# **Competitive Memetic Algorithms for Arc Routing Problems**

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**Abstract.** The Capacitated Arc Routing Problem or CARP arises in applications like waste collection or winter gritting. Metaheuristics are tools of choice for solving large instances of this NP-hard problem. The paper presents basic components that can be combined into powerful memetic algorithms (MAs) for solving an extended version of the CARP (ECARP). The best resulting MA outperforms all known heuristics on three sets of benchmark files containing in total 81 instances with up to 140 nodes and 190 edges. In particular, one open instance is broken by reaching a tight lower bound designed by Belenguer and Benavent, 26 best-known solutions are improved, and all other best-known solutions are retrieved.

Keywords: Capacitated Arc Routing Problem, CARP, metaheuristic, memetic algorithm

#### 1. Introduction

Contrary to the well-known *Vehicle Routing Problem* (VRP), in which goods must be delivered to client nodes in a network, the *Capacitated Arc Routing Problem* (CARP) consists of visiting a subset of edges. CARP applications include for instance urban waste collection, winter gritting and inspection of power lines. From now on, to make the paper more concrete without loss of generality, examples are inspired by municipal refuse collection.

The basic CARP of literature tackles undirected networks. Each edge models a two-way street whose both sides are treated in parallel and in any direction (*zigzag collection*), a common practice in residential areas with narrow streets. A fleet of identical vehicles of limited capacity is based at a depot node. Each edge can be traversed any number of times, with a known traversal cost. Some edges are required, i.e., they have a non-zero demand (amount of waste) to be collected by a vehicle. The CARP consists of determining a set of vehicle trips of minimum total cost, such that each trip starts and ends at the depot, each required edge is serviced by one single trip, and the total demand processed by a trip fits vehicle capacity.

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The CARP is NP-hard, even in the single-vehicle case called Rural Postman Problem (RPP). Since exact methods are still limited to 20–30 edges (Hirabayashi, Saruwatari, and Nishida, 1992), heuristics are required for solving large instances, e.g., Augment-Merge (Golden and Wong, 1981), Path-Scanning (Golden, DeArmon, and Baker, 1983), Construct-and-Strike (Pearn's improved version, 1989), Augment-Insert (Pearn, 1991) and Ulusoy's tour splitting algorithm (Ulusoy, 1985).

The first metaheuristic for the CARP, a simulated annealing procedure, was designed by Eglese (1994) for solving a winter gritting problem. Several tabu search (TS) algorithms are also available, both for particular cases like the undirected RPP (Hertz, Laporte, and Nanchen-Hugo, 1999) or the mixed RPP (Corberán, Martí, and Romero, 2000) and for the CARP itself (Eglese and Li, 1996; Hertz, Laporte, and Mittaz, 2000). All these metaheuristics and classical heuristics may be evaluated thanks to lower bounds, generally based on linear programming formulations, see (Benavent, Campos, and Corberán, 1992; Belenguer and Benavent, 1998; Amberg and Voß, 2002). On most instances, the best-known lower bound is obtained by a cutting-plane algorithm (Belenguer and Benavent, 2003).

Compared to the VRP, the CARP has been relatively neglected for a long time but it attracts more and more researchers: successful applications are reported (Mourão and Almeida, 2000) and extensions are now investigated, for instance the directed RPP with turn penalties (Benavent and Soler, 1999), the multi-depot CARP (Amberg, Domschke, and Voß, 2000) and the CARP with intermediate facilities (Ghianni, Improta, and Laporte, 2001).

This paper presents powerful memetic algorithms (MA) for an extended CARP. Compared to an earlier genetic algorithm (GA) for the mixed CARP with forbidden turns (Lacomme, Prins, and Ramdane-Chérif, 2001), they handle other objectives, like the makespan or the number of vehicles used, and extensions like parallel arcs, turn penalties, a maximum trip length and a limited fleet. Several possible bricks for each MA step are designed and tested, e.g., a generational approach and a partial replacement procedure. The best resulting MA is twice faster, it improves 26 best-known solutions and tackles larger instances with 140 nodes and 190 edges.

The extended problem (ECARP) is presented in section 2. Three classical constructive heuristics are extended to the ECARP in section 3 to provide good initial solutions. Section 4 describes possible components for each step of memetic algorithms. Section 5 is devoted to computational evaluations: the best MA structure is defined after a preliminary testing and results are reported for three sets of benchmark instances.

## 2. Extended CARP model (ECARP)

## 2.1. Extensions considered and street modelling

For the sake of clarity, this subsection presents without mathematical symbols our extended problem and the modelling technique for the streets of a real network. Subsections 2.2–2.4 are respectively devoted to the required notation, to some complications

raised by forbidden turns, and to the representation of solutions. The ECARP tackles the following extensions:

- (a) mixed multigraph with two kinds of links (edges and arcs) and parallel links,
- (b) two distinct costs per link (deadheading and collecting),
- (c) prohibited turns (e.g., U-turns) and turn penalties (e.g., to penalize left turns),
- (d) maximum trip length (an upper limit on the cost of any trip).

Like in the basic CARP, the depot is unique, the fleet is homogeneous, no demand exceeds vehicle capacity, and no split collection is allowed. The number of vehicles is a decision variable. To get feasible solutions, the maximum trip length allows a vehicle to reach any required link, collect it, and return to the depot. The cost of a trip comprises collecting costs (for each link collected) and deadheading costs (for each link traversed without collection). The goal is to find a set of trips of minimum total cost, covering all required links.

A mixed graph allows to model two kinds of non-required streets and three kinds of required streets. A non-required street is modelled either as one arc (1-way street) or two opposite arcs (2-way streets). A required street can be: (i) a 2-way street with zigzag collection (giving one edge, like in the basic CARP), (ii) a 2-way street with sides collected separately (giving two opposite arcs), and (iii) a 1-way street (modelled as one arc). We use a *mixed multigraph* to handle more complicated cases: for instance, two parallel arcs can model a one-way street too wide for zigzag collection and requiring two traversals, one for each side.

## 2.2. Graph transformation and notations

To ease algorithmic design, the mixed multigraph  $\Gamma$  is transformed into a fully directed graph in which each edge is replaced by two arcs with opposite directions. Only one of these arcs must be collected in any feasible solution. To ensure this, both arcs are linked by a pointer variable: when an algorithm selects one direction, both arcs can be marked "collected."

More precisely, the mixed multigraph is coded as a *fully directed graph*, G = (N, A). N is a set of n nodes including a depot node s with identical vehicles of capacity W. A is a set of m arcs identified by indexes from 1 to m instead of pairs of nodes, to avoid ambiguities for parallel arcs. Each arc  $u \in A$  begins at node b(u), ends at node e(u) and has a deadheading cost c(u).

The  $\tau$  required links of  $\Gamma$  (also called *tasks*) comprise  $\varepsilon$  required edges or *edge-tasks* and  $\alpha$  required arcs or *arc-tasks*. They correspond in A to a subset R with  $\rho = 2\varepsilon + \alpha$  arcs. Each arc  $u \in R$  has a demand q(u), a collecting cost w(u) and a pointer inv(u). All costs and demands are non-negative integers. Each arc-task of  $\Gamma$  is coded in R by one arc u with inv(u) = 0, while each edge-task gives two opposite arcs u and v, such that inv(u) = v, inv(v) = u, q(u) = q(v), c(u) = c(v) and w(u) = w(v). To simplify, in the sequel, an arc in R is also called *task*, knowing that u and inv(u) stands

Mixed multigraph $\Gamma$	Ι	Data for each arc u	ľ	Miscellaneous
$\tau = \varepsilon + \alpha$ no. of tasks in $\Gamma$ $\varepsilon$ , $\alpha$ no. of edge-tasks, no of arc-tasks $G = (N, A)$ directed encoding of $\Gamma$ $n$ no. of nodes in $N$ $m$ no. of arcs in $A$ $R \subseteq A$ subset of arcs coding the tasks $\rho = 2 \cdot \varepsilon + \alpha$ no. of arcs in $R$	` '	begin node end node demand deadheading cost service cost pointer to opposite arc set of successor-arcs	s K W L pen(u, v) D P	depot node fleet size (variable) vehicle capacity maximum trip length penalty for turn $(u, v)$ $m \times m$ distance matrix $m \times m$ predecessor matrix

Table 1 Glossary of mathematical symbols.

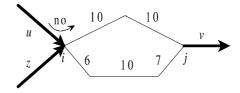


Figure 1. Notion of feasible deadheading path (see text).

for the same task (a required edge) if  $inv(u) \neq 0$ . Table 1 provides a quick reference for the remainder of the paper. From now on, having built the directed encoding of the mixed multigraph, all related algorithms work in that directed equivalence.

## 2.3. Forbidden turns, turn penalties and distance matrix

This subsection shows how to make forbidden turns transparent. In figure 1, the shortest path from node i to node j is the upper one (cost 20). However, the best path for a vehicle coming from arc u is the lower one (cost 23), because it is forbidden to turn left after u.

Such ambiguities can be avoided by considering paths between two arcs. Let suc(u) be the set of allowed successor-arcs for arc u, i.e.  $v \in suc(u)$  if e(u) = b(v) and the turn (u, v) is allowed. Given two arcs u and v, we define a *feasible deadheading path* from u to v as a sequence of arcs  $\mu = (u = u_1, u_2, \ldots, u_k = v)$ , such that  $u_{i+1} \in suc(u_i)$  for  $i = 1, \ldots, k-1$ . Its deadheading cost  $c(\mu)$  is defined by equation (1). Note that the costs of u and v are not included.

$$c(\mu) = pen(u, u_2) + \sum_{i=2}^{k-1} (c(u_i) + pen(u_i, u_{i+1})).$$
 (1)

Dijkstra's algorithm (e.g., see Cormen, Leiserson, and Rivest, 1990) can be adapted to pre-compute a shortest feasible path between all pairs of arcs, in two  $m \times m$  matrices D and P. D(u, v) is the cost of the shortest path found from arc u to arc v and P(u, v) is the predecessor of v on this path. For instance, in figure 1, D(u, v) = 23 and D(z, v) = 20.

```
for v := 1 to m do D(u,v) := \infty; fix(v) := false endfor for each v in suc(u) do D(u,v) := pen(u,v); P(u,v) := u endfor for count := 1 to m do v := argmin\{D(u,z):fix(z)=false\} fix(v) := true for each z in suc(v) with D(u,v) + c(v) + pen(v,z) < D(u,z) do D(u,z) := D(u,v) + c(v) + pen(v,z) P(u,z) := v endfor endfor
```

**Algorithm 1.** Algorithm for shortest paths from one given arc u to all other arcs.

Paths from/to the depot s are handled by putting in A one fictitious loop  $\sigma$  with  $b(\sigma) = e(\sigma) = s$ .

Algorithm 1 computes row u of D and P. It must be called m times with u = 1, 2, ..., m to fill D and P. An arc v is said fixed when a shortest path from u to v is obtained. Initially, no arc is fixed and all paths from u have an infinite cost. Each iteration of the third for loop determines the destination arc v with the smallest path cost, among the arcs not yet fixed. This arc is fixed and each successor-arc z is checked to see if the provisional path from u to z can be improved.

The algorithm runs in  $O(m^2)$ . A heap data structure (Cormen, Leiserson, and Rivest, 1990) allows an  $O(h \log m)$  version, with h the total number of allowed turns in G. So, D and P can be computed in  $O(mh \log m)$  by calling the algorithm m times. For real street networks with  $m \approx 4n$  and  $h \approx 4m \approx 16n$ , D and P are computed very quickly, in  $O(n^2 \log n)$ .

#### 2.4. Implementation of trips and solutions

A solution T is a list  $(T_1, \ldots, T_K)$  of K vehicle trips (K is a decision variable). Each trip  $T_i$   $(i = 1, 2, \ldots, K)$  is a list of tasks  $(T_{i1}, T_{i2}, \ldots, T_{i,|T_i|})$ , with a total demand  $load(T_i) \leq W$  and a total cost  $cost(T_i) \leq L$  defined by equations (2)–(3). Implicitly,  $T_i$  starts and ends at the depot and shortest feasible paths are assumed between two tasks and between one task and the depot loop  $\sigma$  (cf. 2.3). The total cost of a solution T is the sum of its trip costs. Each arc-task appears once in T and each edge-task occurs as one of its two opposite arcs. So, T requires a space proportional to the number of tasks  $\tau$ .

$$load(T_i) = \sum_{j=1}^{|T_i|} q(T_{ij}),$$
 (2)

$$cost(T_i) = D(\sigma, T_{i1}) + \sum_{j=1}^{|T_i|-1} (w(T_{ij}) + D(T_{ij}, T_{i,j+1})) + D(T_{i,|T_i|}, \sigma).$$
 (3)

#### 3. Constructive heuristics for the ECARP

This section extends three classical CARP heuristics to the ECARP: Path-Scanning (Golden, DeArmon, and Baker, 1983), Augment-Merge (Golden and Wong, 1981) and Ulusoy's heuristic (Ulusoy, 1985). The extended versions are used in 4.5 to initialize the population of our MA. The main difference with classical versions is to use D, the arc-to-arc distance matrix of 2.3, to handle forbidden turns.

## 3.1. Extended Path-Scanning (EPS)

This heuristic builds one trip at a time. In constructing each trip, the sequence of tasks is extended by joining the task looking most promising, until capacity W or maximum trip length L are exhausted. For a sequence ending at a task u, the extension step determines the set M of tasks closest to u, not yet collected, and feasible for W and L. Five rules are used to select the next task v in M: (1) maximize the distance  $D(v, \sigma)$  to the depot loop  $\sigma$  (cf. 2.3), (2) minimize  $D(v, \sigma)$ , (3) maximize the yield q(v)/w(v), (4) minimize this yield, (5) use rule (1) if the vehicle is less than half-full, else use rule (2).

Once selected, v must be flagged as "collected," to avoid reselection in subsequent iterations. If v belongs to an edge-task, inv(v) must be flagged too. EPS builds one solution per criterion and returns the best one. It can be implemented in  $O(\tau^2)$ , i.e.,  $O(n^2)$  for a real street network with  $\tau \le \rho \le m \approx 4n$ . In spite of its great simplicity, EPS gives good results in practice, thanks to compensation effects among criteria: the five solutions are never simultaneously bad.

#### 3.2. Extended Augment-Merge (EAM)

The original version is illustrated in figure 2.  $\tau$  trips are built (one per task) and sorted in decreasing cost order. For each trip  $T_i$  ( $i=1,2,\ldots,\tau-1$ ), the *augment phase* scans each smaller trip  $T_j$  ( $j=i+1,i+2,\ldots,\tau$ ). If the unique task u of  $T_j$  is on a deadheading path of  $T_i$  and if  $load(T_i) + q(u) \leq W$ ,  $T_j$  is absorbed. The cost of  $T_i$  does not vary because deadheading and service costs are equal in the basic CARP. However, the total cost decreases by  $cost(T_i)$ . Then, the *merge phase* evaluates the concatenation

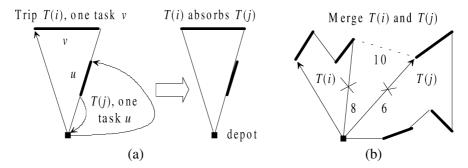


Figure 2. Principle of *augment* (a) and *merge* (b). Thick lines correspond to edge-tasks, thin lines to shortest deadheading paths.

of any two trips, subject to W: e.g, in the figure, concatenating  $T_i$  then  $T_j$  yields a saving of 4. *Merge* concatenates the two trips with the largest positive saving. The process is repeated until no such concatenation is possible.

In the ECARP, each task u has two distinct costs c(u) and w(u). In augment, if trip  $T_i$  absorbs trip  $T_j$  with its task u, the total saving is now  $cost(T_j) + c(u) - w(u)$  and is not always positive like in the basic CARP. In fact, some testing shows that augment can be suppressed without affecting average solution costs. So, we actually removed it. Moreover, matrix D is generally asymmetric for mixed networks and a trip is no longer equivalent to its mirror trip obtained by inverting the sequence of tasks. This gives up to 8 ways of concatenating two trips  $T_i$  and  $T_j$ :  $T_i$  then  $T_j$  or  $T_j$  then  $T_i$ , with each trip inverted or not. Note that a trip cannot be inverted if it contains arc-tasks, non-invertible. The extended heuristic EAM can be implemented in  $O(\tau^2 \log \tau)$ , i.e.  $O(n^2 \log n)$  for real street networks.

## 3.3. Extended Ulusoy's heuristic (EUH)

The original heuristic for the basic CARP temporarily relaxes vehicle capacity W to compute a least-cost giant tour S covering the  $\tau$  edge-tasks. If all edges are required, this sub-problem is an easy undirected Chinese postman problem. If not, it is an NP-hard rural postman problem that can be solved heuristically. Then, this tour is optimally split into capacity-feasible trips.

Figure 3 depicts the splitting procedure (*Split* in the sequel) for a giant tour S = (a, b, c, d, e) with demands in brackets and deadheading costs, assuming W = 9. *Split* builds an auxiliary graph H with  $\tau + 1$  nodes indexed from 0 onward. Each subsequence  $(S_i, \ldots, S_j)$  corresponding to a feasible trip is modeled by one arc (i - 1, j), weighted by the trip cost. A shortest path from node 0 to node  $\tau$  in H (bold) indicates the optimal

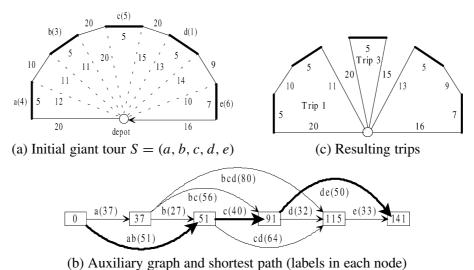


Figure 3. Principle of Split.

splitting: 3 trips and a total cost 141. Note that H is an artificial construction standing in no relation with the CARP graph G.

In the ECARP version EUH, W but also the maximum trip cost L are relaxed to compute a good giant tour S in a mixed multigraph with forbidden turns and turn penalties, modelled by the directed multigraph G. We solve this mixed rural postman problem approximately, by running EPS (cf. 3.1) with a big value of W and L. For better results, we keep the 5 tours obtained by the 5 criteria of EPS, split them, and return the best solution. Split computes the load and cost of  $(S_i, \ldots, S_j)$  using equations (2) and (3) and creates (i-1, j) only if W and L are respected. Forbidden turns are entirely hidden in the arc-to-arc matrix D used in equation (3).

We now analyze complexity, missing in Ulusoy's paper. Path-Scanning (cf. 3.1) returns an initial giant tour in  $O(\tau^2)$ . Then, by construction, H is topologically sorted and contains  $O(\tau^2)$  arcs. Bellman's algorithm (Cormen, Leiserson, and Rivest, 1990) can compute the shortest path in  $O(\tau^2)$ . The global complexity is then  $O(\tau^2)$ , i.e.  $O(n^2)$  for a real street network with  $\tau \le \rho \le m \approx 4n$ . If the minimal demand  $q_{\min}$  is large enough, a trip contains at most  $\omega = \lfloor W/q_{\min} \rfloor$  tasks, H contains  $O(\omega \tau)$  arcs and Split becomes faster, in  $O(\omega \tau)$ .

# 4. Components for memetic algorithms

This section describes the main features of our memetic algorithms: chromosome structure, chromosome evaluation, crossover operators, mutation by local search, population structure and initialization, population management. It describes several possible implementations for certain features. No computational evaluation is performed here: the best assembly of components is determined in section 5.

#### 4.1. Chromosomes: representation, evaluation and generation

Most genetic algorithms for routing problems use quasi-direct representations of solutions, as sequences of tasks. A natural idea for the multi-vehicle case is to use subchromosomes (one per trip), separated by special symbols called *trip delimiters*. In that case, crossovers generally require a repair operator because children may contain overloaded trips. This technique is used for instance by Potvin and Bengio (1996) for the VRP with Time Windows. In our MAs, a chromosome S simply is a sequence of  $\tau$  tasks, *without trip delimiters*, and with implicit shortest paths between consecutive tasks (see figure 4, presented later).

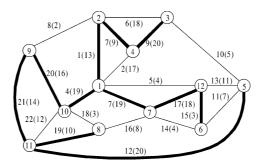
Clearly, S does not directly represent an ECARP solution but can be viewed as a giant trip ignoring capacity W and maximum trip cost L. The Split procedure to be described for Ulusoy's heuristic (cf. 3.3) is applied to S to get an ECARP solution. The fitness F(S) of S is the total cost of this solution. Two good properties hold in our case: (1) chromosomes are optimally evaluated with respect to their sequence, (2) there exists at least one optimal chromosome, i.e., one giving an optimal solution after evaluation (consider one optimal solution and concatenate its trips).

These properties, yet trivial, are not always respected in the GAs whose chromosomes (*genotypes*) are abstractions of actual solutions (*phenotypes*). For instance, most GAs for scheduling problems use lists of operations, because it is sometimes complicated to design crossovers that directly combine two schedules. These lists can be converted into actual schedules, using a scheduling engine. The engines that build non-delay schedules are very good on average, but it is well known that the set of non-delay schedules contains no optimal solution for some instances (French, 1982). In that case, the second property does not hold.

A chromosome is created either by random generation (initial population), by crossover, or by converting an existing ECARP solution  $T = (T_1, \ldots, T_K)$ . In the third case, the trips are concatenated from left to right and the fitness is recomputed with *Split*, i.e. we forget cost(T). There are two main reasons for this policy. First, the solution computed by *Split* is at least as good as T. Second, reproduction is based on a fitness-biased selection of parents (cf. 4.6): to be coherent, all chromosomes must be evaluated in the same way.

Compared to traditional local search, a genetic algorithm works on a population of solutions and its crossovers based on two solutions define larger neighbourhoods. This gives a spatial dimension to the search, often called *intrinsic parallelism*. Thanks to the two properties of our chromosome system, this parallelism is expected to find one optimal ECARP solution.

Figure 4 shows a basic CARP with W=5, 22 edges, and  $\tau=11$  edge-tasks with unit demands (bold) and costs (in brackets). The underlying directed graph G with



Rank :	1	2	3	4	5	6	7	8	9	10	11 =	τ
Cut at :					Ī	)=6		d=8				
						$\downarrow$		$\downarrow$				
Parent P1:	31	21	20	17	15   0	7 0	3	12	23	19	26	
Parent P2:	34	09	29	20	41 2	26 4	3	25	15	39	23	
Child C1:	09	20	41	26	43   0	7 0	3	12	15	39	23	
P1 split :	(31,	21,	20,	17,	15),(0	7),(0	3,	12,	23,	19,	26),	F(P1) = 318
P2 split :	(34,	09,	29),	(20,	41, 2	26),(4	3,	25,	15,	39,	23),	F(P2) = 324
C1 split :	(09,	20,	41,	26),	(43, 0	7),(0	3,	12,	15,	39,	23),	F(C1) = 311

Figure 4. A basic CARP instance with 11 tasks and an example of LOX crossover. Each edge is given with the arc index u for direction (i, j), i < j. The opposite arc is inv(u) = 22 + u.

m=44 is not shown but each edge [i,j] is given with the arc index (i,j) such that i < j, e.g., 7 for (2,4). The index for (j,i), not shown, is by convention 22 + u, e.g., 29 for (4,2). Three chromosomes P1, P2 and C1 are given, for the LOX crossover explained in 4.3. The three last lines give the trips and solution costs found by *Split*. Note that some tasks are treated in two different directions by P1 and P2, e.g., edge [3,4] is collected as (3,4) in P2 (arc index 9) but as (4,3) in P1 (index 31).

## 4.2. Efficient splitting procedures for two objective functions

Algorithm 2 is an  $O(\tau^2)$  version of *Split* working on a given chromosome (giant tour) S. It minimizes total cost (subject to the sequence of tasks defined by the chromosome) and, as a secondary objective, the number of vehicles. It runs in  $O(\tau)$  space only, by avoiding an explicit generation of the auxiliary graph H. Two labels are used for each node i of H:  $V_i$  (cost of the shortest path from 0 to i in H) and  $N_i$  (number of arcs on that path, i.e. number of trips).

Thanks to two indexes i and j, the algorithm enumerates all subsequences  $(S_i, \ldots, S_j)$  of S that correspond to feasible trips and computes their loads and costs using equations (2) and (3). Instead of creating one arc (i-1, j) for a trip composed of tasks  $S_i$  to  $S_j$ , like in 3.3, the labels of j are immediately updated. At the end, the total cost F(S) and the minimum number of vehicles K for that cost can be read in  $V_T$ 

```
procedure split(S)
 V(0), N(0) := 0
 for i := 1 to \tau do V(i) := \infty endfor
 for i := 1 to \tau do
   load, cost := 0; j := i
   repeat
     load := load + q(S(j))
     if i = j then
       cost := D(\sigma, S(i)) + w(S(i)) + D(S(i), \sigma)
       cost := cost - D(S(j-1), \sigma) + D(S(j-1), S(j)) + w(S(j)) + D(S(j), \sigma)
     endif
     if (load \leqslant W) and (cost \leqslant L) then
       VNew := V(i-1) + cost
       if (VNew < V(j)) or ((VNew = V(j))) and (N(i-1) + 1 < N(j)) then
         V(j) := VNew
         N(j) := N(i-1) + 1
       endif
       j := j + 1
     endif
   until (j > \tau) or (load > W) or (cost > L)
 endfor
endproc
```

**Algorithm 2.** *Split* procedure minimizing total cost and number of vehicles.

```
procedure split_for_makespan(S)
 build the auxiliary graph H explicitly
 K, V(0), V2(0), N(0), N2(0) := 0
 for i := 1 to \tau do V(i), V2(i) := \infty endfor
 repeat
   K := K+1; stable := true
   for i := 0 to \tau - 1 do
     for each successor j of i in H with max(V(i),Z(i,j)) < V2(j) do
       V2(j) := \max(V(i), Z(i,j))
       N2(j) := N(i)+1
       stable := false
     endfor
   endfor
   V := V2; N := N2
 until stable or (K = Kmax)
endproc
```

Algorithm 3. Split procedure, version minimizing makespan subject to a limited fleet.

and  $N_{\tau}$ . If required, the ECARP solution can be extracted by tracing the shortest path back.

Algorithm 3 implements *Split* for an interesting ECARP version found in waste collection. The fleet is this time limited. The number of trips K is free but cannot exceed the fleet size  $K_{\rm max}$ . All costs are times and the goal is to minimize makespan (longest trip duration). Note that the problem is trivially solved without  $K_{\rm max}$ , by treating each task by a separate trip. The algorithm uses the same labels as algorithm 2. It computes a min-max path from node 0 to node  $\tau$  in the auxiliary graph H, here explicitly generated.  $Z_{ij}$  is the weight of arc (i, j) in H.

Each iteration of the *repeat* computes in  $O(\tau^2)$  shortest paths in H with at most K arcs. It scans the arcs of H and stores improved label values in V2 and N2. V2 and N2 are copied into V and N at the end of the iteration. The algorithm stops when all labels are stable (checked with the boolean *stable*) or when  $K = K_{\text{max}}$ . The chromosome S is *infeasible* if  $V_{\tau} = \infty$ . If not, the minimal makespan for S and the number of trips actually used are given by  $V_{\tau}$  and  $N_{\tau}$ . Since a shortest path from node 0 to node  $\tau$  in H may have up to  $\tau$  arcs, the algorithm runs in  $O(\min(\tau, K_{\text{max}}) \cdot \tau^2)$ . Note that algorithm 3 can be simply adapted to minimize total cost (but subject to a limited fleet, contrary to algorithm 2), by replacing  $\max(V_i, Z_{ij})$  by  $V_i + Z_{ij}$ .

#### 4.3. Crossovers

Our chromosomes without trip delimiters can undergo classical crossovers for permutation chromosomes. The resulting children are immediately evaluated with *Split*. We tried the linear order crossover LOX and the order crossover OX (e.g., see Oliver, Smith, and Holland, 1987). LOX is designed for linear chromosomes (chromosomes coding objects that clearly have one begin and one end, like hamiltonian paths), while OX rather

```
function OX(P1, P2):chromosome
 //initialise miss to false: no task is already in C
 for u := 1 to \tau do miss(pack(u)) := true endfor
 //draw the substring P1(p)...P1(q) and copy it into C
 draw p in [1,\tau]
 if p = 1 then draw q in [1, \tau -1] else draw q in [p, \tau] endif
 for i := p to q do
   C(i) := P1(i)
   miss(pack(P1(i))) := false
   if inv(P1(i)) \neq 0 then miss(pack(inv(P1(i)))) := false
 endfor
 //browse P2 circularly, starting after the substring, to complete C.
 //{\tt C} is also completed circularly, starting after the substring.
 i,j
         := (q \mod \tau) + 1 //for LOX: i, j := 1
 istart := i
 repeat
   if miss(pack(P2(i))) then
     //for LOX, replace this comment by: if j = p then j := q + 1 endif
     C(j) := P2(i)
     miss(pack(P2(i))) := false
     if inv(P2(i)) \neq 0 then miss(pack(inv(P2(i)))) := false
     j := (j \mod \tau) + 1
   endif
   i := (i \mod \tau) + 1
 until i = istart
 return C
endfunc
```

**Algorithm 4.** Adaptation of the OX crossover for the ECARP.

concerns circular permutations (like TSP tours). Intuitively, the best choice that will be confirmed in section 5 should be OX, because the chromosome before splitting may be viewed as a circular object (giant trip).

Given two parents P1 and P2 with  $\tau$  tasks, both crossovers draw two cutting sites p and q with  $1 \le p \le q \le \tau$ . To get the first child C1, LOX copies P1(p), . . . , P1(q) into C1(p), . . . , C1(q). P2 is then swept from left to right and the tasks missing in C1 are used to fill C1(1), . . . , C1(p - 1) then C1(q + 1), . . . , C1( $\tau$ ). In OX, the sequence for C1 is P1(p), . . . , P1(q) followed by P2(q + 1), . . . , P2( $\tau$ ), P2(1), . . . , P2(p - 1), with restriction that tasks from P2 are taken only if missing in C1. However, C1 is interpreted as a circular list and the result stored such that C1(p) = P1(p). For both crossovers, the other child C2 is obtained by exchanging the roles of P1 and P2.

In the ECARP, a task u is "missing" in C1 if both u and inv(u) are not yet in C1. Algorithm 4 is an ad-hoc version of OX, implemented as a function of two parents P1, P2 that returns a child C (the other child is obtained by swapping the parents in the call). An  $O(\tau)$  complexity is achieved using a table pack that maps the indexes of arcs of R (in  $1, \ldots, m$ ) into  $1, \ldots, \tau$ . Pack is built once for all in O(m), when initializing the MA.

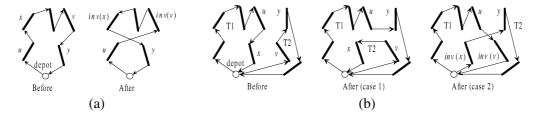


Figure 5. 2-opt moves on one trip (a) and two trips (b). Thick lines correspond to edge-tasks, thin lines to shortest deadheading paths.

The boolean vector *miss* records the tasks missing in C. The algorithm avoids p=1 and  $q=\tau$  at the same time, to ensure  $C \neq P1$ . The only two lines to be changed for LOX are indicated in comments.

#### 4.4. Mutation by local search

In combinatorial optimization, it is well known that the basic GA (Holland, 1975) with simple mutations cannot compete with simulated annealing (SA) and tabu search (TS). To be effective, the generic GA must be hybridized with a local search, giving a *hybrid GA* or *memetic algorithm* (MA) (Moscato, 1999). With a given probability, each child in our MA is converted into an ECARP solution to undergo a local search LS.

LS performs successive *phases* that scan in  $O(\tau^2)$  all pairs of tasks (u, v) to try the following moves, in which x (resp. y) is the task serviced after u (resp. v) in the trip of u (resp. v). Each phase ends by performing the first improving move detected or when all pairs (u, v) are examined. LS stops when a phase reports no improvement. The final ECARP solution is converted into a chromosome, as explained in 4.1. Here are the types of moves examined:

 $N_1$ : invert task u in its trip if it is an edge-task, i.e., replace u by inv(u) in the trip,

 $N_2$ : move task u after task v, or before v if v is the first task of its trip,

 $N_3$ : move adjacent tasks (u, x) after task v, or before v if v is the first task of its trip,

 $N_4$ : swap tasks u and v,

 $N_5$ : two-opt moves (explained in *figure 5*).

Each move type involves *one trip* or *two distinct trips*. Moreover, when moving an edgetask in  $N_1$ – $N_4$ , its service direction may be inverted or not. For instance,  $N_4$  comprises in fact four swapping cases: u and v may be replaced by v and u, inv(v) and u, v and inv(u), or inv(v) and inv(u). In  $N_5$ , some moves may require the inversion of a substring of tasks (cf. figure 5): they are discarded if the substring contains arc-tasks (not invertible).

#### 4.5. Population structure and initialization

The population is implemented as an array  $\Pi$  of nc chromosomes, kept sorted in increasing cost order to ease the selection process described in 4.6. In traditional GAs, identical

solutions or *clones* may appear, leading to a premature convergence. The phenomenon worsens in MAs because the local search quickly compresses  $\Pi$  in a reduced cost interval. A possible remedy is to forbid clones. Exact clone detection can be performed efficiently, e.g., using hashing techniques (Cormen, Leiserson, and Rivest, 1990). We adopted an approximate but faster system in which all individuals have *distinct costs*. Let UB be an upper bound on solution costs and *used* a boolean vector, indexed from 0 to UB, such that used(c) = true iff  $\Pi$  contains an individual of cost c. We know in O(1) if a new chromosome S can be added to  $\Pi$  by checking that used(F(S)) = false. A crossover is said unproductive if its children cannot be kept because of duplicate costs. This concerns a minority of crossovers if nc is not too large (cf. section 5).

 $\Pi$  is initialized with random chromosomes. Because of clones, when nc is too large or the problem very small, many draws may be required to add a new distinct chromosome to the k ones already generated,  $k=1,2,\ldots,nc-1$ . In practice, we try a fixed number of times mnt to generate the next chromosome and truncate  $\Pi$  to nc=k if all draws fail. It is also possible to include in  $\Pi$  a few good heuristic solutions, for instance computed by EPS, EAM or EUH (cf. section 3). These solutions must be converted into chromosomes, as explained in 4.1.

#### 4.6. Incremental memetic algorithms

The basic iteration of an *incremental GA* selects two chromosomes to undergo crossover and mutation. The resulting children immediately replace some existing chromosomes in  $\Pi$ . In a *generational GA* (4.7), the basic iteration (called *generation*) performs a massive reproduction involving all chromosomes. The resulting children are either stored in another population array used for the next generation, or added to  $\Pi$ . In the second case,  $\Pi$  is reduced from 2nc to nc individuals at the end of each generation, by eliminating the nc worst solutions.

We designed incremental versions with two types of selection. The first type (Reeves, 1995) selects the rank i of P1 with probability 2(nc - i + 1)/(nc(nc + 1)). Since  $\Pi$  is sorted in increasing cost order (4.5), the probability of drawing an individual with median cost is roughly 1/nc, the probability of drawing the fittest  $\Pi_1$  is doubled 2/(nc + 1), while the probability of drawing the worst individual  $\Pi(nc)$  is only 2/(nc(nc + 1)). The rank of P2 is drawn uniformly with a probability 1/nc. The second type is *binary tournament*. Two chromosomes are randomly selected and the least-cost one is kept for P1. The process is repeated to get P2.

An OX or LOX crossover (4.3) is applied to (P1, P2). One child C is selected at random and undergoes a mutation by local search (4.4) with a given probability. Two replacement strategies were tested: C replaces either the worst individual  $\Pi(nc)$  or one  $\Pi(k)$  above the median cost, i.e., with  $k \ge \Pi(\lfloor nc/2 \rfloor)$ . Note that both methods preserve the best solution. If no duplicate cost appears, the child mutated or not enters  $\Pi$  and one productive iteration is counted. If not, the child is rejected and the iteration is unproductive.

```
procedure partial_replacement(\Pi,nc,nrep)
 done := 0 //number of solutions currently replaced
 repeat
   generate a population \Omega with nrep distinct costs not present in \Pi
   sort \Omega in increasing cost order
   k := 0
   repeat
     k := k + 1
      if F(\Omega(k)) < F(\Pi(nc)) then
        \Pi(\text{nc}) := \Omega(k); done := done + 1; re-sort \Pi
      else
        cross \Omega(k) with each individual of (\Pi \cup \Omega) \setminus {\Omega(k)}
        C := best child with a cost not present in \Pi
        if F(C) < F(\Pi(nc)) then
          \Pi(\text{nc}) := C; \text{ done } := \text{done } + 1; \text{ re-sort } \Pi
        endif
      endif
    until (done = nrep) or (k = nrep)
 until done = nrep
endproc
```

**Algorithm 5.** Partial replacement procedure used in restarts.

Our incremental MAs perform a main phase stopped after a given number of productive crossovers, after a given number of productive crossovers without improving  $\Pi_1$ , or when reaching a lower bound LB (in that case,  $\Pi_1$  is of course optimal). More instances are solved by adding a fixed number of short restarts, based on a *partial replacement procedure* (Cheung, Langevin, and Villeneuve, 2001). Each restart stops after a fixed number of crossovers or by reaching LB. In section 5, the same number of restarts and the same length per restart are allocated to all instances. Since LB is reached in the main phase for a majority of standard instances, restarts are not always used. Section 5 clearly indicates the number of allowed restarts, the number of crossovers allowed per restart, and the numbers of restarts and crossovers actually performed.

Algorithm 5 adapts Cheung's procedure to replace nrep chromosomes in our population  $\Pi$ . Each set  $\Omega$  is built by generating random chromosomes until nrep distinct cost values not already present in  $\Pi$  are obtained. This process never fails in practice because nc is small (see 5.2) and nrep is even smaller (typically,  $nrep = \lfloor nc/5 \rfloor - \lfloor nc/3 \rfloor$ ). Compared to a blind replacement, the procedure preserves the best solution and never degrades the worst cost. According to Cheung et al., it gives better final solutions for a given CPU time.

## 4.7. Generational memetic algorithms

We also tested a generational scheme described by Hartmann (1998). Each generation randomly partitions  $\Pi$  into pairs. Each pair undergoes a crossover. All children are added to  $\Pi$ , giving 2nc chromosomes, and  $\Pi$  is reduced by keeping the nc best so-

lutions. This method was adapted as follows for populations with distinct costs: the enlarged population is sorted in increasing cost order, and one representative is kept for the *nc* smallest costs. When several chromosomes have the same cost, better diversity is achieved by keeping the most recent one.

## 5. Computational evaluation

## 5.1. Implementation and benchmarks used

All algorithmic components are implemented in the Pascal-like language Delphi 5 and tested on a 1 GHz Pentium-3 PC under Windows 98. The computational evaluation uses three sets of benchmark problems downloadable at  $http://www.uv.es/\sim belengue/carp.html$ .

The first set (*gdb* files) contains 25 undirected instances built by DeArmon (1981), with 7–27 nodes and 11–55 edges. Instances 8 and 9 are never used because they contain inconsistencies. The second set (*val* files) contains 34 undirected instances designed by Belenguer and Benavent (2003) to evaluate a cutting plane algorithm. These files have 24–50 nodes and 34–97 edges. In these two first sets, all edges are required: each instance is in fact a UCPP (*Undirected Capacitated Chinese Postman Problem*), a special case of the CARP.

The third set (*egl* files) provides 24 undirected instances built by Belenguer and Benavent (2003). They are called Eglese instances by these authors, because they are based on the road network of the county of Lancashire (UK), used by Eglese (1994) for a winter gritting problem. Belenguer and Benavent have generated 12 files per area, by varying the vehicle capacity W and the percentage of required edges These instances are very interesting for their realism, their large size (77–140 nodes, 98–190 edges), and also because they contain true CARPs and not only UCPPs like in *gdb* and *val* sets.

## 5.2. Best components, standard setting of parameters and stopping criteria

The best selection of components was determined during an initial testing phase on gdb files. We started from a basic incremental MA, with a population of nc=50 random chromosomes without clones, the Reeves selection, the LOX crossover and a local search rate  $p_m=0.02$ . Like in 4.6, one child is randomly selected after each crossover to replace one chromosome drawn above the median cost. This MA stops when a lower bound is reached or after 5000 crossovers. The list of experiments and resulting decisions are summarized in table 2.

As pointed out by Barr et al. (1995), an acceptable testing of metaheuristics must distinguish "standard" results, reported for one setting of parameters, and "best results" found using various combinations of parameters. The standard setting is important for comparisons with other methods and to give an idea about performance in operational conditions, e.g., when an executable file with frozen parameters is used or when it is too long to try different settings. Our *standard setting* (table 3) has also been found during the preliminary testing. It is the one giving the best average solution values when applied

Table 2 Experiments for selecting best components.

No	Experiment	Impact on solution costs	Decision
1	Inhibit local search LS	Increase	Keep local search
2	Clones always accepted (used ignored)	Increase	Clones never accepted
3	Test various combinations $(nc, p_m)$	Best one is $nc = 30$ , $p_m = 0.1$	Use $nc = 30$ , $p_m = 0.1$
4	Switch to a generational MA	Slight increase	Keep incremental MA
5	Binary tournament selection	Slight decrease	Use binary tournament
6	OX crossover	Slight decrease	Use OX
7	Keep 2 children after OX, not one	Slight increase	Keep one child C
8	C replaces worst solution only	Increase	Not adopted
9	C rejected if $F(C) > a$ given cost	Increase + high rejection rate	Not adopted
10	EPS, EAM, EUH in initial Π	Slight decrease	Use EPS, EAM, EUH
11	Apply LS to initial Π	Increase	No LS on initial Π
12	Add restarts	Decrease	Restarts added

Table 3 Standard setting of parameters.

Name	Role	Value
nc	population size	30
UB	largest cost used (dimension of vector <i>used</i> defined in 4.5)	50000
mnt	max no. of attempts to get each initial random chromosome	50
$p_m$	local search rate in main phase	0.1
mnpi	max no. of productive crossovers in main phase	20000
mnwi	max no. of productive crossovers without changing $\Pi(1)$ , in main phase	6000
mnrs	max no. of restarts	20
nrep	no. of solutions replaced in each restart (partial replacement procedure)	8
$p_r$	local search rate in restarts	0.2
rnpi	max no. of productive crossovers per restart	2000
rnwi	max no. of productive crossovers without changing $\Pi(1)$ , per restart	2000

to all gdb instances. The size of used (see 4.5), UB = 50000, corresponds to the largest cost found in the initial populations of all instances (around 33000 for some egl files), multiplied by a security factor 1.5.

Algorithm 6 illustrates the structure of the best resulting MA and the *stopping criteria*. The procedure *initialize* builds the initial population. The main phase is a call to the procedure *search* (MA basic loop) with a local search rate  $p_m$ . This phase ends after *mnpi* productive iterations (crossovers), after *mnwi* non-improving crossovers, or when a lower bound LB is reached. The MA stops there if  $F(\Pi(1)) = LB$ . If not, it executes a restart loop limited to *mnrs* iterations. Each restart calls the replacement procedure of algorithm 5 and the procedure *search*, but this time with the stronger local search rate  $p_r$  and the reduced numbers of crossovers *rnpi* and *rnwi*. *Search* and the restart loop may stop at any time by reaching LB.

## 5.3. Results for gdb files

Table 4 gathers the results for gdb files. Firstly, we describe the table format, shared by the three sets of benchmarks. After the file name, the number of nodes n and the

```
main program
  initialize(\Pi,nc,used,UB,mnt)
                                                          //initialize population
 if F(\Pi(1)) > LB then begin
                                                          //if LB not reached
   \verb|search|(\Pi, \verb|nc|, \verb|used|, \verb|LB|, \verb|pm||, \verb|mnpi|, \verb|mnwi|)
                                                          //perform main phase
   restarts := 0
                                                          //initialize restart counter
   while (restarts < mnrs) and (F(\Pi(1)) > LB) do
                                                          //perform restarts
     restarts := restarts + 1
                                                          //count one restart
     partial\_replacement(\Pi,nc,nrep)
                                                          //cf. algorithm 5
     search(Π,nc,used,LB,pr,rnpi,rnwi)
                                                          //intensive short phase
   endwhile
  endif
endmain
procedure initialize (\Pi, nc, used, UB, mnt)
  for k := 1 to UB do used(k) := false endfor
                                                          //cost values used, cf. 4.5
                                                          //no of chromosomes built
 get solutions of EPS, EAM and EUH as H(1), H(2), H(3)
                                                          //heuristics of Section 3
 for i := 1 to 3 do
                                                          //try to put solutions in \Pi
   convert H(i) into a chromosome S; split(S)
                                                          //reevaluate, see why in 4.1
   if not used(F(S)) then
                                                          //if cost not duplicated
     k := k + 1; \Pi(k) = S; used(F(S)) := true
                                                          //add S to \Pi
   endif
  endfor
                                                          //generate random solutions
  repeat
   try := 0
                                                          //initialize no of attempts
   repeat
                                                          //loop on attempts
     try := try + 1
                                                          //count one attempt
     generate S at random; split(S)
                                                          //build a random chromosome
   until (not used(F(S))) or (try = mnt)
                                                          //until OK or failed
   if not used(F(S)) then
                                                          //if cost not duplicated
     k := k + 1; \Pi(k) = S; used(F(S)) := true
                                                          //add S to \Pi
 {\bf until} (k = nc) or (used(F(S))
                                                          //\Pi filled or fail
  if used(F(S)) then nc := k endif
                                                          //actual population size
 sort \Pi in increasing cost order
                                                          //sort for replacements
//pls: LS rate, mpi: max. no of productive Xovers, mwi: idem, without
improvement procedure search(\Pi,nc,used,LB,pls,mpi,mwi)
                                                          //productive crossovers
 npi := 0
 nwi := 0
                                                          //idem, without improvement
 repeat
                                                          //MA search loop
   select parents P1, P2 by binary tournament
                                                          //selection, cf. 4.6
   C := OX(P1, P2)
                                                          //crossover, cf. 4.3
                                                          //evaluation (algorithm 2)
   split(C)
   select k at random in [[nc/2],nc]
                                                          //\Pi(k) to be deleted (4.6)
                                                          //local search LS required?
//apply LS, cf. 4.4
   \textbf{if} \text{ random < pls } \textbf{then}
     M := LS(C)
                                                          //{\rm reevaluate}, see why in 4.1
     split(M)
     //\Pi(k) will be replaced in priority by M if M is not a clone,
     //otherwise by the child before its mutation
     if (not used(F(M))) or (F(M) = F(\Pi(k))) then C := M endif
   endif
   if (not used(F(C))) or (F(C) = F(\Pi(k))) then
                                                          //accept replacement
                                                          //count one productive xover
     npi := npi + 1
     if F(C) < F(\Pi(1)) then nwi := 0 else nwi := nwi + 1 endif
     used(F(\Pi(k))) := false; used(F(C)) := true
                                                          //update costs in use
     \Pi(k) := 0
                                                          //perform replacement
     re-sort Π
                                                          //\text{keep}\ \Pi sorted
   endif
 until (npi = mpi) or (nwi = mwi) or (F(\Pi(1)) = LB)
endproc
```

**Algorithm 6.** Best MA structure with initialization and search procedure.

Table 4
Computational results for *gdb* files.

	В																							233*			21	- 1
	Time	0.00	0.44	0.0	0.00	0.11	0.17	0.05	39.82	7.09	0.0	1.26	9.78	7.42	0.05	0.00	0.0	0.05	0.11	0.00	0.33	0.17	3.35	51.19	5.29	51.19		
	$\operatorname{Time}^*$	0.00	0.44	90.0	0.00	0.11	0.17	0.05	99.0	7.09	90.0	1.26	90.0	7.42	0.05	0.00	90.0	0.05	0.11	0.00	0.33	0.17	3.35	51.19	3.16	51.19		
	Xovers	8	2303	179	0	779	1183	0	47187	10708	236	2000	46397	24466	48	0	81	0	141	4	1150	257	3046	24567	7162	47187		
	Rstrts	0	0	0	0	0	0	0	70	_	0	0	20	10	0	0	0	0	0	0	0	0	0	10	2.7	70		
	Std MA	316*	339*	275*	287*	377*	298*	325*	350	303*	275*	395*	458	536*	*001	28*	127*	91*	$164^{*}$	55*	121*	156*	200*	233*	0.15%	1.78%	21	1
gdb files.	EUH	330	353	297	320	407	318	330	388	358	283	413	537	552	104	28*	132	93	172	63	125	162	207	239	6.4%	19.3%	_	4
esults tor	EAM	349	370	319	302	423	340	325*	393	352	300	449	569	260	102	09	129	91*	174	63	129	163	204	237	8.4%	24.2%	C	1
omputational r	EPS	350	366	293	287*	438	324	363	463	354	295	447	581	563	114	09	135	93	177	57	132	176	208	251	10.4%	34.6%	_	
Comp	Best-known	316* abcd	339* abd	275* abcd	287* abcd	377* abd	298* abd	325* abcd	348 bd	303* d	275* abcd	395* abd	458 abd	538 d	$100^*$ abcd	58* abcd	127* abcd	91* abcd	164* abcd	55* abcd	121* abd	156* abcd	200* abcd	233* c	0.14%	1.78%	20	21
	Time	3.15	5.17	0.07	0.0	5.59	0.85	0.00	61.00	53.91	1.55	2.29	20.63	2.42	0.48	0.00	1.70	0.00	0.28	0.20	9.50	1.13	3.38	34.37	9.03	61.00		
	Carpet	316*	339*	275*	287*	377*	298*	325*	352	317	275*	395*	458	<b>5</b> 4	100*	28*	127*	91*	*49I	55*	121*	156*	500*	235	0.48%	4.62%	<u>×</u>	0.7
	TB	316	339	275	287	377	298	325	344	303	275	395	450	536	100	28	127	91	164	55	121	156	200	233				
	1	22	56	22	19	56	22	22	46	51	25	45	23	28	21	21	28	28	36	11	22	33	4	55				
	и	12	12	12	11	13	12	12	27	27	12	22	13	10	7	7	∞	∞	6	∞	11	11	11	11				
	File	gdb1	gdb2	gdb3	gdb4	gdb5	gdb6	gdb7	gdb10	gdb11	gdb12	gdb13	gdb14	gdb15	gdb16	gdb17	gdb18	gdb19	gdb20	gdb21	gdb22	gdb23	gdb24	gdb25	Average	Worst	Optima	)

Lower bounds from Benavent and Belenguer (to appear), except for gdb14 (Amberg and Voß, 2002).

Heuristics cited for published best-known solutions: (a) Belenguer, Benavent, and Cognata (1997), (b) Hertz, Laporte, and Mittaz (2000), (c) Pearn (1989), (d) Lacomme, Prins, and Ramdane-Chérif (2001).

Asterisks denote proven optima, grey cells indicate solutions improved compared to the preliminary GA of Lacomme, Prins, and Ramdane-Chérif (2001), boldface show new best solutions.

Running times in seconds on a 1 GHz Pentium-III PC. Original times for Carpet have been scaled. Improvement for gdb10 in "Best MA" column obtained by using LOX crossover instead of OX.

number of tasks  $\tau$ , the 4th column gives the bound obtained by Belenguer and Benavent (2003), except for gdb14 where it is improved by Amberg and Voß (2002). The two next columns Carpet and Time show the cost reached with standard parameters by Carpet, the best TS heuristic available for the CARP (Hertz, Laporte, and Mittaz, 2000) and the running time in seconds, scaled for the 1 GHz Pentium-III PC used for the MAs. According to SPEC (2001), the power index for the 195 MHz SGI Indigo-2 workstation used by Carpet is 8.88 for integer computations. SPEC does not report benchmarks beyond 500 MHz for the Pentium-III, but we found 41.7 for a 866 MHz at http://you.genie.co.uk/peterw/service/compare.htm, corresponding approximately to 48.2 for 1 GHz. So, we have divided the original Carpet times by 48.2/8.88 = 5.43.

The best-known solutions before this paper are listed in column *Best-known*. The *EPS*, *EAM* and *EUH* columns report solution costs computed by the extended versions of Path-Scanning, Augment-Merge and Ulusoy's heuristic (cf. section 3). Note that this is the first evaluation of Ulusoy's method on standard benchmarks. Then, the table provides the costs obtained by the MA with standard parameters (*Std MA*), the number of restarts used *Rstrts*, the overall number of productive crossovers *Xovers*, the running time until last improvement *Time\**, the overall running time *Time*, and the best cost found using various settings (*Best MA*).

Asterisks denote proven optima, grey cells signal solutions that are improved compared to the GA of Lacomme, Prins, and Ramdane-Chérif (2001), and boldface indicate new best solutions. The last four rows indicate for each column: (a) the average value, given as a deviation to LB in % when the column concerns solution costs (*Average*), (b) the worst value (*Worst*), (c) the number of proven optima (*Optima*) and (d) the number of best-known solutions found (*Best*).

EUH outperforms the other basic heuristics EPS and EAM. The standard MA is at least as good as Carpet in all cases. Compared to Carpet, four instances are improved (10, 11, 15, 25), the average and worst deviations to LB are more than halved and the average running time is 40% smaller. Compared to our first GA (Lacomme, Prins, and Ramdane-Chérif, 2001) needing 21 seconds at 500 MHz on average, the MA runs twice as fast and improves two instances (15, 25). Instance *gdb15* is optimally solved for the first time. Note that these excellent results are achieved without restarts for 18 out of 23 instances (i.e., the lower bound is reached during the main phase). Using several settings (*Best MA*), the MA improves only its solution to *gdb10* but finally finds all best solutions. These results show that *gdb* instances are no longer hard enough for testing CARP metaheuristics.

## 5.4. Results for val files

Table 5 reuses the format of table 4 to present the results for *val* files. The best lower bounds are all obtained by Belenguer and Benavent (2003). Note that the bound 138 they gave to Hertz, Laporte, and Mittaz (2000) for instance *val3c* was not correct, due to an error in their lower bound procedure. The correct value is in fact 137. The *val* files seem empirically harder than *gdb* files: the average deviations to *LB* grow for all

Computational results for val files.

Best MA	173*	245	227*	259*	457	81* 81*	*2	138	*00	412*	428*	530	425÷	440 6440	4,4 1,4 1,4	223*	233	$\frac{317}{670*}$	*672	783 737	386 386	395*	527	323*	326*	332*	391 ***	436* 436*	446*	278	0.54%	4.20% 22	34
IIme	0.00	28.67	0.05	0.22	$\frac{21.76}{2.25}$	0.05	0.00	28.23	0.72	1.21	19.11	103.26	1.86	1.04	90.74	0.17	67.34	52.23	1.97	101	0.66	9.95	71.46	18.29	29.39	71.19	25.115	4.67	17.30	215.04	38.35	413.04	
TITIE	0.00	0.27	0.05	0.22	8.08	0.05	0.00	0.49	0.72	1.21	19.11	6.37	1.86	1.04 4.24	1.32	0.17	6.48	52.23	4.66	60.44	0.66	9.95	62.83	18.29	29.39	71.19	20.07	4.67	17.30	182.85	18.61	182.83	
VOVCIS	0	46554	128	375	50933	34	S	47291	235	503	10391	50746	908 431	4/T	56410	89	53508	53548	126	15/	777	7038	51659	4156	7832	14512	00800	199	4200	52710	56410	6/107	
KSILIS	0	20	0	0	70	0	0	70	0	0	7	50 0	0	)	200	90	20	50 0	0	) (	0	0	20	0	0	4 5	90		0	20	7.3	77	
Std MA	173*	245	227*	259*	45/	<u>81</u> *	*2	138	*00	412*	428*	541	423÷	044	4 5 1 2 4	223*	233	317	2/9* 200*	587	*986	395*	527	323*	326*	332*	391 *907	470 436*	*446 **	530	0.61%	4.20% 22	32
	173*	280	255	281	515	82	66	153	436	468	486	809	104	480 707	504	243	253	367	293	295 281	432	439	603	345	350	368	407 757	457	496	589	10.9%	20.9% 1	_
LAIVI	190	262	238	275	533	98	96	152	443	487	483	631	400 100 100	/8 <del>1</del>	606	241	247	365	300	314 387	412	426	604	348	358	368	450	460	478	590	11.4%	%1.c7 0	C
LFC	186	331	259	284	5 <u>16</u>	× ×	66	158	451	487	539	656	0/4/	208 244	450	271	274	381	326	303 304	433	455	296	358	352	394	492 253	474 474	503	614	16.8%	40.9% 0	C
Best-known	173* abd	245 bd		$259^*$ abd	_					412* abd	428* ad	530 d	423" abd	446 abd	581 d	223* abd	233 ad	317 bd	2/9* abd	283° abd	386* abd	395* abd	527 d	323* abd		332* abd	*			530 d	0.55%	4.20% 22	33
Time	0.02	93.20	0.17	13.02	31.66	0.77	2.79	41.66	28.32	75.66	70.06	233.56	3.80	41.40	23.27 224 11	3.89	26.94	85.18	6.59	0.07 121 14	3.84	81.46	147.40	28.51	59.89	56.44	55.70	18.43	93.47	156.31	63.87	223.28	
Carpet	173*	245	227*	260	464 465	<u>81</u> *	*/*	138	*00	416	453	556	. 574 	4 4 8 7 7	6/4	223*	241	329	*6/2	587 273	386*	401	533	323*	329	332	405 ***	436*	451	145	1.90%	8.37%	17
ΓŖ	173	235	227	259	455	81	87	137	400	412	428	520	577	0440	404 777	223	231	311	6/2	283 233	386	395	517	323	326	332	700	436	446	524			
2	39	39	34	37	45	35	35	35	69	69	69	69	60	65	36	20	50	30	9 (	200	38	63	63	92	92	225	760	07	07	97			
и	24	12	24	43	4.	4	7	7	41	41	4	45	4 5	4 5 4 5	¥ 4	3.5	31	31	<del>5</del> 5	9€	£ 6	30	30	50	20	20	3 6	86	86	20			
rile	valla	valle	val2a	val2b	vaľ2c i2	val3a	val3b	val3c	val4a	val4b	val4c	val4d	valSa	val50	valse valsd	val6a	val6b	val6c	val/a	val/b	val8a	val8b	val8c	val9a	val9b	val9c	valyd vel10e	vall0a vall0b	val10c	val10d	Ayerage	worst Optima	Best

algorithms and 15 instances out of 34 require restarts. The average running time is now 38 seconds, but the last improvement is found earlier (compare 18.61/38.35 = 49% for val-files against 3.16/5.29 = 60% for gdb files). Among the constructive heuristics, EUH performs better than EPS and EAM. Again, compared to Carpet, the standard MA provides identical or better solutions, divides by two the average and worst deviations to LB and runs 40% faster.

Using several settings, the MA yields all best solutions, improves the preliminary GA of Lacomme, Prins, and Ramdane-Chérif (2001) three times, and finds a new best solution for val10d. This solution was obtained by trying an exact clone detection: two chromosomes  $S_1$  and  $S_2$ , converted by Split into ECARP solutions, are clones if each trip of  $S_1$  can be retrieved in the trips of  $S_2$  with the same sequence of tasks or, for undirected instances, with an inverted sequence. However, the simpler clone detection based on distinct costs was kept because it provides slightly better results on average. The reason probably resides in the integer costs and in the relatively small range of solution values in the instances tested, favouring a better dispersal of solutions.

#### 5.5. Results for egl files

Table 6 shows the results for these files constructed by Belenguer and Benavent from Eglese's data. The number of edges m/2, often greater than  $\tau$ , is now mentioned. The *Carpet* column reports unpublished results of Carpet, computed by Mittaz on behalf of Belenguer and Benavent. The running times are unknown. Since Carpet is here the only heuristic compared with the MA, the redundant *Best-known* column is removed.

The *egl* files seem much harder than the previous files: the average deviation to *LB* augments for all algorithms and *LB* is never reached. Of course, the reason is perhaps inherent to the bound and/or to the heuristics. E.g., according to Belenguer and Benavent, the partial graph of required edges is sometimes disconnected and their bound does not exploit this property. All in all, EUH remains the best simple heuristic, the standard MA outperforms Carpet 19 times and the best MA improves all solution values, proving that Carpet finds no optimal solution. The price to pay is a larger average running time (9 min): the instances are bigger and, since *LB* is never reached, the MA performs in all cases its 20 restarts.

#### 5.6. Makespan minimization

The flexibility of the memetic algorithm is illustrated here by minimizing a different objective function for gdb files: the duration of the longest trip (makespan), subject to a limited number of vehicles. The two main changes in the MA are to replace algorithm 2 by algorithm 3 (see 4.2) for Split and to use the new objective function in the local search LS. Let  $q_{tot}$  be the total demand. The fleet size  $K_{max}$  (see 4.2) is set to the smallest possible value  $\lceil q_{tot}/W \rceil$ . This bound is tight for gdb files, since it is always reached by the MAs minimizing total cost.

A relatively simple lower bound LB2 to the optimal makespan can be computed as follows. The duration of a trip containing only one task u is  $cost(u) = D(\sigma, u) + w(u) + v(u) + v(u)$ 

Computational results for egl files. Table 6

	7/11	2	$\Gamma B$	Carpet	FFS	EAIM	EOD EOD	Std MA	KSILIS	Xovers	$\operatorname{Time}^*$	Time	Best MA
egl-e1-A 77	86 /	51	3515	3625	4115	4605	3952	3548	20	47325	1.48	74.26	3548
	86 /	51	4436	4532	5228	5494	5054	4498	20	46249	48.39	69.48	4498
	86 /	51	5453	2663	7240	6429	9919	5595	20	46712	39.98	71.18	5895
_	86 /	72	4994	5233	6458	6253	5716	5018	20	47290	20.60	152.58	5018
	86 /	72	6249	6422	7964	7923	7080	6340	20	57657	22.19	153.41	6340
		72	8114	8603	10313	10453	9338	8415	20	00009	27.52	129.63	8395
		87	6985	2907	7454	7350	6723	2898	20	54522	24.44	242.00	2898
		87	7646	7921	0066	9244	8713	7822	20	29860	173.18	255.35	7816
egl-e3-C 77		87	10019	10805	12672	12556	11641	10433	20	52412	111.50	206.35	10369
_		86	6372	6489	7527	7798	7231	6461	20	48101	275.50	291.87	6461
		86	6088	9216	10946	10543	10223	9021	20	00009	291.49	312.85	9021
egl-e4-C 77	86 /	86	11276	11824	13828	13623	13165	11779	20	51186	77.83	252.38	11779
egl-s1-A 140	) 190	75	4992	5149	6382	6143	5636	5018	20	52115	15.88	208.61	5018
egl-s1-B 140	) 190	75	6201	6641	8631	7992	9802	6435	20	54924	21.42	208.77	6435
egl-s1-C 140	) 190	75	8310	8687	10259	10338	9572	8218	20	49068	160.38	165.55	8518
egl-s2-A 140	) 190	147	08/6	10373	12344	11672	11475	9995	20	26992	795.10	874.36	9995
egl-s2-B 140	) 190	147	12886	13495	16386	15178	14845	13174	20	00009	641.58	760.50	13174
egl-s2-C 140	_	147	16221	17121	20520	19673	19290	16795	20	29606	743.69	746.93	16715
egl-s3-A 140	_	159	10025	10541	13041	11957	11956	10296	20	50853	651.03	1070.50	10296
egl-s3-B 140	) 190	159	13554	14291	17377	15891	15663	14053	20	27886	1043.58	1064.01	14028
egl-s3-C 140	) 190	159	16969	17789	21071	19971	20064	17297	20	00009	622.58	874.30	17297
egl-s4-A 140	) 190	190	12027	13036	15321	14741	13978	12442	20	00009	1529.57	1537.59	12442
egl-s4-B 140	_	190	15933	16924	19860	19172	18612	16531	20	00009	1184.52	1430.26	16531
egl-s4-C 140	) 190	190	20179	21486	25921	24175	23727	20832	20	00009	1464.26	1495.02	20832
Average				4.74%	26.4%	22.8%	15.4%	2.47%	20	54685	416.15	526.99	2.40%
Worst				8.61%	39.2%	31.0%	19.3%	4.46%	20	00009	1529.57	1537.59	4.46%
Best				0	0	0	0	19					24

Table 7
Makespan optimization subject to a limited fleet for *gdb* files.

File	n	τ	LB2	β	$q_{ m tot}/W$	γ	EUH	Std MA	Rstrts	Xovers	Time*	Time
gdb1	12	22	64	63	4.40	64	84	66	10	30152	1.43	10.32
gdb2	12	26	59	59	5.20	57	81	60	10	31479	3.13	17.03
gdb3	12	22	59	59	4.40	55	74	60	10	26110	3.24	10.38
gdb4	11	19	72	64	3.80	72	98	74	10	26487	0.11	6.04
gdb5	13	26	64	64	5.20	63	88	69	10	26054	8.85	13.02
gdb6	12	22	64	64	4.40	60	75	68	10	26476	0.22	10.44
gdb7	12	22	65	57	4.40	65	81	68	10	29679	1.31	11.20
gdb8	27	46	38	38	9.22	35	54	40	10	26791	11.43	33.12
gdb9	27	51	37	37	9.56	31	69	39	10	26224	28.67	48.28
gdb10	12	25	69	39	3.70	69	86	73	10	26477	0.22	10.43
gdb11	22	45	79	43	4.48	79	98	82	10	30544	26.47	44.98
gdb12	13	23	93	93	6.06	64	124	96	10	26490	0.11	6.97
gdb13	10	28	128	128	5.98	90	178	140	10	27866	1.37	19.39
gdb14	7	21	20	15	4.24	20	27	21	10	26158	0.16	24.99
gdb15	7	21	15	8	3.03	15	16	15*	0	3626	4.83	4.83
gdb16	8	28	26	14	4.83	26	40	27	10	27132	0.94	22.14
gdb17	8	28	19	9	4.10	19	22	19*	0	148	0.49	0.49
gdb18	9	36	33	19	4.14	33	40	34	10	27915	3.90	46.90
gdb19	8	11	19	17	2.44	19	24	21	10	26026	0.06	25.00
gdb20	11	22	31	20	3.96	31	45	32	10	27860	1.15	16.75
gdb21	11	33	26	15	5.70	26	50	29	10	26264	0.38	30.65
gdb22	11	44	25	12	7.59	25	45	28	10	30963	17.13	107.71
gdb23	11	55	24	13	9.85	24	39	30	10	31342	21.59	117.38
Average							39.0%	5.9%	9.1	25577	5.96	27.76
Worst							92.3%	25.0%	10.0	31479	28.67	117.38

See section 5.6 for comments.

 $D(u, \sigma)$ , according to equation (3). So, the minimum duration  $\delta(u)$  of a trip reduced to one task u is either cost(u), if u is an arc-task, or  $min\{cost(u), cost(inv(u))\}$ , if u is an edge-task. A first bound to the makespan is obtained by computing the maximum of these costs for all tasks:  $\beta = max\{\delta(u) \mid u \in A, q(u) > 0\}$ . A second bound is  $\gamma = \lceil LB/K_{max} \rceil$ , where LB is the lower bound for the total cost whose values are listed in table 4. Finally, LB2 =  $max\{\beta, \gamma\}$ .

The results are summarized in Table 7. All MA parameters are taken from table 3, except the maximum number of restarts *mnrs*, now set to 10. The only heuristic used for the initial population is the extended Ulusoy's method (EUH, see 3.3). Path-Scanning and Augment-Merge are discarded because they often lead to infeasible solutions. Two optima are found and the average deviation to LB2 is nearly 6%. This gap probably comes from the weakness of the bound: the last improvement is obtained early (5.96 seconds on average) compared to the overall running time (27.76 seconds), indicating that other solutions could be optimal.

DeArmon 23 pbs Benavent 34 pbs Eglese 24 pbs Criterion Carpet MA Carpet MA Carpet MA 0.48 Avg. dev. to LB % 0.15 1.90 0.61 4.74 2.47 Max. dev. to LB % 4.62 1.78 8.57 4.26 8.61 4.46 No. of proven optima 18 21 15 22 0 0 No. of best solutions 19 22 17 32 0 19 Avg. running time (s) 9.02 5.29 63.87 38.35 unknown 526.99

Table 8 Comparison between the standard MA and Carpet.

# 5.7. Performance overview

Table 8 compares performance criteria between the memetic algorithm and Carpet, executed with their respective standard parameters: the average and worst deviations to LB, the number of proven optima (when LB is reached), the number of best-known solutions retrieved, and the average running time on a 1 GHz Pentium-III PC. The standard setting of parameters seems to be *extremely robust*: it gives the best average results for the three sets of instances and its solutions are improved only 8 times out of 81 by trying different settings. Finally, 26 best-known solutions are improved and all other best-known solutions are retrieved.

The MA confirms the interest of a GA template already applied successfully to the open-shop scheduling problem by Prins (2000). Indeed, this earlier GA shares some common features with our MA for the ECARP: a small population with distinct solutions, a few good solutions in the initial population, an improvement procedure used as mutation operator. This shows that powerful hybrid genetic algorithms can be designed thanks to a synergic effect between several simple improvement ideas.

#### 6. Conclusion

The best memetic algorithm for the CARP presented in this paper outperforms all known heuristics on three sets of benchmarks available in public, even when it is executed with one single setting of parameters. This excellent performance results from a combination of several key-features. In spite of simple chromosomes (without trip delimiters) and crossovers, each child is optimally evaluated thanks to the *Split* procedure and can be strongly improved by local search. Small populations of distinct solutions avoid a possible premature convergence. A few good initial solutions are computed via classical heuristics. The incremental management of population and the partial replacement technique used for restarts accelerate the decrease of the objective function value. The absence of complicated techniques must also be underlined.

Moreover, the MA is already designed for tackling several extensions like mixed networks, parallel arcs and turn penalties. We just checked its correct execution on a few instances constructed by hand from a city map. It is too early to provide a computational evaluation for these extensions: more instances must be prepared, appropriate lower

bounds must be developed, while no other algorithm is available for comparison. All these tasks are in progress, in particular a random generator for large scale realistic street networks.

More generally, this paper reflects the increasing research activity on arc routing problems. This activity is confirmed by two very recent articles, published during the last revision of this paper, which bring two other new metaheuristics for solving the CARP: Beullens et al. (2003) describe a guided local search, while Greistorfer (2003) proposes a tabu search procedure that makes use of the scatter search paradigm.

## Acknowledgments

The authors wish to thank the two anonymous referees who performed an extremely detailed analysis of the paper and suggested many useful improvements.

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