



UNIVERSIDAD NACIONAL DE COLOMBIA

Kalman-based schemes for mobile nodes localization in ad-hoc networks

Esquemas de localización de nodos móviles en redes *ad-hoc* basados en filtros de Kalman

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Dedicated to

To my beloved ones, my Mother Aracelly and my Brother Alejo, because they are the strength that makes me get up every day and their support is always there for me...

To my Granny Consuelo, who taught me to have faith and to know that good things happens when you believe...and by extension to all the Murillo's Family...

And to Marley (just be patient, Honey!)

Dedicatoria

A mis seres más queridos, mi Madre Aracelly y mi hermano Alejo, porque ellos son la fuerza que me hace levantarme cada día y su apoyo siempre está disponible para mí...

A mi mamita Consuelo, quien me enseñó a tener fe y a saber que cosas buenas pasan cuando creemos...y por extensión a toda la familia Murillo...

Y a Marley (teneme un poquito de paciencia!)

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Resumen

Esta tesis enfrenta el problema de la determinación de la posición de nodos móviles en redes inalámbricas *ad hoc*, con base en las mediciones del indicador de potencia de la señal recibida (RSSI). Las características de movilidad de los nodos se modelan a través de un sistema no lineal representado por un modelo de Giro Coordinado (CTM). La localización de la posición de los nodos se lleva a cabo mediante multilateración integrada con diferentes esquemas para refinar la estimación basados en las técnicas de Kalman: Filtro Extendido de Kalman (EKF); Filtro de Kalman "Unscented" (UKF); y dos Filtros de Múltiples Modelos Interactuantes (IMM), los cuales consisten en un conjunto de Filtros de Kalman Extendidos (IMM-EKF) y un banco de Filtros de Kalman "Unscented" (IMM-UKF). Se alcanza a estimación de los estados del modelo de movilidad, los cuáles comprenden la posición, la velocidad y, en algunos casos, el parámetro de Tasa de Giro del nodo objetivo móvil. El desempeño de los diferentes esquemas basados en Kalman se compara mediante el seguimiento de dos trayectorias a través de simulaciones Monte Carlo.

Palabras clave: Redes *ad hoc*, Localización de nodos, Seguimiento de trayectorias, Estimación, Filtros de Kalman, Filtro de Kalman Extendido, Filtro de Kalman "Unscented", Múltiples Modelos Interactuantes.

Abstract

This thesis addresses the problem of position localization of mobile nodes in ad hoc wireless networks based on received signal strength indicator (RSSI) measurements. Node mobility is modelled as a non-linear system driven a Coordinated Turn Model (CTM). Self-localization of mobile nodes is performed via multilateration integrated with a different collection of Kalman based schemes for estimation refinement: Extended Kalman Filter (EKF); Unscented Kalman Filter (UKF); an two Interacting Multiple Model Filter consisting of a bank of Extended Kalman Filters (IMM-EKF) and Unscented Kalman Filters (IMM-UKF). Estimation of the mobility state, which comprises the position, speed and, in some cases, the Turn Rate parameter of the mobile node is accomplished. The performance of the Kalman based filters is compared through the tracking of two different trajectories by Monte Carlo simulation.

Keywords: *Ad hoc* networks, Node Localization, Trajectory Tracking, Kalman Filter, Extended Kalman Filter, Unscented Kalman Filter, Interacting Multiple Model

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List of Abbreviations

Abbreviation	Term
<i>AhLoS</i>	Ad Hoc Localization System
<i>ANOVA</i>	Analysis of Variance
<i>AoA</i>	Angle of Arrival
<i>APS</i>	Ad Hoc Positioning System
<i>BER</i>	Bit Error Rate
<i>CWPA</i>	Continuous Wiener Process Acceleration (<i>PVA</i>)
<i>CWPV</i>	Continuous Wiener Process Velocity (<i>PV</i>)
<i>DE</i>	Distance Error
<i>EKF</i>	Extended Kalman Filter
<i>GPS</i>	Global Positioning System
<i>IMM</i>	Interacting Multiple Model
<i>KF</i>	Kalman Filter
<i>MANET</i>	Mobile Ad Hoc Network
<i>MDS – MAP</i>	Multi-Dimensional Scaling
<i>MinMax</i>	Minimum-Maximum Algorithm
<i>MSE</i>	Mean Squared Error
<i>P</i>	Position Model
<i>RMSE</i>	Root Mean Squared Error
<i>RSS</i>	Received Signal Strength
<i>RSSI</i>	Received Signal Strength Indicator
<i>ToA</i>	Time of Arrival
<i>TDoA</i>	Time Difference of Arrival
<i>UKF</i>	Unscented Kalman Filter
<i>WMN</i>	Wireless Mesh Networks
<i>WSN</i>	Wireless Sensor Networks

1 Introduction

1.1. Mobile nodes self-localization in ad hoc networks

The movement patterns of mobile users play an important role in performance analysis of wireless computer and communication networks in which nodes may move freely within an area. The structure of *ad hoc* wireless networks can change dynamically over time; this fact complicates the network control and management tasks. It is very important for this kind of network to localize the node positions and movement [17] [64] [45] as the transmitter range is generally fairly small with respect to the size of the area. Self localization [35] involves the combination of absolute location information (e.g., obtained from a Global Positioning System (GPS)) with relative distance information (e.g. distance measurements between sensors) over regions of the network. It is also desirable to minimize the amount of inter-sensor communications.

There are many methods for self-localization, one class is based on signal measurements and their statistical models (see, e.g., the surveys [17] [45] [35]). These methods rely upon the signal time-of-arrivals, time difference of arrivals, angle-of-arrivals or received signal strengths and they vary in their complexity and accuracy. In this thesis we consider the self-localization of mobile nodes in wireless *ad hoc* networks using received signal strength indicator (RSSI) measurements. Node mobility is modelled with a linear dynamic model, with multiple acceleration modes, which are driven by a discrete Markov process. Due to the fact that the control process of the mobile node is unknown and we have multiple acceleration levels, the Interacting Multiple Model approach [5] is suitable for the considered problem. We implemented it in combination with an Extended Kalman Filter (requires linearization of the non-linear measurement equations) and an Unscented Kalman Filter (affords avoiding linearization of the highly non-linear measurement equations). The IMM Kalman-based filters are compared with EKF and UKF algorithms for mobile nodes self-localization.

Previous approaches to mobile nodes localization depend on the type of the ad hoc network [35]: indoor or outdoor. In the indoor sensor network the localization can be performed with beacons (fixed or moving) or it can be beacon free. In outdoor applications GPS systems are mainly used. Many localisation techniques rely on Kalman Filtering [36] [59] [37] and Monte Carlo techniques [20], including and knowledge of the connectivity between the nodes. Some works [45] consider the case when nodes can communicate among each other which is not

always possible because communications are energy-consuming.

1.2. Motivation

The absence of a fixed infrastructure in wireless *ad hoc* networks make them suitable for use in emergency situations as well as for low cost commercial communication systems. However, the flexibility of the highly dynamic *ad hoc* networks complicates important control and management tasks such as routing, flow control, and power management. For example, traffic routes change over time, subject to the movement of the mobile nodes. The effectiveness of any routing algorithm depends heavily on the accuracy and timelines of the available network topology information. In *ad hoc* networks, knowledge of the network topology can be inferred from the mobility of the nodes. The absence of a fixed infrastructure in wireless *ad hoc* networks make them suitable for use in several applications of target tracking as well as for low cost commercial communication systems.

Although different algorithms for self-localization and tracking have been proposed in the literature, this is an open and active research area, where many questions from theoretical and practical point of view remain unsolved. The difficulties are coming from the changeable network topology, the need of communications between the nodes under limited resources (energy, bandwidth), noisy data and the challenge of overcoming losses.

1.3. Objectives

General Objective

The aim of this thesis is to present a performance comparison of a collection of Kalman based tools capable of target tracking in *ad-hoc* networks by solving one or both target motion and measurement uncertainties.

Specific Objectives

- The development of tracking algorithms for *ad-hoc* networks based on Kalman filters and IMM techniques using distance measurements and lateration.
- The implementation of a simulation scheme for the evaluation of the developed localization algorithms.
- The selection of a performance metric for the comparison of the simulation results of the studied methods.

1.4. Chapter by chapter overview

This thesis is organized in 5 chapters. A brief description of each chapter is given as follows.

- **Chapter 2:** This chapter presents an introduction of the Ad Hoc Wireless Sensor Networks, with a brief definition and an illustration of its properties, applications and the current design challenges.
- **Chapter 3:** This chapter presents an overview of the problem of self-localization of nodes in *ad hoc* networks, and it shows the stages of the localization process. Also, it describes the state of the art of the localization and tracking solutions for *ad hoc* networks found in literature.
- **Chapter 4:** This chapter describes the mobility models used to represent the dynamics of the mobile target, and it also shows the measurement model for extracting position information from the RSSI data over the *ad hoc* network.
- **Chapter 5:** This chapter presents a detailed description of the theoretical tools used to develop the localization algorithms, such as lateration and Kalman Filter and its derivatives (EKF - UKF). We also introduce the Interacting Multiple Models tools for the estimation of the target trajectory. Later, some algorithms for localization and tracking using Kalman Filter in *ad hoc* networks are presented.
- **Chapter 6:** In this chapter, the simulation results of the studied methods are presented.
- **Conclusions:** This dissertation finishes with some concluding remarks and outlines possible future work areas on this matter.

2 Ad Hoc Wireless Sensor Networks

There is a major difference between wireless and wired networks: the former makes the transmission of the information through air and the latter through physical cables. Therefore, the main difference between wired and wireless is the infrastructure. At the wired networks: wires, switches, hubs and all the other communication means are responsible for transporting the information while at wireless networks signals and waves propagate through the air. Figure 2-1 presents a wireless networks classification according to [30]. As it is the subject of this thesis, the following sections introduce the characteristics of the Ad Hoc Wireless Sensor Networks.

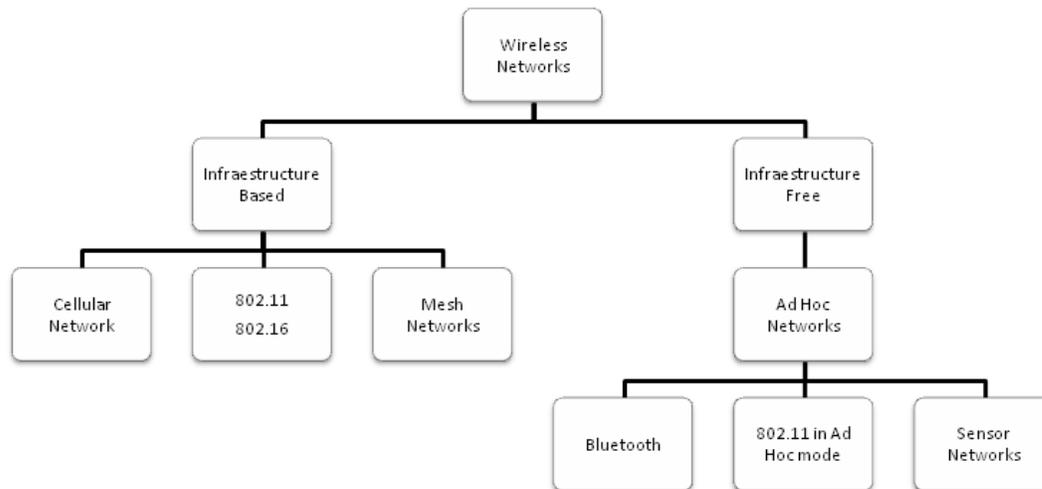


Figure 2-1: Wireless Networks Classification

2.1. What is a Wireless Ad hoc Network?

A wireless ad hoc network is a decentralized wireless network with a dynamical structure changing over time, which complicates the network control and management tasks. Data can be forwarded from any node to any other node through zero, one or more intermediate nodes. Other characteristics are:

- Multi-hop routing is a common feature in ad hoc networks: data from a source to destination may go through certain intermediate nodes.

- Each node can send data, receive data and also act as a router forwarding data for peers.

Wireless cellular networks need a fixed infrastructure where an access point manages communication among nodes. In contrast, a wireless ad hoc network requires minimal prior configuration and can be quickly deployed: the devices self-organize among themselves to form a network on the fly. The emphasis here is on the *ad hoc* character of these kinds of networks; it implies that network deployment and configuration happens with no need of previous preparation, and assumes no existing infrastructure in the area of deployment. The dynamic aspect of these kind of networks constitutes its most interesting and more challenging characteristic.

2.1.1. Types of Wireless Ad hoc Networks

Ad hoc Network (MANET)

A MANET is a self-organizing wireless network of mobile devices. The topology of the network (arrangement of the devices) changes dynamically with time. An example of this kind of networks are the Vehicular Ad hoc Networks(VANET) [63]. A type of MANET that is fast gaining wide research focus, VANETs are used for communication among vehicles (equipped with wireless devices) and between vehicles and roadside equipment.

Wireless Mesh Networks (WMN)

A WMN is an ad hoc network where all nodes are generally static and mobility is not an issue. A WMN is composed of mesh clients (laptops, cell phones and etc.) and mesh routers that forward traffic towards and from the mesh clients to the gateways that connect to the Internet. The mesh topology provides multiple paths from the source to the destination and helps in quick reconfiguration of paths when an existing path fails. The WMN nodes are normally not resource constrained. An example of a WMN could be a mesh formed by placing wireless relaying equipment on the top of houses [2].

Wireless Sensor Networks (WSN)

Being the main subject of this thesis, this kind of ad hoc networks will be reviewed in the next section with more detail.

2.2. What is an Ad Hoc Wireless Sensor Network?

Generally speaking, an ad hoc wireless sensor network consists of a large number of self-sufficient nodes with sensing capabilities. The nodes can perform simple computations and

communicate with each other through a self-deployed and self-healing multihop communication network [1]. The sensors are distributed over a region to cooperatively monitor physical or environmental conditions (such as temperature, pressure, pollutant concentration, forest fire, etc.) at different locations and report the data to a central base station (commonly called sink).

Each node in a sensor network is typically equipped with one or more sensors; a radio transceiver for communication; a micro-controller for computation and decision making and a battery. The lifetime of each individual sensor may not as significant but lifetime of the entire network may be an important performance metric.

2.2.1. Applications of Ad Hoc Wireless Sensor Networks

The claim of wireless sensor network proponents is that this technological vision will facilitate many existing application areas and bring into existence entirely new ones. This claim depends on many factors, but a couple of the envisioned application scenarios shall be highlighted. Apart from the need to build cheap, simple to program and network, potentially long-lasting sensor nodes; a crucial and primary ingredient for developing actual applications is the actual sensing and actuating capabilities of a sensor node. For many physical parameters, appropriate sensor technology exists that can be integrated in a node of a WSN. Some of the few popular ones are temperature, humidity, visual and infra-red light (from simple luminance to cameras), acoustic, vibration (e.g. for detecting seismic disturbances), pressure, chemical sensors (for gases of different types or to judge soil composition), mechanical stress, magnetic sensors (to detect passing vehicles), potentially even radar [13]. But even more sophisticated sensing capabilities are conceivable.

The rest of this section presents a number of applications of ad hoc wireless sensor networks found in the literature.

Environmental applications

Although there are some other techniques to monitor environmental conditions, random distribution and self organization of WSNs make them suitable for environmental monitoring. WSNs can be used to control the environment, for example, with respect to chemical pollutants: a possible application are garbage dump sites. Another example is the surveillance of the marine ground floor; an understanding of its erosion processes is important for the construction of offshore wind farms. Closely related to environmental control is the use of WSNs to gain an understanding of the number of plant and animal species that live in a given habitat (biodiversity mapping). The main advantages of WSNs here are the long-term, unattended, wire-free operation of sensors close to the objects that have to be observed; since sensors can be made small enough to be unobtrusive, they only negligibly disturb the

observed animals and plants. Often, a large number of sensors is required with rather high requirements regarding lifetime. Some of the specific ongoing study cases over this field are:

- Planetary exploration [3]
- Geophysical monitoring [61]
- Habitat monitoring [33]
- Bio-complexity mapping [24]
- Oceanography [7]

Transportation Systems

Cars equipped with sensors can form local communication networks sharing information on weather and road conditions, plan efficiently chosen routes, avoid traffic and identify their position in areas where GPS signals are unavailable [9].

Structural Monitoring

Sensors are placed on structures to detect structural weaknesses, the presence of hazardous materials and, their reaction to extreme weather and physical phenomena [29]. These sensors can form local communication networks and forward the information they gather to central points where suitable simulation models can then decide on the need for structural changes and help prevent accidents.

Medical and Physiological Monitoring

Wireless sensors can be used in hospitals to efficiently locate doctors and patients and when attached to medications, they can minimize the danger of issuing the wrong medication to patients [28]. Small sensors, each with different sensing capabilities, can be attached to different parts of the body of a patient to measure physiological signatures, detect infections and automatically regulate her medication or alert doctors [57].

Intelligent buildings

Buildings waste vast amounts of energy by inefficient usage of Humidity, Ventilation and Air Conditioning (HVAC). WSN can be integrated with building automation systems [44], in such a way that employees are equipped with sensors that form local networks and efficiently locally regulate light, temperature and humidity, based on predefined individual preferences. Every time an employee enters or leaves a room, the network is updated (restructured) and the light, temperature and humidity settings of the room are updated accordingly.

Military and Security

As it also occurs into other fields of technology research, the military plays a central role in the advancement of ad hoc wireless sensor networks by initiating and supporting a variety of related projects. The rapid deployment, self-organization and fault tolerance are characteristics of sensor networks that make them a very appealing promising technique for military applications. These area also includes alerting systems and disaster relief implementations. Some examples of use are:

- Surveillance and battle-space monitoring [58] [23]
- Urban warfare [23]
- Self-healing minefields [6]
- Disaster detection and rescuing operations [1]
- Commercial and residential security [22]

Inventory Control

Even in almost fully automated distribution centers, due to their dynamic nature, inventory management still requires high-cost manual inventory control operations. When items in the warehouse are equipped with low-cost tiny sensors, these operations become obsolete; stock can be easily located in the warehouse and inventory management can become a fully automated process [1].

Robot Swarms

Most of the issues that concern the design and implementation of ad hoc wireless sensor networks are also relevant when multirobot formations are considered. Robots are equipped with sensors; they communicate with each other and form local networks to accomplish various tasks. Note that the problem of network localization that we are here concerned with, plays an important role in multirobot formations and is still an active field of research [32].

Social Networks Analysis

There are fields that cannot be considered as direct applications of ad hoc wireless sensor networks, but still share much of their structure and characteristics, and can benefit from their development. One can name, for example, the study of Social Networks, where the importance of concepts as centrality, network density and visualization (and relative positioning) of such networks plays an important role. Actors in such networks have sensing capabilities, communicate with other actors and interact with their environment in such a way that can be easily modeled as ad hoc wireless sensor networks [19].

2.2.2. Properties of Ad Hoc Wireless Sensor Networks

The increasing number and variety of applications and designs of wireless sensor networks make it difficult to speak of a typical wireless sensor network in a strict sense [49]. Roughly speaking, there is a different kind of network depending on the application purpose. However, it is assumed that ad hoc wireless sensor networks possess certain characteristics that permit them to make the task at hand.

- ***Ad Hoc***: The network can deploy itself without depending on the existence of any external infrastructure, without the requirement of previous planing, and has the ability to self-assemble and self-organize and perform its assigned tasks unattended [1].
- ***Self-sufficiency***: The network is able to manage its own resources: energy, bandwidth and processing power in an optimal way with the goal of maximizing its lifetime.
- ***Communication Network***: The sensors are able to communicate with each other or with an external (central) computational unit, and form a “communication network”. A communication network can involve all the nodes in the network uniformly (each node is connected and communicates with all its neighbors), or can use local *clustering* (each node is associated with a cluster head that they then can communicate with each other).
- ***Sensing Capabilities*** The nodes are equipped with sensing devices and can be specialized to detect certain environmental effects such as motion, heat or sound [1]. By using algorithms that feature emergent behavior, the network can demonstrate sensing capabilities without the cost of having all nodes being able to sense all effects that are of interest.
- ***Robustness***: The network should be in the position to reorganize itself when additional nodes become available or in the case of node failure. When mobile nodes are considered, the network should be in the position to adapt to structural changes due to node movement.
- ***Small size/low cost***: When ad hoc wireless sensor networks are considered, we refer to networks of homogeneous nodes that are small in physical size and of low economic cost; nodes have to be deployed in large numbers and, if needed, left behind.
- ***Location awareness***: Nodes need to be aware of their position, either in an absolute way or in reference to an internal coordinate system [35].

2.2.3. Challenges in the Design of Ad Hoc Wireless Sensor Networks

There are many challenges that designers of hardware and software for wireless sensor networks have to face before they can make wireless sensor networks attractive for large scale

commercial use. Most of these challenges have to do with the large number of nodes that are used and the unpredictable behavior that accompanies such large scale systems [8].

The hardware challenges at hand include extending the lifetime of the networks (energy), building efficient ways for the networks to adapt easily to varying environmental constraints, and building adaptive data collection mechanisms that are fast and robust [15]. On the software side, information flow and localization seem to be the most challenging aspects of the design of wireless sensor networks. Energy constraints play an important role here.

Information Flow

Information should be able to travel through the network. The most popular way of information broadcasting in (large scale) ad-hoc networks is Flooding. With Flooding [56], a message is broadcasted by all receiving nodes to their neighbors until the entire network is reached. In a clustered network, flooding is performed between the cluster heads which then forward the message to all nodes that belong to their cluster. An empirical study [56] has shown that communication in large networks can be hugely affected by small environmental effects and deviations from the norm. These may result in "dark areas" in the network; areas that the broadcasted information is unable to reach.

The network should in some cases be able to communicate with a "base station" or "sink node" outside of its deployment area [18]. When a number of the available nodes are able to transmit and receive messages to and from the base station, flooding can be used to communicate information between the base station and the nodes. When a message from the base station needs to be communicated to some, or all, of the nodes in the network, it is enough if it can reach some of its nodes, since it can then be forwarded to all the nodes using their internal communication network. In the reverse case, nodes that need to communicate information to the base station can simply forward it through the existing communication network to the nodes that are in the position to reach the base station.

In some scenarios the transmitters and receivers of the nodes have different angular spreads, resulting in an asymmetrical communication network; some nodes might receive messages from neighbors that they will not be able to transmit to. Information flow can also be affected by structural changes in the communication network such as node failures, and message collisions. Noisy environments can sometimes be destructive to the information flow in the network [18] as they can severely affect the quality of the transmission.

Recent research has led to the creation of a number of alternative schemes, such as Gossiping, SPIN, SAR, LEACH, SMECN and Directed diffusion (see [1] for an extensive study of their properties and behavior), to address the inefficiencies of simple Flooding communication

protocols (i.e., duplicate messages sent to the same node, multiple nodes sharing the same observing region, resource blindness, etc.).

Energy

It is assumed that ad hoc wireless sensor networks should be low-cost and unattended. It is therefore expected that node failure will be a common event [18]. Accordingly, communication and computational demands should be minimized to prevent node-failure due to energy overconsumption.

Localization

The problem of localization, that we address in the present work, concerns the estimation of the position of wireless sensor nodes in an ad hoc network setting, in reference to a coordinate system that may be internal or external to the network. Localization is one of the main challenges of an ad hoc wireless sensor network set up and there are many technical issues that make this problem a very difficult. Some of these issues are:

- **Non-Line of Sight Problem (NLOS):** The existence of *physical obstacles* can prevent neighboring nodes from detecting each other [48]. This in turn, can lead to losses of important connectivity information.
- **Sparse Node Problem:** The connectivity information available to the network, especially in networks of low density, may not be enough for the construction of a unique solution [48].
- **Geometric Dilution of Precision (GDOP):** In practice, the distance measurements used to compute node coordinates almost always have some error. These measurement errors get reflected in the computed node coordinates. The magnitude of the final computed error depends on both the value of the measurement error and the true geometry of the structure induced by the nodes and edges. The contribution due to geometry is called the *Geometric Dillution of Precission* (GDOP), and is defined as the ratio between the computed coordinate error and the measurement error [48].
- **Range Error Problem:** Localization is based either on distance or connectivity measurements, and this information is prone to error. Distance information may contain errors as large as 50 % of the measurement value [53].

Summary

This section presented the concept of Ad Hoc Wireless Sensor Networks, a type of Ad Hoc Network composed of many self-sufficient nodes with sensing capabilities. The applications,

properties and challenges in the field of Ad Hoc Wireless Sensor Networks were also presented. Between the challenges, the localization problem is highlighted because it is the central topic of this thesis.

3 Self Localization in Mobile Ad Hoc Networks

The tendency of networks, both wired and wireless, is to continuously grow in size and complexity. There is a need to make automatic process of the self-configuration and monitoring tasks, and this need increases along with the network growth. Therefore, the problem of making the nodes localize themselves an constitutes an issue of primary concern in a WSN (and in general in a MANet). Information about the position and orientation of the individual nodes is useful to the target application and service. For instance, routing and querying are services that can be controlled according to position. At the application level, we require position information in order to log the reported data in a sensor network. Consider a meteorological application where it is necessary to make geographical sense of the observed data e.g. labelling a map with temperature or humidity.

3.1. Localization

When sensors are deployed in an ad-hoc manner, a priori knowledge of their location is not possible [13]. There are, therefore, many reasons why solving the localization problem is crucial, and while GPS could be employed in some outdoor settings and only for some types of networks, it is not suitable for wireless sensor networks because of the following reasons [27], [41]:

- GPS cannot be used indoors, where satellite signals are not available.
- Most applications need location information of high accuracy that cannot be offered by GPS.
- Battery constraints in most wireless sensor networks settings make the use of GPS highly inefficient.
- When networks of thousands of disposable nodes are considered, the installation of GPS devices would be forbiddingly expensive.
- The size of the nodes under consideration is, in most cases, much smaller than what would be needed if GPS devices were to be included in the installation.

The problem of localization is an issue that should be addressed, since its importance for ad-hoc wireless sensor networks cannot be neglected. There is a number of reasons why location information is useful and in some cases why it is absolutely necessary:

- **Data Registration:** Sensed data, in many cases, are almost useless when they are not located in space (and sometimes also in time).
- **Efficient targeting:** When sensors are aware of their location, they can either trigger the partial silencing, or the activation of some parts of the network when activities that have to be measured are not present or detected, respectively. This way, sensors can save energy when they are not needed, and the communication becomes more efficient, since the transmitting or receiving of redundant messages is avoided.
- **Target tracking:** When the purpose of the network is to track (possibly moving) targets in its deployment area, node localization is absolutely necessary [28], especially when the network must be able to restructure itself (as in the WSN for Minefield Detection [6]), or to adopt to node failures, target movements or security breaches.
- **Coverage:** When mobile nodes are considered, the network can use different schemes to maximize its coverage of the deployment area, while ensuring the robustness of its communication network. In such a setting, it is assumed that nodes are aware of their position in the deployment area [21].
- **Routing Protocols:** Communication protocols, such as the Location-Aided Routing (LAR) protocol [60] use location information for efficient route discovery purposes. The location information is used to limit the search space during the route discovery process so that fewer route discovery messages will be necessary. In mobile ad hoc networks, such as MANET [30], location information is used to design mobile gateways to increase the information flow inside the network and to create efficient routes to communicate information to stations outside the deployment area. It should though be noted that recent research in Swarm Intelligence [32] has produced routing protocols, such as the Ant-based control (ABC), or the AntNet” method, that can be applied to ad hoc networks and do not depend on location information.

3.2. Technical Challenges

Localization depends on the topology of a network and the quality of the information that is available. As the number of nodes becomes large (currently hundreds of thousands), the dynamic nature and the complexity of the topology of the network become the first major obstacles to localization techniques. Most of the available localization methods depend heavily on proximity and/or distance information between neighboring nodes. Measurement

errors are the second major obstacle to localization. In the following paragraphs, we discuss these two issues in more detail.

3.2.1. Network Topology

The topology of a network is defined by one or more of the following characteristics: size, capacity, diameter, coverage, density, and connectivity.

- **Size:** The size of a network is defined as the number of nodes that are active in it. As the size increases, the complexity of the localization methods can increase in such a way that some of them may become unusable.
- **Capacity:** The capacity of a network is the average rate of data transmission between any two nodes in the network in bits per second [16]. As the size of the network increases this capacity is negatively affected. It has been shown in [16] that for ad hoc networks with n randomly placed nodes, assuming omnidirectional antennas and disregarding link or node failures, the capacity of the network is $O(1/\sqrt{n})$. So, as the size of the network increases, its capacity tends to zero.
- **Diameter:** The diameter of a network is the largest distance between any pair of nodes of the network. It is sometimes estimated by the largest shortest path between any two nodes or by the maximum number of hops [30] that is needed for any two nodes in the network to reach each other.
- **Coverage:** Any node, depending on the capabilities of its sensors, can cover a part of the deployment area of the network. The area that is covered by the whole network is called its coverage [49]. The coverage of a network is closely related to its density and connectivity, properties that influence the quality of localization. In [49], the authors wrote that the coverage of a network can be of three types: *sparse* (when the network coverage is much smaller than its deployment area), *dense* (when the network coverage coincides with its deployment area, or comes close to it) and *redundant* (when multiple sensors cover the same area).
- **Density:** The term density refers to the number of neighbors per node in the network. Density is closely related to connectivity. A network is considered to be connected if there is a path between any pair of its nodes. The density of the network can affect its communication capabilities, its lifetime and the quality of localization. Networks with high-density can be easier localized than networks with low-density, and even poor distance information can be compensated by high connectivity [27]. However, high-density networks are more energy-costly than low-density networks, and have therefore a shorter life expectancy. Nevertheless, research is being done on prolonging the lifetime of high-density networks by making data aggregation more efficient [27].

- **Connectivity:** The connectivity of a network depends largely on the (transmitting) communication range of its sensors, as well as its density. When the range is shorter than the shortest distance between any pair of nodes, the network is completely unconnected. When the range is longer than the diameter of the network, the network is completely connected but energy consumption becomes inefficient in this case and message collisions can diminish the network's communication capabilities. The problem of estimating the minimum range that guarantees that the network stays connected is known as the Critical Transmitting Range problem [27].

3.2.2. Estimating the Distance Between Two Nodes

Localization is based either on connectivity information, or distance information; on case, each node, estimates its distance to its neighbors in a distributed manner. However this information is not free of error. In ad hoc wireless sensor networks, where power consumption constraints and limits on the size of the sensors are common challenges, the influence [30] of the environment (physical obstacles, noise, radio interference) and the transmission system used, render reliable measurements infeasible. Range estimation is usually performed using one of the approaches described below [35], [47].

Received Signal Strength (RSS)

RSS is defined as the voltage measured by the circuit of received signal strength indicator (*RSSI*) in any receiver. It can be equivalently reported as squared magnitude of the signal strength, i.e., the measured power. RSS of Radio Frequency (RF) signals can be estimated by each receiver during normal data communication within the network. This does not require additional bandwidth, energy or hardware. These features of RSS measurements make it relatively inexpensive and simple to implement, and make this techniques appealing to research institutions.

However, RSS measurements are often very noisy, mostly due to the shadowing and fading phenomena. For a distance, d , between a transmitter and receiver, it is often quoted that signal power decays proportional to d^{-2} , but this is not always the case under experimental validation. In practical situations, multiple signals with different amplitudes and phases arrive at the receiver and signals add constructively or destructively as a function of the frequency, causing frequency-selective fading. In addition, the measured RSS is also a function of the transmitter and receiver calibration; RSSI circuits and transmit power vary according to the hardware, and also the battery state.

Time-of-Arrival (ToA) and Time-Difference-of-Arrival (TDoA)

ToA is the measured time at which a signal (RF, acoustic, laser, etc.) first arrives at the receiver. The measured ToA is the time of transmission plus the signal propagation delay. This time delay is a linear function of the transmitter-receiver separation distance if the propagation velocity is previously known. The most difficult aspect of this time-based technique is to accurately estimate the arrival time of the line-of-sight (LOS) signal. This estimation is impaired both by additive noise and multi-path signals; .early-arriving multi-path. and .attenuated LOS. are the main cause of estimation errors. Typically, the ToA is chosen as the time value maximizing the cross-correlation between a known transmitted signal and the received one (Simple Cross-Correlator, SCC). If the nodes of an wireless *ad hoc* network are accurately synchronized, then the time delay is determined by subtracting the known transmit time from the measured ToA.

For an asynchronous network, a common practice is to use either two-signals or round-trip techniques. In the first a sensor transmits two signals simultaneously with two different propagation velocities e.g. RF and ultrasound. The receiver computes the time difference between the two signals. In the second method a node transmits a signal to a second node, which immediately replies with its own signal. At the originating transmitter, the measured delay between transmission and reply is twice the propagation time plus the known delay of the second sensor. It should be noted for a given bandwidth and signal-noise ratio (SNR), the time delay estimate can only achieve a given accuracy. A more in depth study is provided in [47].

Another alternative is the *Time Difference of Arrival* (TDoA) technique. This is similar to ToA, but based on processing signals transmitted simultaneously by different terminals.

Angle-of-Arrival (AoA)

AoA involves the measurement of the direction towards neighboring nodes rather than the distance to them. This kind of information is complementary to RSS and ToA. The normal way in which nodes measure AoA is to use a sensor array, or alternatively two (or more) directional antennas.

In the sensor array case, each sensor node is comprised of two or more individual sensors (microphones for acoustic signals or antennas for RF signals) whose locations with respect to the node center are known. The AoA is estimated from the differences in arrival times for a transmitted signal at each of the sensor array elements employing techniques called .array signal processing". In case of multi-directional antennas, the beams of antennas pointed in different directions overlap and can be used to estimate the AoA from the ratio of their individual RSS values.

Both of the above methods require hardware that considerably increases the device cost and size. AoA measurements are still hampered by additive noise and multi-path. Since it is not likely that sensors will be placed with known direction, localization reliability may depend on the sensor orientation.

Connectivity and Bit-Error-Rate (BER)

This technique determines whether two devices can communicate or not. Two nodes are not regarded as connected simply based on the distance between them. They are connected if the receiver node can successfully demodulate packets from the transmitter. The connection quality is usually correlated with the received signal strength (RSS) and therefore the bit error rate (BER) of the received packets. BER increases as RSS decreases, and demodulation of packets can easily fail if the RSS is too low. BER and connectivity for ranging is generally applied in conjunction with the other methods previously reported since it is a crude technique.

3.3. A brief literature survey about ad-hoc localization schemes

Localization algorithms can be implemented in one of two distinct ways: if each node in a network is able to localize itself it is referred to as distributed localization; otherwise, if the data collected are relayed to a powerful central point (*sink*) it is called centralized localization. A further approach can be identified, similar to the distributed case but each node can only make use of local data (in other words communication is limited), this is called localized localization. Localization approaches can be further categorized as range based or range free protocols. The first bases the estimates upon absolute distance measures, whilst the latter uses relative position and connectivity information.

Generally speaking, a localization process is made up of three phases [31]:

1. **Ranging:** to determine the distances among the nodes that form the network and the reference points.
2. **Positioning:** to give a first evaluation (which in some cases can be very approximate) about the nodal positions from its distances to reference nodes. Several methods exist to do this and depend on the application, available budget and network typology.
3. **Refinement:** to refine the outcome of the previous two phases using nodes information in addition to the range (distance). This is derived from the network and the nodes which form it.

References [35], [47] report complete overviews about the different techniques used in localization.

3.3.1. Ranging

All the methods used to measure relative distances among nodes can be referred to as "*ranging techniques*". The most commonly encountered were presented in subsection 3.2.2.

3.3.2. Positioning

As previously discussed *Positioning* is a stage which follows after *Ranging*. There are several well-known methods to implement positioning but the most commonly encountered are presented in the following subsections.

The first three positioning techniques below (Lateration, MinMax, RocRSSI+) use RSS as a ranging technique. They are attractive because of the low cost and complexity. These methods work on the assumption that anchor nodes or landmark devices with known positions are available and can act as reference points for the nodes with unknown position. Nevertheless, in [38] the authors investigate algorithms for localization in sensor networks where there is a lack of absolute reference. References [35], [34], [36] and [41] list several other possible ways to achieve the positioning stage. In [35], particularly, there is a classification of positioning algorithms and the authors discuss "*one hop*" or "*multi-hop*" solutions.

One hop algorithms assume nodes receiving positioning information are within the connectivity range (*hop*) of the device (landmark, satellite, anchor or beacon) acting as the reference point and providing the localization service.

"*Multihop*" solutions are those where localization algorithms have to consider nodes which may not be in direct contact with reference nodes. They do not necessarily measure ranges to anchor nodes directly. This group of algorithms is more applicable in large *ad hoc* networks. In [41], the positioning algorithms are further classified through the ranging technique used (distance measurements or connectivity derived from the graph associated with the network) and the kind of coordinate systems provided. These include: *absolute*, *local* or *relative* positioning. An *absolute* coordinate system requires global coherence of the estimation. This is desirable for most applications but is expensive in communication cost, especially in mobile environments, since all new positions must be coherent with the previous ones. A *relative* coordinate system, provides a relative position which is local to the network; the coordinates could be arbitrary but network wide coherence is still guaranteed. A *local* coordinate system provides just local coherence between singular clusters of nodes, with respect to each other.

As the reader can infer, clearly implementation costs, and hardware/software challenges vary considerably according to application purposes and needs. A literature review of the available algorithms is presented in the following subsections.

Lateration

This technique is based on the idea of locating a node using overlapping circles (at least three) with rays equal to the estimated distances among nodes. This is discussed fully in section 5.1.1.

Minimum Maximum Algorithm (MINMAX)

A simpler method than lateration is presented in [53] as part of the N-hop multilateration approach. This algorithm (also known as *Bounding Box*) is an easy method and it is used in some systems for locating initial estimates quickly but not very carefully. For all mobile nodes, an anchor node measures the power received, and the computation of the distance allows a square to be drawn around the fixed node, with sides equal to twice the estimated distance. The anchor node assumes that the mobile is within this region. The same thing happens for all other mobile nodes. By overlapping (*bounding*) the square regions (*boxes*) it is possible to identify a mobile node within a rectangle whose area gets smaller as the number of anchor nodes increases. The output of the algorithm is the centroid of this rectangle. The main advantage of this technique relies on its ease of implementation, as the intersection of all bounding boxes can be easily computed without any need for floating point operations. However, the minimum maximum algorithm is characterized by a bias. If an anchor node assumes that the mobile position is represented by a square region, instead of a circumference around itself (as in *lateration*), there is an additive estimation error [36]. As a consequence the bias impairs the output of the algorithm, even in cases where there is no fading or shadowing.

Ring Overlapping Algorithm (ROCRSSI+)

The general idea of ROCRSSI+ is that each sensor node uses a series of overlapping rings to narrow down the possible area in which it resides. This is achieved by comparing the received power from the anchor nodes. Therefore it does not need an accurate estimate of the distance among nodes. Each anchor node can draw circles around itself by measuring the received power to any other fixed node. The circumferences of these circles are centered on an anchor node and ideally, i.e. in a channel which only contains path loss, will intersect the others. In this way each anchor node draws boundaries around itself that identify different regions of the space. By comparing the received signal strength from the mobile node (for all the mobile nodes) with the one coming from the other anchor nodes, each anchor node can apportion the mobile node to one of these regions. Repeating the same principal for all

anchor nodes, we arrive at a space comprising of several overlapping regions, i.e. a slice. The output of this algorithm is the centroid of the region slice where the majority of anchor nodes have established the mobile to be. The power of this algorithm comes from the fact that it is ranging free”, i.e., it does not need to estimate the distances among nodes and just compares the received powers. However, it has been shown in [36] that the localization precision of this method changes significantly according to the anchor topology and the mobile position.

Convex Optimization

This is a centralized method, simple to implement with most types of ranging techniques. It can use both linear programming and semidefinite programming (*SDP*) since there are efficient computational methods available for most convex programming problems. In *SDP*, the optimality is achieved at the cost of centralization and the need to manage large data structures. The complexity is at least quadratic with the number of connections, and it uses connectivity information or graph theory as the ranging technique. In the centralized algorithm of Doherty *et al* [12], a method for estimating unknown node positions in a sensor network based exclusively on connectivity-induced constraints is described and solved as a convex optimization problem through *SDP*.

Multi-Dimensional Scaling (MDS-MAP)

This is another centralized algorithm that makes use of connectivity only to provide positions in a network with or without landmarks [54]. This method operates in three stages: first the shortest path between all pairs of nodes is computed, then a relative map is built for the network nodes, finally this map is converted into an absolute map if three (or more) landmarks are available. The advantage of this method is its wide range of applicability; it can use both connectivity and distance measurement ranging techniques and provides both absolute and relative positioning. The complexity required is quite high and is at least cubic with respect to the number of nodes, but, unlike the previous algorithm, it has a theoretical bound.

Global Position System (GPS)

Like the previously cited *lateration*, *MinMax*, and *RocRSSI+* methods, is a “one hop” positioning algorithm. A dominant technical requirement for GPS is line of sight (LOS). The positioning service is available, through multilateration, when at least four satellites are visible; three to find the coordinate of the receiver, and a fourth one to determine the clock bias on the receiver. So, in environments such as: indoors, under foliage or in the shadow of buildings it is not possible to get a position. Furthermore, when very small sensors are required, size and cost prohibit the inclusion of a GPS receiver. Besides this, the integration of a GPS

device to the sensor nodes will increase the energetic consumption of the sensor, decreasing the operational time of the node [35].

Lighthouse Location System

This is a curious one hop positioning method that achieves the positioning of an entire field of sensors using a "lighthouse" base station (BS). Using a parallel beam that rotates at a constant speed the BS sweeps over the entire set of nodes. By knowing the rotational speed and the width of the beam, and measuring the length of time it sees the light, each node (equipped with a clock and a photo detector) is able to independently estimate its range to the BS. This effectively places it on a circle around the lighthouse. Collecting three of these estimations independently of each other (using three different lighthouses or mutually perpendicular light beams from the same lighthouse) it is possible for a node using trilateration to determine its position [50].

Ad Hoc Localization System (AHLoS)

This is a distributed algorithm [52] consisting of several types of multilateration: atomic, iterative and collaborative. In atomic multilateration, anchor node density is high such that a node has enough neighbours to apply basic trilateration. Nodes that manage to obtain a location begin to behave as landmarks starting the iterative multilateration. However even after applying these two methods there may still be nodes unable to estimate their position, and the problem must be considered in a collaborative fashion. The disadvantage of AHLoS is that it requires a high percentage of anchor nodes in order to achieve an acceptable number of resolved nodes. The overriding advantage is that given a good ranging method, it is likely to produce high-quality position estimates.

Ad Hoc Positioning System (APS)

This is a hybrid distributed, multi-hop algorithm [42] comprising of two ideas: distance vector (DV) routing and beacon based positioning (like GPS). What makes it similar to DV is that information is forwarded hop by hop. What makes it similar to GPS is that eventually each node estimates its own position based on the landmark readings it gets. The advantage of the APS method is that it is both distributed and localized. It supports limited mobility and has variable duty cycles for power-saving oriented applications. It is suitable for the majority of ranging techniques. The disadvantages are: it requires a fairly uniform distribution of anchor nodes, and, due its DV nature, it will face increasing costs in the face of high mobility.

Self Positioning Algorithm (SPA) - Local Positioning System (LPS)

These algorithms simply find positions within a coordinate system of an identified group of nodes called the "location reference group" (LRG). In an SPA a node achieves relative positioning with respect to its neighbors, by iteratively exchanging tables which contain particular information. In LPS a node uses capabilities such as RSS, AoA or magnetic compasses to estimate its position in a local coordinate system. This method allows positioning to be achieved only for nodes participating in packet forwarding, thereby reducing traffic load in the network.

Once each node independently establishes its local coordinate system, an alignment procedure initiated by the LRG aligns all other coordinate systems to the reference group. The advantage of both these methods is that they provide a network-wide coherence in position without the need for landmarks. SPA faces high cost as the mobility increases, whilst LPS cannot provide a globally coherent coordinate system.

3.3.3. Refinement

Finally the refinement stage aims to minimize the estimation error variance. In a centralized approach to localization, the central processing entity can parse data coming from the whole network, and the optimization can be treated as a convex programming problem as seen for the convex optimization. Alternatively in distributed localization, the refinement follows the positioning stage and can be performed either independently by each node or collaboratively. The following subsections identify the commonly used refinement methods.

Optimization of a global cost function

Some algorithms try to find the optimum of a global cost function, e.g., Least Squares (LS), Weighted Least Squares (WLS) or Maximum Likelihood (ML). These methods are very accurate, but they involve a heavy computational load that is not always achievable in sensor nodes. References [59] and [46] illustrate some applications of this approach.

Bayesian Estimators

These algorithms try to maximize the a posteriori probability of the estimation, given measured data (positioning measurements), using Bayes rule [39] to incorporate some a priori information about the model, e.g. density of the nodes. An example of this kind of estimator is the DT/A Algorithm, proposed in [40] and based on a Hidden Markov Model (HMM). The implementation allows the joint tracking of both a mobile terminal and the visibility condition in LOS/NLOS indoor environment. Another application of these methods can be found in [39], where the algorithm acts as a refinement of an existing sampling method denominated as progressive correction.

Cooperative Localization

After each sensor has estimated its location, it then transmits the assertion to its neighbours, which must then recalculate their location and transmit again, until convergence occurs. This technique is required, to implement the AHLoS positioning algorithm [53], but can also be found in [43] and [45].

Particle Filters - Monte Carlo Localization

In the article [20], Hu and Evans present a range-free localization algorithm for mobile sensor networks based on the Sequential Monte Carlo method [4]. The Monte Carlo method has been extensively used in robotics [4] where a robot estimates its localization based on its motion, perception and possibly a pre-learned map of its environment. Hu and Evans extend the Monte Carlo method as used in robotics to support the localization of sensors in a free, unmapped terrain. The authors assume a sensor has little control and knowledge over its movement, in contrast to a robot. They target an environment where there is no hardware for obtaining ranging information, the topology of the network is unknown and most likely irregular, the density of anchors is low and both anchors and sensor nodes can move in an uncontrollable manner. The only assumption that is made is that the sensors or anchors move with a known maximum speed and that the radio range is common to the sensors and anchors, or is distributed together with other messages. This latter point, however, is not described by the authors.

Using the sequential Monte Carlo Localization (MCL), Hu and Evans want to take advantage of mobility to improve the accuracy of localization and reduce the number of anchor nodes that are required in the network. The key idea of the sequential Monte Carlo Localization is to represent the posterior distribution of the possible locations of a node using a set of weighted samples. Localization happens in two steps. First, the prediction step leads to choosing a set of samples representing the belief of the node regarding its location. During the prediction step, a node picks random locations within the deployment area, possibly constrained by its maximum speed and the previous location samples. Second, the filtering step aims at removing the impossible locations from the set of samples. The filtering is done using information obtained from the environment, such as the location of the anchors in the case of a sensor node or the detection of landmarks in the case of a mobile robot. The process repeats and the sensor or robot is able to update its position estimation.

Although this approach has been widely studied, there are some drawbacks for its implementation [62]. First, a sufficient number of anchors are required for the algorithms. The position estimation of MCL depends on local anchor information; hence, the location error could be large when the density of anchor nodes is low. Although MCL is an extension approach that includes information about the neighbours, the problem is not solved completely. Accord-

ing to the simulation results of MCL [20], MCL needs more than one anchor in a one-hop transmission range to obtain reasonable accuracy, and this number of anchors is relatively small but not sufficient to apply the algorithm in a real environment. As a second problem, the particle filter based algorithms assume that the fixed radio transmission range is known. In a real environment, however, the radio range changes due to residual battery, geometric characteristics, and many other factors.

Kalman Filters

The Kalman Filter is a digital filter that provides an efficient recursive means to estimate the state of a process, minimizing the mean of the squared error. We implemented some derivations of this refinement method to improve the performance of the positioning algorithms considered, as it is well suited to the tracking application. In section 5.1.2 we report a detailed description of this filter. Other examples of how the Kalman filter can be used to track mobile terminals can be found in references [36], [59], [37].

Summary

Localization is the process of determining a target position at a particular moment in time. This is one of the main challenges for the Ad Hoc wireless sensor networks, because of both the network topology and the difficulty of the estimation of the distance between nodes. The three phases of a localization process are:

- Ranging
- Positioning
- Refinement

In spite of the extensive search of a solution for the localization problem, this remains as one of the open research issues for wireless networks. There are many proposed schemes to solve this problem, but there is no complete solution because every proposed system has its own drawbacks.

4 Mobility and observation models for mobile node localization

For the successful tracking of a moving target it is essential to extract the maximum useful information about the target state from the available observations. Both good models to describe the target dynamics and sensor will certainly help this information extraction. As the knowledge of information on the kinematics of the target and sensor characteristics are generally known, most of the tracking algorithms base their performance on the a priori defined mathematical model of the target which are assumed to be sufficiently accurately. This section, addresses the problem of describing the target motion model and establishes a good compromise between accuracy and complexity.

4.1. Localization of mobile nodes

We consider the two-dimensional (2-D) problem of cooperative localization [45] of mobile nodes. The vector $\mathbf{X} = \{(x_1, y_1), (x_2, y_2), \dots, (x_n, y_n)\}$ of positions of n mobile nodes is estimated given r reference coordinates $\mathbf{X}^r = \{(x_{n+1}, y_{n+1}), (x_{n+2}, y_{n+2}), \dots, (x_{n+r}, y_{n+r})\}$ previously known and measurements $\{X_{i,j}\}$, where $X_{i,j}$ is a measurement between devices i and j . In this work, besides the mobile nodes positions, we can also estimate their speeds and accelerations depending on the mobility models used. We focus our attention on cases in which mobile sensor nodes receive distance measurements from a subset of reference sensor nodes in the network. This includes applications in which each sensor is equipped with a wireless transceiver and the distance is estimated by received signal strength indicators or time delay of arrival between sensor locations. We consider a sensor deployment architecture as described in [37] (see Figure 4-1). Received signal strength indicators are processed in a central information processor (CIP) through a low-power communication network. We allow unknown-location devices to make measurements with known-location references. In cooperative localisation, unknown-location devices might be allowed to make measurements with other unknown-location devices. The additional information gained from the measurements between pairs of unknown-location devices might enhance the accuracy and robustness of the localisation system. However, apart from the increased energy necessary for communications, the complexity of the localisation algorithms also increases. In this work we limit our consideration to the case without communications between the mobile nodes.

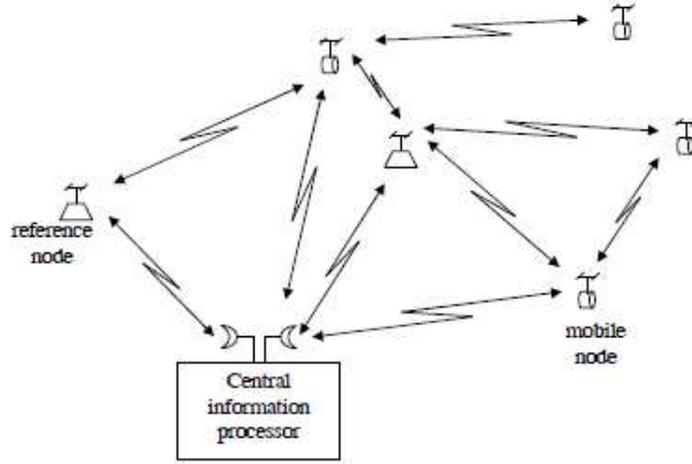


Figure 4-1: *Ad-hoc* sensor network architecture

4.2. Discrete-Time State Space Models

In this work we will consider only discrete-time state space models. The models are defined recursively in terms of distributions as shown in equation (4-1).

$$\begin{aligned} x_k &\sim p(x_k | x_{k-1}) \\ y_k &\sim p(y_k | x_k) \end{aligned} \quad (4-1)$$

In (4-1),

- $x_k \in R^n$ is the *state* of the system at time step k .
- $y_k \in R^m$ is the *output* or *measurement* of the system at time step k .
- $p(x_k | x_{k-1})$ is the dynamic model which characterizes the dynamic behaviour of the system. Usually the model is a probability density (continuous state), but it can also be a counting measure (discrete state), or a combination of them, if the state is both continuous and discrete.
- $p(y_k | x_k)$ is the model for measurements, which describes how the measurements are distributed given the state. This model characterizes how the dynamic model is perceived by the observers.

A system defined this way has the so called *Markov*-property, which means that the state x_k given x_{k-1} is independent from the history of states and measurements, which can also be expressed with the following equality:

$$p(x_k | x_{1:k-1}, y_{1:k-1}) = p(x_k | x_{1:k-1}). \quad (4-2)$$

The past events do not depend on the future given the present, which is the same as

$$p(x_{k-1} | x_{k:T}, y_{k:T}) = p(x_{k-1} | x_k). \quad (4-3)$$

The same applies also to measurements, meaning that the measurement y_k is independent from the histories of measurements and states, which can be expressed with the equality

$$p(y_k | x_{1:k}, y_{1:k-1}) = p(y_k | x_k). \quad (4-4)$$

In actual application problems, we are interested in predicting and estimating the state of a dynamic system given the measurements obtained so far. In probabilistic terms, we are interested in the predictive distribution for the state at the next time step

$$p(x_k | y_{1:k-1}), \quad (4-5)$$

and in the marginal posterior distribution for the state at the current time step

$$p(x_k | y_{1:k}). \quad (4-6)$$

The formal solutions for these distribution are given by the following recursive Bayesian filtering equations (e.g. [51]):

$$p(x_k | y_{1:k-1}) = \int p(x_k | x_{k-1})p(x_{k-1} | y_{1:k-1})dx_{k-1}, \quad (4-7)$$

and

$$p(x_k | y_{1:k}) = \frac{1}{Z_k} p(y_k | x_k) p(x_k | y_{1:k-1}), \quad (4-8)$$

where the normalization constant Z_k is given as

$$Z_k = \int p(y_k | x_k) p(x_k | y_{1:k-1}) dx_k. \quad (4-9)$$

In many cases we are also interested in smoothed state estimates of previous time steps given the measurements obtained so far. In other words, we are interested in the marginal posterior distribution

$$p(x_k | y_{1:T}), \quad (4-10)$$

where $T > k$. As with the filtering equations above also in this case we can express the formal solution as a set of recursive Bayesian equations (e.g. [51]):

$$\begin{aligned} p(x_{k+1} | y_{1:k}) &= \int p(x_{k+1} | x_k) p(x_k | y_{1:k}) dx_k \\ p(x_k | y_{1:T}) &= p(x_k | y_{1:k}) \int \left[\frac{p(x_{k+1} | x_k p(x_{k+1} | y_{1:T}))}{p(x_{k+1} | y_{1:k})} \right] dx_{k+1}. \end{aligned} \quad (4-11)$$

4.2.1. Linear state space estimation

The simplest of the state space models considered in this documentation are linear models, which can be expressed with equations of the following form:

$$\begin{aligned} x_k &= A_{k-1}x_{k-1} + q_{k-1} \\ y_k &= H_k x_k + r_k \end{aligned} \tag{4-12}$$

where

- $x_k \in R^n$ is the *state* of the system on time step k .
- $y_k \in R^m$ is the *state* of the system on time step k .
- $q_{k-1} \sim N(0, Q_{k-1})$ is the process noise on time step $k - 1$.
- $r_k \sim N(0, R_k)$ is the measurement noise on time step k .
- A_{k-1} is the transition matrix of the dynamical model.
- H_k is the measurement model matrix.
- The prior distribution for the state is $x_0 \sim N(m_0, P_0)$, where parameters m_0 and P_0 are set using the information known about the system under study.

The model can also be equivalently expressed in probabilistic terms with distributions

$$\begin{aligned} p(x_k | x_{k-1}) &= N(x_k | A_{k-1}x_{k-1}, Q_{k-1}) \\ p(y_k | x_k) &= N(y_k | H_k x_k, R_k). \end{aligned} \tag{4-13}$$

4.2.2. Discretization of continuous-time linear time-invariant systems

Often many linear time-invariant models are described with continuous-time state space equations of the following form:

$$\frac{dx(t)}{dt} = Fx(t) + Lw(t), \tag{4-14}$$

where

- the initial conditions are $x(0) \sim N(m(0), P(0))$,
- F and L are constant matrices, which characterize the behaviour of the model,
- $w(t)$ is a white noise process with a power spectral density Q_c .

To be able to use the Kalman filter defined in the next section the model 4-14 must be discretized somehow, so that it can be described with a model of the form 4-12. The solution for the discretized matrices A_k and Q_k can be given as [51] [5]:

$$A_k = \exp(F\Delta t_k) \quad (4-15)$$

$$Q_k = \int_0^{\Delta t_k} \exp(F(\Delta t_k - \tau)LQ_cL^T \exp(F(\Delta t_k - \tau))^T d\tau, \quad (4-16)$$

where $\Delta t_k = t_{k+1} - t_k$ is the stepsize of the discretization. In some cases the Q_k can be calculated analytically, but in cases where it is not possible, the matrix can still be calculated efficiently using the following matrix fraction decomposition:

$$\begin{bmatrix} C_k \\ D_k \end{bmatrix} = \exp\left(\begin{bmatrix} F & LQ_cL^T \\ 0 & -F^T \end{bmatrix} \Delta t_k\right) \begin{bmatrix} 0 \\ I \end{bmatrix}. \quad (4-17)$$

The matrix Q_k is then given as $Q_k = C_k D_k^{-1}$.

4.3. Mobility Models

4.3.1. CWPA Model

Consider a case where we track an object moving on a two-dimensional space with a sensor, which gives measurements of the target position in Cartesian coordinates x and y . In addition to, position target also has state variables for its velocities and accelerations toward both coordinate axes, \dot{x} , \dot{y} , \ddot{x} and \ddot{y} . In other words, the state of a moving object on time step k can be expressed as a vector

$$X_k = [x_k \ y_k \ \dot{x}_k \ \dot{y}_k \ \ddot{x}_k \ \ddot{y}_k]^T. \quad (4-18)$$

In continuous case the dynamics of the target motion can be modelled as a linear, time-invariant system

$$\frac{dx(t)}{dt} = \begin{bmatrix} 0 & 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 & 0 & 0 \end{bmatrix} x(t) + \begin{bmatrix} 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 0 & 0 \\ 1 & 0 \\ 0 & 1 \end{bmatrix} w(t), \quad (4-19)$$

where $x(t)$ is the target state on the time t and $w(t)$ is a white noise process with power spectral density

$$Q_c = \begin{bmatrix} q & 0 \\ 0 & q \end{bmatrix}. \quad (4-20)$$

As can be seen from the equation the acceleration of the object is perturbed with a white noise process and hence this model has the name continuous Wiener process acceleration (CWPA) model. There are also other similar models, for example, the continuous white noise acceleration (CWNA) model [5], where the velocity is perturbed with a white noise process.

To be able to estimate this system, the differential equation defined above must be discretized to get a discrete-time state equation of the form of (4-12). It turns out, that the matrices A and Q can be calculated analytically with equations (4-15) and (4-16) to give the following:

$$A = \begin{bmatrix} 1 & 0 & \Delta t & 0 & \frac{1}{2}\Delta t^2 & 0 \\ 0 & 1 & 0 & \Delta t & 0 & \frac{1}{2}\Delta t^2 \\ 0 & 0 & 1 & 0 & \Delta t & 0 \\ 0 & 0 & 0 & 1 & 0 & \Delta t \\ 0 & 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 \end{bmatrix} \quad (4-21)$$

$$Q = \begin{bmatrix} \frac{1}{20}\Delta t^5 & 0 & \frac{1}{8}\Delta t^4 & 0 & \frac{1}{6}\Delta t^3 & 0 \\ 0 & \frac{1}{20}\Delta t^5 & 0 & \frac{1}{8}\Delta t^4 & 0 & \frac{1}{6}\Delta t^3 \\ \frac{1}{8}\Delta t^4 & 0 & \frac{1}{6}\Delta t^3 & 0 & \frac{1}{2}\Delta t^2 & 0 \\ 0 & \frac{1}{8}\Delta t^4 & 0 & \frac{1}{6}\Delta t^3 & 0 & \frac{1}{2}\Delta t^2 \\ \frac{1}{6}\Delta t^3 & 0 & \frac{1}{2}\Delta t^2 & 0 & \Delta t & 0 \\ 0 & \frac{1}{6}\Delta t^3 & 0 & \frac{1}{2}\Delta t^2 & 0 & \Delta t \end{bmatrix} q. \quad (4-22)$$

4.3.2. CWPV Model

In the *Continuous Wiener Process Velocity* (CWPV) model, the state of the target at time step k consists of the position in two dimensional Cartesian coordinates x_k and y_k and the velocity toward those coordinate axes, \dot{x}_k and \dot{y}_k . Thus, the state vector can be expressed as

$$X_k = [x_k \quad y_k \quad \dot{x}_k \quad \dot{y}_k]^T. \quad (4-23)$$

Established in an analogue way as the CWPA model of the previous section, the dynamic of the target is modelled as a linear, discretized Wiener velocity model

$$X_k = \begin{bmatrix} 1 & 0 & \Delta t & 0 \\ 0 & 1 & 0 & \Delta t \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{bmatrix} \begin{bmatrix} x_{k-1} \\ y_{k-1} \\ \dot{x}_{k-1} \\ \dot{y}_{k-1} \end{bmatrix} + \mathbf{q}_{k-1}, \quad (4-24)$$

where \mathbf{q}_{k-1} is Gaussian process noise with zero mean and covariance

$$Q = \begin{bmatrix} \frac{1}{3}\Delta t^3 & 0 & \frac{1}{2}\Delta t^2 & 0 \\ 0 & \frac{1}{3}\Delta t^3 & 0 & \frac{1}{2}\Delta t^2 \\ \frac{1}{2}\Delta t^2 & 0 & \Delta t & 0 \\ 0 & \frac{1}{2}\Delta t^2 & 0 & \Delta t \end{bmatrix} q \quad (4-25)$$

where q is the spectral density of the noise.

4.3.3. Position Model

In the *Position* (P) model, the state of the target at time step k consists only of the position in two dimensional Cartesian coordinates x_k and y_k . Thus, the state vector can be expressed as

$$X_k = [x_k \quad y_k]^T. \quad (4-26)$$

Established in an analogue way as *Random Walk* model, the dynamic of the target is modelled as a linear, discretized Wiener model

$$X_k = \begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix} \begin{bmatrix} x_{k-1} \\ y_{k-1} \end{bmatrix} + \mathbf{q}_{k-1}, \quad (4-27)$$

where \mathbf{q}_{k-1} is Gaussian process noise with zero mean and covariance

$$Q = \begin{bmatrix} \Delta t & 0 \\ 0 & \Delta t \end{bmatrix} q \quad (4-28)$$

where q is the spectral density of the noise.

4.3.4. Coordinated Turn Model

A common way of modelling a turning object is to use the *coordinated turn model* [5]. The idea is to augment the state vector with a turning rate parameter ω , which has to be estimated along with the other system parameters (the position and the velocity of the target). Thus, the joint system vector can be expressed as

$$X_k = [x_k \quad y_k \quad \dot{x}_k \quad \dot{y}_k \quad \omega_k]^T. \quad (4-29)$$

The dynamic model for the coordinated turns is

$$X_{k+1} = \begin{bmatrix} 1 & 0 & \frac{\sin(\omega_k \Delta t)}{\omega_k} & \frac{\cos(\omega_k \Delta t) - 1}{\omega_k} & 0 \\ 0 & 1 & \frac{1 - \cos(\omega_k \Delta t)}{\omega_k} & \frac{\sin(\omega_k \Delta t)}{\omega_k} & 0 \\ 0 & 0 & \cos(\omega_k \Delta t) & -\sin(\omega_k \Delta t) & 0 \\ 0 & 0 & \sin(\omega_k \Delta t) & \cos(\omega_k \Delta t) & 0 \\ 0 & 0 & 0 & 0 & 1 \end{bmatrix} X_k + \begin{bmatrix} 0 \\ 0 \\ 0 \\ 0 \\ 1 \end{bmatrix} v_k, \quad (4-30)$$

where $v_k \sim N(0, \sigma_\omega^2)$ is univariate white Gaussian process noise for the turn rate parameter. This model is, despite the matrix form, non-linear.

4.4. Measurement Model

In a wireless *ad hoc* network, the distance between the mobile and a reference (reachable) node can be inferred from the received signal strength indicator (RSSI) or pilot signal of the node. The RSSI, measured in dB, received at the mobile node N_i from the node N_j with coordinates $(x_{j,k}, y_{j,k})$ at time k is given by [37]

$$z_{ij,k} = \kappa_j - 10\gamma \log_{10}(d_{ij,k}) + v_{i,k}, \quad (4-31)$$

where κ_j is a constant determined by the transmitted power, wavelength, and gain of the node N_j , γ is the slope index (typically $\gamma = 2$ for highways and $\gamma = 4$ for microcells in a city), $v_{i,k}$ is the logarithm of the shadowing component which is found to be a zero mean, stationary Gaussian process with standard deviation σ_v (from 4-8 dB), and $d_{ij,k}$ is the distance between the mobile nodes N_i and N_j

$$d_{ij,k} = \sqrt{(x_{j,k} - x_{i,k})^2 + (y_{j,k} - y_{i,k})^2}. \quad (4-32)$$

All nodes in the group send their pilot signal strength measurements (reference strength indicator signals) to the reference nodes. In the case with n mobile nodes, and r fixed nodes having known positions and without communications between the fixed nodes, the overall observation vector has the form

$$z_k = \{z_{ij,k}\}_{ij=1}^L, \quad (4-33)$$

where $L = n * r$ is the number of measurements. In the developed solutions we do not consider communications between mobile nodes, hence we exclude the measured distances between them. The vector form of the observation equation (4-31) is

$$z_k = H_k x_k + v_k, \quad (4-34)$$

where the noise v_k consists of the path loss and the shadowing component.

Summary

This chapter presented the mathematical fundamentals for the study of localization problem in Ad Hoc wireless sensor networks. Starting from a discrete-time state space representation, the Wiener Process and the Coordinated Turn models are discussed. Also, the measurement model based on the signal strength indicator was established.

5 The application of Kalman-based schemes to localization

5.1. Theoretical tools

5.1.1. Lateration

This algorithm is based on a very simple idea. For all the mobile nodes, each anchor node measures the power received from the mobile and draws a circle around itself by computing the distance (inversion of the path loss model formula). The anchor node assumes that the mobile should be at one of the points on the circumference of the circle. If the measurement of the distances is exact (i.e. there is no fading and no shadowing) it is possible to discover where is the mobile node by the intersection of three circles, given three anchor nodes.

To make the implementation easier, pairs of anchor nodes were considered rather than triplets. In such a manner, for each pair of anchor nodes it should be possible to calculate two points where the circles overlap: one of this is correct, the other one is false. Then, for each pair of anchor nodes it is possible to collect two estimated positions for all the mobile nodes considered. Referring to the collected estimates of an individual mobile node, the objective is to find which ones are correct and which are false.

The main problem associated with the lateration positioning is that in the presence of fading and shadowing this method may produce no results at all, i.e. the error in the distance estimation can prevent the circles overlapping.

5.1.2. Kalman Filter

In the 1960s R.E. Kalman published a paper entitled *A new Approach to Linear Filtering and Prediction Problems* (see [26]). In this work the author presented a novel filtering method to process data for solving problems that deal with measurements hampered by random biases, with the goal of providing a recursive solution for estimation of *linear* discrete-time dynamic systems. A Kalman Filter is *an optimal recursive data processing algorithm*; each of these words composes the magical formula that makes one aware of what this filter is:

- "data processing algorithms", this simply means that the Kalman Filter is a set of

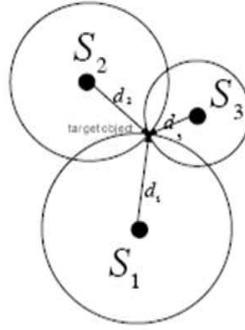


Figure 5-1: The *lateration* process

mathematical formulas implemented in software, and the filtering process is therefore digital;

- ”**recursive**”, that is, the filter repeats at each iteration (i.e. at each measurement done) the same operations;
- ”**optimal**”, namely, the Kalman filter achieves the best performance if we assume the observed process is linear and the measurement noise is white and gaussian.

The purpose of the filter is to estimate a quantity by both the measurement and the *a priori* knowledge about the observed phenomenon. The principle followed is to collect and wisely process the available information. The data process is made up of two stages:

- **Prediction:** a function estimates the state of the Kalman Filter, projecting the quantity for future estimates. During this stage the measurement value of the quantity is also collected; these steps translate to equations as follows

$$\begin{aligned} m_k^- &= A_{k-1}m_{k-1} \\ P_k^- &= A_{k-1}P_{k-1}A_{k-1}^T + Q_{k-1}. \end{aligned} \quad (5-1)$$

- **Update:** the data collected at the previous stage is weighted by an appropriate coefficient called the *Kalman Gain*, determining the estimation for the observed value. The equations for this stage are:

$$\begin{aligned} v_k &= y_k - H_k m_k^- \\ S_k &= H_k P_k^- H_k^T + R_k \\ K_k &= P_k^- H_k^T S_k^{-1} \\ m_k &= m_k^- + K_k v_k \\ P_k &= P_k^- - K_k S_k K_k^T, \end{aligned} \quad (5-2)$$

where

- m_k^- and P_k^- are, respectively, the predicted mean and covariance of the state on the time step k before taking into account the current measurement.
- m_k and P_k are, respectively, the estimated mean and covariance of the state on time step k after taking into account the current measurement.
- v_k is the innovation or the measurement residual on time step k .
- S_k is the measurement prediction covariance on the time step k .
- K_k is the filter gain, an indicator of how much the predictions should be corrected on time step k .

The measurement function is responsible for projecting the state of the filter into the future and contains information about the observed process. This function describes a mathematical model of the measurable monitored phenomenon. In this way we can interpret the measurement obtained dynamically and coherently. The *Kalman Gain* is composed of quantities that are related to uncertainty both of the measurements and of the model used. We are assuming an environment impaired by white gaussian noise; therefore we can take into account both measurement and projection uncertainties referring to the noise and the covariances of the model respectively.

In certain cases a Gaussian assumption is too primitive, but a Gaussian process only requires the first and second order statistics (i.e. mean and covariance) to be statistically described. The Gaussian assumption often allows complex problems to become more tractable. Furthermore, because the real data are influenced by many variables (e.g., scatter noise, multi-path channel, measurement noise, used device bias, etc.) we can utilize the *Central Limit Theorem*, which is at the core of probability theory. This theorem [36], proves that when a number of independent random variables are added together, their overall effect can be described well using a Gaussian probability density.

The assumption of white noise leads to some contradictions. This supposition implies noise values are not correlated in time, e.g. a channel can change from one state to another one abruptly. The assumption of whiteness also implies that all frequencies have the some power, i.e. a noise with infinite power. Despite this, the white-noise model is still useful both for treatment simplicity and because we can overcome the above problems using some "tricks". For instance, any physical system of interest has a bandpass frequency response; we can just disregard the frequencies outside this band to get away from the requirement to consider infinite noise power.

Note that in this case the predicted and estimated state covariances on different time steps do not depend on any measurement, so that they could be calculated off-line before making any measurements provided that the matrices \mathbf{A} , \mathbf{Q} , \mathbf{R} and \mathbf{H} are known on those particular time steps. It is also possible to predict the state of system as many steps ahead as wanted just by looping the predict step of Kalman filter, but naturally the accuracy of the estimate decreases with every step.

5.1.3. Kalman Filter for Tracking Problems

A tracking problem, particularly in a wireless environment, is a complex task indeed. On one hand the power measurements, which are required to determine the mobile positions, are hampered by more complex noise than white-gaussian; and on the other the algorithm and the hardware used can introduce bias that is difficult to control. Furthermore trying to model animated motion is a difficult challenge since feelings, needs, thoughts, and instincts can make the movements (and of the mobile carried) completely unpredictable. Despite this, the trade off between simplicity and reliability leads to investigate the Kalman filter capabilities to track a mobile terminal along an unknown path. The use of the Kalman Filter to assist in tracking mobile nodes has been proposed before; see [36], [59], [37].

5.1.4. Kalman Filters for Non-linear State Estimation

In many cases interesting dynamic systems are not linear by nature, so the traditional Kalman filter cannot be applied in the estimation of the state of such systems. In these kind of systems, one or both the dynamics and the measurement processes can be non-linear. In this section we describe two extensions to the traditional Kalman filter, which can be applied for estimating non-linear dynamical systems by forming Gaussian approximations to the joint distribution of the state x and measurement y . First we present the Extended Kalman filter (EKF), which is based on Taylor series approximation of the joint distribution, and then the Unscented Kalman filter (UKF), which is based on the unscented transformation of the joint distribution.

5.1.5. Extended Kalman Filter (EKF)

The extended Kalman filter (see [5], [51]) extends the scope of Kalman filter to non-linear optimal filtering problems by forming a Gaussian approximation to the joint distribution of state x and measurements y using a Taylor series based transformation. A first order EKF is presented using linear approximation; higher order filters are also possible, but not presented here. The filtering model used in the EKF is

$$\begin{aligned} x_k &= f(x_{k-1}, k-1) + q_{k-1} \\ y_k &= h(x_k, k) + r_k, \end{aligned} \tag{5-3}$$

where $x_k \in \mathfrak{R}^n$ is the state, $y_k \in \mathfrak{R}^m$ is the measurement and $q_{k-1} \sim N(0, Q_{k-1})$ is the process noise, $r_k \sim N(0, R_k)$ is the measurement noise, f is the (possibly non-linear) dynamic model function and h is the (again possibly non-linear) measurement model function. The first order extended Kalman filter approximate the distribution of state x_k given the observations $y_{1:k}$ with a Gaussian:

$$p(x_k | y_{1:k-1}) \approx N(x_k | m_k, P_k). \quad (5-4)$$

Taylor Series Based Linear Approximation

A linear approximation for the distribution of variable y , which is generated with a non-linear transformation of a Gaussian random variable x , is obtained as follows:

$$\begin{aligned} x_k &\sim N(m, P) \\ y_k &= g(x), \end{aligned} \quad (5-5)$$

where $x \in \mathfrak{R}^n$, $y \in \mathfrak{R}^m$ and $g : \mathfrak{R}^n \mapsto \mathfrak{R}^m$ is a general non-linear function. Solving the distribution of y in a formal way is in general not possible, because it is non-Gaussian for all by linear g , so in practice it must be approximated somehow. The joint distribution of x and y can be formed with, for example, linear and quadratic approximations [5].

The linear Gaussian approximation of the joint distribution of variables x and y defined by equations (5-5) is given as

$$\begin{pmatrix} x \\ y \end{pmatrix} = N \left(\begin{pmatrix} m \\ \mu_L \end{pmatrix}, \begin{pmatrix} P & C_L \\ C_L^T & S_L \end{pmatrix} \right), \quad (5-6)$$

where

$$\begin{aligned} \mu_L &= g(m) \\ S_L &= G_x(m) P G_x^T(m) \\ C_L &= P G_x^T(m), \end{aligned} \quad (5-7)$$

and $G_x(m)$ is the Jacobian matrix of g with elements

$$[G_x(m)]_{j,j'} = \left. \frac{\partial g_j(x)}{\partial x_{j'}} \right|_{x=m}. \quad (5-8)$$

Extended Kalman Filter Procedure

Like Kalman filter, also the extended Kalman filter is separated into two steps. The steps for the first order EKF are:

- **Prediction:**

$$\begin{aligned} m_k^- &= f(m_{k-1}, k-1) \\ P_k^- &= F_x(m_{k-1}, k-1) P_{k-1} F_x^T(m_{k-1}, k-1) + Q_{k-1}. \end{aligned} \quad (5-9)$$

■ **Update:**

$$\begin{aligned}
 v_k &= y_k - h(m_k^-, k) \\
 S_k &= H_x(m_k^-, k)P_k^- H_x^T(m_k^-, k) + R_k \\
 K_k &= P_k^- H_x^T(m_k^-, k)S_k^{-1} \\
 m_k &= m_k^- + K_k v_k \\
 P_k &= P_k^- - K_k S_k K_k^T,
 \end{aligned} \tag{5-10}$$

where the matrices $F_x(m, k-1)$ and $H_x(m, k)$ are the Jacobians of f and h , with elements

$$[F_x(m, k-1)]_{j,j'} = \left. \frac{\partial f_j(x, k-1)}{\partial x_{j'}} \right|_{x=m} \tag{5-11}$$

$$[H_x(m, k)]_{j,j'} = \left. \frac{\partial h_j(x, k)}{\partial x_{j'}} \right|_{x=m}. \tag{5-12}$$

Note that the difference between first order EKF and KF is that the matrices A_k and H_k in KF are replaced with Jacobian matrices $F_x(m_{k-1}, k-1)$ and $H_x(m_k^-, k)$ in EKF. Predicted mean m_k^- and residual of prediction v_k are also calculated differently in the EKF.

The Limitations of EKF

As discussed in [25], the EKF has some serious drawbacks, which should be kept in mind when it is used:

1. The linear transformation produces reliable results only when the error propagation can be well approximated by a linear function. If this condition is not achieved the performance of the filter can be extremely poor. At worst, its estimates can diverge altogether.
2. The Jacobian matrices (and Hessian matrices with second order filters) need to exist so that the transformation can be applied. However, there are cases where this is not true. For example, the system might be jump-linear and the parameters can change abruptly [25].
3. In many cases the calculation of Jacobian and Hessian matrices can be a very difficult process, and its also prone to human errors (both derivation and programming). These errors are usually very hard to debug, because it its hard to see which parts of the system produces the errors by looking at the estimates, especially as usually we do not know which kind of performance we should expect.

5.1.6. Unscented Kalman Filter

The *unscented Kalman filter* (UKF) [25] makes use of the *unscented transform* described in section 5.1.6 to give a Gaussian approximation to the filtering solutions of non-linear optimal filtering problems of the form (same as eq. (5-3), but restated here for convenience)

$$\begin{aligned} x_k &= f(x_{k-1}, k-1) + q_{k-1} \\ y_k &= h(x_k, k) + r_k, \end{aligned} \quad (5-13)$$

where $x_k \in \mathfrak{R}^n$ is the state, $y_k \in \mathfrak{R}^m$ is the measurement and $q_{k-1} \sim N(0, Q_{k-1})$ is the Gaussian process noise, and $r_k \sim N(0, R_k)$ is the Gaussian measurement noise.

Unscented Transform

Like Taylor series based approximation presented above the *unscented transform* (UT) [25] can be used for forming a Gaussian approximation to the joint distribution of random variables x and y , which are defined with equations (5-5). In UT we deterministically choose a fixed number of sigma points, which capture the desired moments (at least mean and covariance) of the original distribution of x exactly. After that we propagate the sigma points through the non-linear function g and estimate the moments of the transformed variable from them.

The advantage of UT over the Taylor series based approximation is that UT is better at capturing the higher order moments caused by the non-linear transform, as discussed in [25]. Also the Jacobian and Hessian matrices are not needed, so the estimation procedure is in general easier and less error-prone.

The unscented transform can be used to provide a Gaussian approximation for the joint distribution of variables x and y of the form

$$\begin{pmatrix} x \\ y \end{pmatrix} = N \left(\begin{pmatrix} \mathbf{m} \\ \mu_{\mathbf{U}} \end{pmatrix}, \begin{pmatrix} \mathbf{P} & \mathbf{C}_{\mathbf{U}} \\ \mathbf{C}_{\mathbf{U}}^{\mathbf{T}} & \mathbf{S}_{\mathbf{U}} \end{pmatrix} \right). \quad (5-14)$$

In Eq. 5-14, $\mathbf{m} = E[x]$, with $E[\cdot]$ denoting the expected value. Also, $\mathbf{m} \in \mathbb{R}^n$ and $\mathbf{P} \in \mathbb{R}^{n \times n}$. The transformation is done as follows:

1. Compute the set of $2n + 1$ sigma points from the columns of the matrix $\sqrt{(n + \lambda)\mathbf{P}}$:

$$\begin{aligned} \chi^{(0)} &= \mathbf{m} \\ \chi^{(i)} &= \mathbf{m} + \left[\sqrt{(n + \lambda)\mathbf{P}} \right]_i, \quad i = 1, \dots, n \\ \chi^{(i+n)} &= \mathbf{m} - \left[\sqrt{(n + \lambda)\mathbf{P}} \right]_i, \quad i = n + 1, \dots, 2n \end{aligned} \quad (5-15)$$

and the associated weights:

$$\begin{aligned}
 W_m^{(0)} &= \lambda / (n + \lambda) \\
 W_c^{(0)} &= \lambda / (n + \lambda) + (1 - \alpha^2 + \beta) \\
 W_m^{(i)} &= 1 / \{2(n + \lambda)\}, \quad i = 1, \dots, 2n \\
 W_c^{(i)} &= 1 / \{2(n + \lambda)\}, \quad i = 1, \dots, 2n.
 \end{aligned} \tag{5-16}$$

Parameter λ is a scaling parameter, which is defined as

$$\lambda = \alpha^2(n + \kappa) - n. \tag{5-17}$$

The positive constants α , β and κ are used as parameters of the method.

2. Propagate each of the sigma points through the non-linearity as

$$y^{(i)} = g(x^{(i)}), \quad i = 0, \dots, 2n. \tag{5-18}$$

3. Calculate the mean and covariance estimates for y as

$$\mu_U \approx \sum_{i=0}^{2n} W_m^{(i)} y^{(i)} \tag{5-19}$$

$$S_U \approx \sum_{i=0}^{2n} W_c^{(i)} (y^{(i)} - \mu_U)(y^{(i)} - \mu_U)^T. \tag{5-20}$$

4. Estimate the cross-covariance between x and y as

$$S_U \approx \sum_{i=0}^{2n} W_c^{(i)} (x^{(i)} - m)(y^{(i)} - \mu_U)^T. \tag{5-21}$$

The square root of positive definite matrix \mathbf{P} is defined as $\mathbf{A} = \sqrt{\mathbf{P}}$, where

$$\mathbf{P} = \mathbf{A}\mathbf{A}^T. \tag{5-22}$$

To calculate the matrix \mathbf{A} we can use, for example, lower triangular matrix of the Cholesky factorization.

The Matrix Form of UT

Let $\mathbf{Y} = [m \ \dots \ m]$, $\mathbf{Y} \in \mathbb{R}^{n \times 2n+1}$. The unscented transform described above can be written conveniently in matrix form as follows (see [51] for a proof):

$$\begin{aligned}
 \mathbf{X} &= [\mathbf{Y}] + \sqrt{c} [0_{n \times 1} \quad \sqrt{\mathbf{P}} \quad -\sqrt{\mathbf{P}}] \\
 \mathbf{Y} &= g(\mathbf{X}) \\
 \mu_U &= \mathbf{Y}\mathbf{w}_m \\
 S_U &= \mathbf{Y}\mathbf{W}\mathbf{Y}^T \\
 C_U &= \mathbf{X}\mathbf{W}\mathbf{Y}^T,
 \end{aligned} \tag{5-23}$$

where \mathbf{X} is the matrix of sigma points, function $g(\mathbf{X})$ is applied to each column of the argument matrix separately, $c = \alpha^2(n + \kappa)$, and vector \mathbf{w}_m and matrix \mathbf{W} are defined as follows:

$$\mathbf{w}_m = \left[W_m^{(0)} \quad \dots \quad W_m^{(2n)} \right]^T. \quad (5-24)$$

$$\begin{aligned} \mathbf{W} &= \mathbf{I} - \begin{bmatrix} w_m & \dots & w_m \end{bmatrix} \\ &\times \text{diag}(W_c^{(0)} \quad \dots \quad W_c^{(2n)}) \\ &\times (\mathbf{I} - \begin{bmatrix} w_m & \dots & w_m \end{bmatrix})^T \end{aligned} \quad (5-25)$$

UKF Procedure

Using the matrix form of UT described above the prediction and update steps of the UKF can be computed as follows:

- *Prediction:* Compute the predicted state mean m_k^- and the predicted covariance \mathbf{P}_k^- as

$$\begin{aligned} \mathbf{X}_k^- &= \begin{bmatrix} m_{k-1} & \dots & m_{k-1} \end{bmatrix} + \sqrt{c} \begin{bmatrix} 0 & \sqrt{\mathbf{P}_{k-1}} & -\sqrt{\mathbf{P}_{k-1}} \end{bmatrix} \\ \hat{\mathbf{X}}_k &= f(\mathbf{X}_{k-1}, k-1) \\ \mathbf{m}_k^- &= \hat{\mathbf{X}}_k \mathbf{w}_m \\ \mathbf{P}_k^- &= \hat{\mathbf{X}}_k \mathbf{W} \left[\hat{\mathbf{X}}_k \right]^T + \mathbf{Q}_{k-1}. \end{aligned} \quad (5-26)$$

- *Update:* Compute the predicted mean μ_k and covariance of the measurement \mathbf{S}_k , and the cross-covariance of the state and measurement \mathbf{C}_k :

$$\begin{aligned} \mathbf{X}_k^- &= \begin{bmatrix} m_k^- & \dots & m_k^- \end{bmatrix} + \sqrt{c} \begin{bmatrix} 0 & \sqrt{\mathbf{P}_k^-} & -\sqrt{\mathbf{P}_k^-} \end{bmatrix} \\ \mathbf{Y}_k^- &= \mathbf{h}(\mathbf{X}_k^-, k) \\ \mu_k &= \mathbf{Y}_k^- \mathbf{w}_m \\ \mathbf{S}_k &= \mathbf{Y}_k^- \mathbf{W} \left[\mathbf{Y}_k^- \right]^T + \mathbf{R}_k \\ \mathbf{C}_k &= \mathbf{X}_k^- \mathbf{W} \left[\mathbf{Y}_k^- \right]^T. \end{aligned} \quad (5-27)$$

Then compute the filter gain \mathbf{K}_k and the updated state mean \mathbf{m}_k and covariance \mathbf{P}_k :

$$\begin{aligned} \mathbf{K}_k &= \mathbf{C}_k \mathbf{S}_k^{-1} \\ \mathbf{m}_k &= \mathbf{m}_k^- + \mathbf{K}_k [\mathbf{y}_k - \mu_k] \\ \mathbf{P}_k &= \mathbf{P}_k^- - \mathbf{K}_k \mathbf{S}_k \mathbf{K}_k^T. \end{aligned} \quad (5-28)$$

5.1.7. Multiple Model Systems

In many practical scenarios it is reasonable to assume that the the model of the system can change through time somehow. For example, a fighter airplane, which in normal situation flies with stable flight dynamics, might commence rapid maneuvers when approached by a hostile missile, or a radar can have a different SNR in some regions of space than in others, and so on. Such varying system characteristics are hard to describe with only one certain model, so in estimation one should somehow take into account the possibility that the model of the system might change.

We now consider systems whose current model is one from a discrete set of n models, which are denoted by $M = \{M^1, \dots, M^n\}$. We assume that for each model M^j we have some prior probability $\mu_0^j = \{M_j^0\}$. Also the probabilities of switching from model i to model j in next time step are assumed to be known and denoted by $p_{ij} = P\{M_k^j | M_{k-1}^i\}$. This can be seen as a transition probability matrix of a first order Markov chain characterizing the mode transitions, and hence systems of this type are commonly referred as *Markovian switching systems*. The optimal approach to filtering the states of multiple model system of this type requires running optimal filters for every possible model sequences, that is, for n models n^k optimal filters must be ran to process the k -th measurement. Hence, some kind of approximations are needed in practical applications of multiple model systems.

In this section we describe the Interacting Multiple Model (IMM) filter [5] [14], which is a popular method for estimating systems, whose model changes according to a finite-state, discrete-time Markov chain. IMM filter can also be used in situations where the unknown system model structure or its parameters are estimated from a set of candidate models, and hence it can be also used as a method for model comparison. Initially we start with linear models, and after that we review the EKF and UKF based non-linear extensions to the standard IMM-filter.

Linear Systems

We can now modify the equations of linear systems described in (4-12) to have the form

$$\begin{aligned} x_k &= A_{k-1}^j x_{k-1} + q_{k-1}^j \\ y_k &= H_k^j x_k + r_k^j \end{aligned} \tag{5-29}$$

where now we have denoted by j the model (or mode) acting during the time step $k - 1$. Conditioned on the currently active model we can use the classical Kalman filter (section 5.1.2) for estimating the state of the system on each time step. However, the active model of the system is not usually known, so we must also estimate it.

Interacting Multiple Model (IMM) Filter

IMM-filter [5] is a computationally efficient and in many cases well performing suboptimal estimation algorithm for Markovian switching systems of the type described above. Basically it consists of three major steps: interaction (mixing), filtering and combination. In each time step we obtain the initial conditions for certain model-matched filter by mixing the state estimates produced by all filters from the previous time step under the assumption that this particular model is the right model at current time step. Then we perform standard Kalman filtering for each model, and after that we compute a weighted combination of updated state estimates produced by all the filters yielding a final estimate for the state and covariance of the Gaussian density in that particular time step. The weights are chosen according to the probabilities of the models, which are computed in filtering step of the algorithm. The equations for each step are as follows:

- *Interaction*: The mixing probabilities $\mu_k^{i|j}$ for each model M^i and M^j are calculated as

$$\bar{c}_j = \sum_{i=1}^n p_{ij} \mu_{k-1}^i, \quad (5-30)$$

$$\mu_k^{i|j} = \frac{1}{\bar{c}_j} p_{ij} \mu_{k-1}^i, \quad (5-31)$$

where μ_{k-1}^i is the probability of model M^i in the time step $k-1$ and \bar{c}_j a normalization factor. Now we can compute the mixed inputs (that is, means and covariances) for each filter as

$$\mathbf{m}_{k-1}^{0j} = \sum_{i=1}^n \mu_k^{i|j} \mathbf{m}_{k-1}^i, \quad (5-32)$$

$$\mathbf{P}_{k-1}^{0j} = \sum_{i=1}^n \mu_k^{i|j} \times \left\{ \mathbf{P}_{k-1}^i + [\mathbf{m}_{k-1}^i - \mathbf{m}_{k-1}^{0j}] [\mathbf{m}_{k-1}^i - \mathbf{m}_{k-1}^{0j}]^T \right\}, \quad (5-33)$$

where \mathbf{m}_{k-1}^i and \mathbf{P}_{k-1}^i are the updated mean and covariance for model i at time step $k-1$.

- *Filtering*: Now, for each model M^i the filtering is done as

$$[\mathbf{m}_k^{-,i}, \mathbf{P}_k^{-,i}] = KF_p(\mathbf{m}_{k-1}^{0j}, \mathbf{P}_{k-1}^{0j}, \mathbf{A}_{k-1}^i, \mathbf{Q}_{k-1}^i), \quad (5-34)$$

$$[\mathbf{m}_k^i, \mathbf{P}_k^i] = KF_u(\mathbf{m}_{k-1}^{-,i}, \mathbf{P}_{k-1}^{-,i}, \mathbf{y}_k, \mathbf{H}_k^i, \mathbf{R}_k^i), \quad (5-35)$$

where we have denoted the prediction and update steps (equations ((5-1) and (5-2)) of the standard Kalman filter with $KF_p(\cdot)$ $KF_u(\cdot)$, correspondingly. In addition to mean and covariance we also compute the likelihood of the measurement for each filter as

$$\Lambda_k^i = \mathbf{N}(\mathbf{v}_k^i; 0, \mathbf{S}_k^i), \quad (5-36)$$

where \mathbf{v}_k^i is the measurement residual and \mathbf{S}_k^i is the covariance for model M^i in the KF update step. The probabilities of each model M^i at time step k are calculated as

$$c = \sum_{i=1}^n \Lambda_k^i \bar{c}_i, \quad (5-37)$$

$$\mu_k^i = \frac{1}{c} \Lambda_k^i \bar{c}_i, \quad (5-38)$$

where c is a normalizing factor.

- *Combination*: In the final stage of the algorithm the combined estimate for the state mean and covariance are computed as

$$\mathbf{m}_k = \sum_{i=1}^n \mu_k^i \mathbf{m}_k^i, \quad (5-39)$$

$$\mathbf{P}_k = \sum_{i=1}^n \mu_k^i \times \left\{ \mathbf{P}_k^i [\mathbf{m}_k^i - \mathbf{m}_k] [\mathbf{m}_k^i - \mathbf{m}_k]^T \right\}. \quad (5-40)$$

Non-linear Systems

The non-linear versions of IMM filter reviewed in previous section can be obtained simply by replacing the Kalman filter prediction and update steps (in eqs. (5-34) and (5-35)) by their extended Kalman filter or unscented Kalman filter counterparts, which were reviewed in sections 5.1.5 and 5.1.6. These algorithms are commonly referred as IMM-EKF and IMM-UKF. Naturally, this approach introduces some error to the estimations of an already suboptimal IMM, but it can still provide sufficient accuracy with suitable models.

5.2. Algorithm Description

Using the theoretical concepts presented in the previous section, four localization techniques were implemented in software for simulation purposes. These schemes employ the observed RSSI measurements as the way for obtaining the distance of a mobile target with respect to four reference points. With the distances, the trilateration equations are formed. However, the RSSI measurements contain uncertainty and noise. So, we integrate the Kalman Filter to the localization process through four different approaches, using the concepts illustrated in the previous section.

The analyzed schemes are:

- **IMM-EKF**: EKF based Interacting Multiple Model filter integrated to localization. The procedure is presented in table **5-1**.

- **IMM-UKF**: UKF based Interacting Multiple Model filter integrated to localization. The procedure is presented in table **5-2**.
- **EKF**: Extended Kalman filter integrated to localization. The procedure is presented in table **5-3**.
- **UKF**: Unscented Kalman filter integrated to localization. The procedure is presented in table **5-4**.

Table 5-1: IMM-EKF Localization Algorithm Procedure

- | |
|---|
| <ol style="list-style-type: none"> 1. Measure of RSSI values 2. Calculate distance values from RSSI observed values 3. Calculate the mixing probabilities for each model (eq. (5-30) and (5-31)) 4. Calculate the mixed initial conditions for the states m_{k-1}^{0j} and .. the covariances P_{k-1}^{0j} for the models used (5-32) and (5-33) 5. Perform mode-matched filtering with EKF and calculate the corresponding .. likelihood function (see equations (5-34), (5-35) and (5-36)) 6. Update model probabilities (eq. (5-37) and (5-38)) 7. Extract combined estimate for the state mean $x_k = m_k$ and covariance P_k .. (eq. (5-37) and (5-38)) |
|---|

Table 5-2: IMM-UKF Localization Algorithm Procedure

- | |
|---|
| <ol style="list-style-type: none"> 1. Measure of RSSI values 2. Calculating distance values from RSSI observed values 3. Calculate the mixing probabilities for each model (eq. (5-30) and (5-31)) 4. Calculate the mixed initial conditions for the states m_{k-1}^{0j} and .. the covariances P_{k-1}^{0j} for the models used (5-32) and (5-33) 5. Perform mode-matched filtering with UKF and calculate the corresponding .. likelihood function (see equations (5-34), (5-35) and (5-36)) 6. Update model probabilities (eq. (5-37) and (5-38)) 7. Extract combined estimate for the state mean $x_k = m_k$ and covariance P_k .. (eq. (5-37) and (5-38)) |
|---|

Table 5-3: EKF Localization Algorithm Procedure

1. Measure of RSSI values in dB
2. Calculate distance values from RSSI observed values (eq. (4-31))
3. Initialize values for the states x_0 and the covariance P_0 for the model used
4. Apply EKF procedure using $m_0 = x_0$, P_0 , the distance values and
.. the model parameters (see equations (5-9) and (5-10))
5. Extract the estimated values of the states $x_k = m_k$ at the k instant

Table 5-4: UKF Localization Algorithm Procedure

1. Measure of RSSI values in dB
2. Calculate distance values from RSSI observed values (eq. (4-31))
3. Initialize values for the states x_0 and the covariance P_0 for the model used
4. Calculate *sigma* points and the associated weights (eq. (5-15) and (5-16))
5. Apply UKF procedure using $m_0 = x_0$, P_0 , the distance values and
.. the model parameters (see equations (5-26), (5-27) and (5-28))
6. Extract the estimated values of the states $x_k = m_k$ at the k instant

5.3. Hypothesis Test: ANOVA Analysis

Analysis of variance (ANalysis Of VAriance - ANOVA) is a general method for studying sampled-data relationships [10]. The method analyzes the difference between two or more sample means, achieved by subdividing the total sum of squares. One way ANOVA is the simplest case. The purpose is to test for significant differences between class means, and this is done by analysing the variances. Incidentally, if we are only comparing two different means then the method is the same as the *t- test* for independent samples [11]. The basis of ANOVA is the partitioning of sums of squares into between-class (SS_b) and within-class (SS_w). It enables all classes to be compared with each other simultaneously rather than individually; it assumes that the samples are normally distributed. The one-way analysis is calculated in three steps. First, the sum of squares for all samples, then the within-class cases and, at last, the between-class cases. For each stage the degrees of freedom df are also determined, where df is the number of independent "pieces of information" that go into the estimation of a parameter. These calculations are performed via the Fisher statistic to analyse the null hypothesis. The null hypothesis states that there are no differences between means of different classes, suggesting that the variance of the within-class samples should be identical to that of the between-class samples (resulting in no between-class discrimination capability). It must however be noted that small sample sets will produce random fluctuations due to the assumption of a normal distribution. If d_{ij} is the sample for the i^{th} class and j^{th} data

point then the total sum of squares is defined as:

$$SS_t = \sum_{i=1}^S \sum_{j=1}^D (d_{ij} - GM)^2 \quad (5-41)$$

with degrees of freedom:

$$df_t = (S D) - 1 \quad (5-42)$$

where D is the number of data points (assuming equal numbers of data points in each class) and S is the number of classes and GM is the grand mean:

$$GM = \frac{1}{(S D)} \sum_{i=1}^S \sum_{j=1}^D d_{ij}. \quad (5-43)$$

The second stage determines the sum of squares for the within class case, defined as:

$$SS_w = \sum_{i=1}^S \sum_{j=1}^D (d_{ij} - M_i)^2, \quad (5-44)$$

where M_i is the i^{th} class mean determined by:

$$M_i = \frac{1}{D} \sum_{j=1}^D d_{ij}, \quad (5-45)$$

and the within class df is:

$$df_w = S(D - 1). \quad (5-46)$$

The sum of squares for the between class case is:

$$SS_b = \sum_{i=1}^S D (M_i - GM)^2, \quad (5-47)$$

with the corresponding df of:

$$df_b = S - 1. \quad (5-48)$$

Defining the total degrees of freedom df_t and the total sum of squares SS_t as:

$$df_t = df_b + df_w, \quad (5-49)$$

$$SS_t = SS_b + SS_w. \quad (5-50)$$

Finally if MSS_b is the mean square deviations (or variances) for the between class case, and MSS_w is the reciprocal for the within class case, then:

$$MSS_b = \frac{SS_b}{df_b} \quad ; \quad MSS_w = \frac{SS_w}{df_w}. \quad (5-51)$$

It is now possible to evaluate the null hypothesis using the Fisher statistic, defined as:

$$F = \frac{MSS_b}{MSS_w}. \quad (5-52)$$

If $F \gg 1$, then it is likely that differences between class means exist. These results are then tested for statistical significance or P -value, where the P -value is the probability that a variate would assume a value greater than or equal to the value observed strictly by chance. If the P -value is small (eg. $P < 0,01$ or $P < 1\%$) then this implies that the means differ by more than would be expected by chance alone. By setting a limit on the P -value, (i.e. 1%) a critical F value can be determined. The critical value F_{crit} is determined (via standard lookup tables) through the between-class (df_b) and within-class (df_w) df values. Values of F greater than the critical value denote the rejection of the null hypothesis, which prompts further investigation into the nature of the differences of the class means. In this way ANOVA can be used to prune a list of features [55].

Summary

This section presented the application of four different Kalman-based schemes to the localization problem in Ad Hoc wireless sensor networks. All the proposed schemes employ the multilateration technique as the basis for the unknown position determination. The schemes are distinguished by the Kalman-scheme implemented, and by the use of a single mobility model or a set of interacting mobility models. The proposed schemes were:

- **EKF**: Employs Extended Kalman Filter and a single movement model.
- **UKF**: Employs Unscented Kalman Filter and a single movement model.
- **IMM-EKF**: Employs Extended Kalman Filter and the interacting multiple model technique for the target movement representation.
- **IMM-EKF**: Employs Unscented Kalman Filter and the interacting multiple model technique for the target movement representation.

THIS section enden with the description of the ANOVA analysis, an statistical tool for studying the relationships between different sets of sampled data.

6 Simulation results

In this chapter, the simulation results of the different localization algorithms are shown and discussed. Two examples are exhibited with the objective to demonstrate the capability of the algorithms of tracking a target undergoing different trajectories.

6.1. Wiener Process Models Performance Comparison

In this section, the performance of the P , PV , and PVA models with Extended Kalman Filter for trajectory tracking will be compared. The benefit of one over the other depends upon the characteristics of the motion of the object. A system modeled using just P will work when the position is mostly constant and the velocity can be treated as noise. In the case of a PV model, it will tend to work better when velocity is mostly constant, and the acceleration can be treated as noise. PVA on the other hand works better when the acceleration is mostly constant [?].

6.1.1. Simulation Description

At first, we will explore the performance of the P , PV , and PVA models with an Extended Kalman Filter. We simulate the two dimensional motion of an object by collecting distance measurements of the mobile node while physically moving it in a network composed of 8 wireless sensors (see figure **6-1**). The range measurement employed is Received Signal Strength Indicator (RSSI); however, as was discussed in previous sections, the use of RSSI introduces noise. The distance measurements returned by the sensors fluctuate often and these measurements are the inputs to the Kalman filters.

$$RSSI = \kappa_j - 10\gamma \log_{10}(d_{ij,k}) + v_{i,k}, \tag{6-1}$$

To investigate the positioning capabilities of the implemented RSSI based algorithms the simulation was set up as reported below. We chose the parameters in order to fit the real node features and the characteristics of the environment reported in [36]. The distance measurements were obtained from eq. (6-1) (see section 4.4 for further explanation). All the algorithms have been tested in 100 scenarios, with randomly created noises, and equal for all algorithms; the results obtained can therefore be interpreted statistically. We have repeated

Table 6-1: Simulation parameters for Wiener Process Models Performance Comparison

T	Sampling time [s]	0,08s
n	Number of samples	50
κ_j	Transmission power [dB]	90
γ	Path loss index	2,1060
n_{sim}	Number of Monte Carlo runs	100
σ_v	Covariance of the gaussian noise v [dB]	2,0873

the experiments with the same simulation values, but modifying the process noise parameter q over the Extended Kalman Filter procedure, to gauge its influence on the overall estimation with the different models.

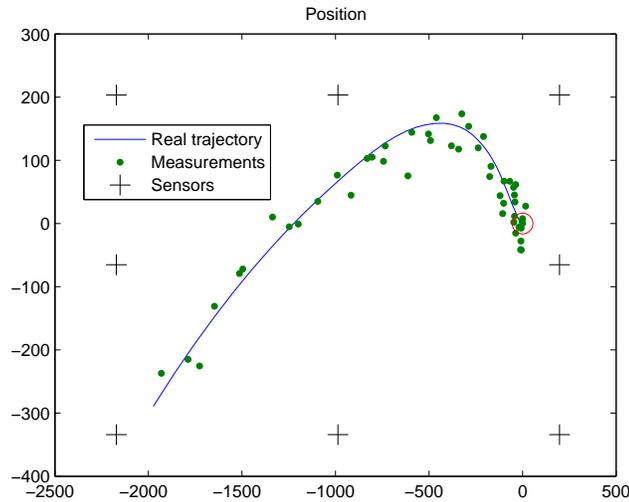


Figure 6-1: Simulation testbed for the comparison of the trajectory estimation for P, PV, and PVA models with an Extended Kalman Filter. The green dots denote a set of the noisy position observations randomly generated.

The sensors or reference points were placed at fixed coordinates (denoted by +) of the grid as shown in **6-1**. For each of the 50 tested position (see the green dots), distances measurements were collected from the eight beacons. The measurement noise for each of the sensors was assumed to be independent of the others. The measurement noise was obtained by taking the average difference between the actual and estimated distances from the beacons to the nodes. The process noise matrix Q is more difficult to obtain. Approximate behavior such as the standard deviation of the position for the P model about the estimated movement of the object can be used. Of course, the standard deviation of velocity would be applied

for the *PV* model, and the standard deviation of the acceleration for the *PVA* model. Determination of the correct process noise parameters are key to accurate localization. Here, we use $Q = \text{diag}\{q\}$, with q taking values from the vector $q = [50 \ 5 \ 0,5 \ 0,05]^T$.

6.1.2. Discussion of Results

We have implemented EKF P, PV, and PVA tracking algorithms and performed experiments extensively. Several metrics were used to compare the performances of the Kalman filters. The first is the average distance error (DE) in localization per estimate, as defined by equation (6-2). Another metric that was used is the Root Mean Square Error (RMSE). The benefit of the RMSE given in equation (6-3) is that the error in localization of the X and Y coordinates is available. The X and Y RMSE values can be combined through equation (6-4) resulting in the Net RMSE that describes the net error. An interesting characteristic of the RMSE is that it is biased towards large errors. A large error make a larger contribution in RMSE than in average distance error.

$$\text{DE} = \frac{\Sigma \sqrt{(x_{\text{actual}} - x_{\text{est}})^2 + (y_{\text{actual}} - y_{\text{est}})^2}}{\text{number of estimates}} \quad (6-2)$$

$$\text{RMSE} = \sqrt{\frac{\Sigma(\text{actual} - \text{estimated})^2}{\text{number of estimates}}} \quad (6-3)$$

$$\text{Net RMSE} = \sqrt{x_{\text{RMSE}}^2 + y_{\text{RMSE}}^2} \quad (6-4)$$

We ran one hundred simulations of the position per value of the noise process q . In table **6-2**, the average error metrics for each case are presented. As table **6-2** indicates, the PV model has the least distance error per estimate and the least Net_{RMSE} and hence the best localization performance in every case, except when the value of the noise process is $q = 0,05$. This is followed closely by the PVA model, and finally the P model. Figure **6-2** shows the best estimation reached by each model for the different q values evaluated. Figure **6-3** plots the average Distance Error of the estimation with each model for the different values of q . Over the performance of the models, we can highlight the following comments:

- The value of the process noise parameter q has a deep effect on the estimation capabilities of each model. When the q value is high, meaning that the uncertainty over the process model is high, all three estimation models have a similar performance. However, as the q value decreases, the models with more complete information prevail. The PV and PVA models, with more rigid kinematic equations, tend to perform better. This fact is corroborated when the PVA model (more descriptive) reaches the lowest error values when the noise process parameter has a minimum value of $q = 0,05$

- The P model estimation decays significantly when the values of the noise parameter decrease; in fact, the P model only works well when $q = 50$. This is due to the assumption that velocity and acceleration are just random noise; when the uncertainty in the trajectory model decrease (lower q), the ability of the P model to keep up with the reference trajectory is completely reduced.

There are two cases when seems that there is no difference between the estimation with the different models. For these specific cases, we performed an ANOVA analysis over the average DE of the estimations. We choose as the null hypothesis the statement that there are no significant statistical differences between the estimations, and we use a critical value or $F_{crit} = 0,05$. Remember, if the P -value of the ANOVA is lower than F_{crit} , we reject the null hypothesis. Conversely, if P -value $> F_{crit}$, then the null hypothesis is accepted. See the results in the paragraphs below.

Table 6-2: Average of the error metrics for the trajectory estimation with P , PV and PVA models for different values of q

<i>Model</i>	<i>DE</i>	<i>x_{RMSE}</i>	<i>y_{RMSE}</i>	<i>Net_{RMSE}</i>	<i>q</i>
P	33.6596	28.4742	25.5553	38.3858	50
PV	32.4624	26.9093	25.3647	37.1278	
PVA	34.2068	27.3339	27.8196	39.1788	
P	38.3033	33.9419	32.6505	47.2804	5
PV	27.2986	23.4787	20.0342	30.9742	
PVA	29.9712	25.1152	22.7514	34	
P	96.3363	106.2014	79.9403	132.9739	0.5
PV	26.3663	20.3114	21.6824	29.9136	
PVA	27.0517	22.2004	20.942	30.6959	
P	226.2218	306.2388	93.8838	320.331	0.05
PV	46.6378	34.3279	45.2519	56.9912	
PVA	25.0649	20.2785	19.5381	28.4024	

- **Case 1 - Estimation when $q = 50$:** Figure 6-4 depicts the boxplots and the P -value of the ANOVA for the different set of estimations with $q = 50$. Even if the estimation with the P, PV, and PVA models look very similar from figures 6-2 and 6-3, the P -value indicates that there is a significant statistical difference among the average Distance Error of the estimations. However, the comparison of P and PVA data showed no difference between them, with a P -value = 0,25994 $>$ $F_{crit} = 0,05$. These results favor the PV estimation as the best possible result obtained whit $q = 50$.
- **Case 2 - Estimation with PV and PVA models when $q = 0,5$:** Figure 6-5 depicts the boxplots and the P -value of the ANOVA for only PV and PVA set of estimations with $q = 0,5$. As we can see from figures 6-2 and 6-3, the P model performance is lower than the other two, so we let this data out of the analysys. In this case, the ANOVA analysis threw a P -value = 0,10608 $>$ $F_{crit} = 0,05$. Therefore, we accept the null hypothesis, and we conclude that there is no significant statistical difference between the average of Distance Error estimation for PV and PVA models with $q = 0,5$

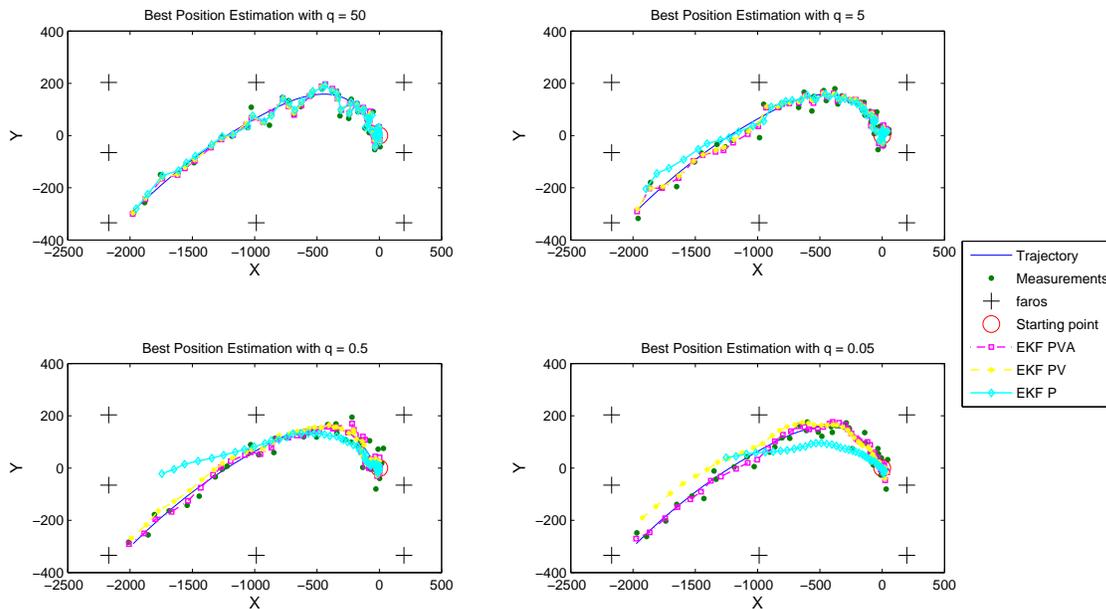


Figure 6-2: Best trajectory estimation with each model for different values of the q parameter

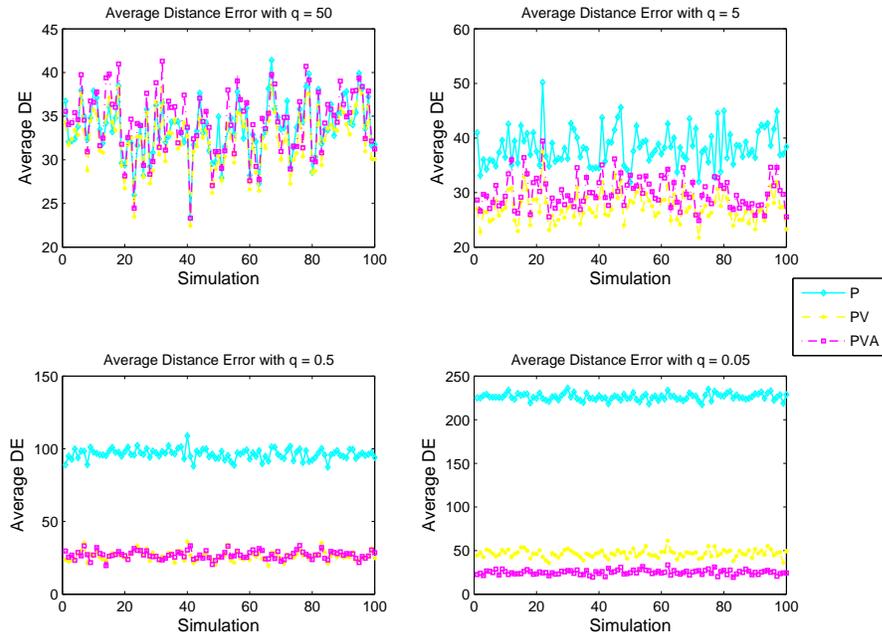


Figure 6-3: Average Distance Error with the different values of q

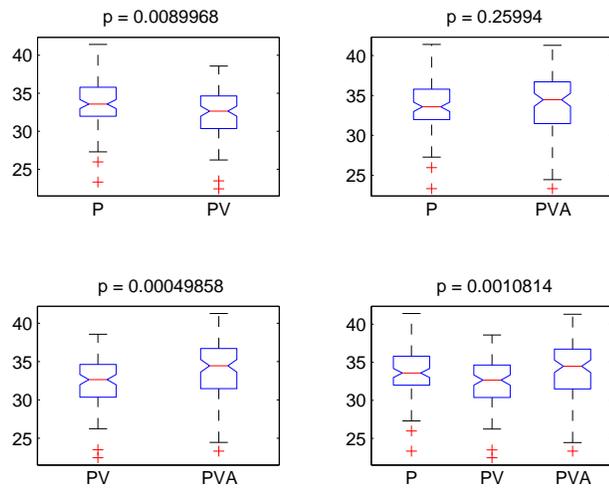


Figure 6-4: Boxplot and p -value of ANOVA analysis for the Distance Error data of the different estimations with $q = 50$

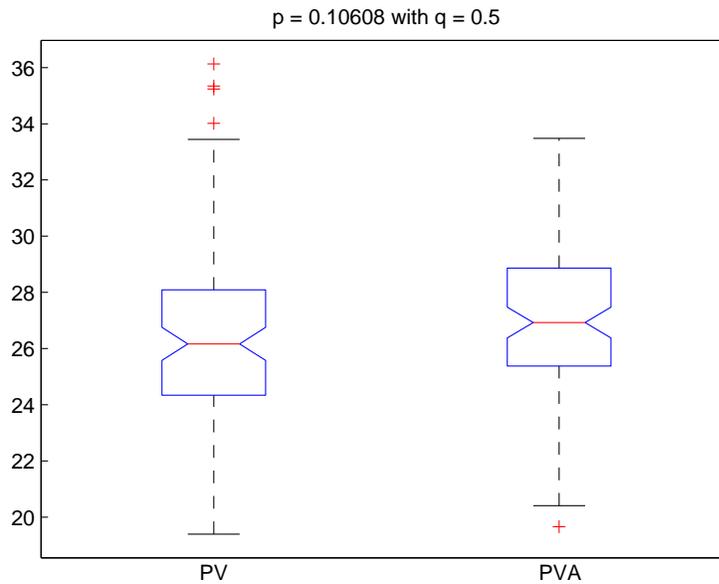


Figure 6-5: Boxplot and p -value of ANOVA analysis for the Distance Error data of PV and PVA estimations with $q = 0,5$

As a summary, we can conclude that the performance of the three Wiener Process models evaluated was very similar when the noise process was high; i.e., there is a high uncertainty over the motion target model. However, as the noise process decreased, the performance of the different estimation got better for the models including more kinematic parameters (P and PVA), as this kind of models have a more detailed descriptive level.

6.2. Target Tracking

Next we review a classical filtering application (see, e.g., [?]), in which we track a moving object with sensors, which measure only the bearings (distances) of the object with respect positions of the sensors. There is a one moving target in the scene and four sensors for tracking it. Solving this problem is important, because often more general multiple target tracking problems can be partitioned into sub-problems, in which single targets are tracked separately at a time [51].

6.2.1. Simulation Description

Similarly to the scenario presented in section 6.1, we consider one mobile target with four reference nodes. Respective pilot signal strengths are used between the mobile nodes and

reference nodes. Two sample mobile trajectories were generated using the Coordinated Turn Model for the mobile target.

The four Kalman-based algorithms described in the previous chapter are used for locating the mobile target. The target and the reference nodes transmit RSSI signals, and every node can receive the signals of the others (a fully communicated one-hop network). The distance measurements were obtained from the RSSI data and the position of the object was extracted directly with the additive noise, in the same way as explained in section 6.1. Over 100 Monte Carlo simulations were performed for each algorithm and trajectory, with the RSSI values randomly generated in each experiment.

In the estimation we use the following models:

1. Standard Wiener process velocity model with process noise variance $q_1 = 0,05$, whose purpose is to model the relatively slow turns.
2. A combination of Wiener process velocity model and a coordinated turn model described above. The variance of the process noise for the velocity model is set to $q_2 = 0,01$ and for the turning rate parameter in the turning model to $q_\omega = 0,15$. The estimation is now done with both the EKF and UKF based IMM filters as the turning model is non-linear. In both cases the model transition probability matrix is set to

$$\Phi = \begin{bmatrix} 0,9 & 0,1 \\ 0,1 & 0,9 \end{bmatrix}, \quad (6-5)$$

and the prior model probabilities are

$$\mu_0 = [0,9 \quad 0,1]. \quad (6-6)$$

The simulation parameters are described in table **6-3**. The position of the object is estimated with the following methods:

- **EKF**: Extended Kalman filter using the Wiener process velocity model.
- **UKF**: Unscented Kalman filter using the same model as the EKF.
- **IMM-EKF**: EKF based IMM filter using a combination of Wiener process velocity model and a coordinated turn model.
- **IMM-UKF**: UKF based IMM filter using the same models as IMM-EKF.

The combined position root-mean-square error (RMSE) shown in the equation 6-7 is used to assess the closeness of the estimated trajectory \hat{x}, \hat{y} to to the actual trajectory x, y averaged over the number of data points n .

$$RMSE = \sqrt{\frac{1}{n} \sum_{i=1}^n [(\hat{x}_i - x_i)^2 + (\hat{y}_i - y_i)^2]}. \quad (6-7)$$

Table 6-3: Simulation parameters for Target Tracking

T	Sampling time [s]	0,1s
n	Number of samples	200
κ_j	Transmission power [dB]	90
γ	Path loss index	2,1060
$nsim$	Number of Monte Carlo runs	100
σ_v	Covariance of the gaussian noise v [dB]	2,0873

6.2.2. Results for trajectory 1

Trajectory 1 is generated from a combination between PV Wiener Process Model and a Coordinated Turn Model as described in table 6-4. The trajectory is depicted in figure 6-6. We ran 100 Monte Carlo simulations for this trajectory. A sample result of one estimations is plotted in Figure 6-7. The estimates seem to be very similar for the four Kalman-based algorithms.

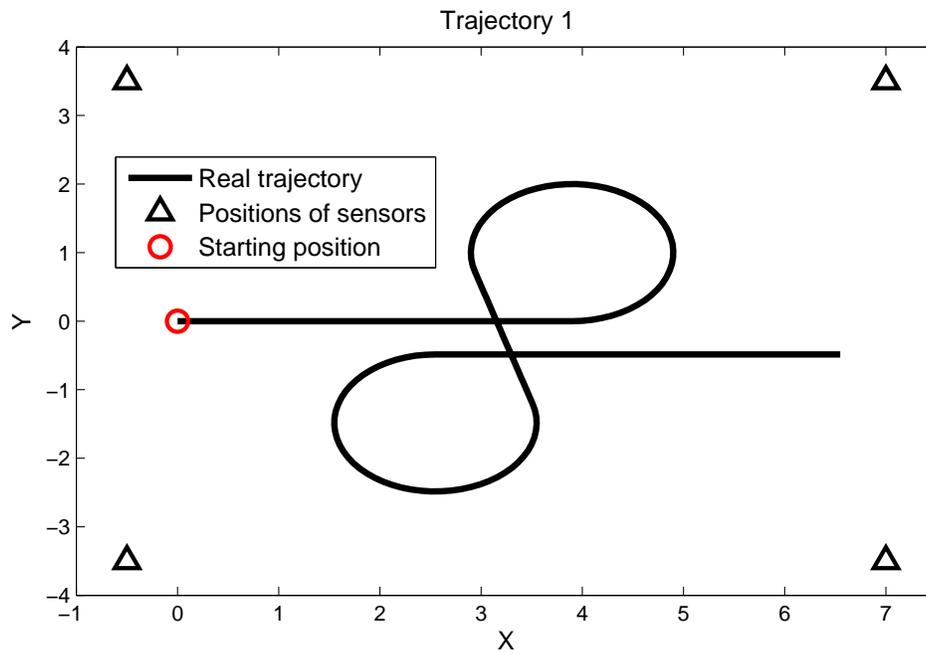
In table 6-5 we have listed the average mean square errors of position estimates over 100 Monte Carlo runs. The histogram of average RMSE of position estimation for each simulation is plotted in figure 6-8. Also, The average RMSE of velocity estimation for each simulation is shown in figure 6-9. It can be observed that the estimates of EKF and UKF are identical in practice. The difference between IMM-UKF and IMM-EKF is trending in the favor of IMM-UKF. To corroborate this observations, we perform an ANOVA analysis like is going to be shown in the next subsection.

Table 6-4: Description of Trajectory 1

1.	Object starts from origin with velocity $(\dot{x}, \dot{y}) = (1, 0)$.
2.	At 4s object starts to turn left with rate $\omega = 1$.
3.	At 9s object stops turning and moves straight for 2 seconds with a constant total velocity of one.
4.	At 11s objects starts to turn right with rate $\omega = -1$.
5.	At 16s object stops turning and moves straight for 4 seconds with the same velocity.

Table 6-5: Average of RMSE values for Trajectory 1

Algorithm	Position RMSE
EKF	0,0587
UKF	0,0586
IMM-EKF	0,0483
IMM-UKF	0,0401

**Figure 6-6:** Representation in 2D of Target trajectory 1.

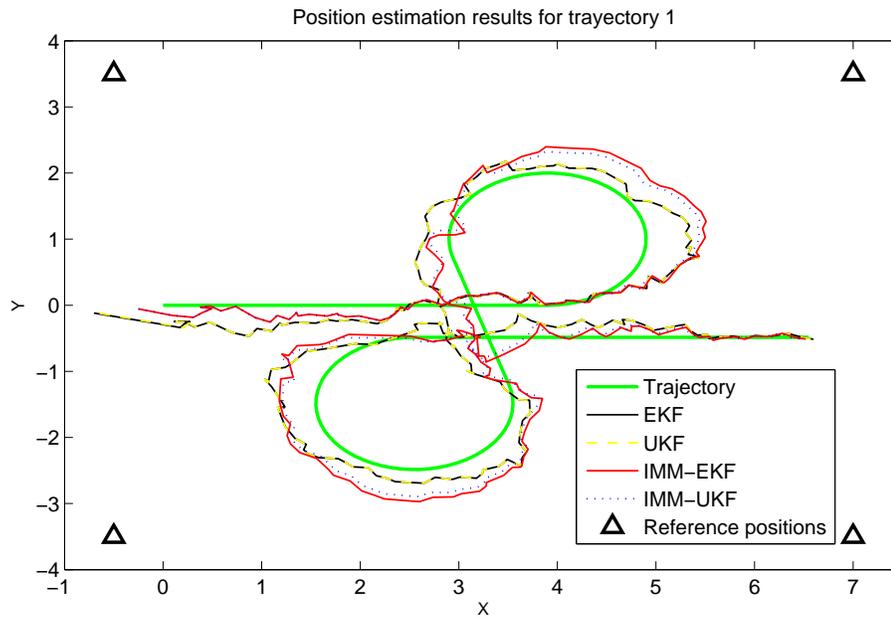


Figure 6-7: A sample of position estimation results for trajectory 1

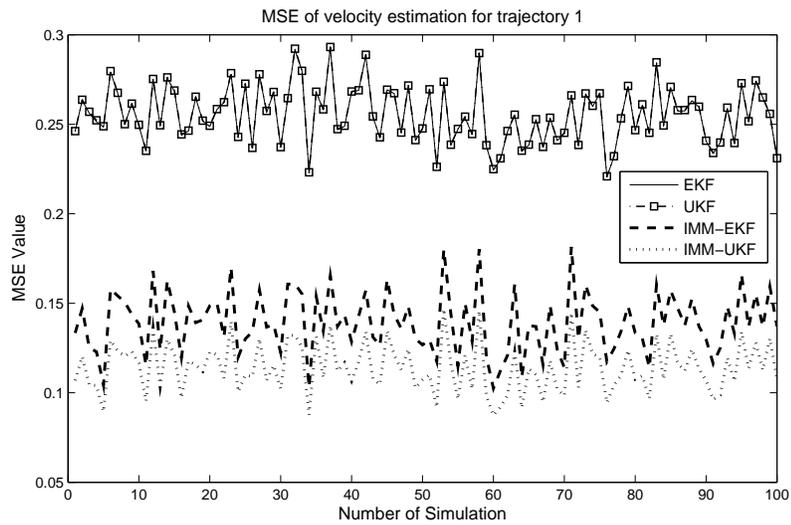


Figure 6-9: MSE of velocity estimation for trajectory 1

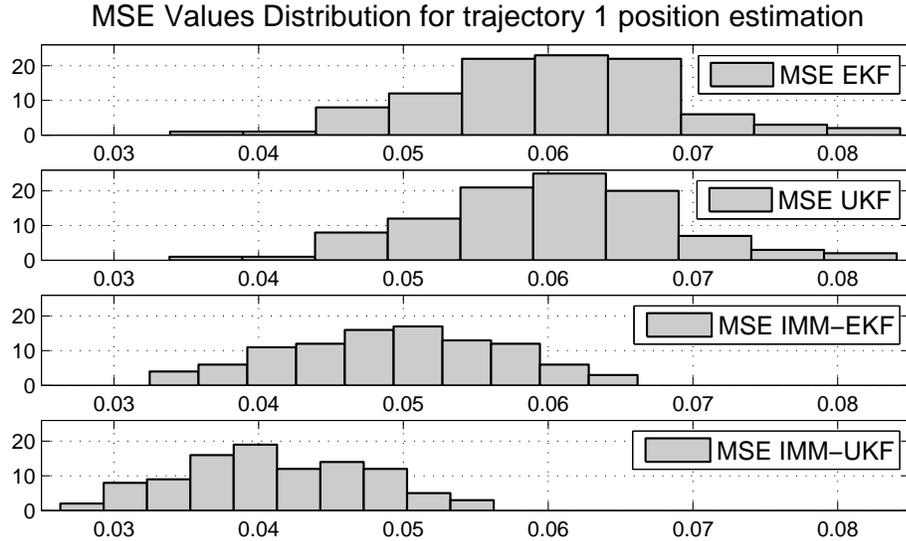


Figure 6-8: MSE of position estimation for trajectory 1

ANOVA analysis for trajectory 1

From figures 6-7 and 6-8, we observe that the position estimation through EKF and UKF is almost identical. Also, the average error for the IMM-EKF and IMM-UKF seems to be very close. For these specific cases, we performed an ANOVA analysis over the average DE of the estimations. As we did in the section 6.1, we choose as the null hypothesis the affirmation that there are no significant statistical differences between the estimations, and we use a critical value or $F_{crit} = 0,05$. Remember, if the P -value of the ANOVA is lower than F_{crit} , we reject the null hypothesis. Conversely, if $P\text{-value} > F_{crit}$, the null hypothesis is accepted. See the results in the paragraphs below.

Figure 6-10 depicts the boxplots and the P -value of the ANOVA for the different set of estimations with the four different filters for trajectory 1. Even if the estimations between subsets look very similar, the P -value for the different combinations of datasets indicates that there is a significant statistical difference among the average error of the estimations. However, the comparison of EKF and UKF data showed no difference between them, with a $P\text{-value} = 0,96013 > F_{crit} = 0,05$. These results favor the IMM-UKF estimation as the best possible result obtained for trajectory 1.

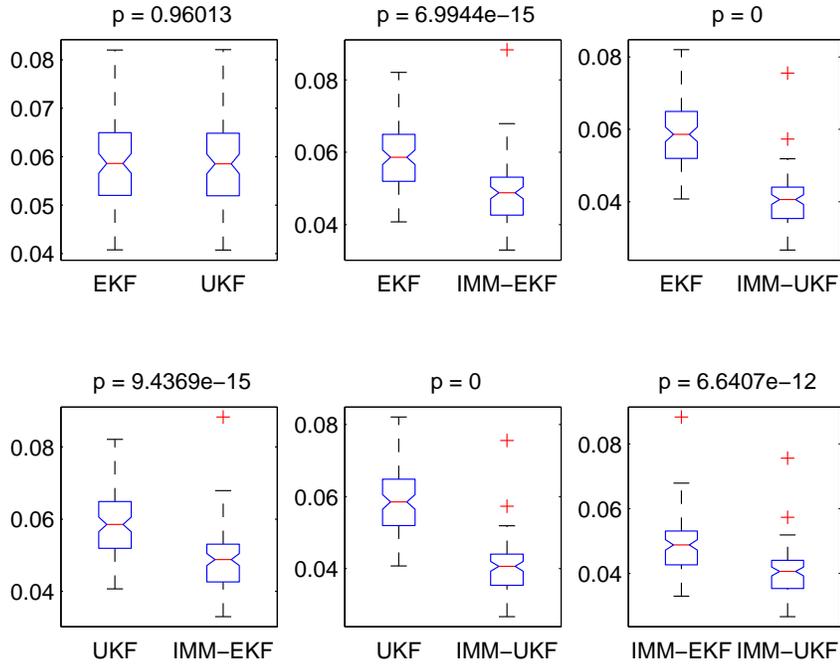


Figure 6-10: Boxplots of MSE distributions for ANOVA analysis of position estimation for trajectory 1

6.2.3. Results for trajectory 2

Trajectory 2 is also generated from a combination between PV Wiener Process Model and a Coordinated Turn Model as described in table 6-6. The trajectory is depicted in figure 6-11. We ran 100 Monte Carlo simulations for this trajectory. A sample result of one estimations is plotted in Figure 6-12. The estimates are very similar for the four Kalman-based algorithms, even more than the previous case. In fact, the results of each method are barely discriminated one from the others.

Table 6-6: Description of Trajectory 2

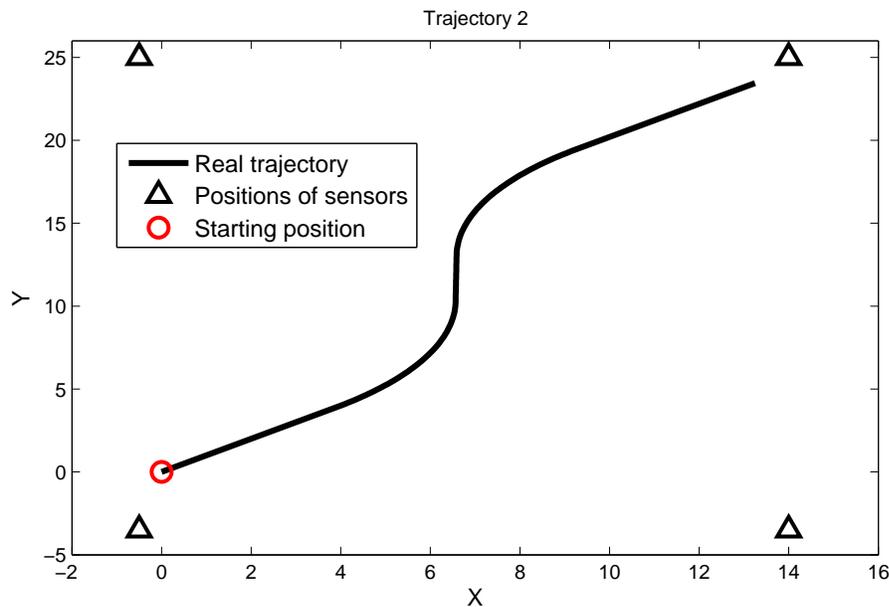
1.	Object starts from origin with velocity $(\dot{x}, \dot{y}) = (1, 1)$.
2.	At 4s object starts to turn left with rate $\omega = 0,1555$ and velocity $(\dot{x}, \dot{y}) = (-1, -1)$.
3.	At 9s object stops turning and moves straight for 2 seconds with a constant total velocity of one.
4.	At 11s objects starts to turn right with rate $\omega = -0,1555$.
5.	At 16s object stops turning and moves straight for 4 seconds with the same velocity.

Table 6-7: Average of RMSE values for Trajectory 2

Algorithm	Position RMSE
EKF	0,0339
UKF	0,0339
IMM-EKF	0,0259
IMM-UKF	0,0249

In table **6-7** we have listed the average mean square errors of position estimates over 100 Monte Carlo runs. The histogram of average RMSE of position estimation for each simulation is plotted in figure **6-8**. Also, the average RMSE of velocity estimation for each simulation is shown in figure **6-9**.

For the states representing the velocity over each axis, the IMM based methods hold a clearly advantage over EKF and UKF. This could be expected, as the IMM techniques rely on the interaction of more mobility models, and thus can extract more information. As in the previous section, it can be observed that the estimates of EKF and UKF are identical in practice. However, there is no appreciable difference between IMM-UKF and IMM-EKF, although the performance of the latter two methods is better than the former two. Again, to corroborate this observations, we perform the ANOVA analysis that is being shown shown in the next subsection.

**Figure 6-11:** Representation in 2D of Target trajectory 2.

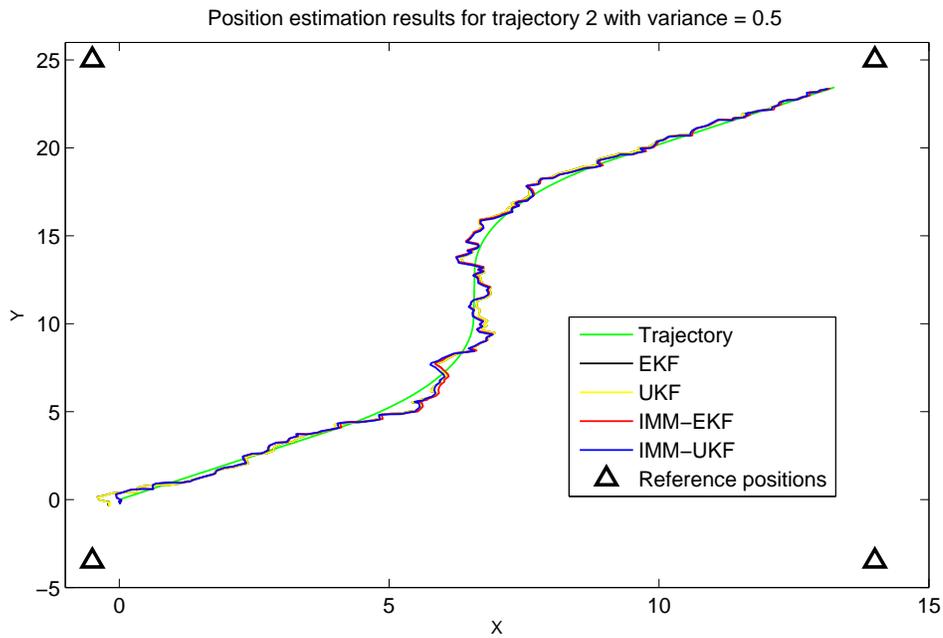


Figure 6-12: A sample of position estimation results for trajectory 2

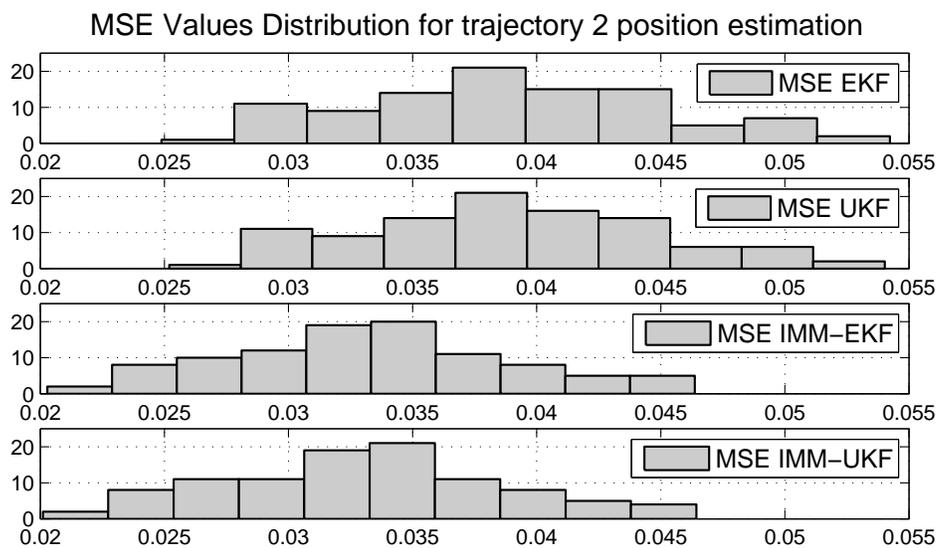


Figure 6-13: MSE of position estimation for trajectory 2

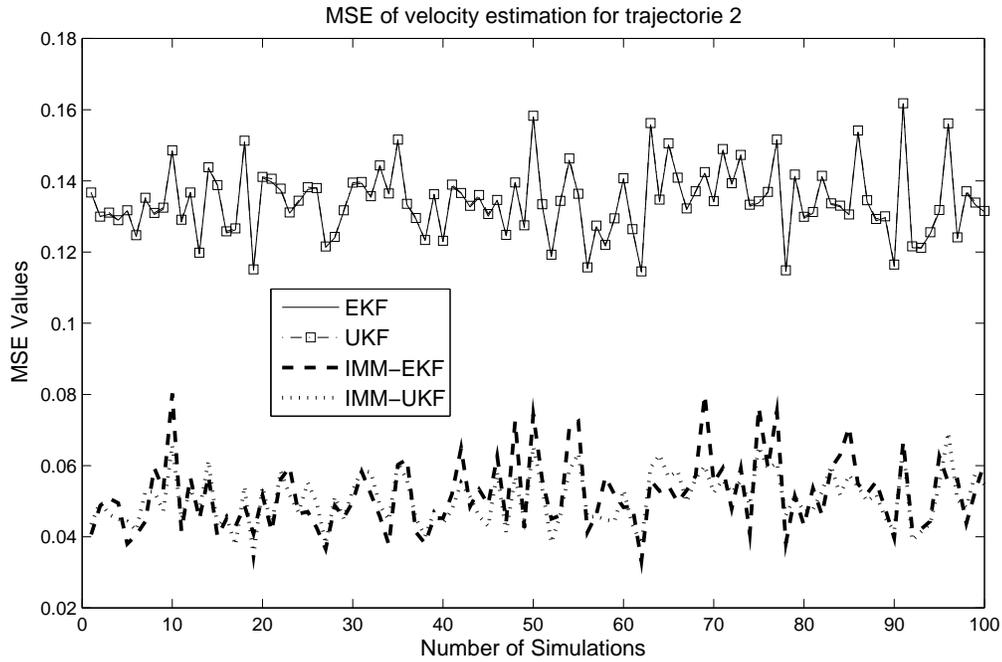


Figure 6-14: MSE of velocity estimation for trajectory 2

ANOVA analysis for trajectory 2

From figures 6-12 and 6-13, we observe that the position estimation trough EKF and UKF is almost identical. Also, the average error for the IMM-EKF and IMM-UKF seems to be equal. For these specific cases, we performed an ANOVA analysis over the average error of the estimations. Again, we choose as the null hypothesis the affirmation that there are no significant statistical differences between the estimations, and we use a critical value or $F_{crit} = 0,05$.

Figure 6-15 depicts the boxplots and the P -value of the ANOVA for the different set of estimations with the four different filters for trajectory 2. Like the previous case, a P -value of 0,9606 shows that there is no significant statistical difference between the estimation with EKF and UKF for this case. Besides this, the null hypothesis is accepted also for the comparison of IMM-EKF and IMM-UKF, with a P -value = 0,82111, meaning the estimation results for these two models is statistically identical. So, there is no clear-cut better method for estimation of trajectory 2: both IMM-based models perform better than the others.

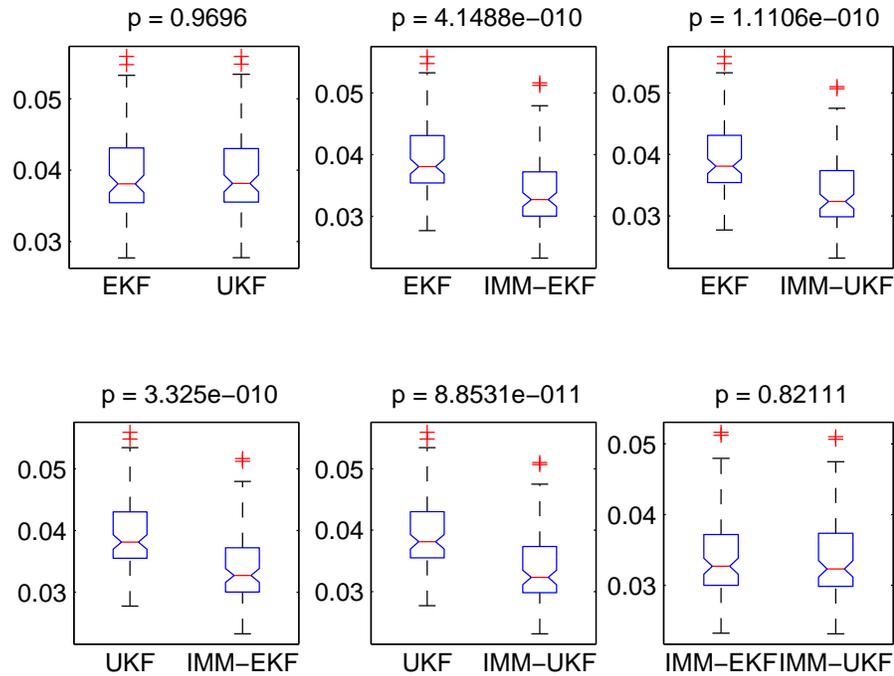


Figure 6-15: Boxplots of MSE distributions for ANOVA analysis of position estimation for trajectory 2

Summary

It has been shown that the four Kalman-based schemes present similar estimation results for the two different trajectories studied. A case could be made for either method, but the choice of the best one depends on the designer's criteria. The IMM-UKF yields the best results, and the Multiple Model algorithms performed better than those with one-model. Besides this, it should be noted that the performance of each tested method could be tuned by optimizing their parameters (e.g. variance of process noise of dynamic models, values of model transition matrix in IMM, etc.) more carefully, so the performance differences could change radically. Still, it is clear that the Kalman filter does actually work also with (at least some) non-linear dynamic and measurement models, and should be considered as a standard location estimation method for multiple model systems. Also, one should prefer IMM-UKF over IMM-EKF as the performance is (at least in these examples) better, although its computation could be higher in some cases.

7 Conclusions and future work

The aim of this thesis was to present an overview of a collection of tools used in the problem of target tracking. Under this thesis scope, several algorithms for target tracking were studied. This chapter will thus emphasize the main results drawn from the algorithms studied, bearing in mind that the validity of these conclusions may be limited to the scenarios simulated.

7.1. Conclusions

- The Kalman filter is capable of localizing using noisy distance measurements in sensor networks. The PV Wiener Process model with Extended Kalman filter was found to have the best performance over the other Wiener Process model in the studied examples. However, depending on the motion of the tracked object the PV or PVA model could be better.
- The process noise parameters allow the Kalman filter to project the position of the object in the next time instant. If the proper process noise parameters are not used, the performance of the Kalman filters will be severely affected.
- The Kalman-based localization algorithms all revealed capable trackers under target motion uncertainty. The first couple of Kalman-based algorithms (EKF and UKF) are the most limited, since they assume the target motion is governed by a constant pattern, i.e., the algorithms only use an unique model to estimate the mobile target trajectory . The other two schemes (IMM-EKF and IMM-UKF) have more complex concepts, even though they are not computationally more complex (when using a low number of models). Their tracking capability appears to be better, since they assume the target can have multiple maneuver motions.
- Regarding the performance comparison shown in chapter 6, this results represent only an example since the algorithms may have different responses for different target motions. In the simulations performed in this thesis, there is no sensible difference between EKF and UKF. Given the availability of more resources (because they had multiple mobility models), the IMM-based algorithms got better results than EKF and UKF. This suggests that there is a sensible advantage for the use of this Multiple Model techniques in localization schemes for Ad-Hoc sensor networks over the one-model based techniques.

- Kalman filtering is an interesting and powerful method to optimally and dynamically process data hampered by white gaussian noise. Filtering also allows one to interpret measurements coherently with a mathematical model of the observed phenomenon. However, as a downside, the performance of the Kalman Filter depends on tuning parameters and on the model used to describe the process of interest. It is possible using Kalman filtering to realize tracking applications, achieving a good performance after a specific delay to achieve the best filter state.

7.2. Future Work

The additional research that can be performed in this area is vast. There are several additional topics to extend this thesis work, they are:

1. **Test further and study even in more depth the algorithms in simulation environment:** more simulations could be realized, specially for the IMM Kalman-based algorithms, but with the intervention of other kind of ranging signals different than RSSI. Also further algorithm development could be done in order to study its performance in the presence of the phenoms affecting the RSSI signals, just like interferences, fading and multipath effects.

Regarding this performance improvement, it would be interesting to analyze possible mathematical designs of the model probability thresholds, characteristic of the IMM algorithms, and study whether they should be static or adaptive, for instances, if there is a high uncertainty on the state estimate, the thresholds can be relaxed.

2. **Analyze the algorithms performance in a real life implementation:** Since all the discussed algorithms, are studied in environments that try to simulate a real life situation, it would be valuable to analyze their performances in a real life application. It would be interesting to implement the algorithms using real sensor measurements and real targets, and test their capability and robustness.
3. **Study alternative algorithms:** there are a lot to choose from, more or less documented and more or less studied. An interesting subject with plenty of room for exploration is the study of more alternative techniques such as the Monte Carlo techniques known as Particle Filtering [20], which have drawn the attention of the tracking community.

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