

In situ calibration of low-cost instrumentation for the measurement of ambient quantities: evaluation methodology of the algorithms and diagnosis of drifts

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Étalonnage in situ de l'instrumentation bas coût pour la mesure de grandeurs ambiantes : méthode d'évaluation des algorithmes et diagnostic des dérives

Thèse de doctorat de l'Institut Polytechnique de Paris préparée à École Polytechnique

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To facilitate the understanding of this part by the people it concerns, French is used here.

Pour commencer ces remerciements, je souhaite tout d'abord m'adresser au jury qui a évalué ces travaux. Merci à Gilles Roussel d'avoir accepté d'examiner cette thèse et d'en présider le jury. Je remercie également Jean-Luc Bertrand-Krajewski et Romain Rouvoy pour avoir accepté de rapporter mes travaux de thèse dans le contexte sanitaire actuel qui bouleverse les activités des enseignants-chercheurs. Merci pour vos rapports et les questions qui y ont figuré, m'aidant à préparer la soutenance. Merci également à Nathalie Redon d'avoir examiné mes travaux et enfin à Éric Peirano et Matthieu Puigt d'avoir pris part au jury en tant qu'invités. J'ai particulièrement apprécié nos échanges, notamment à travers les questions que vous avez chacun pu avoir sur mes travaux.

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In various fields going from agriculture to public health, ambient quantities have to be monitored in indoors or outdoors areas. For example, temperature, air pollutants, water pollutants, noise and so on have to be tracked. To better understand these various phenomena, an increase of the density of measuring instruments is currently necessary. For instance, this would help to analyse the effective exposure of people to nuisances such as air pollutants.

The massive deployment of sensors in the environment is made possible by the decreasing costs of measuring systems, mainly using sensitive elements based on micro or nano technologies. The drawback of this type of instrumentation is a low quality of measurement, consequently lowering the confidence in produced data and/or a drastic increase of the instrumentation costs due to necessary recalibration procedures or periodical replacement of sensors.

There are multiple algorithms in the literature offering the possibility to perform the calibration of measuring instruments while leaving them deployed in the field, called *in situ* calibration techniques.

The objective of this thesis is to contribute to the research effort on the improvement of data quality for low-cost measuring instruments through their *in situ* calibration.

In particular, we aim at 1) facilitating the identification of existing *in situ* calibration strategies applicable to a sensor network depending on its properties and the characteristics of its instruments; 2) helping to choose the most suitable algorithm depending on the sensor network and its context of deployment; 3) improving the efficiency of *in situ* calibration strategies through the diagnosis of instruments that have drifted in a sensor network.

Three main contributions are made in this work. First, a unified terminology is proposed to classify the existing works on *in situ* calibration. The review carried out based on this taxonomy showed there are numerous contributions on the subject, covering a wide variety of cases. Nevertheless, the classification of the existing works in terms of performances was difficult as there is no reference case study for the evaluation of these algorithms.

Therefore in a second step, a framework for the simulation of sensors networks is introduced. It is aimed at evaluating *in situ* calibration algorithms. A detailed case study is provided across the evaluation of *in situ* calibration algorithms for blind static sensor networks. An analysis of the influence of the parameters and of the metrics used to derive the results is also carried out. As the results are case specific, and as most of the algorithms recalibrate instruments without evaluating first if they actually need it, an identification tool enabling to determine the instruments that are actually faulty in terms of drift would be valuable.

Consequently, the third contribution of this thesis is a diagnosis algorithm targeting drift faults in sensor networks without making any assumption on the kind of sensor network at stake. Based on the concept of rendez-vous, the algorithm allows to identify faulty instruments as long as one instrument at least can be assumed as non-faulty in the sensor network. Across the investigation of the results of a case study, we propose several means to reduce false results and guidelines to adjust the parameters of the algorithm. Finally, we show that the proposed diagnosis approach, combined with a simple calibration technique, enables to improve the quality of the measurement results. Thus, the diagnosis algorithm opens new perspectives on *in situ* calibration.

Dans de nombreux domaines allant de l'agriculture à la santé publique, des grandeurs ambiantes doivent être suivies dans des espaces intérieurs ou extérieurs. On peut s'intéresser par exemple à la température, aux polluants dans l'air ou dans l'eau, au bruit, etc. Afin de mieux comprendre ces divers phénomènes, il est notamment nécessaire d'augmenter la densité spatiale d'instruments de mesure. Cela pourrait aider par exemple à l'analyse de l'exposition réelle des populations aux nuisances comme les polluants atmosphériques.

Le déploiement massif de capteurs dans l'environnement est rendu possible par la baisse des coûts des systèmes de mesure, qui utilisent notamment des éléments sensibles à base de micro ou nano technologies. L'inconvénient de ce type de dispositifs est une qualité de mesure insuffisante. Il en résulte un manque de confiance dans les données produites et/ou une hausse drastique des coûts de l'instrumentation causée par les opérations nécessaires d'étalonnage des instruments ou de remplacement périodique des capteurs.

Il existe dans la littérature de nombreux algorithmes qui offrent la possibilité de réaliser l'étalonnage des instruments en les laissant déployés sur le terrain, que l'on nomme techniques d'étalonnage *in situ*.

L'objectif de cette thèse est de contribuer à l'effort de recherche visant à améliorer la qualité des données des instruments de mesure bas coût à travers leur étalonnage *in situ*.

En particulier, on vise à 1) faciliter l'identification des techniques existantes d'étalonnage in situ applicables à un réseau de capteurs selon ses propriétés et les caractéristiques des instruments qui le composent ; 2) aider au choix de l'algorithme le plus adapté selon le réseau de capteurs et son contexte de déploiement ; 3) améliorer l'efficacité des stratégies d'étalonnage in situ grâce au diagnostic des instruments qui ont dérivé dans un réseau de capteurs.

Trois contributions principales sont faites dans ces travaux. Tout d'abord, une terminologie globale est proposée pour classer les travaux existants sur l'étalonnage *in situ*. L'état de l'art effectué selon cette taxonomie a montré qu'il y a de nombreuses contributions sur le sujet, couvrant un large spectre de cas. Néanmoins, le classement des travaux existants selon leurs performances a été difficile puisqu'il n'y a pas d'étude de cas de référence pour l'évaluation de ces algorithmes.

C'est pourquoi dans un second temps, un cadre pour la simulation de réseaux de capteurs est introduit. Il vise à guider l'évaluation d'algorithmes d'étalonnage in situ. Une étude de cas détaillée est fournie à travers l'évaluation d'algorithmes pour l'étalonnage in situ de réseaux de capteurs statiques et aveugles. Une analyse de l'influence des paramètres et des métriques utilisées pour extraire les résultats est également menée. Les résultats dépendant de l'étude de cas, et la plupart des algorithmes réétalonnant les instruments sans évaluer au préalable si cela est nécessaire, un outil d'identification permettant de déterminer les instruments qui sont effectivement fautifs en termes de dérive serait précieux.

Dès lors, la troisième contribution de cette thèse est un algorithme de diagnostic ciblant les fautes de dérive dans les réseaux de capteurs sans faire d'hypothèse sur la nature du réseau de capteurs considéré. Basé sur le concept de rendez-vous, l'algorithme permet d'identifier les instruments fautifs tant qu'il est possible de supposer qu'un instrument n'est pas fautif dans le réseau de capteurs. À travers l'analyse des résultats d'une étude de cas, nous proposons différents

moyens pour diminuer les faux résultats et des recommandations pour régler les paramètres de l'algorithme. Enfin, nous montrons que l'algorithme de diagnostic proposé, combiné à une technique simple d'étalonnage, permet d'améliorer la qualité des résultats de mesure. Ainsi, cet algorithme de diagnostic ouvre de nouvelles perspectives quant à l'étalonnage *in situ*.

List of publications

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In Journals

- F. Delaine, B. Lebental and H. Rivano, "In Situ Calibration Algorithms for Environmental Sensor Networks: A Review," in *IEEE Sensors Journal*, vol. 19, no. 15, pp. 5968-5978, 1 Aug.1, 2019, DOI: 10.1109/JSEN.2019.2910317.
- F. Delaine, B. Lebental and H. Rivano, "Framework for the Simulation of Sensor Networks Aimed at Evaluating In Situ Calibration Algorithms," in *Sensors*, vol. 20, no. 16, 2020, DOI: 10.3390/s20164577.

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F. Delaine, B. Lebental and H. Rivano, "Methodical Evaluation of In Situ Calibration Strategies for Environmental Sensor Networks," Colloque CASPA, Paris, April 2019.

Online repositories

F. Delaine, B. Lebental and H. Rivano, "Example case study applying a "Framework for the Simulation of Sensor Networks Aimed at Evaluating In Situ Calibration Algorithms",", Université Gustave Eiffel, 2020, DOI: 10.25578/CJCYMZ.

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List of Acronyms and Abbreviations

```
VIM
              International Vocabulary of Metrology
   ARMA
              Autoregressive moving average
   SM-SVD Subspace model + Singular value decomposition (calibration strategy used in Chapter
487
              3)
488
              Subspace model + Least squares (calibration strategy used in Chapter 3)
   SM-LS
   SM-TLS Subspace model + Total least squares (calibration strategy used in Chapter 3)
   AB-DT
              Average Based + Difference trigger (calibration strategy used in Chapter 3)
   AB-KF
              Average Based + Kalman filter (calibration strategy used in Chapter 3)
   SVR-KF Support Vector Regression + Kalman filter (calibration strategy used in Chapter 3)
   K-KF
              Kriging + Kalman filter (calibration strategy used in Chapter 3)
   PCA
              Principal components analysis
   WGLI
              Weekly gain linear increase (drift model defined in Chapter 3 Section 3)
   RGOI
              Random gain and offset increase (drift model defined in Chapter 3 Section 5.3.1)
497
   CGOI
              Continuous gain and offset increase (drift model defined in Chapter 3 Section 5.3.1)
```

503

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506

The notations are organised into several categories to quickly identify a set of notations used 500 in a particular context. 501

Note from the author: Conflicts exist due to identical notations that are used for different objects depending on the context. It concerns notably the measurand models (2D Gaussian Model and Gaussian Plume model) and the parameters of the calibration algorithms to remain consistent with the usual notations in the literature.

Common objects

507	t	Instant of time
508	Δt	Time duration
509	ω	Angular frequency
510	E	Expectation
511	$\mathcal{N}(lpha,eta)$	Normal law of mean α and standard deviation β .
512	$\mathcal{U}(lpha,eta)$	Uniform law on the range $[\alpha; \beta]$
513	μ	Average
514	σ	Standard deviation

515	Measuring in	nstruments
516	s_i, s_j	Measuring instruments (or systems if it is mentioned)
517	$c(s_i)$	Accuracy class of s_i
518 519	S	Set of measuring instruments. S^k , S^{k+} and S^{k-} are respectively the sets of measuring instruments where $c(s_i) = k$, $c(s_i) \ge k$ and $c(s_i) \le k$
520	$m(s_i,t)$	Measurement result of s_i at t
521	$M(s_i,(t,\Delta t))$	Set of measurement results for s_i , over the time range $[t - \Delta t; t]$
522	$v(s_i,t)$	Measured value of s_i at t
523	$V(s_i,(t,\Delta t))$	Set of measured values for s_i , over the time range $[t - \Delta t; t]$
524	$v_{true}(s_i,t)$	True value that should be measured by s_i at t if it were ideal
525 526	$\Delta v(s_i, t)$	Measurement uncertainty of the value measured by s_i at t . It can be a constant or not

Relative measurement uncertainty of the value measured by s_i at t.

 $v_{min}(k)$ is the detection limit of instruments of class k

529 $\zeta(s_i,t)$ Indication of s_i at t

Function representing the measuring chain of an instrument

Estimate of the measuring chain of an instrument

Inverse of the estimate of the measuring chain of an instrument

Vector of values of influence quantities $\mathbf{q}(t)$

Estimate of the vector of values of influence quantities

Relationship built during an $in \, situ$ calibration

Calibration relationship derived from an in situ calibration

Fault models

537

G Gain drift

539 O Offset drift

540 arepsilon Noise

541 ψ Spike value

542 p_{ψ} Spike probability

2D Gaussian model

C Concentration

545 x, y, z Coordinates

A Amplitude

Standard deviation

548 FWHM Full width at half maximum

Gaussian Plume model

Emission rate at the source

551 V_w Wind speed

Horizontal and vertical dispersions

H Pollutant effective release

$$H = h_s + \Delta h(t)$$

with h_s , the pollutant source height, and Δh :

$$\Delta h(t) = \frac{1.6F^{\frac{1}{3}}x^{\frac{2}{3}}}{V_w}$$

with F such as

$$F = \frac{g}{\pi} D\left(\frac{\theta_s - \theta(t)}{\theta}\right)$$

g gravity constant

D volumetric flow

 θ ambient temperature

 θ_s source temperature

Parameters of the calibration algorithms

Number of components of the PCA for the algorithms SM-X

Parameter for the algorithms SM-X

Parameter for the algorithms X-KF

Penalty parameter for the algorithm SVR-KF

Kriging parameters for the algorithm K-KF

Usual metrics

556

557

MAE Mean absolute error

MAPE Mean absolute percentage error

Root-mean-square error

Pearson correlation coefficient

 R^2 Coefficient of determination

Error model

574

 \mathcal{F} Function

Slope Slope

572 b Intercept

573 ε Additive error

Metrics for the evaluation of the performances of the diagnosis algorithm

TN True negative

False negative

NDN Non-determined negative

FP False positive

TP True positive

NDP Non-determined positive

Number of positives P

582	N	Number of negatives			
583	Prev	Prevalence			
584	TPR	True positive rate			
585	TNR	True negative rate			
586	FPR	False positive rate			
587	FNR	False negative rate			
588	NDPR	Non-determined positive rate			
589	NDNR	Non-determined negative rate			
590	NDR	Non-determined rate			
591	PPV	Positive predictive value			
592	FDR	False discovery rate			
593	NPV	Negative predictive value			
594	FOR	False omission rate			
595	ACC	Accuracy			
596	$\Delta \mathcal{D}(s_i)$	Delay of first positive detection for an instrument s_i			
	Diagnosis algorithm				
597	Diagnosis al	$\operatorname{gorithm}$			
597 598	Diagnosis alg	gorithm Set of diagnosis procedures			
598	D	Set of diagnosis procedures			
598 599	D d	Set of diagnosis procedures Diagnosis procedure			
598599600601	D d t_d	Set of diagnosis procedures			
598599600601602	D d t_d $\Delta c_{min}^{\mathcal{D}}(k)$	Set of diagnosis procedures			
598599600601602603	D d t_d $\Delta c_{min}^{\mathcal{D}}(k)$ $c_{min}^{\mathcal{D}}(s_i)$	Set of diagnosis procedures			
598 599 600 601 602 603 604	D d t_d $\Delta c_{min}^{\mathcal{D}}(k)$ $c_{min}^{\mathcal{D}}(s_i)$ $\Omega(s_i,t)$	Set of diagnosis procedures Diagnosis procedure Instant of the diagnosis procedure d Minimal relative difference of class required for instruments of class k with their diagnoser instruments Minimal class of sensors allowed to diagnose s_i . True status of s_i at t , equal either to NF (non-faulty) or F (faulty) Predicted status of s_i at t , equal either to NF (non-faulty), F (faulty) or A			
598 599 600 601 602 603 604 605 606	D d t_d $\Delta c_{min}^{\mathcal{D}}(k)$ $c_{min}^{\mathcal{D}}(s_i)$ $\Omega(s_i,t)$ $\widehat{\Omega}(s_i,t)$	Set of diagnosis procedures Diagnosis procedure Instant of the diagnosis procedure d Minimal relative difference of class required for instruments of class k with their diagnoser instruments Minimal class of sensors allowed to diagnose s_i . True status of s_i at t , equal either to NF (non-faulty) or F (faulty) Predicted status of s_i at t , equal either to NF (non-faulty), F (faulty) or A (ambiguous) Actualised status of s_i at t , equal either to NF (non-faulty), F (faulty) or			
598 599 600 601 602 603 604 605 606 607 608	D d t_d $\Delta c_{min}^{\mathcal{D}}(k)$ $c_{min}^{\mathcal{D}}(s_i)$ $\Omega(s_i,t)$ $\widehat{\Omega}(s_i,t)$	Set of diagnosis procedures Diagnosis procedure Instant of the diagnosis procedure d Minimal relative difference of class required for instruments of class k with their diagnoser instruments Minimal class of sensors allowed to diagnose s_i . True status of s_i at t , equal either to NF (non-faulty) or F (faulty) Predicted status of s_i at t , equal either to NF (non-faulty), F (faulty) or A (ambiguous) Actualised status of s_i at t , equal either to NF (non-faulty), F (faulty) or A (ambiguous)			
598 599 600 601 602 603 604 605 606 607 608	D d t_d $\Delta c_{min}^{\mathcal{D}}(k)$ $c_{min}^{\mathcal{D}}(s_i)$ $\Omega(s_i,t)$ $\widehat{\Omega}(s_i,t)$ $\widetilde{\Omega}(s_i,t)$	Set of diagnosis procedures Diagnosis procedure Instant of the diagnosis procedure d Minimal relative difference of class required for instruments of class k with their diagnoser instruments Minimal class of sensors allowed to diagnose s_i . True status of s_i at t , equal either to NF (non-faulty) or F (faulty) Predicted status of s_i at t , equal either to NF (non-faulty), F (faulty) or A (ambiguous) Actualised status of s_i at t , equal either to NF (non-faulty), F (faulty) or A (ambiguous) Set of sensors to diagnose			

```
M^{\approx}(s_i,(t,\Delta t))
                               Set of measurement results coherent with true values for s_i, over the time
613
                               range [t - \Delta t; t]
614
     M^+(s_i,(t,\Delta t))
                               Set of measurement results upper non-coherent with true values for s_i, over
615
                               the time range [t - \Delta t; t]
616
     M^-(s_i,(t,\Delta t))
                               Set of measurement results lower non-coherent with true values for s_i, over
617
                               the time range [t - \Delta t; t]
618
                               Set of measurement results for s_i, over the time range [t - \Delta t; t] that are
     M^*(s_i,(t,\Delta t))
619
                               metrologically valid
620
     \varphi(s_i \to s_j, t)
                               Rendez-vous between s_i and s_j at t
621
     \Phi(s_i \to s_j, (t, \Delta t))
                              Set of rendez-vous between s_i and s_j over the time range [t - \Delta t; t]. \Phi(s_i \to t)
622
                               S_i(t,\Delta t) is the set of rendez-vous between s_i and any other instrument of
623
                               S over the time range [t - \Delta t; t]. It is also denoted \Phi(s_i, (t, \Delta t))
624
     \Phi^{\approx}(s_i \to s_i, (t, \Delta t)) Set of coherents rendez-vous between s_i and s_i over the time range [t - \Delta t; t]
625
     \Phi^+(s_i \to s_j, (t, \Delta t)) Set of upper non-coherents rendez-vous between s_i and s_j over the time
626
                               range [t - \Delta t; t]
627
     \Phi^-(s_i \to s_j, (t, \Delta t)) Set of lower non-coherents rendez-vous between s_i and s_j over the time
628
                               range [t - \Delta t; t]
629
     \Phi(s_i \to S, (t, \Delta t))|_k Set of rendez-vous between s_i and all s_j \in S such as c(s_j) \ge k over the time
630
                               range [t - \Delta t; t]
631
     \mathbf{Mat}_{\mathbf{\Phi}}(D, \Delta t)
                               Matrix of the minimal number of rendez-vous between two instruments on a
632
                               duration \Delta t and over a set of diagnosis procedures D
633
     \Phi_v(s_i \to S, (t, \Delta t))
                              Set of rendez-vous between s_i and s_i \in S over the time range [t - \Delta t; t] that
634
                               are valid
635
     |\Phi_v|_{min}
                               Minimal number of valid rendez-vous required to conduct successfully a
636
                               diagnosis
637
    r_{\varphi_v}^{\approx}(s_i,(t,\Delta t))
                               Rate of coherent rendez-vous in the set of valid rendez-vous of s_i over the
638
                               time range [t - \Delta t; t]
639
    r_{\varphi_v}^+(s_i,(t,\Delta t))
                               Rate of upper non-coherent rendez-vous in the set of valid rendez-vous of s_i
640
                               over the time range [t - \Delta t; t]
641
    r_{\varphi_v}^-(s_i,(t,\Delta t))
                               Rate of lower non-coherent rendez-vous in the set of valid rendez-vous of s_i
642
                               over the time range [t - \Delta t; t]
643
    (r_{\varphi_n}^+)_{max}
                               Maximal tolerated value for r_{\varphi_n}^+(s_i,(t,\Delta t)) for any t
644
     (r_{\varphi_n}^-)_{max}
                               Maximal tolerated value for r_{\varphi_n}^-(s_i,(t,\Delta t)) for any t
645
     (r_{\varphi_v}^+ + r_{\varphi_v}^-)_{max}
                               Maximal tolerated value for (r_{\varphi_v}^+ + r_{\varphi_v}^-)(s_i, (t, \Delta t)) = 1 - r_{\varphi_v}^{\approx}(s_i, (t, \Delta t)) for
646
                               any t
647
     r_{true}(s_i, (t, \Delta t))
                               Rate of measurement results coherent with true values of s_i over the time
648
                               range [t - \Delta t; t]
649
```

w

654

655

Rate of measurement results coherent with true values in the set of rendezvous of s_i over the time range $[t - \Delta t; t]$

 $r_{true}(s_i,(t,\Delta t))|_{\Phi_v}$ Rate of measurement results coherent with true values in the set of valid rendez-vous of s_i over the time range $[t-\Delta t;t]$

Weight operator for rendez-vous such as:

$$w(\varphi(s_i \to s_j, t)) = \frac{1}{|\Phi(s_i, t)|}$$

with
$$\Phi(s_i, t) = \Phi((s_i, S), t|0)$$

Practical definition of rendez-vous

Position of s_i at t

Maximal distance between two instruments to consider them in rendez-vous (spatial condition)

(spatial condition)

Maximal difference between the timestamps of the measurement results of two instruments to consider them in rendez-vous (temporal condition)

Representativity area of the instrument s_i at t

1 Context of the thesis

The monitoring of ambient quantities is a fundamental operation for the digitisation of our environment [59]. From the level of precipitation [32] to the concentration of pollutants in the air [111], more and more quantities need to be monitored. The information brought by measurements is used in various fields from agriculture [117] to public health [46] for instance. Such a monitoring may have a major impact because these quantities can have significant effects on human lives. For instance, heatwaves, corresponding to high temperatures at both day and night, induce an abnormal mortality rate that can be extremely high [81]. Combined with drought, they can also be disastrous for agriculture and generate water stress [159]. Thus, the comprehension of these phenomena is particularly important to explain and predict their evolution over time but also to react to them.

This activity of monitoring has evolved over time, following the progress in science and technologies. Nowadays, the observation of ambient quantities is carried out with high quality measuring systems, at least for strategic or regulatory applications like meteorology [174] or air quality monitoring [158]. They can be deployed on the ground, in the sea, in the air or even in space via satellites. Regulatory-grade instruments are usually expensive. Regarding those on the ground, they are often deployed at fixed positions. Depending on the target area to monitor, on the spatial variability of the measurand and on the spatio-temporal resolution desired, hundreds of devices may be required, as shown for instance by [21] and [101] in the field of air quality monitoring. Thus, it may not be possible to have a high spatial resolution with such instruments as few of them are deployed together most of the time due to their cost [78]. For instance, about 30 measuring stations are deployed across Paris and its inner suburbs to monitor air pollution [2] as shown in the figure page 2. Moreover, all of them are not monitoring the same quantities. For instance, only a third of these stations are monitoring O_3 . Consequently, the spatial resolution is even smaller for some measurands.

Following up the scientific advances in micro and nano-electronics [90, 135], it is now possible to imagine a dense monitoring of ambient quantities with a fine granularity, both temporally and spatially, notably for air quality monitoring [110, 135] which is the practical context chosen for this thesis. Indeed, small low-cost measuring instruments have been designed in recent years. Fostered by the emergence of the Internet of Things, the interest for such devices has been growing significantly because they open up new possibilities for environmental sensing [62]. Without actually replacing regulatory grade instruments, it is hoped they can complement existing networks of measuring instruments. As they are affordable, they can also be owned by everybody. In this way, low-cost sensors are tools that could be used for individual awareness and education, particularly concerning both indoor and outdoor air quality [27].

Nevertheless, these low-cost sensing technologies are still young, and several challenges are yet to be tackled [62, 127]. One of them concerns the quality of the measurement results they produce. The particular issue degrading the results addressed in this thesis is the instrumental drift which these instruments are prone to. The classical approach to solve this problem is to recalibrate the instruments in a calibration facility. This requires taking the instrument out of service, bringing



Symbol	Definition		
	Permanent urban station		
	Permanent traffic station		

Map of permanent air pollution monitoring stations of Airparif for Paris and its inner suburbs (December 2018)

it to a facility, calibrating it, bringing it back to the field and putting it back in service. This operation has both economically and technically a definite cost. Therefore, when hundreds of measuring instruments are deployed, performing this task is particularly challenging because it has to be carried out more frequently than with high-quality instruments. However, because numerous instruments are deployed and forming a sensor network, there may be relationships between their measured values that could be exploited to manage this issue.

In the literature, multiple solutions have been proposed to mitigate the impact of the drift for environmental low-cost sensors by performing what we call an *in situ* calibration. It means the calibration of measuring instruments while leaving them in the field, preferably without any physical intervention. To avoid these physical interventions, the idea is to use the values provided by other instruments deployed.

2 Contributions

The objective of this thesis is to contribute to the research effort on the improvement of data quality for low-cost measuring instruments through their *in situ* calibration. In particular, we aim at:

- facilitating the identification of existing *in situ* calibration strategies applicable to a sensor network depending on its properties and the characteristics of its instruments.
- helping to choose the most suitable algorithm depending on the sensor network and its context of deployment.
- improving the efficiency of in situ calibration strategies through the diagnosis of instruments

that have drifted in a sensor network.

Toward this goal, three main contributions are made in this work. First, a unified terminology is proposed to classify the existing works on *in situ* calibration. Indeed there is no shared vocabulary in the scientific community enabling a precise description of the main characteristics of the algorithms. The review carried out based on this taxonomy showed there are numerous contributions on the subject, covering a wide variety of cases. Due to the type of sensor network deployed, in terms of properties of the measuring instruments or how these devices can interact between them, different approaches may be considered. Nevertheless, the classification of the existing works in terms of performances was difficult as there is no reference case study for the evaluation of these algorithms.

Therefore in a second step, a framework for the simulation of sensor networks is introduced. It is aimed at evaluating *in situ* calibration algorithms. It details all the aspects to take into account when designing a case study. A detailed case study is provided across the evaluation of *in situ* calibration algorithms for blind static sensor networks. An analysis of the influence of the parameters and of the metrics used to derive the results is also carried out. As the results are case specific, and as most of the algorithms recalibrate instruments without evaluating first if they actually need it, an identification tool enabling to determine the instruments that are actually faulty in terms of drift would be valuable.

Thus, the third contribution of this thesis is the design of a diagnosis algorithm targeting drift faults in sensor networks without making any assumption on the kind of sensor network at stake. Based on the concept of rendez-vous, the algorithm allows identifying faulty instruments as long as one instrument at least can be assumed as non-faulty in the sensor network. Across the investigation of the results of a case study, we propose several means to reduce false results and guidelines to adjust the parameters of the algorithm. Finally, we show that the proposed diagnosis approach, combined with a simple calibration technique, enables to improve the quality of the measurement results.

3 Organisation of the manuscript

The manuscript is organised as follows.

In Chapter 1, concepts related to measuring instruments are defined and the performances of low-cost measuring instruments in the context of air quality monitoring are reviewed. Then, the issue of data quality for these devices and how calibration can mitigate drift problems are discussed. The concept of *in situ* calibration is introduced, followed by the interesting properties of sensor networks for this application. At the end of the chapter, the problem statement is recalled and detailed.

In Chapter 2, the taxonomy for the classification of *in situ* calibration strategies is introduced, followed by the review of existing works.

Chapter 3 introduces the framework for the simulation of sensor networks designed for the evaluation of *in situ* calibration algorithms. A case study concerning blind static sensor networks is developed, and multiple derived cases are investigated to determine the influence of parameters and of the metrics used on the interpretations of the results.

Afterwards, the diagnosis algorithm for drift faults in sensor networks is introduced in Chapter 4. A case study is developed and means to improve its results are discussed. The sensitivity of the diagnosis algorithm to the choices made during the design of the case study is investigated. It is followed by the combination of the diagnosis algorithm with a simple calibration approach that exploits information build during the diagnosis, applied to the preceding case study.

Finally, the contributions are summarised in the last chapter and a general conclusion is provided, along with perspectives regarding this work.

Several appendices are also provided in this manuscript. Appendix A provides variations and extensions of the diagnosis algorithm presented in Chapter 4. They require further studies but the basis are introduced towards future work. Appendix B extends the discussion on the main assumption made for the diagnosis algorithm. Lastly, Appendix C presents an additional case study related to the sensitivity of the diagnosis algorithm.

Chapter

Low-cost Measuring Instruments for Air Quality Monitoring: Description, Performances and Challenges

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Introduction

This chapter is aimed at recalling several concepts related to measuring instruments first. The question of the measurement of ambient quantities is detailed by defining low-cost instruments and by reviewing their performances in the context of air quality monitoring. Afterwards, the challenges for these instruments are defined. The threats to the data quality of measuring instruments are discussed before investigating how calibration can mitigate drift problems. Finally, the concept of *in situ* calibration is introduced, followed by the definition of a sensor network and its interesting characteristics for our problem. At the end of the chapter, the problem statement is detailed based on the previous developments.

1 Measurement of ambient quantities with low-cost instruments

In this section, the definition of a measuring instrument is recalled first. Then, a general description of how it makes measurements is provided, followed by the definition of low-cost instruments, their challenges and a review of their actual performances in metrological terms.

1.1 Definition of a measuring instrument

In environmental monitoring, we aim at tracking one or several quantities through measurements of their values over time. Such quantities are for instance temperature, relative humidity, pressure, concentration of chemical components in a gas or a liquid, noise... These quantities we aim at monitoring are called measurands.

- Definition 1 (Quantity [14]). A quantity is the "property of a phenomenon, body, or substance, where the property has a magnitude that can be expressed as a number and a reference".
- Definition 2 (Quantity value [14]). A quantity value is a "number and reference together expressing the magnitude of a quantity".
- Definition 3 (Measurement [14]). A measurement is the "process of experimentally obtaining one or more quantity values that can reasonably be attributed to a quantity".
- Definition 4 (Measurand [14]). A measurand is the "quantity intended to be measured".
 - To carry out measurements, one or several measuring instruments are used. Multiple devices may be required to perform a measurement and they can be embedded into a single device as a measuring system. Such a system may be used to track several measurands.
- Definition 5 (Measuring instrument [14]). A measuring instrument is a "device used for making measurements, alone or in conjunction with one or more supplementary devices".
- Definition 6 (Measuring system [14]). A measuring system is a "set of one or more measuring instruments and often other devices, including any reagent and supply, assembled and adapted to give information used to generate measured values within specified intervals for quantities of specified kinds".

1.2 Measuring chain of an instrument

In a measuring instrument (or system), several components may be necessary to perform a measurement, forming a measuring chain. It usually starts by one or several sensors¹ and provides as an output one or several indications. For example, consider a mercury-in-glass thermometer (Figure 1.1.1). The sensor is the mercury. Depending on the variation of the volume of the liquid following temperature, its level in a tube of glass is the indication. With an appropriate scale of value, this level can be converted back to a temperature through direct reading.

¹In ordinary use, the term sensor may refer to a measuring instrument. This is a slight misuse of language.

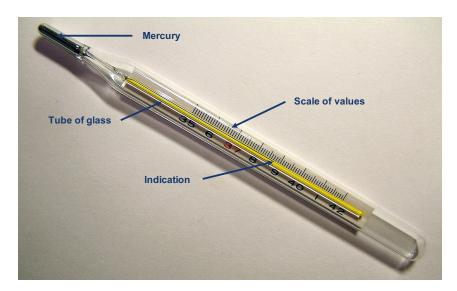


Figure 1.1.1: Description of a mercury-in-glass thermometer²

Definition 7 (Measuring chain [14]). A measuring chain is a "series of elements of a measuring system constituting a single path of the signal from a sensor to an output element".

Definition 8 (Sensor [14]). A sensor is the part of a measuring instrument "that is directly affected by a phenomenon, body, or substance carrying a quantity to be measured".

Definition 9 (Indication [14]). An indication is a "quantity value provided by a measuring instrument or a measuring system".

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As shown in Figure 1.1.1, the raw indication provided by a measuring instrument is not necessarily of the same kind as the measurand. There is often a quantity conversion that is performed at the end of the measuring chain so that the measurand and the indications provided are of the same kind to facilitate the readings. In this case, the indication provided by a measuring instrument is a measured value. For the conversion step, values from different instruments may be involved in the case of a measuring system.

Definition 10 (Measured value [14]). A measured value is a "quantity value representing a measurement result".

The values provided by measuring instruments are not necessarily equal to the actual values of the measurand affecting the sensor. The actual values of measurand are called true quantity values.

Definition 11 (True quantity value [14]). A true quantity value, or true value, is a "quantity value consistent with the definition of a quantity".

Thus, an uncertainty is usually associated with a measured value to give an idea of its precision. Its value can be determined through various means (see [72]), the components of the measuring chain having an influence.

Definition 12 (Uncertainty³ [14]). An uncertainty is a "non-negative parameter characterizing the dispersion of the quantity values being attributed to a measurand, based on the information used."

²Modified image. Original source: Menchi (Licence CC BY-SA), https://commons.wikimedia.org/w/index.php?curid=51236

³The concept of uncertainty is different from the concept of measurement error because even after evaluating the components of the error that are known or supposed, an uncertainty always remains [72]

This is notably why a measured value is a value representing a measurement result but not a measurement result itself. Usually, a measurement result is expressed by a measured value and an uncertainty. A measurement result composed of only a measured value does not allow to estimate how close it is to the true value. In fact, any important information for the interpretation of a measurement result should be added to it. It can concern the operating conditions, or data helping to identify when or where a result was obtained, like a timestamp or GPS coordinates.

Definition 13 (Measurement result [14]). A measurement result is a "set of quantity values being attributed to a measurand together with any other available relevant information".

All along the measuring chain, influence quantities may act on the output of the instrument.

Definition 14 (Influence quantity [14]). An influence quantity is a "quantity that, in a direct measurement, does not affect the quantity that is actually measured, but affects the relation between the indication and the measurement result".

The measuring chain of an instrument can be more or less long depending on the measurement method. For a digital thermometer, based on a thermocouple for instance, the sensor provides a voltage, as an indication of the temperature. Several processings are usually carried out afterwards like amplification and filtering, which adds steps in the measuring chain compared to the case of a mercury-in-glass thermometer.

The description of the measuring chain of an instrument is summarised in Figure 1.1.2.

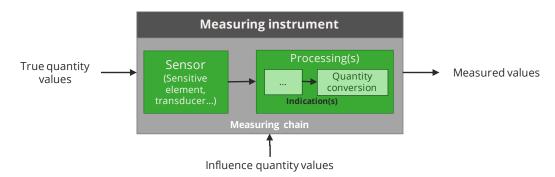


Figure 1.1.2: Schematic description of the measuring chain of a measuring instrument

1.3 Low-cost instruments

1.3.1 Definition

For any measurand, there is a wide range of measuring instruments notably in terms of cost. Their price is related to the capabilities of instruments in terms of measurement range, sensitivity, uncertainty and so on. In the General Introduction, we exposed that reference instruments were sometimes too expensive to enable a dense environmental monitoring of different quantities, taking as an example the context of air quality monitoring. Consequently, low-cost instruments are necessary to help carry out this task.

Definition 15 (Reference measuring instrument). A reference (or regulatory-grade) measuring instrument is a measuring instrument implementing a reference measurement procedure or equivalent. Such an instrument is usually expensive.

Definition 16 (Low-cost measuring instrument). A low-cost measuring instrument is a measuring instrument significantly affordable compared to reference instruments.

Depending on the application, what a low-cost measuring instrument is may be very different [27, 110, 138]. In the context of air pollution monitoring, we speak of devices that cost US\$ 100 or less to few US\$ 1,000 [71, 85, 110] at most, considering that reference instruments can be worth from US\$ 10,000 to more than US\$ 100,000.

Low-cost instruments are often characterised by a small size, thanks to the use of micro and nano technologies [183]. They are also frequently autonomous in terms of energy: they are powered by batteries or small solar panels. Most of the time they have wireless communications capabilities. They can be deployed at fixed positions or be mobile and even wearable [66].

1.3.2 Challenges of environmental monitoring

Few years ago, Hart et al. [62] and Rundel et al. [127] exposed various challenges for measuring devices used in environmental monitoring:

- **Energy supply and management** This subject concerns mainly autonomous devices. In this case, it aims notably at minimising the energy consumption so that size of the system of energy storage and charge can be reduced. In addition, the lifespan of the energy supply components has to be the longest as possible to limit the maintenance operations and its environmental impact.
- Hardware standardisation To have upgradeable systems, interoperability of the hardware must be met, with user-friendly manageability.
 - **Software standardisation** In the same way that hardware has to be standardised, the software of measuring instruments should also be normalised to facilitate the usage of devices from different manufacturers in the same deployment.
- Data standardisation To help data management and the use of values, data formats must be standardised.
 - **Data quality improvement** This challenge concerns different aspects. First, in conjunction with data standardisation, the data quality has to be ensured by providing for each measurement all the necessary and meaningful information regarding its realisation. Secondly, the accuracy of the measurements has to be improved through the upgrade of the hardware but also with the help of data processing algorithms.
 - **Data usage reinforcement** To make sure the most is taken out of the acquired data, it is necessary to facilitate its management and use. It would also be valuable to exploit the measurements in the largest number of possible applications.
 - **Security reinforcement** As systems are more and more connected, they become vulnerable to attacks that can affect data integrity. Protocols must be established to prevent tampering and protect privacy.

Among all these challenges, data quality is an important one for low-cost instruments in particular and is still a major issue [119]. Indeed, if a satisfying data quality cannot be achieved, the responses to any of the other challenges cited will have a limited impact.

1.3.3 Performances of low-cost instruments in the literature for air quality monitoring

While the concepts introduced and methods developed in this work are generic, the examples and illustrations provided are taken from the field of air quality monitoring. Indeed, regulatory-grade instruments as used by air quality surveillance agencies are extremely expensive. Therefore,

the interest for low-cost devices is major in this field [27]. To evaluate the severity of data quality issues for these devices, a review of their actual performances is conducted in this section.

In the literature, there are numerous works proposing an evaluation of the performances of low-cost sensors for air quality monitoring. In a review from 2019 [73], the authors reported 105 references "that include quantitative comparison of sensor data with reference measurements" from 2004 to April 2019, with 91 between January 2015 and December 2018. We consider here the following works: [15, 16, 71, 73, 85, 121, 141], some of them being aggregates of previous publications [73, 121]. They tackle the most common measurands regarding to air quality: PM_{10} [15, 16, 73, 121, 141], $PM_{2.5}$ [15, 16, 71, 73, 85, 121, 141], PM_1 [73, 141], PM_2 [15, 16, 71, 73, 85, 121, 141], PM_3 [15, 16, 71, 73, 85, 121] and PM_3 [15, 16, 71, 73, 121, 141], PM_3 [15, 16, 71, 73, 85, 121] and PM_3 [15, 16, 71]. In these studies, the considered low-cost measuring instruments or systems are co-located to reference, e.g. regulatory-grade, measuring instruments at a given location, for instance an urban environment [15, 141], a suburban environment [71] or a mountain site [85].

These studies report a wide range of performances for the measuring instruments or systems they consider: from poor to suitable at least for indicative measurements, according to the Data Quality Objectives (DQO) set by the European Union [158], and this for all the measurands. Minimum and maximal values of the coefficient of determination R^2 reported in the cited publications for PM_{10} , $PM_{2.5}$, PM_1 , O_3 , NO_2 , NO, CO and SO_2 are listed in Table 1.1.1. This metric indicates the quality of the linear regression between the values of an instrument and a reference.

Publication	PM_{10}	$PM_{2.5}$	PM_1	O_3	NO_2	NO	CO	SO_2
1 donouvon	min – max	min – max	$\min - \max$	min – max	min – max	min – max	min – max	min – max
Borrego et al. [15]	0.13-0.36	0.07-0.27	n/a	0.12-0.77	0.02-0.89	0.34-0.80	0.53-0.87	0.09-0.20
Jiao et al. [71]	< 0.2	5-0.45	n/a	< 0.25 – 0.94	< 0.25 – 0.57	0.77 – 0.87	< 0.25 – 0.68	< 0.25
Li et al. [85]	n/a	0.82	n/a	0.62	n/a	n/a	0.68	n/a
Spinelle et al. [141]	0.01 - 0.81	0.00 - 0.86	0.26 - 0.95	0.04 - 0.96	0.35 - 0.88	n/a	n/a	n/a
Rai et al. [121]*		0.07 – 0.99		0.01 – 0.99	< 0.10 – 0.89	n/a	< 0.10-0.99	n/a
Karagulian et al. [73]*	0.00 – 1.00	0.00 – 1.00	<0.600.95<	0.00 – 1.00	0.00 – 1.00	< 0.10 – 1.00	0.25 – 1.00	n/a

Table 1.1.1: Minimum and maximal values of the coefficient of determination R^2 reported in the cited publications for PM_{10} , $PM_{2.5}$, PM_1 , O_3 , NO_2 , NO, CO and SO_2 . Publications marked with a "*" are surveys aggregating results from other works. Sometimes the values were given for several types of PM instruments. The values of this table concern different types of devices and do not relate the variability that can be observed with several instruments of the same manufacturer [141].

Several authors report that their results should be considered with caution. On the first hand, those who obtained results only at one site, under a certain type of operating conditions, explain that their work should be repeated with different conditions [141]. On the other hand, those who aggregated results point out the diversity of the metrics used [73] and of the conditions under which the results were obtained [121], which makes the comparisons difficult. Karagulian *et al.* [73] chose to rely on the coefficient of determination R^2 to compare the highest number of studies, but they explained that this metric is not the most appropriate. They call for standardised protocols for the evaluation of low-cost measuring instruments. Such a protocol for the evaluation for low-cost gas sensors for air quality monitoring has been proposed [138].

In all of these works, the calibration of the instruments is a main concern. In Section 1.2, we mentioned that in general, measuring instruments provide indications not necessarily of the same kind as the measurand and that a quantity conversion is often carried out to facilitate the reading of measurement results. Determining the relationship to obtain measurements results from indications is what calibration allows to achieve. This operation is carried out on a measuring instrument for the first time right after its manufacturing or before its first use [67]. Generally,

it is performed in a dedicated facility. However, it is only mandatory to have an individual calibration for regulatory-grade instruments. In most cases, several instruments are calibrated altogether. Sometimes it also happens that instruments are not calibrated at all.

In the publications cited, out-of-the-box measurement results were quite inaccurate most of the time and some instruments had to be compensated following influence quantities like temperature, relative humidity or chemical components depending on the measurand and on the measuring principle at stake [16, 71, 73, 121]. Jiao *et al.* [71] showed for instance an improvement of the coefficient R^2 :

- from 0.36 to 0.43, from 0.45 to 0.60 and from 0.43 to 0.50 for three different PM instruments
- from 0.88 to 0.94 for an O_3 instrument
- from 0.57 to 0.82 for a NO_2 instrument

In conclusion, all the authors report that the quality of measurement results can be strongly improved with the use of data treatment and processing tools. To better understand how data processing algorithms can improve the quality of the measurement results, the issues causing their degradation must be identified.

2 Threats to data quality for measuring instruments

2.1 Introduction

 In Section 1.3.2, the challenge of data quality was briefly presented from different perspectives. First, it can be related to the quality of the information composing a measurement result, e.g. the measured value, its uncertainty, the conditions under which the measurement was made and so on

In another way, it can be considered as a lack (or not) of measurement accuracy compared to the one of reference instruments in terms of sensitivity or resolution for instance. Such issues can be known at the design or prototyping stage of measuring instruments depending on their components. Therefore, to improve the quality of the measurements, the performances of the instruments have to be upgraded. This can be achieved along the advances during the conception and manufacturing of its components.

The data quality issue can also be observed as a challenge regarding the ability of a measuring instrument to deliver results consistent with its specifications during its life. In the field of dependability, this can be considered as a question of reliability. It is the continuity of correct service delivered by a system [5]. For an instrument, the correct service is the achievement of measurements consistent with its metrological specifications. It is often observed that the quality of low-cost devices decays with time, even under regular operating conditions [33]. This is the general issue we focus on in this work.

The threats to the dependability of a system, and consequently the threats to its reliability, are **failures**, **errors** and **faults** [5]. A failure is an event making the instrument no longer able to deliver a correct service regarding its specifications or requirements. This event is triggered by a deviation of the normal state of the system, called an error, and its potential cause is a fault. The fault itself may be explained by the failure of another part of the system and so on. Thus, there is a relationship of causality between these concepts [5].

In our case, the failure we are interested in for a measuring system is the event when the measurement results it delivers are not correct regarding the actual values of the measurand and the results that should be provided by the device. To improve the quality of data provided by low-cost measuring instruments, it is necessary to determine the faults that are causing this failure. The faults of a measuring system can be studied from different perspectives depending on the case [115]. Consider an instrument that is in failure mode from the point of view of data

due to the absence of measurement results. With a data-centric point of view, the fault that generated this failure could be called "missing data". However, with a system-centric approach, this "missing data" fault can be explained in several ways, for instance a failure from a component of the measuring chain, a power failure or a communication failure, and these failures could be explained by a wide variety of faults depending on the system considered as illustrated in Figure 1.2.1. Therefore, a point of view must be chosen.

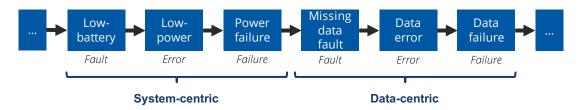


Figure 1.2.1: Example of situation where a fault from a system-centric point of view generates a fault and then a failure from a data-centric perspective.

We propose to take a data-centric point of view. Indeed, from the review of Section 1.3.3, data processing techniques are seen as promising means to improve the quality of the results of measuring instruments. Consequently, we choose to define the faults that we are interested in from what can be observed from the output of the measuring chain of an instrument, e.g. its indications or its measured values, because they are the inputs of data processing algorithms. The identification of the cause of the faults, whether they concern the measuring chain of the instruments or not, will not be studied here.

2.2 Faults

Several taxonomies have been proposed to differentiate faults of measuring instruments and sensor networks [98, 114, 115, 125].

In this work, we propose to use the following taxonomy, inspired from a previous work conducted by Ni et al. [115] that focuses on the relationships between faults taken from different points of view.

Eight different types of faults conducting to a possible data failure were identified. The list of possible faults is illustrated in Figure 1.2.2. In this figure, the fault-less signal is $v_{true}(t) = \sin(\omega t)$ with $t \in [0; 100]$ and $\omega = 2\pi \times 0.01$.⁴

Noise Disturbances present in the indications and which may be correlated or not to the true signal. Usually, it is modelled as a Gaussian white noise, e.g. a random variable following a normal law \mathcal{N} of null average and with a given standard deviation, that is added to the true values. This fault is usually permanent but is also intrinsic to a digital measuring instrument. In fact, noise is considered as a fault when it exceeds the specifications of the instrument. Its severity may vary over time.

In Figure 1.2.2a, the measured values v(t) are equal to $v(t) = v_{true}(t) + \varepsilon(t)$ with $\varepsilon(t) \sim \mathcal{N}(0, 0.3)$.

Spikes The measured values at some instants are very different from the ones expected. This observation is based on the previous and following values, if there is no physical cause related to the measurand that could explain this variation. They can be of two types: single-point (only one point differs in the data set) or multi-points (a few points differ in the data set). The spike is characterised by its amplitude and, for multi-points, by its

⁴The unit of time is arbitrary.

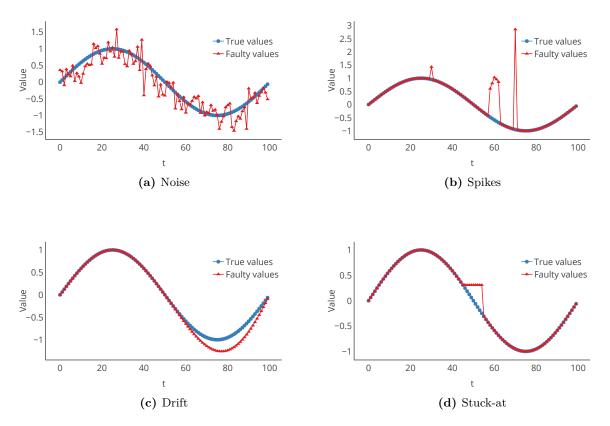


Figure 1.2.2: Considered faults for measuring instruments in the manuscript (continued)

shape, which both may vary. A probability law is associated to determine the frequency of spike occurrence. This fault is transient, but its frequency and severity may vary over time.

In Figure 1.2.2b, two one-point spikes appear at t=30 and t=70 with a respective amplitude of 50% and 250% to the related values, and one multi-points spike appears at t=58 for 5 time steps with a maximal amplitude of 75% of the absolute value of the signal and with a triangular shape.

Drift The drift fault is "a continuous or incremental change over time in indication, due to changes in metrological properties of a measuring instrument" [14]. Usually, this fault is permanent and irreversible, except if it is due to influence quantities which the effect is not compensated by the measuring system and is reversible.

In Figure 1.2.2c, there is a drift of the gain G added, so that $v(t) = (1 + G(t)) \cdot v_{true}(t)$ It is defined as:

$$G(t) = \begin{cases} 0 & \text{if } t < 50\\ 0.01(t - 50) & \text{if } t \ge 50 \end{cases}$$

Stuck-at Several consecutive values are identical when they should not. This fault can be transient or permanent, e.g. it may last for a certain amount of time or forever when it appears.

In Figure 1.2.2d, this fault appears at t = 45 and lasts for 10 time steps.

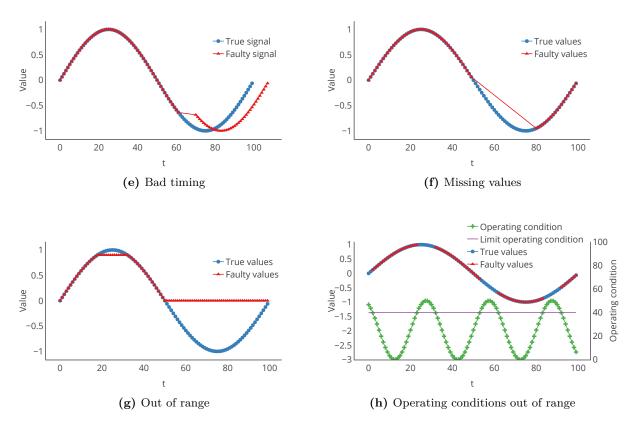


Figure 1.2.2: Considered faults for measuring instruments in the manuscript

Bad timing A measurement result was obtained but has a timestamp differing from the one expected. It can be a transient fault, for instance due to a momentary latency of the instrument, or a permanent fault due to the drift of the clock of the device.

In Figure 1.2.2e, values expected at $t \in [60; 100]$ are recorded with a delay of 8 units of time.

Missing values Measured values have been lost. As for the stuck-at fault, it can be a transient or a permanent fault. In Figure 1.2.2f, values at $t \in [50; 80]$ are missing.

Measured values out of range The recorded measured values are out of the measurement range of the sensor.

The instrument may behave differently depending how it was designed to manage such fault. It may detect this fault and do not record the value, detect it and record it anyway, or it does not have a detection system and record the values normally. If recorded, the value may be equal to one of the bounds of the measurement range, simulating as it is stuck-at, or to another one, more or less related to the true value of the measurand.

In Figure 1.2.2g, we considered the instrument having a range of [0; 0.90] and recording the values, behaving as it is stuck-at when values are out-of-range.

Operating conditions out of range The environment of measurement is out of range of the sensor's normal operating conditions. Like for the fault of measured values out of range, the possibly wrong values may be managed differently depending on the instrument. They can be recorded normally, or recorded but marked as unreliable or not recorded at all.

In Figure 1.2.2h, the instrument is not recording values anymore when the operation condition signal exceeds 40.

2.3 Discussion

This fault taxonomy is a list of the most common data faults that can be observed in sets of measurement results from measuring instruments. Although several faults may be related to each other⁵, they provide a reasonable overview of the threats to the data quality of measuring instruments, in terms of accuracy of the results. The way they appear and how they manifest in the data can be discussed in detail for each type of device but is out of the scope of this work.

These different faults can be differentiated in terms of how an instrument may be able to deal with them. Indeed, low-cost environmental measuring instruments have computing capabilities, though sometimes minimal. Thus, they can be self-aware of their characteristics and may be able to pre-treat their measured values and potentially identify some faults.

For instance, concerning the "out of range" fault, a device can be able to detect when it produces a value higher than its maximal possible output if it knows its measurement range. The instrument may then raise an error indicating it made an incorrect measurement, either because the value of its own measurand was indeed too high compared to its properties or because the measuring instrument is producing an erroneous response to the value of the measurand for instance. This reasoning can be extended to the faults "bad timing", "missing value" or even "operating conditions out of range" if a measuring instrument can access to information on its operating conditions, with the help of other devices. Note that these three faults do not concern directly the measured value but more how it was obtained. On the contrary, "noise", "stuck-at", "spike" and "drift" faults concern directly the measured value like the "out of range" fault. To detect them with a data-centric approach, an analysis of the measured values has to be carried out. "Noise", "stuck-at" and "spike" are faults having a characteristic time shorter than a "drift" fault which usually affects a measuring instrument on a longer time basis. For them it is possible to consider studying the variance or the gradient of the measured signal in a first step [112], although more elaborated techniques have been developed [114]. Be that as it may, self correcting its faults is not straightforward for measuring instruments.

In fact, drift is one of the most problematic faults because it is a continuous change over time in the indications of the instruments. In this way, the relationship between the true values of the measurand and the indications evolves. Consequently, the relationship determined through a previous calibration to derive measured values from true values has to be determined again. This is why calibration is an operation also conducted during the maintenance of an instrument and it must be carried out periodically. To follow the traditional guidelines, it requires to take it out of service and to bring it to a calibration facility. After this operation and the adjustment of the instruments, it can be brought back to its deployment site and put back in service. However, in the context of large deployments of measuring instruments, carrying out this maintenance operation in this manner can have a significant cost, both economically and technically. It is all the more important that low-cost devices are particularly prone to declining performances, often sooner than expected [33].

3 Calibration of measuring instruments

In Section 1.3.3, we briefly introduced the concept of calibration. It was reported as a main concern in studies on the performances of low-cost instruments. Thus, we provide here its formal definition and develop why it is a challenging procedure for large deployments of measuring

⁵For instance, a "spike" fault may trigger a "measured value out of range" fault which itself may be represented by a "stuck-at". However, this is not always the case: a "measured value out of range" fault can happen due to the normal evolution of a measurand.

instruments.

3.1 Definition

The formal definition of calibration is the following:

Definition 17 (Calibration [14]). Calibration is the "operation that, under specified conditions, in a first step, establishes a relation between the quantity values with measurement uncertainties provided by measurement standards and corresponding indications with associated measurement uncertainties and, in a second step, uses this information to establish a relation for obtaining a measurement result from an indication".

Definition 18 (Measurement standard [14]). A measurement standard is the "realisation of the definition of a given quantity, with stated quantity value and associated measurement uncertainty, used as a reference".

In practice, it means in a first step that the calibration of a measuring instruments consists into observing its response to known values of the measurand. From these observations it is possible to determine a relationship giving the indication provided by a measuring instrument following the value of the measurand. To obtain the measured values from the indications, this relationship has to be inverted. This is summarised in Figure 1.3.1

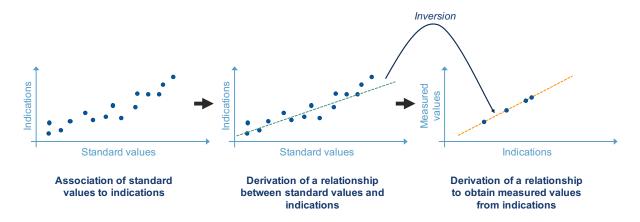


Figure 1.3.1: Schematic description of the steps of a calibration operation

As an example, consider the case of a mercury-in-class thermometer once again. Without its scale, we do not read a temperature value (measured value), we read a level of mercury (indication). Thus, the calibration of this instrument consists into putting the device into several environments for which the conditions are controlled, and the value of the temperature known with its uncertainty. For each value of temperature, the level of mercury is read. Then, it is possible to determine the level of mercury as a function of the temperature. Building the scale for the thermometer consists into inverting this relationship, which is a simple task in this case because it can be approximated as a linear relationship.

Following the VIM [14], "calibration may be expressed by a statement, calibration function, calibration diagram, calibration curve, or calibration table. In some cases, it may consist of an additive or multiplicative correction of the indication with associated measurement uncertainty", a correction being a "compensation for an estimated systematic effect".

The proper achievement of calibration for a measuring instrument is notably meant to ensure its trueness.

3.2 Analysis

Let us have a short analysis of calibration from a mathematical perspective. In this case, the result of calibration is expressed as a function. Consider that $\zeta(s_i, t)$ is the indication provided by s_i at t and that the measuring chain of the measuring instrument is represented as a function \mathcal{F} .

Therefore,

$$\forall t, \zeta(s_i, t) = \mathcal{F}(v_{true}(s_i, t), \mathbf{q}(t))$$

where v_{true} are the true values of the measurand and \mathbf{q} is a vector of the values of the influence quantities of the considered instrument.⁶

The first step of calibration aims at determining $\widehat{\mathcal{F}}$ which is an estimate of \mathcal{F} . Then, the purpose is to derive $\widehat{\mathcal{F}}^{-1}$ such as:

$$v(s_i, t) = (\mathcal{F} \circ \widehat{\mathcal{F}}^{-1})(v_{true}(s_i, t), \mathbf{q}(t))$$

where v are the measured values. Ideally, $v(s_i, t) = v_{true}(s_i, t)$.

However, deriving this function may be challenging because $\widehat{\mathcal{F}}$ depends on $\mathbf{q}(t_{calibration})$, where $t_{calibration}$ is the instant of the calibration. If there is no influence quantity, $\widehat{\mathcal{F}}^{-1}$ can be determined without any problem. On the opposite case, $\widehat{\mathcal{F}}^{-1}$ should be determined on the entire space of \mathbf{q} , so that with $\widehat{\mathbf{q}}$, an estimator of the values of the influence quantities, we can have:

$$v(s_i, t) = \widehat{\mathcal{F}}^{-1}(\zeta(s_i, t), \widehat{\mathbf{q}}(t)) = \widehat{\mathcal{F}}^{-1}(\mathcal{F}(v_{true}(s_i, t), \mathbf{q}(t)), \widehat{\mathbf{q}}(t))$$
(1.1)

Obtaining this relationship over the entire space of \mathbf{q} is difficult, without considering that the uncertainties also have to be derived. This is why there are two steps in a calibration procedure, e.g. the determination of $\widehat{\mathcal{F}}$ and then of $\widehat{\mathcal{F}}^{-1}$, the first step alone being often perceived as calibration. $\widehat{\mathcal{F}}$ and $\widehat{\mathcal{F}}^{-1}$ are often determined under a given finite set of operating conditions. The same reasoning can be extended to any type of expression chosen for the results of calibration (curve, diagram, table...).

3.3 In situ calibration

Calibration is an operation having a major impact on the quality of the measurements performed by an instrument. Indeed, the output of calibration is a relationship enabling to derive measurement results from indications, e.g. measured values and their associated uncertainties. Through this step, making the budget of the uncertainties on the indications is a prerequisite, and if there are influence quantities at stake, they must be taken into account. Thus, from its definition and all its constraints, it is justified to usually carry out calibration operations in dedicated facilities. This is manageable for few tens of instruments and corresponds to what is done by air quality monitoring agencies for instance [3] but it may be less the case in a deployment involving hundreds of nodes, particularly if they are low-cost. Indeed, these nodes are more prone to drift, requiring more frequent calibrations. Also, it was observed in some works [27, 124] that even if a laboratory calibration is performed, such instruments may behave differently once deployed in the field due to the uncontrolled operating conditions. In this context, maintaining a dense deployment of instruments for environmental sensing, with a significant number of low-cost devices is a tremendous task with classical approaches.

This is why research works emerged on *in situ* calibration algorithms, mostly in the past decade.

Definition 19 (In situ calibration algorithms). In situ calibration algorithms aim at calibrating measuring instruments while leaving them in the field, preferably without any physical

⁶Quantities involved in the operating conditions of the calibration can be influence quantities too.

intervention.

They are also called *field* [64], in place [24] or online [150] calibration algorithms.⁷

The literature on the subject studies under which hypotheses, in which manner and with which performances the measurement accuracy of one or several instruments can be improved with such algorithms. Notably in the context of a dense deployment of measuring instruments, comparisons of the measurement results of different instruments could be carried out for instance to perform this task and this is exactly what these algorithms aims at exploiting in practice.

Most of the time, in situ calibration strategies do not take indications as an input but measured values. Mathematically speaking, considering the developments presented in Section 3.2, it means it is not the function $\widehat{\mathcal{F}}$ that is derived but a function \mathcal{H} mapping what are considered as standard values and potentially influence quantity values to measured values so that:

$$\mathcal{H}: (v_{standard}(s_i, t), \mathbf{q}(t)) \to v(s_i, t)$$

Then \mathcal{H}^{-1} is derived and, in the same way that we expressed concerns regarding the determination of $\widehat{\mathcal{F}}^{-18}$, the corrected values $v_{corr}(s_i, t)$ that we obtain after an *in situ* calibration can be expressed as:

$$v_{corr}(s_i, t) = \widehat{\mathcal{H}}^{-1}(v(s_i, t), \widehat{\mathbf{q}}(t))$$
(1.2)

If the in situ calibration is perfect, then $v_{corr}(s_i, t) = v_{true}(s_i, t)$.

In summary, in situ calibration strategies aim at giving relationships to obtain **corrected** measured values from measured values. To do so, it uses standards defined with the measured results of other measuring instruments. The measurement uncertainties associated with the measured value measured both by the instruments to calibrate and the ones used to provide standard values can be used to derive the uncertainties of the corrected values. Finally, the operating conditions may be known through instruments measuring other quantities. Thus, in situ calibration is a procedure conceptually close to the formal definition of calibration.

3.4 Discussion

Deploying a large number of instruments in an environment may offer opportunities to overcome the problem of calibration for low-cost measuring instruments and more generally the issue of data quality. Indeed, the devices deployed can be related in a network, allowing them to share information like their measurement results.

However, regarding the formal definition of calibration and depending on the point of view, using results from instruments obtained in an uncontrolled environment to calibrate other devices may not be considered as equivalent to employing measurement standards or to being under controlled conditions. This is particularly true if results from an instrument of an equivalent accuracy class than another is used to calibrate another device.

Definition 20 (Accuracy class [14]). The accuracy class (or class) is the "class of measuring instruments or measuring systems that meet stated metrological requirements that are intended to keep measurement errors or instrumental measurement uncertainties within specified limits under specified operating conditions".

Indeed, values in which a high confidence can be put are used to calibrate measuring instruments. Thus, the sources of the measurement standards and the conditions under which an algorithm is applied are critical for *in situ* calibration strategies.

⁷Calibration without any adjective is also used but it may be confusing regarding the definition of calibration in the VIM [14].

⁸e.g. it should be carried out on the entire space of \mathbf{q} , with an estimator $\hat{\mathbf{q}}$ of the values of the potential influence quantities.

4 Sensor networks

We define in this section what a sensor network is. Note that we do not discuss (wireless or not) sensor networks in general here. We focus on its characteristics of interest from the perspective of *in situ* calibration of measuring instruments.

4.1 Definition

Definition 21 (Sensor network [62]). A sensor network⁹ is a set of measuring systems spatially deployed in order to periodically measure one or more quantities in an environment and that are exchanging information between them through a communication layer that can be wired or wireless.

The measuring systems of the network are also called **nodes**. A node may be static or mobile. The network can be either meshed, with device to device communications, or a collection of stars centred on gateways for instance. A sensor network may target a single or multiple measurands.

For a given measurand, instruments in a same network are not necessarily of the same accuracy class. In this work, those which are known to be more accurate than the others of the network are called **reference instruments**.

It is also supposed that the nodes have sufficient computing and communication capabilities to carry out the operations proposed by algorithms to overcome the issue of their *in situ* calibration.

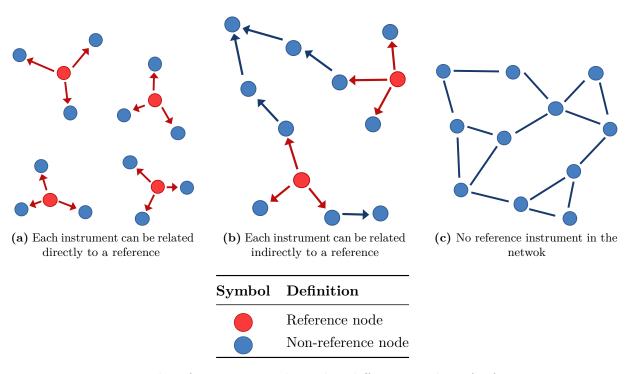


Figure 1.4.1: Examples of sensor networks with a different number of reference instruments and different relationships between the instruments.

4.2 Characteristics of interest in this work

Regarding the issue of data quality for low-cost instruments, there are two characteristics of sensor networks particularly interesting to tackle this issue: the presence of reference instruments

⁹Like sensor being used to refer to a measuring instrument, there is also a slight misuse of language here. However, this expression is widely used.

for each measurand and the mobility of the nodes.

4.2.1 Presence of references

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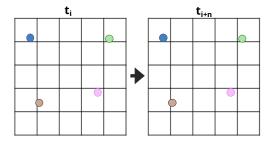
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In sensor networks, reference instruments may be present in different manners. First of all, it is possible that each non-reference node can compare its measurement results directly to those of a reference one. It is illustrated in Figure 1.4.1a where the edges represent the possible comparisons. They can be defined following a criterion of distance between the nodes.

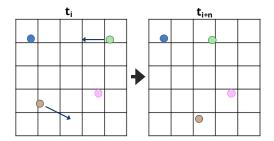
Another possibility is that each instrument can be related to a reference node but in this case there may be intermediary nodes for some instruments as shown in Figure 1.4.1b.

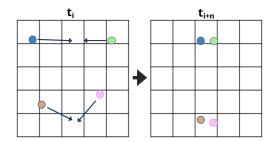
Finally, the last possible case is that there is no reference instrument in the network. This is the example of Figure 1.4.1c. Edges were drawn representing relationships between non-reference instruments.

In the case of nodes that are measuring systems which do not measure the same quantities, it is possible that the graph of relationship between instruments is different depending on the measurand.



(a) Each instrument is static





(b) There are both static and mobile instruments

(c) Each instrument is mobile

Figure 1.4.2: Illustration of the principle of static, static and mobile, and mobile sensor networks.

4.2.2 Mobility

We already mentioned that measuring instruments deployed for environmental monitoring can remain at a fixed position, e.g. they are static, or they can change of position over time, e.g. they are mobile.¹⁰ Thus there are three different situation that can happen, regardless of the accuracy class of the instruments:

- all the nodes of the network are static.
- all the nodes are mobile.

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 $^{^{10}}$ The case of we arable instruments is considered as a particular case of mobile instruments

• there are both static and mobile nodes in the network.

This is illustrated in Figure 1.4.2.

Note that in the presence of mobile nodes, the relationships between reference and non-reference instruments as exposed in the previous section may change overtime.

4.3 Conclusion

In this section, we defined what a sensor network is in a general manner. Two characteristics of sensor network were presented respectively regarding the presence of reference instruments in the network and the mobility of the nodes. The different possible cases for each property were exposed. They are interesting for the problem we want to address because the configuration of the sensor network can be exploited to drive the design of *in situ* calibration algorithms.

5 Problem statement

Based on the previous sections, we now detail the problem statement of this thesis briefly exposed above in the General Introduction.

5.1 Motivations

To perform a dense measurement of ambient quantities in the context of environmental monitoring activities such as air quality, the use of low-cost instruments is necessary to achieve it at a reasonable cost economically speaking.

Among all the challenges for the measurement of ambient quantities in the context of environmental monitoring, the need for an improvement of data quality appears to be major, particularly for low-cost devices. Indeed, they are prone to faults, one of them being the instrumental drift.

The traditional solution to mitigate this fault is to regularly perform calibration, as part of maintenance operations. However, the cost of these tasks may limit the feasibility of dense deployments of measuring instruments in practice. Indeed, it usually requires taking out the node of the service, bringing it to a calibration facility, calibrating it and putting it back in service. Therefore, there is a need for tools that allow recalibrating measuring instruments while leaving them on the field.

Because the aim of environmental monitoring with low-cost instruments is notably to have a high density of instruments in the monitored area, it is tempting to use the measurement results of other instruments to detect and potentially correct drift faults. The sensor network thus constituted may offer opportunities to overcome this issue of data quality. In addition, the presence of reference instruments in the network, or the mobility of the nodes are characteristics that can facilitate the *in situ* calibration of measuring instruments.

Fortunately, researchers have begun to tackle this issue and there are already methodologies exploiting the properties of sensor networks to enable an *in situ* calibration of measuring instruments.

5.2 Objectives of the thesis

As stated in the General Introduction, the objective of this thesis is to contribute to the research effort on the improvement of data quality for low-cost measuring instruments through their *in situ* calibration.

In the first place, the goal is to identify the existing techniques in the literature enabling such a calibration. In Section 4, we reported there are sensor networks with different characteristics in terms of presence of reference instruments and of mobility of the nodes. Thus, the same *in situ* calibration algorithm may not be applicable to each type of sensor network for instance. This is

why a taxonomy is developed for the description of such algorithms, facilitating the identification of the strategies that could be applied to a given sensor network depending on its properties and the characteristics of its instruments.

In the case where several *in situ* calibration algorithms are adapted to a sensor network, it is possible that each strategy does not yield the same performances. In other words, they may not all allow correcting instruments as well as possible. Being able to evaluate the performances of different *in situ* calibration strategies on the same use case would be valuable. The second objective of this thesis is to provide a means to achieve this task.

Finally, before calibrating measuring instruments, it is necessary to determine if an instrument needs to undergo such an operation. In publications introducing *in situ* calibration algorithms, they mainly focus on how to perform the calibration itself and less on how to identify the faulty instruments in a sensor network. Identifying these instruments is the last objective of this thesis.

Chapter 2

In Situ Calibration Algorithms for Environmental Sensor Networks

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Introduction

In Chapter 1, we recalled that calibration is a procedure which is part of the maintenance operations that must be carried out periodically on measuring instruments. To follow the traditional guidelines for a device deployed in the field, it requires to take it out of service and to bring it to a calibration facility. After this operation and the adjustment of the instruments, it can be brought back to its deployment site and put in service again. This is manageable for few tens of instruments and corresponds to what is done by air quality monitoring agencies for instance [3]. However, notably in the particular case of environmental sensing, the increase of the number of measuring instruments deployed to ensure a higher spatial coverage may not be feasible due to the cost of regulatory-grade instruments. This is why low-cost instruments are necessary to achieve this goal. In the literature, it is reported they are suffering from multiple issues, one of them being a faster drift than usual instruments. Regarding calibration, it forces to increase the frequency of this maintenance operation and thus its cost, economically and technically. Also, it was observed in some works [27, 124] that even if a laboratory calibration is performed, once deployed in the field, the instruments may behave differently due to the uncontrolled operating conditions. In this context, maintaining a sensor network for environmental sensing and composed of a significant number of low-cost instruments with classical approaches is a tremendous task.

This is why research works emerged on *in situ* calibration algorithms, mostly in the past decade. The literature on the subject is abundant. This is expected as the topic of sensor networks is a very hot research subject due to the trend for Smart Cities and the Internet of Things. More practically, it is also predictable due to the variety of sensor networks as described in Chapter 1, Section 4. Indeed, one can easily suspect that an algorithm may not be applicable to any type of sensor network, on top of the assumptions surrounding it.

A major gap is the absence of a unified terminology to describe the field of application of *in situ* calibration algorithms, regardless of the quantities measured by the sensor network. The large variety of different techniques are reported under different terms in the literature, such as "blind calibration", "multi-hop calibration", "macro calibration" and so on. The expression describing an algorithm, notably in the title or the abstract of publications, is sometimes not unique and is often incomplete. For instance, Balzano *et al.* [7] or Wang *et al.* [168] present their strategies as "blind calibration" algorithms but it does not indicate if it is applicable to static sensor networks, mobile ones or both. Consequently, it is not crystal-clear for an end user from the first lines of an article if its content matches his/her needs.

Thus, through the literature review carried out on *in situ* calibration algorithms in this thesis, a synthetic taxonomy for their classification is proposed in the following sections.

1 Scope of the taxonomy

To build the taxonomy, several guidelines were adopted. First of all, it should help to identify relevant algorithms for any type of sensor network.

Also, the terms must be independent from the kind of measurand of the sensor networks: the groups of categories described are relevant for any environmental phenomenon.

The attributes should also be limited to the properties involved in the definition of *in situ* calibration, e.g. the definition of the measurement standards, how the relationship between measured values and standard values is established and the relationship correcting the measured values.

There are other considerations that could matter and interact with calibration, such as data integrity, security or even privacy, in particular when crowdsensing platforms involving citizens are developed [102]. The way the calibration relationship is computed, e.g. in a centralised or decentralised manner is also a major subject. As a matter of fact, the care for the management of these questions at an early step of the design of the sensor networks is crucial. The way these

concepts are implemented with respect to the system architecture may have a significant impact on the effectiveness of some calibration methods.

Nevertheless, they are not attributes that are directly driving the result of the calibration procedure compared to the network architecture and the algorithmic principles underlying the calibration method. Thus they are not in the scope of the presented taxonomy but defining characteristics that should be clarified in a second step to describe an *in situ* calibration strategy as part of a "datasheet" could be valuable for future works.

2 Taxonomy for the classification of the algorithms

The following subsections introduce the proposed taxonomy for the classification of *in situ* calibration strategies for sensor networks.

They can be divided into two categories: network architecture characteristics, namely the nature of instruments and their potential mobility, and the algorithmic principles of the calibration techniques, namely the mathematical structure of the calibration relationship and to which point the algorithm can be distributed.

Each subsection represents a primary level group of categories that may have others nested. Categories for each group are in bold font. Figure 2.2.1 represents all the primary levels of the taxonomy and the defined categories.

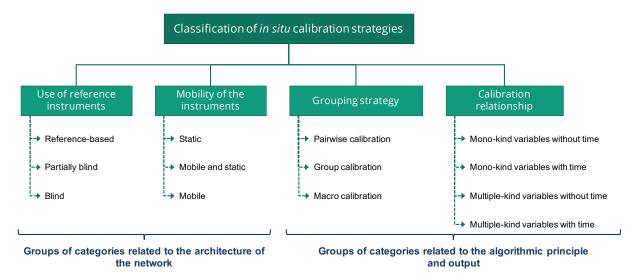


Figure 2.2.1: Proposed taxonomy for the classification of in situ calibration algorithms

2.1 Use of reference instruments

One of the first criteria of classification is how the calibration method assumes the presence of reference instruments within the network. This is why this characteristic was introduced as a major feature of sensor networks for our problem in Chapter 1 Section 4.2.1. Three categories of *in situ* calibration strategies are derived from this property with the examples from Figure 1.4.1 in Chapter 1 to illustrate them.

The calibration of measuring instruments using a sufficient number of reference measurement instruments is called **reference-based calibration**. It means the network is composed of both reference and non-reference instruments and that all the non-reference instruments can be calibrated using at least one reference instrument. The approach postulates the existence of a calibration relationship between each non-reference instrument and at least one reference instrument because they are close enough for instance.

The calibration of measuring instruments in the absence of reference values is called **blind calibration**. It means the network is composed of only non-reference instruments. These various methods may or may not assume the existence of a correlation between the instrument outputs.

The hybrid situation is called **partially blind calibration**. In this setting, the network may gather both reference and non-reference instruments, but a reference-based calibration is not achievable, e.g. when some of the non-reference instruments can never be compared to a reference instrument. It also captures the cases where a cluster of non-reference instruments is considered as good enough to approximate a reference instrument.

2.2 Mobility of the instruments

The second significant aspect of the network architecture is the potential mobility of nodes. As for the presence of references, this characteristic was also introduced as a major feature of sensor networks in Chapter 1 Section 4.2.2. In the same way, three categories of *in situ* calibration strategies are considered to illustrate how the methods exploit this feature. Figure 1.4.2 in Chapter 1 provides examples also illustrating these categories.

A first category of methods addresses networks with **exclusively static** nodes. A second one addresses networks with **exclusively mobile** nodes. The corresponding methods rely often strongly on the mobility of the nodes to achieve calibration. A last group of methods addresses heterogeneous networks with **both mobile and static** nodes. In such cases the mobility of the nodes is not systematically exploited in the calibration strategy.

2.3 Calibration relationships

In Chapter 1 Section 3, the purpose of calibration was defined as the establishment of a mathematical relationship between the indications of the instrument and the standard values of the measurand and then the derivation of a relationship for obtaining measurement results from indications. Based on the mathematical developments in Chapter 1 Section 3.2, it is possible to have different types of relationships, for instance if there are influence quantities or not.

The categories identified in this section are first based on the number of kinds of quantities as input variables in the relationship: the measurand, the indications, the influence quantities, and so on. In terms of algorithmic principles of the calibration methods, it implies the variety and quantity of data to exchange as well as the computational effort necessary to achieve a target accuracy.

The most straightforward relationships are called **mono-kind variables without time**. They only take a single quantity as input variable and do not depend on time.

The second category of relationships gathers the ones that have **mono-kind variables with time**. It accounts for a relationship with mono-kind variables which is influenced by time, for instance in case of an instrument drifting due to ageing [160]. This is the reason why time is considered as a particular influence quantity.

The relationships with **multiple-kind variables without time** account for two or more quantities as variables but remain independent from time. These models are mainly used to include the effect of influence quantities (except time) in the calibration relationship. In these cases, the networks include instruments measuring the influence quantities. They are not systematically reference instruments and therefore their calibration may also be included in the calibration strategy.

Finally, this last approach may be extended into relationships with **multiple-kind variables** with time when appropriate.

For each of these categories, subcategories can be defined based on the kind of mathematical expression used for the calibration relationship. Popular examples are polynomials with constant coefficients [7], gain-phase [13], variable offset [168], neural networks [36].

These categories are particularly important regarding *in situ* calibration strategies. Indeed, considering Equations 1.1 and 1.2 in Chapter 1 Section 3, we have then:

$$v_{corr}(s_i, t) = \widehat{\mathcal{H}}^{-1}(\widehat{\mathcal{F}}^{-1}(\zeta(s_i, t), \widehat{\mathbf{q}}(t)), \widehat{\mathbf{q}}(t))$$

$$= \widehat{\mathcal{H}}^{-1}(\widehat{\mathcal{F}}^{-1}((\mathcal{F}(v_{true}(s_i, t), \mathbf{q}(t)), \widehat{\mathbf{q}}(t)), \widehat{\mathbf{q}}(t)))$$
(2.1)

For example, and recalling that in fine the purpose of in situ calibration is to determine the relationship $\widehat{\mathcal{H}}^{-1}$, assume that the calibration relationship $\widehat{\mathcal{F}}^{-1}$ of an instrument is defined so that:

$$v(s_i, t) = G \cdot \zeta(s_i, t) + O + \kappa \cdot \theta(t)$$

where G, O and κ are constant and G is a correction of the gain of the instrument, O a correction of the offset, and $\kappa \cdot \theta(t)$ a temperature compensation standing for $\hat{\mathbf{q}}(t)$ in the general expression of $\hat{\mathcal{F}}^{-1}$.

If the behaviour of the measuring instrument changes over time, $\widehat{\mathcal{F}}^{-1}$ cannot ensure that $v(s_i,t) = v_{true}(s_i,t)$. If the calibration relationship $\widehat{\mathcal{F}}^{-1}$ cannot be adjusted, the relationship $\widehat{\mathcal{H}}^{-1}$ obtained with an *in situ* calibration strategy may compensate the change of behaviour. Consider $\widehat{\mathcal{H}}^{-1}$ such as:

$$v_{corr}(s_i, t) = G_{corr} \cdot v(s_i, t) + O_{corr} + \kappa_{corr} \cdot \theta(t)$$

Therefore:

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$$v_{corr}(s_i, t) = G_{corr}G \cdot \zeta(s_i, t) + (G_{corr}O + O_{corr}) + (G_{corr}\kappa + \kappa_{corr}) \cdot \theta(t)$$

Here, $\widehat{\mathcal{H}}^{-1}$ allows determining a calibration relationship equivalent to the type of $\widehat{\mathcal{F}}^{-1}$. However, if $\widehat{\mathcal{H}}^{-1}$ is such as:

$$v_{corr}(s_i, t) = G_{corr} \cdot v(s_i, t) + O_{corr}$$

Then:

$$v_{corr}(s_i, t) = G_{corr}G \cdot \zeta(s_i, t) + (G_{corr}O + O_{corr}) + G_{corr}\kappa \cdot \theta(t)$$

In this case, it may not be possible to correctly compensate the drift of the instrument s_i because $G_{corr}G$ and $G_{corr}\kappa$ are coupled: it is not possible to correct them independently.

This is why the type of relationship, both considering the quantities at stake and the kind of relationship between them, is a key parameter of *in situ* calibration algorithms.

2.4 Instrument grouping strategies

While the previous categories are mostly driven by operational constraints (deployment strategy, properties of the measurand and of the selected instruments), the present paragraph considers the number of nodes involved in each calibration step and to which point the algorithm can be distributed.

A first approach is **pairwise calibration**. Two instruments are used, one providing standard values for the other. It is classically applied between a reference instrument (or an approximation of reference) and each of the nodes related to it. It can be a distributed or even localised algorithm. This case is illustrated in Figure 2.2.2a.

A macro calibration strategy consists in calibrating the network as a whole, e.g. the values of all the instruments are used to calibrate each node of the network. Even if they exist, node-to-node relationships may not be exploited directly. A centralised algorithm might be necessary with this grouping strategy. This case is illustrated in Figure 2.2.2c.

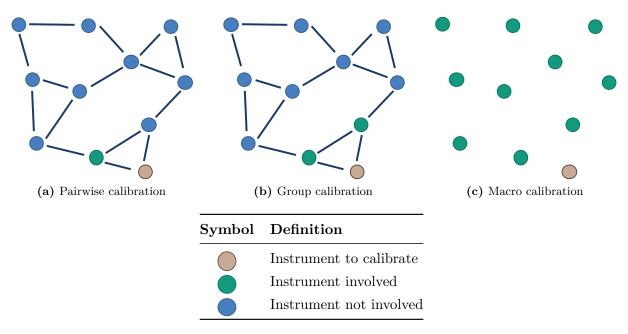


Figure 2.2.2: Examples of cases for grouping strategies from the point of view of an instrument to calibrate. It shows which instruments are used to calibrate it. In these figures, the sensor networks are equivalent in terms of the number of instruments and relationships between the instruments. The relationships between the instruments in Figure 2.2.2c are not represented because they may not be exploited directly.

Group calibration is an intermediate approach consisting in carrying calibration operation among groups of measuring instruments among the whole network. In this case, the criteria defining these groups become essential. This approach may be used when pairwise calibration induces significant errors, while macro calibration approaches are not fine adjusted enough. This category includes notably strategies where groups are composed of instruments measuring additional quantities besides the main target quantity. These additional quantities are often included as influence quantities in the calibration relationship. These algorithms can be at least partially distributed, e.g. the computation is concentrated on an elected group leader, or fully distributed at the cost of messages broadcasting. This case is illustrated in Figure 2.2.2b.

3 Comparison to other taxonomies

During the development of the presented taxonomy, two notable works reviewing calibration algorithms for sensor networks were conducted and published within a few months.

The first publication is from Maag et al. [95] and focus on sensor networks for air pollution monitoring, addressing operational concerns regarding to calibration. Maag et al. build their classification of the algorithms around the terms "blind calibration", "collaborative calibration", "blind and collaborative calibration" and "transfer calibration". They add precision regarding the mobility of the instruments in the publications they cite. Without going into details of the definitions of the terms, it is a taxonomy with few attributes: blind group calibration and blind macro calibration strategies cannot be distinguished with this taxonomy for instance. This information on how the instruments are grouped to perform a calibration is important, notably if a distributed algorithm is targeted. This is why this feature was added in the presented taxonomy.

The second publication is from Barcelo-Ordinas et al. [8]. The taxonomy they developed

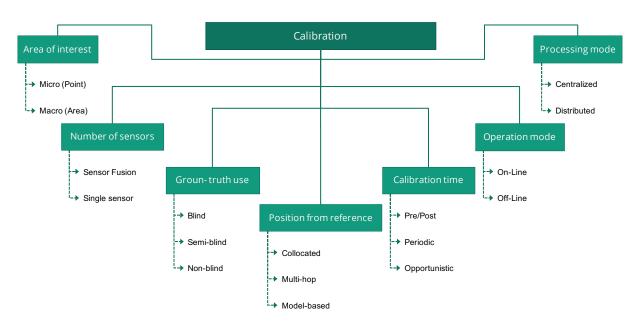


Figure 2.2.3: Taxonomy for the classification of self-calibration algorithms according Barcelo-Ordinas *et al.* [8]

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is more detailed. All the categories of attributes they proposed are represented in Figure 2.2.3. Some of them are common with the presented taxonomy, like the use of reference instruments which is called the "ground-truth use". There are also differences. First, two groups of categories are close to what was defined here as the grouping strategy in the presented taxonomy: the "area of interest" and the "number of instruments". However, the associated definitions and the number of categories in these groups do not strictly match the ones given in Section 2. Also the idea of using reference instruments is refined in this taxonomy with a "position from reference" group of attributes, meaning the authors do not consider a relationship between the use of reference instruments and the category "position from reference". Barcelo-Ordinas et al. also provided other groups of attributes like the "calibration time", the "operation mode" or the "processing mode". The processing mode was stated as not in the scope of the presented taxonomy in Section 1. The "calibration time" (pre/post, periodic, opportunistic) and the "operation mode" (on-line, off-line) can have an influence on the data quality of the instruments. Indeed, two algorithms with the same attributes except one that performs a periodic calibration and the other an opportunistic one may not give the same results. However, we estimate that such features should be considered in a second phase while identifying relevant algorithms considering a given sensor network. For this reason, these features were not added to the presented taxonomy, contrary to Barcelo-Ordinas et al. Finally, contrary to the work of Maag et al. and the work presented here, Barcelo-Ordinas et al. did not consider the mobility of the instruments of the network as an attribute of in situ calibration algorithms.

In both of these works, the mathematical expression of the calibration relationship is not directly an attribute of the algorithms. This topic is discussed but separately from others like the presence of reference in the network. In addition, calibration relationships are considered from the point of view of the mathematical expression only in both publications. The question of the calibration relationship is more developed in the presented taxonomy. The classification of calibration relationships is decomposed in two levels in our work: first the kinds of quantities involved in the relationships and then their mathematical expression. In this way, the proposed taxonomy provides a description capturing more information on the calibration relationships in in situ calibration algorithms.

To conclude, the presented taxonomy is halfway between the one of Maag et al. which is based on few groups of attributes, and the one of Barcelo-Ordinas et al. that is more extensive but with information excluded from the scope of this work. Our taxonomy is limited to what is seen as the essential attributes of an in situ calibration algorithm and designed with an effort to avoid any redundancy or misunderstanding from the terms used, the taxonomy of Barcelo-Ordinas et al. being sometimes confusing due to some groups of categories that are semantically close. Nevertheless, all these works are interesting basis towards the definition and use a shared terminology.

4 Review of the literature based on this classification

An application of the classification is provided here with highlights on the existing literature. A large number of *in situ* calibration studies according to this classification are sorted in Table 2.4.1. Some rows refer to multiple papers as they are related somehow (same technique or same authors) and consist in developments of the same initial paper. The current section focuses on a description of the methods. The topic of performance comparison between methods is addressed in the next section. The addressed measurands cover a wide range of environmental quantities: temperature [7], pressure [182], noise [129], air pollutants [63], light [163]... Most of the reported studies have generic approaches that can be transposed to other measurands.

4.1 Overview

Regarding pairwise strategies, relatively few papers address methodological issues related to reference-based pairwise strategies, as this approach is the closest to a "traditional" calibration approach with measurement standards and features fewer challenges. Partially blind and blind pairwise calibration methods (often focusing on mobile nodes) are more complex as they require to define calibration relationships not only between reference and non-reference nodes, but also between non-reference nodes only. This translates into error propagation issues.

Macro calibration approaches were initially developed to address the absence of reference instruments in a network and thus are mostly blind or partially blind. In the absence of reference, there is a strong challenge in defining valid calibration relationships based on values from non-reference instruments, which explains the major consideration for these methods.

Group strategies have been generating strong interests as they appear to outperform both pairwise and macro strategies with or without reference instruments.

Most methods are based on relationships with mono-kind variables without time and with a linear expression, but more complex models are progressively appearing to better address the complexity of environmental sensing. Moreover, a particular attention is given in most publications to the calibration relationship used compared to the other attributes of *in situ* calibration that we identified. This can be explained by the significance of this information for readers, e.g. it allows determining if the algorithm is able to provide a correction corresponding to the way an instrument drifts.

Likewise, while most works initially focused on static networks, there are now many interests for mobile nodes as they allow for physical rendez-vous between nodes. Henceforth, calibration methods are less impacted by the physical variability of the phenomena.

Finally, an underlying question addressed is the ability to distribute the computation of calibration relationships [23, 55, 107, 129, 143, 144, 177]. The topic is of strong interest when considering privacy preservation issues [102]. The capability to decentralise is linked to the grouping strategy: pairwise and group strategies foster more naturally decentralised computation, under the condition that the nodes are capable of individual procession and of bidirectional communication. On the contrary, macro-calibration strategies tend to be centralised, except when the characteristics of the parameter identification methods allow for partially or fully

Table 2.4.1: In situ calibration strategies for sensor networks R: Reference-based, PB: Partially blind, B: Blind, P: Pairwise, Gr: Group, Ma: Macro, 1V: Mono-kind variables without time, 1VT: Mono-kind variables with time, MV: Multiple-kind variables with time, MV: Multiple-kind variables with time, M: Exclusively mobile, S: Exclusively static, MS: Mobile and static

Papers	Availability of reference instruments			Instruments grouping strategies			Kinds of variables in calibration relationships				$\begin{array}{c} \text{Mobility} \\ \text{of the} \\ \text{instruments} \end{array}$		
	R	РВ	В	P	Gr	Ma	1V	1VT	MV	MVT	M	S	MS
Ramanathan et al. [123]	√	-	-	✓	-	-	✓	=	-	-	-	√	-
Miluzzo et al. [107]	\checkmark	-	-	-	\checkmark	-	\checkmark	-	-	-	\checkmark	_	\checkmark
Deshmukh et al. [39]	\checkmark	-	-	\checkmark	-	-	\checkmark	-	-	-	-	\checkmark	-
Spinelle <i>et al.</i> [139, 140]	\checkmark	-	-	\checkmark	\checkmark	-	\checkmark	-	\checkmark	-	-	\checkmark	-
Moltchanov et al. [109]	\checkmark	_	_	✓	_	_	\checkmark	_	-	_	_	✓	_
Gao et al. [58]	✓	-	_	\checkmark	_	_	\checkmark	_	\checkmark	-	_	\checkmark	_
Lin et al. [88]	\checkmark	_	_	✓	_	_	\checkmark	_	-	_	_	✓	_
Fang et al. [50, 51]	\checkmark	_	-	_	\checkmark	_	_	_	\checkmark	-	_	✓	_
Martin et al. [103]	√	✓	_	_	✓	_	_	_	✓	_	_	√	_
Sun et al. [147]	√	_	_	_	✓	_	_	_	✓	_	_	√	_
Zimmerman et al. [186]	√	_	_	_	✓	_	_	_	✓	_	_	√	_
Weissert et al. [108, 171]	√	_	_	_	√	_	√	_	_	_	_	√	_
Yu et al. [184]	√	_	_	_	√	_	_	_	✓	_	_	√	_
Topalovic et al. [155]	✓	_	_	✓	✓	_	√	_	✓	_	_	_	_
Wei et al. [169, 170]	✓	_	_	_	✓	_	_	_	✓	_	_	√	_
Mahajan <i>et al.</i> [96]	·	_	_	✓	· ✓	_	√	_	· ✓	_	_	· ✓	_
Malings et al. [100]	·	_	_	_	·	_	_	_	· ✓	_	_	· ✓	_
Qin et al. [120]	·	_	_	_	·	_	_	_	· ✓	_	_	_	✓
Barcelo-Ordinas et al. [9, 10, 53]	✓	-	-	-	✓	-	-	-	✓	-	-	✓	- -
Wang et al. [165]	√	_	_	_	√	_	_	_	✓	_	_	\checkmark	_
Loh et al. [91]	√	_	_	✓	_	_	✓	_	✓	_	_	√	_
Badura et al. [6]	√	_	_	✓	√	_	√	_	✓	_	_	✓	_
Cordero et al. [34]	✓	_	_	_	✓	_	_	_	✓	_	_	✓	_
Tsujita et al. [156]	_	√	\checkmark	✓	_	_	√	_	_	_	✓	_	\checkmark
Tsujita et al. [157]	_	✓	✓	✓	√	_	_	_	✓	_	_	√	_
Y. Xiang et al. [177]	_	✓	\checkmark	✓	_	_	√	_	_	_	✓	_	√
Hasenfratz <i>et al.</i> [63], Saukh <i>et al.</i> [131, 132]	-	✓	✓	✓	-	-	✓	-	-	-	✓	-	✓
Fu et al. [57]	_	√	√	✓	_	_	√	_	_	_	✓	_	√
Maag et al. [93, 94]	_	·	·	_	√	_	-	_	√	_	·	_	·
Arfire et al. [4]	_	·	·	_	·	_	1	✓	· ✓	✓	·	_	· ✓
Markert et al. [102]	_	·	↓	1	, ,	_	· ./	-	-	-	· ./	_	· ./
Kizel et al. [75]	_	·	· -	,	-	_	,	_	_	_	_	1	-
Sailhan et al. [129]	_	-	- ✓	,	1	_	,	_	_	_	1	-	_
Fonollosa et al. [56]	_	1	√	./	_	_	./	_	_	_	_	1	_
Whitehouse et al. [172]	_	-	∨	• -	_	- ✓	./	_	_	_	_	,	_
Thler et al. [65]	_	- ✓	∨ ✓	-		v	√	-	-	_	-	v	-
Taylor et al. [152]	-	٧	∨	-	٧	- ✓	∨ ✓	-	-	-	-	v ./	-

continued on next page

Papers	Availability of reference instruments		Instruments grouping strategies			Kinds of variables in calibration relationships				$\begin{array}{c} \text{Mobility} \\ \text{of the} \\ \text{instruments} \end{array}$			
	R	РВ	В	Р	Gr	Ma	$\overline{1V}$	1VT	MV	MVT	M	S	MS
Tan et al. [151]	-	-	√	-	✓	-	✓	-	-	-	-	√	-
Bychkovskiy et al. [24]	-	\checkmark	\checkmark	\checkmark	-	-	\checkmark	-	-	-	-	\checkmark	-
Balzano et al. [7], Lipor et al. [89], Dorffer et al. [44]	-	✓	✓	-	-	✓	✓	-	-	-	-	✓	-
Takruri <i>et al.</i> [148–150]	-	-	\checkmark	-	-	\checkmark	\checkmark	-	-	-	-	\checkmark	-
Kumar et al. [76, 77]	-	-	\checkmark	-	\checkmark	-	\checkmark	-	-	-	-	\checkmark	-
Ramakrishnan et al. [122]	-	-	\checkmark	-	-	\checkmark	\checkmark	-	-	-	-	\checkmark	-
Buadhachain et al. [23]	-	-	\checkmark	-	\checkmark	\checkmark	\checkmark	-	-	-	\checkmark	-	\checkmark
Bilen et al. [13]	-	-	\checkmark	-	-	\checkmark	\checkmark	-	-	-	-	\checkmark	-
Cambareri et al. [25]	-	-	\checkmark	-	-	\checkmark	\checkmark	-	-	-	-	\checkmark	-
C. Wang <i>et al.</i> [161–163]	-	-	\checkmark	-	-	\checkmark	\checkmark	-	-	-	\checkmark	-	-
C. Xiang <i>et al.</i> [176]	-	-	\checkmark	-	-	\checkmark	\checkmark	-	-	-	\checkmark	-	-
Dorffer et al. [41–43, 45]	-	\checkmark	\checkmark	-	-	\checkmark	\checkmark	-	-	-	\checkmark	-	-
Y. Wang <i>et al.</i> [166–168] Li [87]	-	-	\checkmark	-	-	✓	-	\checkmark	-	-	-	✓	-
Ye et al. [182]	-	-	\checkmark	\checkmark	-	-	\checkmark	-	-	-	\checkmark	-	-
De Vito <i>et al.</i> [36, 47]	-	-	\checkmark	-	\checkmark	-	-	-	\checkmark	-	-	\checkmark	-
Son et al. [137], Lee [82]	-	-	\checkmark	-	-	\checkmark	\checkmark	-	-	-	\checkmark	-	-
Stankovic <i>et al.</i> [143, 144]	-	-	\checkmark	-	-	\checkmark	\checkmark	-	-	-	-	\checkmark	-
Fishbain et al. [55]	-	\checkmark	\checkmark	-	\checkmark	-	\checkmark	-	-	-	-	\checkmark	-
Popoola et al. [118]	-	-	\checkmark	\checkmark	-	-	-	-	\checkmark	-	-	\checkmark	-
Yan et al. [180]	-	-	\checkmark	-	\checkmark	-	-	-	\checkmark	-	-	\checkmark	-
Yang et al. [181]	-	-	\checkmark	-	-	\checkmark	\checkmark	-	-	-	-	\checkmark	-
Mueller et al. [113]	-	-	\checkmark	-	\checkmark	-	-	-	\checkmark	-	-	\checkmark	-

decentralised computation. However, while distributed computing impacts the computational performance of algorithms, there is no report on how it affects calibration performances so far.

4.2 Mobile and static nodes

Static networks are more frequently studied than mobile ones. A wide range of solutions is now available to calibrate them. However, these calibration methods usually require a high spatial density of nodes to overcome the spatial variability of the phenomena, which is not always viable technically or economically. The availability of mobile nodes could alleviate this constraint, as calibration operations exploit physical rendez-vous between nodes. In turn, the methods based on this principle are challenged when the rendez-vous frequency is too low compared to the speed of degradation of the measurement accuracy [132]. In such cases, the addition of a few reference nodes seems to yield satisfying results [45, 63, 146]. Moreover, a challenge of mobile instruments is that they face rapid transients. To address this, methods initially developed for static networks appear promising, such as the work of De Vito et al. [36, 47] which uses dynamic and nonlinear

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Kim et al. [74] Chen et al. [30] Wu et al. [175] Becnel et al. [11] Cheng et al. [31] Sun et al. [146] supervised machine learning tools.

4.3 Calibration relationships

Most reported relationships are of mono-kind variables without time type and based on linear expressions. Nevertheless, there is a rising interest for models with multiple-kind variables, which stems from the observation that there are indeed significant influence quantities for various environmental measurands, notably air pollutant concentrations or temperature and humidity. It often depends on the technology of the instruments used [50, 68, 71, 136, 169, 170]. Such relationships gave very interesting results compared to simpler relationship models:

- for reference-based group calibration in [6, 9, 10, 34, 53, 120, 139, 140, 147, 155]
- for partially blind group calibration strategies in [4, 50, 93, 103], including with timesensitive models in [4, 23]
- for blind strategies, either pairwise or group based, in [47, 113, 118, 180].

On the contrary, relationships with multiple-kind variables were shown to be unnecessary in [22, 64] where the control of the operating temperature of the device was sufficient to perform a pairwise calibration without being influenced by this quantity. Malings *et al.* [100] reached a similar conclusion while also indicating that generalised models, e.g. calibration relationships built on all the values of the different measurands of a measuring system for all the instruments of the device, have several advantages: they reduce the effort required to calibrate the instruments because only one generalised relationship is derived instead of individual ones.

In general, time-dependent approaches are used to address drift issues. Drift is often modelled as an additive random variable with a given probability distribution [167, 177], so that drift-compensation translates as an offset correction.

4.4 Pairwise strategies

Reference-based pairwise

Relatively few papers address methodological issues related to reference-based pairwise strategies, as this approach is the closest to a "traditional" calibration approach with measurement standards. Primarily, reference instruments may be directly co-located in the field with non-reference instruments to achieve their calibration [6, 50, 88, 96, 123, 139, 140, 155].

However, more automated strategies are expected, requiring less the co-location of instruments. Nevertheless, even in the simple case of a relatively dense sensor network, the measurand may spatially vary too much in general to relate a reference instrument at a given location to an instrument at another location for calibration purposes. As an elementary solution to this, Moltchanov et al. [109] proposed to carry out calibration against the reference node using only the data collected during a specific time span based on the postulate that the phenomenon varies less during this time span. This was an idea previously developed by Tsujita et al. [157] including weather conditions used to correct the measured values but not with a reference-based approach. In the context of the Internet of Things, Loh et al. proposed a web-query based framework using machine learning to calibrate portable particulate matter measuring instruments. They reference values are obtained from governmental monitoring stations.

Partially blind pairwise

Partially blind pairwise calibration focuses mostly on mobile nodes. Tsujita et al. [156] tackled it first for mobile nodes by proposing that the device to calibrate should display either the value of a reference node that is close enough, or the average measurements between co-located

nodes if no reference node is available. A calibration parameter is adjusted with these values to correct measurements between rendez-vous.

Y. Xiang et al. [177] later proposed another method. They also distinguish calibration based on the values of a reference instrument and on the values of a non-reference instrument. Their originality relies in the correction of the values that is performed with an estimator of the drift error of the node. This error is recalculated at each calibration by minimising its variance according to a linear combination of the values of the instruments involved in the calibration process.

Hasenfratz et al. [63] addressed by various methods the case of calibration for mobile devices against reference instruments or not. They notably provided dedicated extensions for the case where some devices rarely encounter reference instruments. They also demonstrated a linear dependency between the measurement error and the number of intermediary calibrations between a given node and the reference node it is calibrated against. In [132] and [131], Saukh et al. proposed solutions to this issue of error accumulation by working on the occurrence of rendez-vous between nodes, in view of maximising the opportunities of calibration. An alternative idea was developed by Fu et al. [57] who proposed the optimisation of the deployment of reference instruments to ensure that all nodes can be calibrated against one of the references with a path no longer than k hops. Then Maag et al. [93, 94] and Arfire et al. [4] extended this work to models with multiple-kind variables, with and without time dependency. They showed that the complexity of the model should be adjusted based on the frequency of rendez-vous.

In a similar way, Markert $et\ al.\ [102]$ introduced a calibration strategy based on rendez-vous but with a particular focus on privacy aspects for the exchange of data.

Kizel et al. [75] also proposed a multi-hop calibration method consisting into co-locating two devices for a certain time, one being the reference to the other, and then moving the freshly calibrated device close to another non-calibrated, a reference instrument being introduced in the loop to reset the error that accumulates. The advantage is the error is related to the number of hops that took place like in [63].

Sailhan et al. [129] developed a multi-hop, multiparty calibration scheme with the addition of an assessment protocol for the relevancy of the calibration, based on a weighted directed hypergraph of the network, the weights indicating the quality of the calibration. The presented strategy was applied to blind networks but as in [63], it could be extended to partially blind networks.

In the case where the instruments cannot be mobile, Weissert *et al.* [108, 171] proposed to use a proxy model in their framework for the calibration of sensor networks. It consists into using the mean and standard deviation of the data obtained at a given location that is equivalent in terms of land use compared to the location of the instrument to calibrate.

Fonollosa et al. [56] used various models for calibration of chemical measuring instruments with an approach called "calibration transfer" which is a kind of multi-hop calibration. Indeed, the principle is to calibrate one of the instruments and then apply the same model on other instruments to calibrate, eventually with a transformation to properly map the measurement spaces. This approach was also used by Laref et al. [80]. Such strategies are widely used in the field of spectroscopy [173]. It, however, requires the measuring instruments are more or less behaving the same under identical varying conditions, which is not always the case when dealing with low-cost instruments. Moreover, in the context of environmental sensing with static instruments over a large area, it is very unlikely that the instruments measure the same quantities at the same time. Thus, deriving the correct transformation to transfer the calibration relationship is challenging. Cheng et al. [31] proposed an in-field calibration transfer method to learn this transformation when the distributions of the reference values at the source and target locations are similar, and when this transformation can be assumed as linear.

4.5 Blind macro calibration

Blind calibration strategies for sensor networks were first developed as a way to locate spatially the nodes in a static network. Whitehouse et al. [172] proposed to solve an optimisation problem ensuring that consistency and geometrical constraints were respected. Ihler et al. [65] proposed nonparametric belief propagation instead. Taylor et al. [152] developed an inference technique using a Bayesian filter. However, as they targeted spatial localisation, most of these methods based their algorithms on electromagnetic or acoustic [151] propagation (time delay, intensity loss...), and cannot be applied directly to other calibration problems.

Bychkovskiy et al. [24] proposed a first solution that could be used for any measurand, provided that the existence of relations between instruments of the network is known. It demands first to estimate the parameters of each existing relationship between instruments in the network. Then, the consistency of the derived relationships must be maximised to be resilient to circular dependencies. This technique was applied to a dense static sensor network and has not been yet extended to a mobile sensor network. In theory, it could also be applied to a sensor network with reference nodes but there is no report on the topic.

Later, Balzano et al. [7] developed a blind and partially blind calibration strategy for static sensor network suitable for any measurands, without any prior knowledge on existing relationships between the values of instruments. They tested it notably on temperature, light intensity or CO₂ level measurements. The key postulate is the sensor network is dense enough to oversample the signal of interest. They proposed that the true signal lies in a subspace of the space formed by the measured values. Considering the prior choice of the subspace, the parameters of the calibration relationships for all nodes are then estimated using singular value decomposition or by solving a system of equations using a least square estimator. This method was extended later in [89] to provide a total least square formulation and also in [44] to take into account outliers and separate them from the measurement matrix.

Alternately, Takruri et al. [148–150] addressed calibration as a drift compensation problem. They proposed to proceed recursively: measured values at step n are first corrected with predicted drifts obtained at step n-1, then the next measurements are predicted using support vector regression (SVR). Finally, the predicted values are used to estimate the drifts using a Kalman filter. Kumar et al. [76, 77] replaced SVR by kriging, which is a method of interpolation originally from geostatistics, as a prediction method for next values. In the same vein of using spatial relationship between the measured values, Chen et al. [31] introduced a spatial correlation model for drift correction targeting sensor networks deployed in buildings.

Ramakrishnan et al. [122] proposed another blind calibration strategy based on a gossip-based distributed consensus strategy, with SAGE algorithm used for parameter estimation. Later Buadhachain et al. [23] used an expectation-maximisation algorithm instead. The consensus-based approach is interesting as it can reduce the communication bandwidth.

Bilen et al. [13] extended the problem of blind calibration to the case of sparse input signals, which are measurements with missing/useless information, and exploited compressive sensing to estimate the instruments' corrective gains.

Cambareri et al. [25] proposed a non-convex formulation of the problem and gave a formal criterion of convergence of the calibration method enabling to estimate corrective gains for the values.

For mobile nodes, C. Wang et al. [162] proposed a method which exploits the moments of the measurements, here the average and the variance. They formulate calibration as an optimisation problem minimising the difference between the moments of the true signal and the measured one. The method was extended in [161] and [163] for different expressions of the calibration relationship. The approach was later adjusted by C. Xiang et al. [176] to address specifically mobile crowdsensing, with instruments embedded in mobile phones for instance. Dorffer et al.

[41, 42, 45] considered mobile sensor network calibration with mono-kind variables and linear relationships for spatially sparse signals by using matrix factoring, with an extension regarding non-linear calibration [43].

Wang et al. [168] proposed an extension of [7] relaxing some hypothesis and based their estimation on a Kalman filter with the help of a drift detection strategy. The approach was improved in [87] in terms of drift detection, number of instruments allowed to drift at the same time and pre-processing of the input signal with wavelet denoising. Wu et al. [175] applied the same idea recently but in the case of clustered sensor networks, e.g. when the instruments of the networks can be grouped into several clusters. In this case, it is more a group calibration strategy than a macro calibration.

Wang et al. then extended their work for sparse signals with either a Bayesian approach [167] or a deep learning approach [166].

Yang et al. [181] also based their work on the idea of [7]. They prove that, if the underlying signals follow a first-order auto-regressive process, then the parameters of the linear calibration model are recoverable. They use a nonparametric Bayesian model to do so.

Overall, macro calibration methods do not suffer from error propagation issues unlike pairwise approaches. However, because of the absence of references, they usually require large amount of data, which is typically available in the case of highly dense static networks or mobile networks with high frequency of rendez-vous.

4.6 Group strategies

 On the first hand, group calibration strategies are used to calibrate multiple instruments located at the same place and measuring different quantities that could influence the calibration relationship. It was shown that their corrected values were more accurate when exploiting multiple-kind variables in calibration relationships [9, 10, 47, 53, 100, 120, 139, 140]. Therefore, strategies [63] that were developed for pairwise and mono-kind variables calibration in the first place were extended to group calibration [4, 93, 94]. New methods were also developed.

Based on results like [71, 139, 140], Zimmerman *et al.* [186] introduced a calibration strategy using random forests with multiple-kind variable relationships for measuring systems when the latter are co-located to reference instruments.

Kim et al. [74] recently presented an approach of blind group calibration with prior information on the cross-sensitivities of instruments, known from laboratory experiments, to build calibration relationships with multiple-kind variables.

On the other hand, other works proposed group calibration strategies to reinforce the confidence in the values used as standards.

Miluzzo et al. [107] proposed a calibration technique with multiple reference nodes in a sensor network by formulating a distributed average consensus problem estimating the offset of each non-reference node. The concept of sensing factor was introduced. It refers to the area within which the measurand value can be assumed to be identical for the reference instruments and the instruments to calibrate.

Lee et al. [82, 137] proposed a blind group approach for mobile nodes. The area of deployment is divided into several non-overlapping regions. It is assumed that calibration relationships exist in each of these areas. This is used to formulate the parameter estimation problem as a Laplacian linear equation relating the drift, the measured values and noise.

Stankovic et al. [144] developed a novel methodology based on exploiting the neighbours of each node. Relations in the network are expressed in a matrix and, as there are groups of neighbours, the matrix can be decomposed into blocks. Each block represents a group of relations with parameters to estimate, reducing the problem compared to a macro blind calibration. This approach was extended in [143] to better deal with the case of additive measurement noise.

Fishbain *et al.* [55] proposed a method of aggregation for non-calibrated sensor network relying on a group-consensus strategy.

Fang et al. [50] later introduced reference instruments that are measuring different quantities, in view of analysing the influence quantities in the calibration relationship. More recently [51], they also added an outlier detection and removal strategy before the calibration stage to improve the quality of the input data.

Yu et al. [184] also proposed to derive features from the time series of the measurands to reduce the effect of cross-interference phenomenon between two quantities and then to use a deep calibration model to correct the measuring instruments. The same idea of feature extraction was also developed by Wang et al. [165] in a work extending the idea of piecewise linear calibration of [163] to a group calibration with different types of calibration relationship computed piecewise.

Becnel et al. [11] used the concepts developed in [63, 131, 132] notably to propose a group calibration strategy with a weighted neighbourhood approach. On a same dataset, their approach seems to improve the results by 20 % compared to the previous works they considered.

4.7 Comment regarding other surveys

By comparing the common references in Table 2.4.1 to the references listed in Table 5 of the survey of Barcelo-Ordinas et al. [8], we can observe that some works are not classified in the same way like for instance the work of Balzano et al. [7] that is classified as "micro" for the attribute "area of interest" and "single sensor" for the attribute "number of sensors" whereas it was classified as a macro calibration technique here. This was expected as differences were pointed out between the two taxonomies in Section 2. As the groups "area of interest" and "number of sensors" of Barcelo-Ordinas et al. are similar to the grouping strategy in the presented taxonomy, it appears we did not understand the algorithms in the same way. This underlines the use of a shared terminology is critical to ensure a good comprehension of the algorithms.

5 Conclusion

 In this chapter, a taxonomy for the classification of *in situ* calibration algorithms for sensor networks has been proposed. It is based on four groups of categories capturing both the different network architectures and algorithmic principles: the availability of reference instruments in the network, the mobility of the instruments, the kind of input variables in the calibration relationships, and the instruments grouping strategy (pairwise, macro or by group) used for a calibration procedure.

The review shows that relatively few papers address methodological issues related to reference-based pairwise strategies, as this approach is the closest to a "traditional" calibration approach and features relatively little challenges. Partially blind and blind pairwise calibration methods, which often focus on mobile nodes, are more complex as they require to define calibration relationships between non-reference nodes, which translates into error propagation issues. Macro calibration approaches are mostly used to deal with reference-less situation. Their challenge lies in defining valid calibration relationships between non-reference instruments. Group strategies appear to improve on performance of both pairwise and macro strategies with or without reference instruments. Most methods are based on calibration relationships with mono-kind variables and a linear expression, but more complex models are progressively appearing to better address the complexity of environmental sensing. Likewise, while most work initially focused on static networks, there is now a strong interest for mobile nodes as they allow for physical rendez-vous between nodes, which reduces the impact of the physical variability of the phenomena between static distant nodes.

In a general manner, the main situations and issues for the *in situ* calibration of environmental sensor networks have been addressed multiple times in the literature. To identify the improvements

that could be brought, the question of the performances of the existing algorithms must be tackled, which is the subject of the next chapter.

The content of this chapter is based on the following publication:

F. Delaine, B. Lebental and H. Rivano, "In Situ Calibration Algorithms for Environmental Sensor Networks: A Review," in *IEEE Sensors Journal*, vol. 19, no. 15, pp. 5968-5978, 1 Aug.1, 2019, DOI: 10.1109/JSEN.2019.2910317.

The review presented in this thesis was extended to the papers published after this survey.

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Chapter 3

Framework for the Simulation of Sensor Networks Aimed at Evaluating In Situ Calibration Algorithms

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1 Challenges for the comparison of *in situ* calibration algorithms

In Chapter 2, after proposing a taxonomy for the classification of *in situ* calibration algorithms, a review of the existing works was conducted. The literature on the subject is abundant and numerous contributions were reported, covering all the possible types of sensor networks. Nevertheless, no comparison of the *in situ* calibration strategies in terms of performances was proposed in this survey. Indeed, though the presented algorithms in publications are operational from a mathematical perspective, a formal quantification of the improvement is never provided. Instead, one or several case studies are systematically provided to demonstrate how the algorithm is put in practice and its efficiency on these specific cases. This step is either performed through simulations [45], laboratory experiments or field experiments [109], or both [7].

The principle of such studies consists in comparing the measured values, both before and after calibration, to the values that should have been measured by the instruments if they were ideal. The latter values are obtained with the help of simulations or with higher quality instruments co-located to the instruments to recalibrate.

Regarding approaches by simulation, different ones are proposed to compute numerically the values of the measurand. It can be based on 2D Gaussian fields [41] or ARMA processes [166] for instance. Disturbances like noise or drifts are added to the reference values in order to generate the measured values. However, the models can be more or less realistic. Thus, the model used may raise the question of the relevance of the results obtained, notably with respect to experiments.

Concerning experimental approaches, there are lots of datasets produced and used in various studies (for instance the Intel Lab Data [79], a deployment at James Reserve used in [7], the OpenSense Zurich Dataset [86]). They may be representative of particular types of deployments, for instance indoor (Intel Lab Data) or outdoor (James Reserve, OpenSense) measurements, with static and dense sensor networks (Intel Lab Data, James Reserve) or mobile sensor networks (OpenSense). Moreover, whereas repeating a simulation with different network settings (number of instruments, types of instruments, positions...) but identical phenomenon is feasible, it is almost impossible with experiments except under laboratory conditions. Although there are existing facilities that could allow it [38], they are not adapted for the study of sensor networks over large areas. A solution could be the deployment of as many instruments as necessary to produce the desired configurations, but this would drastically increase the cost of the experiment. Moreover, this would have to be reproduced in multiple environments in order to study its influence on the performance of in situ calibration strategies.

More generally, whether we focus on simulation-based or experimental approaches, there is no single case study that is reused a significant number of times in multiple publications by different groups of authors according to the review conducted in Chapter 2. It prevents the comparison of existing works. Such a study is conducted in [35] but as for the work of Karagulian *et al.* [73] reported in Chapter 1 Section 1.3.3, the conclusions must be considered with precautions. In this case, the performances after calibration of measuring systems using instruments from the same manufacturer are compared based on the results published by different groups of researchers. However, the results were not obtained on the same case study. In this way, conclusions may be incorrect.

As an explanation to the absence of one or several main case studies in the community, they are not always easily reproducible. For simulation-based ones, some of their parameters may have not been provided or even mentioned. This is why an identical approach may be used in multiple papers but rarely from different authors. Regarding experiments, datasets are often shared with the community, though they sometimes become unavailable over time or require post-processing that is left to the user. Nevertheless, as for simulation, a dataset is rarely used in publications from authors who did not produce it. This may be due again to the fact that

experiments are tailored for specific cases and calibration strategies.

Another difficulty is the multiplicity of metrics that can be used to assess the performances of the strategies [17]. Such metrics are for instance the mean absolute error, the mean absolute percentage error, the root-mean-square error, the Pearson's correlation coefficient and so on. Original metrics developed in works on the evaluation of the performance of instruments [54] or ones related to official recommendations such as the data quality objective (DQO) in the context of air quality measuring instruments [16, 27, 95] are also used. They are mostly based on prior information concerning the expected corrected values. As shown in the comparison of the existing works carried out by Cross *et al.*[35], authors do not all use the same metrics in their case studies.

Also, in the perspective of comparative studies, while a few authors have shared the code associated to their publications as in [89] or [44], most of the codes are not open-sourced. Therefore, algorithms have to be fully reimplemented in most cases to achieve comparison [168] and due to the efforts it requires, it is carried out in comparative studies only for strategies with features similar to the one they propose–for example the use of machine learning techniques [26, 48, 49] or the use of Kalman filter [168].

Overall, while a new strategy can be overperforming previous ones on a given case study, one rarely knows whether the new strategy is better in general or only for the specific case (Fig. 3.1.1), regardless of whether it is based on simulation or on experiment.

Publication	P_a	\blacksquare P_b	P_c
Algorithm	$A_{\underline{a}}$	A_b	A_c
Case study	I C_a	I C_b	C_c
Conclusion	A_a works on C_a	$A_b > A_a$ on C_b	$\begin{cases} A_c > A_b \\ \text{on } C_c \end{cases}$
Not proved	-	$A_b > A_a$ on C_a	$A_c > A_b$ on $C_{a,b}$ $A_c > A_a$ on $C_{a,b,c}$

Figure 3.1.1: Usual evaluation process performed across papers: consider three publications, P_a , P_b and P_c that are published successively. P_a introduces an algorithm A_a which is evaluated on case study C_a . It concludes that the algorithm is working. Later, P_b presents A_b and evaluate it with A_a on case study C_b , concluding that A_b has better performances than A_a . Finally, P_c brings forward A_c and show it is better than A_b on case study C_c . In the absence of unified validation case study, A_b may still be better than A_c on $C_{a,b}$, and A_a better than A_b and A_c on C_a and $C_{a,b,c}$ respectively.

Thus, there is a strong need for systematic tools and protocols [84, 95] enabling to compare across studies the performances of *in situ* calibration methodologies.

This chapter is addressing this question. The contribution consists in a generic framework to design case studies that can be used for quantitative comparisons of *in situ* calibration strategies. Its objective is to get a better understanding of the factors driving the performances of calibration strategies and the quality of sensor networks. For reproducibility and scaling issues, it is based on numerical simulations of the environmental phenomenon and of the measuring instruments, their metrological properties as well as the faults introduced in the measurements. It is henceforth not a new calibration methodology, but the first methodology to carry out a systematic inter-comparison of the performances of existing methodologies.

We apply this framework on seven calibration algorithms. In a first step, we consider an

elementary case study on a blind static network that is sufficient to get first insights by analysing the performances with commonly used metrics at different scales. Then, we evaluate the impact of the realism of the model of pollution emission and dispersion, the density of the deployment of the measuring instruments, and the error model of the devices. Finally, several engineering insights are deduced from the investigations carried out, validating the relevance of such a framework in the design phase of a practical network of low-cost environmental sensors.

2 Description of the framework

 A generic framework to design case studies is proposed in this section for quantitative comparisons of *in situ* calibration strategies. It aims at yielding a better understanding of the factors driving the performances of calibration strategies, and at providing a protocol as rigorous as possible to conduct such studies. The different steps are described in a generic way to ensure their applicability for most cases.

Like for faults in Chapter 1 Section 2, we adopt a data-centric point of view. We do not target the modelling of the entire measuring chain of an instrument described in Chapter 1 Section 1.2 and illustrated in Figure 1.1.2 in the same chapter. We assume that instruments are grey boxes which provide measured values based on true values of a measurand and possibly also on true values of influence quantities. The grey boxes representing instruments consist into algorithms mimicking the features of real instruments.

In this work, we focus on the metrological performances of *in situ* calibration strategies. Other subjects such as the communication costs, the energy consumption or the computational efforts for instance are not in the frame of this study.

2.1 Simulation-based strategy

The methodology is based on simulation because it enables the following properties that are difficult to get with field or lab experiments

- Ability to perform a study with different operating conditions for a same sensor network.
- Reproducibility: same operating conditions on different sensor networks.
- Knowledge of true values: in experimental operating conditions, true values are not known perfectly—there is always an uncertainty— and having very accurate values requires high-quality instruments which are usually expensive.
- Scalability: the only limit to the density, number and diversity of measuring instruments is the computing time.

To conduct the evaluation of in situ calibration strategies, the framework combines the simulations of the environmental phenomena to produce the true values of the measurand, and of the potential influence quantities, at any position and time, the simulations of the mobility –if any– of the nodes of the network to know where and when measurements are performed, and the simulations of the measuring chain of the instruments to produce the measured values. The realism of the results will depend on the complexity and accuracy of each simulation model. On the other hand, being able to analyse results on a simplistic model can also help to highlight fundamental properties before confronting them to more complex situations. In particular, even if we consider networks of sensors, we neglect to simulate the system and networking aspects. We consider they have a lesser influence on the metrological performances of in situ calibration strategies. One could, however, argue that system and network issues could challenge the robustness of a given implementation of a calibration protocol because of packet loss. It could also be interesting to evaluate the energetic cost or time to converge of such implementation. It

is still possible to extend the framework to these issues that could be covered by tools such as $WSNet^{11}$, $NS-3^{12}$ or other 13 .

The objective of this framework is to help end users in the design of their sensor networks by:

- showing which strategies are applicable to a given use case. Currently only a few strategies are supplemented with strong formal criteria enabling to determine whether they can be used for a given network and in a given situation, for instance in [45] or [7]. Such criterion may, however, be not easy to define and therefore simulation is an interesting solution.
- showing which strategies should give the best performances regarding the assumptions that can be made in practice.
- allowing the optimisation of the settings of the calibration strategies –as in [109] where the time period of calibration is studied. As there are rarely formal applicability criteria, there is also rarely protocols defined for the adjustment of the parameters.

2.2 Functional decomposition

Our methodology, represented in Figure 3.2.1, can be described schematically as follows:

- 1. Build a dataset of ideal measured values.
 - (a) Simulate the quantities involved with a sufficiently high resolution (spatially and temporally).
 - (b) Simulate the positions of the measuring instruments.
 - (c) Combine the two simulations.
- 2. Build measured values from the ideal measured values, e.g. add defects to these values.
- 3. Perform calibration, e.g. determine the correction to apply to the measured values.
- 4. Derive corrected measured values.
 - 5. Compare measured and corrected measured values with ideal measured values. If the algorithm performs well, the corrected values should be closer to the ideal values than the measured ones.

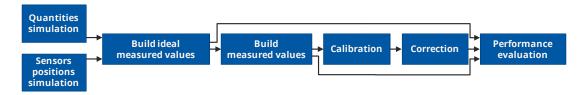


Figure 3.2.1: Schematic diagram of the methodology proposed for *in situ* calibration strategies evaluation

To apply this methodology, input data are the number of instruments involved, the quantities measured, the $in\ situ$ calibration strategies to apply and the metrics considered for the evaluation of the results but also:

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¹¹WSNet: http://wsnet.gforge.inria.fr/

¹²NS-3: https://www.nsnam.org/

¹³OMNet++, OpNet, LoRaSim...

- Information to simulate the quantities involved. Depending on the model, it can be as simple as some parameters of an equation (e.g. a Gaussian model) or a complex description of an urban scene (e.g. http://air.ec-lyon.fr/SIRANE/)
- Specifications on the deployment and possible mobility of the instruments. Since the positions of the instruments is critical for the performances of the sensor network, the number of instruments deployed, and their positions may be optimised [19].
- Descriptions of the measuring chain of the instruments involved, in terms of drift and of their potential influence quantities.
- Configuration of the considered in situ calibration strategies.

As outputs, the framework provides time series of the true values, of the measured values and of the corrected values for each instrument. Values of metrics computed for the comparison of these times series may also be given.

The following subsections give details on each step.

2.2.1 Build ideal measured values

Ideal measured values are measured values that should be measured by instruments if they were perfect. It means for an instrument s_i that its measured values $v(s_i, t)$ are equal to the true value of the measurand, noted $v_{true}(s_i, t)$, for all t.

Generating these ideal measured values for all the instruments of a sensor network can be carried out as follows.

To build ideal measured values, true values of the measurand must be computed. True values of other quantities may also be derived for the simulation such as influence quantities if the instruments to model undergo such effects.

To be able to repeat the study with different sensor networks, the quantities can be simulated independently of the positions of the instruments so that one may build different sets of ideal measured values from a same simulation of the quantities depending on the sensor network.

Therefore, the simulation is performed on the targeted domain of study, e.g. a specific geometry, at a sufficient spatial and temporal resolution, regarding the potential positions of the instruments and their frequency of measurement. The model used for the simulation depends on the application and on the targeted accuracy. Ideally, the simulation should perfectly represent the studied environmental phenomena and the quantities affecting instruments operating under real conditions. This is not always possible and thus the models used may not reflect the real temporal and spatial variability of environmental phenomena. This is particularly true for air pollution monitoring. While high model accuracy can be achieved with advanced computational fluid dynamics models [12], such models require time and resources that are not always acceptable or available for users. In multiple works, case studies were conducted with abstract models [7], enabling to demonstrate that the proposed calibration strategy is functional. However, such models do not enable determining if an algorithm is applicable on real cases.

In addition to the true values of the considered quantities, the sets of the positions where measurements are performed must be built with respect to the position and possible mobility of each instrument. Based on the positions of the instruments over time, true values are extracted and stored from the simulation of the studied quantities at the locations of each instrument at each time step according to their sampling periods.

Afterwards, a time series of ideal measured values is available for each measuring instrument. 14

¹⁴In this section, we considered that the simulation of the quantities and the simulation of the positions of the instruments are carried out separately and then combined. However, it might be possible to determine first the positions of the instruments and then to derive the ideal measured values only at the positions for instance.

After this step, these time series must be validated against the hypotheses of the targeted calibration strategies.

2.2.2 Build measured values from the ideal measured values

Until this step, measuring instruments are assumed ideal. Thus, to obtain realistic measured values, e.g. values that include the real behaviour of measuring instruments, faults must be introduced. The taxonomy presented in Chapter 1 Section 2.2 can be used here to model the measured values. As we study in situ calibration strategies, drift faults are introduced principally. The influence on the performances of in situ calibration strategies of other faults (e.g. noise, spikes, missing data...) can also be studied by introducing them in the measured values (see Section 5.3.2).

2.2.3 Perform calibration and build corrected values

Once measured values are obtained, calibration may then be performed low-cost to correct them. It must be based on algorithms applicable to the specific case study addressed.

Calibration and correction are separated in the methodology for the sake of genericity. Some calibration strategies directly determine the correction to apply to each value [149], while others provide parameters of a mathematical relationship to apply to each measured value [7].

2.2.4 Evaluate

Finally, corrected values, measured values and true values can be compared to assess the performances of the algorithms. Most of the time, usual metrics are employed such as root-mean-square error, mean absolute or Pearson's correlation coefficient for each measuring instrument in the network. Results of case studies may also be presented as the mean of these metrics over the whole network [150]. We discuss the most suitable metrics to use in Section 4.

In the following sections, the framework is applied to the comparison of several strategies for blind static networks on a case study that is simple enough to focus on fundamental issues.

3 Comparison of *in situ* calibration strategies for blind static sensor networks

The purpose of this section is to give a simple example of how to apply the framework.

3.1 Frame of the study

In this section, the case of blind static sensor networks is considered. All the instruments remain at the same position and no reference instrument is present. This type of sensor network offers a particular challenge in terms of calibration. By contrast, when reference instruments are in the network, trustworthy values are available. Secondly, when instruments are mobile, it is possible that two or more instruments are measuring the same true quantity value when they are in a spatio-temporal vicinity. For blind static sensor networks, the availability of standard values may not be assumed, and only the instruments that are deployed close enough can compare their measurements.

We consider seven calibration strategies, from four different groups of authors, that were identified in across the review of Chapter 2. These strategies have never been compared all together on a common case study.

• Balzano et al. [7]: To apply this algorithm, the sensor network must be dense enough to oversample the signal of interest. We refer to this assumption by the term "oversampling" afterwards. With such a sensor network, the true values lie in a subspace of the space formed by the measured values. To calibrate the instruments, this subspace must be

known. The gains and offsets for all nodes are then estimated with the chosen subspace and the measured values using singular value decomposition (SVD) or by solving a system of equations using a least square (LS) estimator. These strategies are respectively called **SM-SVD** and **SM-LS**, SM standing for subspace matching. Note that gains are estimated up to a multiplicative coefficient. Consequently, the gain of one instrument should be known to complete the calibration. Likewise, the offset computation requires either to know some of the offsets, or that an additional hypothesis on the signal is valid. As such, the network is not perfectly blind in truth.

- Lipor et al. [89]: This work is an extent of [7]. It shows first that the solution obtained with SM-SVD is equivalent to the one obtained by solving the system of equations expressed as a total least squares problem (TLS). The latter strategy is called **SM-TLS**. ¹⁶
- Takruri et al. [148–150]: Three contributions from these authors are considered. The first one is based on the idea that the average of the values of neighbouring measuring instruments gives a good estimation of the correct value of an instrument to calibrate. If the absolute difference between the measured value and the average is greater than a given threshold, this difference is added to the instrument's value. We call this strategy **AB-DT** (average-based estimation for difference-based threshold). Two variations are also proposed, a first one in which the difference-based threshold mechanism is replaced by a Kalman filter that estimates the error affecting the instruments (**AB-KF**), and a second one where the average-based part is replaced by support vector regression (**SVR-KF**).
- Kumar et al. [76, 77]: This work, inspired from the previous one, also proposes the use of a Kalman filter to estimate the error of each instrument but instead of performing an average over neighbouring nodes or using SVR for true value estimation, kriging is used (K-KF).

Note that not every strategy that could have been applied to our case study is considered here. For instance, the work of Dorffer et al. [44] extends [7] and [89] for sparse signals, but this is not in the scope of the study.

3.2 Application of the framework

3.2.1 Simulation of the ideal measured values

A space of 1000×1000 m is considered. It is discretised with a step of 10m. At the centre of this area, a NO_2 source is considered. The concentration of NO_2 , C, at the instant t and position (x, y), is modelled as:

$$C(x, y, t) = A(t) \exp\left(-\frac{x^2 + y^2}{\sigma(t)^2}\right)$$

It is a 2D Gaussian function with an equal spread σ for x and y. This model is not very realistic but has been used in other papers [45] for its simplicity of implementation. Examples of pollution maps are represented in Figure 3.3.1.

To facilitate the interpretation of σ , it is expressed as a function of the full width at half maximum (FWHM) of the Gaussian curve:

$$\sigma = \frac{FWHM}{2\ln(2)}$$

¹⁵A usual assumption is that "the signal has a null (or known) average".

¹⁶Note that additional formulations given in the publication, in the case where some gains are known for a subset of instruments, are not considered here as it would be equivalent to have a partially blind sensor network.

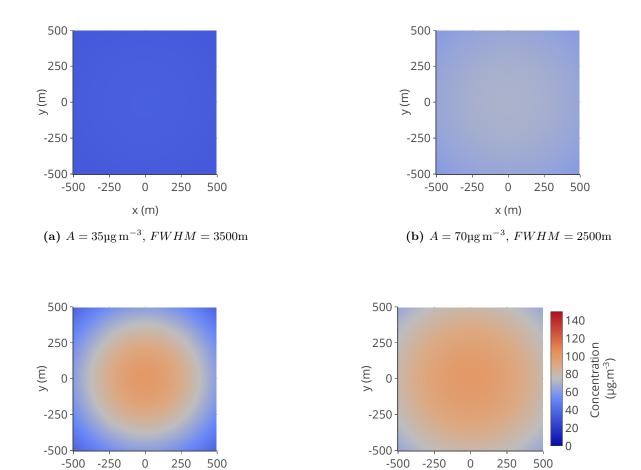


Figure 3.3.1: Examples of maps of concentration C used following $C(x,y) = A \exp \left(-(2 \ln(2))^2 \frac{x^2 + y^2}{FWHM^2}\right)$ for given A and FWHM

x (m)

(d) $A = 100 \mu \text{g m}^{-3}$, FWHM = 1500 m

x (m)

(c) $A = 100 \mu \text{g m}^{-3}$, FWHM = 1000 m

A and FWHM are functions representing respectively the temporal evolution of the amplitude of the 2D Gaussian function and of its FWHM.

To be representative of actually measured concentrations, the function A is based on values measured by a NO_2 monitoring station in Paris between 2017/01/01~00:00 and 2017/12/31~23:00, with an hourly time step (Paris centre station [2]).¹⁷ We consider it represents our pollution source at the coordinates (0,0). Even if the values of the station are not the ones of a real source, we assume it gives a reasonable order of magnitude of variation over time.

The FWHM represents the spread of the pollutant around the source. Two linear piecewise functions are defined to represent the daily evolution of the FWHM: one for weekdays and one for weekends. They are represented in Figure 3.3.2b. Their shape is inspired from the dataset

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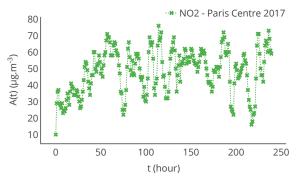
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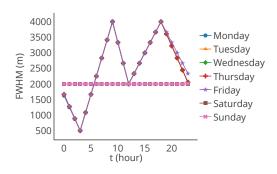
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¹⁷There were missing entries in the time series. Values were first interpolated with a limit of 3 consecutive values to fill, e.g. 3 hours. Then, in the case where more than 3 consecutive values are missing, the interpolation is made based on the values at the same hour of the previous and next day. The values are interpolated with a linear function for both cases.





(a) Partial plot of the values used to model the amplitude A of the phenomenon model

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(b) Profiles designed for the daily evolution of the FWHM per days of the week

Figure 3.3.2: Evolution of A and FWHM for the modelling of the concentration of pollutant

used for the amplitude A. This is not fully realistic, but it provides an order of magnitude of spatial and temporal variations occurring in urban context.

The concentration of pollutant C is simulated for a year at an hourly time-step over the area of study. To avoid the computation of C, at each time step, for the current value of A and FWHM, a catalogue of maps is generated from sets of possible values for the amplitude and the FWHM: respectively from 0 to $150 \mu \text{g m}^{-3}$ with a step of $5 \mu \text{g m}^{-3}$ and from 0 to 4000 m with a step of 500 m. In our case, it required the simulation of 157 maps. To derive the time series, the map generated with the closest allowed values of A and FWHM is picked from the catalogue at each time step. Finally, a time series of 8760 maps is obtained.

A sensor network S of |S|=16 static measuring instruments is considered. It is uniformly deployed spatially. The positions are represented in Figure 3.3.3. Then, the time series of concentration maps and the positions of the instruments are combined to obtain the time series of ideal measured values for all the instruments, noted v_{true} . In this case, we consider that the instruments measure one value per hour.

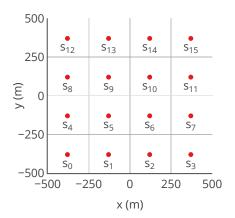


Figure 3.3.3: Positions of the 16 measuring instruments considered in the case study, deployed uniformly in the field

The validity of this set of ideal measured values must be verified for SM-(SVD,LS,TLS)

strategies. The network should satisfy the hypothesis of "oversampling". To verify it, a principal component analysis (PCA) of the matrix formed by the concatenation of the ideal measured values time series of each node is carried out. We assume that the sensor network is satisfying the oversampling hypothesis if the explained variance ratio of the ν first components is greater than 0.999 and $\nu < |S|$. In our case, the condition is met for $\nu = 2 < |S|$. Therefore, the oversampling hypothesis is satisfied.

3.2.2 Building of the measured values

The network is defined as "blind", which can also be expressed by "all the instruments have the same sensing reliability". In a first step, instruments are only drifting. Drift of each instrument is only function of time. We assume all the instruments undergo a linear drift increasing up to 5% of the gain every 7 days. The actual drift is randomly drawn following a uniform law. The value of 5% is set according to what is reported in typical measuring instrument datasheets [105].

The instruments are assumed to be initially calibrated and to remain faultless for eight weeks, which gives a training dataset for the strategies needing it, notably to learn the subspace of the signal with PCA for **SM-(SVD,LS,TLS)**. For four additional weeks, the instruments are kept without drift to study how calibrations strategies behave if the devices are not drifting. Finally, the instruments start drifting after twelve weeks. This instant is noted $t_{\rm start\ drift}$.

Therefore, the gain $G(s_i, t)$ of the instrument s_i at t is:

$$G(s_i, t) = \begin{cases} 1 & \text{if } t < t_{\text{start drift}} \\ G(s_i, t - 1) & \text{if } t \ge t_{\text{start drift}} \text{ and } (t - t_{\text{start drift}}) & \text{mod 7 days } \ne 0 \\ G(s_i, t - 1) + \delta G(s_i, t) & \text{if } t \ge t_{\text{start drift}} \text{ and } (t - t_{\text{start drift}}) & \text{mod 7 days } = 0 \end{cases}$$
with $\forall t, \delta G(s_i, t) \sim \mathcal{U}(0, 0.05)$

This drift model is called "Weekly Gain Linear Increase" (WGLI).

The measured value $v(s_i, t)$ of the instrument s_i at t is expressed from the true value $v_{true}(s_i, t)$ using the relationship:

$$v(s_i, t) = G(s_i, t) \cdot v_{true}(s_i, t)$$

3.2.3 Configuration of the strategies

Each strategy has parameters to set up. They are defined as follows:

- SM-(SVD,LS,TLS): All these strategies share the same parameters, namely the periodicity of calibration and the time interval on which the linear system is solved. Note that the adjustment of the parameters was not discussed by the original authors. We chose to apply the strategy each week and to solve the linear system over the past w = 7 days.
- Kalman filter of (AB, SVR, K)-KF: Variables R and Q are set to 0.0001 and 1 respectively.
- Definition of neighbours for **AB-(DT,KF)**: Instruments are defined as neighbours if they are distant of less than 250m.
- Parameters of the SVR for SVR-KF: Kernel used is 'rbf', the kernel coefficient gamma is set to 10^{-8} and the penalty parameter C is set to 10^{3} .
- Parameters for the weighting for **K-KF**: a, c_0 and c_1 were defined as in [77], therefore a = 12, $c_0 = 0.25$, $c_1 = 0.85$.

We invite the reader to refer to the original publications for a more in-depth understanding of the parameters of these calibration strategies.

3.2.4 Evaluation

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To compare the corrected values of an instrument, obtained with each strategy, and its measured values against its ideal measured values, metrics are needed. We consider the most commonly used metrics in the publications reported in Chapter 2 and described as follows. They can be computed for each instrument or averaged over the network. We discuss the relevance of these statistics in Section 4. In the following, the set x stands for the k measured values of an instrument s_i over $[t - \Delta t; t]$, $V(s_i, (t, \Delta t))$ or the corresponding set of corrected values $V_{corr}(s_i, (t, \Delta t))$ and the set y is the associated k true values $V_{true}(s_i, (t, \Delta t))$.

• Mean absolute error (MAE):

$$MAE(x,y) = \frac{1}{k} \sum_{i=0}^{k} |x_i - y_i|$$

• Root-mean-square error (RMSE):

$$RMSE(x,y) = \sqrt{\frac{1}{k} \sum_{i=0}^{k} (x_i - y_i)^2}$$

• Mean absolute percentage error (MAPE):

$$MAPE(x,y) = \frac{100}{k} \sum_{i=0}^{k} \frac{|x_i - y_i|}{|y_i|}$$

Remark: the values of y must all be different from zero.

• Pearson correlation coefficient (ρ) :

$$\rho(x,y) = \frac{E[(x - E[x])(y - E[y])]}{\sqrt{E[(x - E[x])^2]E[(y - E[y])^2]}}$$

with E[.] being the expectation.

For perfect instruments before correction (or after with v_{corr} instead of v):

- $MAE(V(s_i, (t, \Delta t)), V_{true}(s_i, (t, \Delta t))) = 0.$
- $RMSE(V(s_i, (t, \Delta t)), V_{true}(s_i, (t, \Delta t))) = 0.$
- $MAPE(V(s_i, (t, \Delta t)), V_{true}(s_i, (t, \Delta t))) = 0\%.$
- $\rho(V(s_i, (t, \Delta t)), V_{true}(s_i, (t, \Delta t))) = 1.$

3.3 Results

In Figure 3.3.4, a partial plot of true values, measured values and corrected values obtained with each strategy is displayed for a particular instrument. For the considered device, a visual observation indicates that strategies SM-(SVD, LS, TLS) and SVR-KF provide better results than AB-DT, AB-KF and K-KF.

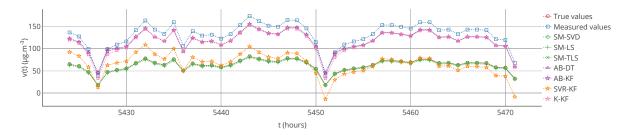


Figure 3.3.4: True values, measured values and corrected values with the strategies considered for a particular instrument $s_1 = 2$ between t = 5424h and t = 5471h. SM-(SVD, LS, TLS) and SVR-KF seem to provide better results than AB-DT, AB-KF and K-KF

	\mathbf{M}	A E	MAPE			RMSE			Pearson		
	μ	σ	μ	ι	σ	_	μ	σ	_	μ	σ
No calibration	18	2	5	3	6		25	2		0.935	0.007
SM- SVD	1	1	5	5	4		2	2		0.994	0.006
$\mathbf{SM}\text{-}\mathbf{LS}$	1	1	3	3	2		2	1		0.995	0.005
$\mathbf{SM}\text{-}\mathbf{TLS}$	1	1	5	5	4		2	2		0.994	0.006
AB-DT	18	1	5	5	6		25	1		0.932	0.004
AB- KF	18	1	5	8	11		25	1		0.933	0.003
$\mathbf{SVR}\text{-}\mathbf{KF}$	16	1	8	3	11		18	1		0.825	0.012
K-KF	18	1	5	3	10		24	1		0.927	0.005

Table 3.3.1: Mean and standard deviation of each metric, computed on the entire time interval of drift, over the 16 nodes of the network. SM-(SVD, LS, TLS) strategies have the best results overall whatever the metric considered. SVR-KF provides corrected values only slightly better than before calibration according to MAE and RMSE but not according to MAPE and Pearson correlation coefficient. AB-(DT, KF) and K-KF do not improve the measured values significantly.

Despite its advantage to easily visualise the results, this representation is not representative of all the instruments of the network.

The computation of the mean and standard deviation of each metric over the whole sensor network, on the entire time interval of study and with the results of each strategy, plus without calibration, is given in Table 3.3.1. With it, we observe that:

- SM-(SVD, LS, TLS) strategies have the best results overall with a small mean error and standard deviation whatever the metric considered
- SVR-KF which seemed to give interesting results in Figure 3.3.4 provides corrected values only slightly better than before calibration according to MAE and RMSE but not according to MAPE and Pearson correlation coefficient. This could be explained by two reasons. The strategy may correct well for high measured values but correct poorly for low ones. It could also be due to errors in the corrected values of particular instruments introduced by the calibration algorithm. The aggregation by averaging makes it impossible to discriminate between these two possible explanations.
- AB-(DT, KF) and K-KF do not improve the measured values

	\mathbf{M}_{L}	MAE		MA	PE	$\mathbf{R}\mathbf{N}$	1SE		Pear	son
	s_1	s_2		s_1	s_2	s_1	s_2	_	s_1	s_2
No calibration	22	18		66	51	30	24		0.924	0.942
$\mathbf{SM}\text{-}\mathbf{SVD}$	1	3		4	10	2	5		0.996	0.983
$\mathbf{SM}\text{-}\mathbf{LS}$	1	2		3	7	2	3		0.997	0.986
$\mathbf{SM}\text{-}\mathbf{TLS}$	1	3		4	10	2	5		0.996	0.983
AB-DT	20	17		61	49	26	23		0.933	0.932
AB- KF	18	17		56	45	25	23		0.935	0.927
SVR- KF	16	16		80	71	19	19		0.828	0.805
K-KF	17	16		51	43	24	23		0.930	0.918

Table 3.3.2: Values of the metrics, computed on the entire time interval of drift, for two particular instruments of the network $s_1 = 2$ and $s_2 = 10$. Strategies AB-(DT, KF) and (SVR, K)-KF are quite equivalent for these two instruments. The improvements are rather small for both instruments. For SM-(SVD, LS, TLS), results are consistent with the observations of Table 3.3.1.

These last two observations invited us to look at results for particular instruments. Table 3.3.2 gathers the results for two instruments selected randomly. We observe that strategies AB-(DT, KF) and (SVR, K)-KF are quite equivalent for these two instruments. The improvements are rather small for both instruments. For SM-(SVD, LS, TLS), results are consistent with our previous observation based on the average and the standard deviation in Table 3.3.1. The standard deviation does give the information of the existence but not the identity of the instruments that are degraded or improved by calibration, which is needed for practical deployments to drive the maintenance operations.

Furthermore, note that the results of Table 3.3.1 were computed over the 12 weeks of drift. The results may be different if computed over a different time range, for instance over each week. Figure 3.3.5 shows the evolution over time of the MAE computed each week for a particular instrument and each strategy. Table 3.3.3 provides statistics on the MAE computed each week but also for the other metrics that were computed in the same way. From the standard deviations of MAE, MAPE and RMSE, it shows that the observations made previously could be locally false, e.g. a strategy is better than others considering a computation of the metrics over the 12 weeks of drift but not always considering a computation of the metrics over each week. This is shown in Figure 3.3.5 where results with MAE for SVR-KF are nearly always worse than those for AB-DT, AB-KF and K-KF until week 24 but are better afterwards. This figure also shows that the performances of SVR-KF could be explained by the presence of a bias, at least according to MAE because it is quite constant in Figure 3.3.5.

3.4 Conclusions

Through this evaluation, we have shown that existing in situ calibration strategies could improve the measured values of a blind static sensor network in the considered case study. Overall, the strategies SM-(SVD, LS, TLS) have the best performances. This can be explained by the fact that the gain of at least one of the instruments in the network has to be known for these methods. In a way, the sensor network is only partially blind, but acquiring this information may not be a major issue in practice for large sensor networks. The other strategies appear to be able to mitigate the drift by few per cents at most only in the best cases, depending on the metric

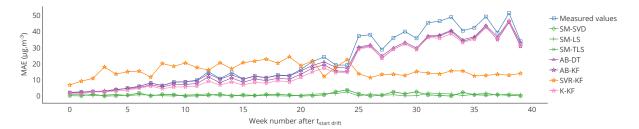


Figure 3.3.5: Evolution of the MAE computed for each week of the drift period between the drifted values and the true values, and between the corrected values for each strategy and the true values for a particular instrument, after the start of drift. MAE for SVR-KF is nearly always worse than those for AB-DT, AB-KF and K-KF until week 24 but are better afterwards. The performances of SVR-KF could be explained by the presence of a bias, at least according to this metrics, as its evolution is quite flat.

	M	MAE		MAP		RMSE		_	Pear	son	
	μ	σ		μ	σ		μ	σ		μ	σ
No calibration	22	16	(66	36		25	17		1.000	0.000
$\mathbf{SM}\text{-}\mathbf{SVD}$	1	1		3	3		1	1		1.000	0.000
$\mathbf{SM}\text{-}\mathbf{LS}$	1	1		3	2		1	1		1.000	0.000
$\mathbf{SM}\text{-}\mathbf{TLS}$	1	1		3	3		1	1		1.000	0.000
AB-DT	20	13	(61	32		22	14		0.999	0.001
AB- KF	18	13	į.	56	31		21	14		0.999	0.000
SVR- KF	16	4	8	80	39		18	4		0.935	0.034
K-KF	17	14	į	51	32		20	15		1.000	0.000

Table 3.3.3: Mean and standard deviation of the weekly mean of each metric for s_1 during the drift period. From the standard deviations of MAE, MAPE and RMSE, the observations made based on Table 3.3.1 could be locally false, e.g. a strategy is better than others considering a computation of the metrics over the 12 weeks of drift but not always considering a computation of the metrics over each week.

considered. The comparison between metrics enables understanding better how they operate. For all algorithms, studying the average values of the metrics over all the network hides strong disparities from instrument to instrument.

However, these results could be challenged based on the following considerations:

- The results may be different with another network (different positions and increased number of instruments, notably for AB-DT, AB-KF and K-KF strategies that use values of neighbour instruments for the prediction of correct values). However, the network is often defined by the use case. Thus, it is not always a criterion that can be optimised.
- The model of the considered phenomenon is not realistic and therefore results may not be practical. It is, however, sufficient to highlight issues of the algorithms under test. A more realistic model may not yield better result but would certainly make the analysis more complex.

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- Only a drift of the gain was added which does not fully relate the behaviour of a real instrument. Adopting a finer-grained model of the behaviour also increases the complexity of the case study.
- Other parameters of the calibration strategies may change the results. We empirically explored the parameters of the algorithms, keeping in this study the values that give the best results for each strategy. There is, however, no claim for optimality here, but the framework can be used to optimise the parameters of a given strategy.

In Section 5, the impact of these four separate aspects on the algorithms is tested. Before moving on to this, we discuss in the next section the metrics used and how additional ones, or other ways to visualise results, can give an enhanced comprehension of the behaviour of the calibration algorithms on a sensor network.

4 Evaluation of measurements after correction

4.1 Problem statement

To evaluate the performances of in situ calibration strategies, various metrics are commonly used like MAE, MAPE, RMSE and Pearson's correlation coefficient notably. We presented the results as statistics of these metrics over the entire network or values for particular instruments. The metrics were computed on the entire time interval of study or on subintervals. Results were displayed as tables and plots. These choices enabled carrying out an analysis at multiple scales (entire network versus individual instruments). This approach underlines disparities of efficiency of the algorithms on particular instruments for instance.

Multiple surveys exist on metrics [17]. They aim at providing a clear reading on the differences and similarities among the wide spectra of metrics used. While metrics are numerous, groups of them are based on the same idea (difference, absolute difference, square difference...). In some works, such as in [54], an integrated performance index was developed to combine the values of several metrics into a single one. However, some metrics are interdependent [154], which can exaggerate or hide information when they are combined. In the literature regarding in situ calibration strategies, the choice of metrics is not always clearly justified, nor is the choice of the values presented and how they are computed (values for particular instruments, statistics of the metric over the network and so on). Moreover, Tian et al. [154] showed it is possible to have equal values of multiple metrics for two sets of values very different from each other. They justified it by the interdependence, underdetermination and incompleteness of usual metrics, in addition to the fact that they assume the error to be linear. This issue is major as metrics are used in official recommendations such as the data quality objectives (DQO) in the context of air quality measuring instruments [16, 27, 95]. In this way, we suggest in the next section to use an error model as introduced by Tian et al. [154]. It is considered as a satisfying means to deal with these issues.

Another question lies in the way a metric is computed or in how statistics are derived from it. As shown in the previous section, looking at values of metrics computed over the full-time interval of drift or computed over each week may yield different interpretations. The choice of the time interval on which it is derived can be guided by requirements on the minimal duration for which the mean error of an instrument must be under a given threshold for instance.

Regarding the use of statistics, it concerns particularly networks with hundreds of nodes, for which interpreting the values of metrics for each node may not be easily feasible. However, statistics derived over the entire network do not enable discriminating easily between instruments differently affected by an algorithm. The use of statistics is perfectly appropriate if the quality of the calibration of the instruments belonging to a sensor network should be evaluated at the level of the network, e.g. when it is the collective performance that matter. Otherwise, if each

instrument is expected to be perfectly calibrated after an *in situ* calibration, statistics may only give a first idea of the performances of the algorithm. At the same time, evaluating the performances of an algorithm on a very large sensor network by looking at each instrument individually may not be feasible. Thus, the scale at which the performances of the instruments are evaluated must be specified in the description of a study, to make a proper use of statistics and to determine if the analysis is manageable in practice.

In addition, the way the results are displayed is important. We discuss this topic in Section 4.3 and propose the use of an easily readable representation, compatible with the error model we use hereafter.

4.2 Evaluation with an error model

We investigate in this section an approach which captures better individual instrument behaviours. It applied to the previous case study for the sake of comparison. We follow the error model approach proposed by Tian *et al.*[154]. It consists in finding a function \mathcal{F} associating each measured value of an instrument s_i , noted $v(s_i,t)$, to the true values of the measurand $v_{true}(s_i,t)$ so that $v(s_i,t) = \mathcal{F}(v_{true}(s_i,t))$. They illustrated this with a linear additive error model expressed as:

$$v(s_i, t) = a \cdot v_{true}(s_i, t) + b + \varepsilon(t)$$

with a and b being respectively a constant gain and a constant offset, and ε a random error, following a given probability law depending on the case. If the actual distribution of ε is not determined, it is equivalent to the fitting a linear trend model between v and v_{true} . A linear regression can be used to determine the slope a and the intercept b in this case. a=1 and b=0 are the ideal results, indicating a perfect calibration. The linear regression between v and v_{true} is appropriate to describe the performances before calibration. Comparing the slope and intercept of this regression to the ones obtained with the linear regression between the corrected values after calibration v_{corr} and the true values v_{true} enables determining the improvement brought by the calibration algorithm. This approach with linear regression is used in the literature on calibration of individual instruments [139]. In our case, this error model is perfectly appropriate to evaluate the remaining error after calibration, considering the drift model defined in Section 3.

The case study of Section 3 is analysed here again with the error model. The results associated with all the tables and figures presented are provided. First, the overall results obtained with this approach for the slope and intercept of the linear model, plus the score of the regression, e.g. the coefficient of determination of the regression, are in Table 3.4.1. Then, results are reported in Table 3.4.2 for the same particular instruments considered in Section 3.From these tables, the observations that can be made are consistent with those made with Table 3.3.1 and Table 3.3.2.

These tables also seem to confirm a bias for the strategy SVR-KF: the standard deviation of the slope and intercept among instruments is small, the mean slope is close to 1 but the mean intercept is equal to 10. However, the poor mean score of the regression invites to be careful.

Table 3.4.3 shows the mean and standard deviation of a, b and of the score computed weekly for a particular instrument. In this case, the means are different from those of Table 3.4.2 for s_1 (a linear regression is not a linear operation). We observe that the slope for SVR-KF is varying significantly. More importantly, the mean intercept is equal to zero with a standard deviation of 17, which is not the case for the other strategies. Therefore, the observation made with Table 3.4.2 about SVR-KF is not valid locally and can explain why the mean score is poor compared to the ones obtained with the other strategies.

In conclusion, the use of an error model allows to make the same conclusions as in Section 3. However, clearer information can be derived such as the remaining error after calibration,

¹⁸This idea of error model is close to the concept of a calibration relationship. The inverse function of the error model is the relationship to apply to an instrument so that calibration is perfect.

2398 expressed in terms of gain and offset.

	Slope		Inte	ercept	Sco	re
	μ	σ	μ	σ	μ	σ
No calibration	1.62	0.07	-2	0	0.87	0.01
SM- SVD	1.02	0.02	1	0	0.99	0.01
$\mathbf{SM}\text{-}\mathbf{LS}$	0.97	0.02	1	0	0.99	0.01
$\mathbf{SM}\text{-}\mathbf{TLS}$	1.02	0.02	1	0	0.99	0.01
AB-DT	1.62	0.05	-2	1	0.87	0.01
AB-KF	1.62	0.04	-1	1	0.87	0.01
SVR-KF	1.05	0.03	10	1	0.68	0.02
K-KF	1.62	0.04	-2	1	0.86	0.01

Table 3.4.1: Mean (μ) and standard deviation (σ) of the parameters of the error model and the regression score, computed on the entire time interval of drift, over the 16 nodes of the network. The observations that can be made with these values are identical to those of Table 3.3.1. This table also seems to confirm that there is a bias for SVR-KF as the standard deviation of the slope and intercept among instruments is small, with mean slope close to 1 but a mean intercept equal to 10. However, the poor mean score of the regression invites to be careful with this statement.

	Slope		_	Intercept		Score	
	s_1	s_2		s_1	s_2	s_1	s_2
No calibration	1.77	1.58		-2	-1	0.85	0.89
$\mathbf{SM}\text{-}\mathbf{SVD}$	1.01	1.05		0	1	0.99	0.97
$\mathbf{SM}\text{-}\mathbf{LS}$	0.97	0.94		0	1	0.99	0.97
$\mathbf{SM}\text{-}\mathbf{TLS}$	1.01	1.05		0	1	0.99	0.97
AB-DT	1.62	1.55		0	-1	0.87	0.87
AB- KF	1.63	1.56		-1	-3	0.87	0.86
SVR- KF	1.09	1.01		9	12	0.69	0.65
K-KF	1.62	1.56		-2	-4	0.87	0.84

Table 3.4.2: Values of the parameters of the error model and the regression score, computed on the entire time interval of drift, for two particular instruments of the network $s_1 = 2$ and $s_2 = 10$. The observations that can be made with these values are identical to those of Table 3.3.2.

	Slope		In	tercept	Sco	re
	μ	σ	μ	σ	μ	σ
No calibration	1.66	0.36	0	0	1.00	0.00
$\mathbf{SM}\text{-}\mathbf{SVD}$	1.03	0.04	0	0	1.00	0.00
$\mathbf{SM}\text{-}\mathbf{LS}$	0.99	0.03	0	0	1.00	0.00
$\mathbf{SM}\text{-}\mathbf{TLS}$	1.03	0.04	0	0	1.00	0.00
AB-DT	1.53	0.28	2	1	1.00	0.00
AB-KF	1.54	0.30	0	1	1.00	0.00
SVR- KF	1.35	0.28	0	17	0.88	0.06
K-KF	1.51	0.32	0	0	1.00	0.00

Table 3.4.3: Mean and standard deviation of the parameters of the error model and the regression score for s_1 , computed on each week of the drift period. The slope for SVR-KF varies significantly and the mean intercept is equal to zero with a standard deviation of 17. This is not the case for the other strategies. Thus, the observation made with Table 3.4.2 on SVR-KF is not valid locally and can explain why the mean score is poor compared to the ones obtained with the other strategies.

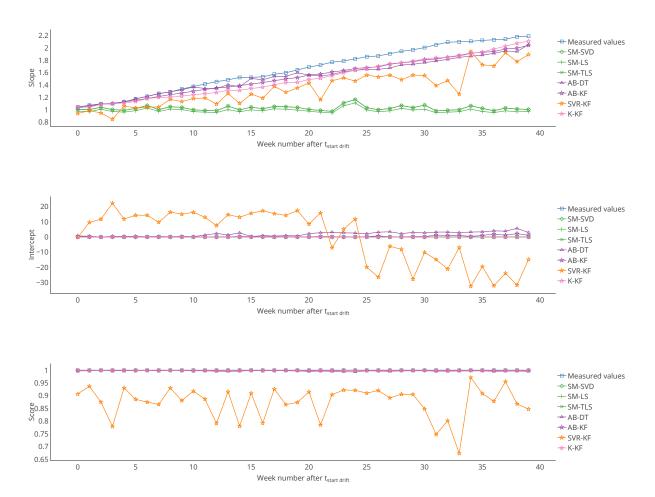


Figure 3.4.1: Evolution of the slope, intercept and score of the error model of s_1 , computed on each week of the drift period. Both the slope and intercept are locally poor for SVR-KF, although this is not visible when computing the error model on the entire time interval of study. This explains the poor values of the score compared to the other strategies.

4.3 Means of visualisation

Choosing other ways of visualising results may help to better understand what happens to each instrument. Various representations (Figures 3.4.2, 3.4.3 and 3.4.4) are discussed in terms of suitability for the comparison of methods.

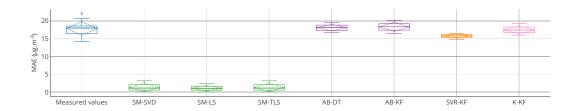


Figure 3.4.2: Box plots of the MAE, computed on the entire time interval of drift, of the 16 nodes of the network without calibration and with **SM-SVD**. It is a graphical representation of the information displayed in Table 3.3.1 for MAE. Supplementary information is provided compared to the table: quartiles, minimum and maximal values.

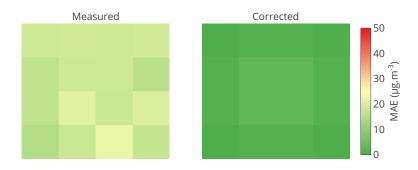


Figure 3.4.3: Matrix of the MAE, computed on the entire time interval of drift, of the 16 nodes of the network without calibration and with **SM-SVD**. It shows exactly the information for each instrument. The colour scale can help to identify problematic instruments.

Figure 3.4.2 is a boxplot-based representation of the values of MAE computed on the entire time interval of study. A boxplot is provided for each strategy. This representation is a graphical illustration of the information displayed in Table 3.3.1 for MAE. In addition, it provides supplementary information compared to the table: quartiles, minimum and maximal values. However, this representation does not show exactly the results for each instrument. One graph is needed for each metric. The plot can also show statistics on the evolution of a metric for a particular instrument like Table 3.3.3 but again it can only display information about one metric.

To display values regarding each instrument of a network, a matrix of values can be a solution. A colour scale can make the values more easily readable. Figure 3.4.3 is a matrix of the mean of MAE on the entire time interval of study. This figure shows exactly the information for each instrument. The colour scale helps to identify problematic instruments. It is appropriate for real-time display. Nevertheless, if multiple metrics (or variables if an error model is considered) are used, one matrix must be plotted for each one. The study of multiple calibration strategies is not possible on a single plot.

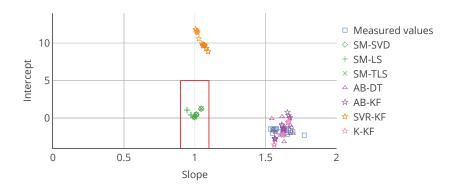


Figure 3.4.4: Target plot of the 16 nodes as a function of their slope and intercept in the error model for each calibration strategy. The slope and intercept are computed on the entire time interval of drift. In this case, it allows to locate an instrument according to the slope and intercept of its associated error model. The red rectangle is an example of area that can be defined to quickly identify instruments that are not satisfying a requirement, here a slope in [0.9; 1.1] and an intercept in [-5; 5].

To overcome the challenge of displaying information for multiple instrument on a single chart, we propose to use target plots. They enable locating a point along multiple axes (two for a 2D diagram, three for a 3D diagram). Boundaries within which the point should be ideally can be depicted. Each axis can be used for a different metrics. A colour scale can be used to depict a third one. In the previous section, we proposed to use an error model with two variables. Therefore, a target plot is perfectly adapted. Figure 3.4.4 is a target plot of the parameters of the error model, each point's coordinates being the values of the slope and the intercept of the error model for each instrument, computed on the entire time interval of study. Indeed, this diagram shows exactly the information for each instrument. Moreover, statistics like standard deviation can be added with error bars. The target diagram is also appropriate for real-time display and results for multiple calibration strategies can be plotted. Thus, it is a very powerful representation, although it is limited to two or three variables, respectively for a 2D plot and a 3D plot (potentially three or four with a colour scale on the markers) to allow a graphical representation, and only one statistic per axis can be added with error bars.

4.4 Conclusion

In this section, we investigated further issues regarding the evaluation of performances of *in situ* calibration strategies in terms of metrics and display. The conclusions are as follows:

- In the case of an evaluation based on usual metrics, the study with multiple metrics is strongly recommended, with a careful regard on the purpose of each metric. Possible interdependence and underdetermination between them 19 may mislead interpretations. We recommend using an error model-based approach, which allows to better capture the correction brought by a calibration algorithm, for instance in terms of gain and offset.
- The way metrics (or the parameters when an error model is used) are computed (time range on which the calculation is performed, periodicity of the computation) should be consistent with the details of the case study.

¹⁹Note that the set of metrics that is used in this work (or a subset) is not optimal itself regarding these questions. See [154]

- Computing statistics of metrics over the whole network is advantageous for large sensor networks, but may hide important information regarding to individual instrument (for instance, a specific device degrading the overall performances of a given method due to outliers)
- Graphical visualisations of metrics (or parameters of an error model) can enable a better
 analysis, notably with a target plot of the model's parameters when an error model with
 two or three parameters is used.

5 Sensitivity of the calibration algorithms to the specificities of the case study

In this section, we discuss the influence of the solutions chosen for the implementation of each step of the framework on the performances of the algorithms studied. First the function used to model the concentration of pollutant is modified. Then, the number of instruments is changed. Afterwards, the influence of the drift model is investigated. Finally, the impacts of a change of parameters for the calibration algorithms are studied. For comparison purposes, the metrics chosen in Section 3 are used, completed with the error model and the target plot proposed in Section 4 when appropriate.

5.1 Using a more realistic model of the true values

The model used in Section 3 is simplistic and can legitimately raise questions on the performances of each calibration strategy with a more representative model. In this section, consider the Gaussian Plume model [61, 145], a refined model with meteorological parameters, taken from a real-world data set. Otherwise, the space of study and its discretisation, the pollutant source location and duration of study, the sensor network, the instrument drift model, and the parameters of the calibration strategies are kept similar to the previous case study.

We suppose that at each time step, the pollutant dispersion to be in steady state. Concentration C at instant t and position (x, y, z) for a wind direction following x is expressed

$$C(x, y, z, t) = \frac{Q}{4\pi V_w(t)\sigma_y \sigma_z} e^{-\frac{y^2}{4\sigma_y^2}} \left(e^{-\frac{(z-H)^2}{4\sigma_z^2}} + e^{-\frac{(z+H)^2}{4\sigma_z^2}} \right)$$

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- σ_y and σ_z : respectively horizontal and vertical dispersion coefficients
- Q: Emission rate at the source
- V_w : Wind speed
- H: Pollutant effective release. $H = h_s + \Delta h(t)$ where

$$\Delta h(t) = \frac{1.6F^{\frac{1}{3}}x^{\frac{2}{3}}}{V_w} \text{ with } F = \frac{g}{\pi}D\left(\frac{T_s - T(t)}{T_s}\right)$$

- $-h_s$: pollutant source height
- -g: gravity constant
- -D: volumetric flow
- $-T_s$: source temperature
- -T: ambient temperature

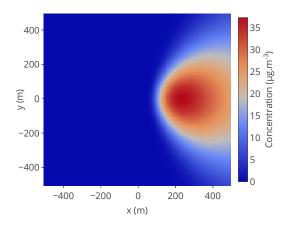


Figure 3.5.1: Example of concentration map used and based on the Gaussian plume model. The pollutant source is at coordinates $(x_s,y_s)=(0,0),\ h_s=25\mathrm{m},\$ the map is observed at $z=2\mathrm{m}.$ Other parameters are : $T=25^{\circ}\mathrm{C}, V_w=10\mathrm{m\,s^{-1}}, T_s=30^{\circ}\mathrm{C}, g=9.8\mathrm{m\,s^{-2}}, D=1.9\times10^{-9}\,\mathrm{m^3\,s^{-1}}, Q=5\times10^{-3}\,\mathrm{kg\,s^{-1}}, \sigma_y=1.36|x-x_s|^{0.82}, \sigma_z=0.275|x-x_s|^{0.69}.$ Wind direction is equal to 0° here.

To allow for wind direction changes, the cylindrical coordinate system is used to rotate the plume.

Emission rate, temperature of the source and dispersion coefficients are supposed constant over the simulation range. Ambient temperature, wind speed and wind direction are extracted from METAR weather reports collected over a year. An example of pollution map obtained for a given altitude z is represented in Figure 3.5.1.

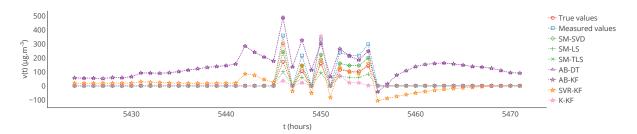


Figure 3.5.2: True values, measured values and corrected values with the strategies considered for a particular instrument between $t=5424\mathrm{h}$ and $t=5471\mathrm{h}$ with the Gaussian plume model. Note that the profiles of the curves are very different from those of Figure 3.3.4 as instruments are not necessarily exposed to the plume of pollutant due to the wind direction. In this case it is exposed to the pollutant between $t=5445\mathrm{h}$ and $t=5456\mathrm{h}$.

Figure 3.5.2 is a temporal plot for a particular instrument like in Section 3. From this plot we observe that the true signal appears to vary more quickly from high to low values. This is due to the fact that only one source of pollutant is considered in the area, and that the wind varies quickly from one direction to another in this dataset. Moreover, as we assumed that the concentration at each time step could be represented following the equation of a stationary Gaussian Plume, the pollutant quickly spreads at some points if the wind direction changes a lot. In practice the pollutant may not disperse as fast.

	\mathbf{M}	AE	MA	APE	RN	ISE	Pearson	
	μ	σ	μ	σ	μ	σ	μ	σ
No calibration	13	6	17	6	38	21	0.98	0.00
$\mathbf{SM}\text{-}\mathbf{SVD}$	2.30×10^3	6.12×10^3	2.28×10^5	6.13×10^5	1.39×10^4	3.73×10^4	0.55	0.43
$\mathbf{SM}\text{-}\mathbf{LS}$	26	24	952	2.42×10^3	93	138	0.85	0.34
SM-TLS	2.30×10^3	6.12×10^3	2.28×10^5	6.13×10^5	1.39×10^4	3.73×10^4	0.55	0.43
AB-DT	36	3	2.11×10^3	529	54	7	0.62	0.19
AB-KF	36	3	2.11×10^3	529	54	7	0.62	0.19
SVR- KF	249	357	1.51×10^4	2.15×10^4	270	380	0.54	0.18
K-KF	47	22	2.18×10^3	1.31×10^3	85	36	0.11	0.43

	Slope		Inter	Sco	re	
	μ	σ	μ	σ	μ	σ
No calibration	1.56	0.08	-2	1	0.97	0.01
$\mathbf{SM}\text{-}\mathbf{SVD}$	-21.11	60.40	2.49×10^3	6.99×10^3	0.48	0.42
$\mathbf{SM}\text{-}\mathbf{LS}$	0.34	0.32	10	28	0.83	0.33
$\mathbf{SM}\text{-}\mathbf{TLS}$	-21.11	60.40	2.49×10^3	6.99×10^3	0.48	0.42
AB-DT	0.82	0.31	19	13	0.42	0.24
AB-KF	0.82	0.31	19	13	0.42	0.24
SVR- KF	0.83	0.19	234	357	0.32	0.19
K-KF	0.26	0.66	30	15	0.18	0.27

Table 3.5.1: Mean and standard deviation of each metric, computed on the entire time interval of drift, over the 16 nodes of the network with the Gaussian plume model. No calibration strategy allows to obtain less error than without calibration, even for SM-(SVD, LS, TLS) strategies that had the best performances in Section 3.

Results with usual metrics and statistics on the parameters of the linear error model are presented in Table 3.5.1 on the entire network. From this table, we observe first that no calibration strategy provides less error than without calibration, even for SM-(SVD, LS, TLS), the strategies that had the best performances in Section 3. This is also the case in Table 3.5.2 for the same particular instruments considered in Section 3, despite apparently satisfying results according to MAE but not according to the slope and intercept notably. The results for each instrument are depicted in detail in Figure 3.5.3 with the target plot of the parameters of the error model for all the instruments.

Using a more realistic model of dispersion obviously produces a more complex signal that may require a higher density of instruments to capture the phenomenon. Indeed, the subspace size for SM-(SVD, LS, TLS) is equal to 11 (it was 2 in Section 3.2.1). This is still fitting the oversampling condition required for these algorithms, but it could explain the actual results: there may not be enough measuring instruments to capture all the features of the measurand. The same reasoning applies for SVR-KF which is also model based. For the other strategies, for which the prediction of true values is based on neighbouring nodes, the quality of the results which were already poor in Section 3 is even worse.

In conclusion, the influence of the model used to derive the true values of the measuring instruments is important. Considering cases on urban air pollution monitoring, the dispersion of

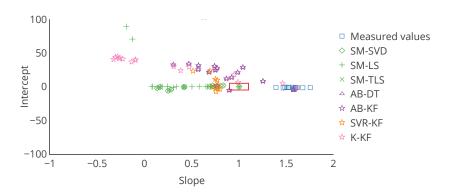


Figure 3.5.3: Target plot of the 16 nodes as a function of their slope and intercept in the error model, computed on the entire time interval of drift, for each calibration strategy and with the Gaussian plume model. The instruments are not all present, notably for those after the correction by SVR-KF. Axis were truncated to keep the plot readable. The red rectangle depicts the same requirement as in Figure 3.4.4 e.g. a slope in [0.9; 1.1] and an intercept in [-5; 5]. Few instruments are inside this area whatever the *in situ* calibration strategy considered.

pollutant for cities with different street layout may lead to different concentration fields despite identical pollutant sources and experimental conditions. Therefore, even with an identical sensor network with instruments drifting in the same way, *in situ* calibration strategies could produce corrections of a variable quality.

	M	AE	MA	PE	RI	MSE	Pea	Pearson		
	s_1	s_2	s_1	s_2	s_1	s_2	s_1	s_2		
No calibration	12	12	19	13	36	35	0.98	0.98		
$\mathbf{SM}\text{-}\mathbf{SVD}$	6	33	108	201	24	269	0.83	0.06		
SM-LS	8	25	55	86	20	57	0.97	0.97		
$\mathbf{SM}\text{-}\mathbf{TLS}$	6	33	108	201	24	269	0.83	0.06		
AB-DT	37	32	2.59×10^3	1.61×10^3	51	45	0.57	0.79		
AB- KF	37	32	2.60×10^3	1.61×10^3	51	45	0.57	0.79		
SVR- KF	24	387	1.31×10^3	2.60×10^4	34	415	0.69	0.34		
K-KF	27	63	818	3.41×10^3	55	100	0.60	-0.25		

	Slope			Intercept		Sco	re
	s_1	s_2	-	s_1	s_2	s_1	s_2
No calibration	1.75	1.47		-1	-2	0.96	0.97
$\mathbf{SM}\text{-}\mathbf{SVD}$	0.83	0.25		2	-6	0.68	0.00
$\mathbf{SM}\text{-}\mathbf{LS}$	0.58	0.17		0	-0	0.94	0.93
$\mathbf{SM}\text{-}\mathbf{TLS}$	0.83	0.25		2	-6	0.68	0.00
AB-DT	0.77	0.92		27	14	0.33	0.63
AB-KF	0.77	0.92		28	14	0.33	0.63
$\mathbf{SVR}\text{-}\mathbf{KF}$	0.78	0.98		3	382	0.48	0.12
K-KF	0.95	-0.26		18	42	0.36	0.06

Table 3.5.2: Values of the metrics, computed on the entire time interval of drift, for two particular instruments of the network $s_1 = 2$ and $s_2 = 10$, with the Gaussian plume model. No calibration strategy allows to obtain less error than without calibration, even for SM-(SVD, LS, TLS) strategies that had the best performances in Section 3, despite apparently satisfying results according to MAE but not according to the slope and intercept notably.

5.2 Density of the sensor network

The density of the network of sensors is crucial to capture fine spatial features of the phenomenon. In this section, we investigate the impact on the performances of the calibration strategies.

First, we replay the initial case study (2D Gauss dispersion model) with regular grid networks with one instrument every $\frac{1000}{n}$ m, $n \in [2..10]$. The size of the network, $|S| = n^2$, grows from 4 to 100. It was equal to 16 Section 3. All other parameters are kept.

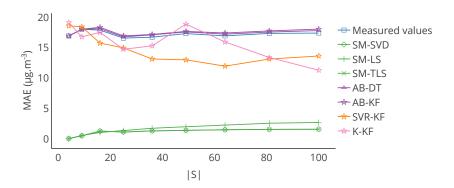


Figure 3.5.4: Evolution of the mean of the MAE, computed on the entire time interval of study, of all the nodes of the network, as a function of the number of nodes |S|, for the 2D Gauss model

Evolution of the mean of the MAE of the network with |S| is shown in Figure 3.5.4. From it, we observe that increasing the number of instruments improves the results with strategies K-KF and SVR-KF. However, the results are not significantly impacted by |S| for strategies AB-(DT, KF) and SM-(SVD, LS, TLS) compared to the baseline case (|S|=16). It can be explained by the simplicity of the model used to derive the true values. Moreover, for these strategies, slightly better results are obtained with a smaller number of instruments than with |S|=16. It seems that such a number (4 or 9 nodes) already enables capturing enough information to calibrate instruments for SM-(SVD, LS, TLS). For AB-(DT, KF), the prediction based on averaging values of neighbour instruments cannot be improved by increasing the number of instruments. Indeed, the network is blind, and instruments are all drifting at the same time, following the same model, with different values of drift. For SVR-KF, issues identified in Sections 3 and 4 seem to prevail despite an improvement of the performances with more instruments (but not with a lower number of nodes). However, in a general manner for K-KF, the more instruments are deployed, the more the kriging seems to provide accurate predictions. Results still have to be improved for |S|=100 though.

Thus, the number of instruments has an influence, though limited, on the results of the *in situ* calibration strategies. We also conjecture that the way instruments are deployed can have an influence, as it is crucial for an efficient reconstruction of the signal [20].

We also replay Gaussian Plume case study with a regular grid network of 100 instruments. We conjectured in Section 5.1 that increasing the density of instruments could improve on the poor calibration performances witnessed. Unexpectedly, our results contradict this intuition: the increased number of instruments does not improve the performances of the considered strategies.

Results with usual metrics and statistics on the parameters of the linear error model are presented in Table 3.5.3 for the entire network. From this table, we indeed observe that no calibration strategy allows obtaining less error than without calibration. In Table 3.5.4 results are quite equivalent to the previous ones in orders of magnitude. SM-(SVD, LS, TLS) strategies

	\mathbf{M}_{\cdot}	MAE		PE	RM	ISE	Pearson	
	μ	σ	μ	σ	μ	σ	μ	σ
No calibration	13	7	17	6	38	25	0.98	0.00
$\mathbf{SM}\text{-}\mathbf{SVD}$	1.04×10^{12}	1.04×10^{13}	1.04×10^{14}	1.04×10^{15}	6.32×10^{12}	6.32×10^{13}	0.55	0.37
SM-LS	4.77×10^8	4.77×10^{9}	4.77×10^{10}	4.77×10^{11}	2.91×10^{9}	2.91×10^{10}	0.91	0.26
$\mathbf{SM}\text{-}\mathbf{TLS}$	8.67×10^{11}	8.67×10^{12}	8.67×10^{13}	8.67×10^{14}	5.29×10^{12}	5.29×10^{13}	0.54	0.37
AB-DT	37	3	2.09×10^3	580	57	8	0.63	0.20
AB-KF	37	3	2.10×10^3	578	57	8	0.63	0.20
SVR- KF	412	208	2.68×10^4	1.33×10^4	432	232	0.51	0.18
K-KF	39	21	1.66×10^3	973	71	41	0.10	0.44

	Slope			Inter		Scor		
	μ	σ	-	μ	σ		μ	σ
No calibration	1.55	0.07		-2	1	0	.97	0.01
SM- SVD	-1.01×10^{11}	1.01×10^{12}		1.22×10^{12}	1.22×10^{13}	0	.43	0.38
$\mathbf{SM}\text{-}\mathbf{LS}$	-4.63×10^7	4.63×10^{8}		5.63×10^{8}	5.63×10^{9}	0	.89	0.26
$\mathbf{SM}\text{-}\mathbf{TLS}$	-8.43×10^{10}	8.43×10^{11}		1.02×10^{12}	1.02×10^{13}	0	.42	0.38
AB-DT	1.02	0.71		17	14	0	.43	0.25
AB-KF	1.02	0.71		17	14	0	.43	0.25
SVR-KF	0.82	0.14		415	208	0	.29	0.16
K-KF	0.19	1.38		23	11	0	.20	0.30

Table 3.5.3: Mean and standard deviation of each metric, computed on the entire time interval of study, over the 100 nodes of the network with the Gaussian plume model. Again, no calibration strategy allows to obtain less error than without calibration.

have abnormal performances, notably due to outliers in the results for some instruments. Figure 3.5.5 depicts the target plot of the parameters of the error model for all the instruments.

For the strategies AB-DT, AB-KF and K-KF, which perform the prediction of true values based on neighbouring nodes, the fact that instruments are drifting altogether may have a more important effect than the density of the network. For SM-(SVD, LS, TLS), the subspace of the signal, which is at the heart of these algorithms, was chosen considering the same criteria on the sum of the explained variance ratios. In this case, 13 components of the PCA with the highest variance ratios were kept (11 in Section 5.1). Thus, setting that a subspace is satisfying if the sum of the explained variance ratios of its components is greater than a threshold may not be sufficient to define it properly. The same reasoning applies for SVR-KF which is also model based.

	MAE		MA	PE	R	RMSE Pearso		
	s_1	s_2	s_1	s_2	s_1	s_2	s_1	s_2
No calibration	10	10	22	22	23	21	0.98	0.99
$\mathbf{SM}\text{-}\mathbf{SVD}$	5	3	36	25	13	6	0.92	0.99
$\mathbf{SM}\text{-}\mathbf{LS}$	6	4	42	30	11	6	0.99	1.00
$\mathbf{SM}\text{-}\mathbf{TLS}$	5	3	36	25	13	6	0.92	0.99
AB-DT	37	37	2.16×10^3	2.08×10^3	53	54	0.59	0.63
AB- KF	38	37	2.17×10^3	2.09×10^3	53	54	0.59	0.63
$\mathbf{SVR}\text{-}\mathbf{KF}$	373	309	2.22×10^4	1.75×10^4	37	4 311	0.63	0.68
K-KF	9	10	44	62	18	22	0.95	0.90

	Slope		Intercept		Score		
	s_1	s_2	s_1	s_2	s_1	s_2	
No calibration	1.59	1.54	-1	-1	0.97	0.97	
$\mathbf{SM}\text{-}\mathbf{SVD}$	0.84	0.87	-0	-0	0.85	0.98	
$\mathbf{SM}\text{-}\mathbf{LS}$	0.71	0.84	-0	-0	0.99	0.99	
$\mathbf{SM}\text{-}\mathbf{TLS}$	0.84	0.87	-0	-0	0.85	0.98	
AB-DT	1.11	1.28	20	16	0.35	0.39	
AB- KF	1.11	1.28	20	16	0.34	0.39	
SVR- KF	0.89	1.02	375	309	0.39	0.46	
K-KF	1.31	1.29	1	1	0.91	0.81	

Table 3.5.4: Values of the metrics, computed on the entire time interval of study, for two particular instruments of the network $s_1 = 2$ and $s_2 = 10$, with the Gaussian plume model. For these instruments, the results are quite equivalent to those of Table 3.5.2 in orders of magnitude

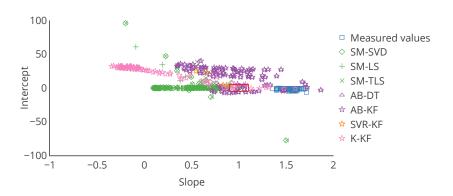


Figure 3.5.5: Target plot of the 16 nodes as a function of their slope and intercept in the error model, computed on the entire time interval of study, for each calibration strategy and with the Gaussian plume model. The instruments are not all present: axes were truncated to keep the plot readable. The red rectangle depicts the same requirement as in Figure 3.4.4 and 3.5.3 e.g. a slope in [0.9; 1.1] and an intercept in [-5; 5]. More instruments are inside this area compared to Figure 3.5.3 but most of them are outside whatever the *in situ* calibration strategy considered.

5.3 Instrument modelling

The modelling of the instruments used combines a simple drift model and the assumption that there are no other faults occurring to instruments than drifting. This hypothesis is not realistic, in particular when considering low-cost instruments. In this section, we introduce a more realistic drift model and evaluate the robustness of calibration strategies when the measurements are corrupted by other faults.

The model for the simulation of the true values, the sensor network and the parameters of the calibration strategies used in Section 3 are considered.

5.3.1 Drift model

To determine the influence of the drift complexity, two models are considered, with $t_{\text{start drift}}$ being the instant when the instrument begins to drift:

• Random gain and offset increase (**RGOI**): Gain $G(s_i, t)$ and offset $O(s_i, t)$ drift of instrument s_i are computed at each time step following:

$$G(s_i, t) = \begin{cases} 1 & \text{if } t < t_{\text{start drift}} \\ G(s_i, t - 1) + \delta G(s_i, t) & \text{if } t \ge t_{\text{start drift}} \end{cases}$$
with $\forall t, \delta G(s_i, t)$ drawn following $\mathcal{U}(0, \delta G_{max})$

$$O(s_i, t) = \begin{cases} 0 & \text{if } t < t_{\text{start drift}} \\ O(s_i, t - 1) + \delta O(s_i, t) & \text{if } t \ge t_{\text{start drift}} \end{cases}$$
with $\forall t, \delta O(s_i, t)$ drawn following $\mathcal{U}(0, \delta O_{max})$

with δG_{max} and δO_{max} being respectively the maximal gain and offset possible increase per time step.

• Continuous gain and offset increase (CGOI): A constant gain $\delta G(s_i)$ and offset $\delta O(s_i)$ increase of instrument s_i are drawn from uniform laws at the beginning of the drift and are added to respectively to the gain and offset of the instrument at each time step.

$$G(s_i, t) = \begin{cases} 1 & \text{if } t < t_{\text{start drift}} \\ G(s_i, t - 1) + \delta G(s_i) & \text{if } t \ge t_{\text{start drift}} \end{cases}$$
with $\delta G(s_i)$ drawn following $\mathcal{U}(0, \delta G_{max})$

$$O(s_i, t) = \begin{cases} 0 & \text{if } t < t_{\text{start drift}} \\ O(s_i, t - 1) + \delta O(s_i) & \text{if } t \ge t_{\text{start drift}} \end{cases}$$
with $\delta O(s_i)$ drawn following $\mathcal{U}(0, \delta O_{max})$

with δG_{max} and δO_{max} being respectively the maximal gain and offset possible increase per time step.

For both models, measured values are expressed following:

$$v(s_i, t) = G(s_i, t) \cdot v_{true}(s_i, t) + O(s_i, t)$$

In this case, $\delta G_{max} = 6 \times 10^{-5}$, and $\delta O_{max} = 0.03$.

There is an offset in RGOI and CGOI drift models. The estimation of the offset could not be carried out with SM-(SVD, LS, TLS) because the phenomenon does not have a null average on the give time span of calibration. Hence, the offsets are expected to be wrong for these strategies.

Results with usual metrics and statistics on the parameters of the linear error model are presented in Table 3.5.5 and 3.5.6. From these tables and compared to Tables 3.3.1 and 3.4.1, we observe that very different results can be obtained depending on the drift model and notably for SM-(SVD, LS, TLS). For these strategies, the measured values with RGOI or CGOI drift models are less improved by the calibration strategies even if we take into account that we expected to have incorrect offsets. Indeed, the average slope is worse than the one in Table 3.4.1, with an equivalent standard deviation. For both drift models, SVR-KF seems to improve the measured values according to MAE, MAPE, RMSE and the Pearson correlation coefficient but not according to the error model with a poor slope (0.76) and a high intercept (20) in average. AB-(DT,KF) and K-KF give corrected values equivalent to the measured values with the RGOI model. They are slightly improved if the measured values are built with the CGOI model and more notably for K-KF.

Thus, the drift model used to derive the measured values can drastically change the corrected values obtained after calibration depending on the algorithm used.

	MAE		MA	PE	$\mathbf{R}\mathbf{N}$	RMSE Pear			son
	μ	σ	μ	σ	$\overline{\mu}$	σ		μ	σ
No calibration	56	0	273	46	65	0		0.65	0.00
$\mathbf{SM}\text{-}\mathbf{SVD}$	50	3	248	32	56	3		0.67	0.01
$\mathbf{SM}\text{-}\mathbf{LS}$	46	1	232	43	52	1		0.68	0.00
$\mathbf{SM}\text{-}\mathbf{TLS}$	50	3	248	32	56	3		0.67	0.01
AB-DT	56	0	274	48	65	0		0.65	0.00
AB-KF	56	1	276	54	65	1		0.64	0.00
SVR- KF	14	0	86	13	16	1		0.84	0.01
K-KF	56	1	276	54	65	1		0.64	0.00

	\mathbf{Slope}		Intercept		\mathbf{Score}		
	μ	σ	μ	σ	μ	σ	
No calibration	1.41	0.00	43	1	0.42	0.01	
M-SVD	1.24	0.01	42	2	0.45	0.01	
$\mathbf{SM}\text{-}\mathbf{LS}$	1.18	0.04	40	0	0.46	0.00	
$\mathbf{SM}\text{-}\mathbf{TLS}$	1.24	0.01	42	2	0.45	0.01	
AB-DT	1.41	0.01	43	1	0.42	0.00	
AB- KF	1.41	0.03	43	1	0.42	0.00	
SVR- KF	0.76	0.02	20	1	0.71	0.02	
K-KF	1.41	0.03	43	1	0.41	0.00	

Table 3.5.5: Mean and standard deviation of each metric, computed on the entire time interval of study, over the 16 nodes of the network with the RGOI drift model. Results are very different with this model compared to Tables 3.3.1 and 3.4.1, notably for SM-(SVD, LS, TLS) even if we take into account that we expected to have incorrect offsets.

	MAE		$\mathbf{M}A$	PE		RMSE Pears			son	
	μ	σ	μ	σ	-	μ	σ	_	μ	σ
No calibration	44	29	210	131		51	34		0.75	0.15
$\mathbf{SM}\text{-}\mathbf{SVD}$	33	21	162	99		38	24		0.78	0.14
$\mathbf{SM}\text{-}\mathbf{LS}$	26	16	133	89		31	19		0.77	0.13
$\mathbf{SM}\text{-}\mathbf{TLS}$	33	21	162	99		38	24		0.78	0.14
AB-DT	46	1	227	44		54	1		0.70	0.01
AB-KF	47	1	228	46		54	1		0.70	0.00
SVR- KF	14	0	85	13		16	0		0.84	0.01
K-KF	20	1	95	22		23	1		0.89	0.00

	Slope		Inte	ercept	S	Score		
	μ	σ	μ	σ	μ	ļ.	σ	
No calibration	1.34	0.20	33	23	0.5	58	0.23	
$\mathbf{SM}\text{-}\mathbf{SVD}$	1.18	0.12	27	17	0.6	63	0.21	
$\mathbf{SM}\text{-}\mathbf{LS}$	1.07	0.13	21	14	0.6	61	0.20	
$\mathbf{SM}\text{-}\mathbf{TLS}$	1.18	0.12	27	17	0.6	63	0.21	
AB-DT	1.35	0.03	35	1	0.4	19	0.01	
AB- KF	1.35	0.03	36	1	0.4	19	0.00	
$\mathbf{SVR}\text{-}\mathbf{KF}$	0.76	0.02	20	1	0.7	71	0.02	
K-KF	1.18	0.02	14	1	0.7	79	0.01	

Table 3.5.6: Mean and standard deviation of each metric, computed on the entire time interval of study, over the 16 nodes of the network with the CGOI drift model. Results are very different with this model compared to Tables 3.3.1 and 3.4.1, notably for SM-(SVD, LS, TLS) even if we take into account that we expected to have incorrect offsets.

5.3.2 Other faults

As stated in Chapter 1 Section 2.2, measuring instruments are subject to other faults but drift. Even if calibration algorithms are not capturing these faults, the quality of input data impacts the performances of calibration. In [50], it was shown that the outliers contained in the time series of measured values have a more important influence on the calibration results than the average quality of the data. As an early work on the topic, Ramanathan *et al.* [123] observed that the measured values were sometimes corrupted, even with an individual field calibration for each instrument performed with a mobile chemistry laboratory at the time of deployment. In this section, we investigate the influence of other faults on the calibration results.

Let us assume that in addition to a WGLI drift, all the instruments are undergoing after $t_{\rm start\ drift}$:

Noise We model it as a random variable following a normal law. For each instrument s_i , the value of the noise $\varepsilon(s_i, t)$ is drawn from $\mathcal{N}(0, \varepsilon_{max})$ at each time step, with $\varepsilon_{max} = 20$ here

Spikes In this study, a spike occurs depending on the random variable $p_{\psi}(s_i, t)$ that follows $\mathcal{U}(0, 1)$. If $p_{\psi}(s_i, t) < 0.05$, a spike is added to the measured value $v(s_i, t)$. The value of the spike is equal to $(\psi \cdot v)(s_i, t)$, with $\psi(s_i, t)$ following $\mathcal{U}(-1, 1)$.

Thus, the measured values of an instrument s_i are equal to:

$$v(s_i, t) = \begin{cases} G(s_i, t) \cdot v_{true}(s_i, t) & \text{if } t < t_{\text{start drift}} \\ G(s_i, t) \cdot v_{true}(s_i, t) + \varepsilon(s_i, t) & \text{if } t \ge t_{\text{start drift}} \text{ and } p_{\psi}(s_i, t) \ge 0.05 \\ (G(s_i, t) + \psi(s_i, t)) \cdot v_{true}(s_i, t) + \varepsilon(s_i, t) & \text{if } t \ge t_{\text{start drift}} \text{ and } p_{\psi}(s_i, t) < 0.05 \end{cases}$$

Results with usual metrics and statistics on the parameters of the linear error model are presented in Table 3.5.7. While SM-(SVD, LS, TLS) strategies worked very well for instruments suffering from drift only, their performances degrade significantly after introducing these faults in the measured values. For the other strategies, noise and spikes do not seem to degrade the results although the performances were fairly unsatisfying from the start like in Section 3. It is expected since this kind of fault is not considered by the calibration strategies. It is henceforth less a matter of performance than an issue of robustness of each method against abnormal values. In practice, one should try to identify and correct as many faults as possible, e.g. by filtering spikes and averaging noise, before applying an *in situ* calibration algorithm.

	MAE		\mathbf{M}	APE	F	RMSE P			Pear	Pearson	
	μ	σ	$\overline{\mu}$	σ		ι	σ		$\overline{\mu}$	σ	
No calibration	25	2	100	5 14	3	3	2		0.79	0.01	
$\mathbf{SM}\text{-}\mathbf{SVD}$	60	50	198	3 144	8	2	70		0.23	0.33	
$\mathbf{SM}\text{-}\mathbf{LS}$	28	6	93	10	3	3	7		0.35	0.57	
$\mathbf{SM}\text{-}\mathbf{TLS}$	60	50	198	3 144	8	2	70		0.23	0.33	
AB-DT	21	1	78	17	2	8	2		0.88	0.01	
AB- KF	21	1	78	17	2	8	2		0.88	0.01	
SVR-KF	18	0	92	15	2	2	0		0.75	0.01	
K-KF	25	1	10'	7 18	3	2	1		0.79	0.00	

	\mathbf{Slope}		In	tercept	Sco	\mathbf{Score}		
	μ	σ	$\overline{\mu}$	σ	μ	σ		
No calibration	1.62	0.07	-2	0	0.62	0.01		
SM-SVD	0.18	0.69	-13	B 14	0.16	0.25		
$\mathbf{SM}\text{-}\mathbf{LS}$	0.15	0.26	-0	1	0.43	0.08		
$\mathbf{SM}\text{-}\mathbf{TLS}$	0.18	0.69	-13	B 14	0.16	0.25		
AB-DT	1.63	0.04	-2	1	0.77	0.02		
AB- KF	1.63	0.04	-1	1	0.78	0.02		
SVR- KF	1.06	0.03	10	1	0.56	0.01		
K- KF	1.61	0.04	-2	1	0.62	0.01		

Table 3.5.7: Mean and standard deviation of each metric, computed on the entire time interval of study, over the 16 nodes of the network with the WGLI drift model, spikes and noise. The results with SM-(SVD, LS, TLS) are significantly influenced by the spikes and noise. The other strategies do not seem to provide more degraded results due to the noise and spikes although the performances were fairly unsatisfying from the start in Section 3.

5.4 Parameters of calibration strategies

Finally, the last step that can influence the results is the configuration of the calibration strategies. Indeed, the strategies considered in Section 3 have multiple parameters.

Multiple studies can be performed for each parameter of each strategy we considered. As an example, we only investigate the case of SM-(SVD, LS, TLS). The influence of the frequency of computation and of the duration of the time range used to select the measured values used is studied. Note that this aspect was not treated in the original publication [7]. We consider here that the period of computation and the duration of the time range for the selection of the measured values are equal and denoted w. The same environment, sensor network and drift model as in Section 3 are considered.

Figure 3.5.6 represents the evolution of the mean of MAE over the network for the considered strategies as a function of w. The best result is obtained with w = 7, which was our original value in Section 3, considering the average of the error for the three strategies with the same w. Nevertheless, changing the value of w may lower the improvement brought by the $in\ situ$ calibration for low and high w but the corrected values are still closer to the true values compared to the measured values.

We also changed the criteria to determine the signal subspace. The sum of the explained variance ratio of the components obtained by PCA must be greater $1-1\times 10^{-30}$ instead of $1-1\times 10^{-3}$ previously. This resulted in a change of the subspace with four components considered instead of two. The same study with varying w was conducted and results are shown in Figure 3.5.7.

We still observe that the best results are obtained for w=7: for most of the values of w, the algorithms degrade the quality of the measurements. This is even more significant for SM-SVD and SM-TLS. We, however, do not claim that w=7 is an optimal parameter since we only showcase this feature of the framework: it enables a full parametric optimisation of the methods. Calibration methods are indeed very sensitive to the adjustment of several parameters and a comparative study should include a parametric optimisation.

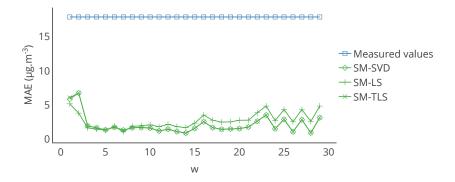


Figure 3.5.6: Evolution of the mean of MAE, computed on the entire time interval of study, over the network for the strategies SM-SVD, SM-LS and SM-TLS as a function of w. The best result is obtained for w=7, which was our original value in Section 3, considering the average error of the three strategies for the same w.

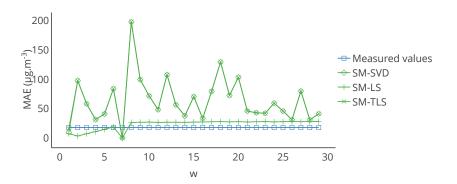


Figure 3.5.7: Evolution of the mean of MAE, computed on the entire time interval of study, over the network for the strategies SM-SVD, SM-LS and SM-TLS as a function of w with a different signal subspace than in Figure 3.5.6. The best results are still obtained for w=7 but for most the values of w, the algorithms degrade the quality of the measurements and more significantly for SM-SVD and SM-TLS.

5.5 Summary of the results

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In this section, we changed several parameters independently: the model of the phenomenon, the number of instruments, the faults introduced, notable the drift model, and the parameters of the calibration strategies for SM-(SVD, LS, TLS) that gave the best results previously.

Method	Baseline case	With Gaussian plume model	Varying number of nodes	Drift RGOI	Drift CGOI	Drift WGLI + faults	Varying calibration parameters
SM-SVD							
SM-LS							
SM-TLS							
AB-DT							
AB-KF							
SVR-KF							
K-KF							
Values improved Values not significantly Improved or degraded Values degraded Case not studied							

Figure 3.5.8: Summary of the results observed by changing the parameters of the baseline case study which was built with a 2D Gaussian function to model the measurand, with the WGLI drift model applied to the measured values of the instruments, with 16 nodes in the network and the values indicated in Section 3 for the parameters of the *in situ* calibration algorithms. The improvements are judged relatively to results of the baseline case, recalled in the first column. Thus, this table does not indicate that the corrected values are very close to the true values when the measured values are improved.

The results obtained are reported in Figure 3.5.8. From a general perspective, we observed that: 2652

- increasing the complexity of the model of the phenomenon drastically worsen the results for all the strategies
- changing the number of instruments between |S| = 4 and |S| = 100 in our case does not have a determining effect.
- the model of the drift strongly impacts the results
- the presence of other faults in the measured values can degrade the results
- the (correct) adjustment of parameters of calibration strategies may have a significant effect

Therefore, comparing in situ calibration strategies requires multiple precautions on how the corrected values were obtained and on what the measured values were based to provide fair and correct conclusions.

6 Discussion and conclusion

 In this chapter, a framework for the simulation of sensor networks, which enables a systematic comparison of *in situ* calibration strategies with reproducibility, and scalability, was presented. We have showcased this methodology by comparing several calibration strategies for blind and static sensor networks. Our results provide several engineering insights on the design of a sensor network.

Based on the methodology described in Section 2, we have proposed a case study in Section 3 concerning blind and static sensor networks. Although we have shown that several strategies have better performances than the others, these results have to be tempered as the conclusions depends on how the evaluation is conducted. We have proposed solutions to conduct a balanced assessment of the performances in terms of metrics used in Section 4. It is still challenging for sensor networks with a high number of instruments that require multiple scale analysis. In addition, we have shown in Section 5 that results may not be the same depending on choices made during the design of the case study regardless of how the evaluation is carried out. We have developed variations of our initial case study by changing the model computing the true values of the measurand, the number of instruments, the fault model of the instruments and parameters of calibration strategies.

Our results highlight the dependence of the performances of *in situ* calibration algorithms to the case study. Besides, these performances seem to be often limited, even on the relatively simple environmental models discussed here. The latter explains why more complex and accurate environmental models were not tested here. Finding the good trade-off between simplicity and realism of the simulation models is important. At first sight, the goal could be to consider models as realistic as possible everywhere. However, the accuracy of simulation-based environmental models remains challenging to establish, and costly to run, while simplified models such as those used in the paper may be sufficient when the goal is to compare methods to single out the most promising ones among several for a predefined case study. This is also true for the models to simulate the measuring chain and the deployment of the instruments.

It allows to point out an additional advantage to the framework proposed in this paper: its suitability as a design tool for sensor networks. For instance, after having determined the number of instruments (and their positions) required to cover a given area with a static sensor network, the best *in situ* calibration strategy and its parameters could be identified among a set of algorithms applicable to this particular network. In such cases, however, a stronger focus on the accuracy of the environmental model should be put whenever possible. Beyond the quality of the models, the adjustment of the strategies is crucial for an objective evaluation. The reproducibility provided by the framework is an asset to conduct a thorough parametric optimisation and compare the best results of each calibration.

To conclude, we can point out the following engineering insights on the design of a sensor network and the evaluation of *in situ* calibration strategies.

- There is actually no method that is universally outperforming the other. The best calibration strategy to apply depends on the deployment of the instruments and the exploitation of the measurements.
- Even on the same scenario, two strategies can outperform each other depending on which metrics of performance is considered and the way they are computed. The metrics and their means of visualisation have to be chosen accordingly to the target exploitation of the sensor network, to focus on the relevant features. We also advise using an error model in order to get meaningful details on the performances.
- Increasing the density of the deployment of instruments does not always lead to better performances of a calibration strategy.

• In some case, the quality of data can be degraded by a calibration strategy. Besides, other faults than drift will happen, in particular when considering low-cost instruments with low-quality electronics. To cope with robustness issues of the calibration strategy, it seems relevant to process these errors upstream.

The content of this chapter is based on the following publication:

F. Delaine, B. Lebental and H. Rivano, "Framework for the Simulation of Sensor Networks Aimed at Evaluating In Situ Calibration Algorithms," in *Sensors*, vol. 20, no. 16, 2020, DOI: 10.3390/s20164577.

Codes and data files used to produce the results of Sections 3, 4 and 5 are available online under licence AGPL 3.0 and ODbL 1.0 respectively in [37]

Chapter 4

Diagnosis of Drift Faults in Sensor **Networks**

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1 Motivations

Across the study of all the *in situ* calibration strategies in Chapter 2, and particularly regarding those considered in Chapter 3, most strategies perform the calibration of all instruments without evaluating first whether they actually need it. Moreover, we observed the algorithms may not improve, or may even degrade, the quality of the measurements. Two guidelines for *in situ* calibration algorithms can be established based on that. First, instruments providing measurement results agreeing with their specifications, should not be affected by an *in situ* calibration strategy. Secondly, the need for a calibration operation must be explicitly determined. This is particularly important if the execution of the algorithm is expensive, e.g. in terms of energy consumption in the case of autonomous nodes. It would add more confidence in the algorithms, namely it could inform it recalibrated an instrument because a criterion on its measurement results was met.

Some calibration algorithms are actually providing the identification of instruments needing a calibration, notably in [87, 168]. In these two publications, the algorithm determines first whether an instrument has drifted since the last time step. If it is the case, it gives the identity of the instrument that has drifted and correct it with the help of all the other instruments. However, if more than one instrument is drifting at a time, the algorithm raises an error. This is problematic as such a case is very likely in practice.

In the same vein of assumptions that are difficult to meet in practice, multiple strategies are designed based on the hypothesis that the sensor network is "dense". It means that each instrument is within few tens of meters to one or several nodes of the network. The need for this assumption to be verified can be explained as follows. To correct drifting instruments, their correct values—or the values of the drift—must be estimated. To do so, mainly two approaches are considered: the estimation from values measured by neighbouring instruments or the estimation from a model of the measurand. This is the case whether the network is static or not and blind or not.

Regarding the first approach, instruments should be physically close enough to have comparable values between them. This is something that can be easily achieved in practice and the reason why the assumption of a dense sensor network is often proposed. However, if the distance between the instrument to recalibrate and the instruments used to estimate its value is too important compared to the characteristic spatial length of variation of the measurand, and if it cannot be reduced, it may not be possible to have an accurate estimation with the desired uncertainty. Thus, as complex quantities are targeted in environmental monitoring, comparisons may be feasible for mobile sensor networks, but it is less likely to be the case for static sensor networks. Indeed, with mobile instruments can sometimes meet whatever the number of nodes in the network. For static ones, the number of instruments may have to be significantly increased depending on the application so that values of different devices may be compared but with an associated cost that may be significant.²⁰

Therefore, a model-based approach for estimating the correct values of instruments may be preferred when meeting the required density of instruments is not possible, e.g. with static instruments close enough or with mobile nodes meeting often. With a model-based approach, it is assumed possible to model satisfyingly the measurand to estimate its value at several locations, with the help of a priori information and sometimes with measured values, preferably from

²⁰The case where neighbouring instruments are used to evaluate the values of an instrument could be extended to the case where the nodes of the network embed several instruments with the same characteristics and measuring the same quantity. In this case, the set of instruments of a node measuring the same quantity can be seen as a measuring system, which would be equivalent to having a single instrument of a higher quality. Indeed, it can be legitimately expected that the average of the values of these instruments is more accurate than their values considered individually. This comes down to the case described in this paragraph. For the sake of clarity, this particular situation is not discussed in detail in this manuscript.

reference instruments. Acquiring such information may not be easy or may require a long training period to capture enough features of the phenomena. The latter may not be compatible with the lifespan of the instruments, especially if they are low cost. The deployment of expensive reference instruments during the deployment phase is a solution but has a major cost impact, notably if a high number of instruments is required. Additionally, using a model for the measurand raises the question of the interest of deploying the sensor network itself. If the model is able to derive at the required spatiotemporal resolution the values of the considered phenomenon, even with the help of a few high-quality instruments, this may be sufficient for most applications.

Overall, these elements point out that most approaches for *in situ* calibration can be challenged by the fact that the reference values used for calibration, e.g. standard values, need to meet certain conditions. They are not necessarily met in *in situ* calibration algorithms. The same precautions should be taken regarding the identification of instruments that have drifted, whatever the approach chosen to determine them.

In this chapter, we investigate the feasibility of drift detection for measuring instruments with the help of measurement results from other devices. The concept of **rendez-vous** is used to achieve this goal and is formally defined in Section 3, based on existing works on the subject [40, 132]. The study of these rendez-vous is expected to help to determine if an instrument needs to be recalibrated as long as one of the instruments involved in the rendez-vous can be considered as a standard, e.g. an instrument of which results can legitimately be trusted. To remain the most generic as possible, **no assumption is made for this algorithm on the type of sensor network, on the density of instruments and on the measurand** such as a priori known information regarding it.²¹

Thus, a formal definition of the concept of rendez-vous is proposed first based on existing works. Which values of instruments may be compared to each other in practice is briefly discussed. Then, a diagnosis algorithm, that is a way to identify the instruments that need a correction of their calibration relationship, is provided. The approach is based on rendez-vous. Afterwards, a case study is presented to illustrate the application of the diagnosis algorithm. The means to improve the results, in terms of false results, are discussed followed by guidelines on the adjustment of the parameters of the algorithm. The sensitivity of the algorithm to the parameters of the case study is studied. Finally, the algorithm is combined to a simple calibration strategy showing promising results. Before moving to these contributions, a literature review is conducted regarding fault diagnosis applied to sensor networks and in particular regarding the detection of drift

2 State-of-the-art of fault diagnosis applied to sensor networks

2.1 Overview of fault diagnosis for sensor networks through existing reviews

In the last decade, while sensor networks gained in popularity, the question of their dependability became a major subject. In particular, multiple contributions were made on fault diagnosis as reported in various surveys [98, 106, 114, 185].

In these publications, the research works are classified by the way the diagnosis is performed. First, the place where the decision is computed, e.g. if it is a centralised, decentralised or hybrid approach, is considered. Then, the general mathematical approach used for the diagnosis (testing, comparisons, majority voting, statistics, probabilities, machine learning algorithms, fuzzy logic, clustering...) is studied.

Faults are generally characterised by "hard" or "soft" and as "permanent", "transient" or "temporary" [98, 106, 185]. Only Muhammed *et al.* [114] used a taxonomy similar to the one of Ni *et al.* [115] that inspired the one used in this manuscript. In particular, these authors report

²¹It could be for instance an actual correlation between the true values at two very distant points that could be exploited in a specific situation.

that the existing diagnosis algorithms are often targeting multiple faults. For those addressing the drift fault²², stuck-at or out-of-range faults are also detected by the same algorithm in the 15 associated references they cited.

In the following surveys [98, 114, 185], the diagnosis approaches reported concern mainly static sensor networks. Mahapatro et al. [98] listed eight algorithms that could be applied to static and mobile sensor networks.²³ Zhang et al. [185] reported two references explicitly dealing with mobile sensor networks [1, 28]. They also stated that approaches designed for static networks behave poorly when applied to mobile ones. This indicates that algorithms exploiting specificities related to networks with mobile nodes in the diagnosis approach have not been deeply investigated so far.

In addition, other subjects such as energy consumption, communication range needed, communication failures, or the amount of required neighbour nodes are taken into account in the surveys on fault diagnosis for sensor networks [98, 106, 114, 185]. Zhang et al. [185] underline it is challenging to design a diagnosis approach having satisfying performances or properties regarding all the features of interest for such an algorithm, e.g. for instance diagnosing faults well with low energy consumption, few communications between nodes and so on.

2.2 Existing methods of diagnosis compatible with the concept of rendez-vous addressing any fault

2.2.1 Introduction

 The methods used to carry out the diagnosis of any fault in sensor networks are diverse in terms of concepts and tools on which they are based on. According to the surveys cited [98, 106, 114, 185], different main ideas come out like the use of comparisons between instruments, the use of statistics and probabilities, as well as the use of machine learning models (regressors, classifiers...). As the goal in this chapter is to propose a diagnosis algorithm for sensor networks based on rendez-vous between the nodes, we focus here on approaches compatible with this idea, e.g. it consists at least in comparison of values measured by different instruments.

2.2.2 Related works

Chen et al. [29] introduced a faulty sensor detection algorithm consisting of four steps of evaluation. First, each instrument compares its measured value at the instant of the diagnosis with the measured values of its neighbours. If the difference between two instruments is higher than a first threshold, the deviation of the difference since the last comparison is computed. If it is higher than a second threshold, the test between the two instruments is marked as positive. In a second step, each instrument determines if it is likely good or faulty depending on the number of positive tests and its number of neighbours. This information is shared between the instruments so that based on its neighbours that are likely good, an instrument determines if it is actually good and share this result. Then, for the remaining undetermined instruments, their statuses are decided if all their neighbours are non-faulty and if all the initial tests gave the same results. If ambiguities remains, they are removed based on the likely statuses of the instruments. Xu et al. [179] proposed an extension of this work dealing with the particular case of tree-like networks to reduce the number of communications. To avoid intermittent faults, they also proposed to compute the initial test result based on multiple comparisons of values between two instruments, instead of a unique comparison. In the same way, Saha et al. [128] developed a similar algorithm but with comparisons for multiple quantities (measurand and remaining

²²In the publication of Muhammed *et al.* [114], calibration, gain, and offset faults correspond to what is notably a drift fault in this thesis.

²³Note that the subject of mobile sensor networks was new at the time when Mahapatro *et al.* [98] published their survey.

energy).

Ssu et al. [142] proposed an approach to diagnosis faults between sources and sinks. The main idea is to send a request through two paths and to compare the results obtained at the sink node. If the results are different, then at least one faulty path exists. The algorithm tries to identify it with the help of a third path and a majority voting procedure, but the faulty paths cannot always be identified. Such a method is particularly relevant to diagnose communication or hardware-related faults.

Lee et al. [83] presented a distributed algorithm to isolate faulty nodes. Initially, the nodes are all assumed as faulty. A comparison is made between the measured values of neighbour nodes and if the result is higher than a threshold, then the test is positive. If less than a predefined number of positive tests has been obtained, or if the test with a non-faulty node is negative, then the diagnosed instrument is non-faulty. The algorithm works either based on single comparisons between neighbours or with comparisons repeated multiple times.

Mahapatro et al. [97, 99] introduced a clustering-based diagnosis algorithm. The clustering part is used for the definition of the neighbours around cluster heads, the cluster heads being the instruments with the highest residual energy levels. They also compare the measured values between instruments of a cluster and a majority voting strategy is used to determine the state of the nodes.

Chanak et al. [28] developed a comparison-based scheme using a reference mobile sink node moving between the static nodes of the network. When it is close to a node, several diagnoses are performed to detect hardware and software faults. This also enables a low consumption of energy for the transmission of the measurement results as it is no longer necessary to communication with a distant gateway. The core of this contribution lies in the determination of an optimal path to meet with each node.

Luo et al. [92] proposed an approach using the concept of average consensus. Each instrument estimates first its status regarding the average consensus measured value build with the values of its neighbours. Each instrument also estimates the statuses of its neighbours. Then, the decisions of all the instruments are merged to make the final decision.

In reaction to contributions computing the average value of the neighbours of an instrument by weighting their values with the inverse of the distance between them, Xiao et al. [178] argued that the distance does not control alone the relationship between the values of two instruments, notably if a closer one is actually faulty. Thus, they propose to take into account an indicator of trustworthiness computed for each node of the network. This confidence value is used in voting procedures that are used to determine the status of instruments.

Ji et al. [69] also developed their algorithm around a weighted average of the values measured by an instrument. The weights are representing a confidence level associated to each instrument. The difference between the measured value of an instrument and the average is compared to a threshold. If it is greater than the threshold, the confidence level of the instrument is decreased and once it reaches zero, the instrument is reported as faulty.

This idea of trust between instruments has been extensively studied and extended, as did for instance Jiang et al. [70]. Their robust trust model is based on direct comparisons but also on third parties and recommendations, with concerns about the security of the communications between the instruments. Their trust framework is also adapted for mobile sensor networks. Wang et al. [164] exploited Petri nets to introduce a trust-based formal model aimed at detecting faults in sensor networks.

Sharma et al. [134] also exploited the idea of confidence in a method similar to the one of Chen et al. [29]. The main difference with Chen et al. is that in a first step each instrument analyses its own behaviour and that a confidence level is associated to the determined statuses.

Feiyue et al. [52] proposed to combine the information resulting of a self-evaluation and of comparisons with the neighbours of each instrument. In a first step, the reliable nodes are

determined through a majority voting procedure. Then, various metrics are computed and combined to make the decisions for the instruments that were not considered reliable in the first step.

2.2.3 Remarks

 In this section, the way the computation is conducted (centralised, decentralised, hybrid) was not discussed on purpose. Indeed, this topic impacts the number of communications required and the energy consumption of the method more than the actual result of the diagnosis procedure.

Another aspect is how the faults are managed after detection in a diagnosis algorithm. There are different visions. Considering the diagnosis of outliers for instance, Ottosen *et al.* [116] adopted a gap-filling strategy. Fang *et al.* [51] preferred a fault removal approach because this procedure is a data processing carried out before the calibration of the instruments of the network. As shown for the calibration of measuring instruments in Chapter 3, it may also be preferable to correct measurement results prior to the diagnosis of drift if any other fault may be present.

Finally, in the works presented here, the diagnosis is performed in most cases based on single values from a given set of instruments measured at the same time. Taking a decision on a short time range may not be correct, especially if drift is targeted. Indeed, drift is a fault increasing little by little and thus a diagnosis on a long sequence of measured values over time could be more appropriate.

2.3 Positioning of the contribution

From this review, we observed that the literature on fault diagnosis for sensor networks is rich and abundant. Static sensor networks have been widely studied while mobile sensor networks have been partially covered. Multiple types of faults are addressed, concerning various aspects of measuring systems, but drift is not specifically targeted in existing works. Various methods have been used to tackle the challenge of fault diagnosis, notably based on comparisons between instruments.

However, the validity of the comparisons is rarely discussed from a metrological perspective. It would be valuable to develop a new approach with considerations on the quality of the measurement results. Moreover, exploiting the concept of rendez-vous allows to take into account that it happens in a spatiotemporal vicinity. This concept is more sophisticated than the simple idea of comparison of measurement results.

In addition, a diagnosis algorithm not requiring any assumptions on the type of sensor network and its structure would be interesting. Indeed, the existing approaches are often targeting static sensor networks, which may imply a dense network in the case of a comparison-based strategy, and rarely networks with mobile nodes. Regarding rendez-vous, mobility is an important feature that can facilitate their occurrence. Because mobile nodes can meet spatially with other mobile or static nodes, a diagnosis approach based on these meeting points would allow more meaningful comparisons of values if a very high density of measuring instruments is not feasible. In this way, if the algorithm works whether the sensor network is static or has mobile node leaves to the users the choice of the most appropriate type of sensor network for their application.

3 Definition of concepts for a drift diagnosis algorithm based on rendezvous

Before the presentation of an algorithm for the diagnosis of drift in sensor networks, this section introduces the definitions of the concepts of validity and compatibility of measurement results, and of rendez-vous. A sensor network composed of a set of instruments S is considered here and in the following sections without any other assumption on its type and on the characteristics of its nodes.

3.1 Validity of measurement results

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In Chapter 1 Section 1.2, we recalled that measuring instruments cannot exactly provide true values. In addition, we indicated they may be different accuracy class, e.g. of different qualities, in Chapter 1 Section 3.

On top of these ideas, we introduce here the notion of the **validity of measurement results**. For instance, a measuring instrument cannot measure all the values of a quantity. It has a bounded measuring interval which is the set of values of a quantity "that can be measured by a given measuring instrument or measuring system with specified instrumental measurement uncertainty, under defined conditions" [14]. In addition, it cannot work properly under any operating conditions. Therefore, a measurement result with a value outside an instrument's measuring interval or which has been obtained outside of the normal conditions of operation of a device, may not be **valid**.

The conditions under which measurement results are valid are specific to each type of instrument. In practice, they are usually provided in the technical documentation or it is possible to determine them experimentally. In this chapter, the conditions under which results are considered as valid are explained in the case study.

The notation $M^*(s_i, (t, \Delta t))$ refers to the set of measurement results for s_i , over the time range $[t - \Delta t; t]$ that are metrologically valid, e.g. $v(s_i, t)$ in the measuring interval and so on depending on the definition of what a valid result is metrologically speaking.

3.2 Compatibility of measurement results

Definition 22. Consider $s_i \in S$. A measurement result $m(s_i, t)$ is compatible with true value if $v_{true}(s_i, t) \in [v(s_i, t) - \Delta v(s_i, t); v(s_i, t) + \Delta v(s_i, t)]$.

Definition 23. Consider $s_i \in S$. A measurement result $m(s_i, t)$ is upper non-compatible with true value if $v_{true}(s_i, t) < v(s_i, t) - \Delta v(s_i, t)$.

Definition 24. Consider $s_i \in S$. A measurement result $m(s_i, t)$ is lower non-compatible with true value if $v_{true}(s_i, t) > v(s_i, t) + \Delta v(s_i, t)$.

The set of measurement results compatible with true values of s_i obtained during $[t - \Delta t; t]$ is noted $M^{\approx}(s_i, (t, \Delta t))$. $M^+(s_i, (t, \Delta t))$ and $M^-(s_i, (t, \Delta t))$ are defined for upper and lower non-compatible measurement results respectively. It can be noticed that

$$M(s_i,(t,\Delta t)) = M^{\approx}(s_i,(t,\Delta t)) \cup M^+(s_i,(t,\Delta t)) \cup M^-(s_i,(t,\Delta t))$$

These definitions can be extended for the comparison of measurement results between different instruments.

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Definition 25. Consider s_i and s_j \in S. m(s_i,t) is compatible with m(s_j,t') if v(s_j,t') - \Delta v(s_j,t') \le v(s_i,t) - \Delta v(s_i,t) \le v(s_j,t') + \Delta v(s_j,t') or v(s_j,t') - \Delta v(s_j,t') \le v(s_i,t) + \Delta v(s_i,t) \le v(s_i,t') + \Delta v(s_i,t'). It is noted m(s_i,t) \approx m(s_j,t').
```

Definition 26. Consider s_i and $s_j \in S$. $m(s_i,t)$ is upper non-compatible with $m(s_j,t')$ if $v(s_i,t) - \Delta v(s_i,t) \geq v(s_j,t') + \Delta v(s_j,t')$. It is noted $m(s_i,t) > m(s_j,t')$.

Definition 27. Consider s_i and $s_j \in S$. $m(s_i,t)$ is lower non-compatible with $m(s_j,t')$ if $v(s_i,t) + \Delta v(s_i,t) \leq v(s_j,t') - \Delta v(s_j,t')$. It is noted $m(s_i,t) < m(s_j,t')$.

The sets of measurement results of s_i obtained during $[t - \Delta t; t]$ associated to these definitions are respectively noted $M^{\approx}(s_i \to s_j, (t, \Delta t)), M^+(s_i \to s_j, (t, \Delta t))$ and $M^-(s_i \to s_j, (t, \Delta t))$.

More restrictive criteria can be set to determine if measurement results are compatible either with their associated true values or with results of other instruments. For instance, it

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could be a minimal level of overlapping between $[v(s_i,t) - \Delta v(s_i,t); v(s_i,t) + \Delta v(s_i,t)]$ and $[v(s_j,t') - \Delta v(s_j,t'); v(s_j,t') + \Delta v(s_j,t')]$. However, the influence of these definitions are not studied afterwards as it is more a matter of specification for the quality of measurements than a matter of diagnosis.

3.3 Rendez-vous

3.3.1 Formal and practical definitions

A rendez-vous between two instruments is defined as follows.

Definition 28 (Rendez-vous). Two instruments s_i and $s_j \in S$ are considered in rendez-vous when they are in a spatiotemporal vicinity so that their measurement results can be compared. In practice, it means for both their measurement results $m(s_i, t)$ and $m(s_j, t')$ that the instruments were spatially close enough and the difference between the instants of measurement t and t' was small enough to actually measure the same quantity value.

 s_i being in a rendez-vous at t with s_j is noted $\varphi(s_i \to s_j, t)$. Respectively s_j being in a rendez-vous at t' with s_i is noted $\varphi(s_j \to s_i, t')$.

The concept of "close enough" instruments will be explained later in the present section. Regarding the instants of measurement, two different ones are defined, t and t', because s_i and s_j may not have the same measurement frequency or may not have synchronous clocks. Consequently, in general, the measurement results of two instruments in rendez-vous are not obtained exactly at the same time.

The set of the rendez-vous encountered by s_i with s_j during $[t - \Delta t; t]$ is noted $\Phi(s_i \to s_j, (t, \Delta t))$. By extension, $\Phi(s_i \to S, (t, \Delta t))$ is the set of rendez-vous between s_i and any other instrument of S during $[t - \Delta t; t]$.

First of all, with this definition of a rendez-vous, there is no reason that any of the instrument should be mobile or static. As long as the conditions regarding the spatiotemporal vicinity are respected, rendez-vous can happen.²⁴ Thus, the use of the concept of rendez-vous in practice requires the definition of what an acceptable spatiotemporal vicinity is for a rendez-vous.

A first possible definition can be based on a maximal distance Δl_{φ} allowed between the instruments and on a maximal acceptable time difference Δt_{φ} between the measurement results of the two instruments. Therefore, s_i is in rendez-vous with s_j at t if

$$\exists t' \in \mathbb{R}^+$$
 such as $\exists m(s_j, t')$ and $|t - t'| \leq \Delta t_{\varphi}$ and $|l(s_i, t) - l(s_j, t')| \leq \Delta l_{\varphi}$

with $l(s_i, t)$ and $l(s_j, t')$ being the positions of s_i at t and s_j at t'. The temporal condition allows multiple rendez-vous with a single instrument for a same measurement result if t' is not unique. In this case, t' can be chosen as the value reaching the minimum of |t - t'| or of $|l(s_i, t) - l(s_j, t')|$. This choice of a unique measurement result for s_j is necessary to ensure that the measurement result of s_i is compared to the result the most probable to have been obtained from the same quantity value.

In practice, Δt_{φ} can be related to the characteristic time of variation of the measurand whereas Δl_{φ} may be related to the spatial homogeneity of this quantity.

Another definition of the spatial closeness can be designed based on the concept of representative area of measurement results [126], noted $a(s_i,t)$ for s_i at t. It is the area around the location of a measurement result in which the measurement result is representative of any other result that would have been obtained at another position within the area. This definition may be more usable in practice than the one based on a norm as it considers the geometry of the

²⁴If no instrument is mobile, it may be necessary to have less restrictive requirements on the spatiotemporal vicinity to observe rendez-vous.

area in which the instruments are. There are buildings and crossings for instance in an urban environment. Thus, with this definition, s_i is in rendez-vous with s_i at t if for instance:

$$\exists t' \in \mathbb{R}^+$$
 such as $\exists m(s_j, t')$ and $|t - t'| \leq \Delta t_{\varphi}$ and $a(s_i, t) \cap a(s_j, t') \neq \emptyset$

An alternate expression could be the following:

$$\exists t' \in \mathbb{R}^+$$
 such as $\exists m(s_i, t')$ and $|t - t'| \leq \Delta t_{\varphi}$ and $|a(s_i, t) \cap a(s_i, t')| \geq \Delta a_{\varphi}$

where Δa_{φ} is the minimal size of the intersection of the representativity area of both measurement results.

In conclusion, the concept of rendez-vous can be applied according to different definitions in practice. In the following sections, the diagnosis algorithm is designed without any assumption on the practical definition of a rendez-vous.

3.3.2 Compatible rendez-vous

Consider a rendez-vous $\varphi(s_i \to s_j, t)$. By associating the definitions of compatible measurement results, a rendez-vous of s_i with s_j at t is stated as **compatible** if $m(s_i, t) \approx m(s_j, t')$. Otherwise it is stated as **upper non-compatible** if $m(s_i, t) > m(s_j, t')$ or **lower non-compatible** if $m(s_i, t) < m(s_j, t')$. Consequently, $\Phi^{\approx}(s_i \to s_j, (t, \Delta t))$, $\Phi^+(s_i \to s_j, (t, \Delta t))$ and $\Phi^-(s_i \to s_j, (t, \Delta t))$ are the respective associated sets of such rendez-vous.

4 Algorithm for the diagnosis of calibration issues in a sensor network

4.1 General idea

The general purpose of a diagnosis algorithm is to determine whether if a system is **faulty** (F) or **non-faulty** (NF). In the present case, the predicted status of an instrument s_i at t is noted $\widehat{\Omega}(s_i,t)$. Ideally, $\widehat{\Omega}(s_i,t)$ is equal to the true status of the instrument, $\Omega(s_i,t)$, which is unknown in practice.

To determine if instruments are correctly calibrated, the proposed approach consists in using the concept of rendez-vous introduced in Section 3.3. When two instruments s_i and s_j are in rendez-vous, it means that, according to chosen spatiotemporal conditions, they were both measuring the same quantity value at the same time and the same place. Suppose that one of the instruments, s_j for instance, was diagnosed as non-faulty and consider $m(s_i,t)$ and $m(s_j,t')$, the measurement results involved in a rendez-vous $\varphi(s_i \to s_j,t)$. If both results are metrologically valid, it means that $m(s_j,t')$ can be seen as a standard value for s_i : it is a known value with an associated uncertainty. Thus, if $m(s_i,t) \approx m(s_j,t')$, it denotes that s_i is correctly calibrated, at least according to s_j , and its status can be predicted as non-faulty. Otherwise, it is predicted as faulty. As s_j helps to predict the status of s_i , it is called a **diagnoser** of s_i .

Predicting the status of an instrument based on only one value may easily lead to false predictions because drift is a fault usually having a longer characteristic time than others like spike faults. For instance, if spikes are not correctly removed before carrying out a diagnosis, it is possible that $m(s_i,t) > m(s_j,t')$ or that $m(s_i,t) < m(s_j,t')$, e.g. the measurement results $m(s_i,t)$ and $m(s_j,t')$ are not compatible, whereas the true status of s_i , $\Omega(s_i,t)$ is actually equal to non-faulty from a calibration perspective. Hence, the use of multiple rendez-vous with s_j and also with other instruments to predict the status of s_i can reduce the impact of these particular

²⁵We recall that more restrictive criterion can be set to determine if measurement results are compatible with the results of other instruments but it is not studied in this work. See Section 3.2 for more details.

²⁶We recall that the measurement results $m(s_i, t)$ and $m(s_j, t')$ have two different timestamps t and t' respectively because they may not have been obtained exactly at the same time (see Section 3.3).

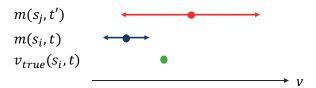


Figure 4.4.1: Example where the measurement results $m(s_i, t)$ and $m(s_j, t')$ are compatible with each other but where $m(s_i, t)$ is not compatible with its true value

cases. Moreover, if the class²⁷ of s_i is such as $c(s_i) \gg c(s_j)$, predicting the status of s_i based on s_j , even if $\widehat{\Omega}(s_j,t') = NF$, may not be advisable. Indeed, it is possible that $m(s_i,t) \approx m(s_j,t')$ but with $m(s_i,t)$ not compatible with its true value as shown in Figure 4.4.1. A solution to that is to allow only instruments of a higher class than the one of s_i to take part in the prediction of its status. This minimal class is noted $c_{min}^{\mathcal{D}}(s_i)$ and is defined by:

$$c_{min}^{\mathcal{D}}(s_i) = c(s_i) + \Delta c_{min}^{\mathcal{D}}(c(s_i))$$

where $\Delta c_{min}^{\mathcal{D}}(k)$ is the relative difference of class required between instruments of class k and the instruments allowed to be their diagnosers.²⁸

Thus, to determine the status of an instrument s_i at t_d , the following set of **valid rendez-vous** $\Phi_v(s_i \to S, (t_d, \Delta t))$ is used:

$$\Phi_{v}(s_{i} \to S, (t_{d}, \Delta t)) = \{ \varphi(s_{i} \to s_{j}, t) \in \Phi(s_{i} \to S, (t_{d}, \Delta t)), \text{ such as}$$

$$s_{j} \notin S^{\mathcal{D}}, // s_{j} \text{ is not an instrument to diagnose}$$

$$\widehat{\Omega}(s_{j}, t') = NF, // s_{j} \text{ is non-faulty}$$

$$c(s_{j}) \geq c_{min}^{\mathcal{D}}(s_{i}), // \text{ The class of } s_{j} \text{ is higher or equal to the minimal}$$

$$class \text{ allowed to diagnose } s_{i}$$

$$m(s_{i}, t) \in M^{*}(s_{i}, (t_{d}, \Delta t)), // \text{ The measurement result of } s_{i} \text{ is valid}$$

$$m(s_{j}, t') \in M^{*}(s_{j}, (t_{d}, \Delta t)) \} // \text{ The measurement result of } s_{j} \text{ is valid}$$

$$(4.1)$$

It is also possible that the input information of a diagnosis algorithm is not sufficient to choose between F and NF statuses. Thus, a third option for the predicted status of a system is **ambiguous** (A).²⁹ In the present case, it is when $\Phi_v(s_i \to S, (t_d, \Delta t))$ does not contain enough rendez-vous to allow a prediction with enough confidence. This minimal size for a set of valid rendez-vous for any instrument is noted $|\Phi_v|_{min}$.³⁰

Based on the definition of $\Phi_v(s_i \to S, (t_d, \Delta t))$, an algorithm can be designed to determine the statuses of all the instruments in a sensor network.

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$$|\Phi_v|_{min} = \min_{k \in [0..c_{max}]} |\Phi_v|_{min}(k)$$

 $^{^{27}\}mathrm{See}$ Definition 20 in Chapter 1 Section 3, page 18.

 $^{^{28}\}Delta c_{min}^{\mathcal{D}}(k)$ can be identical for each class e.g. $\Delta c_{min}^{\mathcal{D}}(k) = \Delta c_{min}^{\mathcal{D}}(S) \forall k \in [0..c_{max}].$

²⁹The true status of an instrument cannot be ambiguous. It is either faulty or non-faulty strictly.

 $^{^{30}}$ It is possible to define different $|\Phi_v|_{min}$ for each class of instrument like for $\Delta c_{min}^{\mathcal{D}}(k)$. $|\Phi_v|_{min}$ was introduced as a common constant for the diagnosis of all the instruments because the amount of valid information to perform a binary diagnosis has no major reason to change from a class to another and in the worst case, $|\Phi_v|_{min}$ can be set as:

4.2 Procedure for the diagnosis of all the instruments in a sensor network

Consider a diagnosis procedure d that occurs at t_d , and Δt such as $[t_d - \Delta t; t_d]$ is the time range on which the diagnosis procedure is carried out. The measurement results obtained and the rendez-vous that occurred during $[t_d - \Delta t; t_d]$ are the data that will be used to predict the statuses of the instruments of S. This forms a scene as defined by [40].

All the instruments in $S^{c_{max}}$ are assumed as non-faulty. ³¹ Thus, at the beginning of a diagnosis procedure, the set of instruments to diagnose $S^{\mathcal{D}}$ is equal to $S \setminus S^{c_{max}}$ where $S^{c_{max}}$ is the set of instruments of class c_{max} .

Then, the predicted statuses $\widehat{\Omega}(s_i, t_d)$ for each $s_i \in S$ are initialised. The predicted statuses of the instruments in $S^{c_{max}}$ are set to non-faulty and those of the instruments in $S^{\mathcal{D}}$ are initially set to ambiguous.

Afterwards the predicted statuses of all the instruments, noted $\widehat{\Omega}(S, t_d)$, are actualised. These actualised statuses are noted $\widetilde{\Omega}(S, t_d)$.

They are determined as follows for each instrument $s_i \in S^{\mathcal{D}}$. First, if $|\Phi(s_i \to S, (t_d, \Delta t))| < |\Phi_v|_{min}$, it means that during $[t_d - \Delta t; t_d]$, s_i did not meet other instruments enough times to be able to diagnose its status, whatever the predicted statuses of the other instruments. Therefore, it is not possible to actualise its predicted status with another value than ambiguous, which is already the value of $\widetilde{\Omega}(s_i, t_d)$, and s_i is removed from $S^{\mathcal{D}}$.

Otherwise, $\Phi_v(s_i \to S, (t_d, \Delta t))$ is determined. If its size is lower than $|\Phi_v|_{min}$, the actualised predicted status of s_i stays equal to ambiguous and s_i remains in $S^{\mathcal{D}}$ as the size of $\Phi_v(s_i \to S, (t_d, \Delta t))$ may change in a future iteration.

If $|\Phi_v(s_i \to S, (t_d, \Delta t))| \ge |\Phi_v|_{min}$, the actualised predicted status can be determined between non-faulty and faulty.

To do so, the rates $r_{\Phi_v}^{\approx}(s_i \to S, (t_d, \Delta t)), r_{\Phi_v}^+(s_i \to S, (t_d, \Delta t))$ and $r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t))$ are computed. They are defined as follows:

 $r_{\Phi_v}^{\approx}(s_i,(t_d,\Delta t))$ is the rate of compatible rendez-vous in the set of valid rendez-vous of s_i over the time range $[t_d - \Delta t; t_d]$. It is equal to

$$r_{\Phi_v}^{\approx}(s_i, (t_d, \Delta t)) = \frac{|\Phi_v^{\approx}(s_i, (t_d, \Delta t))|}{|\Phi_v(s_i, (t_d, \Delta t))|}$$

 $r_{\Phi_v}^+(s_i,(t_d,\Delta t))$ is the rate of upper non-compatible rendez-vous in the set of valid rendez-vous of s_i over the time range $[t_d - \Delta t; t_d]$. It is equal to

$$r_{\Phi_v}^+(s_i, (t_d, \Delta t)) = \frac{|\Phi_v^+(s_i, (t_d, \Delta t))|}{|\Phi_v(s_i, (t_d, \Delta t))|}$$

 $r_{\Phi_v}^-(s_i,(t_d,\Delta t))$ is the rate of lower non-compatible rendez-vous in the set of valid rendez-vous of s_i over the time range $[t_d - \Delta t; t_d]$. It is equal to

$$r_{\Phi_v}^{-}(s_i, (t_d, \Delta t)) = \frac{|\Phi_v^{-}(s_i, (t_d, \Delta t))|}{|\Phi_v(s_i, (t_d, \Delta t))|}$$

Based on $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$, which are the maximal tolerated values associated to the rates $r_{\Phi_v}^{\approx}(s_i, (t_d, \Delta t))$, $r_{\Phi_v}^+(s_i, (t_d, \Delta t))$ and $r_{\Phi_v}^-(s_i, (t_d, \Delta t))$, if one of the following conditions is true:

•
$$r_{\Phi_v}^+(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^+)_{max}$$

³¹This assumption is discussed in Section 6.

• $r_{\Phi_n}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_n}^-)_{max}$

• $(1 - r_{\Phi_v}^{\approx}(s_i \to S, (t_d, \Delta t))) > (r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$

then $\widetilde{\Omega}(S, t_d)$ is equal to faulty. Otherwise $\widetilde{\Omega}(s_i, t_d)$ is set to non-faulty. In the end, s_i is removed from $S^{\mathcal{D}, 32}$

The use of three different rates, $r_{\Phi_v}^{\approx}(s_i \to S, (t_d, \Delta t))$, $r_{\Phi_v}^+(s_i \to S, (t_d, \Delta t))$ and $r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t))$, is useful because it brings an additional information on the status of the instrument if it is diagnosed as faulty. If this prediction was triggered by the condition on $r_{\Phi_v}^+(s_i \to S, (t_d, \Delta t))$, it indicates that the instrument is overestimating its measured values for instance.

After all the instruments in $S^{\mathcal{D}}$ are treated, if $\widehat{\Omega}(S, t_d) = \widetilde{\Omega}(S, t_d)$, it means that no statuses of the instruments in $S^{\mathcal{D}}$ changed. Consequently, the statuses of the instruments of the network at t_d , $\widehat{\Omega}(S, t_d)$, are determined and the diagnosis procedure ends. Otherwise $\widehat{\Omega}(S, t_d)$ takes the values of $\widehat{\Omega}(S, t_d)$ and the actualised statuses are determined again for each instrument $s_i \in S^{\mathcal{D}}$.

The pseudo-code of this algorithm is provided in Algorithm 1.

4.3 Improvements and extensions of the presented algorithm

The considered algorithm can be easily improved in terms of efficiency. Appendix A Section 1 provides a version with several changes that may reduce the number of iterations of the main loop of Algorithm 1.

Moreover, it is possible to add more confidence into the predicted statuses of the instruments by restricting the set of the diagnosers of an instrument s_i to the one of the diagnosers of class $k > c_{min}^{\mathcal{D}}(s_i)$, with k being the highest class such as the predicted status $\widehat{\Omega}(s_i, t_d)$ can be determined as non-faulty. This formulation is provided in Appendix A Section 2.

In addition, the formulation of Algorithm 1 invites to compute it in a centralised manner. Appendix A Section 3 introduces elements allowing an application of the principles of the diagnosis algorithm in a decentralised manner.

Finally, Appendix A Section 4 gives insights on how the diagnosis procedure could be extended to the case of a sensor network measuring different quantities and for which some of its measurands are influence quantities for several instruments in the network.

For the sake of clarity, Algorithm 1 is the version used in the following sections of this chapter.

4.4 Conclusion

In this section, the principle of a diagnosis algorithm for sensor networks aiming at detecting drift faults of measuring instruments was introduced. It is strongly based on the concepts of rendez-vous between measuring instruments and of class of a measuring instrument. As a major assumption, the instruments with the highest class must be assumed as non-faulty. Several parameters drive the algorithm and are studied in the following sections. In a first step, the next section gives an example of case study applying the presented algorithm.

 $^{^{32}}$ Alternate manners to determine the predicted status of an instrument according to compatible and non-compatible valid rendez-vous could be chosen. For instance, the value of the predicted status could be chosen according to the number of compatible and non-compatible valid rendez-vous. Thus $(r_{\Phi_v}^+)_{max}, \ (r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$ would be replaced by maximal numbers of tolerated non-compatible rendez-vous regarding $|\Phi_v^+(s_i \to S, (t_d, \Delta t))|, \ |\Phi_v^-(s_i \to S, (t_d, \Delta t))|$ and their sum.

Algorithm 1: Algorithm of the diagnosis procedure proposed for the detection of drift in sensor networks

```
Data: S, \Phi(S \to S, (t_d, \Delta t)), |\Phi_v|_{min}, \Delta c_{min}^{\mathcal{D}}
Result: \widehat{\Omega}(S, t_d)
/* Initiate the set of instruments to diagnose and the predicted statuses
      of the instruments
S^{\mathcal{D}} \leftarrow S \setminus S^{c_{max}}
\widehat{\Omega}(S^{c_{max}}, t_d) \leftarrow \text{NF}
\widehat{\Omega}(S^{\mathcal{D}}, t_d) \leftarrow \mathbf{A}
\widetilde{\Omega}(S^{\mathcal{D}}, t_d) \leftarrow \widehat{\Omega}(S^{\mathcal{D}}, t_d)
                                                                          /* Initiate the actualised statuses */
/* Ignore the instruments that cannot have enough valid rendez-vous
for s_i \in S^{\mathcal{D}} do
   if |\Phi(s_i \to S, (t_d, \Delta t))| < |\Phi_v|_{min} then
     S^{\mathcal{D}} \leftarrow S^{\mathcal{D}} \setminus \{s_i\}
end
/* Predict the status of each instrument to diagnose
                                                                                                                                                      */
repeat
      \widehat{\Omega}(S, t_d) \leftarrow \widehat{\Omega}(S, t_d)
                                                    /* The actualised statuses are now the predicted
       statuses */
      for s_i \in S^{\mathcal{D}} do
           /* Build the current set of valid rendez-vous for s_i
                                                                                                                                                      */
            c_{min}^{\mathcal{D}}(s_i) \leftarrow c(s_i) + \Delta c_{min}^{\mathcal{D}}(c(s_i))
            \Phi_v(s_i \to S, (t_d, \Delta t)) \leftarrow \{\varphi(s_i \to s_j, t) \in \Phi(s_i \to S, (t_d, \Delta t)), \text{ such as } \}
                  s_j \notin S^{\mathcal{D}}, \ \widehat{\Omega}(s_j, t') = NF, \ c(s_j) \ge c_{min}^{\mathcal{D}}(s_i),
                  m(s_i, t) \in M^*(s_i, (t_d, \Delta t)) \text{ and } m(s_i, t') \in M^*(s_i, (t_d, \Delta t))
           /* If s_i have enough valid rendez-vous, then compute the different
                  rates to actualize its status
                                                                                                                                                      */
           if |\Phi_v(s_i \to S, (t_d, \Delta t))| \ge |\Phi_v|_{min} then
                r_{\Phi_v}^{\infty}(s_i \to S, (t_d, \Delta t)) \succeq \frac{|\Psi_v|_{min} \text{ then}}{|\Phi_v^{\infty}(s_i \to S, (t_d, \Delta t))|}
r_{\Phi_v}^{\infty}(s_i \to S, (t_d, \Delta t)) \leftarrow \frac{|\Phi_v^{\infty}(s_i \to S, (t_d, \Delta t))|}{|\Phi_v(s_i \to S, (t_d, \Delta t))|}
r_{\Phi_v}^{+}(s_i \to S, (t_d, \Delta t)) \leftarrow \frac{|\Phi_v^{+}(s_i \to S, (t_d, \Delta t))|}{|\Phi_v(s_i \to S, (t_d, \Delta t))|}
r_{\Phi_v}^{-}(s_i \to S, (t_d, \Delta t)) \leftarrow \frac{|\Phi_v^{-}(s_i \to S, (t_d, \Delta t))|}{|\Phi_v(s_i \to S, (t_d, \Delta t))|}
                 /* If a condition on the different rates is met, then the
                        actualised status of s_i is set to faulty, otherwise, it is set
                        to non-faulty
                 if r_{\Phi_v}^+(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^+)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or }
                   (1-r_{\Phi_n}^{\approx}(s_i \to S, (t_d, \Delta t))) > (r_{\Phi_n}^+ + r_{\Phi_n}^-)_{max} then
                    \Omega(s_i, t_d) \leftarrow F
                  else
                   \Omega(s_i, t_d) \leftarrow \text{NF}
                  S^{\mathcal{D}} \leftarrow S^{\mathcal{D}} \setminus \{s_i\} /* s_i is diagnosed so it can be removed from S^{\mathcal{D}} */
            \mathbf{end}
      end
until \Omega(S,t_d)=\Omega(S,t_d) /* Repeat until there is no difference between the
  predicted and actualised statuses */
```

5 Application of the algorithm to a first case study

To illustrate the presentation of the diagnosis algorithm, this section provides a first case study based on simulation to appreciate its performances.

5.1 Definition of the case study

5.1.1 Sensor network

A sensor network of 10 instruments is considered. The class of one of them is equal to 1, the others being of class zero. Thus, $c_{max} = 1$ in this case.

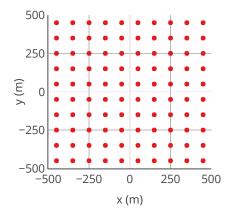


Figure 4.5.1: Map of the 100 positions available for the instruments in the case study.

The instruments of class k=0 move randomly in a discrete space of 100 positions at a discrete time. The positions of the instruments are geometrically defined according to a grid of 10×10 with a step that represents 100m, and centred on (x,y)=(0,0) as shown in Figure 4.5.1. At each time step a new position is chosen randomly for each instrument following a uniform law. Instruments may remain in position and multiple instruments may share a position. Two instruments are in rendez-vous when they are at the same position at the same time. In addition, the instrument of class c_{max} is static. Its position is randomly drawn among the 100 positions.

As the case study is based on simulation, the time step has no actual physical meaning. To ease the comprehension of the study, the time step of the simulation represents 10 min and the case study lasts 265 days.

5.1.2 Instruments

Instruments are assumed to be initially calibrated. In the first place, the instrument of class c_{max} is assumed as perfect, e.g. it does not drift. The instruments of class zero all follow the same drift model. The model chosen corresponds to the RGOI model defined in Chapter 5 Section 5.3.1, page 71. As a reminder, gain $G(s_i,t)$ and offset $O(s_i,t)$ drift of instrument s_i are computed for this model at each time step following:

$$G(s_i, t) = \begin{cases} 1 & \text{if } t < t_{\text{start drift}} \\ G(s_i, t - 1) + \delta G(s_i, t) & \text{if } t \ge t_{\text{start drift}} \end{cases}$$
with $\forall t, \delta G(s_i, t)$ drawn following $\mathcal{U}(0, \delta G_{max})$

$$O(s_i, t) = \begin{cases} 0 & \text{if } t < t_{\text{start drift}} \\ O(s_i, t - 1) + \delta O(s_i, t) & \text{if } t \ge t_{\text{start drift}} \end{cases}$$

with $\forall t, \delta O(s_i, t)$ drawn following $\mathcal{U}(0, \delta O_{max})$

with δG_{max} and δO_{max} being respectively the maximal gain and offset possible increase per time step.

Measured values are expressed following:

$$v(s_i, t) = G(s_i, t) \cdot v_{true}(s_i, t) + O(s_i, t)$$

The instruments start to drift at t = 0

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Each measured value $v(s_i, t)$ is associated to a constant relative uncertainty $\Delta_r v(c_i)$ depending on the class of the instrument such as

$$\frac{\Delta v(s_i, t)}{v(s_i, t)} = \Delta_r v(c_i)$$

A detection $\lim^{33} v_{min}(c_i)$ is also defined to balance the effect of high uncertainties for low measured values. This detection limit is used to determine if a measurement result $m(s_i, t)$ is metrologically valid, e.g. all the values below this detection limit are not considered valid.

All the values of the parameters are listed in Table 4.5.1.

Parameter		Value	Unit
Class 1			
	$\Delta_r v$	1	%
	v_{min}	0.752	$ m \mu gm^{-3}$
Class 0			
	δG_{max}	2/(24*6*30)	$\%/10 \mathrm{min}$
	δO_{max}	18.8/(24*6*30)	$\mu g/m^3/10 min$
	$\Delta_r v$	30	%
	v_{min}	37.6	$\mu \mathrm{g}\mathrm{m}^{-3}$
True values	model		
	A_{max}	200	$ m \mu gm^{-3}$
	σ_{max}	3826	m

Table 4.5.1: Values of the parameters of the case study. The values of δG_{max} and δO_{max} are displayed as a fraction of (24*6*30) min for the sake of clarity, e.g. the numerator is the maximal drift of the gain and of the offset per 30 days.

5.1.3 True values

To model the true values of the instruments, the model used in Chapter 3 Section 3.2.1 is considered. However in this case, instead of using the data that was considered to determine the values of A(t) and $\sigma(t)$, A(t) and $\sigma(t)$ are drawn randomly at each time step following the

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³³The detection limit is the value "obtained by a given measurement procedure, for which the probability of falsely claiming the absence of a component is β , given a probability α of falsely claiming its presence" [14]. In practice, it is recommended that $\beta = \alpha = 0.05$ for information.

uniform laws $\mathcal{U}(0, A_{max})$ and $\mathcal{U}(0, \sigma_{max})$ respectively. The values of A_{max} and σ_{max} are reported in Table 4.5.1.

The impact of this choice on the results is discussed in Section 5.6 and 9.3.

5.2 Configuration of the diagnosis algorithm

We aim at performing a diagnosis **every 15 days**. Thus, for $d \in D$, $t_d = (d+1) \times 15$ days, with D = [0..16].

The goal of this diagnosis is to detect when an instrument has provided at least 25% of measurement results non-compatible with their true values over the past 15 days. According to this specification, we choose $\Delta t = 15$ days and the rate thresholds $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$ are all set to 25%.

The minimal number of valid rendez-vous to conclude with a predicted status different from ambiguous $|\Phi_v|_{min}$ is set initially to 15.

5.3 Definition of the true status of an instrument

As it is a simulation, it is possible to know the true status of an instrument. Indeed its measured values and the true values are accessible, and therefore, it can be determined if 25% of measurement results non-compatible with their true values were obtained or not during a period of 15 days.

 $r_{true}(s_i, (t, \Delta t))$ is the rate of measurement results compatible with true values of s_i over $[t - \Delta t; t]$. It is equal to

$$r_{true}(s_i, (t, \Delta t)) = \frac{|M^{\approx}(s_i, (t, \Delta t))|}{|M(s_i, (t, \Delta t))|}$$

In this case, if $r_{true}(s_i, (t, \Delta t)) < 0.75$, then $\Omega(s_i, t) = F$. Otherwise, $\Omega(s_i, t) = NF$.

Note that the true status of an instrument is based on all its measurement results against the true values on the considered time range whereas its predicted status is based on the measurement results of its valid rendez-vous with other devices. Thus, it is expected that instruments which are actually non-faulty may be predicted as faulty and *vice versa* because the sets of measurement results used to compute the true and predicted status of an instrument are different.

5.4 Metrics for the evaluation of performances of the diagnosis algorithm

To estimate the performance of the algorithm, vocabulary and metrics that are used to evaluate binary classifiers [130, 153] are appropriate but must be adapted as the predicted statuses can take three values instead of two.

Regarding the possible true statuses and predicted statuses, there are six cases:

- A non-faulty status predicted for an instrument which true status is non-faulty is a true negative (TN)
- A faulty status predicted for an instrument which true status is non-faulty is a **false negative** (**FN**)
- An ambiguous status predicted for an instrument which true status is non-faulty is a non-determined negative (NDN)
- A non-faulty status predicted for an instrument which true status is faulty is a **false** positive (FP)
- A faulty status predicted for an instrument which true status is faulty is a **true positive** (**TP**)

3265 3266 • An ambiguous status predicted for an instrument which true status is faulty is a **non-determined positive** (NDP)

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To these primary metrics, P and N are added, which are respectively **the number of positives** (faulty true status) and **the number of negatives** (non-faulty true status).

The relationships between these metrics are summed up in Table 4.5.2

True status

		Non-faulty (N)	Faulty (P)
	Non-faulty	True negative (TN)	False negative (FN)
Predicted status	Ambiguous	Non-determined negative (NDN)	Non-determined positive (NDP)
	Faulty	False positive (FP)	True positive (TP)

Table 4.5.2: Contingency table of the different primary metrics

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For these metrics, only the instruments in $S \setminus S^{c_{max}}$ are considered. The instruments of class c_{max} being always assumed as non-faulty and being not drifting in the following sections, taking them into account would bias the metrics defined afterwards in Table 4.5.3.

Table 4.5.3: Metrics derived from the metrics P, N, TP, TN, FP, FN, NDP and NDN

Name	Expression	Role
Prevalence	$Prev = \frac{P}{P+N}$	Indicates the proportion of positive cases among all the cases
True positive rate	$TPR = \frac{TP}{P}$	Indicates the proportion of positive cases correctly detected
True negative rate	$TNR = \frac{TN}{N}$	Indicates the proportion of negative cases correctly detected
False positive rate	$FPR = \frac{FP}{N}$	Indicates the proportion of positive cases incorrectly detected
False negative rate	$FNR = \frac{FN}{P}$	Indicates the proportion of negative cases incorrectly detected
Non-determined positive rate	$NDPR = \frac{NDP}{P}$	Indicates the proportion of positive cases detected as ambiguous
Non-determined negative rate	$NDNR = \frac{NDN}{N}$	Indicates the proportion of negative cases detected as ambiguous
Non-determined rate	$NDR = \frac{NDP + NDN}{P + N}$	Indicates the proportion of all cases detected as ambiguous
Positive predictive value	$PPV = \frac{TP}{TP + FP}$	Indicates the proportion of correctly detected positive cases per positive call
False discovery rate	$FDR = \frac{FP}{TP + FP}$	Indicates the proportion of incorrectly detected positive cases per positive call
Negative predictive value	$NPV = \frac{TN}{TN + FN}$	Indicates the proportion of correctly detected negative cases per negative call
False omission rate	$FOR = \frac{FN}{TN + FN}$	Indicates the proportion of incorrectly detected negative cases per negative call

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Name	Expression	Role
Accuracy	$ACC = \frac{TP + TN}{P + N}$	Indicates the proportion of correct detection

Another metric added is the **delay of first positive detection for an instrument** s_i , noted $\Delta \mathcal{D}(s_i)$. It is the difference between the index of the diagnosis procedure d where $\Omega(s_i, t_d)$ changes from NF to F and the index of the diagnosis procedure d' where $\widehat{\Omega}(s_i, t_{d'}) = F$ for the first time. Thus:

$$\Delta \mathcal{D}(s_i) = d - d'$$

This value can be positive or negative as there can be early or late positive detection. In the following results, the average and the standard deviation of this metric over $S \setminus S^{c_{max}}$ are considered.

5.5 Results

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5.5.1 Observations from the results of different instruments

In Figure 4.5.2, the evolution of the true and predicted statuses over time are represented for different instruments. It shows that the diagnosis algorithm can produce different results over the instruments:

- the status of the instrument at each diagnosis procedure is often correctly predicted, except for a delay right after it becomes faulty (Figures 4.5.2a and 4.5.2c)
- there are predicted statuses equal to ambiguous, once only in Figure 4.5.2a, but it can happen frequently considering Figure 4.5.2b
- Changing decisions happen, e.g. an instrument predicted as faulty may be predicted as non-faulty later (Figure 4.5.2b)

In all these particular cases, false results are observed, e.g. for instance an instrument is predicted as non-faulty when it is faulty and so on. In Section 5.3, we stated that these behaviours were expected. We investigate the reasons behind these incorrect predictions in Section 5.6.

5.5.2 Overall appreciation

Globally, the values of the metrics computed over all the diagnosis procedures are listed in Table 4.5.4.³⁴ In this study, the prevalence of positive cases, e.g. the proportion of cases where the instruments are actually faulty over all the diagnosis procedures, is equal to 76%, for 153 cases (9 instruments diagnosed 17 times each).

From the number of TP, FP, TN, FN, NDP and NDN, it shows that:

- Most of the predicted statuses are true negatives (36 cases) and true positives (78 cases)
- Few predicted statuses are false negatives (20 cases) and non-determined positives (19 cases)
- No non-determined negative and false positive are predicted

Consequently, the TNR is equal to 1 and the FPR is equal to zero whereas the TPR is important (0.67) and the FNR is moderate (0.17). Moreover, the NDNR is equal to zero and

 $^{^{34}}$ We remind that the metrics do not take into account the instruments of class c_{max} in this Section.

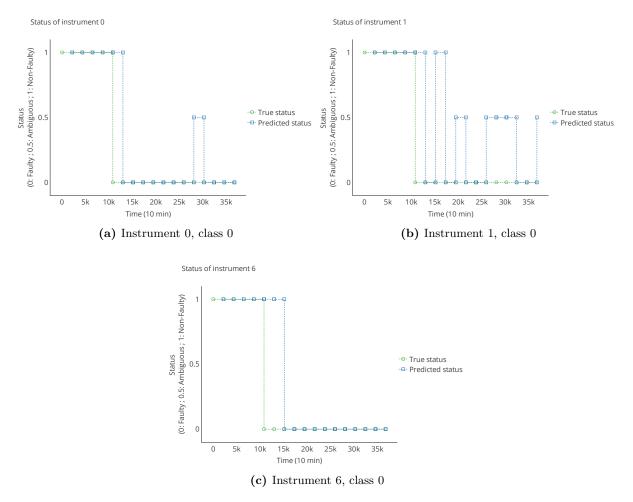


Figure 4.5.2: Evolution of the true status and of the predicted status for several instruments

the NDPR is equal to 0.16, thus NDR = 16% of ambiguous predictions were made which is quite important.

That being said, it appears that the algorithm is correct when it predicts that instruments are faulty (PPV=1, FDR=0). It is often not correct when it predicts that instruments are non-faulty, about one third of the time (NPV=0.64, FOR=0.36). According to Figure 4.5.2, there is a delay when an instrument actually becomes faulty and the first prediction as faulty. It confirmed by the statistics of the delay of positive detection $\Delta \mathcal{D}$ in Table 4.5.5: the average value is negative and in fact it is always lower or equal to zero according to the maximal value of this metric.

Nevertheless, the accuracy is equal to 75% which is acceptable for a first case study without a fine adjustment of the diagnosis algorithm's parameters.

5.5.3 Evolution over time of the metrics

The values of the metrics given in the previous subsection are representative of the results for all the diagnosis 17 procedures. They do not allow to understand how this overall result was build. Therefore, we study here their evolution.

Figure 4.5.3 shows the evolution of the metrics as a function of the index of the diagnosis procedure. Each curve gives the values obtained for the corresponding metric at the diagnosis d with the results of the nine instruments, where d is the index of the diagnosis procedure.

		True status			
		Non-faulty $N = 36$	Faulty $P = 117$	Prevalence 0.76	Accuracy 0.75
	Non-faulty	TN	FN	\overline{NPV}	FOR
	Ivon-lauity	36	20	0.64	0.36
Dradiated status	A malaimu au a	NDN	NDP		
Predicted status	Ambiguous	0	19		
	Faulty	FP	TP	FDR	PPV
		0	78	0.00	1.00
		TNR	\overline{FNR}		
		1.00	0.17		
		\overline{NDNR}	NDPR	$\overline{}$ NDR	-
		0.00	0.16	0.16	
		FPR	TPR		
		0.00	0.67		

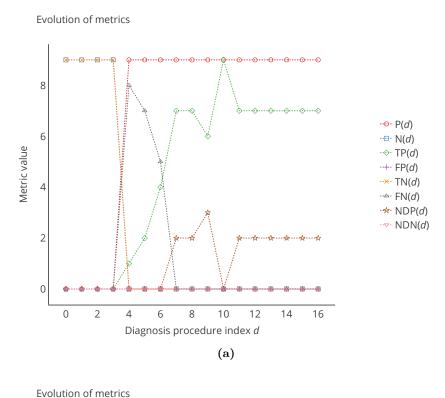
Table 4.5.4: Confusion matrix of the case study (metrics are defined in Section 5.4).

Delay of positive detection						
μ	σ	min	max			
-2.1	1.2	$\overline{-4}$	0			

Table 4.5.5: Statistics of the delay of positive detection $\Delta \mathcal{D}$ for the case study

Figure 4.5.3a shows that first there are only non-faulty instruments (curve of N), and faulty ones appear at the 5th diagnosis (d=4). After the 6th diagnosis, all the instruments of class zero are faulty (curve of P). The transition is quite abrupt but this was expected because all the instruments of class zero follow the same law of drift. The curves of the true positives TP and true negatives TN follow mostly the ones of P and N respectively. The false positives or negatives are in fact occurring around the transition where instruments of class zero become faulty. This behaviour is more clearly represented in Figures 4.5.3b and 4.5.3c with the curves of the TPR and FNR, and the curves of the NPV and FOR respectively. Finally, with Figure 4.5.3d, we observe that the global accuracy, initially equals to one, decreases until the transition phase and then increases again, converging to 0.8, which is explained by the 20% of NDP cases in the last diagnosis procedures (Figure 4.5.3c).

To conclude, this particular study of the evolution of the metrics allowed to understand the global results over the 17 diagnosis procedures: the false results occur mainly when the instruments actually become faulty and the number of predicted statuses equal to ambiguous increases as more as there are instruments actually predicted as faulty. Thus, worse overall results could have been displayed by considering the results of few diagnosis procedures around the transition phase. We conjecture that better results would be obtained by increasing the number of diagnosis procedures in the case study, without changing any parameter of the algorithm, e.g.



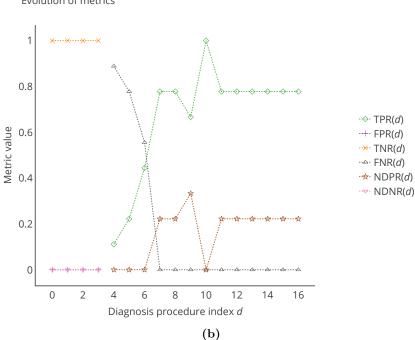
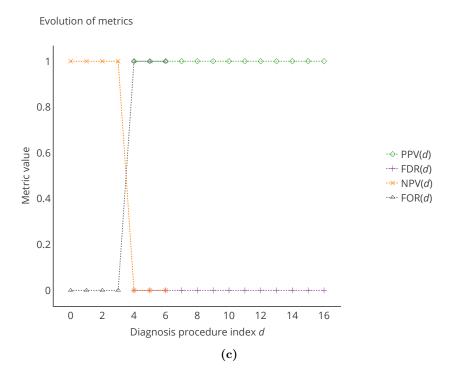


Figure 4.5.3: Evolution of the metrics computed for each diagnosis procedure as a function of the current diagnosis procedure id (continued) (metrics are defined in Section 5.4).

by increasing the duration of the case study. However, this would not reduce the number of false that may happen in absolute terms. The reasons why false results happen are explained in the following section.

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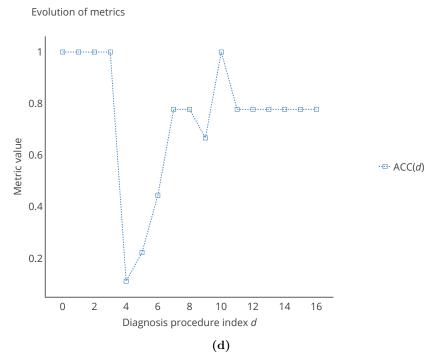
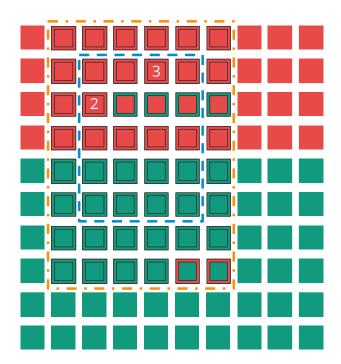


Figure 4.5.3: Evolution of the metrics computed for each diagnosis procedure as a function of the current diagnosis procedure id (metrics are defined in Section 5.4).



Symbol	$m(s_i, t)$ compatible with true value?	$m(s_i,t)$ involved in a rendez-vous?	$m(s_i, t) \approx m(s_j, t')$?
	Yes	No	-
	No	No	-
	Yes	Yes	Yes
	Yes	Yes	No
	No	Yes	Yes
	No	Yes	No
	The number i indicates the in i rendez-vous. Set of the measurements of the measurement $M(s_i, (t_d, \Delta t)) _{\Phi_n}$	results involved in rendez-	vous $M(s_i,(t_d,\Delta t)) _{\Phi}$

Figure 4.5.4: Representation of an example set of measurement results $M(s_i, (t_d, \Delta t))$. First, we observe that the restrictions of the set of measurement results $M(s_i, (t_d, \Delta t))$ to the ones where they are respectively involved in rendez-vous $M(s_i, (t_d, \Delta t))|_{\Phi}$ and involved in valid rendez-vous $M(s_i, (t_d, \Delta t))|_{\Phi}$ do not have the same proportion of measurement results compatible with their true values. Thus, the rates of measurement results compatible with true values $r_{true}(s_i, (t_d, \Delta t))$, $r_{true}(s_i, (t_d, \Delta t))|_{\Phi}$ and $r_{true}(s_i, (t_d, \Delta t))|_{\Phi_v}$ are not equal. Therefore, considering that the predicted statuses are determined with $r_{\Phi_v}^{\approx}(s_i, (t_d, \Delta t))$ for instance, which computation is based on the set $\Phi_v(s_i \to S, (t_d, \Delta t))$, it is possible that $r_{\Phi_v}^{\approx}(s_i, (t_d, \Delta t))$ from $r_{true}(s_i, (t_d, \Delta t))$. Here $r_{true}(s_i, (t_d, \Delta t)) = 0.6$ and $r_{\Phi_v}^{\approx}(s_i, (t_d, \Delta t)) = 0.48$. If a threshold of 0.5 is considered to differentiate faulty instruments from non-faulty ones, then the true status of s_i is $\Omega(s_i, t_d) = NF$ here but its predicted status is $\widetilde{\Omega}(s_i, t_d) = F$: it is a false result.

5.6 Explanations of false results

In this case study, multiple false results, e.g. false positive or false negative, were observed. The reasons why they happen can easily be explained.

Consider as an example a set of measurement results $M(s_i, (t_d, \Delta t))$. This set is composed of 100 measurement results, represented as squares in Figure 4.5.4, such as the rate of compatible measurement results with their true values $r_{true}(s_i, (t_d, \Delta t))$ is equal to $\frac{60}{100} = 0.6$. If a threshold equal to 0.5 for r_{true} is considered to differentiate faulty instruments from non-faulty ones³⁵, then $\Omega(s_i, t_d) = NF$ here.

In the diagnosis algorithm, the predicted status of an instrument is determined based on the measurement results associated to the rendez-vous in the set $\Phi_v(s_i \to S, (t_d, \Delta t))$. These measurement results are the ones in the blue rectangle dashed line in Figure 4.5.4. The question is whether these measurement results are representative of $M(s_i, (t_d, \Delta t))$ and thus allow obtaining a $r_{\Phi_v}^{\approx}(s_i, (t_d, \Delta t))$ equal to $r_{true}(s_i, (t_d, \Delta t))$ or not.

Consider first $M(s_i, (t_d, \Delta t))|_{\Phi}$ which is the restriction of $M(s_i, (t_d, \Delta t))$ to the measurement results associated to rendez-vous that are in the set $\Phi(s_i \to S, (t_d, \Delta t))$. In this case, not all the measurement results of s_i are involved in a rendez-vous, only those with a border in Figure 4.5.4 (grouped in the orange rectangle dashed line in Figure 4.5.4). Computing $r_{true}(s_i, (t_d, \Delta t))|_{\Phi}$, which is defined similarly to $r_{true}(s_i, (t_d, \Delta t))$, gives a value of $\frac{24}{48} = 0.50$ which is worse than $r_{true}(s_i, (t_d, \Delta t))$ but the result is still correct: it allows to say that s_i is non-faulty because $r_{true}(s_i, (t_d, \Delta t))|_{\Phi} \ge 0.5$ according to the threshold previously chosen.

However, the rate used for the prediction of the status of an instrument, $r_{\Phi_v}^{\approx}(s_i, (t_d, \Delta t))$, is not based on comparisons with true values but with other instruments. Different situations can be observed as shown in the table of Figure 4.5.4:

- a measurement result compatible with its true values can be compatible or not with the measurement result of the other instrument involved in a rendez-vous
- a measurement result non-compatible with its true values can be compatible or not with the measurement result of the other instrument involved in a rendez-vous
- measurement results of an instrument may be involved in several rendez-vous.

From the perspective of $\Phi(s_i \to S, (t_d, \Delta t))$ first before moving to $\Phi_v(s_i \to S, (t_d, \Delta t))$, $r_{true}(s_i, (t_d, \Delta t))|_{\Phi}$ may not be equal to $r_{\Phi}^{\approx}(s_i, (t_d, \Delta t))$ and it is the case here as $r_{\Phi}^{\approx}(s_i, (t_d, \Delta t)) = \frac{24-2+4}{48+3} = 0.51 \neq r_{true}(s_i, (t_d, \Delta t))|_{\Phi}$. The difference is small but the reason why false results happen begins to appear. On top of the fact that the compatibility between measurement results of two instruments may not correctly report the true status of an instrument r_{Φ}^{36} , the sets used for the computations of the rates r_{true} and r_{Φ} do not necessarily have the same composition of measurement results.

Considering then $M(s_i, (t_d, \Delta t))|_{\Phi_n}$, we have:

$$r_{true}(s_i, (t_d, \Delta t))|_{\Phi_v} = \frac{8}{20} = 0.40 \text{ and } r_{\Phi_v}^{\approx}(s_i, (t_d, \Delta t)) = \frac{8+3}{20+3} = 0.48$$

Thus, with a threshold $(r_{\Phi}^+ + r_{\Phi}^-)_{max} = 0.5$, the predicted status of s_i , $\widehat{\Omega}(s_i, t_d)$, is equal to faulty (because $r_{\Phi_v}^{\approx}(s_i, (t_d, \Delta t)) < (r_{\Phi}^+ + r_{\Phi}^-)_{max}$) which is different from $\Omega(s_i, t_d)$: it is a false result.

Figure 4.5.5 shows the evolution of the different rates r_{true} and $r_{\Phi_v}^{\approx}$ computed for each block of 15 days for the instrument that was considered in Figure 4.5.2b. With the trigger level in red, the evolution of the green and blue curves in Figure 4.5.2b can be explained following the

³⁵e.g. if $r_{true}(s_i, (t_d, \Delta t)) \geq 0.5$ then $\Omega(s_i, t_d) = NF$

³⁶See Section 4.1

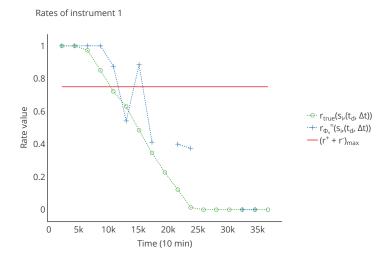


Figure 4.5.5: Evolution of the rates r_{true} and $r_{\Phi_v}^{\approx}$ for s_1 in the case study of Section 5. The missing points for $r_{\Phi_v}^{\approx}$ are due to the ambiguous cases where $r_{\Phi_v}^{\approx}$ could not be computed.

reasoning developed for the presented theoretical example. In this Figure, the diagnosis curve is particularly interesting as the instrument is alternatively diagnosed as faulty and non-faulty, and several times.

Through this development, the reason why false results appear has been explained. It is because the properties of sets of measurement results like $M(s_i, (t_d, \Delta t))$ are not necessarily conserved while building $\Phi_v(s_i \to S, (t_d, \Delta t))$ and it is this latter set that is used to compute the rates r_{Φ_v} . In fact, there is a sort of "sampling" of $M(s_i, (t_d, \Delta t))$ that is performed through the rendez-vous and resulting in $M(s_i, (t_d, \Delta t))|_{\Phi_v}$ in the end. This explains mainly the false results of the diagnosis algorithm. Solving this issue is quite challenging. It might be possible in some situations where the characteristics of the sets of measurement results could be conserved by the "sampling" at stake for the restriction of the sets of measurement results but it would generally require strong assumptions on the measurement results and on the occurrence of rendez-vous, without taking into account the notion of validity of the rendez-vous. However, by adjusting the parameters of the diagnosis algorithm, it may also be possible to reduce the number of false results. Therefore, they are reviewed in the following section.

5.7 On the parameters of the diagnosis algorithm

There are several parameters that drive the proposed algorithm:

- the instant of diagnosis t_d and its periodicity
- the time difference Δt defining the length of the time range on which measurement results selected for the input data of the algorithm
- the minimal relative differences of class required for instruments of class k with their diagnoser instruments $\Delta c_{min}^{\mathcal{D}}(k)$
 - the minimal required number of valid rendez-vous $|\Phi_v|_{min}$
- the conditions for a measurement result to be considered as valid
 - the conditions of spatiotemporal vicinity for two instruments to be in rendez-vous

• the maximal tolerated values $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$ for the rates of upper and lower non-compatible rendez-vous in sets of valid rendez-vous.

It may be possible to improve the results of case study with different rules or values for them. Indeed, values were chosen but may not be the ones allowing obtaining the best results. They may explain the reason why there are many false results in particular. The following paragraphs provide a discussion regarding the choice of the parameters.

Instant of diagnosis and its periodicity

 t_d is the instant of the $d^{\rm th}$ diagnosis. In this case study, we chose to have a periodic diagnosis, e.g. diagnosis procedures are carried out on a regular basis (every 15 days).³⁷ With a different frequency of diagnosis, for instance every week or even every day in the case study, it may be possible to detect sooner that an instrument has become faulty in exchange of a higher computational cost. Consequently, the overall results would be different but this would not change the general behaviour of the algorithm. The choice of the instant of diagnosis and its periodicity is not discussed in detail in this work despite the fact that it may have a significance in practice because it strongly depends on the instruments used and on the context of deployment. Indeed, all the instruments do not have the same characteristic time of drift for instance. This information is often provided as an order of magnitude of drift of the gain and offset over a week or a month in percentage in datasheets [104]. In this way, it would influence the periodicity of the diagnosis. Moreover, the context of deployment can justify to perform the diagnosis at specific times of the day.³⁸ Thus, the characteristics of the instruments used and the context of deployment can help in the choice of the moment of diagnosis and its periodicity.³⁹

Length of the time range on which the algorithm is applied

The length of the time range Δt has an important influence on the results. Indeed, the higher Δt is, the more rendez-vous may be in $\Phi(s_i \to S, (t_d, \Delta t))$ and consequently in $\Phi_v(s_i \to S, (t_d, \Delta t))$, and vice versa. Thus, it influences the contents of $M(s_i, (t_d, \Delta t))|_{\Phi}$ and $M(s_i, (t_d, \Delta t))|_{\Phi_v}$. However, a high value of Δt implies that measurement results obtained at very different instants would be used by the diagnosis algorithm. Considering that old measurement results are more accurate than recent ones in the case of an irreversible drift, the higher Δt is, the less the measurement results at the actual level of the drift will be preponderant in $M(s_i, (t_d, \Delta t))|_{\Phi_v}$. Therefore, choosing a different value for Δt while trying to bring closer the distributions of measured values in $M(s_i, (t_d, \Delta t))$ and $M(s_i, (t_d, \Delta t))|_{\Phi_v}$ could also induce false results for another reason because as shown in Section 5.6, the content of the sets $M(s_i, (t_d, \Delta t))$ and $M(s_i, (t_d, \Delta t))|_{\Phi_v}$ are not necessarily equivalent, regardless of Δt . Moreover, reducing Δt in particular could also have another drawback: on a short time range there may likely be less rendez-vous than on a longer one. Depending on the minimal required number of valid rendez-vous $|\Phi_v|_{min}$, the value of Δt can make the algorithm not functional, e.g. all the instruments

 $^{^{37}}$ An actual day and hour is not associated to t_d because it does not have a real signification here.

³⁸As an example, in [109] for an *in situ* calibration strategy, Moltchanov *et al.* chose to apply their *in situ* calibration algorithm every day after 4:00a.m. and to use the measured values between 1:00a.m. and 4:00a.m.. They proceeded like this because the measuring instruments to calibrate were distant from the reference and at this moment, the measurand was supposed to be homogenous in the measurement area, enabling to compare between them the values measured by the instruments. Therefore, in the same context with our diagnosis algorithm, it may be relevant to apply it at a similar time of the day that was used in [109].

³⁹Note that too high a frequency of diagnosis may be pointless if no rendez-vous occurs between two instants of diagnosis. It would be more relevant to have in this case a formulation of the algorithm so that the diagnosis procedure is carried out when new rendez-vous happen. In this way, Appendix A Section 5 provides a discussion regarding first how to reformulate the algorithm so that it is possible to trigger a diagnosis procedure on an event, e.g. a rendez-vous, and how to use an event-based range on which the algorithm is applied. Then, elements regarding a real-time version of the diagnosis algorithm are given in Appendix A Section 6.

 different from the class c_{max} would be predicted as ambiguous.

In fact, as for the instant of diagnosis and its periodicity, the adjustment of Δt could be defined based on the properties of the instruments used like their characteristic time of drift. Indeed, it can give an order of magnitude of the relevant length of the time range on which the diagnosis algorithm should be applied.

Minimal relative differences of class required

The minimal relative differences of class required $\Delta c_{min}^{\mathcal{D}}(k)$ would usually be set to +1 or 0. Indeed, it is rarely advisable to determine the statuses of instruments with ones of a lower class as shown in Section 4.1. Predicting the status of an instrument with diagnoser instruments of a strictly higher class can give a higher confidence in the diagnosis as their measurement results are standards of a higher quality than the ones provided by diagnoser instruments of an identical class.

Minimal required number of valid rendez-vous

The minimal required number of valid rendez-vous $|\Phi_v|_{min}$ can be a chosen integer but to add up more confidence in the predicted status, it is advisable to make it the highest as possible. However, too high a value can make the diagnosis impossible: all the instruments of class $c_{max}-1$ and lower would be diagnosed as ambiguous. If a training dataset of rendez-vous is available, it is possible to give an upper boundary. This is explained in Section 8.

Conditions for a measurement result to be considered as valid

The conditions to satisfy to state a measurement result of an instrument as valid can be set according to the datasheets of the instruments. Measurement range, detection limit and so on are values usually given. Thus, it is not really a parameter driving the diagnosis algorithm, it is a constraint driven by the choice of the measuring instruments.

Conditions of spatiotemporal vicinity for two instruments to be in rendezvous

This subject was already discussed in Section 3.3. We introduced different definitions for these conditions. In summary, they should be determined based on expert specifications or on characteristics of the measurand and of the measuring instruments.

Maximal tolerated values for the rates of upper and lower non-compatible rendez-vous in sets of valid rendez-vous

Like t_d , its periodicity and Δt , the maximal tolerated values $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ r_{\Phi_v}^-)_{max}$ for the rates of upper and lower non-compatible rendez-vous in sets of valid rendez-vous can be derived according to a requirement. This is what was done in Section 5.2. Nevertheless, these parameters have a major role in the way false results are produced because they are used to make the final decision between a status predicted as faulty or non-faulty. Therefore, their adjustment is discussed in Section 7 where the reduction of false results is tackled.

5.8 Conclusion

In this section, a case study was conducted to put in practice the algorithm of diagnosis presented in Section 4.2. It provides satisfying results according to the initial specification. However, the results are not excellent. There are false results and, in this particular case, a significant number of false negatives. There are also several cases of predicted statuses equal to ambiguous. The reasons behind the false results were explained theoretically.

To improve the results, the adjustment of the parameters of the algorithm was discussed. The brief overview of each one highlighted that the adjustment of some of them can be driven by expert specifications and the characteristics of the measuring instruments composing the sensor network. For $|\Phi_v|_{min}$, $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$, guidelines are provided in Sections 7 and 8 to set their values.

Before that, we discuss in the following section the assumption for the instruments of class c_{max} that are considered as non-faulty-they were even not drifting in the present case study.

6 On the assumption regarding the top-class instruments being always predicted as non-faulty

6.1 Theoretical discussion

The assumption that instruments of class c_{max} are always predicted as non-faulty is a strong one. It is, however, necessary for the proper operation of the presented algorithm. If there is no assumption regarding the status of any instrument of class c_{max} , they are initially considered as ambiguous and thus the sets $\Phi_v(s_i \to S, (t_d, \Delta t))$ are always empty. In this case, the algorithm predicts a status equal to ambiguous for each instrument, which makes it useless.

Actually, the assumption on the instruments of class c_{max} can be relaxed: **one single** instrument of class c_{max} supposed as non-faulty may be sufficient. More practically, depending on the value of $|\Phi_v|_{min}$ and on the configuration of the sensor network, only a subset of $S^{c_{max}}$ may be sufficient to ensure the lowest number of instruments systematically diagnosed as ambiguous. Note that this formulation is also more appropriate in the case of a blind sensor network. Indeed, for this type of sensor network, all the instruments are considered as equivalent from a metrological perspective. Consequently, there is no subset of instruments of a higher class than others.

Following on from this new formulation of the assumption regarding the statuses of the instruments of class c_{max} , we might be tempted to try to relax it to point where the status of only one instrument of any class has to be known. Consider such an instrument s_i . Again, all the instruments $s_j \in S$ such as $c_{min}^{\mathcal{D}}(s_j) > c(s_i)$ cannot be diagnosed differently than ambiguous if the minimal relative differences of class required $\Delta c_{min}^{\mathcal{D}}(k)$ is set to +1 or 0 as recommended in Section 5.7. In particular, if the class of s_i is the lowest class in the network, noted c_{min} , and if $\Delta c_{min}^{\mathcal{D}}(c_{min}) = +1$, then all the instruments of the network, except s_i are systematically diagnosed as ambiguous. In fact, the lower the class of s_i is, the more instruments will always be predicted as ambiguous. Thus, it may be simpler to come back to assuming some instruments of class c_{max} as non-faulty. It is acceptable to guarantee such a hypothesis in real situations.

6.2 Case study with the instrument of class c_{max} drifting

In the case study of Section 5, the instrument of class c_{max} , noted here s_9 , was first assumed as perfect, e.g. not drifting, for the sake of clarity.

Consider here that this instrument is also undergoing drift following the RGOI model but with $\delta G_{max}(s_9) = 0.5/(24*6*30)\%/10$ min and $\delta O_{max}(s_9) = 1.88/(24*6*30)$ µg/m³/10min.⁴² With the previous definition of r_{true} and the same criteria to determine the true status of the instrument $\Omega(s_9,t)$ like the other instruments, the evolution of the true and predicted statuses

⁴⁰Note that if $\Delta c_{min}^{\mathcal{D}}(c_{max}) > 0$, the instruments of class c_{max} which statuses are not assumed as non-faulty, cannot be diagnosed differently than ambiguous. Hence, it may be more relevant to define a sub-network without these instruments on which the diagnosis algorithm is applied so that the results are not biased by the behaviour of the algorithm with these instruments.

⁴¹In the case of a blind sensor network, necessarily $\Delta c_{min}^{\mathcal{D}} = 0$ for all the instruments, otherwise the algorithm cannot work.

⁴²See Table 4.5.1 for the explanation regarding $G_{max}(s_9)$ and $O_{max}(s_9)$ expressed as fractions.

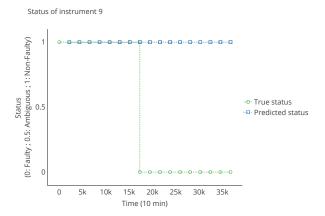


Figure 4.6.1: Evolution of the true status and of the predicted status of s_9 , the instrument of class 1, when it is drifting

		True s	True status		
		Non-faulty $N = 36$	Faulty $P = 117$	Prevalence 0.76	Accuracy 0.72
	Non-faulty	TN	FN	\overline{NPV}	FOR
	11011-1autty	36	28	0.56	0.44
Predicted status	Ambiguous	NDN	\overline{NDP}		
Fredicted status	Ambiguous	0	15		
	Faulty	FP	TP	FDR	PPV
		0	74	0.00	1.00
		TNR	\overline{FNR}		
		1.00	0.24		
		\overline{NDNR}	NDPR	\overline{NDR}	-
		0.00	0.13	0.10	
		\overline{FPR}	TPR		
		0.00	0.63		

Table 4.6.1: Confusion matrix of the case study when the instrument of class 1 is drifting (metrics are defined in Section 5.4).

over time for this particular instrument is given in Figure 4.6.1, the predicted status being again assumed as always non-faulty. We observe that there are consequently false negative if the assumption is kept. However, still without taking into account s_9 in the metrics to evaluate the results, the diagnosis algorithm gives the results of Tables 4.6.1 and 4.6.2. We observe that the values of the metrics do not change significantly from Tables 4.5.4 and 4.5.5. They are slightly worse sometimes: more false negatives (28 instead of 20), less true positives (74 instead of 78), thus NPV is lower and the FOR is higher but the PPV is higher.

Therefore, it appears that even with s_9 drifting, the diagnosis algorithm still provides satisfying results. In this way, assuming that the instruments of class c_{max} are not drifting is acceptable.

Delay of positive detection							
μ	σ	min	max				
-2.3	0.9	$\overline{-4}$	-1				

Table 4.6.2: Statistics of the delay of positive detection $\Delta \mathcal{D}$ for the case study when the instrument of class 1 is drifting

More detailed explanations are provided in Appendix B to show that even when drifting, this instrument is still more accurate than the instruments of a lower class. In fact, the metric used to determine if an instrument is truly faulty or not is not fairly appreciating the status of an instrument when the goal is to compare its measurement results to the ones of another instrument, potentially of a worse quality.

6.3 Conclusion

 In this section, the need for an assumption regarding the instruments of class c_{max} to be considered as non-faulty was explained and relaxed to only require this assumption for one of them at least. The case study of Section 5 was extended by making the instrument of class c_{max} drifting. It showed that the results were only slightly worse in this configuration. In conclusion, assuming that instruments of class c_{max} are always non-faulty is a reasonable hypothesis.

7 Means to reduce false results

In this section, we discuss several means to reduce the number of false results that happen for the reasons explained in Section 5.6.

7.1 Keep the predicted status of instruments unchanged once they are predicted as faulty

A first idea to overcome faults results could be to keep the predicted status of an instrument equal to faulty for the next diagnosis procedures once it is first diagnosed as so. Figure 4.7.1 shows the effect of such choice on the same instrument as in Figure 4.5.2b: here the instrument changes of status only once.

The global performances of the diagnosis algorithm on the case study with this choice of behaviour are shown in Table 4.7.1. The results obtained are improved: there are less false negatives, less non-determined positives and much more true positives. Thus, all the associated metrics based on these metrics are improved. An accuracy of 88% is obtained, while the PPV goes up to 1 for instance. Only the FNR and FOR are still important but this is explained by the late detection of the change of status of the instruments in average (-2.1) as shown in Table 4.7.2.

Therefore, this solution of keeping the predicted status of an instrument equal faulty, once it is predicted as so, avoids observing changes of predicted statuses after being predicted as faulty a first time but it does not solve entirely the problem of false results.

7.2 Alternate definition for rates used to decide of the status of the instruments

In Section 5.6, we observed that measurement results may be considered multiple times in the computation of the rates $r_{\Phi_v}^+$, $r_{\Phi_v}^-$ and $r_{\Phi_v}^{\approx}$ if an instrument encounters multiple other instruments at the same time. It could be preferable to have each measurement result counted only once,

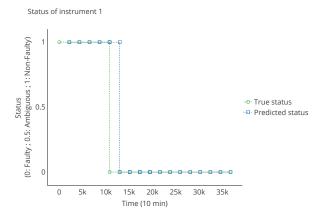


Figure 4.7.1: Evolution of the true status and of the predicted status of s_1 while keeping $\widehat{\Omega}(s_1,t)=F$ for $t\geq t_d$ such as $\widehat{\Omega}(s_1,t_d)=F$ for the first time.

		True status			
		Non-faulty $N = 36$	Faulty $P = 117$	Prevalence 0.76	Accuracy 0.88
Predicted status	Non-faulty	TN	FN	\overline{NPV}	FOR
		36	18	0.67	0.33
	Ambiguous	NDN	NDP		
		0	1		
	Faulty	FP	TP	\overline{FDR}	PPV
		0	98	0.00	1.00
		TNR	FNR		
		1.00	0.15		
		NDNR	NDPR	NDR	
		0.00	0.01	0.01	_
		FPR	TPR		
		0.00	0.84		

Table 4.7.1: Confusion matrix of the case study with the predicted statuses kept as faulty from one diagnosis procedure to another once the status of an instrument is predicted as faulty for the first time (metrics are defined in Section 5.4).

notably if $M(s_i, (t_d, \Delta t))|_{\Phi_v} \sim M(s_i, (t_d, \Delta t))$ in terms of distribution of the measured values. To achieve that, we define $\gamma(s_i \to s_j, t)$, the weight of the rendez-vous $\varphi(s_i \to s_j, t)$:

$$\gamma(\varphi(s_i \to s_j, t)) = \frac{1}{|\Phi(s_i \to S, t)|}$$

with $\Phi(s_i \to S, t) = \Phi(s_i \to S, (t, 0))$ regarding the previous notations of the sets Φ

Delay of positive detection						
μ	σ	min	max			
-2.1	1.2	-4	0			

Table 4.7.2: Statistics of the delay of positive detection $\Delta \mathcal{D}$ for the case study with the predicted statuses kept as faulty from one diagnosis procedure to another once the status of an instrument is predicted as faulty for the first time

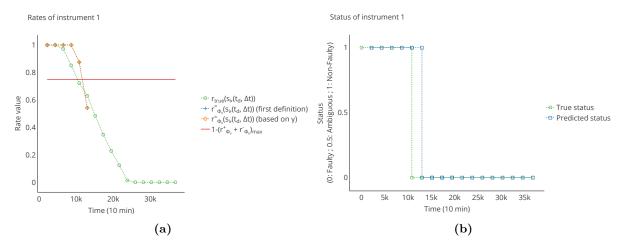


Figure 4.7.2: Evolution for s_1 in the case study of Section 5 (a) of the rates r_{true} and $r_{\Phi_v}^{\approx}$ (both definitions); (b) of its true status and of its predicted status with the alternate definition for $r_{\Phi_v}^{\approx}$; while keeping $\widehat{\Omega}(s_1,t) = F$ for $t \geq t_d$ such as $\widehat{\Omega}(s_1,t_d) = F$ for the first time.

 γ_v can be defined for Φ_v too. With this weight, we have notably:

$$\sum_{\varphi(s_i \to s_j, t') \in \Phi_v(s_i, (t, \Delta t))} \gamma_v(\varphi(s_i \to s_j, t')) = |M(s_i, (t, \Delta t))|_{\Phi_v}|$$

Thus, an alternate definition of $r_{\Phi_v}^{\approx}(s_i,(t,\Delta t))$ is for instance:

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$$r_{\Phi_v}^{\approx}(s_i, (t, \Delta t)) = \frac{\sum_{\varphi(s_i \to s_j, t') \in \Phi_v^{\approx}(s_i, (t, \Delta t))} \gamma_v(\varphi(s_i \to s_j, t'))}{\sum_{\varphi(s_i \to s_j, t') \in \Phi_v(s_i, (t, \Delta t))} \gamma_v(\varphi(s_i \to s_j, t'))}$$

However, Figure 4.7.2a shows that using this definition in this case study does not change the values of $r_{\Phi_v}^{\approx}(s_1, (t, \Delta t))$ compared to the results obtained with the previous definition. Hence, there is no change in the predicted status of the instrument that was considered in Figure 4.7.1 as shown in Figure 4.7.2b. This is also globally the case according to Table 4.7.3 compared to Table 4.7.1.

7.3 Adjustment of the maximal tolerated values for the different rates used to determine the statuses of the instruments

The issue of false results can be seen from another perspective. In fact, the most problematic false results are the false negatives. Indeed, false positives will only trigger an early maintenance

		True status			
		Non-faulty $N = 36$	Faulty $P = 117$	Prevalence 0.76	Accuracy 0.88
Predicted status	Non faulty	TN	FN	\overline{NPV}	FOR
	Non-faulty	36	18	0.67	0.33
	Ambiguous	NDN	NDP		
		0	1		
	Faulty	FP	TP	\overline{FDR}	PPV
		0	98	0.00	1.00
		TNR	\overline{FNR}		
		1.00	0.15		
		\overline{NDNR}	\overline{NDPR}	$\overline{}$ NDR	-
		0.00	0.01	0.01	
		FPR	TPR		-
		0.00	0.84		

Table 4.7.3: Confusion matrix of the case study with the predicted statuses kept as faulty from one diagnosis procedure to another once the status of an instrument is predicted as faulty for the first time and with the alternate definition for $r_{\Phi_n}^{\approx}$ (metrics are defined in Section 5.4).

of the instruments which is preferable in terms of quality of the measurements. The cost of the maintenance has to be taken into account but regarding calibration issues, an *in situ* calibration could be performed.

The parameters that can be adjusted to obtain hopefully more false negatives and less false positives are the maximal tolerated values $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$ for the rates of upper and lower non-compatible rendez-vous in sets of valid rendez-vous. The lower these values are, the less non-compatible result will be accepted.

It is difficult to give a formal procedure to adjust these values and would require additional developments. By the way of a possible solution, an empirical approach could be considered if a training database or test cases are available. While setting first values for $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$ arbitrarily, instruments may be diagnosed and the results studied to see if the predictions were correct. If it is the case, it means that the maximal tolerated values for the rates could be increased if there were only true positives, and decreased if there were only true negatives. In the opposite cases, e.g. if there were only false positives or only false negatives, it means that these values should be increased and decreased respectively. However, if there are both true positives, true negatives, false positives and false negatives, the values for $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$ should be adjusted so that the true positives and true negatives are maximised and the false positives and false negatives minimised. It may be impossible to find values so that there are no false results.

7.4 Conclusion

In this section, two means to reduce false results were presented, one having an interesting effect (keeping the predicted status of instruments unchanged once they are predicted as faulty)

and the other one having no influence on the results (an alternate definition for rates used to decide of the status of the instruments). Then, we discussed the adjustment of the maximal tolerated values for the different rates used to determine the statuses of the instruments because it can help to have more false positives and less false negatives. We explained notably why false positives were less predjudiciable than false negatives.

In conclusion, we can say that despite a careful design of the sensor network and adjustment of the parameters of the diagnosis algorithm, false results may still happen. Therefore, a careful investigation must be carried out if a high rate of false results is observed to understand their causes.

8 Adjustment of the minimal size required for a set of valid rendez-vous to allow a prediction between the statuses faulty and non-faulty

The choice of the value of $|\Phi_v|_{min}$ can make the algorithm non-functional: in the worst situation, all the instruments, except those of class c_{max} , are always diagnosed as ambiguous if the condition $|\Phi_v(s_i, (t, \Delta t))| \ge |\Phi_v|_{min}$ is never satisfied for any s_i to diagnose. Thus, having a method to determine first if the value of $|\Phi_v|_{min}$ is compatible with the considered sensor network and then to give an upper boundary for its value would be interesting.

Note that in the case where the condition $|\Phi_v(s_i, (t, \Delta t))| \ge |\Phi_v|_{min}$ is never satisfied, it is also possible to increase Δt so that more rendez-vous could be included in the set $\Phi_v(s_i, (t, \Delta t))$. We assume here that Δt was defined first to determine a relevant set of rendez-vous based on expert knowledge (Section 5.7) and cannot be changed.

8.1 Algorithm determining an upper boundary for $|\Phi_v|_{min}$

Consider a training database giving the possible rendez-vous between the instruments of a sensor network over a given time range, such as a set D of diagnosis procedures could be carried out on it.

For each $d \in D$, it is possible to determine for each instrument the set of its rendez-vous $|\Phi(s_i \to s_j, (t_d, \Delta t))|$.

Consider the matrix $\mathbf{Mat}_{\Phi}(D, \Delta t) \in \mathbb{N}^{|S| \times |S|}$ where its values are defined by:

$$\phi_{i,j} = \min_{d \in D} |\Phi(s_i \to s_j, (t_d, \Delta t))|$$

Each line of this matrix gives the worst-case scenario of rendez-vous occurring on a range of diagnosis according to the training database.

We will now apply a procedure similar to the diagnosis algorithm to obtain a maximum value for $|\Phi_v|_{min}$ such as all the instruments are diagnosable as non-faulty, noted max $|\Phi_v|_{min}$, under a number of assumptions.

Consider first the set $S^{NF} = S^{c_{max}}$ because the instruments of class c_{max} are assumed as non-faulty by the diagnosis algorithm.⁴³ and the set of instruments to diagnose $S^{\mathcal{D}} = S \setminus S^{c_{max}}$.

For each $s_i \in S^{\mathcal{D}}$, it is possible to compute:

$$|\Phi_v|_{min}(s_i) = \sum_{\substack{j=1\\c(s_j) \ge c_{min}^{\mathcal{D}}(s_i)\\s_i \in S^{NF}}}^{|S|} \phi_{i,j}$$

The maximum of these $|\Phi_v|_{min}(s_i)$ gives a first value for the upper boundary max $|\Phi_v|_{min}$. If

⁴³In the case where only a subset of $S^{c_{max}}$ is assumed as non-faulty (see Section 6), only the instruments of this subset should be in S^{NF} .

we suppose that all the rendez-vous are likely to be valid in the worst case scenario and that all the instruments are likely to be non-faulty, then with $|\Phi_v|_{min} = \max |\Phi_v|_{min}$, all the s_i such as $|\Phi_v|_{min}(s_i) = \max |\Phi_v|_{min}$ would be predicted as non-faulty. Representing this set of instruments by \hat{S}^{NF} , we would have in a following iteration of the diagnosis algorithm $S^{NF} = S^{NF} \cup \hat{S}^{NF}$ and $S^{\mathcal{D}} = S^{\mathcal{D}} \setminus \hat{S}^{NF}$.

 $|\Phi_v|_{min}(s_i)$ can be computed again for the remaining instruments in $S^{\mathcal{D}}$. If the maximum of these new values is lower than the previous value obtained for $\max |\Phi_v|_{min}$, then $\max |\Phi_v|_{min}$ is set to this new maximum so that some of the remaining instruments in $S^{\mathcal{D}}$ can be diagnosed in the worst case scenario of rendez-vous during a second iteration of the algorithm. This procedure is repeated until $S^{\mathcal{D}}$ is empty. The value of $\max |\Phi_v|_{min}$ obtained at the end is the maximal value for $|\Phi_v|_{min}$ ensuring that all the instruments can be diagnosed as non-faulty when they are likely to be so in the worst case scenario of rendez-vous.

However, some instruments can be diagnosed as faulty in practice. In this case, the value $\max |\Phi_v|_{min}$ obtained previously may not be reachable for some instruments.

To overcome this issue, we can compute $\max |\Phi_v|_{min}$ as a function of the number λ of instruments predicted as faulty in the network.

Consider $C^F(S^{(c_{max}-1)-}, \lambda)$ that is the set of all the possible combinations of λ faulty instruments in $S^{(c_{max}-1)-}$. For each set of instruments that are faulty $S^F \in C^F(S^{(c_{max}-1)-}, \lambda)$, it is possible to determine $\max |\Phi_v|_{\min}(S^F)$ following the procedure described previously by initialising $S^{\mathcal{D}}$ to $S \setminus (S^{c_{max}} \cup S^F)$. For $\lambda > 0$, it is possible that some instruments have no rendez-vous with the non-faulty instruments depending on S^F . Thus, the end of the procedure described previously is reached when $S^{\mathcal{D}}$ is empty or when $S^{\mathcal{D}}$ is identical after an iteration.

Then:

$$\max |\Phi_v|_{min}(\lambda) = \min_{S^F \in \mathcal{C}^F(S^{(c_{max}-1)}-,\lambda)} (\max |\Phi_v|_{min}(S^F))$$

is the maximal value allowed for $|\Phi_v|_{min}$ considering λ instruments as faulty.

However, it is possible that $\max |\Phi_v|_{min}(\lambda) > \max |\Phi_v|_{min}(\lambda')$ for $\lambda' < \lambda$.

Thus:

$$\max |\Phi_v|_{min}(\lambda) = \min_{\lambda' \in [0..\lambda]} (\max |\Phi_v|_{min}(\lambda'))$$

Algorithm 2 gives a pseudo-code to determine the values of max $|\Phi_v|_{min}$ as a function of the number of instruments considered as faulty λ .

Algorithm 2: Algorithm to determine the maximal value of the minimal size required for a set of valid rendez-vous to allow prediction between the statuses faulty and non-faulty for an instrument, as a function of the number of faulty instruments in the sensor network

```
Data: S, \{\Phi(S,(t_d,\Delta t)), d \in D\}

Result: \max |\Phi_v|_{min}

/* Compute the matrix of the minimal numbers of rendez-vous encountered on the diagnosis ranges between each couple of instruments */

for i \in [1..|S|] do

| for j \in [1..|S|] do

| \phi_{i,j} \leftarrow \min_{d \in D} |\Phi(s_i \to s_j, (t_d, \Delta t))|
| end

end

/* Determine the maximal value allowed for |\Phi_v|_{min} as a function of the number of faulty instruments */
```

```
for \lambda \in [0..|S^{(c_{max}-1)-}|] do /* For each number of instruments that can be
 faulty
                                                                                                                            */
    \max |\Phi_v|_{min}(\lambda) \leftarrow +\infty
    /* For each combination of \lambda instruments that can be faulty
                                                                                                                             */
     for S^F \in \mathcal{C}^F(S^{(c_{max}-1)-},\lambda) do
          S^{NF} \leftarrow S^{c_{max}}
          S^{\mathcal{D}} \leftarrow S \setminus (S^{c_{max}} \cup S^F)
          \max |\Phi_v|_{min}(S^F) \leftarrow +\infty
         /* Determine the maximal value for |\Phi_v|_{min} according to S^F so that
               all the instruments in S^{\mathcal{D}} can be predicted as non-faulty
          repeat
               \widetilde{S}^{\mathcal{D}} \leftarrow S^{\mathcal{D}}
               \max |\Phi_v|_{min}(S^F)' \leftarrow 0
               for s_i \in S^{\mathcal{D}} do
                   \max |\Phi_v|_{min}(s_i) \leftarrow \sum_{\substack{j=1\\c(s_j) \geq c_{min}^{\mathcal{D}}(s_i)\\c^{NF}}}^{|S|} \phi_{i,j}
                    if \max |\Phi_v|_{min}(s_i) > \max |\Phi_v|_{min}(S^F)' then
                       \max |\Phi_v|_{min}(S^F)' \leftarrow \max |\Phi_v|_{min}(s_i)
                        \widehat{S}^{NF} \leftarrow \{s_i\}
                    else if \max |\Phi_v|_{min}(s_i) = \max |\Phi_v|_{min}(S^F)' then
                        \widehat{S}^{NF} \leftarrow \widehat{S}^{NF} \cup \{s_i\}
               \mathbf{end}
               if \max |\Phi_v|_{min}(S^F)' < \max |\Phi_v|_{min}(S^F) then
                   \max |\Phi_v|_{min}(S^F) \leftarrow \max |\Phi_v|_{min}(S^F)'
               end
               S^{NF} \leftarrow S^{NF} \cup \widehat{S}^{NF}
               S^{\mathcal{D}} \leftarrow S^{\mathcal{D}} \setminus \widehat{S}^{NF}
          until S^{\mathcal{D}} = \varnothing \ or \ S^{\mathcal{D}} = \widetilde{S}^{\mathcal{D}}
          /* If this new maximal value for |\Phi_v|_{min} is lower than the current
               one of \max |\Phi_v|_{min}(\lambda), then it is the new maximal value allowed
               for |\Phi_v|_{min} with \lambda faulty instruments
                                                                                                                             */
          if \max |\Phi_v|_{min}(S^F) < \max |\Phi_v|_{min}(\lambda) then
               \max |\Phi_v|_{min}(\lambda) \leftarrow \max |\Phi_v|_{min}(S^F)
          end
         /* If the maximal value allowed for |\Phi_v|_{min} with \lambda faulty
               instruments is higher than the maximal value allowed for a lower
               number of instruments, then it is this value that is the maximal
               value allowed for |\Phi_v|_{min} with \lambda faulty instruments
         if \lambda > 0 and \max |\Phi_v|_{min}(\lambda - 1) < \max |\Phi_v|_{min}(\lambda) then
              \max |\Phi_v|_{min}(\lambda) \leftarrow \max |\Phi_v|_{min}(\lambda - 1)
          end
     end
end
```

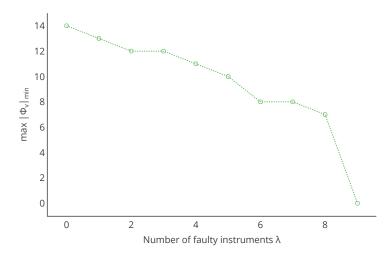


Figure 4.8.1: Evolution of $\max |\Phi_v|_{min}$ as a function of the number of faulty instruments λ in the network, based on the matrix $\mathbf{Mat}_{\Phi}(D, \Delta t)$ derived from the database used for the case study of Section 5

8.2 Application to the case study

Consider the following matrix $\mathbf{Mat}_{\Phi}(D, \Delta t)$ obtained with the database used in the case study of Section 5:

$$\mathbf{Mat}_{\Phi}(D,\Delta t) = \begin{pmatrix} 0 & 13 & 11 & 8 & 4 & 9 & 10 & 8 & 9 & 13 \\ 13 & 0 & 7 & 12 & 7 & 11 & 12 & 10 & 9 & 12 \\ 11 & 7 & 0 & 9 & 10 & 9 & 11 & 11 & 12 & 10 \\ 8 & 12 & 9 & 0 & 9 & 9 & 12 & 9 & 13 & 8 \\ 4 & 7 & 10 & 9 & 0 & 8 & 13 & 11 & 13 & 7 \\ 9 & 11 & 9 & 9 & 8 & 0 & 10 & 8 & 13 & 11 \\ 10 & 12 & 11 & 12 & 13 & 10 & 0 & 14 & 9 & 12 \\ 8 & 10 & 11 & 9 & 11 & 8 & 14 & 0 & 13 & 14 \\ 9 & 9 & 12 & 13 & 13 & 13 & 9 & 13 & 0 & 8 \\ 13 & 12 & 10 & 8 & 7 & 11 & 12 & 14 & 8 & 0 \end{pmatrix}$$

The last row corresponds to the instrument of class $c_{max} = 1$.

Applying the considered algorithm to determine $\max |\Phi_v|_{min}$ as a function of λ gives the Figure 4.8.1.

First of all, note that the overall upper boundary for $|\Phi_v|_{min}$ given by this algorithm is equal to 14, which is lower than the value used in the case study in Section 5: this value was set to 15. This could explain some of the predictions that were observed of instruments diagnosed as ambiguous. Hence, the algorithm we provided allow to identify if a value of $|\Phi_v|_{min}$ is likely to lead to instruments predicted as ambiguous.

Suppose that we want to be still able to diagnose the sensor network without any ambiguous predicted statuses due to a lack of valid rendez-vous when five instruments are predicted as faulty for instance. With Figure 4.8.1, then $|\Phi_v|_{min}$ must be lower or equal to 10 according to the training database used.

		True status			
		Non-faulty $N = 36$	Faulty $P = 117$	Prevalence 0.76	Accuracy 0.91
Predicted status	Non-faulty	TN	FN	\overline{NPV}	FOR
	Non-raurty	36	14	0.72	0.28
	Ambiguous	NDN	NDP		
		0	0		
	Faulty	FP	TP	\overline{FDR}	PPV
		0	103	0.00	1.00
		TNR	FNR		
		1.00	0.12		
		NDNR	NDPR	NDR	
		0.00	0.00	0.00	
		\overline{FPR}	TPR		-
		0.00	0.88		

Table 4.8.1: Confusion matrix of the case study with the predicted statuses kept as faulty from one diagnosis procedure to another once the status of an instrument is predicted as faulty for the first time and $|\Phi_v|_{min} = 10$ (metrics are defined in Section 5.4).

Delay of positive detection				
μ	σ	min	max	
-1.6	1.2	-3	0	

Table 4.8.2: Statistics of the delay of positive detection $\Delta \mathcal{D}$ for the case study with the predicted statuses kept as faulty from one diagnosis procedure to another once the status of an instrument is predicted as faulty for the first time, $|\Phi_v|_{min} = 10$

Consider now $|\Phi_v|_{min} = 10$ in our case study with the predicted status of the instrument kept unchanged once they are equal to faulty (Section 7). The execution of the diagnosis algorithm with this value for $|\Phi_v|_{min}$ gives the results of Tables 4.8.1 and 4.8.2. We observe that the results are significantly improved from what was obtained previously (Tables 4.7.1 and 4.7.2). There are less false negatives, no more non-determined positives, and more true positives. The PPV is now equal to 1 and the NPV increased from 0.67 to 0.72. The accuracy is now equal to 0.91 instead of 0.88. Moreover, the average delay of positive detection has been improved.

Therefore, changing the value of $|\Phi_v|_{min}$ is necessary if the algorithm is not functional at all (e.g. when $\widehat{\Omega}(s_i,t) = A$, $\forall s_i$ and t) but it can also be useful to improve the performances of the algorithm. In our case, a lower value of $|\Phi_v|_{min}$ is interesting.⁴⁴ The algorithm provided to determine max $|\Phi_v|_{min}$ as a function of the number of faulty instruments in the sensor network is adapted to help to find a suitable $|\Phi_v|_{min}$.

⁴⁴Note that a higher (or lower) value than 10 for $|\Phi_v|_{min}$ could give better results. The optimisation of this value was not addressed in this work.

8.3 Conclusion

In this section, an algorithm was provided to give an upper boundary to $|\Phi_v|_{min}$, the minimal size for a set of valid rendez-vous to allow a prediction between the statuses faulty and non-faulty. We defined this upper boundary $\max |\Phi_v|_{min}$ as a function of the number of instruments already diagnosed as faulty in the sensor network to adjust $|\Phi_v|_{min}$ with concerns for the robustness of the diagnosis algorithm when instruments are diagnosed as faulty.

Afterwards, this algorithm was applied on the dataset used in the case studies presented in this chapter. It allowed first to identify that the value of $|\Phi_v|_{min}$ initially chosen was higher than the overall upper boundary given by the algorithm. This could explain the predictions of instruments as ambiguous in previous results. Then, by choosing a lower value for $|\Phi_v|_{min}$ to tolerate five instruments predicted as faulty in previous diagnosis procedures, the results of the case study were improved, showing the interest of the algorithm developed in this section.

9 Sensitivity of the algorithm to changes in the case study

In the previous sections, the diagnosis algorithm was introduced and a case study was presented to illustrate the application of the algorithm. The means to reduce false results were discussed just like the assumption regarding the instruments of class c_{max} . In Chapter 3, the influence on the results of various aspects of the case study was studied for the evaluation of in situ calibration strategies. In this section, the same type of studies are conducted regarding the case study that was presented in Section 5. Indeed, results are likely to be case specific according to the results of the two previous sections where we saw that the choice of the values of the parameters $|\Phi_v|_{min}$, $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$, $(r_{\Phi_v}^+)_{max}$ and so on could change results obtained.

In this section, the influence of the drift values, of the true values and of the model used to build the true values is investigated. The role of the density of instruments is also questioned and a discussion regarding the robustness of the algorithm against other faults is provided.

In the following studies, the parameters of the algorithm that allowed obtaining the best results in Section 8 are used, e.g. with the predicted statuses kept as faulty from one diagnosis procedure to another, once the status of an instrument is predicted as faulty, and with $|\Phi_v|_{min} = 10$. It is explained when the value of a parameter is changed.

9.1 Influence of the values of drift

As the results of Section 8 and the following were obtained for one draw of drift values, the case study was repeated for 100 draws. Table 4.9.1 presents the mean and standard deviation of each metric.⁴⁵ This table shows that:

- the mean values are equal to the values obtained in Table 4.8.1. This indicates that on average, the observations and conclusions do not depend on the variability of the drift.
- the standard deviations of all the metrics are close to zero. This is small considering the mean values and the relative standard deviation varying from zero to at most 16% of the mean value, reached for FN.

Thus, in this case study, the results of the diagnosis algorithm are not dependent on the values of the drift⁴⁶, with this drift model at least, and with this configuration of the algorithm.

⁴⁵The values for the delay of positive detection $\Delta \mathcal{D}$ are not provided here for the sake of clarity.

⁴⁶Note that the range of magnitude of drift does not change here, only the variability is studied. A different range would be equivalent to operating with different measuring instruments in practice.

$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$
Predicted status Ambiguous $ \begin{array}{c ccccccccccccccccccccccccccccccccccc$
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$
FP TP FDR PPV
Faulty μ σ μ σ μ σ μ σ
0.00 0.34 103.06 0.34 0.00 0.00 1.00 0.00
TNR FNR
$\mu \qquad \sigma \qquad \mu \qquad \sigma$
1.00 0.00 0.12 0.01
$egin{array}{cccccccccccccccccccccccccccccccccccc$
μ σ μ σ
0.00 0.00 0.00 0.00 0.00 0.00
$egin{array}{cccc} FPR & TPR \end{array}$
$\mu \qquad \sigma \qquad \mu \qquad \sigma$
$\begin{array}{cccccccccccccccccccccccccccccccccccc$

Table 4.9.1: Confusion matrix for 100 simulations of the case study with drift values drawn again for each one (metrics are defined in Section 5.4).

9.2 Influence of the true values

In the same way, the results of Section 8 were obtained for one draw of true values. Therefore, the case study was repeated for 100 draws of true values. Table 4.9.2 presents the mean and standard deviation of each metric.⁴⁷ This table shows that on average the number of true positives and false negatives are respectively more and less important than with the initial draw. The standard deviations are also close to zero which is small compared to most of the mean values, except for the metrics FP, NDP and FN, plus some the associated metrics derived from them (FNR, FPR, FOR).

Thus, again on average, the results obtained are globally better compared to the specific case study considered previously, in terms of NPV, FOR, accuracy and threat score notably.

In conclusion, the results obtained are dependent of the true values, at least when this model is considered. This is consistent with the studies carried out in Section 5.6 where we showed that the false results depends mostly on the measurement results associated to the rendez-vous.

 $^{^{47}}$ The values for the delay of positive detection $\Delta \mathcal{D}$ are not provided here for the sake of clarity.

Non-faulty Faulty Prevalence Accur μ σ μ σ μ σ μ 36.47 1.38 116.53 1.38 0.76 0.01 0.95	σ 0.02	
	0.02	
36.47 1.38 116.53 1.38 0.76 0.01 0.95		
	R	
TN FN NPV FO		
Non-faulty μ σ μ σ μ σ μ	σ	
35.59 1.19 7.40 2.95 0.83 0.06 0.17	0.06	
$NDN \qquad NDP$		
Predicted status Ambiguous μ σ μ σ		
0.00 0.00 0.10 0.33		
FP TP FDR PP	PPV	
Faulty μ σ μ σ μ	σ	
0.88 2.75 109.03 2.75 0.01 0.01 0.99	0.01	
TNR FNR		
$\mu \qquad \sigma \qquad \mu \qquad \sigma$		
0.98 0.03 0.06 0.03		
$egin{array}{cccccccccccccccccccccccccccccccccccc$		
μ σ μ σ		
0.00 0.00 0.00 0.00 0.00 0.00		
$egin{array}{cccc} FPR & TPR \end{array}$		
$\mu \qquad \sigma \qquad \mu \qquad \sigma$		
0.02 0.03 0.94 0.03		

Table 4.9.2: Confusion matrix for 100 simulations of the case study with true values following a 2D Gaussian model which parameters are randomly drawn for each simulation (metrics are defined in Section 5.4).

9.3 Influence of the model used for the true values

To reinforce the results on the influence of the values, the model used to derive the true values is changed in this section. Because the measurement results of the rendez-vous are samples of all the measurement results acquired during a time range, and because the predicted statuses are determined based on sets without taking into account the actual order of the measurements, we propose to randomly draw the true values at each time step and at each position following the uniform law $\mathcal{U}(0,400)$.

The case study with this model for the true values was repeated for 100 draws of true values. Table 4.9.3 presents the mean and standard deviation of each metric. ⁴⁸ The results of this table are similar to the ones of Table 4.9.2 in Section 9.2, except regarding false positives that are slightly higher. In general, the orders of magnitude are similar for the metrics. Thus, the model chosen for the true values does not seem to have a significant effect on the results. This statement does not take into account the fact that only random models were used here because of the

 $^{^{48}}$ The values for the delay of positive detection $\Delta \mathcal{D}$ are not provided here for the sake of clarity.

Non-faulty Faulty Prevalence Ac	ccuracy	
μ σ μ σ μ σ μ	μ σ	
63.43 0.57 89.57 0.57 0.59 0.00 0.9	92 0.02	
TN FN NPV	\overline{FOR}	
Non-faulty μ σ μ σ μ σ μ	μ σ	
59.97 2.43 8.07 3.39 0.88 0.04 0.3	12 0.04	
$NDN \qquad NDP$		
Predicted status Ambiguous μ σ μ σ		
0.00 0.00 0.15 0.36		
$egin{array}{cccccccccccccccccccccccccccccccccccc$	PPV	
Faulty μ σ μ σ μ σ μ	μ σ	
3.46 3.42 81.35 3.42 0.04 0.02 0.9	96 0.02	
\overline{TNR} FNR		
μ σ μ σ		
0.95 0.04 0.09 0.04		
$egin{array}{cccccccccccccccccccccccccccccccccccc$		
μ σ μ σ		
0.00 0.00 0.00 0.00 0.00 0.00		
$egin{array}{cccc} FPR & TPR \end{array}$		
μ σ μ σ		
$\begin{array}{cccccccccccccccccccccccccccccccccccc$		

Table 4.9.3: Confusion matrix for 100 simulations of the case study with true values randomly drawn following a uniform law (metrics are defined in Section 5.4).

mathematical objects used by the diagnosis algorithm, e.g. sets. Indeed, the order between the measured values is not exploited. We even observed in Section 5.6 that by considering only the measurement results obtained when an instrument is in rendez-vous to predict its status, we were actually performing a sampling of all its measurement results. This already breaks the continuity of the measured signal. In Appendix C, an additional case study is provided to confirm that the realism of the model for the true values does not have a major influence. Therefore, whether or not the relationship between two consecutive measured values is realistic, this has no significance for the diagnosis algorithm: it is aimed at determining a status over a time range and this time range is considered as a whole. In this way, it is the values that have the most significant effect as shown earlier in Section 5.6.

9.4 Influence of the density of instruments

In previous studies, few ambiguous statuses were predicted. This is principally due to the fact that most of the instruments of class 0 have often rendez-vous with the instruments of class 1 so that the size of the set of rendez-vous of an instrument s_i of class 0 with instruments of class 1 $\Phi(s_i \to S^1, (t_d, \Delta t))$ always satisfies $|\Phi(s_i \to S^1, (t_d, \Delta t))| \ge |\Phi_v|_{min}$. In this way, with

at most one or two executions of the main loop of the algorithm, all the instruments can be diagnosed faulty or non-faulty. However, if instruments were meeting less often, we expect that more ambiguous statuses would be produced.

Consequently, to simulate the same sensor network of 10 instruments having less rendez-vous, the number of accessible positions was increased from 100 to 2025^{49} . From Figure 4.9.1, we can observe that with 225 positions the values of the metrics deteriorate compared to the case with 100 positions and are even poorer for 400 positions and more. This can be observed just from the ACC going from more than 0.7 for 225 positions to less than 0.2 afterwards but the other metrics give interesting insights. Indeed, from Figure 4.9.1a, we observe that for more than 400 positions, there are only NDP and NDN. This explains why the TPR, FPR, TNR and FPR are equal to zero, and the PPV, FDR, NPV and FOR are not defined for more than 400 positions, in Figures 4.9.1b and 4.9.1c respectively. This does not only indicate that the diagnosis algorithm provides poor results but also that there is something making the algorithm unable to predict a different status than ambiguous for the measuring instruments.

Therefore, the number of positions has an influence on the performances. This indicates that the number of rendez-vous is at the heart of the diagnosis algorithm. If the number of rendez-vous taking place cannot be increased by performing changes in the sensor network for instance, the parameter that must be adjusted to have better performances is $|\Phi_v|_{min}$. Indeed, if a lower value of $|\Phi_v|_{min}$ can be acceptable, a sensor network with less rendez-vous occurring may be diagnosable.

In our case, while still varying the number of positions from 100 to 2025, $|\Phi_v|_{min}$ is set to 6 instead of 10. Figure 4.9.2 shows that even better results of diagnosis are obtained with 225 positions than with 100 (more TP, less FN, thus a higher TPR and PPV and a lower FNR and FOR, resulting in a higher ACC value notably).

While the diagnosis algorithm still provides acceptable results with 400 positions, although non-determined positives and negatives begin to occur, the performances remain still very poor for a higher number of positions as the number of ambiguous predictions drastically increases according to the evolution of NDP and NDN and the NDR. Thus, the value of ACC deteriorates more slowly than in Figure 4.9.1 but still very quickly.

Therefore, the diagnosis algorithm depends significantly of the density of instruments but more in terms of rendez-vous than of a physical density of instrument. Indeed, in Section 5, we defined the positions of the instruments in an area of 1×1 km. However, we could have chosen any other unit or subunit of length: this would have not changed the results we obtained. In practice, it means the sensor network can cover any area as long as the nodes meet regularly for diagnosis purposes.

⁴⁹The formula used to define these numbers of positions was the following: $((1+0.5\cdot i)\times 10)^2$ with $i\in[0..7]$.

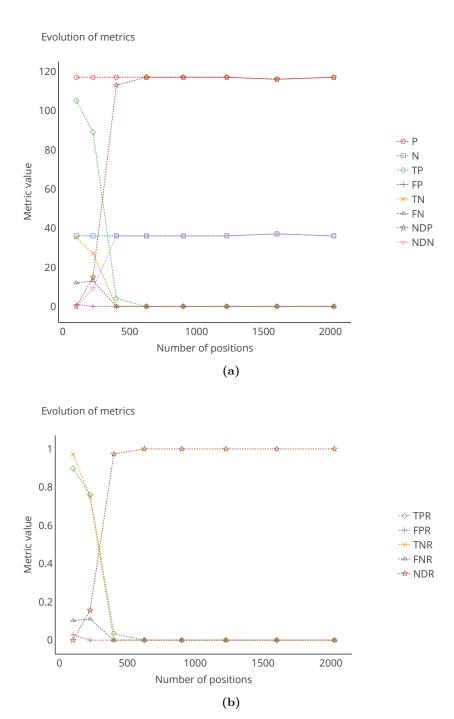
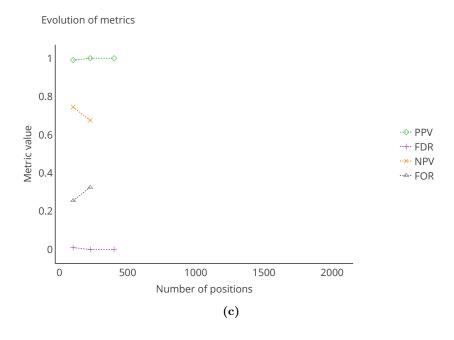


Figure 4.9.1: Evolution of the metrics computed over all the diagnosis procedures as a function of the number of accessible positions with $|\Phi_v|_{min} = 10$ (metrics are defined in Section 5.4). (continued)



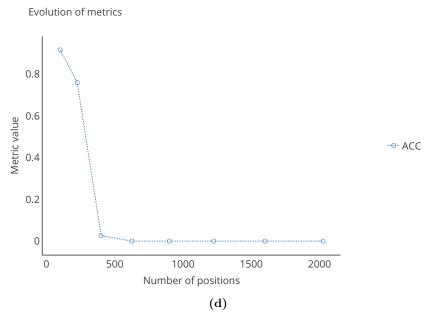


Figure 4.9.1: Evolution of the metrics computed over all the diagnosis procedures as a function of the number of accessible positions with $|\Phi_v|_{min} = 10$ (metrics are defined in Section 5.4).

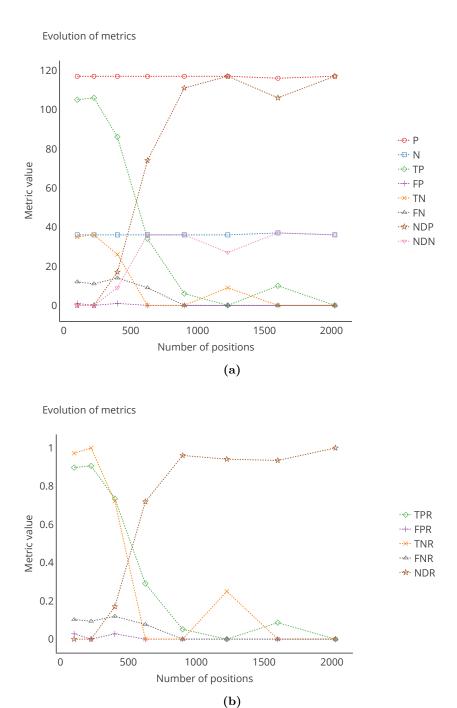
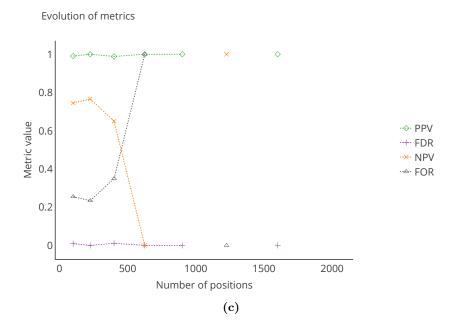


Figure 4.9.2: Evolution of the metrics computed over all the diagnosis procedures as a function of the number of accessible positions with $|\Phi_v|_{min} = 6$ (metrics are defined in Section 5.4). (continued)



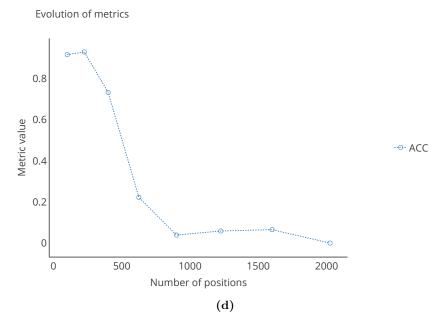


Figure 4.9.2: Evolution of the metrics computed over all the diagnosis procedures as a function of the number of accessible positions with $|\Phi_v|_{min} = 6$ (metrics are defined in Section 5.4).

9.5 Influence of other faults

 Concerning the influence of other faults on the results of diagnosis procedures, a reasoning similar to the one carried out in Section 5.6 for the study of false results can be made. Indeed, measurement results affected by other faults are likely to generate non-compatible measurement results with their true values. We can expect it is also the case regarding other instruments when such measurement results are associated to rendez-vous. Thus, it is very likely that, when they are valid, the rendez-vous of an instrument s_i associated to such measurement results affected by other faults are not belonging to the set of compatible valid rendez-vous of s_i , $\Phi_v^{\approx}(s_i, (t_d, \Delta t))$. Therefore, regarding the rates $r_{\Phi_v}^{\approx}(s_i \to S, (t_d, \Delta t))$, $r_{\Phi_v}^+(s_i \to S, (t_d, \Delta t))$ and $r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t))$ and their maximal associated values, the presence of measurement results affected by other faults in valid rendez-vous for an instrument will tend to generate positive results if they are not treated in advance, e.g. the instruments should be predicted as faulty earlier.

To illustrate this, spikes and noise were added to the measured values following the same drift model and with the same parameters that were used in Chapter 3 Section 5.3.1. The results of the diagnosis are given in Tables 4.9.4 and 4.9.5. From them, we observe that N is lower and P is higher than in Table 4.8.1. Therefore, the instruments should be considered as faulty earlier. Regarding all the other metrics, their values are still good and similar to the results of Table 4.8.1. Thus, it appears that the supplementary faulty measured values are not affecting significantly the results of the diagnosis algorithm. More importantly, despite a slightly higher standard deviation than in Table 4.8.2, the average delay of positive detection is close to zero and even positive. Its maximal value is also positive. It indicates that the prediction of instruments as faulty happens earlier than in Section 8.

To conclude, the proposed diagnosis algorithm targeting initially drift faults is robust to other faults. Although instruments are predicted as faulty more times, regardless if such a prediction is true or false, they are less harmful than false negative results in the context of drift diagnosis. Indeed, it calls for an earlier maintenance of the devices in this case instead of a late one.

9.6 Conclusion

In this section, several studies were conducted to investigate the sensitivity of the diagnosis algorithm to different effects. The following insights were brought:

- the results are not dependent on the values of the drift
- the true values, and thus the measurement results involved in the rendez-vous, have an influence on the results. This is consistent with previous results of Section 5.6
- the model used to derive the true values have less influence on the results than the values themselves
- the density of instruments is critical in terms of frequency of rendez-vous with other instruments
- the algorithm is robust to other faults. Their only effect if they are not managed prior to a drift diagnosis procedure is to make the predictions of instruments as faulty happening earlier. This could generate false positive results, which are less prejudicial than false negatives

		True status			
		Non-faulty $N = 23$	Faulty $P = 130$	Prevalence 0.85	Accuracy 0.90
Predicted status	Non faulty	TN	FN	\overline{NPV}	FOR
	Non-faulty	20	11	0.65	0.35
	Ambiguous	NDN	\overline{NDP}		
		0	2		
	Faulty	FP	TP	\overline{FDR}	PPV
		3	117	0.03	0.97
		TNR	\overline{FNR}		
		0.87	0.08		
		NDNR	NDPR	NDR	
		0.00	0.02	0.01	
		FPR	TPR		-
		0.13	0.90		

Table 4.9.4: Confusion matrix of the case study with the predicted statuses kept as faulty from one diagnosis procedure to another once the status of an instrument is predicted as faulty for the first time, $|\Phi_v|_{min} = 10$ and spikes and noise added to the measured values (metrics are defined in Section 5.4).

Del	ay of p	ositive d	etection
μ	σ	min	max
0.2	1.7	-2	3

Table 4.9.5: Statistics of the delay of positive detection $\Delta \mathcal{D}$ for the case study with the predicted statuses kept as faulty from one diagnosis procedure to another once the status of an instrument is predicted as faulty for the first time, $|\Phi_v|_{min} = 10$ and spikes and noise added to the measured values

10 Combination with a simple calibration approach

The diagnosis algorithm is primarily oriented at detecting when instruments require calibration. We investigate now how this algorithm behaves when calibration is carried out on an instrument after it was predicted as faulty. The diagnosis algorithm is configured as in Section 5, except for $|\Phi_v|_{min}$ which is set to 10 as a consequence of the study in Section 8. Therefore, **the status of an instrument that is predicted as faulty can change** as it is the trigger of a calibration procedure.

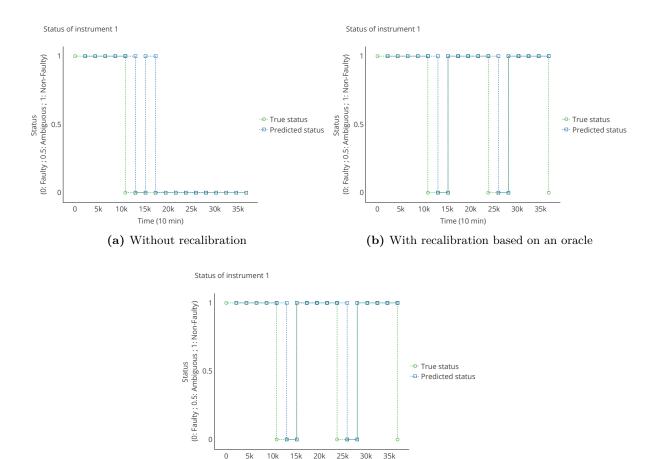
Three cases are considered when an instrument is detected as faulty:

- No calibration is carried out (equivalent to what was done in Section 5)
- Once an instrument is predicted as faulty, its gain and offset are respectively reinitialised to 1 and 0. It is equivalent to having an oracle that knows the gain and offset of each instrument at each instant. That is considered as a perfect calibration.
- A linear regression is carried out based on the couples of measured values of the faulty instruments and its diagnosers that are in the set of valid rendez-vous used for the diagnosis. This technique is usually one of the firsts to be tested for the calibration of measuring instruments [139].

Results for a single instrument are displayed in Figure 4.10.1. The instrument is recalibrated right after being diagnosed as faulty for both recalibration methods. Nevertheless, it is only visible at the next diagnosis procedure. Also from these curves we cannot assess the quality of the calibration, it only shows that it was efficient enough to be again above the threshold mentioned in the specification of the case study.

Table 4.10.1 represents the results for the case without calibration. Tables 4.10.2 and 4.10.3 display the values of the metrics computed on the case study with the oracle-based and linear regression-base calibration approaches respectively. Results are disturbing as the tables give slightly better results without calibration. However this is normal. In the cases where there is actually recalibration performed, instruments are changing of true status several times (this is shown by the values of N, P and the prevalence in each table). In Section 5.5, we observed that most of the false results happen around the moment where an instrument actually becomes faulty. When instruments are recalibrated, they still continue to drift and they may change of status again over time. Therefore, more false results are likely to be obtained when a calibration strategy is applied. For this reason, the performances between the case study for which no calibration is carried out and the ones where it is, based on an oracle or on linear regression here, cannot be compared in terms of metrics without taking into account this fact.

That being said, the behaviour of the instruments is the same in terms of values and drift whether the calibration is performed with an oracle or with linear regression: it is only a different correction that is applied to the instruments. Thus, the results of the diagnosis algorithm can be compared between Tables 4.10.2 and 4.10.3. From these tables, it appears that the calibration based on linear regression provides better results in terms of improvement of the quality of the instruments' measurements: N is greater (consequently P is lower) and there are four less false negative for three more true positive and negative and one more false positive. Therefore, the accuracy with this method of calibration is slightly better. Although that the values of all the metrics are of the same order of magnitude, it is disturbing that the oracle-based method gives less accurate results, yet it corresponds to an ideal recalibration. In fact, like Figure 4.10.1, these tables do not assess the actual quality of the calibration but only the quality of the diagnosis. It only shows that indeed some instruments are corrected so that they are truly non-faulty again—and can be detected legitimately as so.



(c) With recalibration based on linear regression with the set of valid rendez-vous used at the diagnosis instant

Time (10 min

Figure 4.10.1: Evolution of true and predicted status for instrument s_1 for different choices of calibration

To proceed to an evaluation of the impact of the calibration on the measurement results, the error model introduced in Chapter 3 is considered and its parameters are estimated for each instrument considering the entire time interval of study. Table 4.10.4 presents the statistics of the parameters and score of the model and Figure 4.10.3 represents a target plot of the instruments as a function of their slope and intercept. We observe that both the calibration approaches give equivalent results compared to the case without calibration. The error on the slope and on the intercept are reduced of at least 50% and 64% respectively.

However, even with the oracle-based calibration, we observe that the remaining average offset can still be considered as significant. This residual is explained by the fact that the gain and offset of the instruments are corrected at specific moments and not at each time step: between two diagnosis procedures, the instruments continue to drift. Moreover, when an instrument is considered as faulty, the correction is applied on the values that are measured after the moment this prediction is made. This is illustrated in Figure 4.10.2. Thus, it explains why when we compute the parameters of the error model on the entire time range of study, the slope and intercept are not equal to their ideal values (one and zero respectively).

More significantly, the average score of the linear regression on the error model increases from 0.84 to 0.97 at least indicating a good connection between the calibrated values and the

		True status			
		Non-faulty $N = 36$	Faulty $P = 117$	Prevalence 0.76	Accuracy 0.88
Predicted status	Non faulty	TN	FN	\overline{NPV}	FOR
	Non-faulty	36	17	0.68	0.32
	Ambiguous	NDN	NDP		
		0	2		
	Faulty	FP	TP	\overline{FDR}	PPV
		0	98	0.00	1.00
		TNR	\overline{FNR}		
		1.00	0.15		
		NDNR	NDPR	NDR	
		0.00	0.02	0.01	
		FPR	TPR		
		0.00	0.84		

Table 4.10.1: Confusion matrix of the case study on calibration without recalibration (metrics are defined in Section 5.4).

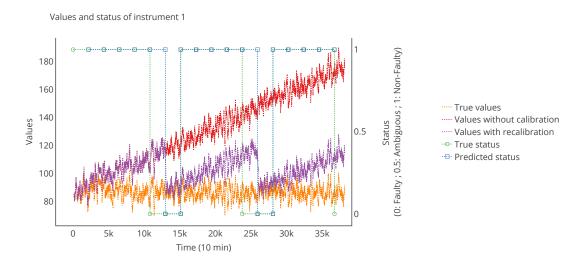


Figure 4.10.2: Evolution of the true values, of the measured values without and with recalibration by linear regression for s_1 . The true and predicted statuses are also plotted to show when recalibrations are triggered.

true values compared to the measured values and the true values if no calibration is performed. Thus, a simple calibration approach, based on the information contained in the rendez-vous used to predict the statuses of the instruments, allows obtaining an interesting improvement of their metrological performances.

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To conclude, this case study shows that, in addition to being able to detect quite accurately when an instrument does not meet a specification, the diagnosis algorithm can also provide useful

		True status			
		Non-faulty $N = 107$	Faulty $P = 46$	Prevalence 0.30	Accuracy 0.83
Predicted status	Non-faulty	TN	FN	\overline{NPV}	FOR
	non-raurty	107	26	0.80	0.20
	Ambiguous	NDN	NDP		
		0	0		
	Faulty	FP	TP	FDR	PPV
		0	20	0.00	1.00
		TNR	FNR		
		1.00	0.57		
		NDNR	NDPR	NDR	
		0.00	0.00	0.00	
		FPR	TPR		-
		0.00	0.43		

Table 4.10.2: Confusion matrix of the case study on calibration with an oracle-based calibration (metrics are defined in Section 5.4).

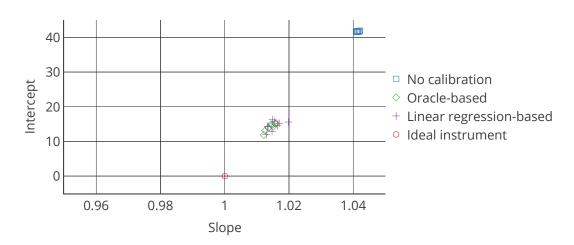


Figure 4.10.3: Target plot of the 9 nodes of class 0 of the network as a function of their slope and intercept following the error model of Chapter 3, computed on the entire time interval of study, for each calibration strategy.

data to perform the recalibration of measuring instruments through the sets of valid rendez-vous used during diagnosis procedures.

		True status			
		Non-faulty $N = 110$	Faulty $P = 43$	Prevalence 0.28	Accuracy 0.85
	Non-faulty	TN 109	FN 22	NPV 0.83	FOR 0.17
Predicted status	Ambiguous	NDN	NDP		
	Faulty	<i>FP</i> 1	<i>TP</i> 21	FDR 0.05	PPV 0.95
		TNR 0.99	FNR 0.51		
		NDNR 0.00	NDPR 0.00	NDR 0.00	
		FPR 0.01	TPR 0.49		

Table 4.10.3: Confusion matrix of the case study on calibration with calibration based on linear regression (metrics are defined in Section 5.4).

Calibration approach	Slope		I	Intercept		Score		
	μ	σ	μ	σ		μ	σ	
No calibration	1.04	0.00	4:	2 0		0.84	0.00	
Oracle-based	1.01	0.00	1	4 1		0.97	0.01	
Linear regression-based	1.02	0.00	1	5 1		0.98	0.01	

Table 4.10.4: Statistics of the slope, intercept and score of the error model of Chapter 3 computed for the nodes of class 0 of the network on the entire time interval of study and for each calibration strategy.

11 Conclusion

In this chapter, a novel algorithm for the diagnosis of drift faults in sensor networks was introduced. It exploits the concept of rendez-vous between measuring instruments. A key property of this algorithm is that it does not require any assumption of the type of sensor network. The only hypothesis needed concern the assumption as always non-faulty of at least a subset of instruments of the highest metrological class in the sensor network.

Based on an initial case study that showed promising results, we discussed the adjustment of the different parameters of the algorithm. While we mostly provided guidelines to adjust them empirically, notably to reduce the number of false results, an algorithm was proposed to facilitate the adjustment of the minimal number of required valid rendez-vous to allow a prediction between the status faulty and non-faulty $|\Phi_v|_{min}$. It gives an upper boundary for the value of this parameter. In our case study, it allowed to identify that the value previously considered was too high to ensure a good diagnosis. Lowering the value of this parameter improved significantly the results. In summary, the recommendations provided across the chapter for all the parameters are summarised in Table 4.11.1.

A study of the parameters of the case study influencing the results of the diagnosis algorithm was conducted. It showed the measurement results of the instruments, in particular those involved in rendez-vous, and the frequency of rendez-vous between instruments are key factors regarding respectively the results of the diagnosis algorithm and the adjustment of its parameters. The algorithm was also shown to be robust to other faults in the measurement results. Indeed, such faulty values are likely to generate false positives, which may be less prejudicial than false negatives.

Finally, the diagnosis algorithm was associated to a simple calibration approach. When an instrument is diagnosed as faulty, the valid rendez-vous that were used to make the prediction of its status are exploited to make a linear regression between the values of the faulty instrument and the values from its diagnosers. This allowed to successfully correct the gain and offset of the faulty measuring instruments. Therefore, the diagnosis algorithm that was presented opens new perspectives on *in situ* calibration as this algorithm does more than indicating which instruments are faulty in a sensor network: it also provides information that can be exploited to correct them.

Parameter	Effect	Recommendation
Conditions of validity of measurement results	Determine if a measurement result can be compared to another.	It should be determined based on the properties of the measuring instruments deployed.
Conditions of spa- tiotemporal vicin- ity to be in rendez- vous	Determine if two measurement results can be compared to each other.	It can be determined based on expert specifications or on characteristics of the measurand and of the measuring instruments.
Instant of diagnosis t_d	Defines the reference time of the time range on which the diagnosis is performed.	Can be determined based on the characteristics of the instruments used and the context of deployment.
Periodicity of the diagnosis	Determines the frequency at which the algorithm is computed.	It can be determined based on the characteristics of the instruments used and the context of deployment.
Length of the time range Δt	Defines the length of the time range on which the diagnosis is performed.	It can be determined based on the characteristics of the instruments used.
Minimal relative difference of class $\Delta c_{min}^{\mathcal{D}}(k)$	Determines which classes of instruments can be the diagnosers of instruments belonging to a given class k .	The higher the relative class is, the more accurate are the diagnosers. It should be set to $+1$ or 0 but not to a negative value.
Maximal tolerated values for the different rates $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$	Defines the triggers to determine if an instrument is faulty or non-faulty.	Too high values are likely to generate false negatives while too low values generate false positives. It should be adjusted to favour false positives instead of false negatives, while minimising their number.
Minimal required number of valid rendez-vous $ \Phi_v _{min}$	Determines if a given set of valid rendez-vous is large enough to determine if an instrument is faulty or non-faulty.	Considering a defined Δt and using the algorithm provided, a value can be chosen based on the number of faulty instruments tolerated in the network. A sufficiently high $ \Phi_v _{min}$ should be chosen so that the content of the sets Φ_v are large enough to likely have representative sets.

 ${\bf Table\ 4.11.1:}\ {\bf Recommendations\ for\ the\ different\ parameters\ of\ the\ diagnosis\ algorithm$

Conclusion and perspectives

1 Conclusion

In this thesis, we tackled the issue of calibration for sensor networks, a question that is particularly critical when low-cost instruments are deployed.

In the first part of this thesis, we reviewed the existing works concerning in situ calibration which are techniques aiming at calibrating measuring instruments while leaving them in the field, preferably without any physical intervention on them. Across this survey, we proposed a concise taxonomy enabling the classification of the algorithms. Four main groups of categories were defined, two of them following properties of the sensor networks that may be exploited, namely the presence of reference instruments in the network and the mobility of the nodes, and the other two concerning respectively the grouping strategy at stake in the algorithm and the kind of mathematical relationship that can be derived. This extensive survey showed that there are numerous contributions on the subject, covering a large spectrum of the cases that can be encountered when deploying sensor networks. However, a quantitative comparison of the performances of the algorithm was shown to be difficult because algorithms are often applied on specific case studies in their original publication. Indeed, there is no standard use cases or protocols to allow such comparisons.

Therefore, a framework for the simulation of sensor networks aimed at evaluating the performances of in situ calibration algorithms was proposed in the second part of this thesis. It was applied to the comparison of seven in situ calibration algorithms for blind and static sensor networks. We independently changed the model used to derive the true values of the measurand, the drift and fault model, the number of instruments and parameters of some calibration algorithms. We showed that even with a particular care to conduct a balanced comparison, different conclusions can be drawn by slightly changing the case study. Thus, in addition to being a protocol enabling the effective comparison of strategies, the proposed framework could be used to choose the most appropriate in situ calibration algorithm during the design of a sensor network.

However, we also observed that some calibration algorithms may correct measuring instruments regardless of their actual performances, particularly concerning those for blind sensor networks. Therefore, it is possible that instruments which measurement results are still correct regarding the metrological specifications of the instruments, have their values corrected. In the worst case, they may have their values degraded. Thus, a diagnosis algorithm for drift faults in sensor networks was proposed in the last part of this thesis. This algorithm is particularly interesting because it is based on rendez-vous between measuring instruments. This concept is more complex than a simple comparison of measured values as it encapsulates the idea of a necessary spatio-temporal vicinity so that measurement results can be relevantly compared. In addition, the algorithm demands no assumption on the type of sensor network used, requiring only that some of the best instruments in the network can be assumed as non-faulty. Across a case study, we showed that the algorithm allowed to successfully identify drifting instruments. In a second step, a better adjustment of the different parameters of the algorithm enabled to improve the results, although our guidelines are mainly empirical. Then, by investigating the influence of the design

choices of our case study, we showed the measurement results involved in rendez-vous and the frequency of rendez-vous between instruments, are key factors regarding respectively the results of the diagnosis algorithm and the adjustment of its parameters. In addition, the algorithm is robust to other faults: faulty values are likely to generate false positives, which may be less prejudicial than false negatives in the context of environmental sensing. Finally, while combining the diagnosis algorithm to a naive calibration approach, it was possible to successfully correct the gain and offset of the faulty measuring instruments based on the measurement results of rendez-vous between instruments, while leaving the measurement results of non-faulty ones unchanged. Indeed, the diagnosis algorithm also provides information that can be exploited to correct measuring instruments.

2 Perspectives

Several perspectives can be considered for the direct pursuit of this work but also for the scientific community and the general public from a larger scope.

Direct perspectives

Extensions of the diagnosis algorithm

In the appendices, several extensions of the diagnosis algorithm are related. They aim at being able to apply the diagnosis algorithm we proposed in any situation and with the same performances. However, their development is not complete and demands supplementary investigations. Moreover, the adjustment of some parameters of the algorithm is empirical and may not be conducted through an automated optimisation procedure. Eventually, being able to ensure that the specification stating when instruments should be considered as faulty or not, is verified, or at least correctly implemented, would be valuable. Therefore, investigating if means could be brought to improve the adjustment of the parameters of the algorithm and the formal qualification of its performances would be of a major interest.

Design of sensor networks

In this work, we proposed a framework that could be used to identify the relevant algorithms for the *in situ* calibration of a sensor network in its environment of deployment. To facilitate the development of sensor networks, there is a need for engineering tools to identify at the design stage what the data processing algorithms are from which a sensor network would benefit, on top of tools already optimising their deployment in the case of static sensor networks for instance [18]. This demands to aggregate the various results obtained by the scientific community from various fields in a dedicated product. Such an integrated tool is not currently available in the literature.

Perspectives for the scientific community and the general public

Standardisation of the vocabulary for the description of $in\ situ$ calibration algorithms

In this work, we proposed a taxonomy for the classification of *in situ* calibration algorithms. We observed that there were also other taxonomies. An intensive work of standardisation of this vocabulary and on the practices to describe and evaluate these algorithms would be valuable in order to structure the work of the scientific community on this subject. Calls were recently published along the same lines [60, 133].

Data processing standardisation

In scientific publications exploiting measurement results from low-cost sensor networks, data processing that are carried out to have a clean dataset are not always fully described. For measuring instruments on the market, even for certified ones, the way the measurements are processed is rarely mentioned or only as a black box. Concerning sensors as components of measuring instruments, technical datasheets are often scarce and imprecise on how they behave over time and depending on their operating conditions. In the scientific community, there are calls for a standardisation of the data processing algorithms in order to determine how far processed measurement results are from raw ones [60, 133]. We extend this call to the need for a clear information on how measuring instruments and their components behave from a metrological perspective, to facilitate the identification and use of adapted data processing algorithms.

Qualification of instruments and their processing

Although experiments under laboratory or field condition conducted during this thesis were not reported in this manuscript, actual measuring instruments and systems were studied in the early months of this thesis through the Sense-City platform [38] as it was done in the studies reported in Chapter 1 Section 1.3.3. Regarding low-cost instruments and considering the cost range we considered in this work (US \$100 to US \$10,000), there is a wide variety of quality among these measuring instruments and systems. In the scientific community, there are reports on the performances of measuring instruments but it is not mandatory for a manufacturer to provide the results of such works with its devices. In fact, it is currently difficult for an end user to identify clearly which instrument suits its needs and at which cost because the information provided is quite poor. Therefore, an effort of qualification of the performances of measuring instruments, with the data processing algorithms they use, and a transparent information of the general public about it when instruments are put on the market would be valuable.

Environmental footprint of low-cost environmental sensor network

Currently, there is little information on the environmental footprint of low-cost measuring instruments [78]. While their interest can be legitimately justified whatever the measurand considered, it has to be counterbalanced by their impact on the environment. We are dealing here most of the time with electronic devices that may be difficult to recycle and may require rare materials. In addition, they also require a computer and telecommunication infrastructure, those being known to have more and more impact on the greenhouse gases emissions around the world. Being able to evaluate the global impact of the entire measuring chain, from the instruments to the datacentres in which measurement results are stored and exploited, would be of a major interest in order to determine the actual benefits of dense low-cost sensor networks compared to what is currently achieved with the information brought by more traditional environmental measuring systems.

Appendix A

Diagnosis Algorithm for Drift Faults in Sensor Networks: Extensions

1 Formulation reducing the number of iterations of the algorithm

The formulation of Algorithm 1 given in Chapter 4 Section 4.2 is not optimised in terms of iterations. Indeed, considering the statuses of instruments are likely predicted based on the results obtained for instruments of a higher class than their own (based on the recommendation to set the minimal relative differences of class required $\Delta c_{min}^{\mathcal{D}}(k)$ to +1 or 0 in Chapter 4 Section 5.7), it would be relevant to browse the set of instruments to diagnose $S^{\mathcal{D}}$ by descending classes of instruments instead of browsing it regardless of the classes of the instruments. In addition, actualising the predicted statuses at the end of the evaluation of the instruments of each class would allow using the instruments predicted as non-faulty at the previous iteration.

Algorithm 3 provides this optimised formulation in terms of iterations.

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Algorithm 3: Algorithm of the diagnosis procedure proposed for the detection of drift in sensor networks with its number of iterations optimised

```
Data: S, \Phi(S \to S, (t_d, \Delta t)), |\Phi_v|_{min}, \Delta c_{min}^{\mathcal{D}}
Result: \Omega(S, t_d)
/* Initiate the set of instruments to diagnose and the predicted statuses
     of the instruments
                                                                                                                                   */
S^{\mathcal{D}} \leftarrow S \setminus S^{c_{max}}
\widehat{\Omega}(S^{c_{max}}, t_d) \leftarrow NF
\widehat{\Omega}(S^{\mathcal{D}}, t_d) \leftarrow \mathbf{A}
\widetilde{\Omega}(S^{\mathcal{D}}, t_d) \leftarrow \widehat{\Omega}(S^{\mathcal{D}}, t_d)
                                                                 /* Initiate the actualised statuses */
/* Ignore the instruments that cannot have enough valid rendez-vous
for s_i \in S^{\mathcal{D}} do
     if |\Phi(s_i \to S, (t_d, \Delta t))| < |\Phi_v|_{min} then |S^{\mathcal{D}} \leftarrow S^{\mathcal{D}} \setminus \{s_i\}
end
/* Predict the status of each instrument to diagnose
                                                                                                                                   */
     \widehat{\Omega}(S, t_d) \leftarrow \widetilde{\Omega}(S, t_d)
                                              /* The actualised statuses are now the predicted
      statuses */
```

```
for k \leftarrow c_{max} - 1 to 0 do /* For each class of instruments in descending
               order
                      \widehat{\Omega}'(S, t_d) \leftarrow \widehat{\Omega}(S, t_d)
                                                                                                                       /* A local instance of the future predicted
                           statuses is created, to take into account the results of the
                          predictions for the previous values of k in this loop */
                       for s_i \in S^{\mathcal{D}} \cap S^k do
                                 /* Build the current set of valid rendez-vous for s_i, using
                                              \widehat{\Omega}'(S,t_d) instead of \widehat{\Omega}(S,t_d)
                                                                                                                                                                                                                                                                                        */
                                 c_{min}^{\mathcal{D}}(s_i) \leftarrow c(s_i) + \Delta c_{min}^{\mathcal{D}}(c(s_i))
                                 \Phi_v(s_i \to S, (t_d, \Delta t)) \leftarrow \{\varphi(s_i \to s_j, t) \in \Phi(s_i \to S, (t_d, \Delta t)), \text{ such as } \}
                                              s_j \notin S^{\mathcal{D}}, \ \widehat{\Omega}'(s_j, t') = NF, \ c(s_j) \ge c_{min}^{\mathcal{D}}(s_i),
                                              m(s_i, t) \in M^*(s_i, (t_d, \Delta t)) \text{ and } m(s_j, t') \in M^*(s_j, (t_d, \Delta t))
                                 /* If s_i have enough valid rendez-vous, then compute the
                                              different rates to actualise its status
                                                                                                                                                                                                                                                                                        */
                                 if |\Phi_v(s_i \to S, (t_d, \Delta t))| \ge |\Phi_v|_{min} then
                                            r_{\Phi_v}^{\approx}(s_i \to S, (t_d, \Delta t)) \leftarrow \frac{|\Phi_v^{\approx}(s_i \to S, (t_d, \Delta t))|}{|\Phi_v(s_i \to S, (t_d, \Delta t))|}
                                            r_{\Phi_v}^+(s_i \to S, (t_d, \Delta t)) \leftarrow \frac{|\Phi_v^+(s_i \to S, (t_d, \Delta t))|}{|\Phi_v(s_i \to S, (t_d, \Delta t))|}
                                            r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) \leftarrow \frac{|\Phi_v^-(s_i \to S, (t_d, \Delta t))|}{|\Phi_v(s_i \to S, (t_d, \Delta t))|}
                                             /* If a condition on the different rates is met, then the
                                                        actualised status of s_i is set to faulty, otherwise, it is
                                                         set to non-faulty
                                             if r_{\Phi_v}^+(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^+)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} \text{ or } r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) > 
                                               (1 - r_{\Phi_n}^{\approx}(s_i \to S, (t_d, \Delta t))) > (r_{\Phi_n}^+ + r_{\Phi_n}^-)_{max} then
                                                       \Omega(s_i, t_d) \leftarrow F
                                              else
                                                       \Omega(s_i, t_d) \leftarrow \text{NF}
                                             S^{\mathcal{D}} \leftarrow S^{\mathcal{D}} \setminus \{s_i\} /* s_i is diagnosed so it can be removed from S^{\mathcal{D}}
                                 end
                       end
           end
until \widehat{\Omega}(S,t_d)=\widehat{\Omega}(S,t_d) /* Repeat until there is no difference between the
   predicted and actualised statuses */
```

2 Diagnosis with the prediction as non-faulty based on the highest sufficient class

In Chapter 4 Section 4.2, the prediction of an instrument s_i as non-faulty or faulty is based on the set of valid rendez-vous $\Phi_v(s_i \to S, (t_d, \Delta t))$ if the condition regarding the minimal number of valid rendez-vous required $|\Phi_v|_{min}$ is satisfied. $\Phi_v(s_i \to S, (t_d, \Delta t))$ is notably defined with the help of the minimal class allowed to be a diagnoser of s_i , $c_{min}^{\mathcal{D}}(s_i)$.

However, suppose that $\exists k > c_{min}^{\mathcal{D}}(s_i)$ such as $|\Phi_v(s_i \to S^{k+}, (t_d, \Delta t))| > |\Phi_v|_{min}$. In this case, the status of s_i can be predicted with a smaller number of instruments. This is interesting

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because instruments of higher class usually have a lower measurement uncertainty. Thus, if the predicted status of s_i , $\widehat{\Omega}(s_i, t_d)$, is determined as non-faulty, a higher confidence in s_i can be granted: it is predicted as non-faulty based on its compatibility with instruments of a higher class than the minimal one required to diagnose it. Consequently, we can estimate that the measurement results of s_i are more accurate than expected. By predicting the statuses of the instruments in this manner, the results of the diagnosis algorithm would be enriched with a useful information.

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Nevertheless, if s_i is predicted as faulty according to $\Phi_v(s_i \to S^{k+}, (t_d, \Delta t))$, it is still possible to obtain $\widehat{\Omega}(s_i, t_d) = NF$ with the set of valid rendez-vous $\Phi_v(s_i \to S, (t_d, \Delta t))$. In this case, it is still preferable to have s_i predicted as non-faulty despite a prediction as faulty with a subset of the diagnoser instruments.

It is also possible to obtain $\widehat{\Omega}(s_i, t_d) = NF$ according to $\Phi_v(s_i \to S^{k+}, (t_d, \Delta t))$ but $\widehat{\Omega}(s_i, t_d) = F$ according to $\Phi_v(s_i \to S, (t_d, \Delta t))$. We estimate the prediction based on $\Phi_v(s_i \to S, (t_d, \Delta t))$ is more likely a false positive as less accurate instruments are considered than in S^{k+} . Such a case may be very rare. Indeed, the less the diagnoser instruments are accurate, the more the predicted statuses have a chance to be equal to non-faulty. This is due to mean used to determine the compatibility between measurement results: it is based on the measurement uncertainty. This time, it is preferable to have s_i predicted as non-faulty with a subset of the diagnoser instruments despite a faulty prediction that can be made with all of them.

Therefore, Algorithm 1 must be adapted to include this idea, expressed as a diagnosis algorithm with negative prediction based on the highest sufficient class. It is described in Algorithm 4^{50} .

Algorithm 4: Algorithm of the diagnosis procedure proposed for the detection of drift in sensor networks with the prediction as non-faulty based on the highest sufficient class

```
Data: S, \Phi(S \to S, (t_d, \Delta t)), |\Phi_v|_{min}, \Delta c_{min}^{\mathcal{D}}
Result: \Omega(S, t_d)
/* Initiate the set of instruments to diagnose and the predicted statuses
      of the instruments
S^{\mathcal{D}} \leftarrow S \setminus S^{c_{max}}
\widehat{\Omega}(S^{c_{max}}, t_d) \leftarrow NF
\widehat{\Omega}(S^{\mathcal{D}}, t_d) \leftarrow \mathbf{A}
\widetilde{\Omega}(S^{\mathcal{D}}, t_d) \leftarrow \widehat{\Omega}(S^{\mathcal{D}}, t_d)
                                                                          /* Initiate the actualised statuses
                                                                                                                                                     */
/* Ignore the instruments that cannot have enough valid rendez-vous
for s_i \in S^{\mathcal{D}} do
     if |\Phi(s_i \to S, (t_d, \Delta t))| < |\Phi_v|_{min} then |S^{\mathcal{D}} \leftarrow S^{\mathcal{D}} \setminus \{s_i\}
end
/* Predict the status of each instrument to diagnose
                                                                                                                                                     */
repeat
      \widehat{\Omega}(S, t_d) \leftarrow \widetilde{\Omega}(S, t_d)
                                                    /* The actualised statuses are now the predicted
     statuses */
for s_i \in S^{\mathcal{D}}
\begin{vmatrix} c_{min}^{\mathcal{D}}(s_i) \leftarrow c(s_i) + \Delta c_{min}^{\mathcal{D}}(c(s_i)) \\ k \leftarrow c_{max} \end{vmatrix}
```

⁵⁰The improvements brought with Algorithm 3 in Section 1 are not added for the sake of clarity.

```
repeat
                    /* Build the current set of valid rendez-vous for s_i with the
                            instruments in S^{k+}
                                                                                                                                                                           */
                    \Phi_v(s_i \to S^{k+}, (t_d, \Delta t)) \leftarrow \{\varphi(s_i \to s_i, t) \in \Phi(s_i \to S^{k+}, (t_d, \Delta t)), \text{ such as } \}
                           s_j \notin S^{\mathcal{D}}, \ \widehat{\Omega}'(s_j, t') = NF, \ c(s_j) \ge c_{min}^{\mathcal{D}}(s_i),
                           m(s_i, t) \in M^*(s_i, (t_d, \Delta t)) \text{ and } m(s_j, t') \in M^*(s_j, (t_d, \Delta t))
                    /* If s_i have enough valid rendez-vous, then compute the
                            different rates to actualise its status
                                                                                                                                                                           */
                    if |\Phi_v(s_i \to S^{k+}, (t_d, \Delta t))| \ge |\Phi_v|_{min} then
                         \begin{split} |\Phi_v(s_i \to S^{\kappa+}, (t_d, \Delta t))| &\geq |\Psi_v| \min \text{ then } \\ r_{\Phi_v}^\approx(s_i \to S^{k+}, (t_d, \Delta t)) &\leftarrow \frac{|\Phi_v^\approx(s_i \to S^{k+}, (t_d, \Delta t))|}{|\Phi_v(s_i \to S^{k+}, (t_d, \Delta t))|} \\ r_{\Phi_v}^+(s_i \to S^{k+}, (t_d, \Delta t)) &\leftarrow \frac{|\Phi_v^+(s_i \to S^{k+}, (t_d, \Delta t))|}{|\Phi_v(s_i \to S^{k+}, (t_d, \Delta t))|} \\ r_{\Phi_v}^-(s_i \to S^{k+}, (t_d, \Delta t)) &\leftarrow \frac{|\Phi_v^-(s_i \to S^{k+}, (t_d, \Delta t))|}{|\Phi_v(s_i \to S^{k+}, (t_d, \Delta t))|} \end{split}
                           /* If a condition on the different rates is met, then the
                                  actualised status of s_i is set to faulty, otherwise, it is
                                  set to non-faulty
                          if r_{\Phi_v}^+(s_i \to S^{k+}, (t_d, \Delta t)) > (r_{\Phi_v}^+)_{max} or r_{\Phi_v}^-(s_i \to S^{k+}, (t_d, \Delta t)) > (r_{\Phi_v}^-)_{max} or (1 - r_{\Phi_v}^{\approx}(s_i \to S^{k+}, (t_d, \Delta t))) > (r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max} then
                              \Omega(s_i, t_d) \leftarrow F
                           else
                                 \widetilde{\Omega}(s_i, t_d) \leftarrow \mathrm{NF}
                           end
                    end
                    k \leftarrow k-1
             until k < c_{min}^{\mathcal{D}}(s_i) or \widetilde{\Omega}(s_i, t) = NF
             if \Omega(s_i, t_d) \neq A then
               S^{\mathcal{D}} \leftarrow S^{\mathcal{D}} \setminus \{s_i\} /* s_i is diagnosed so it can be removed from S^{\mathcal{D}} */
              end
until \widehat{\Omega}(S,t_d)=\widetilde{\Omega}(S,t_d) /* Repeat until there is no difference between the
  predicted and actualised statuses */
```

3 From a centralised to a decentralised computation

From its description provided in Chapter 4 Section 4.2 and its pseudo-code in Algorithm 1, our diagnosis algorithm has to be computed in a centralised manner. Indeed, all the rendez-vous between the instruments must be known, with the associated measurement results and information to determine the validity of the rendez-vous, and along with the predicted and actualised status of the instruments during the different iterations of the main loop in the algorithm.

However, it is possible to consider a distributed calculation, notably if the measurement results are not collected in a centralised manner.

Consider an instrument s_i .

We assume s_i knows $\Delta c_{min}^{\mathcal{D}}(s_i)$, $|\Phi_v|_{min}$, $(r_{\Phi_v}^+)_{max}$, $(r_{\Phi_v}^-)_{max}$ and $(r_{\Phi_v}^+ + r_{\Phi_v}^-)_{max}$.

It is also possible to assume it knows that $\widehat{\Omega}(S^{c_{max}}, t_d) = NF$ due to the assumption of the instruments of class c_{max} as non-faulty required for the algorithm⁵¹.

Finally, s_i knows $\Phi(s_i \to S, (t_d, \Delta T))$. We assume when two instruments meet, they exchange information related on the rendez-vous they have (measurement results and so on).

Consider $S^{\varphi}(s_i, (t_d, \Delta t))$ which is the set of instruments involved in a rendez-vous with s_i between $[t_d - \Delta t; t_d]^{52}$:

$$S^{\varphi}(s_i, (t_d, \Delta t)) = \{s_j \text{ such as } \exists \varphi(s_i \to s_j, t') \in \Phi(s_i \to S, (t_d, \Delta T))\}$$

Based on it, s_i can define a local set of instruments to diagnose $S^{\mathcal{D}} = (s_i \cup S^{\varphi}(s_i, (t_d, \Delta t))) \setminus S^{c_{max}}$.

Depending on $S^{\mathcal{D}}$, s_i can derive $\Phi_v(s_i \to S, (t_d, \Delta T))$ if it knows $\widehat{\Omega}(s_j, t_d)$ for $s_j \in S^{\varphi}(s_i, (t_d, \Delta t)) \setminus S^{\mathcal{D}}$. Then, s_i can determine its actualised status as indicated in Algorithm 1. If $s_i \notin S^{\mathcal{D}}$ after this, its predicted status is known. Otherwise, s_i has to try to determine it again.

In fact, the most challenging step is to access to $\widehat{\Omega}(s_j, t_d)$ for $s_j \in S^{\mathcal{D}} \setminus \{s_i\}$ to actualise $S^{\mathcal{D}}$ between each iteration of the algorithm.

Calls must be sent to these instruments asking if s_j determined if $\Omega(s_j, t_d)$ is equal to NF or F.

There is, however, a need for a synchronisation mechanism between the node. Indeed, because the prediction of each instrument's status depends on exchanges of information with others, some instruments may be able to perform iterations of the algorithm faster or slower than others. To avoid issues by mixing predicted statuses that were not determined after the same iteration of the diagnosis procedure, we define a loop counter $n^{\mathcal{D}}$. This counter is incremented at each repetition of the main loop of the algorithm. When s_i calls the instruments $s_j \in S^{\mathcal{D}} \setminus \{s_i\}$ to ask for $\widetilde{\Omega}(s_j, t_d)$, it also indicates that this call is related to the loop $n^{\mathcal{D}}$.

After receiving the responses, s_i removes the s_j from its local $S^{\mathcal{D}}$ if $\widetilde{\Omega}(s_j, t_d) \neq A$ and it can again try to determine its status.

At some point, $\widetilde{\Omega}(S^{\varphi}(s_i, (t_d, \Delta t), t_d))$ may not change anymore. Thus, like in Algorithm 1, if $\widehat{\Omega}(S^{\varphi}(s_i, (t_d, \Delta t), t_d)) = \widetilde{\Omega}(S^{\varphi}(s_i, (t_d, \Delta t), t_d))$ after an iteration of the diagnosis procedure, the algorithm terminates and $\widehat{\Omega}(s_i, t_d) = A$.

The distributed formulation of the diagnosis algorithm is provided in Algorithm 5. By construction, due to its synchronous behaviour, it is certain that the results are identical to the ones obtained in a centralised manner.

The protocol of communication between the nodes, either when they are in rendez-vous or when they exchange actualised statuses, is not detailed in this pseudo-code. Moreover, we suppose that these communications are managed in a separate thread from the one in which is run the diagnosis algorithm to blockings. Finally, Algorithm 5 does not manage the case where:

⁵¹Or a subset of $S^{c_{max}}$. See Chapter 4 Section 6.

⁵²Therefore, $\Phi(s_i \to S^{\varphi}(s_i, (t_d, \Delta t)), (t_d, \Delta T)) = \Phi(s_i \to S, (t_d, \Delta T)).$

- the diagnosis algorithm run by s_i has terminated after the loop $n^{\mathcal{D}'}$ and $\widehat{\Omega}(s_i, t_d) = A$.
- $\exists s_i$ such as it is still trying to determine its predicted status during a loop $n^{\mathcal{D}} > n^{\mathcal{D}'}$.

Thus, $s_j \in S^{\varphi}(s_i, (t_d, \Delta t))$ during the loop $n^{\mathcal{D}}$. s_i will ask s_j for its actualised status after the loop $n^{\mathcal{D}}$ even if the status of s_j was definitively predicted as ambiguous. We can assume that in practice s_j can tell s_i after the loop $n^{\mathcal{D}'}$ that $\widehat{\Omega}(s_j, t_d)$ will not change anymore and thus avoid useless communications. This would require few changes in Algorithm 5 that were not written for the sake of clarity.

In conclusion, it is possible to apply our diagnosis algorithm in a decentralised manner. The solution proposed is, however, not unique and other protocols could be considered.

Algorithm 5: Algorithm of the diagnosis procedure proposed for the detection of drift in sensor networks written in a distributed manner. Each node of the network runs this pseudo-code.

```
Data: S^{\varphi}(s_i, (t_d, \Delta t)), \Phi(s_i \to S, (t_d, \Delta t)), |\Phi_v|_{min}, c^{\mathcal{D}}_{min}(s_i)
Result: \Omega(S, t_d)
/* Initiate the set of instruments to diagnose and the predicted statuses
       of the instruments
S^{\mathcal{D}} \leftarrow (s_i \cup S^{\varphi}(s_i, (t_d, \Delta t))) \setminus S^{c_{max}}
\widehat{\Omega}(S^{c_{max}}, t_d) \leftarrow \text{NF}
\widehat{\Omega}(S^{\mathcal{D}}, t_d) \leftarrow \mathbf{A}
\widetilde{\Omega}(S^{\mathcal{D}}, t_d) \leftarrow \widehat{\Omega}(S^{\mathcal{D}}, t_d)
                                                                                  /* Initiate the actualised statuses */
/* Ignore the instruments that cannot have enough valid rendez-vous
                                                                                                                                                                       */
      if |\Phi(s_i \to S, (t_d, \Delta t))| < |\Phi_v|_{min} then
         |S^{\mathcal{D}} \leftarrow S^{\mathcal{D}} \setminus \{s_i\}|
end
n^{\mathcal{D}} \leftarrow 0
/* Predict the status of each instrument to diagnose
                                                                                                                                                                       */
repeat
      n^{\mathcal{D}} \leftarrow n^{\mathcal{D}} + 1
      \widehat{\Omega}(S, t_d) \leftarrow \widetilde{\Omega}(S, t_d)
                                                          /* The actualised statuses are now the predicted
        statuses */
      /* Build the current set of valid rendez-vous for s_i
                                                                                                                                                                       */
       \Phi_v(s_i \to S, (t_d, \Delta t)) \leftarrow \{\varphi(s_i \to s_j, t) \in \Phi(s_i \to S, (t_d, \Delta t)), \text{ such as } \}
              s_j \notin S^{\mathcal{D}}, \ \widehat{\Omega}(s_j, t') = NF, \ c(s_j) \ge c_{min}^{\mathcal{D}}(s_i),
              m(s_i, t) \in M^*(s_i, (t_d, \Delta t)) \text{ and } m(s_i, t') \in M^*(s_i, (t_d, \Delta t))
      /* If s_i have enough valid rendez-vous, then compute the different
              rates to actualise its status
                                                                                                                                                                       */
      if |\Phi_v(s_i \to S, (t_d, \Delta t))| \ge |\Phi_v|_{min}

\begin{vmatrix} r_{\Phi_v}^{\approx}(s_i \to S, (t_d, \Delta t)) &\leftarrow \frac{|\Phi_v^{\approx}(s_i \to S, (t_d, \Delta t))|}{|\Phi_v(s_i \to S, (t_d, \Delta t))|} \\ r_{\Phi_v}^+(s_i \to S, (t_d, \Delta t)) &\leftarrow \frac{|\Phi_v^+(s_i \to S, (t_d, \Delta t))|}{|\Phi_v(s_i \to S, (t_d, \Delta t))|} \\ r_{\Phi_v}^-(s_i \to S, (t_d, \Delta t)) &\leftarrow \frac{|\Phi_v^-(s_i \to S, (t_d, \Delta t))|}{|\Phi_v(s_i \to S, (t_d, \Delta t))|} \end{aligned}
```

4 Multiple measurands

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Low-cost sensors are prone to drift due to influence quantities affecting the indication provided by the instrument. Thus, it is frequent that the calibration relationship of such instruments has to take into account their influence quantities. This is easily applicable in a measurement system embedding a main instrument measuring the principal measurand targeted by the system and several other instruments that measure the values of the influence quantities of the main instrument.

Therefore, the sensor network addresses more than one measurand in this case. It can be seen as a "network of sensor networks", each one considering a specific measurand. In this way, some sensor networks are dependent on others to determine their measured values. The diagnosis of the instruments of such systems is discussed in this section.

4.1 General idea

We propose here a general concept of protocol to carry out the diagnosis of drift fault in sensor networks composed of measurements systems in which one or more instruments has influence quantities measured by other instruments of the same system.

The general idea is the following:

- 1. Carry out the diagnosis for each sub-network measuring a same quantity and for which the instruments have no influence quantities following the algorithm described in Chapter 4 Section 4.2.
- 2. Perform the diagnosis for each measuring a same quantity and for which the instruments have influence quantities measured by a sub-network previously diagnosed. The way to carry this diagnosis out is discussed in the following subsection.
- 3. Repeat step 2 until the diagnosis has been performed on each sub-network, e.g. for each measurand.

4.2 Diagnosis of drift faults in a sub-network with instruments having influence quantities

In fact, we aim at using the same algorithm already defined in Chapter 4 Section 4.2. Indeed, the presence of influence quantities questions only the validity of the measurement results.

We remind we assume that the sensor network is composed of measurement systems, in which the influence quantities of one or more of its instruments are measured by other instruments that also belong to the same system.

Consider a sub-network with one influence quantity, a diagnosis procedure d that happens at t_d and suppose we know the predicted statuses at t_d of the instruments measuring the influence quantities.

Consider a rendez-vous between two instruments s_i and s_j , $\varphi(s_i \to s_j, t)$ with $t \in [t_d - \Delta t; t_d]^{53}$, with s_j supposed non-faulty and of class $c(s_j) \ge c_{min}^{\mathcal{D}}$. The instruments measuring the influence quantity are denoted s_i' and $s_j'^{54}$.

Let us discuss the validity of the rendez-vous. First regarding the measured values of s_i , three different cases appear:

 $^{^{53}}$ In this case, we could also speak of a rendez-vous between two systems because all the pairs of instruments measuring the same quantities in the systems i and j are in rendez-vous

⁵⁴We also suppose when two systems are in rendez-vous, we have measurements for all the measurands. This assumption could be discussed in more details as it requires to extend the definition of rendez-vous and to discuss the case of instruments of a same system with different acquisition frequencies. Such particular cases are not treated here

- s'_i is non-faulty: the measured value of s'_i can be used to determine the one of s_i with its calibration relationship. In this case, the measurement result $m(s_i, t)$ is valid and the rendez-vous may be valid too⁵⁵.
- s'_i is faulty but s'_j is non-faulty: the measured value of s'_j can be used to determine the one of s_i with its calibration relationship. In this case, $m(s_i, t)$ is again valid the rendez-vous may be valid too 56 .
- s'_i and s'_j are faulty: the measured value of s_i cannot be determined as no value of its influence quantity can be reliably used with the calibration relationship. In this case, $m(s_i,t)$ is not valid. Therefore, $\varphi(s_i \to s_j,t)$ is **not valid**.

The same reasoning can be applied to s_i . ⁵⁷⁵⁸

In fact, the presence of influence quantities has only an influence on the validity of the measurement results. Thus, the Equation 4.1 from Chapter 4 Section 4.1 used to determine $\Phi_v(s_i \to S, (t_d, \Delta t))$ is still valid. We recall this equation:

```
\Phi_{v}(s_{i} \to S, (t_{d}, \Delta t)) = \{ \varphi(s_{i} \to s_{j}, t) \in \Phi(s_{i} \to S, (t_{d}, \Delta t)), \text{ such as} 
s_{j} \notin S^{\mathcal{D}}, // s_{j} \text{ is not an instrument to diagnose} 
\widehat{\Omega}(s_{j}, t') = NF, // s_{j} \text{ is non-faulty} 
c(s_{j}) \geq c_{min}^{\mathcal{D}}(s_{i}), // \text{ The class of } s_{j} \text{ is higher or equal to the minimal} 
class \text{ allowed to diagnose } s_{i} 
m(s_{i}, t) \in M^{*}(s_{i}, (t_{d}, \Delta t)), // \text{ The measurement result of } s_{i} \text{ is valid} 
m(s_{j}, t') \in M^{*}(s_{j}, (t_{d}, \Delta t)) \} // \text{ The measurement result of } s_{j} \text{ is valid} 
(A.1)
```

This is why, apart from the conditions on the validity of the measurement results that may change, the diagnosis algorithm remains applicable in the context of measuring instruments with influence quantities.

4.3 Conclusion

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In this section, we presented a manner to extend our diagnosis algorithm for drift faults in sensor networks to the case where

However, a formulation enriching the results provided by the diagnosis algorithm could be developed in future work. Indeed, the fact that values of measuring instruments from other systems can be used to determine the status of an instrument in a given system could allow us to determine a status at the level of the measuring system for instance or at least to qualify the predicted statuses in light of how the validity of measurement results is determined.

5 Diagnosis algorithm for drift faults in sensor networks with an eventbased formulation

In the first place, the diagnosis algorithm was presented as being applied at an instant t_d over $[t_d - \Delta t; t_d]$. It means that it is time which triggered the diagnosis and that the determination of

 $^{^{55}}$ Depending on the other conditions defined for the validity of the measurement results.

⁵⁶See note 55

⁵⁷This development can be extended to the case where there are multiple influences quantities but also to the case where more than two instruments meet at the same time, e.g. the values of the influence quantity measured by a third system could be used when s'_i and s'_j are faulty in our case.

⁵⁸The definitions related the compatibility between two measurement results could be extended based on these three cases.

the sets of rendez-vous is notably based on a time range.

However, notably in the perspective of a real-time application, it could be valuable to trigger the diagnosis when rendez-vous occurs. A rendez-vous can be assimilated to an **event** able to trigger a diagnosis procedure. Such rendez-vous is noted $\varphi(...)$, "..." standing for the instruments actually involved in the rendez-vous and the time instant at which it occurred.

In the same way, the contents of the sets of rendez-vous could be determined based on a number of rendez-vous, noted $|\Phi|^{\mathcal{D}}$, that occurred before $\varphi(...)$. It can be useful if the number of rendez-vous on $[t_d - \Delta t; t_d]$ varies a lot between the instruments. For instance, some of them may not be diagnosable with not enough (valid) rendez-vous on this time range. A solution is to increase Δt . However in this case, other instruments may have their statuses not correctly predicted with a greater Δt (see Chapter 4 Section 7). Thus, replacing Δt by $|\Phi|^{\mathcal{D}}$ can help to overcome this situation: all the instruments will be diagnosed based on sets of rendez-vous that are equivalent in terms of size.

In addition to the couple $(t_d, \Delta t)$ defining the moment of the diagnosis and the time period on which it is made, there are other combinations possible to set these characteristics of the algorithm with the help of $\varphi(...)$ and $|\Phi|^{\mathcal{D}}$:

- $(t_d, |\Phi|^{\mathcal{D}})$: the diagnosis occurs at t_d and the last $|\Phi|^{\mathcal{D}}$ rendez-vous are considered for each instrument to predict its status.
- $(t_d, \Delta t, |\Phi|^{\mathcal{D}})$: the diagnosis occurs at t_d and the rendez-vous that happened in $[t_d \Delta t; t_d]$ are considered, up to the $|\Phi|^{\mathcal{D}}$ last ones.
- $(\varphi(...), \Delta t)$: the diagnosis occurs on $\varphi(...)$ and the rendez-vous that happened in $[t_d \Delta t; t_d]$ are considered, with t_d being the instant at which $\varphi(...)$ was generated.
- $(\varphi(...), |\Phi|^{\mathcal{D}})$: the diagnosis occurs on $\varphi(...)$ and the last $|\Phi|^{\mathcal{D}}$ rendez-vous are considered for each instrument to predict its status.
- $(\varphi(...), \Delta t, |\Phi|^{\mathcal{D}})$: the diagnosis occurs at $\varphi(...)$ and the rendez-vous that happened in $[t_d \Delta t; t_d]$ are considered, up to the $|\Phi|^{\mathcal{D}}$ last ones⁵⁹.

The diagnosis algorithm can be easily adapted to consider these alternate definitions of the instant $\alpha \in \{t_d, \varphi(...)\}$ at which is executed and of the duration $\beta \in \{\Delta t, |\Phi|^{\mathcal{D}}, (\Delta t, |\Phi|^{\mathcal{D}})\}$ used to define the sets used by the algorithm. Indeed, the expression " $(t_d, \Delta t)$ " can be replaced by " (α, β) " straightforward in Algorithm 1⁶⁰.

To conclude, the proposed diagnosis algorithm can also be applied with an event-based definition of the instant at which it is executed and of the duration used to define the sets of rendez-vous it uses.

6 Real-time diagnosis algorithm of drift faults in sensor networks

Until then, the diagnosis was considered as a procedure carried out periodically or, with the definitions of Section 5, on events. If it is performed at each time step or on each event, the

- contains less than $|\Phi|^{\mathcal{D}}$ rendez-vous if less than that occurred during $[t_d \Delta t; t_d]$.
- does not contain all the rendez-vous that occurred during $[t_d \Delta t; t_d]$ if there were more rendez-vous than $|\Phi|^{\mathcal{D}}$.

⁵⁹When Δt and $|\Phi|^{\mathcal{D}}$ are combined it is possible that a set of rendez-vous:

⁶⁰In fact, behind "..." in φ (...) lies a " t_d " which is the instant at which the rendez-vous occurred. This t_d is used in practice to determine the sets involved in a diagnosis procedure. Also behind " $|\Phi|^{\mathcal{D}}$ " lies a " t_d " which is the instant when the first of the $|\Phi|^{\mathcal{D}}$ rendez-vous occurred. This t_d is different for each instrument. The definitions given in Chapter 4 Section 3 have to be adapted depending on the choice of α and β .

algorithm is not very efficient: it computes rates based on sets that are not often changing (and probably not a lot) between two time steps. This section aims at providing elements for the design of a more efficient way to perform in real-time the diagnosis of drift faults in a sensor network.

6.1 Choice of an event-based approach

To provide a real-time formulation of the diagnosis algorithm, we choose to rely on an event-based approach, at least regarding α following the notations of Section 5.

This choice is guided by the fact that the sets of rendez-vous $\Phi(s_i \to S, (\alpha, \beta))$ for each $s_i \in S$ do not have to be actualised at each time step but only when a new rendez-vous happens.

However, if $\beta = \Delta t$, it invites to determine the statuses of the instruments based on a sliding window of rendez-vous that changes at each time step. This case is discussed afterwards.

6.2 General idea

Consider two instruments s_i and s_j . When a rendez-vous happens between them, at t for instance, $\Phi(s_i \to S, (\alpha, \beta))$ and $\Phi(s_j \to S, (\alpha, \beta))$ are changed. Thus, the predicted statuses of s_i and s_j have to be actualised.

To determine the statuses of the instruments after this rendez-vous, the principles introduced in Chapter 4 Section 4.1 have to be extended.

Indeed, we must determine which instrument between s_i and s_j is the diagnoser instrument for the other. The set of diagnosers of s_i at t is noted $S_{diagnosers}(s_i, t)$.

At t, in a real-time approach, we have values for $\widehat{\Omega}(s_i,t)$ and $\widehat{\Omega}(s_j,t)$. Depending on the case, the rendez-vous $\varphi(s_i \to s_j,t)$ will be considered either in $\Phi_v(s_i \to S, (\alpha,\beta))$ or $\Phi_v(s_j \to S, (\alpha,\beta))$ (sets of valid rendez-vous).

The following cases are possible:

$\widehat{\Omega}(s_i,t)$	$\widehat{\Omega}(s_j,t)$	Decision
NF	NF	The diagnoser must be chosen between s_i and s_j .
NF	A	s_i is a diagnoser for s_j .
NF	\mathbf{F}	s_i is a diagnoser for s_j .
A	NF	s_j is a diagnoser for s_i .
A	A	The rendez-vous cannot be considered as valid by any of the instruments.
A	\mathbf{F}	The rendez-vous cannot be considered as valid by any of the instruments.
\mathbf{F}	NF	s_j is a diagnoser for s_i .
\mathbf{F}	A	The rendez-vous cannot be considered as valid by any of the instruments.
\mathbf{F}	\mathbf{F}	The rendez-vous cannot be considered as valid by any of the instruments.

Table A.6.1: Possible cases when a new rendez-vous occurs between s_i and s_j for a real-time diagnosis

In fact, there is only one particular case: when $\widehat{\Omega}(s_i,t) = \widehat{\Omega}(s_j,t) = NF$. To choose the one to use as a diagnoser, we propose to rely on the class and on the duration since the instruments were predicted as non-faulty. The diagnoser is the one that has first the highest class and then the one that has been predicted as non-faulty first.

Once this information is determined, e.g. s_j is a diagnoser for s_i for instance, the set of valid rendez-vous $\Phi_v(s_i \to S, (\alpha, \beta))$ is actualised following a derived version of Equation 4.1 from

4282 Chapter 4 Section 4.1:

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\Phi_{v}(s_{i} \to S, (\alpha, \beta)) = \{\varphi(s_{i} \to s_{j}, t) \in \Phi(s_{i} \to S, (\alpha, \beta)), \text{ such as } s_{j} \in S_{diagnosers}(s_{i}, t), // s_{j} \text{ is a diagnoser of } s_{i}
\widehat{\Omega}(s_{j}, t') = NF, // s_{j} \text{ is non-faulty}
c(s_{j}) \geq c_{min}^{\mathcal{D}}(s_{i}), // \text{ The class of } s_{j} \text{ is higher or equal to the minimal } class \text{ allowed to diagnose } s_{i}
m(s_{i}, t) \in M^{*}(s_{i}, (\alpha, \beta)), // \text{ The measurement result of } s_{i} \text{ is valid}
m(s_{j}, t') \in M^{*}(s_{j}, (\alpha, \beta)) \} // \text{ The measurement result of } s_{j} \text{ is valid}
(A.2)
```

Then, the predicted status of s_i can be actualised following the same principles described in Chapter 4 Section 4.2 (rates, criteria...). Because the predicted status of only of the instruments involved in the rendez-vous is actualised, it would be valuable to use a decentralised approach.

6.3 Initialisation of the algorithm

Like in Chapter 4 Section 4.2, all the instruments in $S^{c_{max}}$ are assumed as non-faulty, the set of instruments to diagnose $S^{\mathcal{D}}$ is equal to $S \setminus S^{c_{max}}$ and the predicted statuses $\widehat{\Omega}(s_i, t_d)$ are initialised to non-faulty for $s_i \in S^{c_{max}}$ and to ambiguous for $s_i \in S^{\mathcal{D}}$.

At t = 0, for each $s_i \in S$, $\Phi(s_i \to S, (t, \beta)) = \emptyset$. Overtime, these sets will be filled with rendez-vous. Only when $\Phi_v(s_i \to S, (\alpha, \beta)) \ge |\Phi_v|_{min}$ for $s_i \in S^{\mathcal{D}}$, the predicted statuses can begin to change.

6.4 Allowed changes of status

With the principles presented, the predicted statuses of the instruments can change from one rendez-vous to another like in Chapter 4 Section 4.2. It may be preferable to consider an approach where the predicted status of an instrument is kept equal to faulty once it is first diagnosed as so as suggested in Chapter 4 Section 7 to avoid false results.

6.5 Conclusion

In this section, we provided avenues toward the definition of a real-time formulation of our diagnosis algorithm. We suggested using an event-based approach following the concepts introduced in Section 5. We introduced a general idea to actualise the predicted status of the instruments when a new rendez-vous occurs and discussed the particular cases. We presented the criteria of initialisation which are equivalent to those of Chapter 4 and we recommended using the approach of Chapter 4 Section 7 where the predicted status of an instrument is kept equal to faulty once it is first diagnosed. To validate the ideas presented, it would be necessary to fully develop the algorithm and compared its execution on a same case study where our original diagnosis algorithm⁶¹ would be applied at each time step.

Afterwards, the algorithm giving an upper boundary to $|\Phi_v|_{min}$, the minimal size for a set of valid rendez-vous to allow a prediction between the status faulty and non-faulty could be extended for the cases where the diagnosis is performed in real-time.

 $^{^{61}}$ With the predicted status of the instruments kept equal to faulty once first diagnosed as so.

Appendix **B**

On the Reason Why Assuming the Instruments of Class cmax are not Drifting is Acceptable

In this appendix, more details are briefly given to understand why assuming the instruments of class c_{max} are not drifting is acceptable.

In Chapter 4 Section 6.2, it appeared that even with s_9 drifting, the diagnosis algorithm still provides satisfying results when assuming this instrument as always non-faulty.

Suppose that there was a means to correctly predict the status of s_9 . Following the reasoning that was used to design the algorithm in Chapter 4 Section 4, this instrument should not be considered at some point for the prediction of the statuses of other instruments. Nevertheless, based on the results obtained in the study we conducted, the information provided by this instrument with its measurement results is of a sufficient quality to help the prediction of the other instruments.

In fact, this can be explained by the definition used for r_{true} in Chapter 4 Section 5.3. r_{true} uses the set of measurements results compatible with the true values $M^{\approx}(s_i, (t, \Delta))$. The compatibility of a measurement with its true value is appreciated regarding the uncertainty on the measured value and the true value. For the instruments of class c_{max} here, the uncertainty, absolute or relative, is significantly smaller than the uncertainty of the instruments of lower class, particularly in this case study (see Table 4.5.1 in Chapter 4 Section 5 where $\Delta_r v(s_9) \ll \Delta_r v(s_j)$ for all the other instruments s_j). Thus, with this definition of r_{true} , when two instruments of different classes have their true statuses equal to faulty for the first time, it does not mean that their drift is equivalent. It is possible that an instrument of a given class is still performing well relatively to an instrument of a lower class, which is the case here with s_9 drifting more slowly: $\delta G_{max}(s_9) \ll \delta G_{max}(s_j)$ and $\delta O_{max}(s_9) \ll \delta O_{max}(s_j)$. This is the most important feature and such relationships would likely be valid in practice with instruments of different metrological class.

Therefore, to better appreciate the true status of a given instrument regarding one of a lower class (or not), the compatibility of its measurement results with its true values should be evaluated considering the highest measurement uncertainty between these two devices. In our case, the true status of s_9 is always non-faulty if $M^{\approx}(s_i, (t, \Delta))$ is determined with the relative uncertainty $\Delta_r v$ of the instruments of class zero than with its own $\Delta_r v$ as shown in Figure B.0.1.

In conclusion, assuming that s_9 is not drifting and is always non-faulty is valid from the point of view of all the other instruments of lower class in the network because even when drifting, this instrument is still more accurate than the instruments of a lower class. The metric

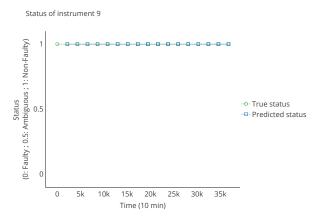


Figure B.0.1: Evolution of the true status of s_9 , computed with the relative uncertainty $\Delta_r v$ on the results of the instruments of class 0 instead of its own $\Delta_r v$, and its predicted status when it is drifting

used to determine if an instrument is truly faulty or not is not fairly appreciating the status of an instrument when the goal is to compare its measurement results to the ones of another instrument, potentially of a worse quality.

Appendix C

Sensitivity of the Diagnosis Algorithm: Case Study with Values Related Over Time for the Parameters of the True Values' Model

1 Introduction

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We extend here the results of Chapter 4 Section 9.3 on the influence of the model used for the true values in the case study developed in Section 5 of the same chapter. More precisely, an additional case study is provided to confirm that the realism of the model for the true values does not have a major influence.

2 Model used

We propose to use for the true values the same model as in Chapter 4 Section 3 but instead of using values drawn randomly for the amplitude and FWHM of the Gaussian curve at each time step, the dataset that was exploited in Chapter 3 Section 3 is considered here to derive these values. Thus, the values of the amplitude and FWHM of the Gaussian curve are related from one time step to another.

In this dataset, the values are given hourly. As we used a time step equivalent to 10min in Chapter 4 Section 3, the dataset has to be resampled. The missing values are interpolated

Apart from the model for the true values, the case study is identical to the one of Chapter 4 Section 8 (case study of Section 3 with improvements regarding false results and the adjustment of $|\Phi_v|_{min}$).

Results

Table C.3.1 gives the values of each metric. Compared to the results of Table 4.8.1 and considering there is a slightly higher prevalence here (0.88 instead of 0.76), e.g. more faulty cases that non-faulty ones, we observe there are less false negative results (10 instead of 14). The number of true negatives TN is equal to the number of negative cases N (18). Thus, there are no false positives. There are also no predicted statuses equal to ambiguous.

Therefore, the TNR and TPR are high (1.00 and 0.93) and so is the PPV (1.00). Only the NPV and FOR are worse than with a random draw of the parameters of the model of the true values (0.64 and 0.36 respectively instead of 0.72 and 0.28). Although, considering the low

		True status			
		Non-faulty $N = 18$	Faulty $P = 135$	Prevalence 0.88	Accuracy 0.93
	Non-faulty	TN	FN	\overline{NPV}	FOR
		18	10	0.64	0.36
Duralistad atataa	Ambiguous	NDN	NDP		
Predicted status		0	0		
	Faulty	FP	TP	\overline{FDR}	PPV
		0	125	0.00	1.00
		TNR	\overline{FNR}		
		1.00	0.07		
		$\overline{}$ $NDNR$	\overline{NDPR}	$\overline{}$ NDR	-
		0.00	0.00	0.00	
		\overline{FPR}	TPR		•
		0.00	0.93		

Table C.3.1: Confusion matrix of the case study with true values based on the 2D Gauss model with values of its parameters derived from a real dataset

number of TN due to the low value of N, this is not major.

The overall performance from the point of view of the accuracy is slightly better than in Table 4.8.1 (0.93 instead of 0.91)

4 Conclusion

 This case study confirms the conclusion from Chapter 4 Section 9.3, e.g. the model chosen for the true values does not seem to have a significant effect on the results. Indeed, the mathematical objects used by the diagnosis algorithm are sets: the order between the measured values is not exploited. Combined to the observation of Chapter 4 Section 7 (the only measurement results considered to the status of a measuring instrument are the ones obtained when it is in rendez-vous. Thus, a sampling of all its measurement results is performed, breaking already the continuity of the measured signal), the relationship between two consecutive measured values has no significance for the diagnosis algorithm: it is aimed at determining a status over a time range and this time range is considered as a whole. Therefore, it is the values that have the most significant effect as shown earlier in Chapter 4 Section 7.

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Titre : Étalonnage in situ de l'instrumentation bas coût pour la mesure de grandeurs ambiantes : méthode d'évaluation des algorithmes et diagnostic des dérives

Mots clés: Réseaux de capteurs, Qualité de mesure, Étalonnage, Evaluation, Diagnostic, Mesures ambiantes

Résumé: Dans de nombreux domaines allant de l'agriculture à la santé publique, des grandeurs ambiantes doivent être suivies dans des espaces intérieurs ou extérieurs. On peut s'intéresser par exemple à la température, aux polluants dans l'air ou dans l'eau, au bruit, etc. Afin de mieux comprendre ces divers phénomènes, il est notamment nécessaire d'augmenter la densité spatiale d'instruments de mesure. Cela pourrait aider par exemple à l'analyse de l'exposition réelle des populations aux nuisances comme les polluants atmosphériques.

Le déploiement massif de capteurs dans l'environnement est rendu possible par la baisse des coûts des systèmes de mesure, qui utilisent notamment des éléments sensibles à base de micro ou nano technologies. L'inconvénient de ce type de dispositifs est une qualité de mesure insuffisante. Il en résulte un manque de confiance dans les données produites et/ou une hausse drastique des coûts de l'instrumentation causée par les opérations nécessaires d'étalonnage des instruments ou de remplacement périodique des capteurs.

Il existe dans la littérature de nombreux algorithmes qui offrent la possibilité de réaliser l'étalonnage des instruments en les laissant déployés sur le terrain, que l'on nomme techniques d'étalonnage *in situ*.

L'objectif de cette thèse est de contribuer à l'effort de recherche visant à améliorer la qualité des données des instruments de mesure bas coût à travers leur étalonnage *in situ*.

En particulier, on vise à 1) faciliter l'identification des techniques existantes d'étalonnage *in situ* applicables à un réseau de capteurs selon ses propriétés et les caractéristiques des instruments qui le composent ; 2) aider au choix de l'algorithme le plus adapté selon le réseau de capteurs et son contexte de déploiement ; 3) améliorer l'efficacité des stratégies d'étalonnage *in situ* grâce au diagnostic des instruments qui ont dérivé dans un réseau de capteurs.

Trois contributions principales sont faites dans ces

travaux. Tout d'abord, une terminologie globale est proposée pour classer les travaux existants sur l'étalonnage *in situ*. L'état de l'art effectué selon cette taxonomie a montré qu'il y a de nombreuses contributions sur le sujet, couvrant un large spectre de cas. Néanmoins, le classement des travaux existants selon leurs performances a été difficile puisqu'il n'y a pas d'étude de cas de référence pour l'évaluation de ces algorithmes.

C'est pourquoi dans un second temps, un cadre pour la simulation de réseaux de capteurs est introduit. Il vise à guider l'évaluation d'algorithmes d'étalonnage in situ. Une étude de cas détaillée est fournie à travers l'évaluation d'algorithmes pour l'étalonnage in situ de réseaux de capteurs statiques et aveugles. Une analyse de l'influence des paramètres et des métriques utilisées pour extraire les résultats est également menée. Les résultats dépendant de l'étude de cas, et la plupart des algorithmes réétalonnant les instruments sans évaluer au préalable si cela est nécessaire, un outil d'identification permettant de déterminer les instruments qui sont effectivement fautifs en termes de dérive serait précieux.

Dès lors, la troisième contribution de cette thèse est un algorithme de diagnostic ciblant les fautes de dérive dans les réseaux de capteurs sans faire d'hypothèse sur la nature du réseau de capteurs considéré. Basé sur le concept de rendez-vous, l'algorithme permet d'identifier les instruments fautifs tant qu'il est possible de supposer qu'un instrument n'est pas fautif dans le réseau de capteurs. À travers l'analyse des résultats d'une étude de cas, nous proposons différents moyens pour diminuer les faux résultats et des recommandations pour régler les paramètres de l'algorithme. Enfin, nous montrons que l'algorithme de diagnostic proposé, combiné à une technique simple d'étalonnage, permet d'améliorer la qualité des résultats de mesure. Ainsi, cet algorithme de diagnostic ouvre de nouvelles perspectives quant à l'étalonnage in situ.

Title: In situ calibration of low-cost instrumentation for the measurement of ambient quantities: evaluation methodology of the algorithms and diagnosis of drifts

Keywords : Sensor networks, Measurement quality, Calibration, Evaluation, Diagnosis, Environmental sensing

Abstract: In various fields going from agriculture to public health, ambient quantities have to be monitored in indoors or outdoors areas. For example, temperature, air pollutants, water pollutants, noise and so on have to be tracked. To better understand these various phenomena, an increase of the density of measuring instruments is currently necessary. For instance, this would help to analyse the effective exposure of people to nuisances such as air pollutants.

The massive deployment of sensors in the environment is made possible by the decreasing costs of measuring systems, mainly using sensitive elements based on micro or nano technologies. The drawback of this type of instrumentation is a low quality of measurement, consequently lowering the confidence in produced data and/or a drastic increase of the instrumentation costs due to necessary recalibration procedures or periodical replacement of sensors.

There are multiple algorithms in the literature offering the possibility to perform the calibration of measuring instruments while leaving them deployed in the field, called *in situ* calibration techniques.

The objective of this thesis is to contribute to the research effort on the improvement of data quality for low-cost measuring instruments through their *in situ* calibration.

In particular, we aim at 1) facilitating the identification of existing *in situ* calibration strategies applicable to a sensor network depending on its properties and the characteristics of its instruments; 2) helping to choose the most suitable algorithm depending on the sensor network and its context of deployment; 3) improving the efficiency of *in situ* calibration strategies through the diagnosis of instruments that have drifted in a sensor network.

Three main contributions are made in this work. First,

a unified terminology is proposed to classify the existing works on *in situ* calibration. The review carried out based on this taxonomy showed there are numerous contributions on the subject, covering a wide variety of cases. Nevertheless, the classification of the existing works in terms of performances was difficult as there is no reference case study for the evaluation of these algorithms.

Therefore in a second step, a framework for the simulation of sensors networks is introduced. It is aimed at evaluating *in situ* calibration algorithms. A detailed case study is provided across the evaluation of *in situ* calibration algorithms for blind static sensor networks. An analysis of the influence of the parameters and of the metrics used to derive the results is also carried out. As the results are case specific, and as most of the algorithms recalibrate instruments without evaluating first if they actually need it, an identification tool enabling to determine the instruments that are actually faulty in terms of drift would be valuable.

Consequently, the third contribution of this thesis is a diagnosis algorithm targeting drift faults in sensor networks without making any assumption on the kind of sensor network at stake. Based on the concept of rendez-vous, the algorithm allows to identify faulty instruments as long as one instrument at least can be assumed as non-faulty in the sensor network. Across the investigation of the results of a case study, we propose several means to reduce false results and guidelines to adjust the parameters of the algorithm. Finally, we show that the proposed diagnosis approach, combined with a simple calibration technique, enables to improve the quality of the measurement results. Thus, the diagnosis algorithm opens new perspectives on *in situ* calibration.

