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(Article begins on next page)

# Lagrangian matheuristics for the Quadratic Multiple Knapsack Problem

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#### Abstract

The Quadratic Multiple Knapsack Problem (QMKP) is a challenging combinatorial optimization problem combining the well-known Quadratic Knapsack Problem with the Multiple Knapsack Problem. After a long stream of research devoted to metaheuristic approaches for large-scale instances, only recently some authors started to study the mathematical properties of the QMKP and proposed exact solution methods for benchmark instances of smaller size. In this paper, we propose the first matheuristic approach for solving the QMKP approximately. The proposed method exploits the strength of a Lagrangian relaxation for the natural quadratic formulation of the QMKP to derive heuristic solutions. Postoptimization local search procedures are embedded in the final framework. Experimental studies show that the resulting deterministic matheuristic approach consistently delivers solutions of very good quality in short computing times.

**Keywords:** Combinatorial optimization; Knapsack problems; Matheuristics; Lagrangian relaxation.

### 1 Introduction

In the Quadratic Multiple Knapsack Problem (QMKP), one is given n items, each having a profit  $p_i$  and a weight  $w_i$  (i = 1, ..., n), m knapsacks, each having a capacity  $c_k$  (k = 1, ..., m), and an additional  $n \times n$  symmetric profit matrix  $[p_{ij}]$ . The QMKP calls for packing m disjoint subsets of items into the knapsacks so that the total weight in any knapsack does not exceed its capacity, with the objective of maximizing the overall profit given by: (i) the profit  $p_i$  of each selected item i, plus (ii) a profit  $p_{ij}$  for each pair of items (i, j) packed into the same knapsack. A natural mathematical formulation of the

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problem is the following binary quadratic program:

$$\max \quad \sum_{i=1}^{n} \sum_{k=1}^{m} p_i x_{ik} + \sum_{i=1}^{n-1} \sum_{j=i+1}^{n} \sum_{k=1}^{m} p_{ij} x_{ik} x_{jk} \tag{1}$$

s.t. 
$$\sum_{i=1}^{n} w_i x_{ik} \le c_k$$
  $(k = 1, \dots, m)$  (2)

$$\sum_{k=1}^{m} x_{ik} \le 1 \qquad (i = 1, \dots, n) \qquad (3)$$

$$x_{ik} \in \{0, 1\} \qquad (i = 1, \dots, n; k = 1, \dots, m), \qquad (4)$$

where  $x_{ik}$  takes the value one iff item *i* is packed into knapsack k (i = 1, ..., n; k = 1, ..., m). Note that the objective function (1) is the sum of a linear term and a quadratic term expressing linear and pairwise profits, respectively. Inequalities (2) express the traditional knapsack capacity constraints, while inequalities (3) forbid an item to be packed into more that one knapsack. We assume throughout the paper that  $p_i$ ,  $w_i$ , and  $c_k$  are positive integers (i = 1, ..., n, k = 1, ..., m), and that  $p_{ij}$  is a nonnegative integer with  $p_{ij} = p_{ji}$  (i, j = 1, ..., n).

The QMKP comes from the combination of two intensively studied combinatorial optimization problems: the Quadratic Knapsack Problem (QKP) and the Multiple Knapsack Problem (MKP). Both the QKP and the MKP are special cases of the QMKP, arising when m = 1 (so constraints (3) become redundant) and when  $p_{ij} = 0$  (i, j = 1, ..., n)(so the quadratic term in (1) vanishes), respectively. When both restrictions hold, the resulting problem is the famous 0-1 Knapsack Problem (01KP). For the vast literature on QKP, MKP, and 01KP, the reader is referred to the classical monographs by Martello and Toth [15] and Kellerer et al. [14], as well as to the recent survey by Cacchiani et al. [3].

While the 01KP is weakly  $\mathcal{NP}$ -hard and can be solved in pseudo-polynomial time, both the QKP and the MKP are strongly  $\mathcal{NP}$ -hard, implying strong  $\mathcal{NP}$ -hardness for the QMKP as well. In addition to its theoretical intractability, the QMKP (which arises in a number of real-world situations like, e.g., project management, capital budgeting, product-distribution system design) is very difficult to solve in practice.

The QMKP was defined in 2006 by Hiley and Julstrom [12], who studied the case (commonly encountered in the literature) in which all capacities are equal. They proposed three heuristic algorithms for its approximate solution and a benchmark set of instances with  $n \in \{100, 200\}$  and  $m \in \{3, 5, 10\}$ . In the following decade, a number of meta-heuristic approaches was computationally tested on these instances. We mention in particular those by Singh and Baghel [18] (genetic algorithm), Sundar and Singh [20] (artificial bee colony), Soak and Lee [19] (memetic algorithm), Garcia-Martinez et al. [10, 11] (strategic oscillation and a Tabu search), Chen et al. [5, 6] (threshold search and evolutionary path relinking). The last approaches were later improved by Qin et al. [16] and Tlili et al. [21]. A local branching approach was recently proposed by Aïder et al. [1].

The study of exact methods for the QMKP started in 2019, when Bergman [2] presented the first exact algorithm for the problem, obtained from an exponentially-sized formulation, solved through Branch-and-Price. He used the generation scheme of [12] to produce new (smaller) instances with  $n \in \{20, 25, 30, 35\}$  and  $m \in \{3, 5, 10\}$ . More recently, Galli et al. [9] studied effective polynomial-size formulations and relaxations of the problem leading to a decomposable structure, and proposed new medium-size benchmark sets (produced with the instance generator of [2]) with  $n \in \{40, 45, 50, 55, 60\}$  and  $m \in \{3, 5, 10\}$ . Computational experiments showed that off-the-shelf solvers like CPLEX applied to the decomposable model are competitive with the Bergman algorithm for small-size instances. A combinatorial branch-and-bound algorithm was very recently proposed by Fleszar [7] for the case where all knapsack capacities are equal, computationally improving on the previous methods. In the following, we consider such a case. (Worth is mentioning that all benchmark instances from the literature consider the same capacity value for each knapsack.)

This paper is devoted to the study of matheuristic methods that can be derived from the mathematical models of the problem. In Section 2, we introduce the Lagrangian relaxation of (1)-(4) and its dual (see [9] for more details), which are used in Section 3 to obtain a matheuristic approach based on the solution of a variant of the set packing problem. A postoptimization stage is introduced in Section 4, and the overall deterministic algorithm is presented in Section 5. The outcome of computational experiments on small and medium size instances is reported in Section 6. Conclusions follow in Section 7.

### 2 Lagrangian relaxation

Let us define a vector  $\lambda$  of nonnegative Lagrangian multipliers  $\lambda_i$  (i = 1, ..., n), and relax in a Lagrangian fashion cardinality constraints (3). We get

$$L(\lambda) = \sum_{\substack{i=1\\n}}^{n} \lambda_i + \max \sum_{\substack{i=1\\k=1}}^{n} \sum_{k=1}^{m} (p_i - \lambda_i) x_{ik} + \sum_{\substack{i=1\\j=i+1}}^{n-1} \sum_{k=1}^{n} \sum_{k=1}^{m} p_{ij} x_{ik} x_{jk}$$
(5)

s.t. 
$$\sum_{i=1}^{n} w_i x_{ik} \le c_k$$
  $(k = 1, ..., m)$  (6)

$$x_{ik} \in \{0, 1\}$$
  $(i = 1, \dots, n; k = 1, \dots, m).$  (7)

Note that the dualization of constraints (3) decomposes the model into m independent (single) QKPs, sharing the same input for what concerns the items, with capacities  $c_1, c_2, \ldots c_m$ , respectively. Further observe that, if two knapsacks have the same capacity value, then the corresponding QKPs are identical. Therefore, for a given vector  $\lambda$  of multipliers, upper bound  $L(\lambda)$  can be obtained by solving a single QKP for each different capacity value, i.e., just one single QKP when all capacities are identical. Also note that the Lagrangian linear profits  $(p_i - \lambda_i)$   $(i = 1, \ldots, n)$  can take negative values.

Coming to the Lagrangian dual problem of finding  $L(\lambda^*) = \min_{\lambda \ge 0} L(\lambda)$ , while the optimal multipliers cannot be computed efficiently, "good" approximations of  $\lambda^*$  can be obtained by traditional subgradient optimization. A more sophisticated technique, known as the *proximal bundle* method (see Frangioni [8]), which can be halted after a prefixed number of iterations, yields either the optimal multiplier vector  $\lambda^*$  or a good approximation of it.

The computer code ProximalBundle, implemented by Frangioni [8] and freely available at the GitLab repository, produces, at each iteration, a possibly improved vector  $\lambda$  of Lagrangian multipliers and the corresponding upper bound  $L(\lambda)$ . We denote as  $\tilde{\lambda}$  the best Lagrangian multiplier vector determined during the execution of the bundle procedure, and with  $L(\tilde{\lambda})$  the corresponding upper bound. In our case, each bundle iteration requires the optimal solution of a QKP: this is obtained by executing the C code quadknap, available at the home page of D. Pisinger. This code implements the QKP algorithm by Caprara et al. [4], and is currently the most effective code for optimally solving the QKP when the linear profits can take negative values. Since the QKP is a strongly  $\mathcal{NP}$ -hard problem, it might be necessary to halt the execution of quadknap after a certain elapsed time (in which case the upper bound provided in the current iteration of the bundle procedure may no longer be valid). In the next section, we show that, in any case, the set of QKP solutions provided by the bundle procedure can be used to derive an effective matheuristic algorithm for the QMKP.

## 3 Matheuristic scheme

In this section, we introduce the basic idea underlying the Lagrangian deterministic matheuristic for the QMKP. Our starting point is that of creating a family  $\mathcal{F}$  of promising subsets of items that can be used to construct a feasible partial solution to the QMKP by filling "some" of the knapsacks. A promising subset is one that is likely to appear in a good quality (possibly optimal) solution to the QMKP. To do that, we take advantage of the subsets of items produced during the bundle iterations (as solutions to the QKP subproblems). In particular, our criterion for inserting a subset in  $\mathcal{F}$  is that the upper bound  $L(\lambda)$  produced at the corresponding bundle iteration be smaller than a given threshold  $UBM = \alpha L(\tilde{\lambda})$  (where  $\alpha$  is a prefixed parameter with  $\alpha > 1$ ). For the case of identical subsets, only one copy is retained. We denote the selected subsets as

$$B_h \subset \{1, \dots, n\}$$
  $(h = 1, \dots, s),$  (8)

where  $s = |\mathcal{F}|$ . Let

$$a_{hi} = \begin{cases} 1 & \text{if item } i \in B_h \\ 0 & \text{otherwise} \end{cases} (h = 1, \dots, s; i = 1, \dots, n)$$
(9)

be the corresponding incidence matrix, and observe that subsets  $B_h$  will normally have nonempty intersections. Note that only "some" knapsacks will be filled according to the subsets in  $\mathcal{F}$ . (In fact, there may not even exist *m* disjoint subsets in  $\mathcal{F}$ .) More precisely, we define a parameter  $r_{\text{max}} < m$  representing the maximum number of (disjoint) subsets of  $\mathcal{F}$  that can be selected to fill the knapsacks. Our approach consists of two phases:

- 1. Select at most  $r_{\max}$  disjoint subsets  $B_h$  from family  $\mathcal{F}$ . Let  $\mathcal{S}^*$  denote the family of selected subsets, and  $\overline{r} = |\mathcal{S}^*|$ .
- 2. Pack the subsets in  $S^*$  into knapsacks  $1, \ldots, \overline{r}$ . Then iteratively pack into knapsacks  $\overline{r}+1, \ldots, m$  subsets of unpacked items obtained by solving a series of (single) QKPs.

### 3.1 Phase 1

This phase requires to solve a combinatorial problem that can be seen as a generalization of the well-known *Set Packing Problem* (SPP), obtained by limiting the number of selected subsets. We denote such problem as the *Cardinality Constrained Set Packing Problem* (CCSPP). The CCSPP can then be formally defined as

(CCSPP) max 
$$\sum_{h=1}^{s} y_h$$
 (10)

s.t. 
$$\sum_{h=1}^{s} a_{hi} y_h \le 1$$
  $(i = 1, \dots, n)$  (11)

$$\sum_{h=1}^{s} y_h \le r_{\max} \tag{12}$$

$$y_h \in \{0, 1\}$$
  $(h = 1, \dots, s),$  (13)

where  $y_h$  takes the value one if subset  $B_h$  is chosen and the value zero otherwise. Constraints (11) ensure that the selected subsets are disjoint, while constraint (12) imposes the maximum cardinality to the family of subsets. The pseudocode implementing Phase 1 is given in Algorithm 1.

Algorithm 1: Procedure Bundle
<b>input</b> : A QMKP instance $I$
<b>output:</b> A family $\mathcal{S}^* \subseteq \mathcal{F}$ of disjoint item subsets
A Lagrangian multiplier vector $\tilde{\lambda}$
1 execute ProximalBundle on $(5)$ - $(7)$ ;
<b>2</b> let $B_h$ $(h = 1, \ldots, s)$ be the subsets in family $\mathcal{F}$ ;
<b>3 solve</b> the CCSPP $(10)$ - $(13)$ ;
4 $y^{\star} \leftarrow \text{CCSPP}$ optimal solution;
5 $S^{\star} \leftarrow \{B_h : h \in \{1, \dots, s\}, y_h^{\star} = 1\}.$

In the special case where  $r_{\text{max}} = s$  the decision version of the CCSPP reduces to the SPP, one of the famous Karp's 21  $\mathcal{NP}$ -complete problems [13]. This proves that the CCSPP is  $\mathcal{NP}$ -hard.

The solution to the CCSPP in procedure Bundle is obtained through an MILP solver (CPLEX 20.0 in our case).

### **3.2** Phase 2

The second phase of the matheuristic approach consists of either one or two steps, depending on the  $\bar{r}$  value produced by the previous phase. If  $\bar{r} < m - 1$ , we iteratively solve  $m - \bar{r} - 1$  (single) QKPs on the unpacked items, with Lagrangian linear profits  $(p_i - \tilde{\lambda}_i)$ . In any case, the *m*-th knapsack is assigned a set obtained by solving a QKP with the original linear profits  $p_i$ . We denote as QKP(T) the (single) QKP induced by a subset T of items with knapsack capacity c and appropriately defined item profits. The corresponding procedure is shown in Algorithm 2.

Algorithm 2: Procedure Pack **input** : A family  $\mathcal{S}^* \subset \mathcal{F}$  of disjoint item subsets A Lagrangian multiplier vector  $\lambda$ **output:** A feasible solution x to the QMKP 1  $\overline{r} := |\mathcal{S}^{\star}|;$ **2** pack the  $\overline{r}$  subsets of  $\mathcal{S}^{\star}$  into knapsacks  $1, \ldots, \overline{r}$ ; **3**  $T \leftarrow$  set of the remaining (unpacked) items; 4 if  $\overline{r} < m - 1$  then for  $\ell := \overline{r} + 1$  to m - 1 do 5 **execute quadknap**(T) with linear profits  $p_i - \widetilde{\lambda}_i \ (i \in T);$ 6  $Q \leftarrow$  set of items in the QKP(T) solution; 7 **pack** the items of Q into knapsack  $\ell$ ; 8  $T := T \setminus Q$ 9 end 10 11 end 12 execute quadknap(T) with linear profits  $p_i$   $(i \in T)$ ; **13**  $Q \leftarrow$  set of items in the QKP(T) solution; 14 pack the items of Q into knapsack m.

The core matheuristic scheme consists thus of consecutively executing procedures Bundle and Pack.

# 4 Postoptimization

In this section, we show how to heuristically improve a feasible solution to the QMKP. We adopt the following notation:

- z(x) = objective function value of a feasible solution x;
- $x^{\text{best}}$  = incumbent solution, UB = upper bound on the optimal solution value. These quantities are assumed to be global variables;
- K(i) = knapsack where item *i* is currently packed (i = 1, ..., n). We assume that an unpacked item *i* is packed into a *dummy knapsack* m + 1;
- X(k) = set of items currently packed into knapsack k, with  $W(k) = \sum_{i \in X(k)} w_i$ (k = 1,...,m) and  $W(m+1) = \infty$ ;

The postoptimization approach makes use of two procedures:

• a procedure Delete which, given a feasible solution, removes, according to two possible criteria, some of the currently packed items;

• a procedure LocalSearch which improves a feasible solution through classical local search moves.

### 4.1 Item removal

For each knapsack k (k = 1, ..., m) and each item  $i \in X(k)$ , we define a global profit  $\overline{p}_i$  given by its profit  $p_i$  plus the overall pairwise profit it currently produces. The items of X(k) can be sorted either according to nondecreasing global profit, or according to global profit per unit weight, and a prefixed percentage  $\gamma$  of the first sorted items is then removed from X(k). The corresponding procedure is provided in Algorithm 3.

Algor	rithm 3: Procedure Delete							
in	<b>uput</b> : A feasible solution $x$ to an instance $I$							
	a parameter $d \in \{1, 2\}$ , a positive parameter	$\gamma < 1$						
01	<b>output:</b> A feasible solution $x^-$ to I with some items removed							
1 x <sup>-</sup>	$\bar{x} := x;$	/* initialize */						
2 fo	or $k := 1$ to $m$ do							
3	foreach $i \in X(k)$ do $\overline{p}_i := p_i + \sum_{j \in X(k) \setminus \{i\}} p_{ij};$	/*global profit*/						
4	if $d = 1$ then sort the items $i \in X(k)$ by nondecrea	asing $\overline{p}_i$ values						
5	<b>else</b> sort the items $i \in X(k)$ by nondecreasing $\overline{p}_i/w_i$ values;							
6	$nd := \lfloor \gamma \cdot  X(k)  \rfloor;$							
7	<b>remove</b> the first $nd$ items $i$ from $X(k)$ and set the	e corresponding $x_{ik}^-$ to 0						
s en	nd							

### 4.2 Local search

In order to improve the solution produced by the Lagrangian matheuristic described in the previous section, we embedded some classical local search basic moves from the literature (item exchanges and relocations). The framework is inspired by a similar one successfully adopted by Qin et al. [16] within a randomized, non-deterministic Tabu search algorithm for the QMKP. The main characteristic of such scheme is that of alternating moves that maintain feasibility and moves that can produce infeasible solutions, and repairing the final solution if infeasible.

The deterministic local search method, presented in Algorithm 4, consists of two main steps:

- Step 1 performs a local search that preserves feasibility. It iteratively invokes two procedures:
  - BestExchange tries to improve a feasible solution by swapping pairs of items;
  - BestReloc tries to improve a feasible solution by moving items to a different knapsack.

Algorithm 4: Procedure Local Search

```
input : A feasible QMKP solution x
    output: A possibly improved QMKP solution x^{\text{best}}
 1 x' := x;
 2 /*Step 1. Feasible Local Search */
 3 d := 1;
 4 while d \leq 2 do
         if d = 1 then \overline{x} \leftarrow \text{BestExchange}(x') else \overline{x} \leftarrow \text{BestReloc}(x');
         if z(\overline{x}) > z(x') then x' := \overline{x}, d := 1 else d := d + 1;
 6
 7 end
 8 x^{\text{best}} := x';
 9 /*Step 2. Infeasible Local Search */
10 Stop := Feas := false;
11 while Stop = false do
          if Feas = true then
12
13
               \tilde{x} \leftarrow \text{BestExchange}(x');
               \overline{x} \leftarrow \text{BestReloc}(x');
\mathbf{14}
               if z(\tilde{x}) > z(\overline{x}) then \overline{x} := \tilde{x};
\mathbf{15}
               if z(\overline{x}) > z(x') then x' := \overline{x}, Feas := true else Feas := false;
\mathbf{16}
               x^{\text{best}} := x'
\mathbf{17}
          else
\mathbf{18}
               \hat{x} := x';
19
               for k := 1 to m do
\mathbf{20}
                    if W(k) < c_k then \overline{x} \leftarrow \text{InfRelocA}(x', k) else \overline{x} \leftarrow \text{InfRelocB}(x', k);
\mathbf{21}
                    if z(\overline{x}) > z(\hat{x}) then \hat{x} := \overline{x}
\mathbf{22}
               end
\mathbf{23}
               if z(\hat{x}) = z(x') then
\mathbf{24}
                    Stop := true
\mathbf{25}
               else
\mathbf{26}
                     Stop := \texttt{false}, x' := \hat{x};
\mathbf{27}
                    if x' is feasible then x^{\text{best}} := x', Feas := true else Feas := false
\mathbf{28}
               end
\mathbf{29}
          end
30
31 end
32 if x' is infeasible then x' \leftarrow \text{Repair}(x');
33 if z(x') > z(x^{\text{best}}) then x^{\text{best}} := x'
```

Step 1 iterates BestExchange until no improvement is obtained, and then switches to BestReloc (possibly returning later to BestExchange);

- Step 2 consists of a main while loop which alternates between a stage that preserves feasibility and a stage that can produce infeasible solutions. The former stage invokes in sequence both procedures above, iterating until no improvement is obtained. The latter stage is then executed by invoking two procedures which can produce an infeasible solution by relocating items:
  - procedures InfRelocA and InfRelocB try to move a currently packed item to a different knapsack, irrespective of a possible capacity violation.

If such relocating procedures are unable to produce an improved (even if infeasible) solution, the local search terminates. Otherwise, one of the two stages is performed, depending on the feasibility of the improved solution. When the **while** loop is ended, if the incumbent solution is infeasible,

procedure Repair removes, exchanges, or relocates items until a feasible solution is obtained.

A detailed description of the local search procedures follows.

#### 4.2.1 Local search procedures

In this section, we provide more details on the inner local search procedures. Despite being based on classical moves from the literature, for the sake of completeness the corresponding pseudocodes are given in the Appendix.

Procedure BestExchange, shown in Algorithm 7, tries to exchange pairs of items (i,j) (with j > i) not packed into the same knapsack, provided the exchange is both feasible and profitable. (Note that one of the two items can currently be unpacked.) An i - j exchange is *feasible* if both capacity constraints of knapsacks K(i) and K(j) are preserved. The profit variation produced by the exchange is:

$$\Delta = \begin{cases} \sum_{\ell \in X(K(j)) \setminus \{j\}} p_{i\ell} + \sum_{\ell \in X(K(i)) \setminus \{i\}} p_{j\ell} - \sum_{\ell \in X(K(i))} p_{i\ell} - \sum_{\ell \in X(K(j))} p_{j\ell} \text{ if } K(i) \le m \text{ and } K(j) \le m \end{cases}$$
$$\Delta = \begin{cases} \sum_{\ell \in X(K(i)) \setminus \{j\}} p_{j\ell} - p_i - \sum_{\ell \in X(K(i))} p_{i\ell} & \text{if } K(i) \le m \text{ and } K(j) = m + 1 \end{cases}$$
$$p_i + \sum_{\ell \in X(K(j)) \setminus \{j\}} p_{i\ell} - p_j - \sum_{\ell \in X(K(j))} p_{j\ell} & \text{if } K(i) = m + 1 \text{ and } K(j) \le m \end{cases}$$

where we assume that  $p_{ii} = 0 \ \forall i$ . An exchange is thus profitable if  $\Delta > 0$ .

Procedure BestReloc, shown in Algorithm 8, shares part of the structure of BestExchange. It tries to move an item i (possibly unpacked) to a different (non dummy) knapsack k, provided the move is both feasible and profitable. In this case, the profit variation produced

by the relocation is:

$$\Delta = \begin{cases} \sum_{\ell \in X(k)} p_{i\ell} - \sum_{\ell \in X(K(i))} p_{i\ell} & \text{if } K(i) \le m \\ p_i + \sum_{\ell \in X(k)} p_{i\ell} & \text{if } K(i) = m + 1 \end{cases}$$

and the move is profitable if  $\Delta > 0$ .

The next two procedures, InfRelocA and InfRelocB, can produce infeasible solutions in which the capacity constraint of some knapsack is violated. Procedure InfRelocA is given in Algorithm 9. It tries to move to a specified knapsack k an item currently packed into a more filled knapsack. The item i (if any) producing the highest *normalized* objective function increase is selected, irrespective of a possible violation of the capacity constraint of knapsack k. The normalization is performed as in Qin et al. [16], i.e., the objective function increase is divided by  $w_i^\beta$  with  $\beta \in (0, 1)$ . Procedure InfRelocB, shown in Algorithm 10, shares part of the structure of InfRelocA, the main difference being that the item is (possibly) moved to a less (or equally) filled knapsack.

The last procedure, **Repair**, shown in Algorithm 11, is only executed if the incumbent solution produced by the previous steps turns out to be infeasible. It tries to iteratively remove an item, exchange a pair of items, or relocate an item with the main objective of reducing the *weight surplus* 

$$\sigma = \sum_{k=1}^{m} \max\{0, W(k) - c\},$$
(14)

where c is the (common) knapsack capacity, breaking ties by the largest profit of the resulting solution. This strategy differs from the one adopted in the repair procedure by Qin et al. [16] which privileges the largest resulting profit. At each iteration, the best move is performed and the execution terminates as soon as a feasible solution is obtained.

### 4.3 Overall postoptimization

The overall postptimization starts by performing a local search on the current solution. It then executes twice the sequence Delete - LocalSearch, for the two removal criteria previously defined. We assume in the following that the incumbent solution  $x^{best}$  is a global variable. The pseudocode is given in Algorithm 5.

### 5 Complete algorithm

In this section, we introduce our complete deterministic algorithmic framework, which combines the matheuristic approach of Section 3, a "weighted" matheuristic variant, and the postoptimization method of Section 4.

The basic idea for an effective implementation of the matheuristic approach is to explore a large number of "high quality" combinations of the item subsets produced by procedure Bundle and to avoid exploring previously generated solutions multiple times. To this end, we define a weighted variant of the CCSPP in which: Algorithm 5: Procedure Postopt

**input** : A feasible solution x to an instance I**output:** A possibly improved solution  $x^{\text{best}}$  to I 1  $x \leftarrow \text{LocalSearch}(x);$ **2** for d := 1 to 2 do del := true;3 while del = true do $\mathbf{4}$  $x^- \leftarrow \texttt{Delete}(x,d);$  $\mathbf{5}$  $x^- \leftarrow \text{LocalSearch}(x^-);$ 6 if  $z(x^{-}) > z(x)$  then  $x := x^{-}$  else del := false  $\mathbf{7}$ end 8 9 end 10 if  $z(x) > z(x^{\text{best}})$  then  $x^{\text{best}} := x$ 11 12 end

- (i) each subset  $B_h \in \mathcal{F}$  is *weighted* by a measure  $g_h$  (h = 1, ..., s) of its "quality" (to be defined later);
- (ii) a *fixed* number r of disjoint subsets of  $\mathcal{F}$  must be selected to fill the knapsacks;
- (iii) a collection  $\mathcal{S}^-$  of subset families is *forbidden*.

The corresponding model is

(CCSPP1) max 
$$\sum_{h=1}^{s} g_h y_h$$
 (15)

s.t. 
$$\sum_{h=1}^{s} a_{hi} y_h \le 1$$
  $(i = 1, \dots, n)$  (16)

$$\sum_{h=1}^{s} y_h = r \tag{17}$$

$$\sum_{h \in S} y_h \le |S| - 1 \qquad (S \in \mathcal{S}^-) \tag{18}$$

$$y_h \in \{0, 1\}$$
  $(h = 1, \dots, s),$  (19)

where points (i)-(iii) above are implemented by equations (15), (17), and (18), respectively.

The overall algorithm, procedure GMRT shown in Algorithm 6,

- starts by generating, through procedure Bundle, a family  $S^*$  of item subsets, and setting  $\overline{r} = |S^*|$ ;
- defines, for each subset  $B_h$  (h = 1, ..., s), two different types of weights:  $g_h^1$ , the cardinality of subset  $B_h$ ,

 $g_h^2$ , the sum of all linear and pairwise profits produced by subset  $B_h$ ;

#### Algorithm 6: Procedure GMRT

```
input : A QMKP instance I
     output: A feasible solution x^{\text{best}} to I
 1 \mathcal{S}^{\star}, \, \tilde{\lambda} \leftarrow \text{Bundle}(I);
 2 \overline{r} := |\mathcal{S}^{\star}|, \ x^{\text{best}} = \emptyset;
                                                                                                                  /* initialize */
 \mathbf{s} for h := 1 to s do
          g_h^1 := |B_h|, \ g_h^2 := \sum_{i \in B_h} p_i + \sum_{i \in B_h} \sum_{j \in B_h : j > i} p_{ij};
                                                                                                                        /* weights */
 4
 5 end
 6 for r := 1 to \overline{r} do
           \mathcal{S}^- := \emptyset, \ \ell := 1 \ ;
                                                                                                                  /* initialize */
 7
           repeat
 8
                 if \ell is odd then
 9
                    g_h := g_h^1 \ (h = 1, \dots, s)
\mathbf{10}
                  else
11
                     g_h := g_h^2 \ (h = 1, \dots, s)
12
                  end
\mathbf{13}
                 solve the CCSPP1 (15)-(19) with an MILP solver;
\mathbf{14}
                  y^{\star} \leftarrow \text{CCSPP1} optimal solution;
15
                  \mathcal{S}^{\star} \leftarrow \{B_h : h \in \{1, \dots, s\}, y_h^{\star} = 1\};
16
                 if \mathcal{S}^{\star} \neq \emptyset then
\mathbf{17}
                        x \leftarrow \operatorname{Pack}(\mathcal{S}^{\star}, \tilde{\lambda});
\mathbf{18}
                        x \leftarrow \mathsf{Postopt}(x);
19
                        if z(x) > z(x^{\text{best}}) then
\mathbf{20}
                             x^{\text{best}} := x;
\mathbf{21}
                              if z(x^{\text{best}}) = L(\tilde{\lambda}) then
\mathbf{22}
                                                                                       /* Optimal solution found */
                                terminate ;
\mathbf{23}
                              end
\mathbf{24}
                        end
\mathbf{25}
                        \mathcal{S}^- := \mathcal{S}^- \cup \mathcal{S}^\star
\mathbf{26}
                  end
\mathbf{27}
                 \ell := \ell + 1
\mathbf{28}
           until \ell = \ell_{\max} or \mathcal{S}^{\star} = \emptyset;
\mathbf{29}
30 end
```

- executes a series of iterations, one for each value  $r = 1, \ldots, \overline{r}$ . Each iteration consists of repeatedly, in turn using weights  $g_h^1$  and  $g_h^2$ ,
  - generating a new QMKP solution from the sequence CCSPP1-Pack, and postoptimizing it through Postopt;
  - adding the new family  $\mathcal{S}^*$  produced by CCSPP1 to the current collection  $\mathcal{S}^-$  of forbidden families.

An iteration ends when either CCSPP1 is no longer able to produce a solution (due to the current  $S^-$ ) or a prefixed maximum number  $\ell_{\text{max}}$  of repetitions is reached:  $S^-$  is then reset to empty, and the next iteration is performed.

The process is terminated if an optimal QMKP solution is found.

Preliminary computational experiments showed that larger instances  $(n \ge 45 \text{ and } m \ge 5)$  can require high computing times for the execution of the quadknap code at each bundle iteration. For this reason, in the final version of the algorithm, we use two time limits within the ProximalBundle code executed by the Bundle procedure invoked by GMRT:

- TLB = time limit for the global execution of the ProximalBundle code. If TLB is reached, the best Lagrangian multiplier vector  $\tilde{\lambda}$  and the best upper bound  $L(\tilde{\lambda})$  found by the bundle code could be improved;
- *TLQ* = time limit for the execution of the **quadknap** code at each iteration of the bundle code. If *TLQ* is reached, the upper bound computed at the current iteration is not a valid upper bound for the current instance.

In addition, it turns out that by executing the GMRT algorithm with different TLB and/or TLQ values, one can obtain different QMKP solutions that are not dominating each other. Therefore, our final solution approach to the QMKP consists of running GMRT twice, with two different pairs of time limits,  $(TLB^1, TLQ^1)$  and  $(TLB^2, TLQ^2)$ , as follows:

- Run 1: during the first run of GMRT, we store in UB1 the best upper bound value (and the corresponding Lagrangian multiplier vector) found at the bundle iterations for which the CPU time of quadknap did not reach time limit  $TLQ^1$ . Note that UB1 represents a valid upper bound for the current instance;
- Run 2: this run of GMRT is only performed when at least one of the two time limits was reached in the first run. We set  $TLB^2$  and  $TLQ^2$  as the new time limits, perform a "warm" start from the best valid upper bound UB1 (and the corresponding Lagrangian multiplier vector) found in **Run 1**, and retain the best QMKP solution found.

# 6 Computational experiments

The deterministic matheuristic algorithm presented in the previous sections was implemented in C++ and computationally evaluated by comparing its results with those produced by the fastest exact algorithm for the QMKP, namely the branch-and-bound algorithm recently proposed by Fleszar [7]. The comparison was performed using small and medium size benchmark instances from the literature. (No computational experiments of other heuristic algorithms on these instances are reported in the literature.) The experiments were performed on a single thread of an  $Intel^{\textcircled{C}}$  Core<sup>TM</sup> i7-8700K running at 3.70GHz with 32 GB RAM.

The following software was used:

- C++ code of the branch-and-bound algorithm by Fleszar [7], available at https: //sites.google.com/view/kfleszar/research;
- C++ code ProximalBundle by Frangioni [8], available at https://gitlab.com/ frangio68/ndosolver\_fioracle\_project;
- C code quadknap, available at http://hjemmesider.diku.dk/~pisinger/codes. html;
- MILP commercial software CPLEX 20.0.

The parameter values used for all our experiments are as follows:

- $-\alpha = 1.02;$
- $-\beta = 0.7;$
- $-\gamma = 0.25;$
- $-\ell_{\rm max} = 300;$

$$-r_{\max} = \begin{cases} \min\{3, \lceil m/2 \rceil\} & \text{if } n \le 35\\ \lceil m/2 \rceil & \text{if } n > 35; \end{cases}$$

- $TLB^1 = 500$  seconds,  $TLQ^1 = 5$  seconds;
- $TLB^2 = 200$  seconds,  $TLQ^2 = 3$  seconds.

The benchmark instances consist of:

- two sets of small-size instances:
  - HJ, based on the generating scheme of Hiley and Julstrom [12]. For each triple (n, m, d) with  $n \in \{20, 25, 30, 35\}$ ,  $m \in \{3, 5, 10\}$ , and  $d \in \{0.25, 0.5, 0.75\}$  (d being the *density*, defined as the fraction of non-zero pairwise profits), five instances were generated, for a total of 180 instances.

The linear and pairwise item profits and the weights are uniformly random integers in (0, 100) and (1, 50), respectively.

The knapsack capacities are set to  $\lfloor 0.8 \sum_{i=1}^{n} w_i / m \rfloor$ ;

- SS, 45 groups of five instances each, based on the generating scheme of Saraç and Sipahioglu [17]. The resulting 225 instances have  $n = 30, m \in \{3, 5, 10\}, d = 0.5$ .

The linear and pairwise profits are randomly generated using a parameter

 $\rho \in \{1, 2, 3\}$  which produces different degrees of correlation with the weights. The item weights are drawn from uniform distributions with five different upper limits depending on a parameter  $\psi \in \{1, 2, 3, 4, 5\}$ . The knapsack capacities are as in **HJ**;

• HJM, one set of medium-size instances generated by Galli et al. [9] according to the HJ scheme, using the instance generator by Bergman [2]. For each triple (n, m, d) with  $n \in \{40, 45, 50, 55, 60\}$ ,  $m \in \{3, 5, 10\}$ , and  $d \in \{0.25, 0.5, 0.75\}$ , five instances were generated, for a total of 225 instances.

Benchmarks HJ and SS by Bergman [2], can be downloaded at the INFORMS address: https://pubsonline.informs.org/doi/suppl/10.1287/ijoc.2018.0840/suppl\_file/ijoc.2018.0840-instances.sm2.zip.

Benchmark HJM by Galli et al. [9] can be downloaded from:

http://or.dei.unibo.it/instances/quadratic-multiple-knapsack-problem-instances.

Tables 1 and 2 provide the results for the small-size instances. The branch-and-bound algorithm by Fleszar [7] was run with a time limit of 3600 seconds. All CPU times are expressed in seconds. Each line refers to a triple (n, m, d) (Table 1) or to a 4-tuple (Table 2) and provides the average values over the corresponding five instances:

- the first group of four columns refers to the Fleszar [7] algorithm: the entries report the best solution value found (Best<sup>F</sup>), the upper bound value (UB), the percentage gap (%Gap<sup>F</sup>) (computed as 100 (UB - Best<sup>F</sup>)/Best<sup>F</sup>), and the CPU time (Time);
- the second group of five columns refers to the "pure" metaheuristic algorithm of Section 3, i.e., before the postoptimization phase: the entries report the best solution value found (Best<sup>M</sup>), the percentage gap (%Gap) (computed as 100 (Best<sup>F</sup> Best<sup>M</sup>)/Best<sup>F</sup>), the CPU time (Time), the upper bound value (UB) computed through the Lagrangian relaxation of Section 2, and the CPU time (T-UB) spent for its computation;
- the third group of three columns refers to the complete algorithm GMRT of Section 5: the entries report the best solution value found (Best<sup>MP</sup>), the percentage gap (%Gap) (computed as 100 (Best<sup>F</sup> - Best<sup>MP</sup>)/Best<sup>F</sup>), and the CPU time (Time) spent in total for its computation.

For these instances, time limits  $TLB^1$  and  $TLQ^1$  were never reached in **Run 1**, so the code never invoked **Run 2**. Tables 1 and 2 show that, on small-size instances, the GMRT algorithm consistently delivers solutions of very good quality in short computing times. In particular, on all **HJ** and **SS** instances, the percentage gap of GMRT is either zero or very close to, and the computing times for  $n \ge 30$  are at least one order of magnitude smaller than those taken by the exact algorithm by Fleszar [7]. The upper bound values produced by GMRT are not very far from the tighter values obtained by such algorithm. A comparison between the percentage gaps of the last two groups of columns reveals the importance of the postoptimization phase, which significantly improves on the matheuristic solution, with a fairly low impact on the overall CPU time.

Instance			Fles	zar			Ma	theuri	stic		Matheu	r.+Po	stopt	
n	m	d	${\tt Best}^{\rm F}$	UB	$ extsf{Gap}^{ extsf{F}}$	Time	${\tt Best}^{\rm M}$	%Gap	Time	UB	T-UB	$\texttt{Best}^{\mathrm{MP}}$	%Gap	Time
20	3	25	2289.20	2289.20	0.00	0.03	2286.00	0.16	0.07	2295.40	0.06	2289.20	0.00	0.09
20	5	25	1916.20	1916.20	0.00	0.07	1911.00	0.23	0.06	1918.20	0.05	1916.20	0.00	0.13
20	10	25	1199.00	1199.00	0.00	0.01	1198.00	0.10	0.05	1199.00	0.04	1199.00	0.00	0.05
20	3	50	3079.80	3079.80	0.00	0.09	3076.60	0.11	0.07	3093.60	0.07	3079.80	0.00	0.10
20	5	50	2296.60	2296.60	0.00	0.21	2268.80	1.16	0.07	2297.00	0.06	2296.60	0.00	0.25
20	10	50	1427.00	1427.00	0.00	0.01	1304.00	7.76	0.05	1433.60	0.04	1427.00	0.00	0.08
20	3	75	3553.80	3553.80	0.00	0.17	3530.60	0.62	0.07	3592.40	0.07	3543.80	0.27	0.12
20	5	75	2640.60	2640.60	0.00	0.16	2633.40	0.28	0.08	2653.80	0.06	2640.60	0.00	0.34
20	10	75	1639.60	1639.60	0.00	0.01	1591.80	2.66	0.05	1639.60	0.04	1639.60	0.00	0.06
	Avg		2226.87	2226.87	0.00	0.08	2200.02	1.45	0.06	2235.84	0.06	2225.76	0.03	0.14
25	3	25	3002.40	3002.40	0.00	0.20	2981.20	0.73	0.14	3014.20	0.13	3000.20	0.08	0.20
25	5	25	2559.00	2559.00	0.00	0.67	2559.00	0.00	0.13	2565.80	0.11	2559.00	0.00	0.32
25	10	25	1781.20	1781.20	0.00	0.08	1771.00	0.60	0.12	1784.40	0.09	1773.80	0.44	0.53
25	3	50	4275.00	4275.00	0.00	1.08	4275.00	0.00	0.18	4294.80	0.17	4275.00	0.00	0.27
25	5	50	3272.80	3272.80	0.00	2.49	3272.80	0.00	0.13	3282.80	0.11	3272.80	0.00	0.33
25	10	50	2093.20	2093.20	0.00	0.14	2087.20	0.29	0.09	2093.20	0.08	2093.20	0.00	0.10
25	3	75	5155.00	5155.00	0.00	3.46	5137.00	0.35	0.26	5178.40	0.25	5155.00	0.00	0.39
25	5	75	3774.80	3774.80	0.00	5.99	3774.80	0.00	0.21	3807.20	0.17	3774.80	0.00	1.26
25	10	75	2420.20	2420.20	0.00	0.16	2408.80	0.47	0.14	2427.00	0.13	2417.40	0.12	2.14
	Avg		3148.18	3148.18	0.00	1.58	3140.76	0.27	0.16	3160.87	0.14	3146.80	0.07	0.61
30	3	25	4194.80	4194.80	0.00	0.73	4161.60	0.79	0.41	4202.80	0.39	4194.80	0.00	0.52
30	5	25	3550.40	3550.40	0.00	6.51	3540.20	0.29	0.26	3557.60	0.22	3550.40	0.00	0.58
30	10	25	2558.40	2558.40	0.00	10.86	2558.40	0.00	0.19	2562.40	0.17	2558.40	0.00	3.34
30	3	50	5858.60	5858.60	0.00	19.09	5829.20	0.49	0.83	5903.20	0.79	5850.80	0.13	1.07
30	5	50	4609.40	4609.40	0.00	21.39	4595.80	0.27	0.28	4611.20	0.27	4609.40	0.00	1.12
30	10	50	2963.20	2963.20	0.00	6.35	2890.80	2.08	0.21	2964.60	0.19	2963.20	0.00	6.61
30	3	75	7210.20	7210.20	0.00	42.83	7207.00	0.04	1.02	7287.80	0.97	7210.20	0.00	1.30
30	5	75	5362.00	5362.00	0.00	30.57	5347.20	0.28	0.35	5375.80	0.26	5362.00	0.00	1.03
30	10	75	3283.40	3283.40	0.00	2.16	3260.40	0.73	0.24	3291.80	0.14	3281.00	0.08	1.02
	Avg		4398.93	4398.93	0.00	15.61	4376.73	0.55	0.42	4417.47	0.38	4397.80	0.02	1.84
35	3	25	5277.80	5277.80	0.00	10.74	5272.40	0.11	2.34	5304.00	2.26	5272.40	0.11	2.75
35	5	25	4463.00	4463.00	0.00	87.35	4416.60	1.06	0.63	4472.80	0.51	4462.80	0.00	1.87
35	10	25	3318.60	3318.60	0.00	45.29	3309.60	0.28	0.51	3323.60	0.27	3317.60	0.03	10.96
35	3	50	7730.60	7730.60	0.00	443.02	7704.00	0.34	3.57	7768.00	3.47	7723.00	0.09	3.87
35	5	50	6000.20	6000.20	0.00	1412.79	5955.60	0.73	0.83	6021.00	0.80	6000.00	0.00	3.02
	10	50		4062.20	0.00	86.65	4028.80	0.84		4068.60	0.26	4041.80	0.50	3.18
35	3	75	9944.40	9944.40	0.00	271.80	9904.60	0.38		9993.40	2.98	9944.40	0.00	3.46
35	5	75		7204.20	0.00	731.79	7189.20	0.20		7226.60	1.03	7202.60	0.02	5.82
	10	75	4482.60	4482.60	0.00	260.15	4449.60	0.73		4498.40	0.40	4479.60	0.07	26.77
	Avg		5831.51		0.00	372.18	5803.38	0.52		5852.93	1.33	5827.13	0.09	6.86
Ov	er. A	vg	3901.37	3901.37	0.00	97.36	3880.22	0.70	0.57	3916.78	0.48	3899.37	0.05	2.36

Table 1:  ${\bf HJ}$  instances. Average values over 5 instances.

Ι	nsta	nce			Fles	zar			Ma	theuri	stic		Matheu	r.+Po	stopt
n	m	$\psi$	$\rho$	${\tt Best}^{ m F}$	UB	$\gamma Gap^F$	Time	${\tt Best}^{\rm M}$	%Gap	Time	UB	T-UB	$\texttt{Best}^{\mathrm{MP}}$	%Gap	Time
30	3	1	1	1610.20	1610.20	0.00	65.83	1609.80	0.03	0.66	1618.00	0.63	1610.20	0.00	1.17
30	3	1	2	1425.60	1425.60	0.00	60.97	1425.00	0.04	0.64	1427.40	0.50	1425.60	0.00	0.89
30	3	1	3	1818.00	1818.00	0.00	135.44	1818.00	0.00	0.98	1821.80	0.82	1818.00	0.00	1.56
30	3	2	1	3530.40	3530.40	0.00	63.08	3529.80	0.02	0.78	3544.00	0.64	3530.40	0.00	1.20
30	3	2	2	3032.00	3032.00	0.00	96.54	3030.00	0.02	0.64	3039.00	0.56	3030.00	0.07	1.16
30	3	$\frac{2}{2}$	$\frac{2}{3}$	3714.00	3714.00	0.00	340.36	3707.60	0.18	1.02	3720.40	0.95	3710.80	0.09	1.83
30	3	3	1	5263.20	5263.20	0.00	159.68	5260.20	0.06	1.02	5289.80	0.98	5263.20	0.00	1.92
30	3	3	2	4188.80	4188.80	0.00	282.75	4188.20	0.00 0.01	0.59	4193.60	0.50 0.54	4188.20	0.00	0.98
30	3	3	$\frac{2}{3}$	5340.80	5340.80	0.00	533.46	5336.00	0.01	0.03 0.77	41 <i>3</i> 5.00 5360.40	0.64	5338.80	0.01	1.65
30	3	4	1	6971.80	6971.80	0.00	69.28	6971.80	0.09	0.98	6990.80	0.08	6971.80	0.00	1.63
$30 \\ 30$	3	4	$\frac{1}{2}$	5964.40		0.00	628.57	5960.60	0.00 0.07	0.98	5975.80	0.90 0.72	5963.60	0.00	$1.08 \\ 1.37$
	3 3	4	$\frac{2}{3}$		5964.40									0.01	
30	3 3			7120.80	7120.80	0.00	436.66	7120.40	0.01	0.74	7143.00	0.69	7120.80		1.35
30		5	1	9039.00	9039.00	0.00	62.49	9033.60	0.06	1.02	9059.20	0.91	9037.00	0.02	1.79
$\begin{array}{c} 30\\ 30 \end{array}$	$\frac{3}{3}$	$\frac{5}{5}$	$\frac{2}{3}$	6566.80	6566.80	0.00	229.49	6565.20	0.02	0.50	6570.20 8274.20	0.46	6565.40 8235.20	0.02	0.65
30			э	8242.80	8242.80	0.00	240.64	8234.80	0.10	0.82		0.73		0.09	1.62
20	Av	-	1	4921.91	4921.91	0.00	227.02	4919.40	0.05	0.80	4935.17	0.71	4920.60	0.02	1.39
$\begin{vmatrix} 30 \\ 20 \end{vmatrix}$	5	1	1	1509.60	1509.60	0.00	354.06	1505.80	0.26	0.46	1514.40	0.41	1509.40	0.01	4.56
30	5	1	2	1382.00	1382.00	0.00	159.25	1381.20	0.05	0.57	1384.20	0.22	1382.00	0.00	1.75
30	5	1	3	1740.00	1740.00	0.00	805.66	1739.60	0.02	2.07	1743.60	0.45	1740.00	0.00	11.05
30	5	2	1	3342.60	3342.60	0.00	637.74	3339.60	0.09	1.92	3355.40	0.39	3340.20	0.07	6.42
30	5	2	2	2933.00	2933.00	0.00	345.75	2929.40	0.13	0.82	2941.00	0.26	2932.80	0.01	3.19
30	5	2	3	3511.60	3511.60	0.00	723.43	3511.20	0.01	0.58	3520.40	0.32	3511.20	0.01	5.92
30	5	3	1	4984.80	4984.80	0.00	594.26	4983.60	0.02	0.51	4996.20	0.41	4983.60	0.02	3.60
30	5	3	2	4041.40	4041.40	0.00	901.17	4040.60	0.02	0.52	4048.40	0.24	4040.60	0.02	2.79
30	5	3	3	5036.80	5036.80	0.00	1877.67	5033.60	0.06	1.18	5048.20	0.37	5036.80	0.00	8.09
30	5	4	1	6546.60	6546.60	0.00	758.69	6530.60	0.23	0.93	6569.80	0.42	6540.80	0.08	6.08
30	5	4	2	5726.60	5736.20	0.16	2127.03	5719.40	0.13	0.95	5741.60	0.28	5720.60	0.11	4.62
30	5	4	3	6704.00	6704.80	0.01	1905.60	6703.20	0.01	0.86	6726.20	0.34	6703.20	0.01	9.43
30	5	5	1	8517.00	8517.00	0.00	1209.14	8508.60	0.11	0.95	8549.80	0.47	8513.60	0.04	7.24
30	5	5	2	6274.80	6287.00	0.20	2129.79	6274.20	0.01	0.49	6295.40	0.24	6274.20	0.01	3.48
30	5	5	3	7762.80	7762.80	0.00	1431.44	7756.40	0.07	1.58	7794.00	0.35	7757.20	0.06	9.38
	Av	g		4667.57	4669.08	0.02	1064.05	4663.80	0.08	0.96	4681.91	0.34	4665.75	0.03	5.84
30	10	1	1	1366.40	1366.40	0.00	286.16	1357.80	0.63	0.80	1369.80	0.25	1365.60	0.05	6.48
30	10	1	2	1290.20	1290.20	0.00	4.49	1288.40	0.14	0.49	1291.00	0.18	1289.80	0.03	4.09
30	10	1	3	1640.00	1640.00	0.00	96.81	1636.80	0.18	1.67	1645.80	0.20	1638.00	0.11	11.80
30	10	2	1	3069.40		0.00	649.28	3063.00	0.21		3071.80	0.22	3068.80	0.02	9.10
30	10	2	2	2746.80	2746.80	0.00	17.91	2746.20	0.02	0.56	2753.80	0.15	2746.20	0.02	8.92
30	10	2	3	3265.60	3265.60	0.00	74.16	3265.60	0.00	0.59	3276.40	0.20	3265.60	0.00	18.75
30	10	3	1	4575.40	4575.40	0.00	187.47	4572.60	0.07	0.46	4582.40	0.29	4575.20	0.00	15.82
30	10	3	2	3776.40	3776.40	0.00	8.13	3774.00	0.06	0.23	3778.60	0.12	3774.60	0.04	4.96
30	10	3	3	4678.00	4678.00	0.00	439.50	4673.20	0.12	0.53	4685.80	0.21	4673.20	0.12	18.63
30	10	4	1	5989.80	5989.80	0.00	509.02	5981.20	0.13	0.28	5994.00	0.22	5989.80	0.00	4.80
30	10	4	2	5319.20	5319.20	0.00	99.11	5316.60	0.05	0.63	5322.20	0.16	5319.20	0.00	12.26
30	10	4	3	6259.20	6259.20	0.00	310.73	6257.60	0.02	0.17	6267.20	0.14	6259.20	0.00	9.59
30	10	5	1	7863.00	7882.00	0.21	902.33	7840.20	0.29	0.87	7883.20	0.27	7860.40	0.03	12.25
30	10	5	2	5866.80	5866.80	0.00	45.38	5866.80	0.00	0.13	5868.60	0.10	5866.80	0.00	2.33
30		5	3		7151.20	0.00	40.58	7127.20	0.30	0.29	7158.00	0.16	7151.20	0.00	14.19
	Av	g		4323.83	4325.09	0.01	244.74	4317.81	0.15	0.56	4329.91	0.19	4322.91	0.03	10.27
0	ver.		g	4637.77	4638.69	0.01	511.93	4633.67	0.09	0.77	4649.00	0.42	4636.42	0.03	5.83

Table 2: SS instances. Average values over 5 instances.

Table 3 provides the results for the medium-size instances **HJM**. In this case too, the algorithm by Fleszar [7] was run with a time limit of 3600 seconds. Each line refers to a triple (n, m, d) and provides the average values over the corresponding five instances:

- the first group of four columns refers to the Fleszar [7] algorithm, and reports the same information as in the previous tables;
- the second group of five columns refers to Run 1. The entries report the best upper bound value UB found and the CPU time T-UB needed to obtain it, the best solution value found (Best<sup>MP</sup>), the percentage gap (%Gap), computed as 100 (Best<sup>F</sup> Best<sup>MP</sup>)/Best<sup>F</sup> and the related CPU time (Time);
- the third group of two columns refers to **Run 2**, and reports the best upper bound value UB found by the two runs and the CPU time T-UB of **Run 2**;
- the fourth and fifth group of three columns each summarize the overall results (best solution value, percentage gap and CPU time) produced by the matheuristic approach before and after postoptimization, respectively.

The results in Table 3 show that on larger HJ instances Fleszar's exact method consistently hits the time limit of 3600 seconds, with few exceptions, mostly on the smaller (n = 40) instances. For such cases, columns  $Best^F$  and UB in the first group provide the best solution and upper bound values, respectively, computed by the branch-and-bound algorithm before the time limit. Already **Run 1** of GMRT (second group of columns) regularly provides solutions of better quality with computing times that are one or two orders of magnitude smaller. The next group of two columns is only reported to examine the effect of **Run 2** on the upper bound computations: we can observe that: (i) the second run is only executed for relatively large values of n; (ii) a moderate increase of computing time occasionally allows to improve the upper bound values.

The overall performance of the proposed approach is summarized in the last two groups of columns, before and after the postoptimization phase, respectively. The results show that: (i) *before* postoptimization the best solutions found are worse than those obtained by Fleszar's algorithm; (ii) *after* postoptimization, the overall algorithm favorably compares with Fleszar's algorithm (negative percentage gap values), providing good solutions in CPU times of few hundred seconds; (iii) the execution of **Run 2** further improves the results obtained after the execution of **Run 1**, with a moderate increase of CPU time.

### 7 Conclusions

Recent literature on the QMKP shows an increasing attention by different authors to develop exact methods for this challenging combinatorial optimization problem. We propose the first deterministic Lagrangian matheuristic for solving the QMKP approximately. Our method relies on the strength of the Lagrangian relaxation for the natural quadratic formulation of the QMKP, which can be exploited to generate "high quality" subsets of

0.21 0.07 0.05 0.13 0.13 0.13 0.13		$\begin{array}{c ccccccccccccccccccccccccccccccccccc$
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$\begin{array}{rrrr} 1.96 & 2886.33 \\ 4.20 & 2996.54 \end{array}$	3100.39 3600.42 3600.59 3600.59 3600.53 3600.53 3600.17 3594.76 3594.76 3590.17 3600.51 3600.51 3600.63 3600.64 3600.63 3600.64 3600.63 3600.63 3600.63 3600.64 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.63 3600.53 3600.53 3600.53 3600.53 3600.53 3600.53 3600.53 3600.53 3600.53 3600.53 3600.53 3600.50 3700.50 3700.5	$\begin{array}{c} 3100.39\\ 3600.42\\ 3600.59\\ 3600.53\\ 3600.53\\ 3600.53\\ 3600.53\\ 3600.61\\ 3544.76\\ 35600.58\\ 3600.61\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.61\\ 3600.61\\ 3600.63\\ 3600.63\\ 3600.61\\ 3600.61\\ 3600.63\\ 3600.63\\ 3600.61\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.63\\ 3600.61\\ 3600.63\\ 3600$
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0 0 4 7 7		

items to be packed into the knapsacks. Experimental studies show that the method is competitive with the currently best performing exact algorithm by Fleszar [7] on small **HJ** instances with  $n \ge 30$ , and on all **SS** instances, consistently delivering solutions of very good quality in much shorter computing times. On the medium-size **HJM** instances, the proposed method outperforms the algorithm by Fleszar [7] both in terms of solution quality and computing time. The results show that the QMKP is very difficult to solve, especially when n and/or m grow. Directions for future research could thus be the development of faster exact algorithms to solve large problem instances. In particular, we believe that our Lagrangian matheuristic could be embedded within an exact framework to quickly generate high quality solutions and upper bounds.

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# **Appendix: Local Search procedures**

```
Algorithm 7: Procedure BestExchange
    input : A feasible QMKP solution x
    output: A possibly improved QMKP solution x'
  1 Stop := false;
  2 while Stop = false do
        Stop := true;
  3
        for i := 1 to n - 1 do
  4
            \Delta_i := 0;
  \mathbf{5}
           for j := i + 1 to n do
  6
               if K(j) \neq K(i) and the i - j exchange is feasible and profitable
  \mathbf{7}
               then \Delta_i := \Delta, j^* := j;
  8
           end
  9
           if \Delta_i > 0 then
 \mathbf{10}
               exchange items i and j^* and update the relative variables;
 11
               Stop := false;
 12
           end
 13
        end
 \mathbf{14}
15 end
```

Algorithm 8: Procedure BestReloc

**input** : A feasible QMKP solution x**output:** A possibly improved QMKP solution x'1 Stop := false;2 while Stop = false doStop := true;3 for i := 1 to n do  $\mathbf{4}$  $\Delta_i := 0;$  $\mathbf{5}$ for each  $k \in \{1, \ldots, m\} \setminus \{K(i)\}$  do 6 7 if the relocation of item i to knapsack k is *feasible* and *profitable* then  $\Delta_i := \Delta, k^* := k;$ 8 end 9 if  $\Delta_i > 0$  then  $\mathbf{10}$ **relocate** item i to knapsack  $k^*$  and update the relative variables; 11 Stop := false; $\mathbf{12}$ end  $\mathbf{13}$ end  $\mathbf{14}$ 15 end

#### Algorithm 9: Procedure InfRelocA

**input** : A feasible or infeasible QMKP solution x', a knapsack index  $k \in \{1, \ldots, m\}$ output: A possibly improved (feasible or infeasible) QMKP solution  $\overline{x}$ 1  $\overline{x} := x'$ , Maxqain := 0; 2 for  $\kappa := 1$  to m do if  $\kappa \neq k$  and  $W(\kappa) \geq W(k)$  then 3 foreach  $i \in X(\kappa)$  do 4  $\Delta^{\kappa} := \sum_{\ell \in X(\kappa)} p_{i\ell}, \quad \Delta^k := \sum_{\ell \in X(k)} p_{i\ell}, \quad gain := \frac{\Delta^k - \Delta^{\kappa}}{w_i^{\beta}};$  $\mathbf{5}$ if gain > Maxgain then  $Maxgain := gain, i^* := i$ 6 end  $\mathbf{7}$ end 8 9 end 10 if Maxgain > 0 then **relocate** item  $i^*$  to knapsack k and define the relative solution  $\overline{x}$ 11 12 end

#### Algorithm 10: Procedure InfRelocB

**input** : A feasible or infeasible QMKP solution x', a knapsack index  $k \in \{1, \ldots, m\}$ **output:** A possibly improved (feasible or infeasible) QMKP solution  $\overline{x}$ 1  $\overline{x} := x'$ , Maxgain := 0; 2 for each  $i \in X(k)$  do  $\Delta^k := \sum_{\ell \in X(k)} p_{i\ell};$ 3 for  $\kappa := 1$  to m do  $\mathbf{4}$ if  $\kappa \neq k$  and  $W(\kappa) < W(k)$  then  $\mathbf{5}$  $\Delta^{\kappa} := \sum_{\ell \in X(\kappa)} p_{i\ell}, \quad gain := \frac{\Delta^{\kappa} - \Delta^k}{w_i^{\beta}};$ 6  $\textbf{if } gain > \textit{Maxgain } \textbf{then } \textit{Maxgain} := gain, \, i^* := i, \, k^* := \kappa \\$  $\mathbf{7}$ end 8 end 9 10 end 11 if Maxqain > 0 then **relocate** item  $i^*$  to knapsack  $k^*$  and define the relative solution  $\overline{x}$  $\mathbf{12}$ 13 end

Algorithm 11: Procedure Repair **input** : An infeasible QMKP solution x'**output:** A possibly worse feasible QMKP solution  $\overline{x}$ 1 while x' is infeasible do  $\sigma^* := +\infty, \ \Delta^* := 0;$ 2 foreach  $i \in \{1, \ldots, n\} \setminus X(m+1)$  do 3 compute surplus  $\sigma$  and profit variation  $\Delta$  produced by removing *i*; 4 if  $(\sigma < \sigma^*)$  or  $(\sigma = \sigma^* \text{ and } \Delta > \Delta^*)$ 5 then  $\sigma^* := \sigma$ ,  $\Delta^* := \Delta$ , move := remove(i); 6 end 7 foreach item pair (i, j) with  $i, j \in \{1, ..., n\}, j > i$  and  $K(j) \neq K(i)$  do 8 compute surplus  $\sigma$  and profit variation  $\Delta$  produced by exchanging *i* and *j*; 9 if  $(\sigma < \sigma^*)$  or  $(\sigma = \sigma^* \text{ and } \Delta > \Delta^*)$ 10 then  $\sigma^* := \sigma$ ,  $\Delta^* := \Delta$ , move := exchange(i, j);  $\mathbf{11}$ end 12**foreach** item-knapsack pair (i, k) with  $i \in \{1, \ldots, n\} \setminus X(m+1)$  $\mathbf{13}$ and  $k \in \{1, \ldots, m\} \setminus K(i)$ do  $\mathbf{14}$ compute surplus  $\sigma$  and profit variation  $\Delta$  produced by relocating *i* into *k*;  $\mathbf{15}$ if  $(\sigma < \sigma^*)$  or  $(\sigma = \sigma^* \text{ and } \Delta > \Delta^*)$ 16 then  $\sigma^* := \sigma$ ,  $\Delta^* := \Delta$ , move := relocate(i, k);  $\mathbf{17}$ end  $\mathbf{18}$ update solution x' by executing move 19 20 end 21  $\overline{x} := x'$