# Ant colony optimization and the vehicle routing problem 

Tuomas Pellonperä

University of Tampere
School of Information Sciences
Computer Science
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Supervisor: Erkki Mäkinen
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University of Tampere
School of Information Sciences
Computer Science
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#### Abstract

Ant Colony Optimization algorithms are swarm intelligence algorithms, and they are inspired by the behavior of real ants. They are well suited to solving computational problems which involve traversing graphs. The Vehicle Routing Problem is a combinatorial optimization problem which is studied in the field of operations research. Its numerous variants have several real-life applications. In this thesis, I will present how Ant Colony Optimization algorithms have been used to solve a particular variant of the Vehicle Routing Problem - the Vehicle Routing Problem with Time Windows.


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## 1 Introduction

### 1.1 Vehicle routing problem

The gist of the vehicle routing problem (VRP) is how to distribute goods, taking most out of the available resources. More concretely, the VRP is about transporting, in the most optimal and cost-effective way, items between depots and customers by means of a fleet of vehicles [24]. There are many variants of the VRP, and the more complicated ones are able to model accurately many problems arising in real-life situations. Hence, VRP has numerous real-life applications, such as school bus routing, street cleaning, municipal solid waste collection, dial-a-ride systems, routing of salespeople, heating oil distribution, milk delivery, mail pickup and delivery, routing of maintenance units, transportation of disabled people, currency delivery to ATM machines, prisoner transportation between jails and courthouses, beverage delivery to bars and restaurants, and so forth $[9,24,27]$. Its practical benefits can be profound. They can be environmental (reduced exhaust gas emissions due to reduced fuel consumption), or financial (reduced transportation costs, lower consumer prices, and optimal business resource usage).

Ever since the VRP problem was introduced in the late 1950s, it has gained more and more attention in the operations research community. In fact, the growth of the number of published articles in peer-reviewed scientific journals has been exponential. There is already a large enough body of literature for the VRP to allow it be considered a separate and distinct field of knowledge [9]. These facts alone are a convincing evidence of the vitality and the significance of the VRP problem.

### 1.2 Ant colony optimization

Ant colony optimization (ACO) algorithms are swarm intelligence algorithms, and they are inspired by the behavior of real ants looking for good sources of food. First presented in the early 1990s, they were originally used to solve computational problems involving traversal of graphs. There are many ACO algorithms, and the ACO metaheuristic was defined so as to provide a common characterization of a new class of algorithms, and a reference framework for the design of new instances of ACO algorithms [8]. There is a body of knowledge about the
theoretical properties of some members of the ACO algorithm family. There is a dedicated website ${ }^{1}$ which brings the research community together, and a moderated mailing list ${ }^{2}$ on which information regarding the latest developments may be exchanged and discussed [8]. There are a biannual conference on Ant Colony Optimization and Swarm Intelligence ${ }^{3}$, and the IEEE Swarm Intelligence Symposium series, and many journal issues dedicated to the subject [8]. To summarize, the ACO metaheuristic has been a well-established research topic for many years now [8].

The domain of ACO algorithms is vast [7], and they have been applied to dozens of different $\mathcal{N} \mathcal{P}$-hard computational problems, often achieving a state-of-theart performance $[4,8]$. Originally, the first ACO algorithm - the Ant System - was applied to the Traveling Salesman Problem [8]. Since then, they have been applied to, for instance, routing problems, scheduling problems, assignment problems, subset problems, machine learning problems, protein folding, DNA sequencing, and routing packets in telecommunications networks [4, 7, 8]. ACO algorithms may be a particularly viable technique for solving "ill-structured" or highly dynamic problems [8].

All in all, ACO metaheuristic and ACO algorithms are viable and effective tools, and worthy of consideration, in solving hard computational problems.

### 1.3 Road map

In this thesis, I will present a brief overview of the VRP, introducing two fundamental versions of the problem. I will cover the ACO Metaheuristic in detail, presenting three influential and significant members of the ACO algorithm family, and some of their theoretical properties. Finally, I will show how ACO algorithms have been used to solve a particular variant of VRP (i.e., the Vehicle Routing Problem with Time Windows).

The structure of this thesis is as follows. In Chapter 2, some background material on graph theory and computational complexity theory will be presented upon which the material in the subsequent chapters will build. In Chapter 3, the

[^0]vehicle routing problem, and a chosen few of its many variants, will be properly introduced. Having introduced the problem, I will present one technique to solve it. In Chapter 4, the longest chapter of this thesis, the most significant members of the ant colony optimization algorithm family will be introduced. In Chapter 5, I will cover how the problem at hand (VRP) has been solved with our technique of choice (ACO). Finally, in Chapter 6, some conclusions will be drawn.

## 2 Preliminaries

In this chapter, I will make known some theoretical tools which aid in comprehending and clarifying the material to be presented in the upcoming chapters of this thesis.

### 2.1 Graph theory

The problem formulation of the VRP contains the term graph, and the ACO algorithms were developed as a technique for solving computational problems which can be reduced to finding good paths through graphs. Therefore, elementary knowledge of graph theory will help grasp both the problem and its solution on a more accurate level. In what follows, a few very basic definitions of graph theory will be presented. Unless cited otherwise, all the definitions are taken from [15].

A (finite) graph $G$ consists of two finite sets $V$ and $E$. An element $v$ of $V$ is called a vertex ${ }^{1}$, and an element $e$ of $E$, a pair of vertices, is called an edge. If one wants to emphasize that a vertex set $V$ belongs to the graph $G$, one can write $V(G)$; similarly, if one wants to emphasize that an edge set $E$ belongs $G$, one can write $E(G)$. If every edge in $E(G)$ is an unordered pair of two vertices, the graph $G$ is said to be undirected; if every edge of $E(G)$ is an ordered pair of two vertices, the graph $G$ is said to be directed. ${ }^{2}$ The order of a graph is the cardinality of its vertex set; the size of a graph is the cardinality of its edge set.

Assume $e \in E(G)$ and $e=\{u, v\}$ for some $u, v \in V(G) .{ }^{3}$ For notational convenience, $e$ is usually denoted simply as $u v$. The vertices $u$ and $v$ are called the end vertices of $e$, and they are said to be adjacent. Furthermore, it is said that $u$ (similarly for $v$ ) is incident with $e$. If $\{u, v\} \notin E(G)$, then the vertices $u$ and $v$ are non-adjacent. If $u=v$, then $u v$ is called a loop. ${ }^{4}$ The neighborhood of a vertex $v$, denoted by $N(v)$, is the set of vertices adjacent to $v$. The neighborhood of a set $S$, denoted by $N(S)$, is the union of the neighborhoods of the vertices in $S$. The degree of a vertex $v$, denoted by $\operatorname{deg}(v)$, is the number of edges incident

[^1]with $v$. A graph is said to be complete if every vertex is adjacent to every other vertex.

A walk in a graph is a sequence of vertices $v_{1}, v_{2}, \ldots, v_{k}$ such that $v_{i} v_{i+1} \in E$ for all $i=1,2, \ldots, k-1$, and $v_{1}$ and $v_{k}$ are called the end vertices of the walk. If all the vertices in a walk are distinct, the walk is called a path. If all the edges in a walk are distinct, the walk is called a trail. A closed path, or a cycle, is a path $v_{1}, v_{2}, \ldots, v_{k}(k \geq 3)$ together with an edge $v_{k} v_{1}$. A trail which begins and ends at the same vertex is called a closed trail, or a circuit. The length of a walk (or a path, a trail, a cycle, a circuit) is the number of edges, counting repetitions. A graph is connected if every pair of vertices can be joined by a path.

A Hamiltonian path is a path which contains every vertex of $G$. Similarly, a Hamiltonian cycle is a cycle which contains every vertex of $G$ [16]. The graph $G$ is called Hamiltonian if it contains at least one Hamiltonian cycle [19].

Given a graph $G$, the weight function $w$ maps the edges of $G$ to non-negative real numbers. A weighted graph is a pair $(G, w)$.

### 2.2 Computational complexity theory

The purpose of this section is to explain what it basically means for a computational problem to be labelled as $\mathcal{N} \mathcal{P}$-hard. The exposition will be informal.

There are three categories of computational problems - decision problems, search problems, and optimization problems. A model of computation defines the set of allowable operations used in computation and their respective costs. Though there are several alternative models of computation - such as Turing machine, register machine, random-access machine, lambda calculus, combinatory logic the Church-Turing and Cobham-Edmonds theses state that they are fundamentally equivalent in their capabilities and "cost-effectiveness" [14]. Computational problems are further divided into complexity classes, depending on how much computational resources - time and space - their solutions require on a fixed model of computation [22].

Before introducing few specific complexity classes by name, I have to clarify what an "efficient" algorithm means. Briefly, an algorithm the time complexity of which is upper-bounded by a polynomial in the length of the input, is considered efficient. If the time complexity of an algorithm is not bounded by a polynomial
with respect to the length of the input, such an algorithm is considered ineffective or intractable. ${ }^{5}[13,14]$ This is the case when, for instance, an algorithm performs an exhaustive search to come up with a solution to a problem (assuming that the size of the search space is superpolynomial). A problem is efficiently solvable if there is an algorithm which can solve an arbitrary instance of the problem efficiently; a problem has an efficiently checkable solution if there is an algorithm which can check the correctness of a solution efficiently.

Now the two fundamental complexity classes will be introduced. The class $\mathcal{P}$ consists of problems which are efficiently solvable. The class $\mathcal{N} \mathcal{P}$ consists of problems which have an efficiently checkable solution. The most important unsolved problem in computational complexity theory - and in theoretical computer science in general - is what is called the " $\mathcal{P}$ versus $\mathcal{N P}$ problem"; that is, is it true that $\mathcal{P}=\mathcal{N} \mathcal{P}$. The crux of the matter is that if a computer can efficiently check the correctness of a solution, can it efficiently find the solution, too.

A computational problem $\Pi_{A}$ is reducible to another problem $\Pi_{B}$ if an instance of $\Pi_{A}$ can be solved efficiently with the help of a solution to an instance of $\Pi_{B}$. Thus, a reduction transforms one problem into another, preserving the polynomial time bound. A problem $\Pi$ is said to be $\mathcal{N} \mathcal{P}$-complete if $\Pi \in \mathcal{N} \mathcal{P}$ and if every problem in $\mathcal{N} \mathcal{P}$ is reducible to $\Pi$. [13] A problem $\Pi$, with numerical parameters, is said to be $\mathcal{N} \mathcal{P}$-complete in the strong sense if it remains so even when all of its numerical parameters are bounded by a polynomial in the length of the input [12].

One of the most startling discoveries of the complexity theory is the fact that if an efficient algorithm is found to any one of the $\mathcal{N} \mathcal{P}$-complete complete problems ${ }^{6}$, then by definition, every problem in $\mathcal{N P}$ would be efficiently solvable.

A problem $\Pi$ is said to be $\mathcal{N} \mathcal{P}$-hard if every problem in $\mathcal{N} \mathcal{P}$ is reducible to $\Pi$. In addition, a problem $\Pi$ is said to be $\mathcal{N} \mathcal{P}$-hard in the strong sense if a strongly $\mathcal{N} \mathcal{P}$-complete problem has a polynomial reduction to it [12]. It is noteworthy that $\Pi$ need not be a member of $\mathcal{N} \mathcal{P}$. Hence, $\Pi$ may not be efficiently checkable. Basically, $\Pi$ is (computationally) as hard a problem as any problem in $\mathcal{N P}$, and

[^2]possibly even harder. The vehicle routing problem is $\mathcal{N} \mathcal{P}$-hard in the strong sense [27].

To summarize, if a problem is $\mathcal{N} \mathcal{P}$-hard, then there is no known algorithm which can solve an arbitrary instance of the problem efficiently. Given a "large enough" instance of the problem, an algorithm producing an optimal solution will require an infeasible amount of computing resources.

### 2.3 Traveling salesman problem

I will briefly mention the traveling salesman problem (TSP), as it is related both to the vehicle routing problem, and to the ant colony optimization. Besides, it is an important problem in theoretical computer science in its own right.

TSP can be formulated as a graph problem as follows. Given a complete weighted graph - whose vertices represent cities, edges represent roads, and weights represent costs or distances - find a Hamiltonian cycle with the least weight. Stated as a graph problem, the TSP is $\mathcal{N} \mathcal{P}$-hard [19]; stated as a decision problem, the TSP is $\mathcal{N} \mathcal{P}$-complete [13].

The TSP is related to the VRP in that the latter is a generalization of the former [27], whereas it is related to the ACO algorithms in the sense that it was among the first computational problems to be solved with those algorithms, and it is often used as a benchmark to test new ideas in, and new variants of, them [8].

## 3 Vehicle Routing Problem

The structure of this chapter is as follows. First, the basic components of the VRP will be introduced. Then two significant variations of the VRP will be covered in depth one at a time; the other variants will be skimmed over. Finally, other techniques to solve the VRP, besides the ACO metaheuristic, will be mentioned.

### 3.1 Basic components

The term "distributing goods," mentioned in Section 1.1, is at the heart of the VRP, and it is succinctly clarified by Toth and Vigo [27]:

> "The distribution of goods concerns the service, in a given period of time, of a set of customers by a set of vehicles, which are located in one or more depots, are operated by a set of crews (drivers), and perform their movements by using an appropriate road network."

It is carried out by devising a set of routes meeting certain requirements [27]:


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" $[T]$ he solution of a VRP calls for the determination of a set of routes, each performed by a single vehicle that starts and ends at its own depot, such that all the requirements of the customers are fulfilled, all the operational constraints are satisfied, and the global transportation cost is minimized."


Next, each of these basic components will be briefly introduced.

The road network. This component, used for the transportation of goods, is usually described as a weighted graph. The vertices represent "locations" i.e., road junctions, the depot(s), and the customers - whereas the arcs represent connections between locations. Each arc is associated with a cost, and a travel time.

The customers. A customer has several properties:

- the vertex in the road graph where the customer is located;
- the amount of goods (the demand) which must be delivered to, or collected at, the customer;
- the interval during which the customer must be served (the time window);
- the time required to serve the customer (the loading time, or the unloading time);
- the subset of vehicles which can serve the customer.

The vehicles. A vehicle has several properties:

- the home depot;
- the capacity of the vehicle (for instance, the maximum weight the vehicle can carry);
- devices available for loading and unloading;
- the subset of arcs of the road graph which are traversable by the vehicle;
- the cost associated using the vehicle (for instance, per time unit, or distance unit).

The drivers operating the vehicles must satisfy several constraints imposed by the legal system. For example, a driver needs to have a license to drive a particular type of vehicle, and to have proper breaks during the work day and enough rest between work days. From now on, the constraints on drivers will be ignored; instead, it is assumed that they are embedded into the constraints on vehicles.

The routes. A multitude of constraints are imposed on the routes. Their exact nature depends on the type of goods being transported, and on the characteristics of the vehicles and the customers. For instance, the capacity of the vehicle must not be exceeded on any route; customers must be served during their time windows; the driver taking care of the route must not exceed the allowed limit on working hours; the customers may need to be served in a particular order (precedence constraints).

The global transportation cost. This objective of the VRP depends on the global distance traveled, and on the costs arising from the usage of vehicles (their number, their type, etc.).

### 3.2 Graph-theoretic formulation

With the basic vocabulary clarified, I will now formulate the general VRP as a graph-theoretic problem.

Let $G=(V, A)$ be a complete graph, where $V=\{0,1, \ldots, n\}$ is the vertex set and $A$ is the arc set. Vertices $j=1,2, \ldots, n$ correspond to the customers, each with a known non-negative demand, whereas the vertex 0 corresponds to the depot. A non-negative cost, $c_{i j}$, is associated with each $\operatorname{arc}(i, j) \in A$, representing the cost of traveling from vertex $i$ to vertex $j$. If $c_{i j}=c_{j i}$, for all $i, j \in V$, then the problem is said to be a symmetric VRP; otherwise, it is called asymmetric VRP. Usually, the cost matrix satisfies the triangle inequality: $c_{i k}+c_{k j} \geq c_{i j}$, for any $i, j, k \in V$. The VRP consists of finding a collection of $k$ simple circuits, each corresponding to a vehicle route with minimum cost, defined as the sum of the costs of the circuits' arcs such that:
i. each circuit visits the vertex 0 (the depot)
ii. each vertex $j \in V \backslash\{0\}$ is visited by exactly one circuit; and
iii. the sum of vertices' demand visited by a circuit does not exceed the vehicle capacity $C$. [9]

The different variants of VRP extend and build on this formulation, adding some additional constraints to the problem statement.

### 3.3 Capacitated VRP

The least general and the most fundamental version of the VRP is the "Capacitated VRP" (CVRP) [27]. In this version, there are no pickups, only deliveries to customers. The demand of a customer is invariable, and it is known in advance. There is only one depot, and all the vehicles are identical (their capacities are equal). The only constraints are those imposed on the capacities of the vehicles.

More formally, the CVRP may be described as a graph theoretic problem of the following kind. Let $G=(V, A)$ be a complete weighted graph, where $V=$ $\{0,1, \ldots, n\}$. The depot is the vertex 0 ; other vertices are customers. The weight $c_{i j}$ associated with each $\operatorname{arc}(i, j) \in A$ is a non-negative real number, usually called the cost (of traveling along the arc). If $c_{i j}=c_{j i}$, for every $i$ and $j$, the CVRP is called symmetric (SCVRP); otherwise, the it is called asymmetric (ACVRP).

Loops are not allowed in $G$, and they can be prohibited by setting $c_{i i}=\infty$, for all $i \in V$. It is assumed that arcs obey the triangle inequality; that is, $c_{i k}+c_{k j} \geq c_{i j}$, for all $i, j, k \in V$. For each customer $i(1 \leq i \leq n)$, there is an associated nonnegative demand $d_{i}$ which is known a priori and is invariable. For the depot, we set $d_{0}=0$. Given a subset $S$ of customers (i.e., $S \subseteq V \backslash\{0\}$ ), let $d(S)=\sum_{i \in S} d_{i}$ denote the total demand of the customer subset. The fleet of vehicles consists of $K$ identical vehicles, each with a capacity $C$. To make the problem feasible, it is implicitly assumed that $\max \left\{d_{i}\right\} \leq C, i=1,2, \ldots, n$, and that $K$ is large enough to serve all the customers. Neither the time window properties nor the service time properties will be taken into consideration in CVRP.

The CVRP may now be stated formally as follows [27].
Find exactly $K$ simple circuits - i.e., routes - with minimum cost, defined as the sum of the costs of the arcs belonging to the circuits, such that
(i) each circuit visits the depot vertex;
(ii) each customer vertex is visited by exactly one circuit; and
(iii) the sum of the demands of the vertices visited by a circuit does not exceed $C$.

The CVRP is $\mathcal{N} \mathcal{P}$-hard in the strong sense. Therefore, it is not known whether there is an algorithm which can solve it in polynomial time.

It is noteworthy that by setting $K=1$ and assuming that $C \geq d(V)$, the CVRP becomes the TSP. Thus, the CVRP can be regarded as a generalization of the TSP. [27]

There are several variants of the basic CVRP. For example, the vehicles may be heterogeneous; i.e., their capacities may differ. In the so called DistanceConstraint VRP (DVRP), the route capacity constraints are replaced by the constraints on the maximal lengths of the routes. In the so called DistanceConstrained CVRP (DCVRP), both the vehicle capacity constraint and the maximum distance constraint are present. [27]

### 3.4 VRP with Time Windows

The VRP with Time Windows (VRPTW) is an extension of the CVRP in which capacity constraints are imposed and each customer $i$ is associated with a time
interval $\left[a_{i}, b_{i}\right]$ (called a time window), and an additional service time $s_{i}[1,27]$. Each customer must be serviced within the associated time window, and the service itself takes $s_{i}$ time units. The time window requirements essentially impose an implicit orientation on each of the routes. Hence, VRPTW is usually modelled as an asymmetric problem. [27]

The depot is actually represented by two nodes: 0 and $n+1$. Their time windows are equal; namely, $\left[a_{0}, b_{0}\right]=\left[a_{n+1}, b_{n+1}\right]=\left[a_{\alpha}, b_{\omega}\right]$, where $a_{\alpha}$ and $b_{\omega}$ represent the earliest possible departure from the depot and the latest possible return to the depot, respectively. [1] The travel time for each $\operatorname{arc}(i, j) \in A$ is denoted by $t_{i j}$. Feasible solutions exist only if

$$
a_{0}=a_{\alpha} \leq \min _{i \in V \backslash\{0\}}\left\{b_{i}\right\}-t_{0 i}
$$

and

$$
b_{n+1}=b_{\omega} \geq \min _{i \in V \backslash\{0\}}\left\{a_{i}\right\}+s_{i}+t_{i 0}
$$

The demands $d_{0}$ and $d_{n+1}$, and service times $s_{0}$ and $s_{n+1}$, associated with the depot nodes, are all equal to 0 . In case a vehicle arrives at the service location of the customer $i$ ahead of time, it is allowed to wait until the time instant $a_{i}$ before servicing the customer [27].

VRPTW can be stated as follows [27]:

Find exactly $K$ simple circuits with minimum cost, such that
(i) each circuit visits the depot vertex;
(ii) each customer vertex is visited by exactly one circuit;
(iii) the sum of the demands of the vertices visited by a circuit does not exceed $C$; and
(iv) for each customer $i$, the service starts within the time window $\left[a_{i}, b_{i}\right]$, and the vehicle stops for $s_{i}$ time instants.

The VRPTW is $\mathcal{N} \mathcal{P}$-hard in the strong sense, and it generalizes the CVRP: by setting $a_{i}=0$ and $b_{i}=+\infty$, for every $i \in V \backslash\{0\}$, VRPTW turns into CVRP [27]. Interestingly enough, even finding a feasible solution to the VRPTW is an $\mathcal{N} \mathcal{P}$ hard problem [1,24].

### 3.5 Other variants

I will briefly mention other major variants of the VRP, but as they fall outside the scope of this thesis, the coverage will be shallow.

The VRP with Backhauls (VRPB) is an extension of the CVRP in which the customer set is partitioned into two subsets [28]. The first subset, $L$, denotes Linehaul customers, each having a certain demand to be delivered; the second subset, $B$, denotes Backhaul customers, each having a certain demand to be picked up. There is a precedence constraint between the Linehaul and Backhaul customers: a route serving both types of customers must serve all the Linehaul customers before serving any of the Backhaul customers. There are symmetric (VRPB) and asymmetric (AVRPB) versions of the VRPB, and they both are $\mathcal{N} \mathcal{P}$ hard in the strong sense [28], generalizing the SCVRP and ACVRP, respectively. [27]

In the VRP with Pickup and Delivery (VRPPD), each customer $i$ is associated with two quantities $d_{i}$ and $p_{i}$. The quantity $d_{i}$ represents a demand of homogeneous commodities to be delivered, whereas the quantity $p_{i}$ represents a demand of homogeneous commodities to be picked up. [27] A heterogeneous fleet of vehicles, based at multiple terminals, must satisfy a set of transportation requests. Each request is defined by a pickup point, a corresponding delivery point, and a demand to be transported between these locations. [10] The current load of a vehicle, before arriving at a given location, is defined as the initial load subtracted by the demands already delivered, and summed up with the demands already picked up. VRPPD is $\mathcal{N} \mathcal{P}$-hard in the strong sense, and it generalizes the CVRP. [27]

Vehicle routing problems may be further complicated with the presence of heterogeneous vehicle fleets, limitations on customer accessibility, and constraints regarding the order of pickups and deliveries. These variants may be called rich vehicle routing problems. [24] Yet, even these variants may not be able to accurately model highly complex real-life problems. For instance, travel times may be uncertain or dependent on traffic conditions [24]; some of the orders may be received as time progresses [20]. These variants are called Dynamic (or Online) Vehicle Routing Problems (DVRP) [20, 24]. Dynamic problems have arisen due to advances in information and communication technologies which enable information to be obtained and processed in real-time [17].

### 3.6 Other solving techniques

There are various exact algorithms for solving different variants of the VRP, but because of the inherent computational complexity, they do not scale to big, let alone huge, instances. None the less, three different mathematical programming formulations exist for the VRPs presented above: (i) vehicle flow formulations; (ii) commodity flow formulations; and (iii) formulations as a set-partitioning problems [27]. The exact algorithms for the VRP can be classified into three broad categories: (i) direct tree search methods; (ii) dynamic programming; and (iii) integer linear programming [18].

Besides exact algorithms and metaheuristic techniques, there are other methods to solve the VRP, approximation algorithms for instance. However, further examination of these techniques is beyond the scope of this thesis.

## 4 Ant Colony Optimization Algorithms

Ant Colony Optimization is a swarm intelligence technique and a metaheuristic which is inspired by the behavior of real ants in their search for food. A colony is a population of simple, independent, and asynchronous agents that cooperate to find a good solution to the problem at hand [6]. The ants in ACO are stochastic solution construction procedures that probabilistically and iteratively build complete solutions from smaller components [8]. It is this cooperation on a colony level that gives the members of the ACO algorithm family their distinctive flavor.

The structure of this chapter is as follows. First, I will describe the behavior of real ants which the ACO algorithms mimic. Then I will present three highly influential and successful variants of the ACO algorithm family.

### 4.1 Biological inspiration

In the 1940s and 1950s, the French entomologist Pierre-Paul Grassé was among the first scientists to study the social behavior of insects [4,6-8]. His area of expertise was termites. Grassé discovered that they are capable of reacting to what he termed "significant stimuli" - signals which activated a genetically encoded response. He observed further that these reactions could act as a new significant stimuli for both the insect that produced them and other members of the colony. He coined the term stigmergy to describe this indirect communication in which agents are stimulated by the performance they have achieved. The stigmergic communication is characterized by its non-symbolic and local nature; i.e., the non-symbolic information can be accessed only close to the location where it was originally released. [6]

Examples of stigmergy can observed in ant colonies. The "communication medium" is pheromone, a chemical which ants can smell and which they deposit on the ground, and which evaporates at a steady rate. This creates a pheromone trail, and its presence influences an ant's choice of path. When at an intersection of several paths, it has been verified experimentally that an ant tends to choose the path with the most pheromone on it. Goss, Aron, Deneubourg, and Pasteels developed a formula to model this behavior [6]. Assuming that $t$ time units have passed since the beginning of the experiment, and that $m_{1}$ ants have used
the first ${ }^{1}$ bridge and $m_{2}$ ants used the second bridge, the probability $p_{1}$ for the $(m+1)$ th ant to choose the first bridge can be given by

$$
\begin{equation*}
p_{1}=\frac{\left(m_{1}+k\right)^{h}}{\left(m_{1}+k\right)^{h}+\left(m_{2}+k\right)^{h}}, \tag{4.1}
\end{equation*}
$$

where parameters $k$ and $h$ are empirically determined to fit the model to data ${ }^{2}$. Naturally, the probability $p_{2}$ to choose the second bridge is $1-p_{1}$.

As time goes by, shorter paths collect more pheromone than longer ones, which helps ants find good sources of food. There are certain aspects of this foraging behavior which have inspired the behavior of artificial ants. The first one is autocatalysis - exploitation of positive feedback. The second one is called the implicit evaluation of solutions; that is, the shorter paths are completed more quickly, and hence they receive pheromone reinforcement more quickly too. [6, 8] Together, stigmergy, implicit solution evaluation, and autocatalytic behavior give rise to ACO [6]. On one hand, the artificial ants resemble the real ants in a straightforward manner. They traverse graphs, and they lay (artificial) pheromones on problem states. On the other hand, in order to increase the competitiveness of the ACO algorithms in solving demanding computational problems, artificial ants are equipped with capabilities not found in real ants. For example, an artificial ant may possess memory, it can deposit a quantity of pheromone proportional to the quality of the solution produced, and it may use techniques such as look-ahead, local search, and backtracking. [6,8] By and large, new ACO algorithms are less and less biologically inspired and more and more motivated by the aim of improving the algorithmic performance. [8]

### 4.2 ACO metaheuristic

After few ACO algorithms had been published, the ACO metaheuristic was formally defined so as to provide a common characterization of a new class of algo-
${ }^{1}$ Goss et al. conducted what is called a "binary bridge experiment," the purpose of which was to study pheromone trail laying. Ants traversed two bridges of unequal lengths, which connected the ant nest to a food source. In the beginning of the experiment, when no pheromone had accumulated on the paths, the ants showed no bias towards any of the bridges. However, as the pheromone started piling up, the ants clearly favored the shorter path.
${ }^{2}$ Using Monte Carlo simulations, it was discovered that with $k \approx 20$ and $h \approx 2$ a good fit was achieved.
rithms, and a framework for the design of new instances of ACO algorithms. An algorithm which implements the metaheuristic is called an ACO algorithm. [8] The following characteristics are present in all ACO algorithms $[7,8]$ :

- the use of a colony;
- the role of autocatalysis;
- the cooperative behavior mediated by artificial pheromone trails;
- the probabilistic construction of solutions biased by artificial pheromone trails and local heuristic information;
- the pheromone updating guided by solution quality;
- the evaporation of pheromone trails.

What distinguishes one ACO algorithm from another is the implementation of the pheromone update.

In an ACO algorithm, there is an interplay between two techniques commonly found in approximate approaches to hard computational problems; namely, between construction algorithms and local search algorithms. Construction algorithms build solutions to a problem in an incremental way, starting with an empty solution and iteratively adding appropriate components without backtracking until a complete solution is obtained. In contrast, local search algorithms start with a complete initial solution and try to find a better solution in the neighborhood of the current solution (see Code Iterative-Improvement on page 18 for the most basic version of the local search technique). [8] Used in isolation, neither technique is usually able to find a near-optimal solution: a construction algorithm, although fast, does not produce a high quality solution [8], whereas a local search algorithm is often trapped in a local minimum [7]. In an ACO algorithm, the colony of artificial ants is the stochastic construction procedure, which builds a complete solution probabilistically by iteratively adding solution components to a partial solution using as a guide problem-specific heuristic information and artificial pheromone trails. These complete solutions are then optimized with local search algorithms before the pheromone trail is updated.

### 4.2.1 Pseudocode

Before presenting the metaheuristic formally, I will present it in pseudocode (see Algorithm 1) and explain what the different components do.

```
Procedure Iterative-Improvement(s)
    Data: an initial solution \(s \in S\)
    Result: a local optimum \(s_{i}\)
    \(s_{i} \leftarrow\) Improve \((s)\)
    while \(s \neq s_{i}\) do
        \(s \leftarrow s_{i}\)
        \(s_{i} \leftarrow \operatorname{Improve}(s)\)
    end
    return \(s_{i}\)
```

Initialize The parameters are set and the pheromone variables are initialized to the value $\tau_{0}$, which is a parameter of the algorithm.

ConstructAntSolutions A set of $m$ ants constructs solutions. Each ant starts with an empty solution $s_{p}=\emptyset$, and extends it one component at a time.

ApplyLocalSearch Once complete candidate solutions have been obtained, they may be further improved by applying a local search algorithm on them. This is an example of daemon actions, the purpose of which is to implement problem specific or centralized actions which are beyond ants' sphere of responsibility.

UpdatePheromones The purpose of pheromone update is to make solution components of good solutions more desirable for the ants operating in subsequent iterations. The update is implemented using pheromone deposit and pheromone evaporation.

```
Algorithm 1: The ACO Metaheuristic [8]
    Initialize
    while termination condition not met do
        ConstructAntSolutions
        ApplyLocalSearch /* optional */
        UpdatePheromones
    end
```

A slightly different, but nevertheless compatible, version of the ACO metaheuristic is presented in Algorithm 2. The ScheduleActivities construct schedules the three primary components of the algorithm; namely, the management of ants' activity, pheromone update, and daemon actions. It is implementation-dependent how these activities are actually scheduled and synchronized. For instance, they may be executed in parallel, or in a synchronized manner.

```
Algorithm 2: The ACO Metaheuristic [7]
    ScheduleActivities
        ManageAntsActivity()
        EvaporatePheromone()
        DaemonActions() /* optional */
    end
```


### 4.2.2 Problem Representation

I will now present how combinatorial optimization problems (COP) are represented in ACO algorithms. Unless cited otherwise, this section is based on [7] and [6].

A combinatorial optimization problem $\Pi$ is a 3-tuple $(\mathcal{S}, \Omega, f)$, where $\mathcal{S}$ is the search space defined over a finite set of discrete decision variables, $\Omega$ is the set of constraints, and $f: \mathcal{S} \rightarrow \mathbb{R}_{0}^{+}$is the objective function which assigns to each candidate solution $s \in \mathcal{S}$ a cost $f(s, t)^{3}$. The goal is to find a globally optimal solution $s_{\text {opt }}$; i.e., a minimum cost solution satisfying the constraints $\Omega .^{4}$

The search space $\mathcal{S}$ is defined as follows. Given a set of discrete variables $\left\{X_{i}\right\}$ and their respective domains $D_{i}$, where $D_{i}=\left\{v_{i}^{1}, \ldots, v_{i}^{\left|D_{i}\right|}\right\}$, the assignment of value $v_{i}^{j}$ to a variable $X_{i}$ is denoted by $X_{i} \hookleftarrow v_{i}^{j}$. An assignment in which each decision variable is assigned a value from its domain, is called a solution. A solution $s \in \mathcal{S}$ is a feasible solution if it is a complete assignment in which each

[^3]decision variable is assigned a value which satisfies all the constraints in $\Omega$. If the set $\Omega$ is empty, $\Pi$ is said to be unconstrained. If the constraints cannot be violated, they are called hard constraints; if they may be violated (possibly with a penalty), they are called soft constraints.

Some observations and definitions will follow.

- An instantiated decision variable $X_{i} \hookleftarrow v_{i}^{j}$ is called a solution component, and is denoted by $c_{i j}$. The set of all possible solution components is denoted by $\mathcal{C}$.
- Ants traverse what is called the construction graph $G=(V, E)$. The set of components $\mathcal{C}$ may be associated either with the vertices $V$, or with the edges $E$. The graph $G$ is complete and connected.
- The states of the problem are defined in terms of sequences $\left\langle c_{1}, c_{2}, \ldots, c_{k}\right\rangle$ (abbreviated as $\left\langle c_{k}\right\rangle$, for some positive integer $k$ ) of elements of $\mathcal{C}$, and in terms of vertices of $G$. For instance, $\left(\left\langle c_{k}\right\rangle, v\right)$ denotes the state where $v$ is the vertex the ant is at, and the sequence $\left\langle c_{k}\right\rangle$ represents the (partial) solution the ant has built so far. The set of all possible sequences is denoted by $\mathcal{X}$.
- A pheromone trail parameter $T_{i j}$ is associated with each component $c_{i j}$. The value of $T_{i j}$ is denoted by $\tau_{i j}$ and is called pheromone value. ${ }^{5}$ The pheromone values are used and updated during the execution of an ACO algorithm.
- A heuristic value $\eta_{i j}$ may be associated with each component $c_{i j}$. Heuristic values represent a priori knowledge of the problem instance or of the problem domain.
- The set of candidate solutions $\mathcal{S}$ is a subset of $\mathcal{X}$.
- The finite set of constraints $\Omega$ defines the set of feasible states $\overline{\mathcal{X}}$, with $\overline{\mathcal{X}} \subseteq \mathcal{X}$.

[^4]I assume the parameter $t$ to be implicit for the sake of simplicity.

- The set of feasible solutions is denoted by $\mathcal{S}^{\star}$, with $\mathcal{S}^{\star} \subseteq \overline{\mathcal{X}}$, and $\mathcal{S}^{\star} \subseteq \mathcal{S}$.
- The set of optimal solutions is denoted by $\mathcal{S}^{\bullet}$ :

$$
\mathcal{S}^{\bullet}=\left\{s^{\bullet} \in \mathcal{S} \mid f\left(s^{\bullet}\right) \leq f(s), \forall s \in \mathcal{S}\right\} .
$$

Naturally, $\mathcal{S}^{\bullet} \subseteq \mathcal{S}^{\star}$.

- An ant $k$ has a memory $\mathcal{M}_{k}$, which is used for building feasible solutions (i.e., for implementing constraints), for evaluating the solution produced, and in updating pheromones.
- An ant $k$ is assigned a start state $x_{s}^{k}$ and one or more termination conditions $e^{k}$. When at least one termination condition is fulfilled, the construction procedure of the ant $k$ stops.
- When the ant $k$ is in state $\left(\left\langle c_{k}\right\rangle, v\right)$, it tries to move to any vertex $j$ in $N(v)$ (the neighborhood of $v$ ), and in doing so, transitioning to state $\left(\left\langle c_{k+1}\right\rangle, j\right)$. The move is selected using a probabilistic decision rule, which is a function of locally available pheromone trails and heuristic values $\eta$, the private memory $\mathcal{M}_{k}$, and the problem constraints.
- If the ant $k$, after adding a component $c_{i j}$ to its current solution, updates immediately the pheromone trail associated with the component, this is called online step-by-step pheromone update.
- If the ant $k$, after it has finished building its solution, retraces the same path backwards, updating the associated pheromone trails, this is called online delayed pheromone update.

It needs to be emphasized that ants move concurrently and independently. An ant is "intelligent" enough to produce a solution, albeit probably a poor one, on its own, but it is a characteristic of the ACO metaheuristic that good quality solutions emerge as the result of the colony level interaction among the ants via indirect communication.

Now the time is ripe to present three successful ACO algorithms - the Ant System (AS), the Ant Colony System (ACS), and the Max-Min Ant System (MMAS).

### 4.3 Ant System

The Ant System, the first ACO algorithm proposed in the literature [6], is the progenitor of all the research efforts with ant algorithms [5]. As a matter of fact, the AS was originally a set of three algorithms called the ant-cycle, the antdensity, and the ant-quantity [7], but as the ant-cycle turned out to be superior to the other two, performance-wise, it attracted more attention [3]. Thus antcycle would later be called simply the Ant System, whereas ant-density and ant-quantity have gradually fallen into oblivion [8].

The chief characteristic of the AS is that the pheromone values are updated by all the ants which have completed the tour. ${ }^{6}$ After each ant has completed its tour, an ant updates the pheromone values of every solution component which is a part of its tour. More exactly, in ant-cycle, the pheromone $\tau_{i j}$, associated with the edge ( $i, j$ ), is updated as follows:

$$
\begin{equation*}
\tau_{i j} \leftarrow(1-\rho) \cdot \tau_{i j}+\sum_{k=1}^{m} \Delta \tau_{i j}^{k}, \tag{4.2}
\end{equation*}
$$

where $\rho$ is the pheromone evaporation rate, $m$ is the number of ants, and $\Delta \tau_{i j}^{k}$ is the quantity of pheromone laid on the edge $(i, j)$ by the ant $k$ :

$$
\Delta \tau_{i j}^{k}= \begin{cases}Q / L_{k}, & \text { if ant } k \text { used the edge }(i, j) \text { in its tour }  \tag{4.3}\\ 0 & \text { otherwise }\end{cases}
$$

Above, $Q$ is a constant, and $L_{k}$ is the tour length of $k$ th ant.
Ant-density and ant-quantity differ from ant-cycle with regard to how the pheromone trail values are updated. In them, each ant lays its trail at each step; i.e., the pheromone values are updated after an ant has added a new component to its

[^5]tour. Both of them follow Formula 4.2, but they define $\Delta \tau_{i j}^{k}$ differently. In acodensity, it is defined as
\[

\Delta \tau_{i j}^{k}= $$
\begin{cases}Q & \text { if ant } k \text { used the edge }(i, j) \text { in its tour }  \tag{4.4}\\ 0 & \text { otherwise },\end{cases}
$$
\]

whereas in aco-quantity, it is defined as

$$
\Delta \tau_{i j}^{k}= \begin{cases}Q / d_{i j}, & \text { if ant } k \text { used the edge }(i, j) \text { in its tour }  \tag{4.5}\\ 0 & \text { otherwise }\end{cases}
$$

where $d_{i j}$ is the distance between the vertices $i$ and $j$ in the construction graph.
When the ant $k$ is in the component $i$, the transition probability from the component $i$ to the component $j$ is defined as

$$
p_{i j}^{k}= \begin{cases}\frac{\tau_{i j}^{\alpha} \cdot \eta_{i j}^{\beta}}{\sum_{l \in N_{k}(i)} \tau_{i l}^{\alpha \cdot} \cdot \eta_{i l}^{\beta}} & \text { if } j \in N_{k}(i),  \tag{4.6}\\ 0 & \text { otherwise. }\end{cases}
$$

The parameters $\alpha$ and $\beta$ control the relative importance of the pheromone versus the heuristic information $\eta_{i j}$ [6], and $N_{k}(i)$ denotes the neighborhood of the component $i$; i.e., the "allowed" components the ant $k$ can move to when in component $i$ (at the time $t$ ). ${ }^{7}$ Dorigo, Maniezzo, and Colorni [3] call it the tabu list, and tabu ${ }_{k}$ is the tabu list of the ant $k$.

To wrap up, I will quote [3] in order to provide with a summary of the ant-cycle algorithm:
"At time zero an initialization phase takes place during which ants are positioned on different towns and initial values $\tau_{i j}(0)$ for trail intensity are set on edges. The first element of each ant's tabu list is set to be equal of its starting town. Thereafter every ant moves from town $i$ to town $j$ choosing the town to move to with a probability that is a function (with parameters $\alpha$ and $\beta$, see formula (4) [Formula 4.6]) of two desirability measures. The first, the trail $\tau_{i j}(t)$, gives information about how many ants in the past have chosen the same edge $(i, j)$;

[^6]the second, the visibility $\eta_{i j}$, says that the closer a town the more desirable it is. [--]
"After $n$ iterations all ants have completed a tour, and their tabu lists will be full; at this point for each ant $k$ the value of $L_{k}$ is computed and the values $\Delta \tau_{i j}^{k}$ are updated according to formula (3) [Formula 4.3]. Also, the shortest path found by the ants [- -] is saved and all the tabu lists are emptied. This process is iterated until the tour counter reaches the maximum (user-defined) number of cycles $\mathrm{NC}_{\text {max }}$, or all the ants make the same tour."

Although the performance of the AS was promising, it never the less was not on par with the state-of-the-art techniques for solving the TSP; for instance, though it was capable of finding good or optimal solutions only for small instances of TSP, the time required to find such solutions for larger instances was infeasible [5]. Hence further improvements became the focus of research. The first one was called the elitist strategy, which was motivated by the similar idea used in genetic algorithms [3]. The idea is to give the best tour since the start of the algorithm - called $s_{\mathrm{gb}}$ ('gb' standing for "globally-best") - a strong additional weight [8]. In other words, at every cycle the trail laid on the edges belonging to the best-so-far tour is reinforced more than in the standard version in the hope that the trail of the best tour, so reinforced, will direct the search of all the other ants in probability toward a solution composed by some edges of the best tour itself [3]. This additional weight is the quantity $e \cdot Q / L^{*}$, where $e$ is the number of elitist ants, and $L^{*}$ is the length of the best found tour.

Another improvement, called the rank-based Ant System ( $\mathrm{AS}_{\mathrm{rank}}$ ), is a sort of extension of the elitist strategy: it sorts the ants according to the lengths of the tours they have produced and, after each tour construction phase, only the $(w-1)$ best ants and the global-best ant are allowed to deposit pheromone. The $r$ th best ant of the colony contributes to the pheromone update with a weight given by $\max \{0, w-r\}$, while the global-best tour reinforces the pheromone trails with weight $w$. [8]

Although I will present only two improvements and extensions of the AS in the next sections - the Ant Colony System, and the Max-Min Ant System - there are many more of them. Indeed, perhaps the most long-lasting contribution of the AS was to pave the way for, and stimulate, further researchers to develop extensions and improvements [8]. It pioneered the idea of a population of agents
each guided by an autocatalytic process directed by a greedy force [3], and of the synergistic use of cooperation among these simple agents which communicate via a non-symbolic medium [5].

### 4.4 Ant Colony System

The ACS algorithm was expressly devised to improve the efficiency of the AS to solve instances of symmetric and asymmetric TSP [5]. It differs from preceding ant algorithms in three respects: (i) the state transition rule provides a direct way to balance between exploration of new edges and exploitation of a priori and accumulated knowledge of the problem; (ii) the global updating rule is applied only to edges belonging to the best ant tour; and (iii) while ants construct a solution a local pheromone updating rule is applied. [5]

In the following quotation, the developers of the ACS algorithm themselves explain in "layman terms" how the algorithm works [5]:
" $m$ ants are initially positioned on $n$ cities chosen according to some initialization rule (e.g., randomly). Each ant builds a tour (i.e., a feasible solution to the TSP) by repeatedly applying a stochastic greedy rule (the state transition rule). While constructing its tour, an ant also modifies the amount of pheromone on the visited edges by applying the local updating rule. Once all ants have terminated their tour, the amount of pheromone on edges is modified again (by applying the global updating rule). As was the case in ant system, ants are guided, in building their tours, by both heuristic information (they prefer to choose short edges), and by pheromone information: An edge with a high amount of pheromone is a very desirable choice. The pheromone updating rules are designed so that they tend to give more pheromone to edges which should be visited by ants."

The effect of the best ant depositing its pheromone at the end of iteration $t$ is defining a "preferred tour" for the following algorithm iteration $t+1$. During iteration $t+1$, the ants will regard the edges belonging to the best tour as highly desirable and will choose them with high probability. However, thanks to local updating "eating" pheromone away, and the probability mentioned above not being too high, the exploration of new, possibly better tours in the neighborhood of the previous best tour remains likely enough. [5]

The ACS algorithm is presented in pseudo-code in Algorithm 3.

```
Algorithm 3: ACS
    Initialize
    /* At this level each loop is called an 'iteration' */
    while (EndCondition \(\neq\) true)
    \{
        Each ant is positioned on a starting node
        /* At this level each loop is called an 'step' */
        repeat
        \{
            /* Ants incrementally build a solution */
            Each ant applies the state transition rule
            Each ant applies the local pheromone updating rule
        \}
        until (all ants have built a complete solution)
        Apply global pheromone updating rule
    \}
```


### 4.4.1 Components

The state transition rule. The following formula, which appears in the state transition rule, is basically the Formula (4.6) with $\alpha$ set to value 1:

$$
p_{i j}^{k}= \begin{cases}\frac{\tau_{i j} \cdot \eta_{i j}^{\beta}}{\sum_{l \in N_{k}(i)} \tau_{i l} \cdot \eta_{i l}^{\beta}} & \text { if } j \in N_{k}(i),  \tag{4.7}\\ 0 & \text { otherwise. }\end{cases}
$$

An ant $k$ positioned on node $r$ chooses the city $s$ to move to by applying the following rule:

$$
s= \begin{cases}\underset{u \in N_{k}(r)}{\arg \max }\left\{\tau_{r u} \cdot \eta_{r u}^{\beta}\right\} & \text { if } q \leq q_{0} \text { (exploitation) }  \tag{4.8}\\ S & \text { otherwise (biased exploration) }\end{cases}
$$

where $q$ is a random variable uniformly distributed in $[0,1], q_{0}$ is a parameter with $0 \leq q_{0} \leq 1$, and S is a random variable selected according to the probability given in Formula (4.7). The state transition rule arising from the Formulas (4.7)

```
Algorithm 4: ACS-3-opt
    Initialize
    /* At this level each loop is called an 'iteration' */
    while (EndCondition \(\neq\) true)
    \{
        Each ant is positioned on a starting node
        /* At this level each loop is called an 'step' */
        repeat
        \{
            /* Ants incrementally build a solution */
            Each ant applies the state transition rule
            Each ant applies the local pheromone updating rule
        \}
        until (all ants have built a complete solution)
        Bring each ant to local minimum with 3 -opt heuristic
        Apply global pheromone updating rule
    \}
```

and (4.8) is called the pseudo-random-proportional rule. It favors transitions towards nodes connected by short edges and with a large amount pheromone. The parameter $q_{0}$ determines the balance between exploitation and exploration. [5]

The global updating rule. The "peculiarity" of the ACS is that only the ant which has produced the globally best solution so far, is allowed to deposit pheromone. ${ }^{8}$ With the pseudo-random-proportional rule, this has the effect of allocating a greater amount of pheromone on shorter tours, and of making the search more directed: during the iterations to follow, ants search in the neighborhood of the best tour. [5]

The pheromone level update performed at the end of an iteration uses the following formula:

$$
\begin{equation*}
\tau_{i j} \leftarrow(1-\rho) \cdot \tau_{i j}+\rho \cdot \Delta \tau_{i j}, \tag{4.9}
\end{equation*}
$$

[^7]where $\rho$ is the pheromone decay parameter $(0<\rho<1)$,
\[

\Delta \tau_{i j}= $$
\begin{cases}{\left[L_{\mathrm{best}}\right]^{-1},} & \text { if the edge }(i, j) \text { belongs to the best tour },  \tag{4.10}\\ 0 & \text { otherwise },\end{cases}
$$
\]

and $L_{\text {best }}$ is either $L_{\mathrm{gb}}$ (the length of the globally best tour since the start of algorithm execution), or $L_{\mathrm{ib}}$ (the length of the best tour found during the current iteration of the algorithm). [5, 6]

The local updating rule. While building a solution (to an instance of TSP), ants visit edges and change their pheromone level by applying the following rule:

$$
\begin{equation*}
\tau_{i j} \leftarrow(1-\rho) \cdot \tau_{i j}+\rho \cdot \Delta \tau_{i j} \tag{4.11}
\end{equation*}
$$

where $0<\rho<1$ is a parameter. The term $\Delta \tau_{i j}$ may be defined several different ways:

1. $\Delta \tau_{i j}=\gamma \cdot \max _{z \in N_{k}(r)}\left\{\tau_{s z}\right\}$, taking inspiration from one of the reinforcement learning algorithms - the Q-learning algorithm [5] (this version is named as "Ant-Q");
2. $\Delta \tau_{i j}=\tau_{0}$, where $\tau_{0}$ is the initial pheromone level (this version is named as "ACS- $\tau$ " in this thesis); and
3. $\Delta \tau_{i j}=0$ (this version is named as "ACS-0" in this thesis).

According to experiments, ACS-0 performs worse than Ant-Q and ACS- $\tau$. Since Ant-Q is more computation-intensive, $\mathrm{ACS}-\tau$ has become the most popular version. [5]

Besides these three versions, there is yet another version of ACS; that is, not to use the local update rule at all (just like in the AS). However, the absence of the local update rule degrades the overall performance of the ACS algorithm. The performance experiments mentioned above suggest this, as does the following assertion. As the edges are being visited by the ants, the application of the local update rule makes their pheromone diminish, which makes them less and less attractive, thus encouraging the exploration of other edges. Consequently, the ants never converge to a common path (a phenomenon known as "stagnation").

### 4.4.2 Performance

How did ACS succeed in its stated goal of enhancing the performance of the AS algorithm? Dorigo and Gambardella [5] ran two sets of experiments: in the first set, they compared the ACS to other heuristics; in the second one, they tested the ACS on larger problems. In both sets, the problems were instances of the TSP.

Compared to other heuristics - simulated annealing, elastic net, self-organizing map - the ACS almost always offered the best performance. In order to solve large instances of the TSP, the ACS needs to resort to some "tricks of trade"; namely, to using "candidate lists," a common technique in solving demanding instances of the TSP. Indeed, the ACS with a candidate list ${ }^{9}$ was able to produce good results for problem instances comprising of more than 1,500 cities. [5]

Further performance improvements were achieved when the ACS was modified to imitate "ad-hoc heuristics" geared specifically towards solving instances of the TSP. A "tour improvement heuristic" begins from a given, complete tour and attempts to reduce its length by exchanging edges chosen according to some heuristic rule. The most used and well-known tour improvement heuristic is called a "3-opt" [5]. Dorigo and Gambardella added the 3-opt heuristic to the ACS, resulting in what is called "ACS-3-opt." (ACS-3-opt is presented in pseudocode in Algorithm 4.) The results obtained with ACS-3-opt applied to instances of the TSP, both symmetric and symmetric ones, are "impressive" and comparable to other algorithms "considered good" [5].

### 4.5 Max-Min Ant System

### 4.5.1 Motivation

The key to achieving a good performance for ACO algorithms is to exploit the best solutions found during the search to a far greater extent than was done in the AS. This can be justified by analyzing the shape of the search space of many

[^8]combinatorial optimization problems. [26] Next, I will cover this background in a greater detail, relying on [26] as a reference.

Central to the search space analysis is the concept of fitness landscape. It determines the shape of the search space as encountered by the local search algorithm, and it is defined by

1. the set of all possible solutions $S$;
2. an objective function that assigns a fitness value $f(s)$ to every $s \in S$;
3. a neighborhood structure $N \subseteq S \times S$.

The neighborhood structure induces a distance metric on the set of solutions; if $s_{1}$ and $s_{2}$ are solutions, then the distance $d\left(s_{1}, s_{2}\right)$ between them can be defined as the minimum number of moves - or transformations - that have to be performed so as to transform $s_{1}$ into $s_{2}$.

The distribution of local minima and their relative location with respect to global optima are an important determining factor for the effectiveness of adaptive multistart algorithms, such as ACO algorithms. If $s$ is an arbitrary solution, and if $s_{\text {opt }}$ is an optimal solution, then the values $f(s)$ and $d\left(s, s_{\text {opt }}\right)$ and the correlation between them help analyze the shape of the fitness landscape. In the literature on genetic algorithms, this correlation is customarily called the fitness distance correlation (FDC), and it is captured by the correlation coefficient:

$$
\begin{equation*}
\rho(F, D)=\frac{\operatorname{Cov}(F, D)}{\sqrt{\operatorname{Var}(F) \cdot \operatorname{Var}(D)}}, \tag{4.12}
\end{equation*}
$$

where $F$ and $D$ are random variables describing the fitness and the distance of local optima to a global optimum, respectively, Var denotes the variance, and Cov denotes the covariance. If there is a high, positive correlation between the solution cost and the distance to a global optimum (assume the context of a minimization problem), it indicates that the smaller the cost, the shorter the distance to a global optimum. Hence, if a problem exhibits a high FDC, algorithms combining adaptive solution generation with a local search may be expected to perform well. Conversely, if there is no correlation, or if there is a negative correlation, the fitness value provides hardly any guidance towards better solutions. In such a situation, ACO algorithms probably perform poorly.

The symmetric TSP is one of the most widely studied problems with regard to search space analysis. There is a strong empirical indication that the solution cost and the distance from the closest optimal solution are strongly positively
correlated in the TSP. ${ }^{10}$ This, in turn, indicates that locally optimal tours are concentrated around a small region of the entire search space. Therefore, it makes sense to "focus" ants' attention and efforts to the best tours (either iteration-best or globally-best). It is this observation which is behind many of the design choices of MMAS.

### 4.5.2 Algorithm

Stützle and Hoos [26] describe the requirements the MMAS was designed to meet as follows:
"Research on ACO has shown that improved performance may be obtained by a stronger exploitation of the best solutions found during the search and the search space analysis [- -] gives an explanation of this fact. Yet, using a greedier search potentially aggravates the problem of premature stagnation of the search. Therefore, the key to achieve best performance of ACO algorithms is to combine an improved exploitation of the best solutions found during the search with an effective mechanism for avoiding early search stagnation."

There are three key differences between the AS and the MMAS.

1. To exploit the best solutions found during an iteration (iteration-best), or during the run of the algorithm (algorithm-best), only one single ant adds pheromone after each iteration.
2. To avoid search stagnation, the range of possible pheromone trails on each solution component is limited to the interval $\left[\tau_{\min }, \tau_{\max }\right]$.
3. At the start of the algorithm execution, all pheromone trails are initialized to $\tau_{\text {max }}$ in order to achieve higher exploration of the search space.
[^9]Pheromone trail updating. In MMAS, only one single ant will update the pheromone trails after each iteration. The pheromone update rule is given by

$$
\begin{equation*}
\tau_{i j} \leftarrow \rho \cdot \tau_{i j}+\Delta \tau_{i j}^{\text {best }}, \tag{4.13}
\end{equation*}
$$

where $\Delta \tau_{i j}^{\text {best }}=1 / f\left(s_{\text {best }}\right)$ and $f\left(s_{\text {best }}\right)$ denotes the solution cost of either the iteration-best $\left(s_{\mathrm{ib}}\right)$ or the global-best $\left(s_{\mathrm{gb}}\right)$ solution. The difference between the ACS and the MMAS is that the ACS tends to favor $s_{\mathrm{gb}}$, whereas the MMAS tends to favor $s_{\mathrm{ib}}$. Choosing one over the other has an effect on how the search history is exploited. As iteration-best tours tend to vary from one iteration to the other, the designers of the MMAS reasoned that by favoring $s_{\mathrm{ib}}$, the risk of search stagnation is reduced.

It is possible, and maybe worthy of further research, to combine both approaches: for example, $s_{\mathrm{ib}}$ can be the default choice, but $s_{\mathrm{gb}}$ could be used instead at every pre-determined iteration. Or perhaps their weighted sum could be used:

$$
\begin{equation*}
\Delta \tau_{i j}^{\text {best }}=w_{1} \cdot f\left(s_{\mathrm{ib}}\right)+w_{2} \cdot f\left(s_{\mathrm{gb}}\right), \tag{4.14}
\end{equation*}
$$

where $0 \leq w_{1}, w_{2} \leq 1$ and $w_{1}+w_{2}=1$.
Stützle and Hoos [26] ran experiments comparing updating pheromone values, on one hand, with $s_{\mathrm{ib}}$, and on the other hand, with $s_{\mathrm{gb}}$. Their results indicated that the average performance when using the iteration-best solution was superior to that of global-best solution. Thus, they recommend not to use $s_{\mathrm{ib}}$ exclusively for the pheromone trail update. However, they note that for large problem instances, it may be beneficial to alternate $s_{\mathrm{ib}}$ and $s_{\mathrm{gb}}$.

Pheromone trail limits. Regardless of how one chooses between $s_{\mathrm{gb}}$ and $s_{\mathrm{ib}}$, search stagnation can nonetheless take place. This occurs if the pheromone trail on one component is significantly stronger than on the other components, when the ants are to choose their next solution component. In such a situation, the ants construct the same solution over and over again. If pheromone trail differences are prevented from becoming too substantial, search stagnation is less likely to occur. To achieve this goal, the MMAS imposes special limits on pheromone values: $\tau_{\min } \leq \tau_{i j} \leq \tau_{\max }$, for all pheromone trail values $\tau_{i j}$. Assuming that $[x]_{b}^{a}$ is defined as:

$$
[x]_{b}^{a}= \begin{cases}a & \text { if } x>a  \tag{4.15}\\ b & \text { if } x<b, \\ x & \text { otherwise }\end{cases}
$$

then by combining Formula (4.13) with Formula (4.15), the pheromone trail update rule can be expressed as

$$
\begin{equation*}
\tau_{i j} \leftarrow\left[\rho \cdot \tau_{i j}+\Delta \tau_{i j}^{\text {best }}\right]_{\tau_{\min }}^{\tau_{\max }} \tag{4.16}
\end{equation*}
$$

Stützle and Hoos [26] ran empirical tests on four instances of the symmetric TSP, comparing the approach with pheromone trail lower limits to one without such limits. They report that for all instances, the average solution quality is always better if pheromone trail lower limits are used. Hence they conclude that they are definitely advantageous. In addition, they provide some guidelines on how to choose the values $\tau_{\min }$ and $\tau_{\min }$ but I will not cover them here.

Pheromone trail initialization. In the MMAS, the pheromone trails are initialized in such a way that after the first iteration, all pheromone trail values $\tau_{i j}$ are equal to $\tau_{\max }$. The motive behind this is to increase the exploration of solutions - i.e., different paths in the construction graph - during the first iterations of the algorithm.

Stützle and Hoos compared empirically this approach to one involving the initialization of pheromone values to the lower limit $\tau_{\min }$. They found that the proposed trail initialization is superior to the alternative approach, and concluded that the higher exploration of the search space achieved in this way was important to achieving a better solution quality. [26]

Pheromone trail smoothing (PTS). PTS is an additional mechanism to enhance the performance of MMAS. ${ }^{11}$ Before introducing it, some terms need to be defined. A choice point is any point of time during the execution of the algorithm at which an ant has to decide which solution component to add to its partial solution. The notion of convergence for the MMAS is an undesirable situation where, for each choice point, one of the solution components has $\tau_{\text {max }}$ as its pheromone trail value, whereas all the other components have $\tau_{\min }$ as their pheromone trail value. In such a situation, the solution produced by the ants at the end of an iteration usually equals the best solution found by the algorithm. Convergence differs from stagnation in a subtle way: the latter describes the

[^10]situation where all the ants follow the same path; this does not happen in the convergence of the MMAS thanks to the use of pheromone trail limits.

When the MMAS has converged, or is at the verge of converging, the PTS mechanism increases the pheromone trails proportionally to their difference to the maximum pheromone trail limit:

$$
\begin{equation*}
\tau_{i j}^{*}(t)=\tau_{i j}(t)+\delta\left(\tau_{\max }(t)-\tau_{i j}(t)\right), \tag{4.17}
\end{equation*}
$$

with $0<\delta<1$. The point is to increase exploration by increasing the probability of selecting solution components with low pheromone values. The information gathered during the run of the algorithm is not lost, but merely weakened.

To conclude, Stützle and Hoos [26] state that
"PTS is especially interesting if long runs are allowed, because it helps achieving a more efficient exploration of the search space. At the same time, PTS makes MMAS less sensitive to the particular choice of the lower pheromone trail limit."

### 4.6 Performance comparison

Having introduced several major variants of the ACO algorithm family, I will compare their performance on solving some instances of the TSP. Table 4.1 is an abridged version of Table 6 compiled by Stützle and Hoos [26], and their analysis and observations are summarized in what follows. ${ }^{12}$ I have left out two variations of the AS which originally appeared in the comparison mentioned above, as I have not covered them in this thesis. In Table 4.1, "opt" indicates the known optimal solution value for the instance in question; " $\mathrm{AS}_{\text {rank }}+\mathrm{pts}$ " is a modification of the rank-based Ant System, which utilizes pheromone trail smoothing mechanism. ${ }^{13}$

The AS has the worst performance. This is hardly an unexpected discovery, since the sole purpose of other ACO algorithms is to improve the overall performance.

[^11]| Instance | opt | MMAS+pts | MMAS | ACS | AS $_{\text {rank }}$ | AS $_{\text {rank }}+\mathrm{pts}$ | AS |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- | :--- |
| eil51 | 426 | 427.1 | 427.6 | 428.1 | 434.5 | 428.8 | 437.3 |
| kroA100 | 21282 | 21291.6 | 21320.3 | 21420.0 | 21746.0 | 21394.9 | 22471.4 |
| d198 | 15780 | 15956.8 | 15972.5 | 16054.0 | 16199.1 | 16025.2 | 16702.1 |
| ry48p | 14422 | 14523.4 | 14553.2 | 14565.4 | 14511.4 | 14644.6 | 15296.4 |
| ft70 | 38673 | 38922.7 | 39040.2 | 39099.0 | 39410.1 | 39199.2 | 39596.3 |
| kro124p | 36230 | 36573.6 | 36773.5 | 36857.0 | 36973.5 | 37218.0 | 38733.1 |
| ftv170 | 2755 | 2817.7 | 2828.8 | 2826.5 | 2854.2 | 2915.6 | 3154.5 |

Table 4.1: Performance comparison of ACO algorithms on solving instances of the TSP.

By and large, the MMAS achieves the best performance on all the instances except for "ry48p", where it is outperformed by the $\mathrm{AS}_{\text {rank. }} \cdot{ }^{14}$ Overall, the ACS has the second best performance with respect to the solution quality on the symmetric instances; on asymmetric instances, the MMAS and ACS have a similar performance. This performance difference between these two ACO algorithms may attributable to the fact that the MMAS does not center the search around $\mathrm{s}_{\mathrm{gb}}$ as strongly as the ACS does.

Stützle and Hoos combined the MMAS with a local search heuristic in order to investigate its performance on large instances of the TSP, and to compare it to the algorithms with best performance. For symmetric instances, the 3 -opt local search algorithm was used; for asymmetric ones, a restricted form of 3 -opt was used (called reduced 3 -opt). Based on their preliminary experiments, they decided to gradually shift the emphasis from $s_{\mathrm{ib}}$ to $s_{\mathrm{gb}}$ for the pheromone trail update. The goal is to transition smoothly from stronger exploration of the search space (early iterations) to stronger exploitation of the overall best solution (later iterations). Based on their experiments, Stützle and Hoos came to the conclusion that the MMAS was able to find solutions of high quality for all the instances, and that it was able to find the optimal solution at least once for each instance. In addition, they noted that the computational results, and the computation times, were better when local search was used. All in all, they claimed that the MMAS was the best performing ACO algorithm for the TSP at the time of the writing.

[^12]
### 4.7 Theoretical Results

Empirical results have shown that ACO algorithms are efficient at solving a variety of combinatorial optimization problems; however, there used to be very little theory to explain why. Convergence proofs ${ }^{15}$ appeared several years after the first variants of ACO algorithms had been presented. Gutjahr was among the first people to prove the convergence to the globally optimal solution with probability $1-\epsilon$. The downside of his proof was that is was applicable only to a particular ACO algorithm - called the Graph-Based Ant System (GBAS) - and hence it could not be generalized to other ACO algorithms. Moreover, there was no empirical data available about the performance of GBAS. [25]

Fortunately, the limitations of the Gutjahr's proof were overcome by Stützle and Dorigo [25], who discovered a convergence proof which applies directly to the ACS and MMAS. To prove the convergence, they introduced an algorithm class $A C O-\tau_{\text {min }}$, and a particular member of that class, called $A C O-g b-\tau_{m i n}$. They went on to prove the convergence for $A C O-g b-\tau_{\min }$, and then they generalized the proof so that is applies to every member of $A C O-\tau_{\min }$. (Naturally, they proved that MMAS and ACS are members of $A C O-\tau_{\text {min }}$.)

In what follows, I will present the $A C O-g b-\tau_{\min }$ algorithm, define the $A C O-\tau_{\text {min }}$ algorithm class, and present the theorems and propositions of the convergence proof. ${ }^{16}$ Unless otherwise cited, my presentation is based on the article by Stützle and Dorigo [25].

### 4.7.1 $A C O-g b-\tau_{\text {min }}$ and $A C O-\tau_{\text {min }}$

The presentation to follow will make use of the terminology and notation introduced in Subsection 4.2.2.

The procedure 'ACO-Gb-Tau-Min' (page 37) gives a high-level description of $A C O-g b-\tau_{\text {min }}$. In the pseudocode, $\hat{s}$ will denote the best feasible solution found so far, whereas $s_{t}$ will denote the best feasible solution found during the $t$ th iteration of the algorithm.

[^13]```
Procedure ACO-Gb-Tau-Min \(\left(\alpha, \tau_{0}, \tau_{\text {min }}\right)\)
    Initialize \(\left(\tau_{0}\right) \quad / *\) See Algorithm 5. */
    while (EndCondition \(\neq\) true)
    \{
        AntSolutionConstruction \((\alpha)\) /* See Algorithm 6. */
        PheromoneUpdate \(\left(\tau_{\min }\right)\) /* See Algorithm 7. */
    \}
```

After the initialization (Algorithm 5), the algorithms ' 6 ' and ' 7 ' are repeated until the termination condition is true.

```
Algorithm 5: Initialize \(\left(\tau_{0}\right)\)
    Data: For the parameter \(\tau_{0}\) it holds \(\tau_{\min } \leq \tau_{0}<+\infty\).
    \(\hat{s} \leftarrow\) a feasible solution
    \(\tau_{i j} \leftarrow \tau_{0} \quad / * \forall(i, j) \in E(G) \quad\) */
    for (every ant \(a_{i}\) )
    \{
        assign a start vertex \(c_{i}\) for \(a_{i}\)
        \(\mathrm{k} \leftarrow 1\)
        \(x_{k} \leftarrow\) a single element sequence \(\left\langle c_{i}\right\rangle\)
    \}
```

```
Algorithm 6: AntSolutionConstruction( \(\alpha\) )
    Data: \(0<\alpha<+\infty\)
    while (SolutionComplete \(\neq\) true)
    \{
        choose next solution component using Formula (4.6) with \(\eta_{i j}^{\beta}=1\)
    \}
```

During the pheromone value update, all the pheromone values are first decreased by a factor of $\rho$ (the evaporation rate); then the components which are part of $\hat{s}$ have their pheromone values reinforced ("the global best offline pheromone update"); finally, it is made certain that none of the pheromone values are less than the specified limit $\left(\tau_{\text {min }}\right)$.

```
Algorithm 7: PheromoneUpdate \(\left(\tau_{\text {min }}\right)\)
    Data: \(\tau_{\text {min }}>0\)
    \(\tau_{i j} \leftarrow(1-\rho) \cdot \tau_{i j} \quad / * \forall(i, j) \in E(G) \quad * /\)
    if \(\left(f\left(s_{t}\right)<f(\hat{s})\right)\)
    \{
        \(\hat{s} \leftarrow s_{t}\)
    \}
    /* For \(g: \mathcal{S} \rightarrow \mathbb{R}^{+}\)it holds \(f(s)<f\left(s^{\prime}\right) \Rightarrow g(s) \geq g\left(s^{\prime}\right)\). */
    for \((\) every \((i, j) \in \hat{s})\)
    \{
        \(\tau_{i j} \leftarrow \tau_{i j}+g(\hat{s})\)
    \}
    for \((\) every \((i, j) \in E(G))\)
    \{
        \(\tau_{i j} \leftarrow \max \left\{\tau_{i j}, \tau_{\text {min }}\right\}\)
    \}
```


### 4.7.2 Convergence proof

Two theorems will be proved for the $A C O-g b-\tau_{\min }$ algorithm. First, it will be proved that it is guaranteed to find an optimal solution, if given enough time, with a probability which can be made arbitrarily close to 1 . Second, it will be proved that after a fixed number of iterations $t_{0}$ has elapsed since the discovery of an optimal solution, the pheromone trails on the connections of the optimal solution are larger than those on any other connection.

Proposition 1 ([25]). For any $\tau_{i j}$, the following holds:

$$
\lim _{t \rightarrow \infty} \tau_{i j}(t) \leq \tau_{\max }=\frac{1}{\rho} \cdot g(s \cdot)
$$

where $s^{\bullet}$ is an optimal solution.

Proposition 1 states that the maximum possible pheromone level $\tau_{\max }$ is bounded asymptotically.

Proposition 2 ([25]). After an optimal solution has been found, remembering that $\tau_{i j}^{*}(t) \geq \tau_{\text {min }}$ and that the global best pheromone update rule is used, we have

$$
\forall(i, j) \in s^{\bullet}: \lim _{t \rightarrow \infty} \tau_{i j}^{\bullet}(t)=\tau_{\max }=\frac{1}{\rho} \cdot g\left(s^{\bullet}\right)
$$

Proposition 2 states that the pheromone trail values of all the components of $s^{\bullet}$ converge to $\tau_{\text {max }}$.

Theorem 3 ([25]). Let $P^{\bullet}(t)$ be the probability that the algorithm finds an optimal solution at least once within the first $t$ iterations. Then, for an arbitrary small $\epsilon>0$ and for a sufficiently large $t$, it holds that

$$
P^{\bullet}(t) \geq 1-\epsilon,
$$

and asymptotically

$$
\lim _{t \rightarrow \infty} P^{\bullet}(t)=1
$$

Theorem 4 ([25]). Let te be the iteration when the first optimal solution has been found. Then a value $t_{0}$ exists, called the transition period, such that the following holds:
$\forall(i, j) \in s^{\bullet}, \quad \forall(k, l) \notin s^{\bullet}, \quad \forall t>t^{\bullet}+t_{0}=t^{\bullet}+[(1-\rho) / \rho]$.

Theorems 3 and 4 seal the convergence proof. Theorem 3 states that the probability of finding an optimal solution can be made arbitrary large, and that "as time goes by," - i.e., as $t$ becomes arbitrarily large - the algorithm will eventually find an optimal solution. Theorem 4 states that once an optimal solution has been found ${ }^{17}$, the pheromone values on its components will eventually become larger that the pheromone values on the other components. ${ }^{18}$ Thus, the algorithm will converge as all the ants will choose the same paths, producing an identical solution.

Proposition 5 ([25]). Once an optimal solution has been found, then for any $(i, j) \notin s^{\bullet}$ :

$$
\lim _{t \rightarrow \infty} \tau_{i j}(t)=\tau_{\min }
$$

Proposition 5, together with transition probability rules such as Formula (4.6), explicates why the ants gradually steer clear from the non-optimal paths: as the pheromone values of components decrease, so does the probability that an ant will choose that component as a part of its solution. Thus these components do

[^14]not receive any pheromone reinforcement, but are consumed by the pheromone evaporation rate.

Theorem 3 applies to any ACO algorithm for which the probability $P(s)$ of constructing a solution $s \in \mathcal{S}$ always remains greater than a small positive constant $\epsilon$. An ACO algorithm is a member of the class $A C O-\tau_{\text {min }}$ if 1) there is a minimum value $\tau_{\min }$, explicitly set, for the pheromone trail values; 2 ) there is a limit to the amount of pheromone which an ant may deposit after each iteration; 3) and if the pheromone evaporates over time (i.e., $\rho>0$ ). By definition, Theorem 3 holds for any algorithm in $A C O-\tau_{\text {min }}$.

### 4.7.3 ACO algorithms in $A C O-\tau_{\text {min }}$

Of the several ACO algorithms covered thus far, which ones are members of $A C O-\tau_{\text {min }}$ ?

The ancestor of all the ACO algorithms, the AS, is not a member of $A C O-\tau_{\text {min }}$. Neither is the $A_{\text {rank }}$, an elitist variant of the AS. [25]

It is relatively straighforward to show that the MMAS is a member of $A C O-\tau_{\text {min }}$. By Formula (4.16), it is clear that there are minimum and maximum limits to the pheromone values. Moreover, the pheromone evaporation rate is positive. Thus, Theorem 3 holds for the MMAS, and with some minor adaptions, so does Theorem 4. [25]

It is less straightforward to show that the ACS is a member of $A C O-\tau_{\text {min }}$. As the maximum of amount of pheromone is limited, as any feasible solution can be constructed with a nonzero probability, and because an algorithm parameter corresponds to $\tau_{\min }$, it can be observed that Theorem 3 holds for the ACS. Although determining the transition period $t_{0}$ for the ACS is complicated, it can be concluded that Theorem 4 holds for the ACS, too. [25]

As far as employing local search techniques and heuristic information are concerned, they do not have an effect on the convergence [25].

It is unfortunate that the convergence proof does not shed light on the speed with which the algorithms converge towards an optimal solution [8].

## 5 Ant Colony Optimization and Vehicle Routing Problem

In this chapter, I will discuss how the ACO metaheuristic has been applied to solve the VRPTW. First I will present some general guidelines and observations which need to be taken into consideration when applying ACO algorithms to computational problems. Then, an influential ACO algorithm - MACS-VRPTW - will be presented along with two of its variants.

### 5.1 Application Principles

When ACO algorithms are being applied to combinatorial optimization problems other than TSP, few basic issues must be addressed [8].

One issue, of crucial importance, is the definition of the solution components and the pheromone model. To illustrate the differences in the definition of solution components, I will compare the TSP to any job scheduling problem ${ }^{1}$. Although both problems represent solutions as permutations, they define the components differently. In TSP, the two permutations $\pi_{1}=(1,2, \ldots, n)$ and $\pi_{2}=(2,3, \ldots, n, 1)$ are essentially the same; however, in a job scheduling problem, they represent two different solutions. Therefore, in the context of TSP, a solution component $X_{i}$, taking on values $j=1,2, \ldots, n$, could model the fact the node $i$ precedes the node $j$ in the solution to be produced, whereas in the context of a job scheduling problem, it could model the fact the job $i$ is assigned to position $j$ in the job sequence. [8]

Another issue involves balancing the exploitation of the previous search experience and the exploration of unvisited, or relatively unexplored, search space regions $[7,8]$. The tradeoff between them has an effect on the solution quality and on the computation time. Too much exploration and not enough exploitation may render an ACO algorithm into a plain brute-force algorithm, thus improving the solution quality at the cost of increasing the running time; too much exploitation and not enough exploration may cut down on the running time, but at the cost of worsened solution quality, and at a higher risk of search stagnation. Typically,

[^15]ACO algorithms aim to strike this balance through the management of pheromone trails. A stronger exploitation of the previous search experience can be achieved by allowing the best solution(s) found during the search to contribute strongly to pheromone trail updating ${ }^{2}$. To further the exploration of the search space, various methods have been devised. For instance, in ACS, the local pheromone update rule makes already trodden paths less desirable for future ants; in MMAS, the explicit lower pheromone limit guarantees a minimal level of exploration. [8] In addition, it is possible to manipulate the values of $\alpha$ and $\beta$ (see Formula (4.6), for example) to shift the emphasis towards either exploration or exploitation [7].

Yet another issue involves taking advantage of problem and domain-specific knowledge in the form of heuristic information. This information guides the ants' solution construction process. However, it should be noted that the significance of the heuristic information is greatly reduced if local search methods are used to improve solutions, which is usually the case. $[7,8]$

### 5.2 MACS-VRPTW

MACS-VRPTW stands for "Multiple Ant Colony System for Vehicle Routing Problem with Time Windows" [11]. The VRPTW is regarded as multiobjective optimization problem with two objective functions: i) a function to minimize the number of vehicles used; ii) a function to minimize the total travel time. The minimization of the number of vehicles takes precedence over the minimization of the total travel time; i.e., a solution with a smaller number of vehicles is always preferred to a solution with a higher number of vehicles but a shorter total travel time. MACS-VRPTW employs two ant colonies; one colony (called ACS-VEI) to diminish the number of vehicles, and one colony (called ACS-TIME) to optimize the total travel time, and to optimize the solutions produced by ACSVEI. The two colonies are independent, possessing distinct pheromone trails, but they collaborate via a shared variable into which the currently best solution is stored. [11,24]

The three main components of MACS-VRPTW are procedures macsvrp, acstime, and acsvei. The procedure macsvrp ${ }^{3}$ is the entry-point to the algorithm, and after the initialization phase and the computing of a feasible solution, it repeatedly

[^16]activates the procedures acstime and acsvei, waiting for an improved solution to be found. If such a solution is found, it saves it, deactivates the active procedures, and repeats the process until some stopping criterion has been fulfilled.

The procedure acstime ${ }^{4}$ implements a traditional ACS-based colony. As a parameter, it receives a number indicating the smallest number of vehicles for which a feasible solution has been computed [24]. It basically waits for the ants to produce their solutions, and then checks whether any of the feasible ones requires less travel time than the currently best solution. If so, this solution is sent to the procedure macsvrp. The solution serving the most customers is sent to the procedure macsvrp. [11] The procedure acstime serves the purpose of refining the solutions produced by acsvei which has no feel for the time, since its objective function is independent of it [24].

The procedure acsvei ${ }^{5}$ is very similar, but with two differences. First, the parameter it receives is a number indicating the number one less than the smallest number of vehicles for which a feasible solution has been computed [24]; second, rather than paying attention to the travel time of a feasible solution, it keeps an eye on how many customers are served instead. The more customers a solution serves, the better. [11] It bears pointing out that the ants of acsvei usually construct infeasible solutions (not all customers can be visited due to the constraints given) [24].

The constructive step of both ant colonies is based on a procedure which computes the set of all feasible nodes reachable from the current node $i$. A node $j$ is a member of this set if it is unvisited and if the arrival at $j$ is compatible with the time window requirements. In order to take the time windows into account, distance between nodes becomes a function of time: a node appears to be "closer" if the end of its time window is near. Once the ants have finished building their solutions, one algorithm iteration has reached its end. The pheromone trails are updated both locally and globally in accordance with the ACS algorithm. The pheromone trails are updated globally for two different solutions: (i) for the infeasible solution with the highest number of visited customers; and (ii) for the feasible solution with the lowest number of vehicles and the shortest travel time. [24]

[^17]Gambardella et al. [11] state that at the time of writing, MACS-VRPTW was very competitive, in terms of solution quality and computation time, with the best known methods of the time, and that it was able to produce relatively good solutions in a short time. Furthermore, they say that one of the innovations of MACS-VRPTW was to utilize, as a new methodology for optimizing multiobjective problems, multiple ant colonies whose activity is coordinated through information exchange, and each of whom is responsible for optimizing a different objective [11]. Other sources have confirmed the merits of of MACS-VRPTW. For example, Rizzoli et al. [24] state that it "is the most efficient ACO algorithm for the VRPTW and one of the most efficient metaheuristics overall for this problem [VRPTW]".

### 5.3 ANTROUTE

ANTROUTE is an implementation of MACS-VRPTW with two modifications (i) at the start of each tour, an ant chooses the type of vehicle to be used; (ii) a waiting cost was introduced in order to prevent vehicles from arriving too early at the stores. [24]

ANTROUTE was applied to a real-life distribution problem in Switzerland. The task was to distribute 52,000 pallets to 6,800 customers related to a major Swiss supermarket chain in a 20 -day time period. The goods were to be delivered within specified time windows, using three types of vehicles. All the routes had to be performed in one day. [24]

The performance of ANTROUTE was compared to that of human route planners. As the human planners were using a "regional" planning strategy - leading to petal shaped tours - ANTROUTE was run in two configurations: (i) with regional planning mode and with time window constraints enabled (AR-RegTW); (ii) with regional mode and time window constraints relaxed (AR-Free). Every day, ANTROUTE was run on the available set of orders, and it took approximately five minutes to come up with a solution; in comparison, the human planners needed no less than three hours at the minimum. Both configurations of ANTROUTE were able to outperform human planners (see Table 5.1). The shorter the time windows were, the more tours (and hence vehicles) were needed. [24]

|  | Human | AR-RegTW | AR-Free |
| :--- | ---: | ---: | ---: |
| Total number of tours | 2056 | 1807 | 1614 |
| Total km | 147271 | 143983 | 126258 |
| Avg truck loading | $76.91 \%$ | $87.35 \%$ | $97.81 \%$ |

Table 5.1: Comparison between ANTROUTE-generated and man-made tours [24]

### 5.4 Time Dependent MACS-VRPTW

In what follows I will present how a logistical problem was modelled and solved as Time Dependent VRPTW (TDVRPTW) for the city of Padua in Northern Italy.

In TDVRPTW, travel times are time dependent; i.e., the travel time required to traverse a given arc depends on the time of the day [2]. This helps model real-life road networks, where variable traffic conditions are often present. For instance, in many urban environments arterial roads are congested during the rush hours, but free flowing during the other times of day. Taking these conditions into consideration is essential when planning the delivery time windows.

Time dependent MACS-VRPTW is a variant of MACS-VRPTW geared towards solving TDVRPTW [2]. Every arc is associated with information about travel time, so that the travelling time at time $t$ can be deduced. Donati el al. chose to provide this information in the form of a step-like speed distribution $\nu_{i j}(t)$, defined on a time interval $\left[t_{0}, t_{c}\right]$, which partitions the time into periods of time $S_{k}$ defined by intervals $\left[t_{0}^{s}, t_{k}^{s}\right]$. Within an interval, the speed is constant. Hence a single interval can be solved as an instance of classic VRP. The pheromone model associates to each arc $a_{i j}$ a time dependent distribution $\tau_{i j k}$, where the index $k$ refers to the subspace $S_{k}$. Stated differently, the element $\tau_{i j k}$ encodes the convenience of going from $i$ to $j$, departing from $i$ at the time $S_{k}$. The probability distribution used by the ants during the tour construction procedure to select the next customer $j$ out of the feasible set $J$, depends on the departure time from $i$, and is as follows:

$$
p_{j}=\tau_{i j k} \cdot h_{i j}\left(t_{d}\right), \quad j \in J,
$$

where $t_{d} \in T_{k}$ is the departure time from $i$ and $k$ is the corresponding time index,
and $h_{i j}\left(t_{d}\right)$ is the local heuristic function ${ }^{6}$. In the construction procedure, all occurrences of the concept of distance are replaced by the concept of travel time. Hence, the goal is to optimize, not the total length of the solution, but its total travel time. [2]

Dependency on time complicates the local search procedure. The two fundamental operations performed in a local search procedure - namely, insertion and removal of a delivery in a tour - generate a time shift for all the subsequent customers, causing a change in travel and delivery times. Therefore, the local search procedure needs to be modified ${ }^{7}$. [2]

Donati et al. [2] compared the solution to the VRPTW using the time dependent variant with a solution to the same problem in which the travel times on arcs were constant, and depended only on the distances. They ran nine tests, and they reported that the time dependent variant performed $7 \%$ better than the standard VRPTW algorithm [24].

[^18]
## 6 Concluding Remarks

One issue I did not cover in this thesis is how to parallelize ACO algorithms. By their very nature, they lend themselves to be parallelized in the data or population domains [7]. I presume this will be a "hot" research topic in the coming years because of what is happening to computer hardware and processor architectures.

Today, parallel hardware has become ubiquitous [21]. From 1986 to approximately 2002, the performance of microprocessors increased annually by $50 \%$ on average; from 2002 onwards, the single-processor performance improvement slowed down to $20 \%$ per year on average. From 2005 onwards, the major microprocessor manufacturers have sought the performance improvements of microprocessors, not by increasing the speed, but by increasing the number of cores on a microprocessor. [21] Since the first decade of the 21st century, multicore processors - processors with multiple cores - have become the mainstream. For instance, in 2012 quod-core processors - i.e., processors with four cores - are the norm in normal desktop computers, whereas the high-end computers have 10- or 12-core processors [23]. In fact, Moore's law can be stated in a novel way: the number of cores per processor chip will double every 18-24 months [23, 29]. For programmers, these developments and trends mean that in order to fully reap the benefits of modern hardware, programs and algorithms must be made more concurrent and/or more parallel.

The increasing computing power has guided the development of VRP. Research on VRP accelerated during the 1990s, primarily due to, and in parallel with, the rapid growth in processor speeds [9]. Thanks to more powerful computers, researchers were able to model and solve more and more complicated variants and bigger and bigger instances of VRP more easily [9]. Furthermore, enhancements in vehicle tracking, data storage, and exchange media gave rise to dynamic and rich VRPs. As a matter of fact, trends in the VRP literature have shifted from static to more dynamic and fuzzy cases which incorporate real-time data. [9]

It remains to be seen whether the current trends and developments in the processor architecture have a similar boosting impact on bringing parallelism in ACO algorithms to the centre of attention. ACO researchers have studied different parallelization strategies for a long time now, most of which can be classified into fine-grained and coarse-grained strategies [7, 8]. Characteristically, fine-grained parallelization assigns very few individuals to one single processor and emphasizes
frequent information exchange between processors; in contrast, coarse-grained parallelization assigns large subpopulations or entire populations to single processors and de-emphasizes information exchange. Experimental research has concluded that the downside of the fine-grained approach is communication overhead, as the ants end up spending most of their time communicating their modifications to other ants. Indeed, several studies have shown coarse-grained parallelization schemes to be much more promising approach for ACO. [7, 8]

When applied to ACO, coarse-grained schemes run $p$ subcolonies in parallel, where $p$ is the number of available processors [8]. Current development trends in microprocessor architecture allow the value of $p$ to double every year or so. Therefore, it is interesting to see whether the future members of the ACO algorithm family involve dozens or hundreds of ant colonies working in parallel.

Besides processor speeds, other hardware improvements are likely to shape ACO algorithms, too. Similarly to what happened with VRP, the advances in, say, data storage and communication technologies, may shift the focus of ACO algorithms to dynamic and real-time problems, and to problems with multiple objectives.

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[^0]:    ${ }^{1}$ http://www.aco-metaheuristic.org/
    ${ }^{2}$ http://www.aco-metaheuristic.org/mailing-list.html
    ${ }^{3} \mathrm{http}: / /$ iridia.ulb.ac.be/ants/

[^1]:    ${ }^{1}$ A synonym for "vertex" is a "node".
    ${ }^{2}$ Often, a directed edge is called an "arc".
    ${ }^{3}$ As far as the definitions to follow are concerned, it makes no difference whether $e$ is directed or undirected.
    ${ }^{4}$ In general, loops are not allowed in the graphs modelling the VRP.

[^2]:    ${ }^{5}$ The reason why polynomials are used as a divider, is due to their "algebraic" properties; i.e., they are closed under addition, multiplication, and functional composition. These closure properties guarantee the closure of the class of efficient algorithms under natural composition of algorithms. [14]
    ${ }^{6}$ There are thousands of $\mathcal{N} \mathcal{P}$-complete problems. [14, 22]

[^3]:    ${ }^{3}$ The parameter $t$ indicates that the objective function $f$ can be time-dependent. However, in this thesis, I will leave it out for the sake of brevity and assume it is implicitly present.
    ${ }^{4}$ Here we are dealing with a minimization problem. However, it is trivial to turn a maximization problem into a minimization one. Assuming $g$ is the objective function to be maximized, we simply set $f=-g$ to get a minimization problem.

[^4]:    ${ }^{5}$ Note that a pheromone value is generally a function of the algorithm's iteration $t$ :

    $$
    \tau_{i j}=\tau_{i j}(t)
    $$

[^5]:    ${ }^{6}$ Dorigo et al. [3] introduced the AS by applying it to solving instances of TSP. Hence, they used the less general term "tour" to denote a "solution." Also, the choices they made in modeling the problem had an effect on their definitions. They built the construction graph $G$ by associating the locations (i.e., the cities) with the vertices of $G$, and the moves between the locations with the solution components. The pheromone values were associated with the edges $E(G)$. There are other, though perhaps less intuitive, ways to model the problem [6], but as a consequence the formulas, or their explanations, might look different. This subtle observation needs to be kept in mind while reading the formulas to follow, as I will make no attempt to present them in a generalized fashion.

[^6]:    ${ }^{7}$ If the component $j$ is already a part of the (partial) solution built by the ant $k$, then $j \notin N_{k}(i)$.

[^7]:    ${ }^{8}$ In this respect, the ACS is similar to the ant-density and ant-quantity algorithms.

[^8]:    ${ }^{9}$ In this extended version of the ACS, an ant first chooses the city to move to from the candidate list; if none of those cities can be visited, only then will it consider the rest of the cities.

[^9]:    ${ }^{10}$ Not all optimization problems exhibit this behavior. For instance, quadratic assignment problem (QAP) - the problem of assigning a set of facilities to a set of locations with given distances between the locations and given flows between facilities - is one such problem. In the QAP, local minima tend to spread over large $(\mathrm{r})$ parts of the search space than in the TSP. This is probable one of the reasons why the QAP is regarded as one of the hardest optimization problems [26].

[^10]:    ${ }^{11}$ The PTS mechanism has the potential to increase the performance of any elitist version of the AS, such as the the rank-based Ant System $\left(\mathrm{AS}_{\text {rank }}\right)$.

[^11]:    ${ }^{12}$ In Table 4.1, "eil51", "kroA100", and "d198" are symmetric instances of the TSP, whereas "ry48p", "ft70", "kro124p", and "ftv170" are asymmetric instances. For each instance, the average solution quality has been reported.
    ${ }^{13}$ As already noted, the PTS can be added to any elitist version of the AS.

[^12]:    ${ }^{14}$ It needs to be pointed out that the $\mathrm{AS}_{\text {rank }}$ never found the optimal solution for the "ry48p" instance, whereas the MMAS and ACS did.

[^13]:    ${ }^{15}$ A convergence proof means proving that during the execution of an ACO algorithm, the ants will eventually produce the optimal solution, and nothing but the optimal solution.
    ${ }^{16}$ I will omit the proofs of the theorems and propositions; instead, I will refer the reader to [25].

[^14]:    ${ }^{17}$ Note that the length of the transition period $t_{0}$ is bounded.
    ${ }^{18}$ See Proposition 5.

[^15]:    ${ }^{1}$ Essentially, a job scheduling problem is about scheduling $n$ jobs on $m$ identical machines while trying to minimize the total length of the schedule, and subject to possible some constraints.

[^16]:    ${ }^{2}$ This is called elitist strategy.
    ${ }^{3}$ For pseudocode, see Figure 2 in [11].

[^17]:    ${ }^{4}$ For pseudocode, see Figure 3 in [11].
    ${ }^{5}$ For pseudocode, see Figure 4 in [11].

[^18]:    ${ }^{6}$ For the exact formula, see Formula (8) in [2].
    ${ }^{7}$ For more details, see Section 6 in [2].

