A new precedence-based Ant Colony Optimization for permutation problems

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Abstract. In this paper we introduce ACOP, a novel ACO algorithm for solving permutation based optimization problems. The main novelty is in how ACOP ants construct a permutation by navigating the space of partial orders and considering precedence relations as solution components. Indeed, a permutation is built up by iteratively adding precedence relations to a, initially empty, partial order of items until it becomes a total order, thus the corresponding permutation is obtained. The pheromone model and the heuristic function assign desirability values to precedence relations. An ACOP implementation for the Linear Ordering Problem (LOP) is proposed. Experiments have been held on a large set of widely adopted LOP benchmark instances. The experimental results show that the approach is very competitive and it clearly outperforms previous ACO proposals for LOP.

Keywords: Ant Colony Optimization, Permutations Representation, Partial Orders, Linear Ordering Problem

1 Introduction

Ant Colony Optimization (ACO) [6] is a popular meta-heuristic scheme for solving hard combinatorial optimization problems inspired by the foraging behavior of natural ant colonies. Since the seminal work of Dorigo in the early '90s [8], ACO has been extensively and successfully applied to permutation-based optimization problems, i.e., problems where a solution is a permutation of items. See for example the several ACO proposals for the traveling salesman problem [8, 7], the quadratic assignment problem [9] or the permutation flowshop scheduling problem [19].

The typical ACO approach to permutation problems is to consider the items (to be ordered) as the solution components. Therefore, an artificial ant constructs a n-length permutation by starting from the empty sequence and iteratively adding up items until all the n items appear in the sequence exactly once. Usually, the items are sequentially inserted from left to right, though simple variations, where the items can be inserted in arbitrary positions of the (partial) permutation, have been proposed [15].

To the best of our knowledge, this construction scheme is applied by all the ACO proposals for permutation problems available in literature. Basically, they all represent a permutation of n items directly as a n-length sequence without repetitions, thus, during the construction process, a partial solution is a sequence containing empty slots.

Though this representation is very natural, it is not always the most suited for the problem at hand. This is the case of the objective functions defined by means of non-local characteristics such as the precedence relations between the pairs of items contained in a permutation. Indeed, precedences are important in a variety of problems such as some scheduling problems [3, 14] and, in particular, in the Linear Ordering Problem (LOP) [16] where the objective function is defined as the sum of the contributions associated to all the precedences encoded by a permutation. It is evident that, in this case, adding a new item to a partial sequence induces a drastic change to the expected objective value of the permutation under construction.

In this paper, we propose the smoother approach of constructing a permutation by iteratively adding precedences to a, initially empty, partial order until a total order, i.e., a permutation, is formed. Basing on this idea, a new precedence-based ACO approach for permutation problems, namely ACOP, is introduced. Therefore, the main novelty is that ACOP ants represent a partial permutation, i.e., a partial order, as a collection of precedence relations (between items) which are consistent each other. When this collection contains $\binom{n}{2}$ precedences, then a n-length permutation is mathematically guaranteed to be built.

ACOP has been implemented and applied to LOP. Experiments have been held on a large set of widely adopted benchmark instances where ACOP performances have been compared to state-of-the-art results. Moreover, a further experiment has been held to compare ACOP with, as far as we know, the only ACO approach to LOP available in literature, i.e., ACS-IM [5, 18].

2 Ant Colony Optimization

2.1 ACO General Scheme

ACO algorithms [8,9,15] are inspired by the stigmergic foraging behavior of natural ant colonies. When a real ant discover a food source, it walks back to its nest by also leaving pheromone trails on the way, so other ants can sense the trails and reach the food themselves. Analogously, in ACO, artificial ants build up combinatorial solutions component-by-component using a probabilistic construction procedure biased by the artificial pheromone trails deposited on solution components by the best-performing ants of the previous iterations.

Let $f: S \to \mathbb{R}$ be the objective function to be optimized, where S is the set of solutions, then each $s \in S$ is composed by a certain number of components c_1, c_2, \ldots, c_n taken from the set of possible components C. Clearly, S and C are problem dependent. For example, in the traveling salesman problem, C is the set of cities, while S contains all the permutations of C.

ACO aims to optimize f by iteratively probing S by means of N artificial ants. The ants indirectly communicate through a common data structure, called pheromone, which associates a real value τ_c to each solution component $c \in C$. The main scheme of ACO is depicted in Figure 1, where, without loss of generality, maximization is assumed.

```
1: function ACO(N, \alpha, \beta, \rho, \Delta_{\tau}^{ib}, \Delta_{\tau}^{gb})
2: Initialize pheromone values \tau_c for all c \in C
 3:
           s^{gb} \leftarrow \text{null}
           while termination condition is not met do
 4:
                s^{ib} \leftarrow \text{null}
 5:
                for i \leftarrow 1 to N do
 6:
                      s_i \leftarrow \text{BuildSolution}(\alpha, \beta)
 7:
 8:
                      Evaluate f(s_i)
                      if f(s_i) > f(s^{ib}) then s^{ib} \leftarrow s_i
 9:
10:
11:
                       end if
                      if f(s_i) > f(s^{gb}) then
12:
                            s^{gb} \leftarrow s_i
13:
14:
                       end if
                 end for
15:
                 Optionally perform a local search on \boldsymbol{s}^{ib} (and update \boldsymbol{s}^{gb})
16:
                 EvaporatePheromone(\rho)
17:
                 DepositPheromone(s^{ib}, s^{gb}, \Delta_{\tau}^{ib}, \Delta_{\tau}^{gb})
18:
19:
            end while
           return s^{gb}
20:
21: end function
```

Fig. 1. General scheme of ACO

Pheromone values are usually initialized to a constant value. Then, at every ACO iteration, each ant starts from an empty partial solution and builds up a complete solution by iteratively choosing components from C. In many problems, the set C_t of feasible components at construction step t is restricted by the choices done in the previous steps, thus, in general, $C_t \subseteq C$. The choice of a component c from C_t is influenced by its pheromone value $\tau_c \in \mathbb{R}^+$ and a problem dependent heuristic value $\eta_c \in \mathbb{R}^+$ which estimates the contribution of c to the solution quality. Formally, the probability of choosing $c \in C_t$ is

$$p(c) = \frac{\tau_c^{\alpha} \eta_c^{\beta}}{\sum_{k \in \mathcal{C}_t} \tau_k^{\alpha} \eta_k^{\beta}},\tag{1}$$

where the parameters $\alpha, \beta \in \mathbb{R}$ determine the influence of, respectively, pheromone and heuristic values.

The construction process terminates when N complete solutions (one per ant) have been generated. Each solution is evaluated using f and the pheromone is

updated. First, for all $c \in C$, the pheromone value τ_c is evaporated as follows

$$\tau_c \leftarrow (1 - \rho)\tau_c,\tag{2}$$

where $\rho \in [0,1]$ is the evaporation rate parameter. Then, a given amount of pheromone is deposited on the components belonging to the best solutions. Though various deposition strategies are possible [15], the most common ones consider the iteration and global best solutions, respectively, s^{ib} and s^{gb} . Formally, for all $c \in C$, the pheromone value τ_c is updated as

$$\tau_c \leftarrow \tau_c + I\left(c \in s^{ib}\right) \Delta_{\tau}^{ib} + I\left(c \in s^{gb}\right) \Delta_{\tau}^{gb},\tag{3}$$

where: $I\left(c \in s\right)$ is 1 if component c belongs to solution s and 0 otherwise, while $\Delta_{\tau}^{ib}, \Delta_{\tau}^{gb} \in \mathbb{R}^{+}$ are the "awards" of pheromones for the components of, respectively, the iteration and global best solutions.

Finally, note that, before pheromone update, a local search refinement can be optionally applied to a selected set of solutions (usually, the iteration best).

2.2 Pheromone Models for Permutation Problems

While the typical permutation construction procedure of ACO schemes has been described in Section 1, here we provide a brief overview of the different pheromone models for permutation problems available in literature [3, 17].

One simple approach, denoted as PH_{abs} in [3], is to associate pheromone values to pairs composed by an item and an absolute position, in order to indicate the desirability to have a given item at a given position in the permutation.

Two other relevant approaches are PH_{suc} and PH_{rel} [3], which both assign pheromone values to ordered pairs of items. While PH_{suc} aims to encode the desirability of having the two items in consecutive positions of the permutation, PH_{rel} is less stringent and only indicates the desirability of the precedence relation between the two items independently of their distance in the permutation.

We highlight that, as far as we know, all the ACO proposals in literature using the pheromone model PH_{rel} build up the permutation as seen in Section 1, i.e., by iteratively adding items to a incumbent sequence till it becomes a complete permutation.

3 Permutations, partial and total orders

Here, we provide a brief mathematical background useful to describe the representation of the (partial) solutions in ACOP. In particular: we introduce an encoding for generic partial orders of items, we show under which conditions the partial order is also a total order and how to obtain its corresponding permutation.

Let I be a finite set of items that, without loss of generality, can be taken as $I = \{1, ..., n\}$, then a strict partial order relation \prec on I is a binary relation which satisfies the following properties:

- Irreflexivity: $a \not\prec a$, for all $a \in I$;
- Transitivity: if $a \prec b$ and $b \prec c$, then $a \prec c$, for $a, b, c \in I$;
- Anti-symmetry: if $a \prec b$, then $b \not\prec a$, for $a, b \in I$.

As any binary relation, a partial order \prec on I can be represented as the set of pairs $P = \{(a,b) : a,b \in I \text{ and } a \prec b\}$, thus the pair $(a,b) \in P$ indicates the precedence $a \prec b$.

Conversely, given any finite set of precedences $P = \{(a_1, b_1), \ldots, (a_k, b_k)\}$, where $a_i, b_i \in I$ for $i = 1, \ldots, k$, such that the precedences in P do not violate any partial order property, it is possible to find the corresponding partial order \prec_P on I generated by P, i.e., the smallest partial order which respects all the precedences in P. Operatively, \prec_P is the transitive closure P^* of P, which is computed as follows. Let $P_0 = P$, then

$$P_{r+1} = P_r \cup \{(a,b) : \exists c \in I \text{ such that } (a,c), (c,b) \in P_r\}.$$
 (4)

After a finite number s of steps, the sequence of sets $\langle P_r \rangle_r$ stabilizes (i.e., $P_s = P_{s+i}$ for any integer $i \geq 1$), because the maximum number of "compatible" precedences is finite and equal to $\binom{n}{2}$. Hence, $P^* = P_s$ and $a \prec_P b$ if and only if $(a,b) \in P^*$.

Importantly, the partial order \prec_P , represented by a given set of pairs P, can also be seen as the arcs set of the digraph G_{\prec} whose nodes set is I and such that there is an arc $a \to b$ for each precedence $(a,b) \in P$. Therefore, a partial order P can be encoded by the $n \times n$ incidence matrix A of G_{\prec} , whose entries are

$$A_{ab} = \begin{cases} 1 & \text{if } a \prec b \\ -1 & \text{if } b \prec a \\ 0 & \text{otherwise.} \end{cases}$$
 (5)

Furthermore, if \prec also satisfies the property that for all $a, b \in I$, with $a \neq b$, either $a \prec b$ or $b \prec a$, then \prec is a strict total order. For a total order, the matrix A does not contain any 0-entry, except in its main diagonal. Moreover, it contains $\binom{n}{2}$ 1-entries and the same number of -1s.

It is easy to see that there is a bijective correspondence between the set of total orders on I and the set S_n of the permutations of I. Indeed, given $\pi \in S_n$, its corresponding total order \prec_{π} is defined as all the precedences $a \prec_{\pi} b$ such that $a, b \in I$ and a appears before b in π . More formally, $a \prec_{\pi} b$ if and only if $\pi^{-1}(a) < \pi^{-1}(b)$, where π^{-1} is the inverse permutation of π . On the other hand, if \prec is a total order, the corresponding permutation π_{\prec} is recursively defined as: (i) $\pi_{\prec}(1) = a$ where $a \in I$ is the unique item such that $b \not\prec a$ for every $b \in I$, and (ii) $\pi_{\prec}(k) = a$ if $b \prec a$ only for all the $b \in \{\pi_{\prec}(1), \ldots, \pi_{\prec}(k-1)\}$.

Therefore, given a total order encoded by a matrix A (see equation (5)), the corresponding permutation π can be recovered by observing that $\pi(k) = a$ if and only if the a-th row of A has exactly n - k 1s. Hence, by setting $\sigma(a)$ to n minus the number of 1-entries in the a-th row of A, for all $a = 1, \ldots, n$, and observing that σ is a permutation, then $\pi = \sigma^{-1}$.

As a further interpretation, note that a partial order \prec individuates the set of permutations $Q_{\prec} \subseteq \mathcal{S}_n$ such that $\pi \in Q_{\prec}$ if and only if π agrees with \prec , i.e., for all $a, b \in I$, if $a \prec b$ then $\pi^{-1}(a) < \pi^{-1}(b)$.

Finally, given a partial order \prec and a new pair (c,d), with $c,d \in I$ and $c \neq d$, such that $d \not\prec c$, it is possible to extend \prec with the precedence $c \prec d$ by simply computing the transitive closure of the set $P_{\prec} \cup \{(c,d)\}$.

4 ACOP: Ant Colony Optimization on Precedences

ACOP is a new Ant Colony Optimization algorithm for permutation based optimization problems which works on the space of partial orders. Its aim is to optimize an objective function of the form $f: \mathcal{S}_n \to \mathbb{R}$, where \mathcal{S}_n contains all the permutations of a set I of n items.

The main structure of ACOP follows the same ACO general scheme depicted in Figure 1. It handles a colony of N artificial ants and uses the pheromone model PH_{rel} previously described in Section 2.2, i.e., pheromone values are maintained in a $n \times n$ matrix where the entry $\tau_{a,b}$, with $a,b \in I$, is the amount of pheromone assigned to precedence $a \prec b$.

The original parts of ACOP are: (i) the (partial) solution representation, and (ii) the construction procedure performed by the artificial ants, i.e., the implementation of *BuildSolution* (see Figure 1).

Indeed, every ant builds up a permutation by iteratively adding precedence relations to a partial order \prec , which is initially empty, until it becomes a total order. The pseudo-code of the procedure is depicted in Figure 2.

```
1: procedure BuildSolution(\alpha, \beta)
                                                                                             \triangleright All 0s in matrix A
 2:
          A \leftarrow \mathbf{0}
 3:
          np \leftarrow 0
                                                                               \triangleright Number of precedences in A
 4:
          while np < \binom{n}{2} do
               C = \{(a, b) : A_{a,b} = 0 \text{ and } a \neq b\}
 5:
                                                                              ▷ Candidate set of precedences
 6:
               (a, b) \leftarrow \text{ChoosePrec}(C, \tau, \eta, \alpha, \beta)
                                                                                                ▶ See equation (1)
 7:
               Q \leftarrow \{(a,b)\}
 8:
               while Q \neq \emptyset do
                                                    \triangleright Insert (a,b) and compute the transitive closure
9:
                    (a,b) \leftarrow \text{remove an element from } Q
10:
                    A_{a,b} \leftarrow 1
                    A_{b,a} \leftarrow -1
11:
                    np \leftarrow np + 1
12:
13:
                    Q \leftarrow Q \cup \{(a,c) : A_{a,c} = 0 \text{ and } A_{b,c} = 1\}
                              \cup \{(c,b): A_{c,b} = 0 \text{ and } A_{c,a} = 1\}
14:
15:
               end while
          end while
16:
17:
          Return A
18: end procedure
```

Fig. 2. The permutation construction procedure of ACOP

The matrix A encodes the partial order \prec that is initially empty, while the variable np is the number of 1s in A, i.e., the number of precedences of \prec . The loop of lines 4–16 iteratively adds 1-entries to A and terminates when A contains exactly $\binom{n}{2}$ 1-entries, i.e., when A encodes a total order which corresponds to a permutation. At each construction step, the ant chooses a precedence from C (line 5). This precedence can be safely added to \prec . The choice of line 6 is performed by considering pheromone and heuristic values as in the general ACO scheme provided in equation (1). The inner loop at lines 8–15 inserts the selected precedence in \prec , by removing some 0-entries from A and iteratively computing the transitive closure on A as described in Section 3. Finally, though, at the end of the procedure, the matrix A can be converted to a permutation (see Section 3), some objective functions can be directly computed on A, therefore BuildSolution returns the matrix A.

5 Application of ACOP to LOP

The Linear Ordering Problem (LOP) is a classical NP-Hard combinatorial optimization problem [16] and has received considerable attention because of its many applications in diverse research fields such as economy [13], graph theory [4], archeology [10] and computational social choice [12].

LOP can be straightforwardly formulated as a matrix triangulation problem [16]. Given a $n \times n$ matrix H, LOP requires to find a permutation $\pi \in \mathcal{S}_n$ of the row and column indices $\{1, \ldots, n\}$ that maximizes the objective function

$$f(\pi) = \sum_{i=1}^{n} \sum_{j=i+1}^{n} H_{\pi(i),\pi(j)}$$
 (6)

The permutation structure of the LOP solutions allows to apply a variety of meta-heuristics and evolutionary algorithms specifically designed for the permutations search space. See for instance [1, 2, 11, 20]. To the best of our knowledge, the only ACO approach to LOP has been proposed by Pintea et al. in [5] and [18].

An interesting observation is that the objective function of LOP can be directly computed on the permutation representation used by ACOP, i.e., on the matrix A returned by the function BuildSolution depicted in Figure 2. Indeed, it is easy to see that equation (6) can rewritten as

$$f(A) = \sum_{a=1}^{n} \sum_{b=1}^{n} H_{a,b} \cdot \max\{A_{a,b}, 0\}.$$
 (7)

Here, it is evident that the contribution of any single precedence relation (a, b) present in the permutation to be evaluated is exactly $H_{a,b}$. Hence, a simple but effective choice for the heuristic function $\eta_{a,b}$ is to set $\eta_{a,b} = H_{a,b} + \epsilon$, where ϵ is a small positive quantity introduced to avoid null probabilities when $H_{a,b} = 0$.

Further details of the implementation of ACOP for LOP are as follows. Pheromone is deposited using a mix of the iteration-best and global-best strategies as depicted by equation (3). Inspired by [21], the pheromone values are constrained to the interval $[\tau_{min}, \tau_{max}]$, where $\tau_{max} = (\Delta_{\tau}^{ib} + \Delta_{\tau}^{gb})/\rho$ and $\tau_{min} = \tau_{max}/(2n^2-n)$. All the pheromones are initialized to τ_{max} . ChoosePrec (line 6 of Figure 2) has been implemented as in [7], i.e., with probability q_0 the most probable precedence is chosen, otherwise a tournament is performed. Precedence probabilities are computed as in equation (1).

Finally, an enhanced variant of ACOP, called ACOP⁺, has been devised. ACOP⁺ performs a local search refinement on the iteration best solution at the end of every iteration. The local search has been implemented by iteratively applying the best item insertion move till no improvement is observed (see [16]). Moreover, in order to avoid stagnation, ACOP⁺ reinitializes the pheromone values if no improvement to the global best solution has been observed during the last r iterations.

6 Experiments

The ACOP application to LOP has been experimentally investigated on the three widely known benchmark suites LOLIB, SGB and MB³. Therefore, a total of 105 LOP instances, with dimensionalities ranging from 44 to 250, has been considered.

The optima of these instances are known⁴ and they have been used to compute two performance measures: the success rate (SR), and the average relative percentage deviation (ARPD). An algorithm is executed k times per instance, thus SR indicates the percentage of executions that reach the known optimum, while ARPD = $\frac{100}{k} \sum_{i=1}^{k} \frac{opt-run_i}{opt}$ is the average percentage deviation from the known optimum.

ACOP parameters have been experimentally tuned on a subset of 10 selected instances: the (lexicographically) first instances for every dimensionality available in the benchmarks. A set of settings have been individuated by some preliminary experiments, then a full factorial experimental design has been considered in order to choose the best setting. The involved parameters and their values are: $N \in \{20, 50, 100\}, \ \alpha, \beta \in \{1, 2\}, \ \rho \in \{0.05, 0.1, 0.2\}, \ q_0 \in \{0, 0.01, 0.1\}, \ \text{and} \ \left(\Delta_{\tau}^{ib}, \Delta_{\tau}^{gb}\right) \in \{(10, 0), (7.5, 2.5)\}.$ Therefore, a total of 216 settings have been tested by performing 10 executions per instance with a termination criterion of 60 seconds. Then, the average rank of the ARPDs obtained in every instance are computed and the setting with the best average rank is chosen as the reference configuration of ACOP. This setting is $(N=20,\alpha=2,\beta=2,\rho=0.05,q_0=0.1,(\Delta_{\tau}^{ib},\Delta_{\tau}^{gb})=(7.5,2.5)).$

The tuned setting has been used both for ACOP and ACOP⁺. The further parameter r of ACOP⁺ has been set to r = 50. Then, ACOP and ACOP⁺ have

³ The instances are available from http://www.optsicom.es/lolib.

⁴ During the years and using a considerably large amount of computational time, they have been proved to be optima using exact methods [16].

been executed 20 times on every instance. The termination criteria adopted are: 120 seconds for LOLIB instances, 300 seconds for SGB instances, and 600 seconds for the larger MB instances. All the experiments have been run on a homogeneous cluster of computers equipped with Intel Xeon X5650 processors clocking at 2.67GHz. The SR, ARPD and the median time where the global best solution has been found are reported in Tables 1 (LOLIB) and 2 (SGB and MB).

Table 1. Experimental Results on LOLIB instances

Instance	T	ACOP			$ACOP^+$			Instance		ACOP			ACOP ⁺		
	SB	ARPD	_				l	Name		$_{\rm SR}$		Time		ARPD	
-	1			<u> </u>			11								
		0.0004			0	0.133		N-t75d11xx			0.0009	4.145		0	0.089
	1 100		0.319		0	0.038		N-t75e11xx		1-00	0	0.854	100	0	0.046
	1 100		0.303		0	0.061		N-t75i11xx		1-00	0	1.463	100	0	0.227
	1 100		0.140		0	0.024		N-t75k11xx			0		100	0	0.024
N-t59n11xx 44	1 100	0	0.026	100	0	0.020		N-t75n11xx	44	100	0	0.138	100	0	0.022
N-t65b11xx 44	1 0	0.0163	1.003	100	0	0.115		N-t75u11xx	44	100	0	0.196	100	0	0.024
N-t65d11xx 44	1 100	0	0.467	100	0	0.056		N-be75eec	50	100	0	0.553	100	0	0.081
N-t65f11xx 44	1 100	0	0.195	100	0	0.023		N-be75np	50	0	0.0062	18.025	15	0.0002	0.172
N-t65i11xx 44	100	0	0.298	100	0	0.070	ji.	N-be75oi	50	75	0.0003	45.038	80	0.0002	0.090
N-t65l11xx 44	1 100	0	0.007	100	0	0.010	ji.	N-be75tot	50	90	0.0003	1.031	100	0	0.214
N-t65n11xx 44	1 100	0	0.191	100	0	0.051	ii.	N-tiw56n54	56	95	< 0.0001	1.493	100	0	0.780
N-t65w11xx 44	1 100	0	0.345	100	0	0.023	ii.	N-tiw56n58	56	100	0	0.767	100	0	0.312
N-t69r11xx 44	1 100	0	0.067	100	0	0.024	ii.	N-tiw56n62	56	95	0.0011	1.780	95	0.0011	0.183
N-t70b11xx 44	1 100	0	0.295	100	0	0.028		N-tiw56n66	56	100	0	1.618	100	0	0.130
N-t70d11xx 44	1 10	0.0012	0.654	100	0	0.022		N-tiw56n67	56	60	0.0891	8.158	90	0.0218	0.364
N-t70d11xxb 44	1 100	0	0.675	100	0	0.047		N-tiw56n72	56	65	0.0008	28.039	100	0	0.259
N-t70f11xx 44	1 100	0	0.145	100	0	0.043	İ	N-tiw56r54	56	60	0.0017	2.643	95	0.0009	0.637
N-t70i11xx 44	1 100	0	0.227	100	0	0.087	li	N-tiw56r58	56	100	0	1.509	100	0	0.186
N-t70k11xx 44	1 60	0.0041	0.907	100	0	0.039	ii.	N-tiw56r66	56	100	0	1.892	100	0	0.151
N-t70l11xx 44	100	0	0.033	100	0	0.043	ii.	N-tiw56r67	56	55	0.0002	1.610	100	0	0.089
N-t70n11xx 44	100	0	0.150	100	0	0.023	li	N-tiw56r72	56	95	< 0.0001	1.503	100	0	0.101
N-t70u11xx 44	100	0	0.019	100	0	0.019	li	N-stabu70	60	0	0.0242	4.791	80	0.0052	1.670
	1 100		0.322		Õ	0.023		N-stabu74	60	35	0.0160	5.059	100	0	0.913
	1 100		0.464		0	0.023		N-stabu75	60	15	0.0418	4.303			
		0.0010			0	0.046		N-usa79	79	10	0.0306	25.779		0.0051	
							LC	OLIB Avera	ge	80	0.0047	3.432	96	0.0007	0.316

Tables 1 and 2 clearly show that both ACOP and ACOP⁺ obtained remarkable performances throughout all the instances of the benchmark suites considered. Regarding the success rates, ACOP obtained the optimum in at least one execution (SR>0) on about the 57% of the instances, while ACOP⁺ reached the instance optimum on all the 105 instances. Moreover, in 63 cases, ACOP⁺ reached the optimum in all the executions performed (SR=100). Most notably, also when the optimum is not reached, the very small ARPDs clearly show that both ACOP and ACOP⁺ have been able to obtain very high quality solutions. Indeed, the worst ARPD of ACOP, obtained in the N-sgb75.19 instance (see Table 2), is of only the 0.1057%, while for ACOP⁺ it is even smaller, i.e., 0.0218% in N-tiw56n57 (see Table 1). Furthermore, though the computational time to reach the best solution increases with the instance size n, the average times reported at the end of the tables show that the two algorithms are able to provide high quality solutions in a reasonable amount of time.

 $ACOP^+$ Instance ACOP Instance ACOP $ACOP^+$ SR ARPD Time SR Name ARPD Time Name SR ARPD Time SR. ARPD Time N-sgb75.01 75 0.0670 17.789 50 0.0063 2.983 N-r100a2 100 0 0.0155 29.261 0.0005 10.407 75 0.0551 33.936 35 N-sgb75.02 0 0.0002 6.568 N-r100b2 100 0 0.0224 31.812 0.005410.828 N-sgb75.03 0.0197 25.826 100 1.570 N-r100c2 0.0375 40.122 0.005313.498 100 N-sgb75.04 N-r100d2 0.0225 28.953 1.509 0.0003 6.515N-sgb75.05 0 0.0413 41.994 100 1.252 N-r100e2 100 10 0.0015 27.703 0.0002 7.730 N-r150a0 N-sgb75.06 75 0 0.0291 23.84370 0.0001 8.384 150 70 0.0004 84.635 100 29.493 N-sgb75.07 0.0311 18.408 100 0.932 N-r150a1 0 0.0157 137.665 0.0005 58.211 0.0341 28.849 100 N-r150b0 150 8.287 N-sgb75.08 75 2.774 65 0.0002 74.576 N-sgb75.09 75 0 0.0166 69.700 50 0.001668.278 N-r150b1 150 0 0.0051 115.006 0.0019 36.030 N-sgb75.10 75 0 0.0413 54.677 100 N-r150c0 150 55 0.0007 84.107 100 2.720 13.054N-sgb75.11 $0.0841\ 28.972$ 0.0003 48.659 N-r150c1 150 $0.0051\ 129.176$ 95 0.0001 47.817 N-sgb75.12 $0.0272\ 14.141$ 100 2.418 N-r150d0 150 0 $0.0044\ 102.742$ N-sgb75.13 75 0.0125 28 362 30 0.0001 4 894 N-r150d1 150 0 0.0079 140.940 35 0.0011 72.893 N-r150e0 N-sgb75.14 75 0.0258 46.338 3.622 100 71.879 60 0.0001 150 0 100 7.742N-sgb75.15 0.0850 34.499 90 0.0009 N-r150e1 150 0.0097 139.383 0.0001 27.44150.005 N-sgb75.16 75 0.0379 16.061 3.171 N-r200a0 200 $0.0016\ 340.860\ 100$ 100 N-sgb75.17 75 0.0545,44,567 100 0 1.640 N-r200a1 200 0 0.0042.403.498 $95 < 0.0001 \ 140.697$ 0.0435 54.402 0.0001 N-sgb75.18 75 N-r200b0 200 0.0024 335.645 100 70 2.671 178.688 0 $0.1057\ 19.253$ N-sgb75.19 75 < 0.0001 4.057N-r200b1 0.0099 374.202 221.687 10 N-sgb75.20 0.0415 14.087 < 0.0001N-r200c0 200 $0.0029 \ 305.989$ 45 87.9785 70 0.0003 N-sgb75.21 75 0 0.0453 18 634 75 < 0.00016.069 N-r200c1 200 0 0.0036 349 472 100 0 89 4125 0.0839 37.686 100 N-sgb75.22 N-r200d0 200 0.0015 317.184 100 126.250 75 1.579 0 N-sgb75.23 $0.0320\ 73.893$ 0.0005 4.519 N-r200d1 0.0160 459.413 0.0025 N-sgb75.24 75 0.0505 65.582 70 N-r200e0 25 0.0004 276.738 100 0.0001 6.877 200 0.0004 N-sgb75.25 75 0.0613 40.212 80 0.0019 2.638 N-r200e1 200 20 0.0016 306.774 55 140.123 N-r250a0 < 0.0001 148.343 250 15 0.0009 522.939 95 $0.0006\ 528.481$ 75 $< 0.0001\ 369.338$ N-r250c0 250 10 0.0007 528 663 100 0 0.0034 557.957 95 < 0.0001 318.197 N-r250d0 250 N-r250e0 250 0 0.0024 556.871 $80 < 0.0001 \ 361.753$

Table 2. Experimental Results on SGB (left) and MB (right) instances

Finally, a comparison with the ACO algorithm for LOP proposed in [18], namely ACS-IM, has been performed. The results for ACS-IM have been directly taken from its original paper [18], while ACOP and ACOP⁺ have been run for 20 executions on their same set of 49 instances (old non-normalized LOLIB instances⁵). The termination criterion has been set to 50 000 iteration as in [18]. The ARPD results are provided in Table 3.

MB Average | 16 | 0.0060 | 246.602 | 72 | 0.0006 | 101.683

8.792

The comparison reported in Table 3 clearly shows that ACOPs perform largely better than ACS-IM, thus promoting our proposal as the first prominent ACO approach to the linear ordering problem.

7 Conclusion and Future Work

SGB Average | 0 0.0460 35.227 | 79 0.0005

ACOP, a new precedence-based ACO algorithm for permutation problems, has been proposed.

With respect to other proposals in literature, the main novelty of ACOP is to consider a permutation as a total order which is obtained through an incremental refinement of a, initially empty, partial order. The refinement process works by

⁵ Non-normalized LOLIB instances are available at https://www.iwr.uni-heidelberg.de/groups/comopt/software/LOLIB.

Instance $n \mid ACOP \mid ACOP \mid ACS-IM \mid Instance \mid n \mid ACOP \mid ACOP \mid ACS-IM \mid$ t59b11xx 44 | 0.0097 0.08 t75d11xx 44 0.0018 t59d11xx 44 0.0043 0 0.03t75e11xx 44 0.0009 0 0.21t59f11xx 44 0.00610 0.02t75i11xx0.0108o 0.0544 t59i11xx44 0.0001 0 0.06 t75k11xx44 0.0163 o 0.02 t59n11xx 44 0 0.19t75n11xx 44 0 0.040 0 0.0405 t65b11xx 44 0 0.09 t75u11xx 440.00070 0.08 t65d11xx44 0.00660 0.18 be75eec 50 0.00040 0.16 t65f11xx 44 0.00160 0.14 be75np 50 0.0089 0.00020.0004 0.0102 0.0005 t65i11xx0 0.19be75oi 0.004 44 50 0.0023t65l11xx44 0 0 0.03 be75tot 50 0.0037 0.0001 0.12 t65n11xx 44 0.04410 0.16 tiw56n54 56 0.0039 0.13 t65w11xx 440 tiw56n58o 0.0033 0.1456 0.0024 0.15t69r11xx 44 tiw56n62 56 O O 0.410.0068O 0.08t70b11xx 44 0.0043 0 0.03 tiw56n66 560.0057O 0.13t70d11xn 44 0.0095 0 0.05 tiw56n67 56 0.1623 0.450.0002 t70d11xx 44 0.0005 0 0.13 tiw56n72 56 0.0106 0.17 t70f11xx 44 O O 0.16 tiw56r54 56 0.00570.0003 0.19 t70i11xx44 O 0 0.15tiw56r5856 0.00850 0.16t70k11xx 44 0.00840 0.05tiw56r6656 0.0086o 0.09 tiw56r67 0 0.39 t70111xx 44 0 0 0.2456 0.0101 t70n11xx 44 O 0 0.1tiw56r72 56 0.0630 O 0.11 t70u11xx 44 0.00150 0.07stabu1 0.06650.0015 0.26 60 t70w11xx 44 0.00210 0.04stabu2 60 0.07140 0.27 stabu3

Table 3. Experimental Comparison with ACS-IM on non-normalized LOLIB instances

iteratively adding up a selected precedence relations together with the induced precedences. Also the pheromone model and the heuristic values are defined on the precedence relations.

0.0600

0

0.0141 < 0.0001

0.27

0.1454

60

Average

0.02

0.24

This approach is particularly suited for those permutation problems where the precedence relations play an important role, for instance in the linear ordering problem (LOP). An ACOP implementation for LOP is then proposed and experimentally validated on a wide set of popular LOP benchmark instances. ACOP is competitive with the state-of-the-art results and clearly outperform the previous ACO proposals for LOP.

Future research directions are: a thorough investigation of the pheromone update strategies in ACOP, the application to other permutation problems and the proposal of a ACO scheme for problems where the solutions are partial orders.

References

t70x11xx 44

t74d11xx 44 0.0051

0.0026

0

0

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