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# Iterated variable neighborhood search for the capacitated clustering problem



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#### ARTICLE INFO

Article history:
Received 6 January 2016
Received in revised form
16 June 2016
Accepted 6 August 2016
Available online 9 September 2016

Keywords: Capacitated clustering Grouping problem Variable neighborhood search Heuristics

#### ABSTRACT

The NP-hard capacitated clustering problem (CCP) is a general model with a number of relevant applications. This paper proposes a highly effective iterated variable neighborhood search (IVNS) algorithm for solving the problem. IVNS combines an extended variable neighborhood descent method and a randomized shake procedure to explore effectively the search space. The computational results obtained on three sets of 133 benchmarks reveal that the proposed algorithm competes favorably with the state-of-the-art algorithms in the literature both in terms of solution quality and computational efficiency. In particular, IVNS discovers an improved best known result (new lower bounds) for 28 out of 83 most popular instances, while matching the current best known results for the remaining 55 instances. Several essential components of the proposed algorithm are investigated to understand their impacts on the performance of algorithm.

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# 1. Introduction

Given a weighted undirect graph G = (V, E, C, w), where  $V = \{v_1, v_2, ..., v_n\}$  is the set of n nodes, E is the set of its edges,  $C = \{c_{ij}: \{v_i, v_j\} \in E\}$  represents the set of edge weights, and  $w = \{w_i \ge 0: v_i \in V\}$  is the set of node weights, the capacitated clustering problem (CCP) is to partition the node set V into a fixed number p ( $p \le n$  is given) of disjoint clusters (or groups) such that the sum of node weights of each cluster lies in a given interval [L, U] while maximizing the sum of the edge weights whose two associated endpoints locate in the same cluster. In some related literature like Deng and Bard (2011) and Martínez-Gavara et al. (2015), an edge weight  $c_{ij} \in C$  is also called the benefit of the edge  $\{v_i, v_i\}$ , while L and U are called the lower and upper capacity limits of a cluster.

Formally, the CCP can be expressed as the following quadratic program with binary variables  $X_{ig}$  taking the value of 1 if node  $v_i$  is in cluster g and 0 otherwise (Deng and Bard, 2011; Martínez-Gavara et al., 2015):

(CCP) Maximize 
$$\sum_{g=1}^{p} \sum_{i=1}^{n-1} \sum_{j=i+1}^{n} c_{ij} X_{ig} X_{jg}$$
 (1)

Subject to 
$$\sum_{g=1}^{p} X_{ig} = 1, i = 1, 2, ..., n$$
 (2)

$$L \le \sum_{i=1}^{n} w_{i} X_{ig} \le U, g = 1, 2, ..., p$$
(3)

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$$X_{ig} \in \{0, 1\}, i = 1, 2, ..., n; g = 1, 2, ..., p$$
 (4)

$$c_{ij} = 0, \forall \{v_i, v_i\} \notin E \tag{5}$$

where the set of constraints (2) guarantees that each node is located in exactly one cluster (or group) and the set of constraints (3) forces the sum of node weights of each cluster to be at least L and at most U. The set of constraints (5) ensures that the benefit between nodes  $v_i$  and  $v_j$  is 0 if  $\{v_i, v_i\} \notin E$ .

The CCP is closely related to three other clustering problems: the graph partitioning problem (GPP) (Bader et al., 2013; Benlic and Hao, 2011, 2013; Galinier et al., 2011; Soper et al., 2004), the maximally diverse grouping problem (MDGP) (Brimberg et al., 2015; Fan et al., 2010; Gallego et al., 2013; Johnes, 2015; Palubeckis et al., 2015; Rodriguez et al., 2013; Weitz and Lakshminarayan, 1997), and the handover minimization problem (HMP) (Martínez-Gavara et al., 2015; Morán-Mirabal et al., 2013). First, the GPP is a special case of the CCP when the lower and upper capacity limits of the clusters are respectively set to 0 and  $(1 + \epsilon) \left[ \frac{\sum_{i=1}^n w_i}{p} \right]$ , where  $\epsilon (\in [0, 1))$  is a predetermined imbalance parameter. As such, the CCP is also known as the node capacitated graph partitioning problem (Feo and Khellaf, 1990; Ferreira et al., 1996, 1998; Özsoy and Labbé, 2010) or the min-cut clustering problem (Johnson et al., 1993) in the literature. Second, the MDGP is also a special case of the CCP when the given graph is a complete graph and the nodes have a unit weight ( $w_i$ =1, i=1, 2, ..., n) (Martínez-Gavara et al., 2015). Additionally, as discussed in Martínez-Gavara et al. (2015) and Morán-Mirabal et al. (2013), the HMP can be viewed as a practical application of the CCP in the context of mobile networks.

Given that the CCP generalizes the NP-hard MDPG, GPP, and HMP problems, the CCP is at least as computationally difficult as these problems. Moreover, any real-world applications that can be formulated by the MDPG, GPP, or HMP models can be cast as the CCP, such as creation of peer review groups (Chen et al., 2011), parallel computing (Hendrickson and Kolda, 2000), assignment of students to groups (Johnes, 2015), VLSI design (Weitz and Lakshminarayan, 1997), etc.

Given the NP-hard nature of the CCP and its practical importance, a large number of studies have been proposed to investigate the problem and the three related clustering problems. Below, we highlight some most recent approaches on the CCP, while refereeing the reader to two recent papers (Deng and Bard, 2011; Martínez-Gavara et al., 2015) for a comprehensive review of existing studies in the literature.

In 2011, Deng and Bard (2011) proposed a greedy randomized adaptive search procedure with path relinking (GRASP+PR) by hybridizing a construction procedure of initial solution, a randomized variable neighborhood descent method as well as a path-relinking procedure. The reported computational results showed that the proposed GRASP+PR algorithm outperforms the reference algorithms. In 2013, Morán-Mirabal et al. (2013) proposed the following three heuristic algorithms for the HMP problem that is a special case of the CCP: a GRASP method (denoted by GQAP in their paper), an evolutionary path-relinking algorithm combined with GRASP (GevPR-HMP), and a population-based biased random-key genetic algorithm (BRKGA). Their study showed that GevPR-HMP achieved the best performance among the three proposed algorithms. In 2014, Lewis et al. (2014) made a comparison between the linear and nonlinear models for the CCP under the framework of exact methods, and showed that the quadratic model generally outperforms its equivalent linear alternatives.

Recently (2015), Martínez-Gavara et al. (2015) introduced several heuristic algorithms for the CCP, including a new GRASP method, a tabu search (TS) method, and a hybrid algorithm combining the proposed GRASP and tabu search methods (GRASP+TS). The authors also adapted a tabu search algorithm with strategic oscillation (TS\_SO) originally designed for the MDGP to solve the CCP. Their study showed that the proposed GRASP+TS and TS algorithms significantly outperform their reference algorithms, including their GRASP method and TS\_SO, and Deng and Bard's GRASP presented in Deng and Bard (2011), as well as the GevPR-HMP algorithm of Morán-Mirabal et al. (2013). Consequently, the TS and GRASP+TS algorithms proposed in Martínez-Gavara et al. (2015) can be considered as the current best performing approaches for the CCP.

In this paper, we are interested in solving the general CCP problem approximately and propose for this purpose an effective iterated variable neighborhood search algorithm (IVNS). The main contributions of the present work can be highlighted as follows:

- The proposed IVNS algorithm introduces an extended variable neighborhood descent (EVND) method to ensure an intensified local optimization. Contrary to the standard variable neighborhood descent (VND) method (Mladenović and Hansen, 1997), our EVND method focuses on a more balanced exploitation between different neighborhoods, which provides the search with a reinforced diversification effect. Additionally, IVNS integrates two dedicated construction procedures to generate initial solutions and a randomized shake procedure to escape deep local optima (i.e., local optimum solutions which are difficult to attain and difficult to escape for a search algorithm).
- When it is assessed on three sets of 133 benchmark instances of the literature, the proposed IVNS algorithm achieves highly competitive
  performances both in terms of the solution quality and computational efficiency compared to the state-of-the-art results. On the two
  sets of 50 standard instances, IVNS outperforms the state-of-the-art CCP algorithms in the literature. Moreover, for the 83 popular
  benchmark instances of the third set, IVNS improves the best known results (new lower bounds) in 28 cases and matches the best
  known results for the 55 remaining cases.

The rest of the paper is organized as follows. In the next section, our IVNS algorithm and its components are described in detail. Section 3 is dedicated to computational assessments based on the commonly used benchmarks and comparisons with the state-of-the-art algorithms in the literature. In Section 4, several essential components of the proposed algorithm are investigated to shed light on how they affect the performance of the proposed algorithm. Concluding comments are summarized in the last section.

## 2. Iterated variable neighborhood search for the CCP

Variable neighborhood search (VNS) (Hansen et al., 2010; Mladenović and Hansen, 1997) has been applied with success to many combinatorial optimization problems (see for instance, Armas et al., 2015; Brimberg et al., 2015; Mladenović et al., 2016; Urošević, 2014; Villegas et al., 2010). In this work, we follow the general VNS framework and propose the iterated variable neighborhood search (IVNS) method for the CCP which integrates specially designed components to reach a suitable trade-off between intensification and diversification of the search process. Specifically, the proposed IVNS algorithm employs a randomized construction procedure to generate the initial solution, a new local optimization approach called the EVND method (extended variable neighborhood descent method) to discover local optima, and a shake procedure to perturb the incumbent solution. The proposed IVNS algorithm also employs a diversification stage to produce transition states between high-quality local optimum solutions.

#### 2.1. General procedure

Our IVNS algorithm (Algorithm 1) starts from an initial feasible solution that is generated by a randomized construction procedure (Section 2.3) and is improved by the EVND method (lines 1 and 2, Section 2.4.3). Then it enters a 'while' loop in which an iterated local optimization (the inner 'while' loop, lines 5–17) and a diversification phase (the *Shake* call, line 18) are iteratively performed until a cutoff time  $t_{max}$  is reached.

```
Algorithm 1: Main framework of IVNS method for CCP
  Input: Instance I, parameter \beta_{max}, cutoff time t_{max}, \eta shake strength
   \mathbf{Output}: The best solution s^* found during the whole search process
 1 \ s \leftarrow InitialSolution(I)
                                                              /* section 2.3 */
s \leftarrow EVND(s)
                                                            /* section 2.4.3 */
s^b \leftarrow s, s^* \leftarrow s
4 while Time() \leq t_{max} do
      \beta \leftarrow 0
       /* Intensified search: iterated local optimization with
      Shake and EVND
      while \beta < \beta_{max} \land \mathit{Time}() \leq t_{max} do
          s \leftarrow Shake(s^b, \eta)
                                    /* perturb s^b before EVND improvement,
          section 2.5 */
          s \leftarrow EVND(s)
                                    /* local improvement, section 2.4.3 */
          if f(s) > f(s^*) then
10
                                /* update the best solution ever found */
11
          if f(s) > f(s^b) then
12
              s^b \leftarrow s, \beta \leftarrow 0 /* s^b denotes the best solution obtained by
13
             the current inner 'while' loop */
14
          else
             \beta \leftarrow \beta + 1
15
16
          end
17
      end
       /* Diversification: additional Shake to escape deep local
         \leftarrow Shake(s^b, \eta) /* an additional shake of deep local optimum
18
      s^b before next round of iterated local optimization st/
19 end
20 return s
```

The inner 'while' loop aims to find, from a given solution (a local optimum), an improved local optimum by iterating the Shake procedure (line 7) followed by the EVND procedure (line 8). The starting solution is first shaken by making  $\eta$  changes ( $\eta$  is called shake strength, see Section 2.5) which serves as the starting point of the extended variable neighborhood descent procedure (see Section 2.4.3). The outcome of each EVND application is used to update the best solution ever found ( $s^*$ , lines 9–11) and the best local optimum found during the current iterated local optimization phase ( $s^b$ , lines 12–16). The counter  $\beta$  indicates the number of consecutive local optimization (Shake + EVND) iterations during which  $s^b$  is not updated ( $\beta$  is reset to 0 each time an improved local optimum  $s^b$  is discovered). The inner 'while' loop exits when the cutoff time is reached (in which case the whole algorithm terminates) or when  $\beta$  attains a fixed value  $\beta_{max}$  (a large  $\beta_{max}$  thus induces a more intensified search). In the later case,  $s^b$  indicates a deep local optimum that the inner Shake call (line 7) is not sufficient to help EVND to escape. For this reason, we apply an additional Shake call (line 18) to modify  $s^b$  before giving it to the next round of the inner 'while' loop.

Note that with the second Shake call (line 18), the next inner 'while' loop starts the local optimization (EVND) with a doubly shaken starting solution, which diversifies the search strongly and helps escape deep local optima. More generally, the second Shake call may be replaced by other diversification techniques like random or customized restarts. In our case, we simply adopt the same Shake procedure used in the iterated local optimization phase. As shown in Section 3, this technique proves to be suitable and effective for the tested benchmarks. We also provide a study in Section 4.3 about the diversification effect of this second Shake application.

## 2.2. Search space, evaluation function and solution representation

For a given CCP instance that is composed of a weighted graph G = (V, E, C, w), the number p of clusters, and the lower and upper limits L and U on the capacity of clusters, the search space  $\Omega$  explored by the IVNS algorithm contains all feasible solutions, i.e., all partitions of the nodes of V into p clusters such that the weight of each cluster lies between its lower and upper limits.

Formally, let  $\pi$ :  $V \to \{1, ..., p\}$  be a partition function of the n nodes of V to p clusters. For each cluster  $g \in \{1, ..., p\}$ , let  $\pi_g = \{v \in V : \pi(v) = g\}$  (i.e.,  $\pi_g$  is the set of nodes of cluster g). Then the search space  $\Omega$  explored by our IVNS algorithm is given by:

$$\Omega = \left\{ \pi \colon \forall \ g \in \{1, \dots, p\}, L \leq |\pi_g| \leq U, |\pi_g| = \sum_{v \in \pi_g} w(v) \right\}.$$

For any candidate partition  $s = \{\pi_1, \pi_2, ..., \pi_n\}$  in  $\Omega$ , its quality is evaluated by the objective function value f(s) of the CCP:

$$f(s) = \sum_{g=1}^{p} \sum_{v_i, v_j \in \pi_g, i < j} c_{ij}$$
(6)

Given a candidate solution  $s = \{\pi_1, \pi_2, ..., \pi_p\}$ , IVNS employs a n-dimensional vector x (element coordinate vector) to indicate the cluster of each node (or element). That is, if element i belongs to cluster  $\pi_g$ , then x[i] = g ( $i \in \{1, ..., n\}$ ). IVNS additionally uses a p-dimensional vector WC (cluster weight vector) to indicate the weight of each cluster of solution s, i.e.,  $WC[g] = \sum_{v \in \pi_g} w(v)$  ( $\forall g \in \{1, ..., p\}$ ). Moreover, to facilitate neighborhood operations during the search process, the algorithm maintains a  $n \times p$  matrix  $\gamma$  in which the entry  $\gamma[v][g]$  represents the sum of edge weights between the node v and the nodes of cluster g in the candidate solution s, i.e.,  $\gamma[v][g] = \sum_{u \in \pi_g} c_{vu}$ . Consequently, any candidate solution  $s \in \Omega$  can be conveniently indicated by the s and s and s and s and its objective function value s, i.e., s and s and

# 2.3. Initial solution

The proposed IVNS algorithm needs, for each run, an initial solution to start its search. In this work, we devise two randomized construction procedures for this purpose. The first procedure (Algorithm 2) operates in two stages. In the first stage, it iteratively performs a series of insertion operations until all clusters satisfy their lower capacity constraint. Specifically, for each insertion operation, a node v and a cluster g are randomly chosen from the set AN of unassigned nodes and the set AC of clusters whose lower bound constraint is not satisfied, then the node v is assigned to cluster g. In the second stage, the construction procedure performs again a series of insertion operations until all nodes are assigned. Each insertion operation consists of randomly picking an unassigned node v and a cluster g such that  $WC[g] + w(v) \le U$ , and then assigning v to g, where WC[g] and w(v) denote respectively the current weight of cluster g and the weight of node v.

```
Algorithm 2: Initial Solution Procedure
Function InitialSolution()
Input: Instance I
Output: A feasible initial solution (denoted by \langle x[1:n], WC[1:p], \gamma, f \rangle)
\overrightarrow{AN} \leftarrow \{1, 2, \dots, n\}
                                      /* AN is the set of available nodes */
AC \leftarrow \{1, 2, \dots, p\}

AC \leftarrow \{1, 2, \dots, p\}

for g \leftarrow 1 to p do

|WC[g] \leftarrow 0
                                    /* AC is the set of available clusters */
/* WC[g] is the weight of cluster g for the current solution */
 /* x represents the coordinate vector of current solution
    v \leftarrow RandomNode(AN)
                                              /* randomly pick a node from AN */
    g \leftarrow RandomCluster(AC) \ /* \ {\tt randomly \ pick \ a \ cluster \ from \ } AC \ */
    WC[g] \leftarrow WC[g] + w[v]
AN \leftarrow AN \setminus \{v\}
if WC[g] \ge L then
AC \leftarrow AC \setminus \{g\}
    end
end
AC \leftarrow \{1, 2, \dots
while AN \neq \emptyset do |Flag \leftarrow true|
    while Flag do
         v \leftarrow RandomNode(AN) /* randomly pick a node from AN */
         g \leftarrow RandomCluster(AC) /* randomly pick a cluster from AC
        if WC[g] + w[v] \le U then |Flag \leftarrow false
         \quad \text{end} \quad
    end
    x[v] \leftarrow g
WC[g] \leftarrow WC[g] + w[v]
AN \leftarrow AN \setminus \{v\}
_{
m end}
Compute \gamma and f for x
                                                                          /* Section 2.4.2 */
return \langle x, WC, \gamma, f \rangle
```

However, the preliminary experiments showed that it was often difficult to obtain a feasible solution by the above procedure when the upper capacity limit of clusters is very tight. As a result, we modify slightly the above procedure as follows to obtain a second construction procedure. For each insertion operation, instead of randomly picking a node from the set AN of unassigned nodes, we always choose the node v in AN such that v has the largest weight (break ties at random).

Due to the random choices for insertion operations, the construction procedures are able to generate diversified initial solutions which allow the algorithm to start each run in a different area of the search space.

#### 2.4. Local optimization method

Our IVNS algorithm employs the EVND method as its local optimization procedure which extends the standard variable neighborhood descent (VND) method. The detail of the EVND method is described in the following subsections.

## 2.4.1. Neighborhood structures

Our EVND procedure exploits systematically three neighborhoods, i.e., the insertion neighborhood  $N_1$ , the swap neighborhood  $N_2$ , and the 2-1 exchange neighborhood  $N_3$ . Note that although these three neighborhoods have been proposed in previous studies (Deng and Bard, 2011; Martínez-Gavara et al., 2015), they have never been used together like we do in this work.

The insertion neighborhood  $N_1$  is based on the *OneMove* operator. Give a solution (or a partition)  $s = \{\pi_1, \pi_2, ..., \pi_p\}$  in the search space  $\Omega$ , the *OneMove* operator transfers a node v of s from its current cluster  $\pi_i$  to another cluster  $\pi_j$  such that  $|\pi_i| - w(v) \ge L$  and  $|\pi_j| + w(v) \le U$ , where L and U denote respectively the lower and upper limits of capacity of clusters, w(v) represents the weight of node v, and  $|\pi_i|$  and  $|\pi_j|$  denote respectively the weights of the clusters  $\pi_i$  and  $\pi_j$  in s. Let  $\langle v, \pi_i, \pi_j \rangle$  designate such a move and  $s \oplus \langle v, \pi_i, \pi_j \rangle$  be the resulting neighboring solution produced by applying the move to s. Then the neighborhood  $N_1$  of s is composed of all possible neighboring solutions that can be obtained by applying the *OneMove* operator to s, i.e.,

$$N_{1}(s) = \left\{ s \ \oplus < v, \ \pi_{i}, \ \pi_{i} > : \ v \in \pi_{i}, \ |\pi_{i}| - w(v) \ge L, \ |\pi_{i}| + w(v) \le U, \ i \ne j \right\}$$

Clearly, the size of  $N_1$  is bounded by  $O(n \times p)$ .

The neighborhood  $N_2$  is defined by the SwapMove operator. Given two nodes v and u which are located in two different clusters of s, the SwapMove(v, u) operator exchanges their clusters such that the resulting neighboring solution is still feasible. Thus, the neighborhood  $N_2$  of s is composed of all feasible neighboring solutions that can be obtained by applying SwapMove to s, i.e.,

```
N_2(s) = \{s \; \oplus \; SwapMove(v, \, u) \colon v \in \pi_i, \, u \in \pi_i, \, L \leq \{|\pi_i| + w(u) - w(v), \, |\pi_i| + w(v) - w(u)\} \leq U, \, i \neq j\}
```

The size of  $N_2$  is bounded by  $O(n^2)$  and is usually larger than that of  $N_1$ .

The neighborhood  $N_3$  is based on the 2-1 exchange operator (Exchange(v, u, z)). Given the current solution  $s = \{\pi_1, \pi_2, ..., \pi_p\}$  and three nodes v, u and z, where v and u are located in the same cluster  $\pi_i$  and z is located in another cluster  $\pi_j$ , Exchange(v, u, z) transfers the nodes v and u from their current cluster  $\pi_i$  to the cluster  $\pi_j$ , and transfers simultaneously the node z from the cluster  $\pi_j$  to the cluster  $\pi_i$  in such a way that the resulting solution is still feasible. For the current solution s, the neighborhood s0 of s1 is composed of all feasible neighboring solutions which can be obtained by applying the s1 exchange(s2, s3 operator to s3.

```
N_3(s) = \{s \in Exchange(v, u, z): v, u \in \pi_i, z \in \pi_i, L \le \{|\pi_i| - w(u) - w(v) + w(z), |\pi_i| + w(v) + w(u) - w(z)\} \le U, i \ne j\}
```

Since the *Exchange*(v, u, z) operator involves three nodes, the size of  $N_3$  is bounded by  $O(n^3)$  and is usually much larger than that of  $N_2$  and  $N_1$ .

Additionally, it is worth noticing that these three neighborhoods (i.e.,  $N_1$ ,  $N_2$ , and  $N_3$ ) are functionally complementary. Actually, the *OneMove*, *SwapMove*, and *Exchange*(v, u, z) operators transfer at a time 1, 2, and 3 nodes, respectively. As a result, their combined use offers more opportunities for the local search procedure to discover high-quality solutions.

# 2.4.2. Fast neighborhood evaluation technique

Similar to the previous studies for the MDGP (Brimberg et al., 2015; Lai and Hao, 2016; Palubeckis et al., 2015; Rodriguez et al., 2013; Urošević, 2014), our EVND procedure employs an incremental evaluation technique to calculate rapidly the move value (i.e., the change of objective value) of each candidate move. As mentioned in Section 2.2, our procedure maintains a  $n \times p$  matrix  $\gamma$  in which the entry  $\gamma[v][g]$  represents the sum of weights between the node  $\nu$  and the nodes of cluster g for the current solution, i.e.,  $\gamma[v][g] = \sum_{u \in \pi_g} c_{vu}$ . With the help

of matrix  $\gamma$ , the evaluation function value f can be calculated as  $f(s) = \frac{1}{2} \sum_{i=1}^{n} \gamma[i][x[i]]$  for an initial candidate solution s = (x[1], x[2], ..., x[n]). Moreover, the matrix  $\gamma$  is frequently used in the neighborhood search operations (see Algorithms 4–6).

Based on the current solution (or partition)  $s = \{\pi_1, \pi_2, ..., \pi_p\}$ , if a *OneMove* operation  $\langle v, \pi_i, \pi_j \rangle$  is performed, the move value can be easily determined as  $\Delta_f(\langle v, \pi_i, \pi_j \rangle) = \gamma[v][j] - \gamma[v][i]$ , and then the matrix  $\gamma$  is accordingly updated. More specifically, the i-th and j-th columns of matrix  $\gamma$  are updated as follows:  $\gamma[u][i] = \gamma[u][i] - c_{vu}$ ,  $\gamma[u][j] = \gamma[u][j] + c_{vu}$ ,  $\forall u \in V$ ,  $u \neq v$ , where  $c_{vu}$  is the edge weight between the nodes v and u. As such, the evaluation function value f can be rapidly updated as  $f \leftarrow f + \Delta_f$ .

When a SwapMove(v, u) operation is performed, its move value is calculated as  $\Delta_f(SwapMove(v, u)) = (\gamma[v][x[u]] - \gamma[v][x[v]]) + (\gamma[u][x[v]]) - \gamma[u][x[u]]) - 2c_{vu}$ , where x[v] and x[u] are the cluster of nodes v and u in the current solution s. As stated in Section 2.2, x = (x[1], x[2], ..., x[n]) represents the coordinate vector of the incumbent solution s. Since a SwapMove(v, u) operation is composed of two consecutively performed OneMove operations, i.e.,  $s \oplus SwapMove(v, u) = (s \oplus \langle v, x[v], x[u] \rangle) \oplus \langle u, x[u], x[v] \rangle$ , matrix  $\gamma$  is consecutively updated two times according to the OneMove move.

When a Exchange(v, u, z) move is performed, the move value is calculated as  $\Delta_f(Exchange(v, u, z)) = (\gamma[v][x[z]] - \gamma[v][x[v]]) + (\gamma[u][x[z]] - \gamma[u][x[u]]) + (\gamma[z][x[v]] - \gamma[z][x[z]]) + 2(c_{vu} - c_{vz} - c_{uz})$ . Since a Exchange(v, u, z) move is composed of three consecutively performed OneMove moves, i.e.,  $s \oplus Exchange(v, u, z) = ((s \oplus \langle v, x[v], x[z] \rangle) \oplus \langle u, x[u], x[z] \rangle) \oplus \langle z, x[z], x[v] \rangle$ , matrix  $\gamma$  is consecutively updated three times according to the OneMove move.

Matrix  $\gamma$  is initialized at the beginning of each run of the EVND procedure with a time complexity of  $O(n^2)$ , and is updated after each move operation in O(n) for any considered move operator.

#### 2.4.3. Extended variable neighborhood descent

Let  $N_k$  ( $k=1,2,...,k_{max}$ ) be a sequence of neighborhood structures (also called the neighborhood in this section) with respect to a given optimization problem, the standard variable neighborhood descent (VND) method changes in a deterministic way the current neighborhood in order to find a high-quality local optimum solution with respect to all  $k_{max}$  neighborhoods (Deng and Bard, 2011; Mladenović and Hansen, 1997). Specifically, the standard VND method starts with the first neighborhood  $N_1$  (k=1) and makes a complete exploitation of the neighborhood. Then, the VND method switches orderly to the next neighborhood  $N_{k+1}$  ( $k \leftarrow k+1$ ) when the current neighborhood  $N_k$  ( $k=1,2,...,k_{max}-1$ ) is fully explored without finding an improving solution. Moreover, the search process switches immediately to the first neighborhood  $N_1$  as soon as an improving solution is detected in the current neighborhood  $N_k$ , i.e.,  $k \leftarrow 1$ . Finally, the VND method stops if the search process reaches the last neighborhood  $N_{k_{max}}$  and no improving solution can be found in  $N_{k_{max}}$ , and the best solution found during the search process is returned as a result of the VND method. Clearly, the returned solution is a local optimum solution with respect to all  $k_{max}$  neighborhoods.

# 2.4.3 Extended Variable Neighborhood Descent

```
Algorithm 3: Extended Variable Neighborhood Descent (EVND) for CCP
\overline{\text{Function }EVND}(s_0)
Input: Solution s_0
Output: The local optimum solution s
s \leftarrow s_0
repeat
   repeat
       s \leftarrow LSN_1(s)
                                                           /* Algorithm 4 */
       (Flag, s) \leftarrow LSN_2(s)
                                                           /* Algorithm 5 */
   until Flag = false
   (Flag, s) \leftarrow LSN_3(s)
                                                            /* Algorithm 6 */
until Flag = false
return s
```

```
Algorithm 4: Local search with N_1
Function LSN_1(x, WC, \gamma, f)
Input: x[1:n], WC[1:p], \gamma, f
Output: The local optimum solution (denoted by \langle x, WC, \gamma, f \rangle)
/\ast~WC represents the weight vector of clusters
 /* x represents the coordinate vector of current solution
Improve \leftarrow \mathbf{true}
while Improve = true do
    Improve \leftarrow \mathbf{false}
    for v \leftarrow 1 to n do
        for g \leftarrow 1 to p do
            if (x[v] \neq g) \wedge (WC[x[v]] - w[v] \geq L) \wedge (WC[g] + w[v] \leq U) then
                \Delta_f \leftarrow \gamma[v][g] - \gamma[v][x[v]]
                if \Delta_f > 0 then
                    WC[x[v]] \leftarrow WC[x[v]] - w[v]
                    WC[g] \leftarrow WC[g] + w[v]
                    x[v] \leftarrow g
                     f \leftarrow f + \Delta_f
                                                                /* Section 2.4.2 */
                    Update matrix \gamma
                    Improve \leftarrow \mathbf{true}
                end
            end
       end
    end
end
return < x, WC, \gamma, f >
```

```
Algorithm 5: Local search with N_2
Function LSN_2(x, WC, \gamma, f)
Input: x[1:n], WC[1:p], \gamma, f, m
Output: The obtained solution (denoted by \langle x, WC, \gamma, f \rangle) 
/* WC represents the weight vector of clusters
  /* x represents the coordinate vector of current solution
 Improve \leftarrow \mathbf{true}, Flag \leftarrow \mathbf{false}, counter \leftarrow 0
while Improve = true do | Improve \leftarrow false
       \begin{aligned} & \textit{Improve} \leftarrow \text{taise} \\ & \text{for } v \leftarrow 1 \; \textit{to} \; n-1 \; \text{do} \\ & | \; \text{for } u \leftarrow v+1 \; \textit{to} \; n \; \text{do} \\ & | \; \text{if } \; (x[v] \neq x[u]) \wedge (L \leq WC[x[v]] + (w[u] - w[v]) \leq U) \; \wedge \\ & | \; (L \leq WC[x[u]] + (w[v] - w[u]) \leq U) \; \text{then} \\ & | \; \Delta_f \leftarrow (\gamma[v][x[u]] - \gamma[v][x[v]]) + (\gamma[u][x[v]] - \gamma[u][x[u]]) - 2c_{vu} \\ & | \; \text{if } \Delta_f > 0 \; \text{then} \\ & | \; \text{then} \; (x[x], y]) + (x[x], y] + (x[x], y] + (x[x], y]) \end{aligned} 
                                        \begin{array}{l} \Delta_f > 0 \text{ then} \\ WC[x[v]] \leftarrow WC[x[v]] + (w[u] - w[v]) \\ WC[x[u]] \leftarrow WC[x[u]] + (w[v] - w[u]) \end{array}
                                        Swap(x, v, u) 
 f \leftarrow f + \Delta_f
                                         Update matrix \gamma
                                                                                                                             /* Section 2.4.2 */
                                         Flaa \leftarrow \mathbf{ture}
                                         Improve \leftarrow \mathbf{true}
                                         counter \leftarrow counter + 1
                                         if counter \ge m then | return < Flag, x, WC, \gamma, f >
                                                          /* Return the reached solution and stop the
                                                 function */
                                         end
                                end
                       end
                end
       end
end
\mathbf{return} < Flag, x, WC, \gamma, f >
```

```
Algorithm 6: Local search with N_3
Function LSN_3(x, WC, \gamma, f)
Input: x[1:n], WC[1:p], \gamma, f, m
Output: The obtained solution (denoted by \langle x, WC, \gamma, f \rangle) /* WC represents the weight vector of clusters
 /* x represents the coordinate vector of current solution
Improve \leftarrow \mathbf{true}, Flag \leftarrow \mathbf{false}, counter \leftarrow 0
while Improve = true do | Improve \leftarrow false
     for v \leftarrow 1 to n-1 do

for u \leftarrow v+1 to n do

for z \leftarrow 1 to n do
                      \begin{array}{l} WC[x[v]] \leftarrow WC[x[v]] + (w[z] - w[u] - w[v]), \\ WC[x[z]] \leftarrow WC[x[z]] + (w[v] + w[u] - w[z]) \end{array} 
                                 swap \leftarrow x[v], \, x[v] \leftarrow x[z], \, x[u] \leftarrow x[z], \, x[z] \leftarrow swap
                                  f \leftarrow f + \Delta f
                                 Update matrix \gamma
                                                                                     /* Section 2.4.2 */
                                  Flaq \leftarrow \mathbf{ture}
                                 Improve \leftarrow \mathbf{true}
                                 counter \leftarrow counter + 1
                                 \begin{array}{c} \textbf{if} \ \ counter \geq m \ \ \textbf{then} \\ \ \ | \ \ \textbf{return} < Flag, x, WC, \gamma, f > \end{array}
                                                /* Return the reached solution and stop
                                      the function */
                                 end
                           end
                     end
                end
           end
     end
\mathbf{return} < Flag, x, WC, \gamma, f >
```

Our EVND method described in Algorithm 3 extends the standard VND method in the sense that EVND employs a different condition to switch from the current neighborhood  $N_k$  (k > 1) to the first neighborhood  $N_1$ . In the standard VND method, the search process switches back to the first neighborhood as soon as an improving solution is found in the current neighborhood  $N_k$  (even if more improving solutions can be further found in  $N_k$ ). However, our EVND method switches to the first neighborhood  $N_1$  when one of the following two conditions is

satisfied. First, the solution has been updated (or improved) m ( $m \ge 1$ , a parameter called 'the depth of improvement in neighborhood search') times with the current neighborhood  $N_k$ . Second, the solution has been updated (improved) at least one time with the neighborhood  $N_k$  and no improving solution can further be found in the neighborhood  $N_k$ . Clearly, the standard VND method is a special case of our EVND method when m = 1. Note that compared to the standard VND method, our EVND method imposes a stronger condition to move back to the first neighborhood.

Such an extension for the standard VND method is based on two considerations. First, in the standard VND method, the first neighborhood  $N_1$  is explored more often than the other neighborhoods since we move back to  $N_1$  as soon as an improving solution is discovered in the current neighborhood  $N_k$  (k > 1). However, a more balanced exploitation of all the k neighborhoods constitutes another possibility and may help the search process to discover better solutions. Compared to the standard VND method, our EVND method promotes a more balanced exploitation of the neighborhoods  $N_k$  ( $k = 2, 3, ..., k_{max}$ ) relative to the first neighborhood  $N_1$ . Second, the solutions returned by the neighborhoods  $N_k$  ( $k = 2, 3, ..., k_{max}$ ) generally have a larger distance from the local optimum solution produced most recently by the first neighborhood  $N_1$  with our EVND method than with the standard VND method. Thus, compared to the standard VND method, our EVND method creates some diversification effect during its intensified descent process with each neighborhood.

Our EVND method for the CCP exploits three complementary neighborhoods introduced in Section 2.4.1, i.e.,  $N_1$ ,  $N_2$  and  $N_3$  and is described in Algorithms 3–6, where the variables x, WC,  $\gamma$  and f have the same meanings as those in Section 2.2.

From Algorithm 3, one can observe that our EVND procedure consists of two loops and each loop stops as long as the corresponding Flag variable receives the value false. Specifically, for the inner loop the Flag variable will get the value false when no improving neighbor exists in the neighborhood  $N_2$  according to Algorithm 5. Similarly, for the outer loop the Flag variable will get the value false when no improving neighbor exists in the neighborhood  $N_3$  according to Algorithm 6. Consequently, the EVND procedure always stops when no improving neighbor exists in the neighborhoods  $N_2$  and  $N_3$  for the incumbent solution.

# 2.5. Shake procedure

When our IVNS algorithm reaches a local optimum solution, we apply a *Shake* procedure to the reached solution to jump out of the local optimum trap. The *Shake* procedure used by the IVNS algorithm consists of consecutively performing  $\eta$  randomly selected feasible *OneMove* or *SwapMove* moves, where  $\eta$  is a parameter called the shake strength. In other words, from the incumbent solution  $s_0$ , we construct the  $N_1(s)$  and  $N_2(s)$  neighborhoods which include all (feasible) neighboring solutions of s (see Section 2.4.1) and then pick randomly a solution s' from the union of  $N_1(s)$  and  $N_2(s)$  to replace  $s_0$ . We repeat this operation  $\eta$  times. It is clear that a large (respect. small)  $\eta$  value leads to a shaken solution which will be more (respect. less) distant from the input solution. In this work, the value of  $\eta$  is empirically set as  $\eta = \min\{15, \max\{5, 0.02n\}\}$ , where n is the number of nodes in the graph. Note that it is possible to include *Exchange* moves for the Shake operations. Meanwhile, for the reason of simplicity, we only apply *OneMove* and *SwapMove* moves, which proves to be sufficient for our purpose of diversification. The pseudo-code of our *Shake* procedure is given in Algorithm 7.

# Algorithm 7. Shake procedure.

```
1 Function Shake (s_0, \eta) Input: Solution s_0, strength of shake \eta Output: The perturbed solution s, matrix \gamma, objective value f 2 s \leftarrow s_0 3 for l \leftarrow 1 to \eta do 4 | Randomly pick a neighboring solution s' \in N_1(s) \cup N_2(s) 5 | s \leftarrow s' 6 end 7 Compute \gamma and f for s /*Section 2.4.2 */8 return (s, \gamma, f)
```

#### 3. Experimental results and comparisons

In this section, we assess the performance of the proposed IVNS algorithm by showing computational results on well-known benchmark instances and by making a comparison with the state-of-the-art algorithms in the literature.

#### 3.1. Benchmark instances

Our IVNS algorithm was assessed on three sets of 133 benchmark instances commonly used in the literature. These instances are available at http://www.optsicom.es/ccp/, and their details are described as follows.

- RanReal Set (40 instances): This set was originally proposed in Gallego et al. (2013) for the MDGP and adapted to the CCP in Martínez-Gavara et al. (2015) by generating the node weights with a uniform distribution U(0,10). This set is composed of 20 instances with n=240, p=12, L=75, and U=125, and 20 instances with n=480, p=20, L=100, and U=150. For all instances of this set, the edge weights  $c_{ij}$  are a real number which is uniformly and randomly generated in (0, 100).
- DB Set (10 instances): This set was originally proposed by Deng and Bard (2011) for the MDGP in the context of mail delivery, and adapted to the CCP in Martínez-Gavara et al. (2015) by generating the node weights with a uniform distribution U(0,10). These 10 instances are characterized by the following features: n=82, p=8, L=25, and U=75.
- MM Set (83 instances): These 83 synthetic instances were proposed by Morán-Mirabal et al. (2013) for the handover minimization

problem and were widely used in the literature. These instances have the following characteristics:  $n \in \{20, 30, 40, 100, 200, 400\}$ ,  $p \in \{5, 10, 15, 25, 50\}$ , the edge weights  $c_{ij}$  are a real number, and the lower and upper capacity limits of clusters respectively are 0 and a real number depending on each instance.

# 3.2. Parameter settings and experimental protocol

In this section, we show some basic information about our experiments, including the parameter settings of our algorithm, the reference algorithms, the experimental platform, and the termination criterion of algorithms.

First, Table 1 shows the parameter setting of our IVNS algorithm which was achieved by a preliminary experiment. For this preliminary experiment, we used 20 *RanReal* instances with n=240 which were also used in the sensibility analysis of parameters presented in Section 4.4. The computational results indicated that for m and  $\beta_{max}$  the default settings shown in Table 1 are suitable for the algorithm (see Section 4.4). For  $\eta$ , it involves three variables, so we manually tuned its value based on a principle that the strength of shake procedure should be proportional to the size of instance but in an appropriate interval. The computational results on the preliminary experiment indicated that the default setting of  $\eta$  in Table 1 is able to reach an acceptable performance of the algorithm.

Second, according to the previous surveys (Martínez-Gavara et al., 2015; Morán-Mirabal et al., 2013), the TS (Martínez-Gavara et al., 2015), GRASP+TS (Martínez-Gavara et al., 2015), and GevPR-HMP (Morán-Mirabal et al., 2013) algorithms can be considered as the state-of-the-art algorithms for the CCP. Hence, in the present study we use them as the main reference algorithms for our comparative study.

Our IVNS algorithm was programmed in C++. To make a fair comparison with the state-of-the-art algorithms, we also implemented faithfully the GRASP, TS, GRASP+TS algorithms of Martínez-Gavara et al. (2015) which are three state-of-the-art algorithms in the literature. For the GRASP, TS, GRASP+TS algorithms, we adopted the best parameter settings identified in the original paper (Martínez-Gavara et al., 2015). Moreover, all source codes were compiled using g++ compiler with the '-03' flag, and the corresponding experiments were carried out on a computing platform with an Intel E5-2670 processor (2.80 GHz CPU and 2 Gb RAM), running Linux. Following the DIMACS machine benchmark procedure, our machine requires respectively 0.19, 1.17, and 4.54 s for the graphs r300.5, r400.5, r500.5.

For the GRASP, TS, GRASP+TS, and IVNS algorithms, we used a cutoff time  $t_{max} = n$  (in seconds) as the unique stopping criterion where n is the number of nodes of the input graph.

Finally, for the IVNS algorithm, the initial solution was generated by the second initialization procedure for the handover minimization instances due to their tight upper bounds on the capacity of clusters, and by Algorithm 2 for the remaining instances. For GRASP, TS and GRASP+TS, the handover minimization instances are not used in the experiments, since their initial solution procedures cannot guarantee to generate a feasible solution for a part of them due to the tight upper bounds on the capacity of clusters.

#### 3.3. Computational results and comparison on the general CCP instances

The first experiment aims to assess the performance of our IVNS algorithm on the first two sets of instances by comparing its results with those of the state-of-the-art algorithms in the literature. In this experiment, all the compared algorithms (GRASP, TS, GRASP+TS and IVNS) were respectively performed 20 times on each instance, based on the experimental protocol of Section 3.2. The computational results are summarized in Tables 2.3.

In Table 2, the first column identifies the instances, columns 2–4 show respectively the best objective values ( $f_{best}$ ) obtained by the three reference algorithms (GRASP, TS, GRASP+TS), and column 5 reports the best objective values of our IVNS algorithm. Columns 6–9 show respectively the average objective values for the four compared algorithms ( $f_{avg}$ ). The best results among the algorithms in terms of the best and average objective values are indicated in bold. In Table 3, columns 2–5 show the standard deviation ( $\sigma$ ) of the objective values obtained over 20 runs for the compared algorithms, respectively, and columns 6–9 give the average running times (in seconds) of the algorithms to reach their respective objective values ( $time_{avg}$ ). The row #Best of the tables indicates the number of instances for which the corresponding algorithm produces the best results among the compared algorithms. In addition, to verify whether there exists a significant difference between the reference algorithms and our IVNS algorithm on the best and average objective values, as well as the standard deviation of objective values, the p-values from the non-parametric Friedman test are reported in the last row of the tables. Notice that a p-value smaller than 0.05 means that there exists a significant difference between two sets of results compared.

One observes from Tables 2 and 3 that the proposed IVNS algorithm outperforms the reference algorithms. First, IVNS obtained the best result on all 50 instances in terms of the best objective value, whereas the GRASP, TS, GRASP+TS algorithms produced respectively the best result on 10, 7 and 10 instances. Second, when comparing the average objective values, it can be found that the IVNS algorithm yielded the best result on all instances, whereas the GRASP, TS, and GRASP+TS algorithms respectively obtained the best results on only 2, 0, 4 instances. In addition, the small p-values ( < 0.05) confirm the significant differences between the results of IVNS and those of the compared reference algorithms.

Finally, compared to the reference algorithms, the IVNS algorithm produced the smallest standard deviation ( $\sigma$ ) on all tested instances, indicating that IVNS is the most robust algorithm among the compared algorithms, which is also confirmed by the associated small *p*-values.

# 3.4. Computational results and comparison on the handover minimization instances

The second experiment aims to assess the performance of the IVNS algorithm on the set of 83 handover minimization instances with  $n \le 400$ , where for each instance the IVNS algorithm was independently run 20 times. The computational results are summarized in Table 4 for the large instances with  $n \ge 100$ . For very small instances with  $n \le 40$ , the computational results are reported in the appendix

<sup>&</sup>lt;sup>1</sup> The source codes of these algorithms will be available at: http://www.info.univ-angers.fr/pub/hao/ccp.html

<sup>&</sup>lt;sup>2</sup> dmclique, ftp://dimacs.rutgers.edu/pub/dsj/clique

(Table 10) since they are very easy to be solved by the IVNS algorithm (see the appendix for the details). Notice that in the present section all results are given in the form of minimization to make a direct comparison between the results of the IVNS algorithm and those reported in the literature, and that the results of the maximization form can be converted to the minimization form as follows:

**Table 1** Settings of parameters.

Parameters	Section	Description	Values
$\beta_{max}$	2.1	Strength of intensification search	30
m	2.4.3	Depth of improvement in neighborhood search	10
η	2.5	Strength of shake	$\min\{15, \max\{5, 0.02n\}\}$

Table 2
Comparison between the IVNS algorithm and three state-of-the-art algorithms from the literature (i.e., GRASP, TS, GRASP+TS in Mladenović et al., 2016) on the first two sets (RanReal and DB) of CCP instances in terms of the best and average objective function values over 20 independent runs. The best results among the compared algorithms are indicated in bold.

	$f_{best}$				$f_{avg}$			
Instance	GRASP	TS	GRASP+TS	IVNS	GRASP	TS	GRSP+TS	IVNS
Sparse82_01	1342.17	1336.82	1342.17	1342.17	1342.13	1315.21	1342.17	1342.17
Sparse82_02	1306.64	1303.17	1306.64	1306.64	1305.65	1281.57	1306.16	1306.64
Sparse82_03	1353.94	1353.94	1353.94	1353.94	1351.69	1335.89	1353.00	1353.94
Sparse82_04	1291.22	1291.22	1291.22	1291.22	1289.15	1276.85	1290.05	1291.22
Sparse82_05	1352.35	1352.35	1352.35	1352.35	1352.35	1328.15	1352.35	1352.35
Sparse82_06	1354.61	1354.61	1354.61	1354.61	1353.86	1329.86	1354.61	1354.61
Sparse82_07	1266.94	1266.94	1266.94	1266.94	1266.86	1227.01	1266.89	1266.94
Sparse82_08	1393.02	1391.53	1393.02	1393.02	1393.02	1362.54	1393.02	1393.02
Sparse82_09	1294.12	1294.12	1294.12	1294.12	1293.69	1280.97	1293.46	1294.12
Sparse82_10	1356.98	1356.98	1356.98	1356.98	1356.85	1330.79	1356.95	1356.98
RanReal240_01	192,320.30	222,871.98	223,272.87	224,893.92	191,140.45	221,547.29	222,268.07	224,785.27
RanReal240_02	185,612.48	202,356.36	202,344.44	204,608.66	183,435.92	200,582.86	200,356,15	204,415.88
RanReal240_03	179,316.65	196,422.35	196,143.12	198,885.19	176,821.81	194,348.91	194,783.74	198,626.93
RanReal240_04	197,342.26	222,298.86	223,076.30	225,627.16	194,629.26	220,521.18	221,165.75	225,227.11
RanReal240_05	175,967.25	193,358.53	194,115.62	195,440.94	174,465.68	191,234.34	192,076.34	195,228.86
RanReal240_06	192,789.55	214,840.96	215,004.12	216,736.00	188,264.62	212,626.52	213,259.64	216,474.84
RanReal240_07	191,714.87	208,223.05	208,045.67	209,273.70	190,379.19	205,808.95	206,092.48	209,004.05
RanReal240_08	185,930.72	203,595.13	203,168.62	205,246.82	181,699.18	201,102.55	201,519.15	204,958.19
RanReal240_09	189,573.48	207,711.19	207,984.26	209,059.28	186,992.97	206,540.74	206,788.13	208,789.79
RanReal240_10	176,327.61	189,597.87	190,532.75	192,977.28	174,638.77	187,534.34	188,379.13	192,788.59
RanReal240_11	184,198.31	203,109.91	203,037.25	204,722.75	182,673.54	201,113.86	201,701.83	204,523.95
RanReal240_12	181,337.55	199.710.82	199,708.68	201,052.53	180,048.84	198.365.85	198,389.30	200,904.16
RanReal240_13	180,865.29	201,238.18	200,742.90	202,335.99	179,893.27	199,590.02	198,727.63	202,139.55
RanReal240_14	194,009.94	226,813.94	226,621.92	228,844.44	191,947.27	225,362.61	225,721.97	228,512.11
RanReal240_15	173,114.22	188,896.07	188,318.39	191,170.98	171,311.83	187,410.50	187,299.92	190,914.31
RanReal240_15	182,348.50	202,475.44	202,463.27	203,999.02	180,823.13	199,999.63	200,722.29	203,834.68
RanReal240_17	181,270.70	194,155.13	193,835.64	195,242.31	179,607.16	191,978.36	191,908.95	195,114.49
RanReal240_17	174,650.37	192,772.21	193,004.84	195,069.62	173,876.39	190,428.51	190,979.42	194,853.70
RanReal240_19	179,859.43	196,739.04	196,717.17	199,200.03	177,521.56	194,916.62	195,053.62	199,019.23
RanReal240_20	191,936.64	210,399.94	210,365.38	212,264.10	190,110.07	208,970.68	208,887.59	212,046.92
RanReal480_01	463,538.35	543,259.55	544,301.19	555,057.10	460,340.75	539,074.11	539,550.62	554,331.89
RanReal480_01	446,829.64	500,646.59	502,039.41	510,418.44	443,573.33	495,273.27	497,700.18	509,519.84
RanReal480_03	434,854.27	486,379.91	487,561.06	496,641.22	433,059.42	482,624.38	483,908.86	495,847.80
RanReal480_03	455,470.88	510,971.67	513,425.37	521,984.68	450,861.04	504,054.37	507,123.99	520,891.75
RanReal480_05	415,295.30	474,548.74	473,732.77	483,228.99	413,263.81	466,932.68	469,914.28	482,595.19
RanReal480_05	461,624.20	524,191.64	524,520.14	533,762.18	458,049.90	519,002.55	520,607.74	532,888.64
RanReal480_07	461,236.11	537,464.01	537,674.45	545,157.68	456,319.85	530,513.86	532,199.71	544,530.14
RanReal480_07		521,894.07	523,602.31	532,308.09	458,299.68	515,438.23	518,685.30	531,417.94
RanReal480_08	460,756.67 466,977.73	546,057.32	546,394.74	556,478.39		540,459.20	,	555,098.72
_	447,088.04	508,294.13		,	462,121.78		541,615.49	
RanReal480_10			508,168.74	519,456.96	441,810.74	503,161.12	504,750.89	518,612.02
RanReal480_11	451,321.35	515,296.61	515,189.66	523,450.04	448,295.75	510,410.05	509,657.31	522,814.96
RanReal480_12	434,343.91	492,469.23	493,845.25	501,596.63	433,067.26	487,442.12	488,699.18	500,580.84
RanReal480_13	467,130.54	526,936.22	524,825.92	534,638.19	461,557.62	519,201.37	521,740.67	533,763.20
RanReal480_14	428,544.75	500,100.04	508,349.48	513,777.84	426,455.75	495,859.32	499,543.12	512,975.73
RanReal480_15	446,764.19	509,377.07	509,005.17	516,941.11	443,810.07	503,159.99	503,946.63	516,017.98
RanReal480_16	465,499.89	540,493.75	540,840.67	549,371.23	462,345.11	533,502.50	535,129.37	548,276.15
RanReal480_17	460,122.80	531,353.71	529,388.54	537,483.76	458,004.02	523,459.84	524,362.15	536,655.06
RanReal480_18	456,573.73	515,692.20	518,675.12	525,813.39	452,767.79	511,193.51	512,917.14	524,650.86
RanReal480_19	454,922.45	514,503.74	512,339.77	522,158.86	449,744.91	506,858.76	508,311.67	521,180.84
RanReal480_20	443,851.41	510,045.22	509,167.33	518,288.03	440,398.01	503,995.42	505,190.58	517,261.92
# Best	10	7	10	50	2	0	4	50
p-value	2.54e – 10	5.47e – 11	2.54e – 10		4.26e – 12	1.54e – 12	1.18e – 11	

Table 3

Comparison between the IVNS algorithm and three state-of-the-art algorithms from the literature (i.e., GRASP, TS, GRASP+TS Mladenović et al., 2016) on the first two sets (RanReal and DB) of CCP instances in terms of the standard deviation and the average running time to reach its final objective value. Each instance was independently solved 20 times by each algorithm respectively.

	Standard devia	ation $(\sigma)$			$Time_{avg}(s)$			
Instance	GRASP	TS	GRASP+TS	IVNS	GRASP	TS	GRASP+TS	IVNS
Sparse82_01	0.16	23.24	0.00	0.00	14.01	9.31	15.15	0.13
Sparse82_02	0.93	19.99	0.97	0.00	30.82	7.53	29.27	0.82
Sparse82_03	1.24	14.90	1.63	0.00	41.60	2.38	38.78	0.21
Sparse82_04	1.32	12.97	1.04	0.00	30.60	6.36	36.99	4.27
Sparse82_05	0.00	30.33	0.00	0.00	4.97	8.02	5.11	0.07
Sparse82_06	1.39	25.63	0.00	0.00	24.43	4.03	26.04	0.08
Sparse82_07	0.18	22.46	0.10	0.00	32.74	4.75	28.10	0.39
Sparse82_08	0.00	29.24	0.00	0.00	1.00	9.31	0.74	0.03
Sparse82_09	0.40	13.01	0.22	0.00	28.54	5.67	16.31	0.58
Sparse82_10	0.16	24.48	0.07	0.00	29.41	2.08	37.81	0.59
RanReal240_01	864.05	1098.76	601.21	97.88	166.10	15.24	132.87	147.28
RanReal240_02	936.74	1140.64	1886.92	102.77	128.76	17.81	132.97	160.2
RanReal240_03	786.12	1192.60	875.04	170.83	123.40	7.19	125.43	146.88
RanReal240_04	1018.21	1302.30	1060.61	237.09	122.63	13.49	146.27	162.63
RanReal240_05	646.56	1255.73	894.36	91.62	113.10	8.38	129.66	161.77
RanReal240_06	1321.33	1251.75	959.32	169.61	128.25	15.04	139.80	165.88
RanReal240_07	592.59	1696.59	1324.09	120.82	114.48	27.88	143.93	120.4
RanReal240_07	1271.82	1184.11	1096.61	139.71	137.49	19.52	128.97	174.29
_	1271.82	1073,39	873.02	148.77	135.58	27.24	158.97	160.17
RanReal240_09								
RanReal240_10	659.31	1160.27	1080.52	154.18	107.53	7.40	127.40	170.78
RanReal240_11	855.28	1127.48	1041.35	96.57	115.17	36.41	162.40	148.5
RanReal240_12	624.13	1347.08	961.73	126.79	128.65	34.82	151.19	169.3
RanReal240_13	531.24	1095.26	1191.67	155.41	102.67	38.44	131.83	150.7
RanReal240_14	1185.33	1086.76	672.46	170.71	112.17	10.11	124.59	141.2
RanReal240_15	780.67	1026.16	700.91	174.08	120.50	8.78	126.35	144.9
RanReal240_16	645.97	1341.25	1033.99	131.32	97.79	19.38	141.58	147.80
RanReal240_17	679.79	1280.03	936.77	109.08	120.85	25.01	142.44	146.0
RanReal240_18	458.00	1155.66	1050.83	126.17	141.20	7.55	130.71	163.4
RanReal240_19	924.99	1173.30	1357.67	109.79	113.09	8.71	125.27	176.3
RanReal240_20	829.60	908.52	816.33	130.92	99.00	24.49	146.64	149.6
RanReal480_01	1767.54	2915.75	2305.94	415.49	276.78	87.81	270.05	369.2
RanReal480_02	1541.69	3027.29	1982.60	569.71	212.88	50.36	276.25	416.40
RanReal480_03	909.49	2529.01	1999.88	426.86	281.57	55.11	274.68	415.8
RanReal480_04	1426.35	4323.08	2867.38	648.97	230.07	53.23	275.02	380.2
RanReal480_05	1094.52	3383.90	1886.34	497.25	310.87	33.56	271.51	338.9
RanReal480_06	1573.22	3178.92	2393.40	555.89	205.83	58.20	291.32	338.2
RanReal480_07	1926.89	2864.48	2451.06	413.13	198.46	63.44	291.77	388.7
RanReal480_08	1404.82	4291.62	2058.61	555.42	194.59	77.79	291.08	373.4
RanReal480_09	1994.95	3774.05	2163.84	514.64	231.09	78.00	282.38	388.0
RanReal480_10	1667.97	2799.90	2715.16	589.69	296.32	49.75	268.99	403.4
RanReal480_11	1488.69	2771.21	2220.12	402.37	243.62	98.13	277.81	382.5
RanReal480_12	1078.08	3170.40	2713.09	540.16	292.27	80.68	296.84	386.0
RanReal480_13	2379.98	3611.97	2431.74	423.89	264.67	66.61	289.75	420.5
RanReal480_14	1025.07	2482.93	2984.86	408.22	275.91	41.00	276.64	401.0
_		3001.21	3149.15	408.22	2/5.91 241.74	41.00 86.60		
RanReal480_15	1358.30						296.67	375.9
RanReal480_16	1331.41	4488.81	2636.34	590.34	217.53	56.34	279.50	366.1
RanReal480_17	1306.58	3384.98	2538.33	402.38	274.19	56.09	286.52	389.5
RanReal480_18	1343.73	4530.39	2388.17	494.86	248.92	80.52	307.61	379.0
RanReal480_19	2115.36	2949.64	2357.29	550.77	248.76	49.19	282.87	405.6
RanReal480_20	1146.60	3505.30	2635.41	530.58	255.89	91.93	288.15	398.1
# Best	2	0	4	50				
p-value	4.26e - 12	1.54e - 12	1.18e – 11					

 $f_{min} = 2(\sum_{i < j} c_{ij} - f_{max})$ , where  $f_{min}$  and  $f_{max}$  respectively correspond to the results of minimization and maximization forms.

Columns 1 and 2 of Table 4 respectively give the instance name and the best known solution (BKS) published in the literature. Columns 3–5 show the best results of three reference algorithms in Morán-Mirabal et al. (2013): a GRASP method (GQAP), a GRASP embedded within a population-based evolutionary path-relinking algorithm (GevPR-HMP), and a population-based biased random-key genetic algorithm (BRKGA). The results of these reference algorithms were directly extracted from Morán-Mirabal et al. (2013), which correspond to the best outcomes ( $f_{best}$ ) yielded by 5 runs with a cutoff time of 24 h based on a cluster running Intel X5650 processors at 2.67 GHz or a cluster running Intel Xeon E5530 processors at 2.4 GHz (Morán-Mirabal et al., 2013). It is worth noting that the cutoff time of the three reference algorithms is much higher than ours (24 h vs.  $n \le 400 \text{ s}$ ). Columns 6–10 show the results of our IVNS algorithm, including the best objective value ( $f_{best}$ ) over 20 runs, the average objective value ( $f_{agv}$ ), the worst objective value ( $f_{worst}$ ), the standard deviation of objective value ( $\sigma$ ), and the average running time in seconds to reach its final objective value ( $f_{time_{avg}}$ ). The rows Improve, Match denote the

**Table 4**Comparison between the IVNS algorithm and the three reference algorithms in Morán-Mirabal et al. (2013) on the set of handover minimization instances. Each instance was independently solved 20 times by the IVNS algorithm, and the current best results are indicated in bold. The results are given in the form of minimization to make a direct comparison with the results from the literature.

		GevPR-HMP	GQAP	BRKGA	IVNS				
Instance	BKS	$f_{best}$	$f_{best}$	$f_{best}$	$f_{best}$	$f_{avg}$	$f_{worst}$	σ	Time <sub>avg</sub> (s)
100_15_270001	19,000	19,174	19,000	19,000	19,000	19,000.00	19,000	0.00	1.01
100_15_270002	22,686	22,686	22,686	23,288	22,686	22,686.00	22,686	0.00	0.71
100_15_270003	14,558	14,558	14,558	14,616	14,558	14,558.00	14,558	0.00	0.09
100_15_270004	19,700	19,762	19,700	19,882	19,700	19,700.00	19,700	0.00	0.15
100_15_270005	22,746	22,892	22,746	23,092	22,746	22,746.00	22,746	0.00	0.67
100_25_270001	36,412	36,412	36,448	36,752	36,412	36,412.00	36,412	0.00	2.42
100_25_270002	38,608	39,144	38,608	39,256	38,608	38,608.00	38,608	0.00	1.04
100_25_270003	32,686	32,966	32,686	32,708	32,686	32,686.00	32,686	0.00	1.82
100_25_270004	35,322	35,678	35,322	35,954	35,322	35,322.00	35,322	0.00	0.26
100_25_270005	36,878	36,906	36,878	37,100	36,690	36,690.00	36,690	0.00	0.21
100_50_270001	60,922	60,922	61,172	61,554	60,922	60,922.00	60,922	0.00	2.58
100_50_270002	62,022	62,046	62,022	62,524	62,022	62,022.00	62,022	0.00	0.53
100_50_270003	54,596	54,618	54,596	55,192	54,596	54,596.00	54,596	0.00	4.07
100_50_270004	57,894	57,894	57,894	58,208	57,894	57,894.00	57,894	0.00	1.18
100_50_270005	61,088	61,088	61,318	62,784	61,080	61,082.80	61,090	4.31	46.02
200_15_270001	81,558	81,558	82,834	81,558	81,558	81,558.00	81,558	0.00	12.53
200_15_270002	89,810	89,810	90,620	90,506	89,492	90,502.80	91,172	546.53	48.73
200_15_270003	79,232	79,232	80,980	79,548	79,232	79,277.60	80,144	198.77	13.39
200_15_270004	78,324	78,324	80,538	80,026	78,324	78,485.50	79,726	375.08	52.95
200_15_270005	95,998	95,998	98,826	98,830	95,680	96,,137.10	96,986	622.92	21.85
200_25_270001	133,168	133,168	138,454	140,492	133,168	133,168.00	133,168	0.00	51.17
200_25_270002	136,038	136,038	140,066	140,690	133,778	133,859.80	133,926	47.19	68.02
200_25_270003	139,438	139,438	144,120	143,724	136,782	136,795.50	136,812	14.92	67.84
200_25_270004	128,554	128,554	134,054	131,786	128,246	128,246.00	128,246	0.00	53.92
200_25_270005	148,402	148,402	154,260	152,934	147,844	147,844.00	147,844	0.00	10.20
200_50_270001	219,672	221,550	223,096	223,098	215,388	215,531.20	215,572	64.28	67.70
200_50_270002	216,444	218,254	219,910	219,834	212,798	212,864.60	212,912	36.18	92.01
200_50_270002	221,348	221,500	222,404	221,110	214,364	214,413.90	214,426	15.03	61.89
200_50_270004	211,832	212,044	212,544	213,170	206,476	206,509.80	206,590	29.01	61.66
200_50_270005	231,890	231,890	236,136	237,156	229,918	230,050.70	230,082	22.82	76.55
400_15_270001	370,314	372,694	456,158	375,650	369,048	372,055.40	385,786	4442.20	187.66
400_15_270001	370,274	370,274	460,232	383,096	365,878	369,275.90	378,508	4126.82	228.23
400_15_270002	358,684	358,684	448,830	366,314	352,588	356,988.80	365,886	4140.92	135.41
400_15_270003	334,430	334,430	406,834	346,282	331,888	339,169.90	350,388	6361.04	214.63
400_15_270004	361,904	361,904	457,274	377,094	360,422	363,890.10	383,154	4970.80	200.30
400_15_270003	568,830	570,852	663,908	579,130	545,118	546,936.70	549,318	1026.66	208.04
400_25_270001	543,182	544,568	658,440	554,840	528,470	529,087.80	530,118	444.68	201.85
400_25_270002	548,000	548,000	667,982	553,162	524,678	526,220.60	530,438	1695.07	215.17
400_25_270003	501,750	501,750	607,672	516,416	481,568	482,070.00	484,566	722.65	174.58
400_25_270004	556,044	556,044	679,848	585,070	548,100	549,839.10	557,482	2608.26	240.02
400_23_270003	851,412	851,412	951,882	879,438	824,766	825,581.20	826,202	402.47	240.02
400_50_270001	845,496	845,496	949,562	874,226	824,766 823,094	824,239.00	825,442	591.82	216.54
400_50_270002	819,242	819,242	919,140	843,242	823,09 <del>4</del> 801,586	802,672.90	804,310	776.71	266.22
400_50_270003	774,564	774,564	878,912	806,690	760,602	761,370.40	763,076	553.34	275.49
400_50_270004	774,564 854,726	854,726	940,358	882,060	760,602 828,384	829,335.50	830,116	421.66	275.49
	854,726 0	854,726 0	940,358	882,060 0	<b>828,384</b> 28	023,333.30	011,000	421.00	230.40
# Improve	0 17	9	0 11	0 2	28 17				
# Match					17 45				
# Total	45	45	45 5 510 0	45 5 470 11	45				
p-value	1.21e-7	1.97e-9	5.51e-9	5.47e-11					

number of instances for which the associated algorithm improved or matched the best known results in the literature, and row *Total* shows the total number of instances. Note that the current best results are indicated in bold, and other symbols are the same as those in Table 2. Besides, it should be mentioned that this section focuses on the best results produced by the compared algorithms, since the compared algorithms were run on different computers and the cut-off times of the reference algorithms are much longer than that of our IVNS algorithm.

Table 4 clearly discloses that the proposed IVNS algorithm outperforms the three reference algorithms designed for the handover minimization problem. First, the IVNS algorithm improved the best known results for 28 out of 45 instances with  $n \ge 100$ , while matching the best known results for the remaining instances. Second, compared to any of the three reference algorithms, our IVNS algorithm obtained the better or equal objective values for all instances, even if IVNS uses much shorter cutoff times than that of the reference algorithms ( $n \le 400$  seconds vs. 24 h). Third, even the worst objective value produced by the IVNS algorithm is better than the best known result reported in the literature for instances with n=400, and the average computing time  $time_{avg}$  is smaller than 300 seconds for each instance. Finally, one observes that all p-values are smaller than 0.05, implying that there exits a significant difference between the results of the IVNS algorithm and those yielded by

**Table 5**Comparison between the standard VND method and the extended VND (EVND) method on the set of 40 representative instances. Each instance was independently solved 100 times by both algorithms respectively, and better results in the average objective value ( $f_{avg}$ ) between the compared algorithms are indicated in bold.

	$f_{avg}$		time <sub>avg</sub>		
Instance	VND	EVND	VND	EVND	
RanReal240_01	220,221.45	221,035.18	0.12	0.12	
RanReal240_02	199,165.17	199,461.37	0.09	0.10	
RanReal240_03	193,878.39	194,087.07	0.09	0.10	
RanReal240_04	219,242.90	220,280.46	0.08	0.11	
RanReal240_05	190,443.88	190,569.72	0.08	0.10	
RanReal240_06	211,538.07	212,198.32	0.09	0.10	
RanReal240_07	203,850.69	204,429.68	0.10	0.13	
RanReal240_08	200,600.21	200,710.54	0.13	0.12	
RanReal240_09	204,291.07	204,961.02	0.08	0.10	
RanReal240_10	186,995.53	187,202.99	0.09	0.11	
RanReal240_11	199,062.90	199,574.51	0.09	0.11	
RanReal240_12	196,535.07	196,618.48	0.08	0.11	
RanReal240_13	197,326.13	197,585.12	0.10	0.11	
RanReal240_14	224,438.53	224,784.93	0.12	0.12	
RanReal240_15	185,489.25	186,227.47	0.08	0.10	
RanReal240_16	198,794.30	199,277.26	0.10	0.11	
RanReal240_17	189,651.07	190,188.43	0.07	0.10	
RanReal240_18	189,290.28	189,691.65	0.10	0.10	
RanReal240_19	193,267.47	194,274.68	0.08	0.10	
RanReal240_19	207,193.59	207,692.77	0.08	0.10	
RanReal480_01	541,860.57	545,446.40	0.08	0.10	
_	•	,			
RanReal480_02	497,118.68	<b>498,148.99</b>	1.01 0.99	0.95	
RanReal480_03	483,285.49	482,184.26		0.99	
RanReal480_04	507,951.82	509,674.81	0.76	0.90	
RanReal480_05	469,493.30	469,172.22	0.86	0.91	
RanReal480_06	516,732.69	518,796.77	0.96	0.95	
RanReal480_07	528,136.56	533,541.23	0.99	0.93	
RanReal480_08	516,413.17	518,435.89	0.81	0.83	
RanReal480_09	543,150.66	546,057.68	0.88	0.98	
RanReal480_10	507,686.31	508,665.48	0.77	0.86	
RanReal480_11	511,503.09	512,682.60	0.96	0.94	
RanReal480_12	487,411.85	488,100.45	1.04	1.08	
RanReal480_13	517,853.71	521,632.82	0.87	0.88	
RanReal480_14	500,399.35	500,139.40	1.01	0.94	
RanReal480_15	501,547.03	503,165.49	0.88	0.83	
RanReal480_16	536,611.41	537,921.48	1.00	0.86	
RanReal480_17	526,315.11	526,884.43	0.92	0.95	
RanReal480_18	508,572.09	511,441.64	0.88	0.85	
RanReal480_19	509,748.10	509,208.68	0.70	0.84	
RanReal480_20	502,534.92	504,453.71	0.93	0.87	
# Better	4	36			
# Equal	0	0			
# Worse	36	4			
p-value		4.20e – 7		1.96e	

the reference algorithms. In summary, these outcomes indicate that the proposed IVNS algorithm is highly efficient for solving the handover minimization instances compared to the state-of-the-art algorithms in the literature (Morán-Mirabal et al., 2013).

# 4. Analysis and discussions

We now turn our attention to analyze some essential aspects of the proposed IVNS algorithm, including the local optimization procedure (i.e., the EVND method), the influence of the diversification stage on the performance of IVNS algorithm, and a sensitivity analysis of the key parameters. In this section, all experiments were carried out based on the set of *RanReal* instances (20 instances with n=240 and 20 instances with n=480).

# 4.1. Comparison between the standard and extended VND methods

The IVNS algorithm employs the extended VND method (EVND) as its local optimization procedure. Since the EVND method is an extension of the standard VND method, we carried out an experiment to compare both methods. In this experiment, both EVND and VND were respectively run 100 times on each instance. Specifically, for each run, both methods were performed with the same initial solution

**Table 6**Comparison between the IVNS method and its variant (IVNS $^-$ ) in which the neighborhood  $N_3$  is disabled on the set of 40 representative instances. Each instance is respectively solved 20 times by both algorithms, and better results in the average objective value between two algorithms are indicated in bold.

Instance					$time_{avg}$	
	IVNS-	IVNS	IVNS-	IVNS	IVNS-	IVNS
RanReal240_01	223,891.05	224,785.27	179.47	97.88	114.90	147.28
RanReal240_02	203,789.58	204,415.88	160.64	102.77	120.46	160.27
RanReal240_03	198,236.57	198,626.93	96.21	170.83	145.08	146.88
RanReal240_04	224,617.33	225,227.11	86.58	237.09	149.30	162.63
RanReal240_05	195,181.62	195,228.86	82.96	91.62	149.44	161.77
RanReal240_06	215,491.95	216,474.84	126.77	169.61	109.82	165.88
RanReal240_07	208,889.10	209,004.05	74.03	120.82	147.34	120.41
RanReal240_08	204,207.83	204,958.19	230.94	139.71	158.62	174.29
RanReal240_09	208,562.04	208,789.79	50.70	148.77	119.07	160.17
RanReal240_10	192,253.79	192,788.59	100.09	154.18	147.11	170.78
RanReal240_11	203,776.92	204,523.95	109.20	96.57	149.70	148.56
RanReal240_12	200,251.35	200,904.16	145.50	126.79	101.56	169.33
RanReal240_13	201,581.76	202,139.55	143.75	155.41	123.42	150.75
RanReal240_14	228,332.63	228,512.11	228.63	170.71	108.74	141.27
RanReal240_15	190,130.92	190,914.31	190.24	174.08	137.21	144.95
RanReal240_16	203,072.49	203,834.68	197.23	131.32	110.16	147.86
RanReal240_17	194,564.42	195,114.49	92.24	109.08	113.22	146.03
RanReal240_18	194,370.83	194,853.70	121.64	126.17	107.11	163.49
RanReal240_19	198,496.90	199,019.23	144.56	109.79	111.23	176.31
RanReal240_20	211,418.56	212,046.92	130.93	130.92	106.77	149.66
RanReal480_01	550,160.70	554,331.89	777.24	415.49	304.77	369.26
RanReal480_02	507,893.39	509,519.84	457.07	569.71	286.47	416.40
RanReal480_03	493,226.43	495,847.80	584.05	426.86	307.52	415.80
RanReal480_04	518,088.82	520,891.75	462.43	648.97	303.57	380.20
RanReal480_05	481,778.58	482,595.19	436.13	497.25	310.42	338.96
RanReal480_06	529,846.53	532,888.64	537.12	555.89	240.48	338.25
RanReal480_07	542,575.23	544,530.14	675.84	413.13	310.92	388.74
RanReal480_08	529,264.25	531,417.94	464.02	555.42	265.15	373.48
RanReal480_09	551,329.66	555,098.72	547.85	514.64	246.68	388.08
RanReal480_10	516,929.26	518,612.02	466.20	589.69	282.97	403.46
RanReal480_11	520,357.01	522,814.96	424.04	402.37	318.02	382.52
RanReal480_12	498,472.55	500,580.84	410.95	540.16	258.80	386.04
RanReal480_13	530,678.39	533,763.20	628.41	423.89	261.10	420.59
RanReal480_14	509,895.03	512,975.73	685.12	408.22	251.09	401.09
RanReal480_15	513,947.87	516,017.98	645.76	408.07	255.18	375.96
RanReal480_16	544,921.69	548,276.15	933.12	590.34	332.77	366.12
RanReal480_17	533,634.44	536,655.06	498.31	402.38	277.45	389.54
RanReal480_18	521,497.91	524,650.86	775.43	494.86	288.94	379.04
RanReal480_19	519,184.62	521,180.84	716.91	550.77	230.32	405.61
RanReal480_20	514,755.82	517,261.92	500.77	530.58	286.77	398.1
# Better	0	40	500,77	330,30	200,77	550.1
# Equal	0	0				
# Worse	40	0				
p-value	40	2.54e-10				

generated by the first construction procedure presented in Section 2.3.

The computational results of this experiment are summarized in Table 5, including the average objective function value ( $f_{avg}$ ) and the average running time ( $time_{avg}$ ). In addition, the rows *Better*, *Equal* and *Worse* of the table denote the number of instances for which the corresponding algorithm obtained a better, equal, and worse average objective value compared to another one. The p-values from the non-parametric Friedman test are given in the last row of the table.

It can be observed from Table 5 that in terms of the average objective value, the EVND method achieves a better result than the standard VND method for 36 out of 40 instances, whereas both methods consumed a similar computational time for most instances. These outcomes demonstrate the interest of the EVND method compared to the standard VND method.

#### 4.2. Importance of 2-1 exchange neighborhood N<sub>3</sub>

The EVND method employs three complementary neighborhoods i.e.,  $N_1$ ,  $N_2$  and the 2-1 exchange neighborhood  $N_3$ . While  $N_1$  and  $N_2$  are very popular and their effectiveness has been shown on a number of the clustering problems in the literature (Brimberg et al., 2015, 2015; Deng and Bard, 2011; Palubeckis et al., 2015; Rodriguez et al., 2013; Urošević, 2014),  $N_3$  is not well studied and thus less understood. In this section, we assess the influence of  $N_3$  on the performance of the IVNS algorithm. The computational experiment was carried out as follows. We ran our IVNS and IVNS<sup>-</sup> methods 20 times to solve each instance, where IVNS<sup>-</sup> is a variant of IVNS in which  $N_3$  (corresponding to subroutine  $LSN_3$  of Algorithm 6) is disabled while keeping the other algorithmic ingredients unchanged. The experimental results are

**Table 7**Comparative results of the IVNS method with and without its diversified stage (IVNS-D), on the set of 40 representative instances. Each instance was independently solved 20 times by both algorithms respectively, and better results in terms of the average objective value between two algorithms are indicated in bold.

	$f_{a  u g}$		$\sigma$		$Time_{avg}$	
Instance	IVNS-D	IVNS	IVNS-D	IVNS	IVNS-D	IVNS
RanReal240_01	223,986.99	224,785.27	272.57	97.88	69.37	147.28
RanReal240_02	203,614.45	204,415.88	299.81	102.77	113.36	160.27
RanReal240_03	197,731.77	198,626.93	418.98	170.83	100.27	146.88
RanReal240_04	224,424.68	225,227.11	459.32	237.09	93.67	162.63
RanReal240_05	194,298.12	195,228.86	474.64	91.62	113.46	161.77
RanReal240_06	215,609.74	216,474.84	318.55	169.61	90.56	165.88
RanReal240_07	208,341.50	209,004.05	378.79	120.82	122.25	120.41
RanReal240_08	204,211.41	204,958.19	219.57	139.71	93.92	174.29
RanReal240_09	208,092.99	208,789.79	286.18	148.77	109.41	160.17
RanReal240_10	191,828.97	192,788.59	482.02	154.18	133.85	170.78
RanReal240_11	203,921.95	204,523.95	329.82	96.57	75.89	148.56
RanReal240_12	199,971.56	200,904.16	300.91	126.79	105.04	169.33
RanReal240_13	201,224,78	202,139.55	467.84	155.41	87.82	150.75
RanReal240_14	227,825.89	228,512.11	356.96	170.71	103.34	141.27
RanReal240_15	189,814.49	190,914.31	446.62	174.08	81.73	144.95
RanReal240_16	202,951.20	203,834.68	412.54	131,32	87.63	147.86
RanReal240_17	194,328.55	195,114.49	265.17	109.08	165.12	146.03
RanReal240_18	193,915.24	194,853.70	286.06	126.17	104,30	163.49
RanReal240_19	197,900.00	199,019.23	461.25	109.79	101.25	176.31
RanReal240_20	211,284.31	212,046.92	280.17	130.92	96.18	149.66
RanReal480_01	552,979.67	554,331.89	560.30	415.49	375.66	369.26
RanReal480_02	507,974.23	509,519.84	816.89	569.71	337.88	416.40
RanReal480_03	494,000.83	495,847.80	961.92	426.86	372.23	415.80
RanReal480_04	519,475.06	520,891.75	843.47	648.97	367.62	380.20
RanReal480_05	481,098.07	482,595.19	832.34	497.25	315.49	338.96
RanReal480_06	531,667.60	532,888.64	649.14	555.89	360.97	338.25
RanReal480_07	543,160.19	544,530.14	666.03	413.13	345.01	388.74
RanReal480_08	530,127.30	531,417.94	626,52	555.42	369.13	373.48
RanReal480_09	553,832.78	555,098.72	703,40	514.64	383.55	388.08
RanReal480_10	516,798.04	518,612.02	653.19	589.69	327.60	403.46
RanReal480_11	520,835.32	522,814.96	952,45	402.37	359.04	382.52
RanReal480_12	499,078.98	500,580.84	687.26	540.16	334.42	386.04
RanReal480_13	532,308.08	533,763.20	792,43	423.89	390.62	420.59
RanReal480_14	511,808.09	512,975.73	635.87	408.22	408.37	401.09
RanReal480_14	514,720.84	516,017.98	482.10	408.22	365.89	375.96
RanReal480_16	547,296.05	548,276.15	682.57	590.34	371.18	366.12
RanReal480_17	534,975.71	536,655.06	719.18	402.38	353.90	389.54
_	,	•	720.31	402.38 494.86	348.96	379.04
RanReal480_18	523,253.63	524,650.86				
RanReal480_19	519,700.55	521,180.84 517.361.03	564.17	550.77	378.77	405.61
RanReal480_20	515,947.56	517,261.92	567.91	530.58	325.18	398.10
# Better	0	40				
#Equal	0	0				
#Worse	40	0				
p-value		2.54e-10				

summarized in Table 6, including the average objective value  $f_{avg}$ , the standard deviation of objective value ( $\sigma$ ), and the average running time to reach its final objective value ( $time_{avg}$ ), and other symbols are the same as those in the previous tables.

Table 6 shows that without  $N_3$ , the performance of IVNS deteriorates for all instances in terms of average objective value. Moreover, the average computing times indicate that  $N_3$  helps the IVNS algorithm to continue its search for a longer time and thus to attain better solutions. This experiment demonstrates the usefulness of the 2-1 exchange neighborhood for the IVNS algorithm.

# 4.3. Importance of the diversification mechanism

The IVNS algorithm performs an intensified search stage with the iterated local optimization (lines 5–17 of Algorithm 1) and a diversified stage with the Shake procedure (line 18 of Algorithm 1). The diversified stage aims at producing transition states between two high-quality local optima, since these transition states are usually necessary to help the search process to move from a basin of attraction to another basin.

In order to assess the impact of this diversified stage on the performance of the IVNS algorithm, we created a variant of the IVNS method (denoted by IVNS-D) by removing the Shake operation of line 18 of Algorithm 1 while keeping other components of IVNS

**Table 8** Sensitivity analysis of the parameter m. Each instance was independently solved 20 times by the IVNS algorithm for each parameter value in the range  $\{4, 6, 8, 10, 12, 14, 16, 18\}$ , and the average objective values  $(f_{avg})$  over 20 runs are respectively reported.

	$f_{ m avg}$									
Instance/m	4	6	8	10	12	14	16	18		
RanReal240_01	224,747.89	224,743.77	224,750.51	224,736.19	224,734.17	224,784.25	224,725.39	224,718.09		
RanReal240_02	204,400.52	204,403.15	204,425.39	204,459.24	204,400.94	204,418.82	204,448.37	204,449.36		
RanReal240_03	198,657.99	198,717.34	198,685.34	198,691.63	198,679.41	198,658.77	198,661.33	198,685.05		
RanReal240_04	225,204.66	225,140.01	225,146.45	225,190.13	225,152.00	225,194.58	225,242.84	225,239.61		
RanReal240_05	195,275.00	195,287.90	195,237.80	195,259.40	195,256.56	195,199.85	195,246.73	195,246.16		
RanReal240_06	216,533.35	216,475.23	216,454.28	216,497.78	216,513.23	216,465.82	216,515.60	216,491.18		
RanReal240_07	209,098.93	209,062.00	209,060.14	209,043.44	209,092.56	209,024.46	209,021.64	209,029.07		
RanReal240_08	204,914.37	204,947.69	204,883.38	204,950.99	204,938.25	204,962.13	204,943.61	204,930.28		
RanReal240_09	208,826.26	208,832.55	208,883.41	208,833.99	208,799.52	208,794.97	208,822.74	208,745.94		
RanReal240_10	192,736.25	192,755.01	192,815.78	192,698.13	192,811.66	192,776.03	192,640.29	192,796.48		
RanReal240_11	204,478.83	204,444.22	204,464.45	204,448.14	204,470.10	204,495.97	204,478.49	204,454.86		
RanReal240_12	200,921.94	200,878.36	200,830.81	200,867.36	200,808.50	200,783.89	200,887.89	200,824.69		
RanReal240_13	202,094.54	202,050.52	202,042.04	202,105.03	201,987.09	202,076.88	202,094.15	202,098.55		
RanReal240_14	228,474.87	228,487.74	228,483.37	228,478.88	228,531.74	228,528.80	228,525.70	228,557.20		
RanReal240_15	190,924.82	190,924.06	190,907.09	190,958.90	190,939.27	190,933.87	190,944.36	190,913.93		
RanReal240_16	203,763.28	203,789.09	203,798.46	203,856.97	203,816.01	203,850.67	203,746.08	203,801.18		
RanReal240_17	195,115.48	195,125.04	195,150.36	195,147.21	195,078.97	195,106.20	195,132.90	195,141.27		
RanReal240_18	194,947.69	194,852.90	194,893.01	194,756.48	194,875.62	194,852.77	194,866.09	194,836.40		
RanReal240_19	198,962.52	199,040.68	198,966.06	198,984.75	198,934.98	199,008.06	198,988.46	198,974.27		
RanReal240_20	212,074.56	212,092.28	212,056.10	212,032.47	211,998.49	212,046.46	212,001.59	212,105.61		
Average	205,607.69	205,602.48	205,596.71	205,599.85	205,590.95	205,598.16	205,596.71	205,601.96		

**Table 9** Sensitivity analysis of the parameter  $\beta_{max}$ . Each instance was independently solved 20 times by the IVNS algorithm for each parameter value in the range  $\{5, 10, 15, 20, 25, 30, 35, 40\}$ , and the average objective values  $(f_{avg})$  over 20 runs are respectively reported.

	$f_{avg}$							
Instance/ $\beta_{\text{max}}$	5	10	15	20	25	30	35	40
RanReal240_01	224,676.47	224,758.30	224,738.12	224,753.19	224,749.78	224,769.53	224,739.29	224,675.94
RanReal240_02	204,384.58	204,428.48	204,454.16	204,422.07	204,403.51	204,425.57	204,381.89	204,378.07
RanReal240_03	198,621.22	198,612.88	198,694.68	198,622.62	198,700.29	198,600.93	198,693.71	198,489.30
RanReal240_04	225,111.94	225,215.51	225,217.08	225,262.81	225,203.17	225,225.99	225,039.70	225,189.87
RanReal240_05	195,202.16	195,260.55	195,222.28	195,278.64	195,256.13	195,158.88	195,199.11	195,273.42
RanReal240_06	216,421.63	216,540.77	216,524.29	216,511.22	216,500.76	216,531.96	216,491.43	216,397.47
RanReal240_07	208,954.86	209,051.89	209,053.73	209,066.47	209,030.62	209,042.97	209,025.39	208,978.27
RanReal240_08	204,957.91	204,935.06	204,950.87	204,998.23	204,918.30	204,968.93	204,864.60	204,754.79
RanReal240_09	208,775.98	208,872.97	208,803.31	208,762.89	208,814.66	208,793.45	208,786.65	208,819.28
RanReal240_10	192,729.71	192,761.86	192,767.85	192,774.27	192,787.08	192,666.16	192,647.02	192,789.16
RanReal240_11	204,457.23	204,464.00	204,483.40	204,492.54	204,490.35	204,487.99	204,463.07	204,385.61
RanReal240_12	200,897.62	200,913.55	200,930.03	200,918.37	200,868.58	200,824.06	200,865.56	200,857.78
RanReal240_13	202,110.38	202,121.17	202,114.62	202,100.76	202,097.89	202,124.86	202,087.71	202,049.26
RanReal240_14	228,454.78	228,514.51	228,484.05	228,516.60	228,536.93	228,490.37	228,494.82	228,413.24
RanReal240_15	190,869.97	190,940.52	190,895.56	190,944.22	190,966.06	190,940.83	190,868.54	190,804.69
RanReal240_16	203,702.06	203,824.85	203,821.77	203,859.91	203,778.12	203,774.10	203,819.94	203,790.17
RanReal240_17	195,096.28	195,211.71	195,161.74	195,155.80	195,166.45	195,159.81	195,051.22	194,973.69
RanReal240_18	194,861.25	194,879.05	194,882.91	194,900.61	194,900.42	194,861.18	194,881.60	194746.11
RanReal240_19	199,005.21	199,037.12	199,047.45	198,990.69	198,983.59	199,071.70	198,893.53	198,805.32
RanReal240_20	211,997.53	212,060.94	212,059.02	212,049.21	212,060.15	212,052.58	211,986.43	211,926.56
Average	205,564.44	205,620.28	205,615.35	205,619.06	205,610.64	205,598.59	205,564.06	205,524.90

unchanged. We ran IVNS-D and IVNS 20 times to solve each instance. The experimental results are summarized in Table 7, where the symbols have the same meanings as those in the previous tables.

Table 7 indicates that IVNS-D deteriorates the results of IVNS. First, IVNS-D performs worse than IVNS on all instances in terms of the average objective value. Second, concerning the standard deviation ( $\sigma$ ) of the objective value, IVNS obtained a better result for all instances. This experiment confirms the usefulness of the additional diversification stage introduced in line 18 of Algorithm 1.

# 4.4. Sensitivity analysis of parameters

Our IVNS algorithm employs two main parameters, i.e., m and  $\beta_{max}$ . Parameter m is employed in the EVND procedure (Section 2.4.3) to control the exploitation balance between the different neighborhoods, a larger value of m leading to a more balanced neighborhood exploitation. Parameter  $\beta_{max}$  is used to control the strength of intensification search, a larger value of  $\beta_{max}$  implying a stronger intensification for the IVNS algorithm. In this section we show a sensitivity analysis of these two key parameters, which also helps to find an appropriate value for each of them.

In this study, we carried out two additional experiments based on 20 *RanReal* instances with n=240. In the first experiment, we varied the value of m within the range {4, 6, 8, 10, 12, 14, 16, 18} and ran the algorithm 20 times for each value of m and each instance, while keeping other parameters with their default values as shown in Table 1. The computational results are summarized in Table 8, where the second row indicates the values of m, the first column gives the names of instances, the other columns show the average objective function values over 20 independent runs ( $f_{avg}$ ) for each value of m and each instance, and the last row shows the average results over all instances. Similarly, we varied in the second experiment the value of  $\beta_{max}$  within the range {5, 10, 15, 20, 25, 30, 35, 40}. The computational results are summarized in Table 9, where the second row gives the values of  $\beta_{max}$ , and the other entries have the same meanings as those in Table 8.

First, we observe from Table 8 that the performance of the IVNS algorithm is not sensitive to the setting of parameter m. Specifically, for most instances the different values of m leaded to very similar results in terms of  $f_{avg}$ . Indeed, the relative difference between the results yielded by the different parameter values across the 20 instances is very small  $\left( \leq \frac{(205607.69 - 205590.95)}{205607.69} \times 100\% = 0.0081\% \right)$ . Hence, the default value of m was set to 10 in this work. As for  $\beta_{max}$ , Table 9 shows that for most instances the tested values leaded also to similar results in terms of  $f_{avg}$ , with a very small relative difference between the results yielded by the different  $\beta_{max}$  values  $\left( \leq \frac{(205620.28 - 205524.90)}{205620.28} \times 100\% = 0.046\% \right)$ . These outcomes indicate that the IVNS algorithm is not sensitive to the setting of parameter  $\beta_{max}$ . Consequently, to ensure that a lasting intensified search effect when a long computational time is allowed, the default value of  $\beta_{max}$  was set to 30 in this study .

#### 5. Conclusions

The capacitated clustering problem (CCP) is a general and useful model for a number of applications. It also generalizes three well-known NP-hard problems: the maximally diverse grouping problem, the graph partitioning problem, and the handover minimization problem. In this paper, we proposed the iterated variable neighborhood search (IVNS) algorithm for solving the CCP. The proposed algorithm organically combines an extended variable neighborhood descent (EVND) method for intensification and a shake procedure for diversification.

The proposed algorithm was assessed on the 133 instances commonly used in the literature, and the computational results indicated that our IVNS algorithm significantly outperforms the state-of-the-art CCP algorithms both in terms of solution quality and computational efficiency. In particular, the proposed algorithm improved the best known results (new lower bounds) for 28 out of 83 handover minimization instances, while matching the best known results for the 55 remaining instances.

The investigations of several essential components of the proposed algorithm shed light on the following points. First, for the CCP, the EVND method usually outperforms the standard variable neighborhood descent method in terms of the local search ability, and the 2-1 exchange neighborhood  $N_3$  reinforces the intensified search capacity of the EVND method. Second, the diversification stage is essential for the proposed algorithm to reach a suitable trade-off between the diversification and intensification of the search process.

Based on this work, we advance some research perspectives for further improvements. First, within the IVNS algorithm, diversification is ensured by the shake procedure as well as the shake strength. Since different degrees of diversification may be needed at different search stages, it would be interesting to investigate adaptive techniques able to adjust dynamically the shake strength. Moreover, to escape deep local optima, it would also be useful to study other diversification methods like random or adaptive restarts. Second, using the presented EVND method as a local optimization procedure, it may be possible to devise more efficient hybrid evolutionary algorithms for the CCP. Third, the IVNS algorithm only visits feasible solutions. Meanwhile, previous studies like (Chen et al., 2016; Glover and Hao, 2011) showed that tunneling through feasible and infeasible regions can improve the performance of the search process. It would be relevant to study dedicated methods able to explore infeasible regions in a controlled manner. Finally, given that the basic idea of the proposed IVNS algorithm, i.e., integrating organically the EVND method with multiple neighborhoods and a diversified shake procedure, is independent of the CCP, it would be interesting to examine its applicability to other grouping or clustering problems.

# Acknowledgments

We are grateful to the reviewers for their valuable comments which helped us to improve the paper. This work is partially supported by the PGMO project (2013–2015, Jacques Hadamard Mathematical Foundation, Paris, France) and a post-doc grant (for X.J. Lai) from the Region of Pays de la Loire (France).

#### Appendix

See Table 10.

Table 10 Computational results of IVNS on the small handover minimization instances. The results are given in the form of minimization to make a direct comparison with the best known results in the literature.

		IVNS				
Instance	BKS	$f_{best}$	$f_{avg}$	$f_{worst}$	σ	time <sub>avg</sub> (s)
20_5_270001	540	540	540.00	540	0.00	0.00
20_5_270002	54	54	54.00	54	0.00	0.00
20_5_270003	816	816	816.00	816	0.00	0.00
20_5_270004	126	126	126.00	126	0.00	0.00
20_5_270005	372	372	372.00	372	0.00	0.00
20_10_270001	2148	2148	2148.00	2148	0.00	0.00
20_10_270002	1426	1426	1426.00	1426	0.00	0.00
20_10_270003	2458	2458	2458.00	2458	0.00	0.00
20_10_270004	1570	1570	1570.00	1570	0.00	0.00
30_5_270001	772	772	772.00	772	0.00	0.00
30_5_270002	136	136	136.00	136	0.00	0.00
30_5_270003	920	920	920.00	920	0.00	0.00
30_5_270004	52	52	52.00	52	0.00	0.00
30_5_270005	410	410	410.00	410	0.00	0.00
30_10_270001	3276	3276	3276.00	3276	0.00	0.00
30_10_270002	1404	1404	1404.00	1404	0.00	0.00
30_10_270003	2214	2214	2214.00	2214	0.00	0.00
30_10_270004	2150	2150	2150.00	2150	0.00	0.00
30_10_270005	2540	2540	2540.00	2540	0.00	0.00
30_15_270001	6178	6178	6178.00	6178	0.00	0.00
30_15_270002	4042	4042	4042.00	4042	0.00	0.00
30_15_270003	4126	4126	4126.00	4126	0.00	0.00
30_15_270004	3920	3920	3920.00	3920	0.00	0.00
40_5_270001	610	610	610.00	610	0.00	0.00
40_5_270002	136	136	136.00	136	0.00	0.00
40_5_270003	234	234	234.00	234	0.00	0.00
40_5_270004	232	232	232.00	232	0.00	0.02
40_5_270005	774	774	774.00	774	0.00	0.00
40_10_270001	4544	4544	4544.00	4544	0.00	0.00
40_10_270002	2068	2068	2068.00	2068	0.00	0.00
40_10_270003	2090	2090	2090.00	2090	0.00	0.00
40_10_270004	1650	1650	1650.00	1650	0.00	0.01
40_10_270005	4316	4316	4316.00	4316	0.00	0.00
40_15_270001	8646	8646	8646.00	8646	0.00	0.01
40_15_270002	4586	4586	4586.00	4586	0.00	0.03
40_15_270003	5396	5396	5396.00	5396	0.00	0.01
40_15_270004	4800	4800	4800.00	4800	0.00	0.00
40_15_270005	6272	6272	6272.00	6272	0.00	0.00

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