worse than $f(x)$, the value of the current solution $x$ [27]. The SVNS is motivated by the topology of search spaces. The search gets trapped in a local optimum and cannot leave it because all neighboring solutions are worse. Yet, if it opts for a "not-too-close" neighborhood, it may reach the global optimum. This neighborhood should not be too far and the neighbor should not be too much worse can to enable to return to the current neighborhood, if the exploration fails to identify an improving solution.

Table 2.6: Pseudo Code of the SVNS

1 Find an initial solution $x \in X$; set the best solution $x^{*}=x$, and $k=1$.
2 Choose a stopping condition.
3 Do While (Stopping condition is False)
Do While ( $k \leq \bar{k}$ ):
$3.1 \quad x^{\prime} \leftarrow \underline{\text { Shake }}(x, k)$.
$3.2 x^{\prime \prime} \leftarrow$ Local $-\operatorname{Search}\left(x^{\prime}, k\right)$.
3.3 If $f\left(x^{\prime \prime}\right)-\alpha \rho\left(x, x^{\prime \prime}\right)<f(x)$, then set $x=x^{\prime \prime}$ and $k=1$;
If $f\left(x^{\prime \prime}\right)<f\left(x^{*}\right)$ set $x^{*}=x^{\prime \prime}$
else
set $k=k+1$.

The difference between $x$ and $x^{\prime \prime}$, is measured in terms of a distance $\rho\left(x, x^{\prime \prime}\right)$ while the difference between $f\left(x^{\prime \prime}\right)$ and $f(x)$ is considered tolerable if it is less than an
expression pondering the distance $\rho\left(x, x^{\prime \prime}\right)$ by a parameter $\alpha$. That is, the condition $f\left(x^{\prime \prime}\right)<f(x)$ of step 3.3 in Table 2.4 is replaced by $f\left(x^{\prime \prime}\right)-\alpha \rho\left(x, x^{\prime \prime}\right)<f(x)$. A detailed description of the BVNS is provided in Table 2.6.

The utility of the SVNS is demonstrated by Hansen and Mladenovic [27] for the weighted maximum satisfiability of logic problem for which the SVNS performs better than the GRASP (Greedy Randomized Adaptive Search Procedure) and Tabu Search, for medium and large problems, respectively. The choice of $\alpha$ is generally based on an analysis of the behavior of a multiple start VNS.

### 2.3.6 The variable neighborhood decomposition search

The VNDS is used in the particular case, where the set $X$ of feasible solutions is finite. Yet, its extension to the infinite case is possible. It is one of two techniques designed to reduce the computational time of local search algorithms. Even though they investigate a single neighborhood structure, local search heuristics tend to have their runtime increase significantly as the size of the combinatorial problems become large [53]. It is "a BVNS within successive approximations decomposition" scheme [53]. That is, both the VNDS and the BVNS have the same algorithmic steps except that procedures Shake and Local - Search of steps 3.1 and 3.2 of Table 2.4, respectively, are implemented on a partial solution $y$ of free variables, whereas all other variables remain fixed as in $x$ throughout the random selection of $x^{\prime} \in X$ and the local search for an improving solution $x^{\prime \prime} \in X$. A detailed description of the pseudo-code of the VNDS is given in Table 2.7.

Obtaining a random solution $x^{\prime} \in X$ from the current solution $x \in X$ via procedure $\underline{\operatorname{Shake}}(x, k, y)$ entails choosing values for the partial solution $y$, which consists of a set of free variables of $x$, while ensuring that the new value of every

Table 2.7: Pseudo Code of the VNDS

1 Find an initial solution $x \in X$; set the best solution $x^{*}=x$, and $k=1$.
2 Choose a stopping condition.
3 Do While (Stopping condition is False)
Do While $(k \leq \bar{k})$ :
$3.1 \quad x^{\prime} \leftarrow \underline{\text { Shake }}(x, k, y)$.
$3.2 \quad x^{\prime \prime} \leftarrow$ Local - Search $\left(x^{\prime}, y, k\right)$.
3.3 If $f\left(x^{\prime \prime}\right)<f(x)$, then
set $x=x^{\prime \prime}$ and $k=1$;
If $f\left(x^{\prime \prime}\right)<f\left(x^{*}\right)$ set $x^{*}=x^{\prime \prime}$
else
set $k=k+1$.
free variable is different from its current value in $x$. The choice of the free and fixed variables constitutes the heart of the decomposition approach. It can follow some rule of thumb or some logical pattern. Note that the cardinality of $y$ set is equal to $n-k$.

Obtaining a local optimum $x^{\prime \prime} \in X$ from the current solution $x^{\prime} \in X$ via procedure Local $-\operatorname{Search}\left(x^{\prime}, y, k\right)$ entails finding the best values of the partial solution $y$, given that the all other fixed variables of $x^{\prime}$ retain their values in the local optimum $x^{\prime \prime}$. It is equivalent to undertaking a local search on the reduced search space of $y$. It is possible not to apply a local search and to simply implement an inspection approach or exactly solve the reduced problem if the number of fixed variables is very large.

The VNDS is useful in binary integer programming in general [36], and solving
large-scale $p$-median problems [53] in particular. The $p$-median problem involves choosing the location of $p$ facilities among $m$ potential ones in order to satisfy the demand of a set of users at least cost. Lazic et al. [36] provide computational proof that the VNDS performs best on all performance measures of solution approaches including computation time, optimality gap, and solution quality. Hanafi et al. [60] tackle the 0-1 mixed-integer programming problem using a special VNDS variant. This latter exploits the information obtained from an iterative relaxation-based heuristic in its search. This information serves to reduce the search space and avoid the reassessment of the same solution during different replications of the VNDS. The heuristic adds pseudo-cuts based on the objective function value to the original problem to improve the lower and upper bounds; thus, it reducing the optimality gaps. The approach yields the best average optimality gap and running time for binary multi-dimensional knapsack benchmark instances. It is inferred that the approach can yield tight lower bounds for large instances.

### 2.3.7 Comparison of the VNS variants

The VNS was designed for combinatorial problems, but is applicable to any global optimization problem. It explores distant neighborhoods of incumbent solutions in search for global optima. It is simple and requires very few parameters. The main characteristics of its seven variants are summarized in Table 2.8. The VNS has been hybridized at different levels with other heuristics and meta-heuristics. For instance, Kandavanam et al. [20] consider route multi-class network communication planning problem in order to satisfy service quality. They hybridize the VNS with a genetic algorithm and apply the hybrid heuristic to maximizing the residual bandwidth of all links in the network, while meeting the requirements of the expected quality of service.

Table 2.8: Main Characteristics of the VNS Variants

VND : deterministic change of neighborhoods; more likely to reach a global minimum, and if many neighborhood structures are used.

RVNS : useful for very large instances, for which local search is costly; best with $\bar{k}=2$, and analogous to a Monte-Carlo simulation, but more systematic.

BVNS : deterministic and stochastic changes of neighborhoods, and systematic change of neighborhoods.

GVNS : VND is used as a local search within the BVNS; very effective, and useful for low-level hybridization.

SVNS : useful for flat problems, and useful for clustered local optima.
VNDS : a two-level VNS (decomposition at the first level), and useful for large instances.

### 2.4 Parallel VNS

Most sequential heuristic approaches are being implemented as parallel approaches. The increasing tendency towards parallelism is motivated both by the potential reduction of computational time (through the segmentation of the sequential program, and by the expansion of the investigation of the search space (through the provision of more processors and memory for the computing device). The VNS is among the sequential heuristics that were implemented in a parallel computing environment. Four parallelization techniques have been so far proposed [29]. The first two techniques are basic: the parallelization of the neighborhood local search and of the VNS itself by assigning the same VNS algorithm to each thread, and retaining the best solution among all solutions reported by the threads. There is no cooperation among the individual threads. The remaining two techniques on the other hand
utilize cooperation mechanisms to upgrade the performance level of the algorithm. They are more complex and involve intricate parallelization [39, 72] to synchronize communication. A detailed description of these four techniques follows.

The first parallelization technique is the synchronous parallel VNS (SPVNS). It is the most primary parallelization technique [39]. It is designed to shorten the

Table 2.9: Pseudo Code of the SPVNS
1 Find an initial solution $x \in X$; set the best solution $x^{*}=x$, and $k=1$.
2 Choose a stopping condition.
3 Do While (Stopping condition is False) Do While ( $k \leq \bar{k}$ ):
$3.1 \quad x^{\prime} \leftarrow$ Shake $(x)$.
3.2 Divide the neighborhood $N_{k}\left(x^{\prime}\right)$ into $n_{p}$ subsets.
3.3 For every processor $p, p=1 \ldots, n_{p}, x_{p}^{\prime \prime} \leftarrow \operatorname{Local}-\operatorname{Search}\left(x^{\prime}, k\right)$.
3.4 Set $x^{\prime \prime}$ such that $f\left(x^{\prime \prime}\right)=\max _{p=1, n_{p}}\left\{f\left(x_{p}^{\prime \prime}\right)\right\}$
3.5 If $f\left(x^{\prime \prime}\right)<f(x)$, then
set $x=x^{\prime \prime}$ and $k=1$;
If $f\left(x^{\prime \prime}\right)<f\left(x^{*}\right)$ set $x^{*}=x^{\prime \prime}$
else
set $k=k+1$.
runtime through the parallelization of the local search of the sequential VNS. In fact, the local search is the most time-demanding part of the algorithm. The SPVNS splits the neighborhood into $n_{p}$ parts and assigns each subset of the neighborhood to an independent processor, which returns to the master processor an improving neighbor within its subset of the search space. The master processor sets the best among the $n_{p}$ neighbors returned by the $n_{p}$ processors as the current solution. Table
2.9 provides the pseudo code of the SPVNS, adapted from Garcia et al. [39].

The second parallelization technique is the reproduced parallel VNS (RVNS) or replicated parallel VNS or simply a multi-start VNS. It consists of $n_{p}$ parallel independent searches, where $n_{p}$ is the number of parallel threads of the computing device. Each independent search operates an independent VNS on a separate processor. Table 2.10 provides a detailed description of the RPVNS. It can be perceived as a multi-start RVNS where the best solution $n_{p}$ is retained as the best solution except that, the $n_{p}$ replications are undertaken in parallel instead of sequentially [39].

Table 2.10: Pseudo Code of the RPVNS

1 Choose a stopping condition.
2 Do While ( $p \leq n_{p}$ )
2.1 Find an initial solution $x_{p} \in X$; set the best solution $x_{p}^{*}=x_{p}$, and $k=1$.
2.2 Do While (Stopping condition is False)

Do While $(k \leq \bar{k})$

- $x^{\prime} \leftarrow \underline{\text { Shake }}\left(x_{p}\right)$.
- $x^{\prime \prime} \leftarrow$ Local $-\operatorname{Search}\left(x^{\prime}, k\right)$.
- If $f\left(x^{\prime \prime}\right)<f(x)$, then
set $x=x^{\prime \prime}$ and $k=1$;
If $f\left(x^{\prime \prime}\right)<f\left(x_{p}^{*}\right)$ set $x_{p}^{*}=x^{\prime \prime}$;
else
set $k=k+1$.
$3-$ Set $x^{*}$ such that $f\left(x^{*}\right)=\max _{p=1, n_{p}}\left\{f\left(x_{p}^{*}\right)\right\}$.

The third parallelization technique is the replicated shaking VNS (RSVNS) proposed by Garcia et al. [39]. It applies a synchronous cooperation mechanism
through a conventional master-slave methodology. The master processor operates a sequential VNS, and sends its best incumbent to every slave processor, which shakes this solution to obtain a starting point for its own local search. In turn, each slave returns its best solution to the master. This latter compares all the solutions obtained by the slaves, retains the best one and subjects it to its sequential VNS. This information exchange is repeated until a stopping criterion is satisfied. The

Table 2.11: Pseudo Code of the RSVNS

1 Find an initial solution $x \in X$; set the best solution $x^{*}=x$, and $k=1$.
2 Choose a stopping condition.
3 Do While (Stopping condition is False)
Do While ( $k \leq \bar{k}$ )
3.1 For $p=1 \ldots, n_{p}$

- Set $x_{p}=x$
- $x^{\prime} \leftarrow \underline{\text { Shake }}\left(x_{p}^{\prime}\right)$.
- $x_{p}^{\prime \prime} \leftarrow$ Local - Search $\left(x_{p}^{\prime}, k\right)$.
3.2 Set $x^{\prime \prime}$ such that $f\left(x^{\prime \prime}\right)=\max _{p=1, n_{p}}\left\{f\left(x_{p}^{\prime \prime}\right)\right\}$
3.3 If $f\left(x^{\prime \prime}\right)<f(x)$, then
set $x=x^{\prime \prime}$ and $k=1$;
If $f\left(x^{\prime \prime}\right)<f\left(x_{p}^{*}\right)$ set $x_{p}^{*}=x^{\prime \prime}$;
else
set $k=k+1$.
fact that VNS permits changing neighborhoods and types of local search makes this type of parallelization possible. Different neighborhoods or local searches can be performed by independent processors and the resulting information is channeled to the master processor which analyzes its results to obtain the best solution and
conveys it to the other slaves. Maintaining a trace of the last step undertaken by every processor strengthens the search as it avoids the duplication of computational efforts. Table 2.11 provides a detailed description of the RSVNS. It can be perceived as a VNS with a multi-start shaking and local search where the best local solution among $n_{p}$ ones is retained as the best local solution and routed to each of the $n_{p}$ processors [39]. A comparison of the performance of these first three parallel VNS algorithms is undertaken by Garcia et al. [39] for the $p$-median problem.

The fourth and last parallelization technique is the cooperative neighborhood VNS (CNVNS) suggested by Crainic et al. [72] and Moreno-Perez et al. [38]. It is particularly suited to combinatorial problems such as the $p$-median problem. It deploys a cooperative multi-search strategy to the VNS while exploring a centralmemory mechanism. It coordinates many independent VNSs by asynchronously exchanging the best incumbent obtained so far by all processors. Its implementation preserves the simplicity of the original sequential VNS ideas. Yet, its asynchronous cooperative multi-search offers a broader exploration of the search space thanks to the different VNSs being applied by the different processors. Each processor undertakes an RVNS, and communicates with a central memory or the master, every time it improves the best global solution. In turn, the master relays this information to all other processors so that they update their knowledge about the best current solution. That is, no information is exchanged in terms of the VNS itself. The master is responsible for launching and stopping the parallel VNS. In case of Crainic et al. [72], the parallel algorithm is an RVNS where the local search is omitted; resulting in a faster algorithm. Table 2.12 and Table 2.13 gives a detailed description of the CNVNS as originally intended and as described by Moreno-Perez et al. [38]. A comparison of the performance of the four parallel VNS algorithms is undertaken by Moreno-Perez et al. [38] for the $p$-median problem.

Table 2.12: Pseudo Code of the CNVNS: Master's Algorithm

1 Find an initial solution $x \in X$; set the best solution $x^{*}=x$, and $k_{p}=1$, $p=1, \ldots, n_{p}$

2 Choose a stopping condition.
3 Set $x_{p}=x$
4 For $p=1, \ldots, n_{p}$, launch RVNS with $x_{p}$ as its initial solution.
5 When processor $p^{\prime}$ sends $x_{p^{\prime}}^{\prime \prime}$.
5.1 If $f\left(x_{p^{\prime}}^{\prime \prime}\right)<f(x)$, then update current best solution setting $x=x_{p^{\prime}}^{\prime \prime}$ and send it to all $n_{p}$ processors.

6 When processor $p^{\prime}$ requests best current solution, send $x$

Table 2.13: Pseudo Code of the CNVNS: Slave's Algorithm

1 Obtain initial solution $x_{p}$ from master; set the best solution $x_{p}^{*}=x_{p}$, and randomly choose $k \in\{1, \ldots, \bar{k}\}$

2 Do While $(k \leq \bar{k})$
$2.1 x^{\prime} \leftarrow \underline{\text { Shake }}(x)$.
2.2 If $f\left(x_{p}^{\prime}\right)<f\left(x_{p}\right)$, then
set $x_{p}=x_{p}^{\prime}$ and $k=1$;
if $f\left(x_{p}^{\prime}\right)<f\left(x_{p}^{*}\right)$ set $x_{p}^{*}=x_{p}$.
else
send $x_{p}^{*}$ to the master;
receive $x_{p}$ from master;
set $k=k+1$.

### 2.5 Discussion and conclusion

The VNS is a general framework for solving optimization problems. Its successful application to different continuous and discrete problems is advocating for its wider use to non-traditional areas such as neural networks and artificial intelligence. The VNS owes its success to its simplicity and to its limited number of parameters:the stopping criterion and maximal number of neighborhoods. Depending on the specific problem at hand, a VNS variant may be deemed more appropriate than other variants. In fact, a judicious choice of the neighborhood structure and local search strategy could determine the success of an approach. The local search may vary from exact optimization techniques for relaxed or reduced problems to gradient descent, line search, steepest descent, to meta-heuristics like Tabu Search and simulated annealing.

## Chapter 3

## Packing Unit Spheres into the Smallest Sphere Using the VNS and NLP

### 3.1 Introduction

Sphere packing problems (SPP) consist comprise of packing spheres into a sphere of the smallest possible radius. It has many important real-life applications including materials science, radio surgical treatment, communication, and other vital fields. In materials science, random sphere packing is a model to represent the structure of liquids, proteins, and glassy materials. The model is used to study phenomena such as electrical conductivity, fluid flow, stress distribution and other mechanical properties of granular media, living cells, random media chemistry and physics. SPP also entails the investigation of processes such as sedimentation, compaction, and sintering [76]. In radio surgical treatment planning, sphere packing is crucial to

X-ray tomography. In [75] a problem on how to pack minimum number of unequal spheres into a three dimensions bounded region, in connection with radio-surgical treatment planning, is studied. It is used for treating brain and sinus tumor (see Figure 3.1).


Figure 3.1: X-Ray

During the operation, medical unit should not effect other organs. Gamma knife is one of more effective radio-surgical modalities. It can be described as radioactive dosage treatment planning. A high radiation dose is called the shot which can be viewed as a sphere. These shots are ideal equal spheres to be packed into a container, but also they could be of different spheres. The tumor can be viewed as an approximate spherical container. No shots are allowed outside the container, which means that the radiation shots are only hitting the tumor cells and not the healthy ones. Multiple shots are usually applied at various locations and may touch the boundary of the target. They avoid overlapping and touch each other. The stronger the high packing density, the more doses delivered. The target of the search to minimize the number of shots into the container (The tumor). As a result, this ap-
proach has met with certain success in medical applications. Dynamic programming algorithm is being used to find the optimum number of shots.

In digital communication and storage, it emerges in compact disks, cell phones, and the Internet [15, 74]. The most frequent application of minimum sphere packing problem is connected with location of antennas. For example it is crucial in antenna location in some large warehouse or container yards (see Figure 3.2). Each article usually has radio frequency identifier (RFID) or tag. The management of the warehouse wants to locate minimum number of antennas that cover all warehouse such that vehicles, connected with the antenna system, are able to find any article. The radii of each antenna are known in advance. This system provides real-time location visibility, illuminating vital information that is needed [6].


Figure 3.2: Warehouse

Other applications of sphere packing are encountered in powder metallurgy for three-dimensional laser cutting [42]; cutting different natural crystals; layout of computers, buildings, etc.

In this chapter the special case of packing unit spheres into the smallest sphere (PSS) is considered. PSS entails packing $n$ identical spheres, of radius 1
unit, without overlap into the smallest containing sphere $S$. The goal is to search for the best packing of the $n$ spheres into $S$, where the best packing minimizes waste. According to the Typology of Cutting and Packing of Wascher et al. [79], PSS are a three-dimensional variant of the Open Dimension Problem since all small items (which are spheres) have to be packed and the dimension of the large object (which is a sphere) is not given, and has to be minimized. PSS is equivalent to finding the coordinates $\left(x_{i}, y_{i}, z_{i}\right)$ of every sphere $i, i=1, \ldots, n$, the radius $r$ and coordinates $(x, y, z)$ of $S$, such that no pair of spheres $(i, j) \in I \times I$ and $i<j$ overlap. Formally, the problem can be stated as finding the optimal level of the decision variables $r,(x, y, z)$, and $\left(x_{i}, y_{i}, z_{i}\right), i=1, \ldots, n$, such that

$$
\begin{array}{cc}
\text { min } & r \\
\text { subject to } & \left(x_{i}-x\right)^{2}+\left(y_{i}-y\right)^{2}+\left(z_{i}-z\right)^{2} \leq(r-1)^{2} \\
& i=1, \ldots, n, \\
\left(x_{i}-x_{j}\right)^{2}+\left(y_{i}-y_{j}\right)^{2}+\left(z_{i}-z_{j}\right)^{2} \geq 4 & 1 \leq i<j \leq n, \\
r \geq \underline{r}, \tag{3.4}
\end{array}
$$

where $\underline{r}$ is a strictly positive lower bound on $r$. The first set of constraints enforces the containment of every sphere within $S$. The second set enforces the no overlap constraint of any pair of distinct spheres. Finally, the last constraint provides a positive lower bound for the radius of the containing sphere.

PSS is a non-convex optimization problem [32, 75]. It is NP-hard [75], since it is an extension of packing unit circles into the smallest circle, which is in turn NPhard [32]. Thus, the search for an exact local extremum is time-consuming, without any guarantee of a sufficiently good convergence to the optimum [65]. Indeed, PSS is a challenging problem. As the number of unit spheres increases, identifying a reasonably good solution becomes extremely difficult [32]. In addition, PSS has
an infinite number of solutions with identical minimal radii. In fact, any solution may be rotated or reflected or may have free spheres which can be slightly moved without enlarging the radius of the containing sphere. Finally, there is the issue of computational accuracy and precision.

Solving PSS via an off-the-shelf non-linear programming (NLP) solver is generally not successful. In fact, most of these solvers fail to identify global optima because of PSS' difficulty, which gets further amplified as the problem size increases (since the number of local minima increases too). Subsequently, PSS should be solved by a mixture of search heuristics with local exhaustive (exact) search of local minima or their approximations as suggested in $[67,32]$. This chapter follows this line of research. It tackles PSS using a hybrid heuristic which combines these two main components. It applies a variable neighborhood search (VNS) which investigates different neighborhoods of the incumbent solution, and a special purpose NLP solver to locally search for one or more minima within each neighborhood.

Section 3.2 reviews the most prominent literature in the area of sphere packing. Section 3.3 details the proposed approach. Section 3.4 investigates its performance. Specifically, this section demonstrates the superiority of the proposed approach in terms of accuracy and precision, provides new upper bounds for 14 instances, and determines the utility of VNS and the local search. Finally, section 3.5 is a summary.

### 3.2 Literature review

Sphere packing is considered when the container three-dimensional into which the sphere are to be packed is a cube, a parallelepiped, a cylinder, a polytope, or a sphere.

Gensane [22] adapts the Billiard simulation to the problem of identifying the largest radius of $n$ identical spheres that fit inside a unitary cube. Billiard simulation is a stochastic method that mimics the idealized movement of billiard balls inside a domain, with the initial centers of the balls and their directions being randomly fixed. The resulting configuration emanates from probabilistic choices. To improve the convergence of the stochastic algorithm, different types of local searches were considered. These included moving the spheres randomly in a random walk simulation, decreasing the magnitude of the stochastic movements as the size of the spheres increased, and perturbing all spheres simultaneously.

Stoyan and Yaskov [66] focus on finding the minimal height of a parallelepiped $(L, W, H)$ that packs $n$ identical spheres. They model the problem as a NLP with a linear objective function and linear and quadratic constraints. They approximately solve the model using a special search tree in conjunction with a modification of the Zoutendijk method [80] of feasible directions to calculate local minima, and a modification of the decremental neighborhood method to search for an approximation to the global minimum.

Using the same technique, Stoyan and Yaskov [68] tackle the problem of minimizing the height $H$ of a right circular cylinder, of known radius $r$, that packs $n$ identical spheres. The authors obtain the best results to date for $n=498,499$, and 500. Their approach is very effective for $n \leq 500$, and can handle instances for $n \leq 2000$.

Stoyan et al. [77] use techniques similar to those considered in [67] to identify a packing of $n$ non-identical spheres, each of radius $r_{i}, i \in I$, into a parallelepiped of fixed length $L$ and width $W$ but of variable height $H$ with the objective of minimizing $H$. The authors provide numerical results with up to 60 spheres.

Yaskov et al. [21] tackle the problem of maximizing the number of identical spheres that can be packed into a cylindrical composed domain. They construct a mathematical model based on the concept of $\Phi$-functions [78], and design a solution algorithm based on a modification of the optimization method by groups of variables.

Wang [75] formulates mathematically the automated radio surgical treatment planning problem as the packing of spheres into a three-dimensional region with a packing density greater than a given threshold level. He proves that this packing problem is NP-complete and proposes an approximate algorithm to solve it.

Sutou and Dai [70] assimilate the automated radio surgical treatment planning problem to packing non-identical spheres in a three-dimensional polytope with the objective of maximizing the volume of the packed spheres. They formulate the problem as a non-convex optimization one with quadratic constraints and a linear objective function. They propose a variety of algorithms which outperform the generic algorithm for the general non-convex quadratic program. In fact, they incorporate heuristics into the generic algorithm to strengthen its efficiency. They demonstrate its efficiency computationally for limited problem sizes.

Stoyan and Yaskov [69] consider the problem of packing the maximal number of congruent hyper spheres of unit radius into a larger hyper sphere of given radius. They construct a mathematical model, and approximately resolve it by solving a sequence of packing subproblems -whose objective functions are linear- using the Zoutendijk feasible direction method. Starting points are generated in accordance with the lattice packing of hyperspheres translated on various vectors or in a random way. They improve the convergence of their sequential approach by perturbing the lattice packing.

Hales [24] proves the Kepler conjecture which stipulates that the packing
density of identical three-dimensional spheres cannot exceed $\frac{\pi}{\sqrt{18}}$. Bezdek [11] brings forth a strong Kepler conjecture which alleges that the density of at least two equal spheres in a sphere of the three-dimensional space of constant curvature is less than $\frac{\pi}{\sqrt{18}}$.

Birgin and Sobral [13] propose a reduced model for PSS. They aggregate the $\frac{n(n-1)}{2}$ non-overlap constraints into a single constraint:

$$
\begin{equation*}
\sum_{i=1}^{n-1} \sum_{j=i+1}^{n} \max \left\{0,4-\left(x_{i}-x_{j}\right)^{2}-\left(y_{i}-y_{j}\right)^{2}-\left(z_{i}-z_{j}\right)^{2}\right\}=0 \tag{3.5}
\end{equation*}
$$

and implement a reduction approach that avoids computing many of the $\frac{n(n-1)}{2}$ terms of the sum. In fact, only terms relative to spheres that could potentially be touching are included. They solve the resulting model, starting from several initial solutions, using a local solver that is based on an augmented Lagrangian method for smooth general constrained minimization. They report solutions for instances with up to 100 spheres.

Liu et al. [37] adopt an unconstrained model for the problem where they aggregate all the constraints of PSS and "Lagrange" them into the objective function:

$$
\begin{align*}
& \quad \min  \tag{3.6}\\
& \text { where } \quad r+\sum_{i=1}^{n} d_{i 0}^{2}+\sum_{i=1}^{n-1} \sum_{j=i+1}^{n} d_{i j}^{2},  \tag{3.7}\\
& \text { and } \quad d_{i j}^{2}=\max \left\{0,4-\left(x_{i}-x_{j}\right)^{2}-\left(y_{i}-y_{j}\right)^{2}-\left(z_{i}-z_{j}\right)^{2}\right\},  \tag{3.8}\\
& d_{i 0}^{2}=\max \left\{0,\left(x_{i}-x\right)^{2}+\left(y_{i}-y\right)^{2}+\left(z_{i}-z\right)^{2}-(r-1)^{2}\right\} .
\end{align*}
$$

Their model is inspired from quasi physics where spheres that are in contact have extrusive elastic energy, and a system is in equilibrium when all elastic energies are nil. They solve the resulting model using a heuristic which combines the energy landscape paving (ELP) method with an adaptive step length gradient descent pro-
cedure. ELP is an optimization strategy that mimics the Monte Carlo simulation while using concepts from simulated annealing.

Pfoertner [1] posts the best known radii $r_{i}^{*}$ of $n$ identical spheres that can be packed into a containing sphere of radius one. This problem can be stated as: equal sphere packing (EPP)

$$
\begin{equation*}
\max \quad r_{i} \tag{3.9}
\end{equation*}
$$

Subject to $\left(x_{i}-x\right)^{2}+\left(y_{i}-y\right)^{2}+\left(z_{i}-z\right)^{2} \leq\left(1-r_{i}\right)^{2} \quad i=1, \ldots, n$,

$$
\begin{array}{cl}
\left(x_{i}-x_{j}\right)^{2}+\left(y_{i}-y_{j}\right)^{2}+\left(z_{i}-z_{j}\right)^{2} \geq 4 r_{i}^{2} & 1 \leq i<j \leq n(  \tag{3.11}\\
0 \leq r_{i} \leq 1 & i=1, \ldots, n
\end{array}
$$

The EPP model is equivalent model to the PSS model. In fact, its solution value $r_{i}^{*}$ can be mathematically mapped to the PSS optimal solution value $r^{*}$ :

$$
\begin{equation*}
r^{*}=\frac{1}{r_{i}^{*}} . \tag{3.13}
\end{equation*}
$$

### 3.3 Proposed approach

PSS is a difficult problem. It has an infinite number of alternative optima, an exponential number of local optima which are not globally optimal, and an uncountable set of stationary points - i.e., solutions that satisfy the Karush Kuhn Tucker (KKT) conditions but which do not correspond to local optima. Moreover, the solution to a PSS problem is sensitive to the choice of the initial solution (as is the case of non-linear optimization). Finally, the larger the number of spheres to be packed, the larger the number of local minima and stationary points. Thus, the simple multi-start global optimization strategies need more local minimizations to reach a
global minima [13].
PSS is herein solved using an approach that applies Schittkowski local search (LS) as the intensification strategy and variable neighborhood search (VNS) as the diversification strategy. In sections 3.3.1 and 3.3.2, detail LS and VNS, respectively, are detailed.

### 3.3.1 Schittkowski's local search

An approximate solution to PSS can be obtained using an off-the-shelf NLP solver. However, its quality (i.e., its deviation from the optimal radius) depends both on the initial solution fed to the solver and on its proximity to the global optimum. In addition, there is no guarantee that the obtained solution is a local optimum. In fact, the solution can be a saddle point. Moreover, the solver can stop for various reasons such as no improving search direction, errors in evaluating function and gradient values, violation of regularity conditions, etc.

To overcome the shortcomings of NLP solvers, a PSS problem is approximately solved via a better adapted NLP solver. It applies a NonLinear Programming with Non-Monotone and Distributed Line Search (NLPQLP), the sequential quadratic programming (SQP) algorithm to a non-monotone line search of Schittkowski [63]. NLPQLP is designed to resolve smooth NLP problems. When the solver aborts because of computational errors caused by inaccurate function or gradient evaluations, a non-monotone line search is activated. Internal restarts are performed in case of errors when computing the search direction due to inaccurate derivatives. Additional automated initial and periodic scaling with restarts is implemented. Schittkowski [63] reports that the extensive computational investigation of the code on benchmark instances clearly demonstrated the improvements brought
up by the non-monotone search. He further indicates that more than $90 \%$ of the 306 tested instances, were successfully solved, subject to a stopping tolerance of $10^{-7}$.

NLPQLP requires an initial solution and bounds for the variables of the problem. Providing loose bounds on the variables may hinder the convergence of the solver to a feasible direction or induce computational errors. The radius $r$ can be bounded in two ways. A straightforward lower bound $\underline{r}$ assumes that the spheres are not solid, and that $S$ contains no unused volume. Subsequently, the minimal volume of $S$ is the sum of the volumes of the $n$ spheres. It follows that the radius of this utopian sphere is $\underline{r}=\sqrt[3]{n}$. A tighter lower bound uses the Kepler conjecture [24] that the density of the best packing of $n \geq 2$ non-overlapping spheres cannot exceed $\frac{\pi}{\sqrt{18}}$. Consequently, $\underline{r}=\sqrt[3]{\frac{\pi}{\sqrt{18}} n}$.

The upper bound $\bar{r}$ is estimated by assuming that the spheres can be fitted into a cube of side $c=2\lceil\sqrt[3]{n}$, where $\lceil a\rceil$ is a integer such that $\lceil a\rceil-1<a \leq\lceil a\rceil$. The sphere fitting this cube has a radius $\bar{r}=\sqrt{3\left(\frac{c}{2}\right)^{2}}$.

From a mathematical point of view, having $S$ centered at $(0,0,0)$ or $(\bar{r}, \bar{r}, \bar{r})$ yield equivalent configurations since they can be obtained from each other via a simple transformation. However, LS did not handle negative values of the coordinates will and often failed to converge to a feasible solution in the former case. Thus, each of the coordinates $\left(x_{i}, y_{i}, z_{i}\right)$ of a small sphere $i, i=1, \ldots, n$, are bounded by 1 and $2 \bar{r}-1$. Similarly, these bounds apply to the coordinates $(x, y, z)$ of the containing sphere $S$. Tighter bounds can be used, but in many instances it let to infeasible solutions. Our experimental investigation further indicated that most NLP solvers (including LS) face difficulties in restoring containment feasibility if the initial $r$ fed to the solver is too small (even in the absence of overlap). Very loose bounds are counter-productive too. They cause the local search to converge to local optima
that are too far from the global optimum.
It is believed that feeding the solver with a feasible initial solution speeds up the search process and provides a better estimate for the global minima. Meanwhile many newer NLP approaches claim their robustness with respect to the initial solution and its feasibility. Herein, LS is started from both a feasible solution and randomly generated solutions.

When started from an initial feasible solution, the search is denoted SLS. The feasible solution is obtained using a very simple constructive heuristic which packs the spheres in layers in a cube of side $c$. It positions sphere $i, i=1, \ldots, n$, in $\left(x_{i}, y_{i}, z_{i}\right)$, where $x_{i}=2 d+1, y_{i}=2 b+1, z_{i}=2 a+1$, with $a, b, d$ integers satisfying $i=a c^{2}+b c+d$. This heuristic is obviously very simple and does not yield a compact packing of the spheres. On the other hand, it is very fast, causes no truncation or roundoff errors, offers the spheres enough degrees of freedom to move, and gives LS a feasible descent direction. Evidently, the lattice configuration is another feasible configuration. However, it was discarded since it is suspected to correspond to the optimal packing in many instances and represents a local minima for the others. Thus, it may hinder the search for better quality optima.

When started with a random initial solution (which is not necessarily feasible), the search is denoted as RLS. The random configuration is generated by arbitrarily selecting the coordinates $(x, y, z)$ and $\left(x_{i}, y_{i}, z_{i}\right), i=1, \ldots, n$, from the uniform $[1,2 \bar{r}-1]$, and setting $r=\bar{r}$. When RLS is restarted $k_{M}$ times, the search is denoted as M-RLS (for Multistar RLS).

### 3.3.2 Variable neighborhood search

VNS is a metaheuristic which exploits systematically the idea of neighborhood change, both in descent to local minima and in escape from the valleys which contain them [27, 26, 28, 54, 45]. Its success emanates from three main characteristics. First, a local minimum with respect to one neighborhood is not necessarily a local minimum for another neighborhood. Second, a global minimum must be a local minimum with respect to all possible neighborhood structures. Third, in many problems including PSS, local minima with respect to one or more neighborhoods are relatively close to each other.

In VNS the search is initiated with a feasible solution $\mathbf{u}$ for PSS with a radius $r_{\mathbf{u}}$. It sets $k=1$, and randomly generates a solution $\mathbf{u}^{\prime}$ from a neighborhood $N_{k}(\mathbf{u})$ of $\mathbf{u}$. It then applies a local search to $\mathbf{u}^{\prime}$ (SLS or RLS depending on whether $\mathbf{u}^{\prime}$ is feasible or not) to identify a local minimum $\mathbf{u} " \in N_{k}(\mathbf{u})$. If $r_{\mathbf{u}^{\prime \prime}}$ improves the current solution, the best solution is updated, i.e, the current solution $\mathbf{u}$ is set to $\mathbf{u}$ ". That is, if $r_{\mathbf{u}^{\prime \prime}}<r_{\mathbf{u}^{*}}$, then $r_{\mathbf{u}^{*}}$ is reset to $r_{\mathbf{u}^{\prime \prime}}$, and $\mathbf{u}^{*}$ to $\mathbf{u}$ ". Differently stated, the focus of the search is re-centered on $\mathbf{u}$ ". Consequently, $k$ is reset to 1 and the search is restarted by generating a random solution from the neighborhood $N_{1}(\mathbf{u})$. On the other hand, if $\mathbf{u}$ " does not improve the current solution $r_{\mathbf{u}}$, a different neighborhood of $\mathbf{u}$ is investigated. Specifically, $k$ is incremented, and a different random solution $\mathbf{u}^{\prime}$ from the new neighborhood $N_{k}(\mathbf{u})$ of $\mathbf{u}$ is obtained, and the local search is applied to $\mathbf{u}^{\prime}$ to get $\mathbf{u}$ ". It represents the inner loop of the method. When it is completed, then $k$ is set to 1 and the inner loop is started again. It is repeated until the VNS runtime is smaller than the maximum runtime allowed. The detailed algorithm of VNS is given in Algorithm 1.

A solution $\mathbf{u}^{\prime}$ from a neighborhood $N_{k}(\mathbf{u}), k=1, \ldots, \bar{k}$, of $\mathbf{u}$ is generated

## Algorithm 1 Detailed Algorithm of the VNS for PSS

## Input

1 An initial feasible solution $\mathbf{u}=\left(x_{i}, y_{i}, z_{i}\right)_{i=1, \ldots, n}$, and $r_{\mathbf{u}}$ the corresponding radius of $S$.
$2 \bar{t}$, the maximum runtime for VNS.
$3 \bar{k}$, the maximum number of neighborhoods.

## Output

1 An approximate optimal solution $\mathbf{u}^{*}=\left(x_{i}^{*}, y_{i}^{*}, z_{i}^{*}\right)_{i=1, \ldots, n}$, and its associated radius $r_{\mathbf{u}^{*}}$.

## Algorithm

$1 \quad$ Set $\mathbf{u}^{*}=\mathbf{u}$ and $r_{\mathbf{u}^{*}}=r_{\mathbf{u}}$.
2 Do while $t \leq T$
2.1 Set the neighborhood type $k=1$.
2.2 Do while $k \leq \bar{k}$
2.2.1 Apply the procedure $\operatorname{Shake}\left(k, \mathbf{u}, \mathbf{u}^{\prime}\right)$ to obtain a random solution $\mathbf{u}^{\prime} \in$ $N_{k}(\mathbf{u})$, where $N_{k}(\mathbf{u})$ is the $k^{\text {th }}$ neighborhood of $\mathbf{u}$.
2.2.2 Apply the local search with $\mathbf{u}^{\prime}$ as the starting solution, to get the new solution $\mathbf{u}$ " and its associated value $r_{\mathbf{u}}$ ".
2.2.3 If $r_{\mathbf{u}^{\prime \prime}}<r_{\mathbf{u}^{*}}$, set $k=1, \mathbf{u}^{*}=\mathbf{u}$ ", $r_{\mathbf{u}^{*}}=r_{\mathbf{u}}$, and $\mathbf{u}=\mathbf{u}$ "; else, set $k=k+1$.
using procedure $\underline{\operatorname{Shake}}\left(k, \mathbf{u}, \mathbf{u}^{\prime}\right)$. Let $S_{[k]}$ denote the $k^{\text {th }}$ closest sphere to the center of the containing sphere $S$, and $d_{[k]}$ the distance that separates the centers of $S_{[k]}$ and $S$. For $k \leq n$, procedure $\underline{\operatorname{Shake}}\left(k, \mathbf{u}, \mathbf{u}^{\prime}\right)$ translates sphere $S_{[k]}$ by $\delta_{[k]}^{x}, \delta_{[k]}^{y}$ and $\delta_{[k]}^{z}$ in the $x, y$ and $z$ directions, respectively. The translation distances are randomly generated from the Uniform $[-\Delta, \Delta]$, where $\Delta$ is the neighborhood parameter. If this translation results in the violation of the lower (resp. upper) bound of a variable, this latter is reset to its lower (resp. upper) bound. For $n<k \leq 2 n$, procedure
$\underline{\operatorname{Shake}}\left(k, \mathbf{u}, \mathbf{u}^{\prime}\right)$ translates every sphere $S_{\left[k^{\prime}\right]}, k^{\prime}=1, \ldots, k-n$, by $\delta_{\left[k^{\prime}\right]}^{x}, \delta_{\left[k^{\prime}\right]}^{y}$ and $\delta_{\left[k^{\prime}\right]}^{z}$ in the $x, y$ and $z$ directions, respectively. Finally, for $k>2 n$, procedure $\underline{\operatorname{Shake}}\left(k, \mathbf{u}, \mathbf{u}^{\prime}\right)$ translates every sphere $S_{\left[k^{\prime}\right]}, k^{\prime}=1, \ldots, k-2 n$, as long as its $d_{\left[k^{\prime}\right]} \leq 3.0$. This strategy reflects the fact that smaller variations on the positions of spheres closest to the center of $S$ have a more sizeable impact on the final radius of $S$ than do variations on spheres closer to its surface.

The VNS can only be started from a feasible solution u. Herein, a distinction is made between the cases where this feasible solution is obtained from SLS and from RLS, and the two corresponding searches are denoted as VNS-SLS and VNSRLS. Finally, we consider a final version of the search, denoted M-VNS-RLS, where VNS-RLS is restarted with the $k_{M}$ solutions obtained by M-RLS.

From a pure mathematical point of view, PSS can be solved by setting either $(x, y, z)$ or any $\left(x_{i}, y_{i}, z_{i}\right), i \in\{1, \ldots, n\}$, to $(0,0,0)$ or to any other triplet value of interest. In fact, any solution can be translated in a way that results in one of the spheres been centered at the triplet of interest. However, our extensive computational investigation showed that this is not always advisable. Even though this strategy is supposed to reduce the size of the problem and to alleviate some of the difficulties caused by symmetry, it also limits, in many instances, the capacity of the NLP solver to converge to a good quality local optimum. Subsequently, two versions of the proposed hybrid heuristic are considered: without a fixing position (denoted hereafter with a suffix N ) and with a fixing position (denoted hereafter with a suffix F).

Our computational investigation showed that fixing the three coordinates of a sphere is not judicious in most instances. Therefore, it was decided that the fixing should be limited to only one coordinate of one of the spheres. Because of the special structure of the problem, there is no guideline to the choice of the coordinate
to be fixed. Consequently, when fixing is adopted, procedure $\underline{S h a k e}\left(k, \mathbf{u}, \mathbf{u}^{\prime}\right)$ fixes $x_{[1]}$ if $k \leq n, y_{[1]}$ if $n<k \leq 2 n$, and $z_{[1]}$ if $k>2 n$. The coordinate is fixed at its corresponding value in $\mathbf{u}$ whatever that value is. It is obviously not 0 . Since there is no guarantee that using the VNS with fixing yields better results than without fixing, this search option can be used as a diversification strategy.

Subsequently, VNS is run with three types of diversification strategies: fixing versus not fixing a coordinate of one of the spheres, varying the search time $T$ of the VNS, and varying the neighborhood size $\Delta$. The first strategy, as explained in the previous paragraph, is motivated by the fact that fixing one of the variables would reduce the search space by eliminating a large number of equivalent solutions (such as symmetrical ones, or ones obtained by rotation or reflection, etc.). The last two strategies are motivated by the fact that an enlarged neighborhood, or an intensive search within a neighborhood, could lead to a very good quality local minimum that is better than one obtained with smaller values or a less intensive search. However, this local minimum may cause the search to stagnate and prohibit its escape toward a better quality local minimum.

### 3.4 Computational results

The objective of the computational investigation is threefold: to show that the results obtained using the proposed approach are more accurate and precise than those in the literature; to provide new upper bounds for 14 instances; and to investigate the effect of the three diversification strategies, and point out the utility of the VNS and of the multi-start on the performance of LS.

All codes were written in Fortran using double precision computation. The Schittkowski code was modified for two reasons: the stopping tolerance is $10^{-14}$,
and it is automatically restarted with a new random solution if its current run does not yield a feasible solution that satisfies the optimality conditions or if the current solution is not the best solution encountered during the search. Unless otherwise stated, the VNS heuristics use $\Delta=1, \bar{k}=50$, and $\bar{T}=12$ seconds. All instances were run on a Pentium IV dotted with a 3.20 GHz processor and 4 GB of memory. The results are discussed in the following section.

### 3.4.1 Overall performance

Table 3.1 displays the best-known and best-obtained radii. Column 1 indicates the problem size $n$, where $n=1, \ldots, 50$. Columns 2 to 5 display the best reported radii $\hat{r}_{B}, \hat{r}_{0}, \hat{r}_{1}, \hat{r}_{2}$, where $\hat{r}_{B}, \hat{r}_{0}, \hat{r}_{1}, \hat{r}_{2}$ are obtained from [13, 1, 2, 3], respectively. In Columns 3 to 5 , for a given $n$, the minimum of $\hat{r}_{0}, \hat{r}_{1}$, and $\hat{r}_{2}$ is highlighted. Finally, Columns 6 to 10 report $r_{\mathrm{H}}$, the best radius $r$ obtained by heuristic $\mathrm{H}, \mathrm{H} \in \mathcal{H}=$ \{SLS,VNS-SLS, M-RLS, VNS-RLS, M-VNS-RLS\}, and the $\min _{\mathrm{H} \in \mathcal{H}}\left\{r_{\mathrm{H}}\right\}$ is highlighted. For the VNS heuristics, the reported radii are the best over all runs: F and N , $T=6$ and 12 seconds, and $\Delta=0.1,0.2,0.5,0.75,1.0,1.25,1.5$. For M-RLS and M-VNS-RLS, the reported radii are obtained using the same set of $k_{M}=50$ initial random solutions. In fact, for a given $\Delta, T, \bar{k}$, and neighborhood structure, $r_{\mathrm{SLS}} \geq$ $r_{\mathrm{VNS}-\mathrm{SLS}}, r_{\mathrm{M}-\mathrm{RLS}} \geq r_{\mathrm{M}-\mathrm{VNS}-\mathrm{RLS}}$, and $r_{\mathrm{VNS}-\mathrm{RLS}} \geq r_{\mathrm{M}-\mathrm{VNS}-\mathrm{RLS}}$. Some of these relationships may not hold in the values in Table 3.1 since the displayed results are summarized over all $\Delta, T$ and neighborhood structures.

The analysis of Table 3.1 involves the discussion of issues related to accuracy, precision, and convergence. The results of Birgin and Sobral [13] presented in Column 2 of Table 3.1 are most likely inaccurate. The authors admitted the inaccuracy of their results in the two-dimensional case and proposed improved re-
Table 3.1: Best Local Minima

sults for circle packing [12]. Because they used the same algorithm for packing unit spheres, their results are probably erroneous. For example, when packing two unit spheres into the smallest sphere, the authors obtain $r_{B}=1.9999202433$ while the correct exact answer should be 2.0000000000 . Similarly, when packing 13 circles, $r_{B}=2.9997977416$, while the correct answer should be 3.0000000000 . Thus, no comparison will be undertaken with respect to these results.

The radii $\hat{r}_{0}, \hat{r}_{1}$, and $\hat{r}_{2}$ may lack some precision. They were obtained by mapping the solution of equal sphere packing problem (ESPP) to the solution of SPP. ESPP yields, in general, less precise solutions than SPP because of computational errors due to truncation and their faster rate of propagation for values that are less than 1. However, as this chapter makes no attempt to assess these errors, it is assumed that the precision of $\hat{r}_{0}$ and $\hat{r}_{1}$ is $10^{-7}$ (even though it could be suspected that it does not exceed $10^{-6}$ since $\hat{r}_{0}=\hat{r}_{1}=3.0000003$ for $n=13$ ). Let $\hat{r}=\min \left\{\hat{r}_{0}, \hat{r}_{1}, \hat{r}_{2}\right\}$. Evidently, $\hat{r}$ has the same precision as the least precise of $\hat{r}_{0}, \hat{r}_{1}$, and $\hat{r}_{2}$. On the other hand, $r_{\mathrm{H}}, \mathrm{H} \in \mathcal{H}$, are suspected to be precise to $10^{-13}$, as can be inferred from $r_{\mathrm{H}}, \mathrm{H} \in \mathcal{H}$, for $n=13$. A valid comparison of $r_{\mathrm{H}}$ to $\hat{r}$ requires either truncating or rounding off $r_{\mathrm{H}}$ to $10^{-7}$. However, regardless of the adopted comparison method, the proposed approach improves $\hat{r}$ for $n=3-6,9-14,31,32,38$, and 39.

Finally, in many instances, starting LS with a non-feasible solution causes it to stop the search and return an infeasible solution. The number of random initial solutions that needed to be investigated prior to obtaining one that led to a local minima reached 517 in one instance. This number was, however, much less in most instances and averaged 7 for the $k_{M}$ iterations of M-RLS over and 3 per neighbor for the VNS based heuristics over all runs.

Table 3.2 is a summary of Table 3.1. It reports the number of times $r_{\mathrm{H}}$ is
less than $r_{H^{\prime}},\left(\mathrm{H}, \mathrm{H}^{\prime}\right) \in\left(\mathcal{H} \cup\left\{\hat{H}_{1}, \hat{H}_{2}, \hat{H}_{3}\right\}\right)^{2}$, with the diagonals giving the number of times H is the only heuristic where in the best upper bound is obtained, and the number in parentheses is the number of times H reaches the best upper bound but is not the only heuristic where it is attained. It highlights the role of the VNS in the search. Most of the best solutions are obtained using VNS based heuristics. In fact, both SLS and M-RLS rarely outperform the VNS heuristics. Table 3.2 further shows that M-VNS-RLS is the best heuristic; that is, its multi-start strategy coupled with its variable neighborhood search allowed it to reach local minima that are closer to the global minimum than the other proposed heuristics.

Table 3.2: Number of Times $r_{\mathrm{H}}<r_{\mathrm{H}}$,

| $\downarrow r_{\mathrm{H}} \quad r_{\mathrm{H}} \longrightarrow$ | $r_{\mathrm{SLS}}$ | $r_{\mathrm{VNS} \text {-SLS }}$ | $r_{\mathrm{M}-\mathrm{RLS}}$ | $r_{\mathrm{VNS} \text {-RLS }}$ | $r_{\mathrm{M}-\mathrm{VNS} \text {-RLS }}$ | $\hat{r}_{0}$ | $\hat{r}_{1}$ | $\hat{r}_{2}$ |
| :--- | ---: | ---: | ---: | ---: | ---: | ---: | ---: | ---: |
| $r_{\mathrm{SLS}}$ | $0(1)$ | 0 | 6 | 1 | 5 | 4 | 0 |  |
| $r_{\text {VNS-SLS }}$ | 44 | $1(4)$ | 35 | 16 | 3 | 22 | 15 | 10 |
| $r_{\mathrm{M}-\mathrm{RS}}$ | 42 | 7 | $0(0)$ | 4 | 1 | 18 | 12 | 9 |
| $r_{\mathrm{VNS} \text {-RLS }}$ | 47 | 12 | 34 | $0(4)$ | 0 | 24 | 16 | 12 |
| $r_{\mathrm{M}-\mathrm{VNS} \text {-RLS }}$ | 47 | 28 | 37 | 23 | $0(4)$ | 24 | 16 | 13 |
| $\hat{r}_{0}$ | 43 | 26 | 30 | 24 | 24 | $0(3)$ | 0 | 21 |
| $\hat{r}_{1}$ | 44 | 33 | 36 | 32 | 32 | 32 | $13(5)$ | 29 |
| $\hat{r}_{2}$ |  | 35 | 25 | 26 | 23 | 22 | 14 | 4 |

### 3.4.2 Feasibility of the initial solution

Table 3.2 shows that the number of times $r_{\text {M-RLS }}<r_{\text {SLS }}$ is 42 whereas the number of times $r_{\text {SLS }}<r_{\text {M-RLS }}$ is only 6 (corresponding to $n=3,23,25,39,41,43$ ). This discrepancy is expected and should be interpreted with care since $r_{\text {M-RLS }}$ is the best radius over $k_{M}$ runs of LS; though, it remains true that starting LS from a feasible solution is not always necessary. For example, the cases where $n=26$ and 27 require that LS be started from a very good quality solution. In fact, starting them from a feasible solution constructed using our naive approach leads to radii that
are very far from the best-known upper bounds. Our computational investigation show that starting LS with a lattice configuration yields radii that match the bestknown upper bounds for those two cases. However, these results were omitted from Table 3.1 for consistency. On the other hand, the LS failed to converge to a local optimum for those two instances when fed with randomly generated solutions. MRLS was allowed up to 1000 restarts for each of the $k_{M}$ iterations. Subsequently, it is recommended that the LS search be started from multiple solutions, with a subset of them being feasible and a subset being randomly generated. This strategy guarantees the convergence of M-SLS to a feasible local minimum while it allows it to investigate the existence of other local minima that might be closer to the global minimum. Therefore, this hybrid strategy is used for larger value of $n$. We also run our code for $n=26$ and $n=27$, using the starting solution obtained by $n=28$. We get the following results: Tables 3.3, 3.4.

Table 3.3: VNS with initial obtained by hybrid strategy

| For | $r_{\text {M-RLS }}$ | $r_{\text {VNS-RLS }}$ | $r_{\text {M-VNS-RLS }}$ |
| :--- | :--- | :--- | :--- |
| $\mathrm{n}=26$ | 3.747405776527507 | 3.747405776527507 | 3.747405776527507 |
| $\mathrm{n}=27$ | 3.816151203682777 | 3.846600560763699 | 3.816151203682776 |

Table 3.4: VNS with initial point obtained as for $\mathrm{n}=28$ by removing 1 or 2 centers at random

| For | $r_{\text {M-RLS }}$ | $r_{\text {VNS-RLS }}$ | $r_{\text {M-VNS-RLS }}$ |
| :--- | :--- | :--- | :--- |
| $\mathrm{n}=26$ | 3.788751680052301 | 3.747405776527507 | 3.747405776527507 |
| $\mathrm{n}=27$ | 3.817290771971196 | 3.846600560763699 | 3.846600560763699 |

### 3.4.3 Utility of the diversification strategies

The first diversification strategy consists in varying $\Delta$. Figures 3.3 to 3.8 illustrate the impact is running when the VNS heuristics with and without fixing a coordinate of one of the spheres. They display $\frac{1}{r_{\mathrm{H}}}, \mathrm{H} \in\{$ VNS-SLS, VNS-RLS, M-VNS-RLS $\}$, as a function of $n$ for different neighborhood sizes (measured in terms of $\Delta$ ). They confirm the importance of varying $\Delta$ in many instances such as $\mathrm{n}=13,19-21,24$, 31, 37, 40, 45-49 of Figure 3.3. However, there is no general rule of thumb on how to select the "best" $\Delta$ value. Opting for a larger neighborhood is not necessarily better than using a smaller one. Investigating a large neighborhood is equivalent to too much perturbation of the current feasible solution (which itself corresponds to a local minima). In fact, large perturbations induce "large" infeasibility (i.e., overlap). When the VNS heuristics are run without fixing a coordinate of one of the spheres, opting for small $\Delta$ values seems more judicious (in the instances where the size of the neighborhood matters). In such cases, using $\Delta=0.1$ and 0.2 seem reasonable. However, none of these values of consistently dominate the others. This rule holds when the VNS heuristics are run with the coordinate of one of the spheres being fixed in all instances, except for $n=24$ in Figure 3.6.

The second diversification strategy entails in varying $T$. Starting from the same initial solution, each VNS heuristic $\mathrm{H} \in\{$ VNS-SLS, VNS-RLS, M-VNS-RLS\} was run twice: once with $T=6$ and once with $T=12$. Let $r_{\mathrm{H}}^{(6)}$ and $r_{\mathrm{H}}^{(12)}$ denote the respective resulting radii. The results are summarized in Table 3.5; column 1 indicates whether the algorithms are run with or without fixing a coordinate of one of the spheres. Columns 2 and 3 specify the neighborhood size $\Delta$ and the heuristic H. Columns 4-6 display $\eta_{1}, \eta_{2}, \eta_{3}$, the number of times $r_{\mathrm{H}}^{(12)}>r_{\mathrm{H}}^{(6)}, r_{\mathrm{H}}^{(12)}=r_{\mathrm{H}}^{(6)}$, and $r_{\mathrm{H}}^{(12)}<r_{\mathrm{H}}^{(6)}$, respectively. Column 7 reports $\eta_{4}$ the number of times H fails to






identify a feasible solution when $T=6$. Finally, Columns 8 and 9 report the average and maximum deviations of $r_{\mathrm{H}}^{(6)}$ from $r_{\mathrm{H}}^{(12)}$ over the cases with $r_{\mathrm{H}}^{(6)}<r_{\mathrm{H}}^{(12)}$ whereas the last two columns report the average and maximum deviations of $r_{\mathrm{H}}^{(12)}$ from $r_{\mathrm{H}}^{(6)}$ over the cases with $r_{\mathrm{H}}^{(12)}<r_{\mathrm{H}}^{(6)}$.

Results in Table 3.5 suggest that varying $\Delta$ can be successfully used as a diversification strategy of VNS. Using $T=6$ yields more frequent better-quality solutions than using $T=12$. However, the average and maximum improvements brought up by using $T=12$ when $r_{\mathrm{H}}^{(12)}<r_{\mathrm{H}}^{(6)}$ are more sizeable than their counterparts when $r_{\mathrm{H}}^{(12)}=r_{\mathrm{H}}^{(6)}$.

The cases when the magnitude of $T$ does not affect the solution quality are numerous, and the number is more important when H is run with a fixed coordinate than when it is not fixed and for $\Delta=1.0$ than for $\Delta=0.5$. It also depends on the heuristic H (most probably on the quality of the initial solution fed to H ).

When $\Delta=1.0$, the VNS heuristics fail in $35 \%$ of the instances to identify a feasible solution when $T=6$, for either the F or N versions of the algorithm. This percentage is twice as large (16\%) for the F version than for the N version (8\%) when $\Delta=0.5$. This suggests that using $\Delta=1.0$ while fixing a coordinate of one of the spheres may result in infeasibility that cannot be resolved by the VNS within the preset time limit $T=6$; hence, none of the solutions fed to LS would yield a feasible solution.

The third diversification strategy consists in running each of the VNS heuristics twice: starting from the same initial solution, $\mathrm{H} \in\{$ VNS-SLS, VNS-RLS, M-VNS-RLS\}, and run once with and once without fixing a coordinate of one of the spheres; denoted by $\mathrm{H}^{\mathrm{F}}$ and $\mathrm{H}^{\mathrm{N}}$ respectively. The results are summarized by Table 3.6 . Columns 1 and 2 specify the neighborhood size $\Delta$ and the heuristic H .
Table 3．5：Impact of $T$

|  |  |  |  |  | BOB ※䟵 M 응응 000 <br> 下o $\stackrel{\rightharpoonup}{7}$曾名名 $\circ \circ$ － |
| :---: | :---: | :---: | :---: | :---: | :---: |
|  |  |  |  |  |  <br> 청충융 덩융 항 <br>  응 000 |
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|  | $\begin{aligned} & \approx \\ & \stackrel{\circ}{\approx} \\ & \stackrel{y}{c} \end{aligned}$ |  | $\begin{aligned} & 000 \\ & \approx 0 \wedge \\ & \exists \pm 0 \end{aligned}$ | 102010 |  |
|  | I $\triangleleft$ |  |  |  |  |

Columns 3 to 5 display $\eta_{1}, \eta_{2}, \eta_{3}$, the number of times $r_{\mathrm{H}^{\mathrm{N}}}<r_{\mathrm{H}^{\mathrm{F}}}, r_{\mathrm{H}^{\mathrm{F}}}=r_{\mathrm{H}^{\mathrm{N}}}$, and $r_{\mathrm{H}^{\mathrm{N}}}>r_{\mathrm{H}^{\mathrm{F}}}$, respectively. Column 6 reports $\eta_{4}$, the number of times the VNS heuristic fails to identify a feasible solution. Finally, Columns 8 and 9 report the average and maximum deviations $r_{\mathrm{H}^{\mathrm{N}}}-r_{\mathrm{H}^{\mathrm{F}}}$ when $r_{\mathrm{H}^{\mathrm{F}}}<r_{\mathrm{H}^{\mathrm{N}}}$, whereas Columns 10 and 11 give the average and maximum deviations $r_{\mathrm{H}^{\mathrm{F}}}-r_{\mathrm{H}^{\mathrm{N}}}$ when $r_{\mathrm{H}^{\mathrm{F}}}>r_{\mathrm{H}^{\mathrm{N}}}$.

These results show that fixing the coordinate of one of the spheres is neither always preferable nor undesirable. In fact, the maximum deviations of $\left|r_{\mathrm{H}^{\mathrm{F}}}-r_{\mathrm{H}^{\mathrm{N}}}\right|$ can reach 0.012: a very important variation for the problem at hand. Subsequently, in many instances, reducing the search space (with the objective of eliminating symmetry and related issues) is not advisable. Thus, a successful search strategy should opt for testing both approaches of the problem, as is further elucidated by the detailed results for $\mathrm{H}^{\mathrm{F}}$ and $\mathrm{H}^{\mathrm{N}}$ in Table 3.7.

### 3.4.4 Utility of the VNS and the LS

The best radii obtained by $\mathrm{H}^{\mathrm{F}}$ and $\mathrm{H}^{\mathrm{N}}$ when $\Delta=0.5$ and $T=12$ is given in Table 3.7. Column 1 indicates the problem size $n$. Columns 2 and 3 display $r_{\text {SLS }}$ and $r_{\text {M-RLS }}$. Columns 3 to 5 report $r_{\mathrm{H}}^{\mathrm{F}}$, the best radius $r$ obtained by heuristic $\mathrm{H}^{\mathrm{F}}, \mathrm{H}$ $\in\{$ VNS-SLS, VNS-RLS, M-VNS-RLS $\}$, whereas Columns 6 to 8 display the same information but for $\mathrm{H}^{\mathrm{N}}$.

The result in Table 3.7 confirm that initiating LS from a feasible solution always yields a local minimum. This is obviously not the case for infeasible solutions. For instance, in the 48 instances tested above, LS failed to reach a local minimum when started from a large set of randomly generated solutions. It is also showed that the LS obtains a better solution when started with a feasible solution in ten out of 48 instances, with the maximum deviation of $r_{\text {SLS }}$ from $r_{\text {M-RLS }}$ reaching

| One of the Spheres |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: |
|  | $\begin{aligned} & x \\ & y \\ & i \\ & i \\ & z \\ & z \\ & y \end{aligned}$ |  |  |  |
| U |  | \# | - + | ง ง ง |
| $\begin{aligned} & \ddot{0} \\ & \ddot{0} \\ & 0 \\ & 0 \end{aligned}$ |  |  |  | $10 \infty$ <br> $\underset{\sim}{\infty} \because=$ <br> ํㅜN우 |
|  |  | $\triangleleft$ |  |  |

12

\begin{tabular}{|c|c|c|c|}
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\hline \&  \&  \& $\infty$

$\sim$ <br>
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\hline \& ¢ $n$＋上 \&  \& <br>
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\end{tabular}

0.148309392561900 . On the other hand, LS obtains a better solution when started from randomly generated solutions in 36 out of 48 instances with the maximum deviation being 0.392854138549501 . It seems therefore judicious to include an initial feasible solution among the solutions investigated by LS to guarantee its convergence to a local optimum and its improved overall performance.

### 3.4.5 Comparison of the diversification strategies

## VNS-SLS versus SLS

As expected, $r_{\mathrm{SLS}} \geq r_{\mathrm{VNS}-\mathrm{sLS}}$. Applying $\mathrm{VNS}^{\mathrm{F}}$ to the local optimum obtained by SLS improved $r_{\text {SLS }}$ in 44 out of 48 instances. The improvement reaches 0.392855382254420 , and averages 0.055338061287952 over the 48 instances. Similarly, VNS ${ }^{N}$ improves $r_{\text {SLS }}$ in 43 out of 48 instances. The improvement reaches 0.392855382254420 , and averages 0.063762149389844 over the 48 instances.

## VNS-RLS versus VNS-SLS

VNS $^{\mathrm{F}}$-RLS yields better solutions than $\mathrm{VNS}^{\mathrm{F}}$-SLS in 25 out of 48 instances, with the maximum deviation of $r_{\mathrm{VNS} \text {-SLS }}$ from $r_{\mathrm{VNS}-\mathrm{RLS}}$ reaching -0.105594271176310 . The opposite case occurs in 16 instances with the maximum deviation equaling 0.091145444888800 . The overall average deviation is -0.013753483385387 . Similarly, $\mathrm{VNS}^{\mathrm{N}}$-RLS yields better solutions than VNS-SLS in 27 out of 48 instances, with the maximum deviation of $r_{\mathrm{VNS}-\mathrm{SLS}}$ from $r_{\mathrm{VNS}-\mathrm{RLS}}$ reaching -0.097886967409520 . The opposite case occurs in 12 instances with the maximum deviation equaling 0.091145444888900 . The overall average deviation is -0.010181470957522 . This suggests that VNS should be started from a feasible solution when not used in con-
junction with a multi-start strategy.

## VNS-RLS versus M-RLS

The hybridization of VNS and LS can provide a powerful search method if it addresses the two competing goals of meta-heuristics: exploration and exploitation. Exploration allows an extensive search of the solution space in order to determine the part of the space that has a higher chance of containing the global optimum whereas exploitation refines the search and focuses on the part of the space that has a high potential of containing the global optimum. Exploration can be herein obtained via multi-start or via VNS whereas exploitation is obtained via LS. Both the multi-start and VNS strive for global optimization while LS strives for local optimization in the global optimum neighborhood. The effectiveness of using VNS and multi-start as the diversification strategies is discussed here.

The findings presented in Table 3.7 does not provide a clear answer to this dilemma. It suggests using a mixture of the two search strategies. In fact, $r_{\mathrm{VNS}-\mathrm{RLS}}^{\mathrm{F}}<$ $r_{\mathrm{M}-\mathrm{RLS}}^{\mathrm{F}}$ in 22 out 48 instances, while $r_{\mathrm{VNS}-\mathrm{RLS}}^{\mathrm{F}}>r_{\mathrm{M}-\mathrm{RLS}}^{\mathrm{F}}$ in 18 out 48 instances. The maximum deviation in the first case equals -0.148309392561900 while it equals 0.097886967409590 in the second case. Its average is 0.004562699793446 . Similarly, $r_{\mathrm{VNS}-\mathrm{RLS}}^{\mathrm{N}}<r_{\mathrm{M}-\mathrm{RLS}}^{\mathrm{N}}$ in 28 out 48 instances while $r_{\mathrm{VNS}-\mathrm{RLS}}^{\mathrm{N}}>r_{\mathrm{M}-\mathrm{RLS}}^{\mathrm{N}}$ in 15 out 48 instances. The maximum deviation in the first case equals -0.148309392561900 while it equals 0.097886967409520 in the second case. Its average is -0.006120725568965 . These different magnitudes of improvement over all instances of both versions of the algorithms suggest that both diversification and intensification of the search are important; one dose not predominate the other in any case.

## M-VNS-RLS versus VNS-RLS

Comparison of Columns 5 and 6 and Columns 8 and 9 of Table 3.7 highlights the importance of the diversification of the search by starting VNS-RLS from a number of randomly generated solutions. In fact, the additional investigation brings forth sizeable improvements. For $\mathrm{VNS}^{\mathrm{F}}$, an improvement is registered in 32 out of 48 instances with the maximum improvement reaching 0.120916908092260 and averaging 0.021911628194197 over the 48 instances. The same trend is observed for VNS $^{N}$ where an improvement is registered in 37 out of 48 instances with the maximum improvement reaching 0.102674008819120 and averaging 0.019515894278337 over the 48 instances.

## M-VNS-RLS versus M-RLS

Comparison of Column 3 to Columns 6 and 9 of Table 3.7 quantifies the importance of diversifying the search by applying VNS to each of the $k_{M}$ local optima obtained by M-RLS. The larger the discrepancy between $r_{\mathrm{M}-\mathrm{VNS}-\mathrm{RLS}}$ and $r_{\mathrm{M}-\mathrm{RLS}}$, the more important the role of VNS is, in identifying the global optima. For VNS ${ }^{\mathrm{F}}$, an improvement is observed in 36 out of 48 instances with the maximum improvement reaching 0.148309392561900 and averaging 0.014861185920063 over the 48 instances. Similarly, for $\mathrm{VNS}^{\mathrm{N}}$, an improvement is observed in 42 out of 48 instances with the maximum improvement reaching 0.148309392561900 and averaging 0.022481100856624 over the 48 instances.

### 3.5 Conclusion

A variable neighborhood search approach is proposes in this chapter, to solve the problem of packing $n$ unitary spheres into the smallest containing sphere $S$ where the objective is to identify the radius of $S$ and a feasible configuration of the uni-
tary spheres within $S$. The approach follows the recent research trends of combining search heuristics with non-linear programming tools and of balancing exploitation and exploration. Indeed, it ensures exploitation by applying a local search based on a sequential quadratic programming algorithm with a non-monotone line search (as a non-linear programming solver) and exploration by applying a variable neighborhood search and a multi-start strategy. The approach provides accurate results with a precision of $10^{-13}$. It can be extended to the case of non-identical spheres and to the packing of spheres into other 3-dimensional shapes.

## Chapter 4

## Packing Spheres in a Cube

### 4.1 Introduction

In this chapter, the adaptation of the VNS to the problem of packing sphere in a cube is explored. An introduction to the VNS has been given in section 2.1 of chapter 2 , so we will not repeat its rules here.

Sphere packing in a cube is an optimization problem, that could be classified either as continuous and or as discrete. The positions of the spheres are presented in continuous variables, whereas the structure of an optimal configuration is discrete [49]. A successful solution technique should tackle these two aspects simultaneously.

Accordingly, we approximately solve the problem of packing unit spheres into the smallest cube (PSC) using a variable neighborhood search (VNS) that combines these two aspects. A VNS addresses the discrete aspect of the PSC by shaking one or more spheres and its continuous aspect by applying a non-linear programming (NLP) optimizer that identifies a local optimum within the neighborhood of the fed solution. In addition, it yields local optima that are nearer to the global optimum,
in a more straightforward manner than do other meta-heuristics. The VNS has successfully been applied to solve many problems, including integer and continuous optimization problems [52].

For 3D packing of spheres the literature is not as extensive as for solving correspond 2D packing. However, it is clear that mathematical programming formulations could easily be extended from 2D to 3D. As far as we know, there is only one journal paper considering 3D packing of spheres into a cube by Birgin and Sobral [13]. The authors considered the objects in different shapes as triangle, circle, square and strips. More detailed description of their work is given in chapter 3. Other two references that we will use in computational results to compare with, we take from websites [4, 5] that just report the best known values. However, it is not clear what methods they used to report those values.

### 4.2 Mathematical model

PSC comprises of packing $n$ identical spheres, of radius one, without overlap into the smallest containing cube $C$. The goal is to search for the best packing of the $n$ spheres into $C$, where the best packing minimizes $w$, the length of the side of $C$. According to the typology of Wascher et al. [79], PSC is a three-dimensional variant of the Open Dimension Problem, since all small items (which are spheres) have to be packed and the extension of the large object (which is a cube) is not given but has to be minimized. PSC is equivalent to finding the coordinates $\left(x_{i}, y_{i}, z_{i}\right)$ of every sphere $i, i=1, \ldots, n$, and the side $w$ of $C$. There are several equivalent nonlinear programming formulations of PSC. The simplest one is given in Costa et al. [7] where objective is to maximize the small radius of sphere.

$$
\begin{array}{ll}
\text { min } & w \\
\text { s.t. } & \\
& \left(x_{i}-x_{j}\right)^{2}+\left(y_{i}-y_{j}\right)^{2}+\left(z_{i}-z_{j}\right)^{2} \geq w \\
0 \leq x_{i} \leq 1 & 1 \leq i<j \leq n, \\
0 \leq y_{i} \leq 1 & i=1, \ldots, n, \\
& i=1, \ldots, n \\
& i=z_{i} \leq 1 \tag{4.6}
\end{array}
$$

In this chapter we will use the following model [13]

$$
\begin{array}{cc}
\text { min } & w \\
\text { s.t. } & \\
& \left(x_{i}-x_{j}\right)^{2}+\left(y_{i}-y_{j}\right)^{2}+\left(z_{i}-z_{j}\right)^{2} \geq 4 \\
1 \leq x_{i} \leq w-1 & 1 \leq i<j \leq n \\
& i=1, \ldots, n \\
& \\
& i=y_{i} \leq w-1 \\
& \\
& i \leq z_{i} \leq w-1 \tag{4.12}
\end{array}
$$

where $\underline{w}=\sqrt[3]{\frac{4}{3} \pi n}$ is a lower bound on $w[13]$. The first set of constraints ensures there is no overlap of any pair of distinct spheres. Note that this set of constraints could be replaced with a single one as [13]:

$$
\begin{equation*}
\sum_{i=1}^{n-1} \sum_{j=i+1}^{n} \max \left\{0,4-\left(x_{i}-x_{j}\right)^{2}-\left(y_{i}-y_{j}\right)^{2}-\left(z_{i}-z_{j}\right)^{2}\right\}=0 \tag{4.14}
\end{equation*}
$$

The next three sets (4.10-4.12) ensure the containment of every sphere within $C$.
PSC assumes that the bottom leftmost lowest of $C$ coincides with the origin $(0 ; 0 ; 0)$. PSC has an in infinite number of alternative optima caused by symmetrical configurations. Fixing the position of one of the spheres would reduce the search space by eliminating a large number of equivalent solutions [8, 16, 17, 45, 61], without necessarily loosing subspaces including good quality local optima. Herein, we set two of the coordinates of a randomly chosen sphere to 1 so that the solver maintains this sphere adjacent to two edges of the cube during its search. Additional bounding constraints that lexicographically ordered the spheres were investigated but not adopted in this model. Solving PSC via an NLP solver only is generally not successful [5]. Subsequently, PSC is solved by VNS.

### 4.3 Variable neighborhood search-based algorithm for the PSC problem

The proposed VNS is to solve the PSC problem detailed in Algorithm 2. It is in fact very similar to one described in chapter 3 for solving PSS. It starts its search from a feasible solution $\mathbf{u}=\left(x_{1}, y_{1}, z_{1}, \ldots, x_{n}, y_{n}, z_{n}\right)$ for PSC with a side length $w_{\mathbf{u}}$. The coordinates of $\mathbf{u}$ by randomly generating by using uniform distribution from the interval $(1,2 \sqrt[3]{n})$. Any overlapping spheres are moved to out-skirt of $C$. VNS sets the best solution $\mathbf{u}^{*}=\mathbf{u}$, and its upper bound $w_{\mathbf{u}}^{*}=w_{\mathbf{u}}$. It initializes its cumulated
runtime $t$ is 0 and $k=1$.

## Algorithm 2 Detailed Algorithm of the VNS for PSC

## Input

1 A feasible solution $\mathbf{u}=\left(x_{i}, y_{i}, z_{i}\right)_{i=1, \ldots, n}$, and $w_{\mathbf{u}}$ its side length of $C$.
$2 T$, the maximal runtime for VNS .
$3 \bar{k}$, the maximal number of neighborhoods.

## Output

1 A (near-) optimal solution $\mathbf{u}^{*}=\left(x_{i}^{*}, y_{i}^{*}, z_{i}^{*}\right)_{i=1, \ldots, n}$, and its side length $w_{\mathbf{u}^{*}}$.

## Algorithm

$1 \quad$ Set $\mathbf{u}^{*}=\mathbf{u}$ and $w_{\mathbf{u}^{*}}=w_{\mathbf{u}}$.
2 Set the algorithm's cumulated runtime $t$ to 0 .
3 Do while $t \leq T$
3.1 Set the neighborhood type $k=1$.
3.2 Do while $k \leq \bar{k}$
3.2.1 Generate a random solution $\mathbf{u}^{\prime}$ from the $k^{\text {th }}$ neighborhood $N_{k}(\mathbf{u})$ of $\mathbf{u}$.
3.2.2 Starting from $\mathbf{u}^{\prime}$, find a local minimum $\mathbf{u}$ " and its side length $w_{\mathbf{u}}$.
3.2.3 If $w_{\mathbf{u}}$ " $<w_{\mathbf{u}}$,
set $k=1, \mathbf{u}=\mathbf{u} "$, and $w_{\mathbf{u}}=w_{\mathbf{u}}$; if $w_{\mathbf{u}^{\prime}}<w_{\mathbf{u}^{*}}$, set $\mathbf{u}^{*}=\mathbf{u}$ ", and $w_{\mathbf{u}^{*}}=w_{\mathbf{u}}$.
Else
set $k=k+1$.
3.2.4 Update $t$.

The VNS is an iterative procedure with two loops. The outer loop controls the runtime whereas the inner one undertakes the search. At each iteration of the inner loop, the VNS generates a random solution $\mathbf{u}^{\prime}$ from a neighborhood $N_{k}(\mathbf{u})$
of $\mathbf{u}$. It finds a local minimum $\mathbf{u} " \in N_{k}(\mathbf{u})$ by applying NLP to the PSC starting from $\mathbf{u}^{\prime}$. The selected NLP solver is NLPQLP, a sequential quadratic programming algorithm with a non-monotone line search [63]. The NLPQLP is designed for smooth NLP problems. When the solver aborts because of computational errors caused by inaccurate function or gradient evaluations, a non-monotone line search is activated. Internal restarts are performed in case of errors when computing the search direction due to inaccurate derivatives. Additional automated initial and periodic scaling with restarts are implemented [63].

If $w_{\mathbf{u}^{\prime \prime}}<w_{\mathbf{u}}$, the current solution $\mathbf{u}$ is updated; i.e., the focus of the search is re-centered on $\mathbf{u}$ ", and $k$ is reset to 1 . Furthermore, if $w_{\mathbf{u}^{\prime \prime}}<w_{\mathbf{u}^{*}}$, then $w_{\mathbf{u}^{*}}$ is reset to $w_{\mathbf{u}^{\prime \prime}}$, and $\mathbf{u}^{*}$ to $\mathbf{u}$ ".

On the other hand, if $w_{\mathbf{u}}{ }^{\prime \prime} \geq w_{\mathbf{u}}$, the neighborhood of $\mathbf{u}$; i.e., $k$ is incremented. A random solution $\mathbf{u}^{\prime} \in N_{k}(\mathbf{u})$ is chosen, and the local optimum $\mathbf{u}$ " within $N_{k}(\mathbf{u})$ is identified. The neighborhood is enlarged until one of the two stopping criteria is met:

* A solution $\mathbf{u}$ " $\in N_{k}(\mathbf{u})$ such that $w_{\mathbf{u}}$ " $<w_{\mathbf{u}}$ is obtained; in which case, $k$ is reset to 1 and the inner loop is repeated.
* The number of investigated neighborhoods $k$ reaches the threshold $\bar{k}$; in which case the algorithm's total runtime $t$ is updated and the control of the algorithm is transferred to the outer loop of the VNS.

The outer loop is the stopping criterion of the VNS. It compares $t$ to the maximum runtime allowed $T$. VNS stops when $t$ exceeds $T$.

The iterative step generates a solution $\mathbf{u}^{\prime}$ from a neighborhood $N_{k}(\mathbf{u}), k=$ $1, \ldots, \bar{k}$, of $\mathbf{u}$ using a shaking procedure. For $k=1$, this procedure translates one randomly chosen sphere by $\delta^{x}, \delta^{y}$ and $\delta^{z}$ in the $x, y$ and $z$ directions, respectively.

For $1<k<\bar{k}$, the procedure translates the $\min \left\{\frac{n}{2}, \bar{k}\right\}$ randomly chosen spheres in the $x, y$ and $z$ directions, respectively. Finally, when $k=\min \{\bar{k}, n-1\}$, it moves the $k$ closest neighbors of a randomly chosen sphere; thus inducing variations on the position of a cluster of spheres. The parameter $\bar{k}$ is set experimentally to $\min \left\{\frac{n}{2}, 6\right\}$. In all cases, the translation distances are randomly generated from the Uniform $[-\Delta, \Delta]$, where $\Delta$ is the neighborhood parameter. If a translation causes the violation of the lower (resp. upper) bound of a variable, this latter is reset at its lower (resp. upper) bound. The shaking procedure is very useful since it offers multiple starts for NLPQLP; thus, enhances its chances of identifying a (near-) global optimum.

### 4.4 Computational results

In this section we provide computational results obtained by our method for solving PSC problem. This section will be much smaller in size than one in chapter 3, since we use experience from extensive computational analysis from chapter 3. In other words, we use variant of the model with fixing the cube centre, assuring that all variables are nonnegative. As a nonlinear solver, we use NLPQLP [63]. Further, we use here the better initial solution method from there. We generate random initial solution for equivalent maximization of small circle problem. In order to preserve feasibility, the initial value of $r$ is set to 0.001 . The transformation to minimization problem is then performed to be in appropriate form for the model we use. For the parameter $\Delta$ we always use value 0.5 . The value of VNS parameter $T$ is again set to 6 seconds for each independent run. Each test instance is run 10 times and best values reported. The value of $\bar{k}$ is chosen as $\min \{n, 50\}$.

The results are presented at Table 4.1. Column 1 indicates the problem size

Table 4.1: Comparing the Best Local Minima to the Best Known Radii

| $n$ | $\hat{w}_{B}$ | $\hat{w}$ | $\hat{w}_{P}$ | $w^{*}$ |
| :---: | :---: | :---: | :---: | :---: |
| 3 | 3.4141527373 | 3.4142137817 | 3.4142135624 | 3.4142135624 |
| 4 | 3.4141480640 | 3.4142137817 | 3.4142135624 | $\underline{3.4142135624}$ |
| 5 | 3.7887647862 | 3.7888547050 | 3.7888543820 | 3.7888543820 |
| 6 | 3.8855478230 | 3.8856182776 | 3.8856180832 | $\underline{3.8856180832}$ |
| 7 | 3.9977338382 | 3.9978227857 | 3.9978227238 | $\underline{3.9978227238}$ |
| 8 | 3.9999345267 | 4.0000000000 | 4.0000000000 | 4.0000000000 |
| 9 | 4.3093402273 | 4.3094012173 | 4.3094010768 | $\underline{4.3094010768}$ |
| 10 | 4.6666081946 | 4.6666669778 | 4.6666666667 | $\underline{4.6666666667}$ |
| 11 | 4.8159601052 | $\underline{4.8164394712}$ | 4.8164397406 | 4.8164397406 |
| 12 | 4.8280686639 | 4.8284278518 | 4.8284270236 | $\underline{4.8284270519}$ |
| 13 | 4.8282131208 | 4.8284278518 | 4.8284271247 | $\underline{4.8284271247}$ |
| 14 | 4.8282912625 | 4.8284278518 | 4.8284271247 | 4.8284271247 |
| 15 | 5.1997445225 | 5.1999997920 | 5.2000000000 | 5.2000000000 |
| 16 | 5.2964155013 | 5.2967013468 | 5.2967008293 | $\underline{5.2967008293}$ |
| 17 | 5.2996749577 | 5.2998313329 | 5.2998316455 | 5.2998316455 |
| 18 | 5.3279843248 | 5.3282026294 | 5.3282011774 | 5.3282011774 |
| 19 | 5.4586383844 | 5.4589562912 | 5.4589532170 | $\underline{5.4589532170}$ |
| 20 | 5.6048721029 | 5.6051549489 | 5.6051549576 | 5.6051549576 |
| 21 | 5.6431452803 | 5.6427342320 | 5.6427344101 | 5.6430934602 |
| 22 | 5.7710452688 | 5.7712362738 | 5.7712361663 | $\underline{5.7712361663}$ |
| 23 | 5.8199237336 | 5.8201577088 | 5.8201531911 | 5.8201531911 |
| 24 | 5.8633943713 | 5.8637080741 | 5.8637033052 | $\underline{5.8637033052}$ |
| 25 | 5.9589662351 | 5.9593423866 | 5.9593308221 | $\underline{5.9593308221}$ |
| 26 | 5.9952148260 | 5.9914286622 | 5.9914262468 | 5.9955822990 |
| 27 | 5.9998355203 | 6.0000006000 | 6.0000000000 | 6.0000000000 |
| 28 | 6.2421317695 | 6.2425479584 | 6.2425477277 | $\underline{6.2425477277}$ |
| 29 | 6.2423824504 | 6.2426414863 | 6.2426406871 | 6.2426406871 |
| 30 | 6.2424721518 | 6.2426414863 | 6.2426406871 | $\underline{6.2426406871}$ |
| 31 | 6.2425174557 | 6.2426414863 | 6.2426406871 | $\underline{6.2426406871}$ |
| 32 | 6.2425285827 | 6.2426414863 | 6.2426406871 | $\underline{6.2426406871}$ |
| 33 | 6.4689813902 | 6.4680814732 | 6.4680780465 | $\underline{6.4680780465}$ |
| 34 | 6.5735658875 | 6.5738915268 | 6.5738831658 | $\underline{6.5738831658}$ |
| 35 | 6.5931612342 | 6.5933290674 | 6.5933259094 | $\underline{6.5933259094}$ |
| 36 | 6.6970889746 | 6.6975333654 | 6.6944181198 | $\underline{6.6975126034}$ |
| 37 | 6.7083709308 | 6.7086358928 | 6.7086344826 | $\underline{6.7086344826}$ |
| 38 | 6.7093947495 | 6.7096644329 | 6.7096635745 | $\underline{6.7096635745}$ |
| 39 | 6.7739983433 | 6.7742714525 | 6.7742701172 | 6.7742701172 |
| 40 | 6.7998570644 | 6.8000010880 | 6.8000000000 | $\underline{6.8000000000}$ |
| 41 | 6.9039667243 | 6.9042869063 | 6.9042712513 | $\underline{6.9042712513}$ |
| 42 | 6.9906644142 | 6.9909219383 | 6.9907863116 | $\underline{6.9907863116}$ |
| 43 | 7.0610542561 | 7.0595743359 | 7.0595386299 | 7.0595386299 |
| 44 | 7.0991542522 | 7.0992323600 | 7.0992166891 | 7.0992166891 |
| 45 | 7.1269867567 | $\underline{7.1107596816}$ | 7.1107603214 | 7.1107603214 |
| 46 | 7.1396025069 | 7.1302925488 | 7.1302839201 | $\underline{7.1302839201}$ |
| 47 | 7.1447631846 | 7.1449572499 | 7.1449567477 | 7.1449574116 |
| 48 | 7.2254788705 | 7.1449572499 | 7.1449575543 | 7.1449575543 |
| 49 | 7.3396050716 | 7.3299434788 | 7.3299402906 | 7.3312658497 |
| 50 | 7.3606467872 | 7.3554452729 | 7.3554013072 | 7.3598400994 |
| Average | $\mathbf{5 . 8 3 2 1 6 6 5 0 3 9}$ | 5.8296846262 | $\mathbf{5 . 8 2 9 6 1 3 2 0 7 4}$ | 5.8298918455 |
| 51 |  | 7.4070535149 | 7.4061413497 | 7.4263293310 |
| 52 |  | 7.4727301395 | 7.4727224395 | 7.4729829434 |
| 53 |  | $\underline{7.5058057407}$ | 7.5057697640 | 7.5561707846 |
| 54 |  | 7.5967666642 | 7.5641117363 | 7.6046039060 |
| 55 | 7.6497254384 | 7.6402978341 | 7.6207736668 | $\underline{7.6392010852}$ |

$n$. Columns 2,3 and 4 display the side length $\hat{w}_{B}, \hat{w}$ and $\hat{w}_{P}$ obtained by [13], [4] and [5], respectively. Column 5 reports the side length $w^{*}$ obtained by VNS. The configurations corresponding to $\hat{w}_{B}$ are not necessarily feasible. They may
have overlapping spheres or spheres that are not totally contained within the cube. Thus, $\hat{w}_{B}$ will not be used for comparison purposes. For each instance of Table 4.1, the underlined value indicates the tightest upper bound. In column 4 we give the currently best known solutions from web site Packomania. However, we did not compare our results with those from Packomania, since they are obtained by many different people and many different methods. Thus, although we report results obtained by 3 different sources, we compare ours only with those from [4].

The following observations one can get from the results of Table 4.1: (i) VNS improves the best solutions obtained by [4] in 35 out of 48 instances; it represents $73 \%$ of the cases. (ii) The best improvement of an existing upper bound is 0.0010967489 occurring for $n=55$, with the improvement averaging 0.0000391436 over the 35 improved instances. These improvements are important despite their seemingly small magnitude. (iii) Results obtained by [4] are of better quality for large values of $n$. This clearly indicates that another parameter values for our VNS should be chosen. That could be a task for the future work.

### 4.5 Conclusion

In this chapter we apply Variable neighbourhood search (VNS) approach for solving sphere packing problem within smallest containing cube, in 3D. VNS is framework for building heuristics. It starts from initial solution and use different neighbourhoods of that solution in order to improve it. The perturbation or shaking phase of the current solution is obtained by moving $k(k=1, \ldots, \bar{k})$ sphere centres for the $\Delta$ (a parameter) in each dimension. As a local search it is used well-known software for solving nonlinear convex problems [63]. It appears that 35 out of 48 better results are reported when compared with current state of the art. Future work may contain
application of this approach to other packing problems in different containers. In addition, better parameter estimation for large values of $n$ could be performed.

## Chapter 5

## Conclusion

### 5.1 Summary

The Variable Neighborhood Search (VNS) is one of the most efficient general search frameworks. It has several heuristic variations, and can easily be adapted to continuous and discrete combinatorial problems. This thesis presents a tutorial along with a detailed literature review on recent applications of the VNS and its variants. As for other meta-heuristics, the hybridization of the VNS with other approximate or exact search algorithms enhances its efficiency and efficacy. This thesis applies a hybrid VNS to approximately solve the problem of three-dimensional circular packing, where the containing object is spherical or cubical and the items are unit spheres. This problem is relevant to many real-life applications. A successful application of a VNS-based heuristic requires a good definition of the representation of the solution, its neighbourhoods, its moves within a neighbourhood, and its local search. In this context, the neighbourhood has a special structure since it is continuous but disconnected. The proposed VNS implementation represents a solution by the
three-dimensional coordinates of its solutions and its solution value by the radius of the containing sphere or by the length of the side of the containing cube. It generates a neighbouring solution by shaking one or more spheres (depending on the neighbourhood size). It applies a local search by applying the non-linear programming search technique that had obtained a large percentage of best near-global optima for other classes of complex problems while ensuring high precision. The proposed heuristic is restarted a fixed number of times so that it benefits from the diversification induced by different initial starting points. The computational results provide computational proof of the efficiency and efficacy of the VNS-based heuristic. When the containing object is a sphere, our best method is able to improve $60.4 \%$ of best known solutions and matches all other results. When the containing object is a cube, it improves $76.4 \%$ of existing solutions. Many of these solutions are suspected to be optimal, and any improvement is due a large computational precision. Indeed, the proposed approach has a higher precision level than most of the state-of-the-art approaches.

### 5.2 Future research

Future research might include applying the VNS to the simpler two-dimensional shapes or to more complex p-dimensional containers such as rectangular, triangular, pyramidal, and strip-shaped. Different variants of the problem may require different neighbourhood structures and/or different moves. Future and undergoing research concerns the augmentation and reduction of the problem via linearization and reformulation techniques. In real life packing of spheres, optimizes space usage is the most significant goal; yet, different issues are likewise considered. For example, cargo steadiness, multi-drop loads or weight circulations are also pertinent and
should be considered during the packing or loading to ensure cargo security. Thus, adding these constraints into the three-dimensional packing problem is imperative for enhancing the applicability of this research. Such extensions could focus on the efficiency of space usage, and mixing different types of items to fill these voids. Finally, the proposed VNS heuristic could inspire the development of efficient computational heuristics for continuous optimization in other areas such as engineering and econometric.

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