

Knowledge-inspired fusion strategies for the inference of PM2.5 values with a Neural Network

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Abstract. Ground-level concentrations of Particulate Matter (more precisely PM2.5) are a strong indicator of air quality, which is now widely recognized to impact human health. Accurately inferring or predicting PM2.5 concentrations is therefore an important step for health hazard monitoring and the implementation of air quality related policies. Various methods have been used to achieve this objective, and Neural Networks are one of the most recent and popular solutions.

- 5 In this study, a limited set of quantities that are known to impact the relation between column AOD and surface PM2.5 concentrations are used as input of several networks architectures to investigate how different fusion strategies can impact and help explain predicted PM2.5 concentrations. Different models are trained on two different sets of simulated data, namely global scale atmospheric composition reanalysis provided by the Copernicus Atmospheric Monitoring Service (CAMS) as well as higher resolution data simulated over Europe with the Centre National de Recherches Météorologiques ALADIN model.
- 10 Based on an extensive set of experiments, this work proposes several models of knowledge-inspired Neural Networks, achieving interesting results both from the performance and interpretability points of view. Specifically, novel architectures based on BC-GANs (which are able to leverage information from sparse ground observation networks) and on more traditional UNets, employing various information fusion methods, are designed and evaluated against each other. Our results can serve as baseline benchmark for other studies and be used to develop further optimised models for the inference of PM2.5 concentrations
- 15 from AOD at either global or regional scale.

1 Introduction

Particulate Matter (PM2.5), defined as fine airborne particles with an aerodynamic diameter of less than 2.5 micrometers, serves as a critical indicator of air quality. PM2.5 levels are strongly associated with adverse health outcomes, including respiratory and cardiovascular diseases (Bose et al., 2015; Madrigano Jaime et al., 2013; Neophytou et al., 2014). The Global Burden of

20 Disease study has recognized air pollution as the fifth leading risk factor for mortality worldwide (Cohen et al., 2017). Accurate estimation and prediction of PM_{2.5} concentrations are therefore essential for effective health hazard monitoring.

Research on the health effects of PM_{2.5} is fundamental for the development of air pollution management strategies. Access to air pollution exposure data is also critical for assessing the negative health impacts of ambient PM_{2.5}. Historically, regional and national ground monitoring networks have been the primary sources for PM_{2.5} data. However, the establishment and maintenance of such networks are costly, especially on a large scale, and may not be prioritized in some countries. Martin et al. (2019) report how a substantial portion of the world lacks adequate PM_{2.5} monitoring, with only 10% of countries having more than three monitors per million inhabitants and 60% of countries not conducting routine PM_{2.5} monitoring. Furthermore, the scarcity of historical data impedes longitudinal health studies. For instance, China's or India's PM_{2.5} nationwide monitoring networks were only established respectively in late 2012 and 2015, resulting in a lack of data prior to those dates (Ma et al., 2019; Dey et al., 2020).

30 Networks of ground-based sensors monitoring PM_{2.5} concentration at surface level are instrumental but can only provide information for a few sparse locations. Obtaining complete maps of PM_{2.5} values from satellite observations is therefore an interesting and important task. Aerosol Optical Depth (AOD), a metric used to indicate aerosol loading in the vertical column, has strong positive relationships with ground-level PM_{2.5} concentrations (Jill A. Engel-Cox and Haymet, 2004; Wang and Christopher, 2003; Mukai et al., 2006; Xin et al., 2014). In recent decades, advanced space-borne sensors have provided AOD measurements with broad spatial coverage and high spatial resolution. This has enabled the use of satellite derived AOD products for large scale estimate of mass concentration at ground level through more or less complex AOD-PM_{2.5} conversion schemes and models (van Donkelaar et al., 2006; Wu et al., 2016; Hu et al., 2014; Chu et al., 2016; Di et al., 2019; Guo et al., 2021; Ma et al., 2022; Gilik et al., 2022).

40 However, while space-borne observations of aerosol properties, such as AOD and the Ångström exponent, can provide large-scale information, these quantities are not easily nor directly related to PM_{2.5} concentrations near surface level. This is because the PM_{2.5}-AOD relationship can be a multivariate function of a wide range of influencing factors. At first order, AOD and aerosols properties (fine mode fraction, hygroscopicity) are indeed skillful predictors for near-surface PM_{2.5} concentration. The literature, however, points to a wide range of parameters that may also contribute positively to PM_{2.5} statistical prediction. Meteorological variables (wind speed, height of the planetary boundary layer (HPBL), humidity, temperature, rainfall), surface conditions (albedo, normalized difference vegetation index (NDVI)), distance to the ocean, road infrastructure, population density, elevation or calendar month are regularly considered as useful influencing factors (Lary et al., 2015; Son et al., 2018; Reid et al., 2021; Su et al., 2022).

Among the numerous studies aimed at retrieving PM_{2.5} concentrations from satellite, we can generally identify three main categories of methods. The first ones are based on atmospheric chemical transport models (CTMs) and establish a scaling factor between simulated values of AOD and PM_{2.5} (Lyu et al., 2022; Xiao et al., 2022). This factor can then be transferred to estimate ground level PM_{2.5} from satellite derived AOD (van Donkelaar et al., 2006; Geng et al., 2015). This method accuracy heavily depends on the scaling factor spatiotemporal variability and has therefore clear limitations if the variability is not properly accounted for and represented by the scaling model. The second set of methods are directly data-driven and aim at

55 establishing a univariate or multivariate statistical relationship between AOD, other influencing factors and ground-level PM_{2.5}
observed concentrations. While the initial studies proposed to use simple linear or generalized linear regression models, more
complex nonlinear methods, such as neural networks (Gupta and Christopher, 2009) or boosting (Reid et al., 2015), have been
applied later. Machine learning techniques have developed rapidly (Irrgang et al., 2021; Unik et al., 2023) and proved highly ef-
60 ficient for representing the nonlinear relationships between PM_{2.5} and multiple variables (Lee et al., 2022). Yet, performances
of machine learning based methods remain eventually affected by the distribution and density of ground stations used to feed
the regression algorithms (Gupta and Christopher, 2009; Li et al., 2017). Finally, a third type of approaches combined physics
based explicit relations between core aerosol properties (size distribution, hygroscopicity, optical extinction efficiency) and
PM_{2.5} concentrations. While those also rely partly on empirical formulation for establishing some parameters (especially the
link between optical properties and aerosols composition), they tend to provide a better physical interpretability than purely
65 statistical methods and are also more independent of ground stations observations specifics. Combining the interpretability
advantage of semi-physical empirical models with the strength of machine-learning to improve the accuracy of physical pa-
rameters acquisition, opens a clear path to obtain accurate PM_{2.5} concentration from satellite observations as illustrated by
(Jin et al., 2023a).

Machine learning has been increasingly used to develop PM_{2.5} models and deep learning, in particular deep convolutional
70 neural networks (DCNN), has recently revolutionized many prediction-related application areas, including diagnostic. Several
recent and extremely thorough review papers provide clear evidence for the exploding number of studies in the field (Ma et al.,
2022; Unik et al., 2023; Zhou et al., 2024) and also illustrate the need for more standardized comparison methodologies and
metrics (Zhou et al., 2024).

While models tend to perform increasingly well, especially once optimized for a particular region (Chen et al., 2024), they do
75 not necessarily help understand the relative importance of input parameters on final decision. An old and persistent criticism of
neural networks (NN) among physicists is that they do work often at the expense of hiding physical understanding, especially as
NN based models tend to rely on increasingly complex architectures. Not surprisingly the general growing interest in so-called
"explainable AI" is also echoed in sciences (Beckh et al., 2021), including atmospheric sciences, as the use of deep learning
create paradigm shifts in atmospheric modeling. In that respect, the study by Park et al. (2020) provides a valuable approach to
80 evaluate model sensitivity to predictors through layerwise relevance propagation (LRP) (Bach et al., 2015) but remains quite an
exception among the ocean of PM_{2.5} models. Finally while ML actually provides skillful models, there has been little work in
the atmospheric sciences to understand how 2D AOD distribution could actually inform on aerosol properties and be combined
with column properties in order to improve AOD to PM_{2.5} scaling. While some essential parameters are not easily handled
or predictable (boundary layer height, aerosol type Fine Mode Fraction and aerosol vertical profiles) all depend strongly on
85 atmospheric dynamics and geographical location which in turn is somehow translated in the 2D AOD distribution. CNN have
shown excellent generalization capability for dealing with input data that has spatial auto-correlation, like images (Szegegy
et al., 2016) and are therefore potentially well suited in order to extract the information on aerosol properties contained in their
spatial distribution (Marais et al., 2020).

Among the three different approaches often used to estimate PM_{2.5} from AOD observation, we explore here an hybrid
90 method for addressing the scaling approach. We use DCNN (Deep Convolutional Neural Networks) or DC-GAN (Deep Con-
volutional Generative Adversarial Networks) in order to better capture the spatiotemporal heterogeneity of the PM_{2.5}-AOD
relationship. We aim at testing different architectures and information fusion strategies in order to develop a model for PM_{2.5}
whose results and performances can be better explained.

In previous work (Dabrowski et al., 2023), the AOD alone is used for the inference of PM_{2.5}, which leads to promising
95 results, surpassing other methods such as Polynomial Interpolation and the Random Forest Machine Learning algorithm.
However, other variables (such as the surface-level wind speed and direction, temperature, pressure, humidity and Ångström
exponent), could be used as well, as they are known to strongly drive surface PM_{2.5} concentrations (Unik et al., 2023). We
evaluate in the present study if these additional information could enhance the inference performance depending on network
architecture and information fusion strategy.

100 The main contributions of this paper are :

- a study on the interest of several variables (Ångström exponent, wind speed and direction, temperature, pressure, humid-
ity) for the prediction of surface PM_{2.5} concentration when used jointly with the AOD
- a study on the best type of fusion method to use for the prediction of surface PM_{2.5} concentration, depending on the
variables used as input and on the type of model used
- 105 – based on the knowledge from these studies, a model architecture is proposed, along with a selection of additional input
variables to use

The insights this study provides and the knowledge it represents help in building an efficient (and knowledge-inspired)
model. Indeed, based on a performance analysis, we propose a combination of network architectures that appear most suitable
for the application. We note here that our main objective is not to develop an optimized network for a specific application but
110 rather to investigate whether certain types of network architecture or fusion strategies may be more suitable for leveraging
information contained in 2D multi-component atmospheric fields for aerosol characterization.

This paper begins with an explanation of the experimental approach chosen to investigate this problem, in section 2. Then
section 3 proposes a more in-depth description of the data used. Section 4 is dedicated to describing the models and methods
proposed as solutions in this paper. A quick overview of several relevant concepts from the fields of Machine Learning, Deep
115 Learning and Computer Vision is provided in subsection 4.1, followed by a more detailed explanation of the models of interests.
Indeed, some other methods are only used as a baseline for comparison. A precise description of our experiments is realised in
section 5, along with their results and interpretation in section 6. Finally, section 7 gives an overview of the main findings and
proposed solutions deriving from this study.

2 Objective and Approach

120 The purpose of the models designed in this paper is to infer maps representing values of the PM_{2.5} concentration at ground level, from maps (of the same size) representing values of the AOD in conjunction with maps of other atmospheric variables. To do this, NN such as UNets and GANs are used as their convolutional versions showed an ability to take into account the spatial variability of the data they are being presented with. In the case of Convolutional Neural Networks (CNN) these data mainly take the form of images or matrices.

125 As further detailed in section 3, we use different aerosol optical properties in addition to AOD and meteorological quantities that are known to characterize or drive the aerosol concentration as well. These are namely, the wind speed and direction, atmospheric pressure, temperature, relative humidity (all five of these meteorological variables being measured at surface level) and the Ångström exponent.

An important number of experiments are conducted in order to study the impact of these additional variables on the inference
130 performance for each network architecture. Furthermore, as stated in section 4.3, there exists different strategies to leverage several inputs at the same time and within the same model. In this work, we test three different fusion techniques, namely: feature fusion (FF), decision fusion (DF) and channel concatenation (CC) (also called data fusion). These strategies and their implementation are described in section 4.3. Experiments are performed to identify the best fusion strategy depending on network architecture and available input variables considered for inference of PM_{2.5}.

135 More classical solutions, such as the kriging method or even polynomial interpolation, are implemented as well to serve as baseline for comparison of inference performances. The expected outcome of this important number of experiment is a performant NN architecture for the prediction of PM_{2.5} concentration from complete and incomplete maps of the AOD, along with insights on the design process of this type of model.

3 Data

140 In this work, we exclusively use data from simulations (namely from the CAMS and ALADIN models), as it allows us to easily obtain all necessary information. It also maintains the possibility to select or sample data to represent realistic observation scenarii. The CAMS model provides maps representing values of various meteorological quantities and optical measurements, covering the entire world.

The ALADIN model provides the same type of maps, but they cover Europe and the North of Africa instead of the entire
145 world, and at a higher spatial and temporal resolution than CAMS.

Even though we use simulated data, our objective remains to simulate what could be obtained in a real situation. This is why we come up with a scenario in which a part of the data is simulated, and the most recent part is real, measured data. More precisely, in this hypothetical scenario, optical sensors operating from geostationary satellites (Ceamanos et al., 2021) allow us to obtain AOD values in near real-time. These satellites are, namely, two Meteosat Second Generation (MSG) satellites,
150 the Himawari satellite, and two Geostationary Operational Environmental Satellites - New Generation (GOESNG). They respectively cover Europe and Western Asia, Eastern Asia, and the Americas. The cumulated coverage of these geostationary

satellites allows to generate complete (as opposed to sparse) maps of the AOD. The PM2.5 concentration values at surface level would be obtained through photometers, Lidar instruments, optical counting sensors or even filters. Each of these sensors can only provide concentration values for its own geographical location. This is why this network of sensors can only provide
155 sparse maps of the PM2.5 concentration.

This means that, for a real use-case scenario, no complete ground truths are available in the measured data. Instead, sparse ground truths are available. In order to reproduce that scientific obstacle, we produce sparse maps of the PM2.5 concentration using the complete ones, by randomly selecting pixels. For a part of the training set, we consider having only access to these sparse maps instead of the complete ones. This was suggested by the authors of (Dabrowski et al., 2023), as it allowed for some
160 level of control over the sparsity of these sparse maps, and therefore allowed for a study of the impact of the sparsity of these maps over the results. We use this method too in order to be able to compare the results of this paper to ours.

The Aerosol Optical Depth, expressed for a wavelength of $550nm$, is our main input and is used systematically. Apart from it, six other quantities that are either routinely observed or modelled can be used as additional inputs. Five of them are meteorological quantities : wind speed and direction, relative humidity, temperature, and pressure, all of which are measured
165 at surface level. These quantities are known to drive aerosols concentration and their size distribution. The last quantity, the Ångström Exponent (AE), is actually an optical quantity that is derived from AOD at two different wavelengths. It characterizes the spectral variation of the AOD and is related to aerosol particle size distribution such that aerosols with a dominance of fine particles will tend to exhibit larger AE. This is therefore an important parameter in aerosol modeling and potentially an important predictor of PM2.5 (Jin et al., 2023b).

170 Links to the data and code used in this article are available in the code and data availability section, just after section 7.

4 Models

4.1 Background

The task of inferring maps of PM2.5 at ground level from maps of AOD combined with one or several other variables, can be seen as a regression problem, or as what is known in the field of Computer Vision an "image-to-image translation" problem.
175 Indeed, we want to infer an output image from a different input image (or from a number of them). In the litterature, one can find several methods used to solve this kind of task such as the polynomial interpolation, the kriging method (Matheron, 1963), machine learning (Ho, 2018) and deep learning algorithms (Goodfellow et al., 2020). The most relevant algorithms in our context will be briefly described below. Details can be found in section A.

Kriging (Matheron, 1963) is a spatial interpolation and extrapolation method governed by prior covariances. It performs
180 better with important volumes of data, and when estimated values follow a normal distribution. For each inference, a new kriging model is built, which leads to longer inference times compared to other methods used in this paper. This method is described in greater detail in section A.

UNets (Ronneberger et al., 2015) represent a type of Neural Networks (NN) architecture, known for its performances, particularly in the field of Computer Vision. The architecture is typically composed of an Encoder and a Decoder, and the main

185 idea behind UNets is to add skip connections linking the outputs of each layer in the encoder to a corresponding layer in the decoder. This makes it possible to reduce dimensionality without the risk of losing relevant information in the process. Figure 1 gives an example of a model with this type of architecture.

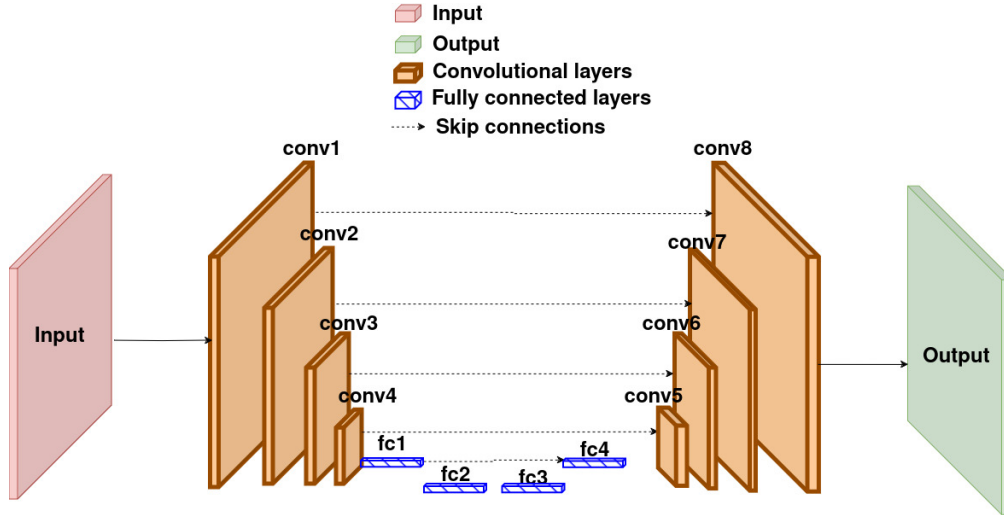


Figure 1. Architecture of a UNet with both Convolutional and Feed-Forward layers. Here, the Encoder and Decoder are symmetrical.

GANs (Goodfellow et al., 2020) are a type of NN that actually consist of two networks. One is called the generator, and the other is the discriminator. The role of the discriminator is to distinguish between real data and data generated by the generator. The purpose of the generator is to produce output close enough to the real data that the discriminator labels them as real. These two networks learn competitively: the higher the loss of the generator, the lower the loss of the discriminator, and inversely. This allows such models to show interesting performances in the context of semi-supervised learning. They are also highly efficient for image-to-image translation tasks. The learning process of the type of GAN that will be used in this paper is described in (Dabrowski et al., 2023), in figure 1 of that article.

195 4.2 Relevant Deep Learning models and architectures

Different types of model are implemented, trained and tested on our data. We use as baseline for comparison the Random Forest algorithm (a Machine Learning algorithm), a polynomial interpolation method (of degree 3), and a kriging algorithm as described in 4.1.

Two Deep convolutional neural networks architectures, UNets and BC-GANs (Dabrowski et al., 2023) are the basic components of the models we propose in this paper. The first one is a purely supervised UNet, and the second one a semi-supervised BC-GAN as described in 2. The architecture of the generator of these BC-GANs will each time correspond to the architecture of the corresponding purely supervised UNet. As for the architecture of the discriminators, they are described in section B.

These models allow to leverage sparse ground truth when complete ones are unavailable, which increases performance (compared to classical GANs) in the context of semi-supervised learning.

205 Indeed, we do not have access to complete ground truths for the whole training set. For a part of it, we only have access to sparse ground truths. The authors of (Dabrowski et al., 2023) were mainly interested about how these sparse ground truths and the information they represent could be leveraged in order to ease the training and obtain better results. They proposed a method to leverage those sparse ground truths based on the literature around Physics-Informed Networks, that implied seeing those sparse ground truths as Boundary Conditions (BC), hence the name of BC-informed GAN. The authors of (Dabrowski
210 et al., 2023) illustrate this method in figure 2 of their article. It includes the design of an additional loss function in order to train the model to respect the BCs. This essentially allows for localised supervision.

4.3 Information fusion strategies

All our models have in common that they use as input one or several variables to produce the same type of output. It is therefore necessary to merge these variables together during this process. This also ensures the production of the output makes use of all
215 these pieces of information.

Fusion strategies are therefore an important aspect of the architecture definition, and eventually the performance, of the model. They represent different methods that can be applied to leverage several sources of data (several inputs) within the same NN. They can be applied on models such as UNets as well as GANs. There are three main different fusion strategies according to (Mangai et al., 2010), namely data fusion, feature fusion and decision fusion. Sections 4.3.1, 4.3.2 and 4.3.3
220 describe (respectively) each of these approaches and the way we use them.

4.3.1 Data fusion (channel concatenation)

The general idea behind this strategy is to use several pieces of raw data to build a new, more complete and useful, piece of raw data.

As our data consists of images, the simplest way to do data fusion is to realise channel concatenation: in other words, to use
225 our different inputs as if they were different channels of one single image. For this reason, this method is also called Channel Concatenation throughout this paper. This approach is the most straightforward of the three.

In terms of architecture of our neural networks, this simply implies using more convolution filters. The UNet’s architecture with data fusion is illustrated by Figure 2. The GAN’s discriminator’s architecture with data fusion is illustrated by Figure B1 in section B.

230 In terms of interpretation, this architecture relies on the local (rather than global) relationships between the different quantities used as input. We believe this architecture to work better if local patterns in one input image correspond to local patterns in other input images.

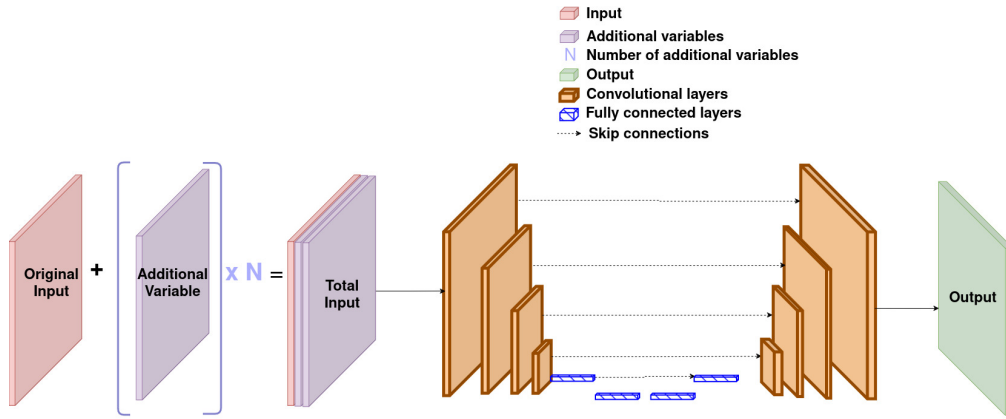


Figure 2. Architecture of the UNet with data fusion approach. Also corresponds to the architecture of the GAN’s generator.

4.3.2 Feature fusion

One of the most common architectures in neural networks for computer vision tasks is the encoder-decoder. The general idea is that the encoder generates what is called a feature. This feature is usually a vector, but can technically also be a matrix (although the idea is for it to be of smaller dimensions than the input). It is supposed to contain all the relevant information from the input with regard to the task at hand. In other words, it describes the input well enough so that only this feature is needed for the task. During the next step, the decoder uses the feature as input and produces the output.

Feature fusion methods are often applied to computer vision tasks, for example when dealing with complex hyperspectral images (Song et al., 2018). The idea behind them is to obtain a feature for each input, and use them to obtain one super-feature (for example by simply concatenating the various features). According to the authors of (Sun et al., 2005), two interesting ways to fuse feature vectors are the serial feature fusion (based on an union vector) and the parallel feature fusion (based on a complex vector), although the same authors actually propose a new method based on canonical correlation analysis (CCA).

This unique feature is then used by the decoder to produce the output. In terms of architecture, this implies having as many encoders as inputs, but only one decoder (since there’s only one output). The UNet’s architecture with feature fusion is illustrated by Figure 3.

The UNet architecture is a specific type of encoder-decoder, in which a specific type of connections, called skip connections, can be found. It is also often symmetric (in the sense the decoder’s layers mirror the encoder’s one). After each layer in the encoder, the obtained feature is sent to the corresponding layer in the decoder. This allows for the decoder to have access to several features rather than simply the smallest one.

Implementing feature fusion with a UNet is therefore non-trivial: among all the features obtained for each input, which ones should be sent to the decoder through skip connections ? We choose to apply what we call multiple feature fusion. The principles of feature fusion are applied to feature of each and every scale, and those merged features are sent to the corresponding layer of the decoder through skip connections.

255 In terms of interpretation, this architecture relies on the global relationships between the inputs. As obtaining features relies on dimension reduction, those features represent the input in a more global way, and do not necessarily represent local patterns. The smaller the scale of these features (and the deeper their corresponding layers are), the truer it is. This architecture relies on the idea that each of the input images contains a global, non-localized piece of information that can be useful in order to estimate the aerosol concentration. Again, the global or non-localized aspect of each of the features actually depends on its
260 scale or depth in the network.

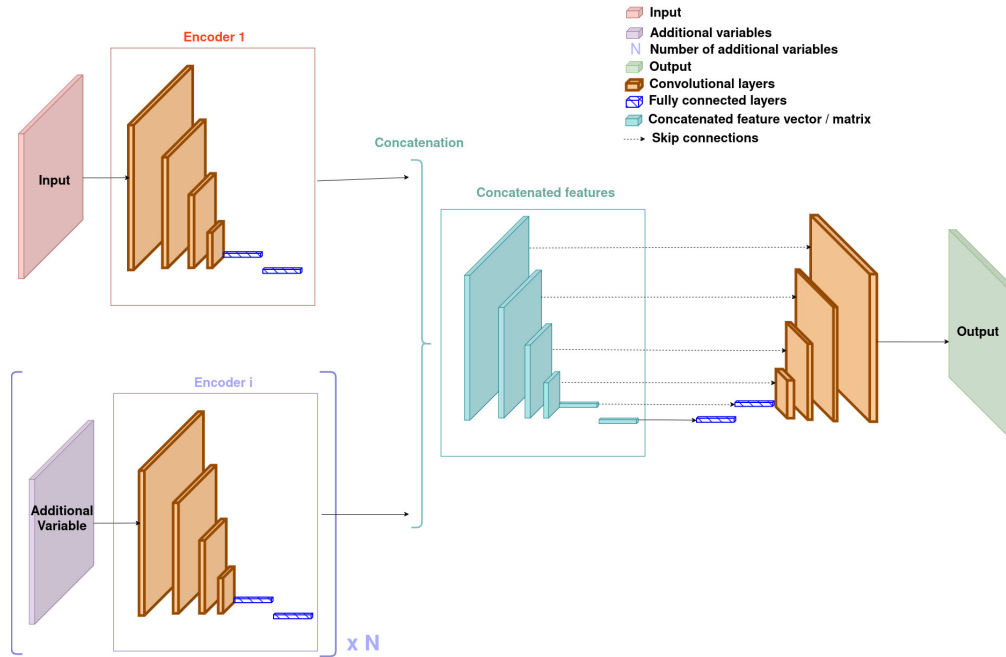


Figure 3. Architecture of the UNet with feature fusion approach. Also corresponds to the architecture of the GAN’s generator.

The GAN’s discriminator’s architecture with feature fusion is illustrated by Figure B2 in section B.

4.3.3 Decision fusion

The idea behind decision fusion is to use a separate model for each input, obtain an output for each of them, and then merge those outputs together to obtain a final decision, supposedly better. In classification tasks, decisions represent the predicted
265 class. In regression tasks, like the one considered in this paper, they represent the estimated quantity. Losses are computed using the final output. The backpropagation process takes place through the entire model (and the smaller models that compose it).

There exists several ways to fuse decisions, such as the linear or log opinion pool (corresponding respectively to a weighted sum or product) (Sinha et al., 2008). Voting algorithms can even be used for classification tasks (Sinha et al., 2008). In this

270 article, a linear opinion pool approach is chosen: we apply a weighted mean of all the outputs to obtain the final one. The weights are learnable parameter, which allows the model to learn which outputs are most relevant.

This principle relies on the idea that each of the inputs can individually be used to produce an estimation of the aerosol concentration, but that these estimation may be flawed, and the best estimation can be obtained through a combination (here a weighted mean) of these flawed estimation. In other words, it is possible to correct the estimation produced with an input using the estimation produced with an other input. Since the model can learn which inputs are the most relevant to produce the desired output, we expect this approach to provide the best inference when all inputs are used.

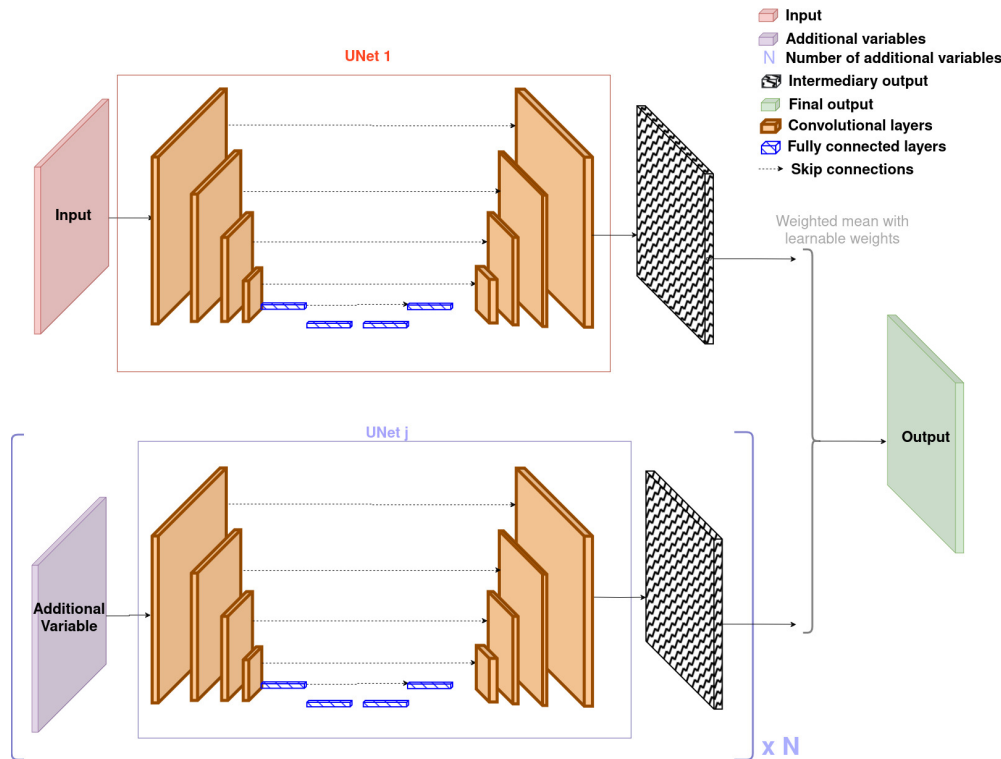


Figure 4. Architecture of the UNet with decision fusion approach. Also corresponds to the architecture of the GAN's generator.

The architecture of the GAN's discriminator with decision fusion is, in principle, very similar to the architecture of our UNet with decision fusion. The main difference is the type of output, as the discriminator outputs a single scalar for each iteration while the UNet outputs images. It is illustrated by Figure B3 in section B.

280 4.3.4 Hybrid fusion models

The physical nature of PM2.5 predictors obviously has an impact on the non-linear function linking AOD and PM2.5. While AOD is directly linked to total column aerosol concentration at a given location, surface pressure can indirectly be linked to PM2.5 through accumulation in the atmospheric boundary layer under stable conditions while wind speed can influence

PM2.5 concentration over longer range in space and time. Therefore we can distinguish "state" variables that can directly link PM2.5 to AOD through an integral expression over the atmospheric column and "indirect predictors" that act on PM2.5 concentrations over different space and time scales. In our current analysis, the Wind variables (speed and direction) stand out as they describe the atmosphere dynamics while the AOD and Ångström exponent are clearly states variables regarding the inference of PM2.5 concentration. Humidity, Pressure and Temperature variables, can be considered primarily as states variables as they strongly impact the particle size distribution through aerosol hygroscopicity but can also indirectly influence near surface PM2.5 concentrations by favoring accumulation under stable atmospheric conditions or on the contrary by removal of atmospheric particles through dry deposition or wet scavenging.

The performance of networks and their robustness to noise is known to be impacted by their architecture and network performances can be improved upon SOTA when the network is well-aligned with the target function (Li et al., 2021). We hypothesize here that for atmospheric applications, the optimal alignment of network architecture with the target function may depend on the nature of variables used as input and on the fusion strategy used for merging information carried by those variables. Through this hypothesis we ask whether there is an advantage in applying different fusion strategies for different types of input variables.

Based on this insight, we propose two hybrid models, using different fusion strategies depending on the variable considered. The idea is to use Data Fusion (Channel Concatenation) for the AOD and all other input variables, except for the Wind and Temperature for which Feature Fusion is used. Figures 5 and 6 describe these two models more precisely. Section 6.5 shows

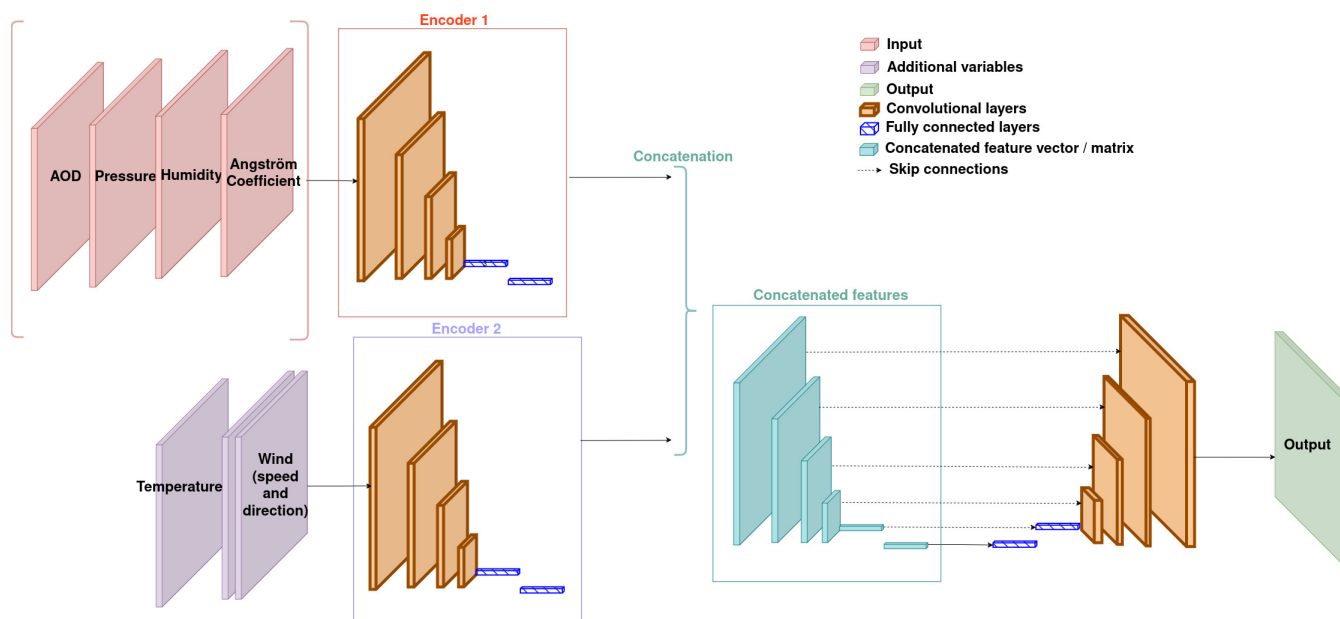


Figure 5. The first proposed hybrid model, using both data fusion and feature fusion.

and discusses the results obtained by these models.

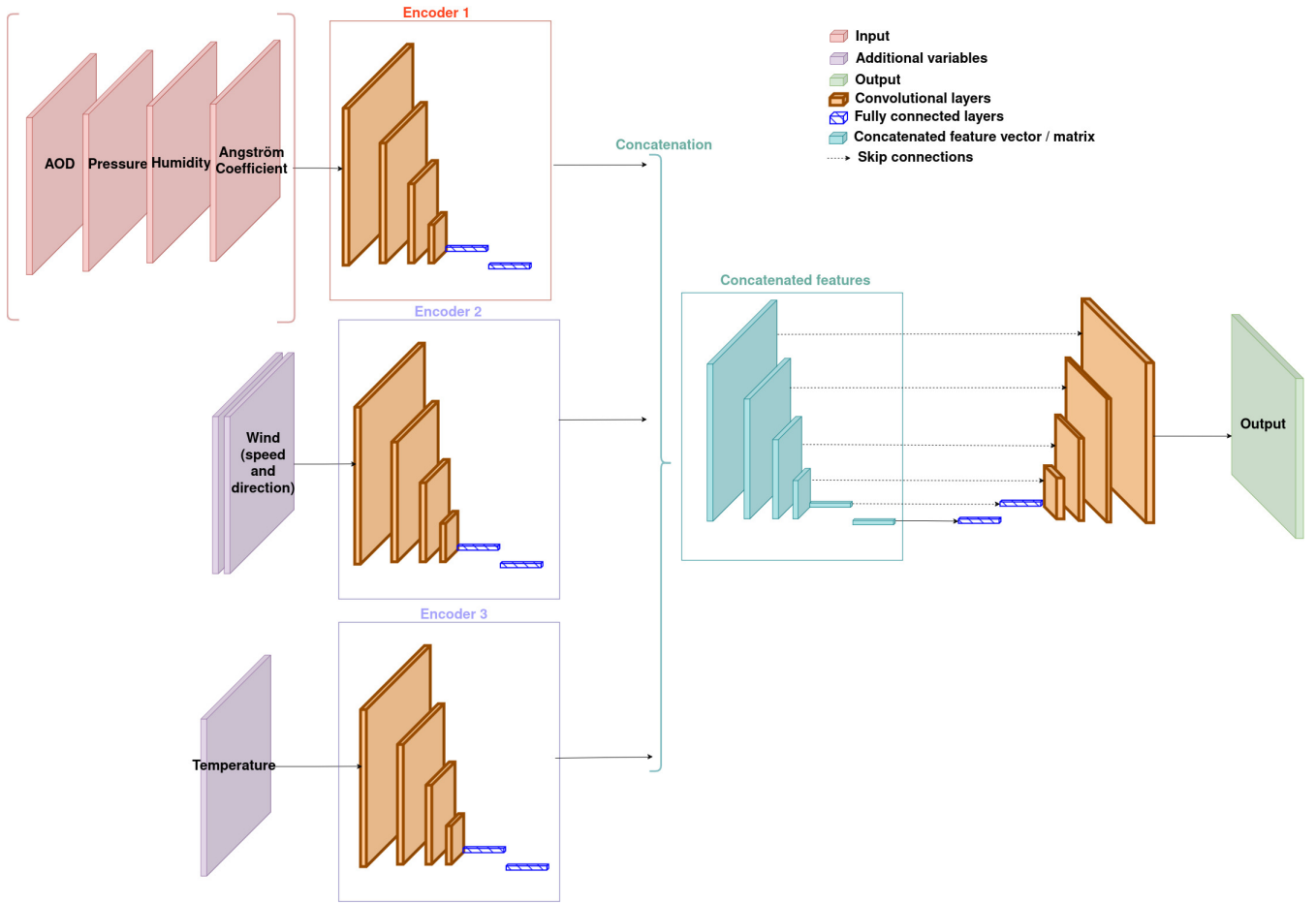


Figure 6. The second proposed hybrid model, using both data fusion and feature fusion.

5 Methodology

In order to find the best way to handle our variety of input quantities, we propose to study the three main fusion approaches described in section 4.3. Each fusion method is experimented on two types of models: a purely-supervised UNet, and a BC-
 305 GAN using sparse measurements of the aerosol concentration at ground level as boundary conditions. Experiments on the hybrid approaches proposed in section 4.3 are realised as well.

The goal is also to understand which input quantities have the most important impact on our results, in other words which additional variables actually help our models to better predict PM2.5 at surface level. This is why, for each distinct model architecture, experiments are realised with different combinations of additional variable, to study their impact on the results.

310 5.1 Learning and validation protocol

Experiments are conducted on CAMS and ALADIN datasets. AOD is systematically used as input in all our experiments. In addition, six other variables are also considered (wind speed and direction, relative humidity, atmospheric pressure, temperature, Ångström exponent) in order to evaluate their impact on the results. It is important to specify that the wind speed and direction are always used jointly to describe the wind state variable. With this rule in mind, experiments are performed with all possible combinations of these six variables (including using none of them and all of them). The AOD is also used in all cases.

For both CAMS and ALADIN dataset, we always consider the same type of scenario, shown by Figure 7. In this scenario, we have access to a dataset with complete ground truths, corresponding to a period of eleven months. We also have access to a second dataset with sparse ground truths, which can therefore only be used in the context of semi-supervised (as opposed to purely supervised) learning, corresponding to a period of one month. These sparse ground truths will be used as Boundary Conditions as stated in section 2. They contain an amount of pixels corresponding to 5% of the pixels available in complete ground truths.

For the CAMS dataset, one sample is generated every three hours. Samples take the form of matrices of size 241x480. Depending on the chosen number of input modalities, each model input can be composed of one to seven of these matrices. The training set therefore contains 2680 of these inputs, the sparse training set (with sparse ground truths) 240 inputs, and the test set 2920 inputs.

For the ALADIN datasets, one sample is generated every hour. The size of the matrices is 405x613. Again, depending on the chosen number of input variables, each model input can be composed of one to seven of these matrices. The training set therefore contains 8040 of these inputs, the sparse training set (with sparse ground truths) 720 inputs. For the test set, we only use one image for every three hours, so that it contains as many samples as the CAMS test set. Therefore it also contains 2920 inputs.

An exhaustive study on our two models (UNets and GANs), three fusion strategies and six input variables (and all possible combinations of these) is conducted on the CAMS dataset. This corresponds to 192 different experiments. Only experiments that lead to the best performances on the CAMS dataset are conducted on the ALADIN dataset. The aim is to provides insight on the impact of the characteristics of each dataset on the results, as shown in section 6.2.

Then, experiments are realised on the hybrid approaches, but only on UNet models, and always with all six additional input variables. These experiments are realised on both datasets as well.

5.2 Data pre-processing

All values of AOD inferior to 0.005 can be considered as noise. They are therefore set to 0 before being used, be it for training or prediction.

In order to speed up the convergence of the models, we equalise the PM2.5 and AOD distributions by applying the function $\ln(1+x)$ to those values. The inverse function $\exp(x) - 1$ is then simply applied to inferred outputs in order to obtain actual concentration values (in $\mu\text{g}/\text{m}^3$ and ease the interpretation of our results.

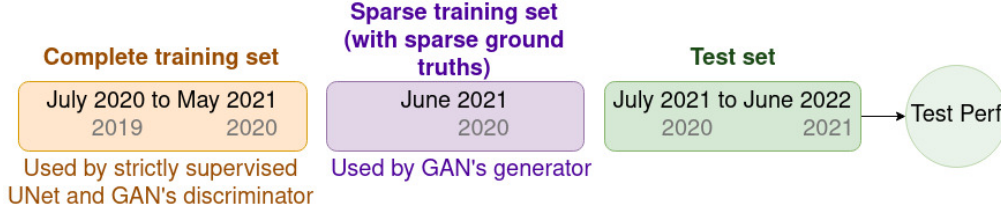


Figure 7. Representation of the datasets used for our experiments. The text in gray specifies the corresponding periods of time for the datasets built with the data from the ALADIN model instead of the CAMS model.

In our context, polluted regions (with high aerosol concentration values) are our areas of main interest. Therefore filtering very low aerosol concentration values in our ground truths and predictions allows us to better evaluate the model performance in these areas. Specifically, values inferior to $1\mu g/m^3$ are set to $0\mu g/m^3$. This is done before computation of the evaluation metrics.

These previous pre-processing protocols are based on the protocols proposed by (Dabrowski et al., 2023).

The data gives us access to values of both Eastward and Northward wind speeds. Instead of using them as such, we apply the transformation described by equation 1 to instead obtain two different matrices. In this equation, U and V respectively represent the Eastward and Northward wind speeds. The first one contains values of the wind speed norm (regardless of direction), and the second one values of the direction (in degrees) of the wind. Each time wind values are used in an experiment, both of these matrices are used.

$$\begin{aligned}
 norm &= \sqrt{U^2 + V^2} \\
 direction &= \arctan\left(\frac{V}{U}\right) \cdot \left(\frac{180}{\pi}\right)
 \end{aligned} \tag{1}$$

Our data does not originally contain values of the Angström exponent, but it is easily possible to compute them using values from the AOD measured at two different wavelengths, and the equation 2. In this equation, λ_1 represents the wavelength of the original AOD, which is $550nm$, and λ_2 is the wavelength of the second AOD, used simply to compute the Ångström exponent. When using data from the CAMS model, this second wavelength is of $865nm$, while with the ALADIN model, it is of $1000nm$.

$$angstrom(\lambda_1, \lambda_2) = -\frac{\ln\left(\frac{AOD(\lambda_2)}{AOD(\lambda_1)}\right)}{\ln\left(\frac{\lambda_2}{\lambda_1}\right)} \tag{2}$$

Pressure values are converted from Pascal to atmospheres, and temperature values from Kelvin to Celsius degrees.

5.3 Metrics and Losses

During training, depending on the type of model, different losses can be used. For our GANs, the adverse loss is used (to train both the generator and the discriminator), as well as the Boundary Condition loss (which is essentially a localised MSE

as stated in section 2). This last loss is described by equation (3), where y is the ground truth, $\hat{y}$ represents the output of the
 365 model, and BC is a matrix containing ones at the location of known values of y and 0 elsewhere.

$$BCloss(BC, \hat{y}, y) = MSE(BC \cdot \hat{y}, BC \cdot y) \quad (3)$$

For our supervised UNets, we use the MSE loss function as well as the FSIM (Zhang et al., 2011), which is also used as a metric and described in this section.

Four metrics are used to evaluate the models during testing: the Mean Absolute Error (MAE), the Mean Bias Error (MBE),
 370 and the Quantized Error (QE) as proposed by (Dabrowski et al., 2023), as well as a metric proposed by (Zhang et al., 2011) called the FSIM. The MAE and MBE are expressed in $\mu g/m^3$ (the unit of aerosol concentration) and in % (for their relative versions).

In the following equations, y represents the ground truth, y_i its elements, and $\bar{y}$ its average. $\hat{y}$ is the output of the model and $\hat{y}_i$ its elements. The number of elements in either matrix is N .

375 – Mean Absolute Error (MAE)

It is the most widely used of these metrics, it represents the error of the models in a general sense (Wang and Lu, 2018). It is described by equation (4) in which y is the ground truth, $\hat{y}$ represents the output of the model, and N represents the number of pixels in y .

$$MAE(\hat{y}, y) = \left| \frac{y - \hat{y}}{N} \right| \quad (4)$$

380 Equation (5) describes the way the $rMAE$ is computed from the MAE , using $\bar{y}$ the average of the ground truth y .

$$rMAE(\hat{y}, y) = \frac{MAE(\hat{y}, y)}{\bar{y}} \quad (5)$$

– Mean Bias Error (MBE)

It represents the model's tendency to overestimate or underestimate the values to predict. A model with poor performance can however still have a low bias, as it is not the only aspect of its performance. This error can sometimes be used to
 385 correct the bias of the model. It is described by equation (6) which uses the same notation as equation (4).

$$MBE(\hat{y}, y) = \frac{y - \hat{y}}{N} \quad (6)$$

Equation (7) describes the way the $rMBE$ is computed from the MBE , using $\bar{y}$ the average of the ground truth y .

$$rMBE(\hat{y}, y) = \frac{MBE(\hat{y}, y)}{\bar{y}} \quad (7)$$

– Quantized Error (QE)

390 This metric is used to quantize the prediction as well as the ground truth before comparing them. The quartiles of the ground truth values distribution ($q1$, $q2$, $q3$, $q4$) are used to define four classes for this quantization process, which is

described by equation (8). In this same equation, $M_{i,j}$ represents a pixel of coordinates (i, j) from the unquantized matrix M . $C_{i,j}$ represents the pixel with these same coordinates in the corresponding quantized matrix C .

$$C_{i,j} = \begin{cases} 0 & \text{if } M_{i,j} \leq q1 \\ 1 & \text{if } M_{i,j} \in]q1, q2] \\ 2 & \text{if } M_{i,j} \in]q2, q3] \\ 3 & \text{if } M_{i,j} > q3 \end{cases} \quad (8)$$

395 This process allows us to obtain a quantized ground truth C_{gt} and a quantized prediction C_{pred} . The Quantized Error is computed from these matrices following the equation (9). N represents the number of pixels in matrix C_{gt} in this equation.

$$QE = \frac{|C_{gt} - C_{pred}|}{N} \quad (9)$$

400 This type of metric is usually better suited for classification or segmentation tasks. However, the air quality is often represented as indexed values, and involves thresholds corresponding to different levels of health hazards (or policy alerts): this metric is therefore more closely related to this representation. It is also more sensitive than the MAE to very localised errors. It is on the other hand less suited to represent tendencies to generally overestimate or underestimate the values to predict than the MBE.

– Feature Similarity Index (FSIM)

405 This metric has been proposed by the authors of (Zhang et al., 2011), as an Image Quality Assessment metric. It relies on the concepts of Phase Congruency and Gradient Magnitude (in the sense of Image Gradient). The Phase Congruency is also used to weigh the contribution of each pixel to the similarity of two images. This leads to a significant weight being given to edges, shapes and other structures in the images.

5.4 Scores

410 As a great number of experiments were realised during this work, we can not easily compare them all in a table. We decide to present an overview of these results through charts, and to select a few experiments to compare in a table. Since a lot of different metrics are used, a protocol is needed to select the experiences to compare. Three different scores, computed from the previously described metrics, are used. This allows us to select experiments that lead to good overall results, rather than experiments that performed well on one metric and poorly on all others, for example. These scores are exclusively used for
415 this selection process and do not intervene further when interpreting and analyzing our results. For this reason, and to avoid overloading our results tables, the score values are not presented in these same results tables.

– Total score

This score is computed using all of our metrics at the same time, including the inference time. It follows equation (10). $rMAE$ and $rMBE$ represent the relative counterparts of MAE and MBE , and are expressed in % instead of $\mu g/m^3$.
 Regarding the inference time, the threshold of $0.05s$ is used as it is the maximum inference time among our experiences with our deep learning models.

$$Total\ score = \frac{\frac{0.05s - Inference\ time}{0.05s} + (1 - rMAE) + \frac{3 - QE}{3} + (1 - rMBE) + (1 - FSIM)}{5} \quad (10)$$

– Timeless score

This score is very similar to the first one, but does not include the inference time metric. This allows to identify the best performing models for a situation in which the inference time is not a predominant factor. It follows equation (11).

$$Timeless\ score = \frac{(1 - rMAE) + \frac{3 - QE}{3} + (1 - rMBE) + (1 - FSIM)}{4} \quad (11)$$

– Reduced score

This score is computed using only $rMAE$ and $FSIM$. These two metrics are the most relevant ones in the Computer Vision domain. This score therefore allows for a comparison of the models and experiments from this point of view only. It also does not make use of the inference time. It follows equation (12).

$$Reduced\ score = \frac{(1 - rMAE) + (1 - FSIM)}{2} \quad (12)$$

5.5 Methodology and data summary

In order to ease the comparison of this work to other models using a scaling approach for inference of $PM_{2.5}$ from AOD, we provide hereafter in Table 1 a summary of our methodology, as well as some characteristics of the dataset and methods used in this paper. It follows the standard proposed by the authors of (Zhou et al., 2024).

Table 1. Characteristics of the dataset, method and experiments used in this paper, in the standard proposed by the authors of (Zhou et al., 2024).

Standard	Indicator	Description
Dataset	Open source	Yes. For more info, see Code and data availability section, just after section 7.
	Data feature	
	Predict step	Single step.
	Time resolution	Every 3 hours for CAMS, every hour for ALADIN.
	Data size	CAMS : 5840 samples. ALADIN : 11 680 samples.
	Data dimensions	Up to eight matrices per sample. Shape for CAMS : 241x480. For ALADIN : 405x613.
	Dataset split	Training set and test set both span over a year. Test set always contains 2920 samples out of the total.
	Pre-processing	
	Missing value	Handled by CAMS and ALADIN models.
	Conversions	Pressure : Pa to atm. Temperature : K to °C.
Method	Filtering	AOD with a threshold of 0.005.
	Normalizing	Function $\ln(1 + x)$ applied to AOD and PM2.5
	Others	Computing the Ångström Exponent out of AOD values.
		Extracting the wind speed norm and direction from its northward and eastward speeds.
	Open source	Yes. For more info, see Code and data availability section, just after section 7.
	Architecture	UNet, inspired from (Ronneberger et al., 2015). GAN, inspired from (Goodfellow et al., 2020). Implementation of Data Fusion, Feature Fusion and Decision Fusion methods.
	Training process	Optimizer : Adam. Loss functions : MSE and FSIM (UNet), adversarial and based on BC (GAN). 500 epochs.
	Visual analysis	Available in figures 10 and 12.
	Novelty	Implementation of a Hybrid Fusion method. Study of the impact of several meteorological variables on the results.
Experiments	Experimental setting	
	Model config	Encoder kernel sizes : 9, 7, 7, 3. Decoder kernel sizes : 3, 7, 7, 9. Size of latent vector : 128.
	Computation setup	GPU Nvidia Tesla A100 with 80 Go of V-RAM.
	Results metrics	MAE, MBE, FSIM and QE (non-classical, see section 5.3).
	Params	Depending on models, between 10 and 250 millions of trainable paramters. See figure C1.
	Comparison with SOTAs	Model outperforms the kriging method, Polynomial Regression of Degree 3, and Random Forest algorithms.
	Ablation study	Yes. Several models with several information fusion methods are tested. Experiments are realised with different numbers of meteorological variables as input.

6 Results

We start with a general overview of our results. The next step is to identify the best performing and most interesting results and models among the experiments realised on the CAMS dataset. This allows us to reproduce these experiments on the ALADIN dataset, in order to compare the results and better understand the impact of the characteristics of these datasets (namely, spatial domain and resolution) on the results.

6.1 Overview

We choose to summarize our results in the form of radar charts, with five metrics represented on these charts: the inference time t , $rMAE$, QE , $rMBE$ and $FSIM$. The values of these metrics were normalized the same way they are when used to compute our scores, which allows us to represent them on the same scale. High values of these normalized versions of our metrics represent high performance. Lines made of cyan dots represent both the maximum and minimum performances for each metric, and are linked by a light cyan area that gives an overview of the performance range on each radar graph. Blue dashes and purple dashdot represent respectively the median and average performance for each metric among all results presented on this radar graph. Out of these same results, the experiments that lead to the obtention of the best total score are represented with a plain red line on the graph. The relative values (in %) of each metric for these experiments (the ones leading to the best score) are also systematically represented in yellow boxes on the graph.

Figure 8 gives an overview of the performance of each couple model-fusion strategy. It shows that, on average, models using Decision Fusion have the longest inference time, while models using Data Fusion are the fastest and those using Feature Fusion are in the middle. This is expected, as it correlates with the number of parameters of each model, as illustrated by figure C1.

It also shows that GANs seem to generally suffer from poorer $rMAE$ and $rMBE$ scores than UNets. Our interpretation is that the proportion of the training set reserved for strictly supervised training is important enough for purely supervised methods to perform well. The interest of GANs lies in their ability to realise semi-supervised training. In our case, it also corresponds to their ability to make use of the portion of our dataset that only contains sparse ground truths. This portion is small, which therefore makes the interest of GANs (and arguably semi-supervised methods) limited in this case. In comparison, the authors of (Dabrowski et al., 2023) show the efficiency of their GANs in a context where only half of the dataset contains complete ground truths. It is interesting to note that the $FSIM$ and QE metrics do not seem to be affected by this in the same way, or at least not as intensely.

Finally, the Data Fusion or Channel Concatenation (CC on the figure) strategy seems to be leading to more stable results than the other two fusion strategies. This may also be linked to the difference in model complexity, as shown in section C.

Figure 9 shows the evolution of the performances of our UNet when we increase the number of input variables. It shows that, when increasing the number of input variables, the average inference time increases. This is expected, as figure C1 shows that the models grow in complexity with the number of input variables. Other metrics, and especially $rMAE$, show on average an increase in performance when adding more input variables. However, this increase in performance is not linear, and when deciding to add an input variable to a given experiment, we are not guaranteed to obtain better results. We can also note that,

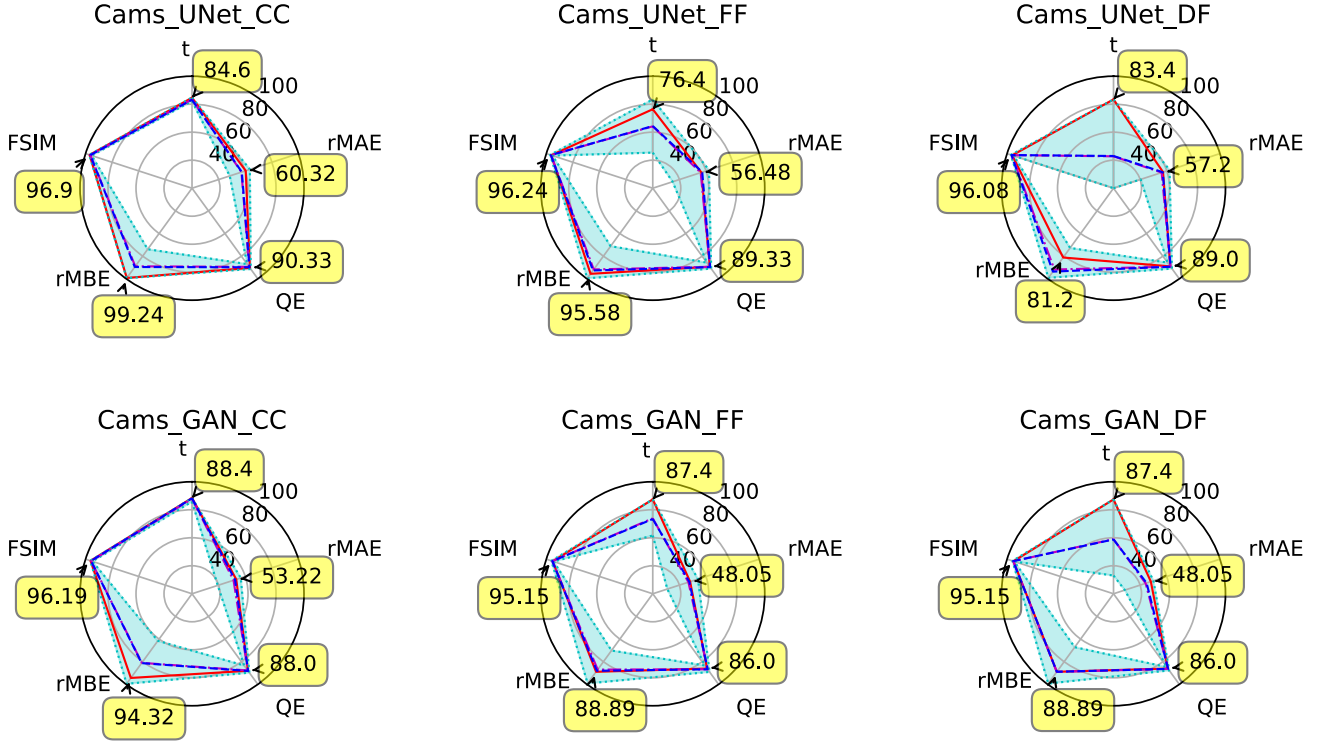


Figure 8. Overview of each model - fusion strategy couple. Regarding the fusion type, 'CC' means 'Channel Concatenation' or Data Fusion, 'FF' means Feature Fusion, and 'DF' means Decision Fusion. Cyan dots: max and min, with a cyan area in between. Blue dashes: median. Purple dashdots: average. Red plain line: best total score (corresponds to annotated values).

when comparing experiments with two additional input variables to three, we observe less stable $rMBE$ values, even though the number of experiments for these two categories is the same. Finally, these charts also show that, regardless of the fusion method used, experiments realised using all five input variables tend to produce the best results (apart from the point of view of the inference time).

6.2 Best results on the CAMS dataset

Table 2 shows the best results according to each of the three scores. First line is the result with the best Total Score, second line is the best Timeless Score, and third line is the best Reduced Score. Figure 10 shows the output of these models for one given sample.

This table shows that using more variables as input seems to generally lead to the best results, except for the Ångström Exponent on the last line. It also shows that Decision Fusion methods suffer from larger inference times, especially compared to Data Fusion. These results lead to three main recommendations depending on the context and the desired performances. If the MBE (or bias) of the output is not an important factor, then the recommended model is a UNet using the Data Fusion

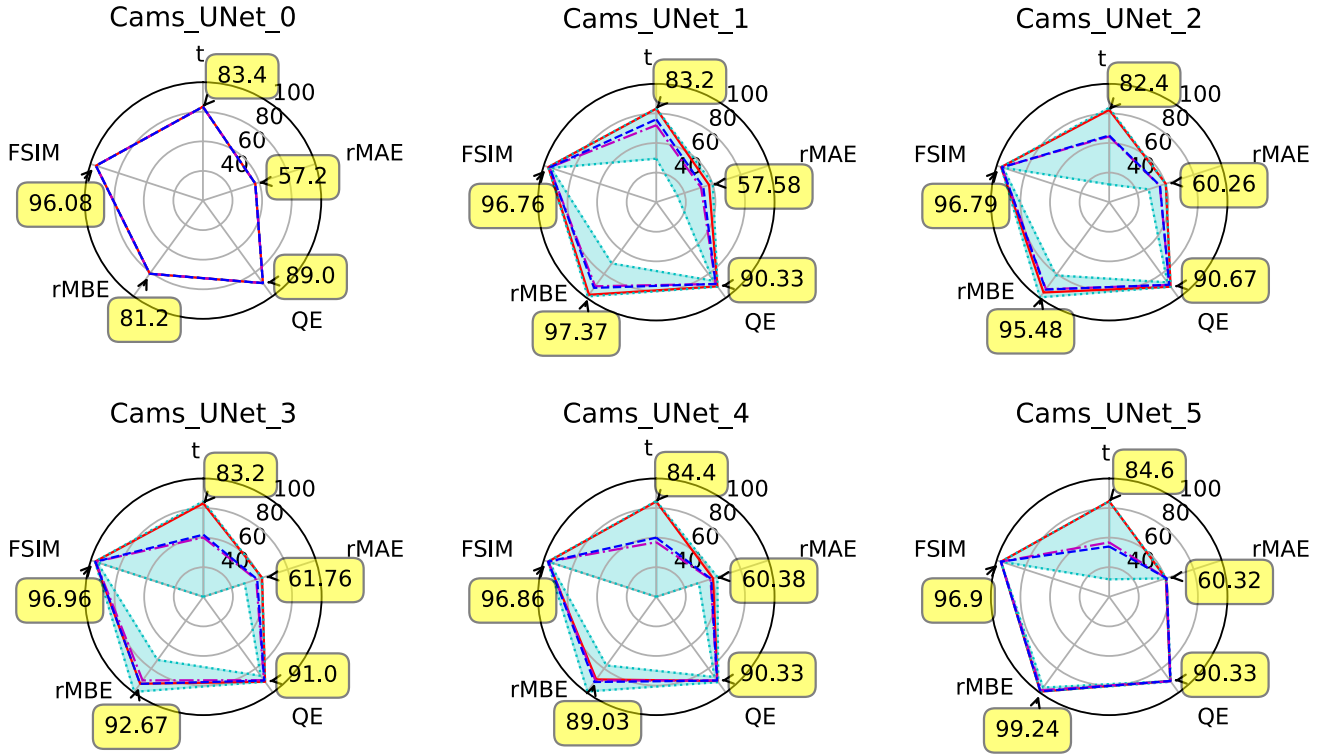


Figure 9. Overview of the evolution of our UNet’s performances when increasing the number of input variables. The wind is counted as one variables even though it contains two channels (one for wind speed and a second one for direction). Cyan dots: max and min, with a cyan area in between. Blue dashes: median. Purple dashdots: average. Red plain line: best total score (corresponds to annotated values).

Table 2. Best results on CAMS Dataset. In the column "variables used", we put the initial of each used variable: W for Wind, H for Humidity, P for Pressure, T for Temperature, A for Ångström Exponent. The AOD is always used as input. The column "fusion type" contains "Data" for Data Fusion, "Feature" for Feature Fusion, "Decision" for Decision Fusion. First line is the result with the best Total Score, second line is the best Timeless Score, and third line is the best Reduced Score.

Model Type	Fusion Type	Variables Used	Inference Time	MAE	QE	MBE	FSIM
UNet	Data	WHPTA	0.0077	4.38	0.29	-0.13	3.10%
UNet	Decision	WHPTA	0.0341	4.33	0.29	0.05	2.95%
UNet	Data	WHPT	0.0086	4.04	0.26	-2	2.8%

strategy, as well all proposed input variables, except for the Ångström exponent. If the Inference Time is not an important factor, then the use of a UNet model with the Decision Fusion strategy and all proposed input variables is advised. Finally, a UNet with the Data Fusion strategy as well as all proposed variables gives the most balanced results.

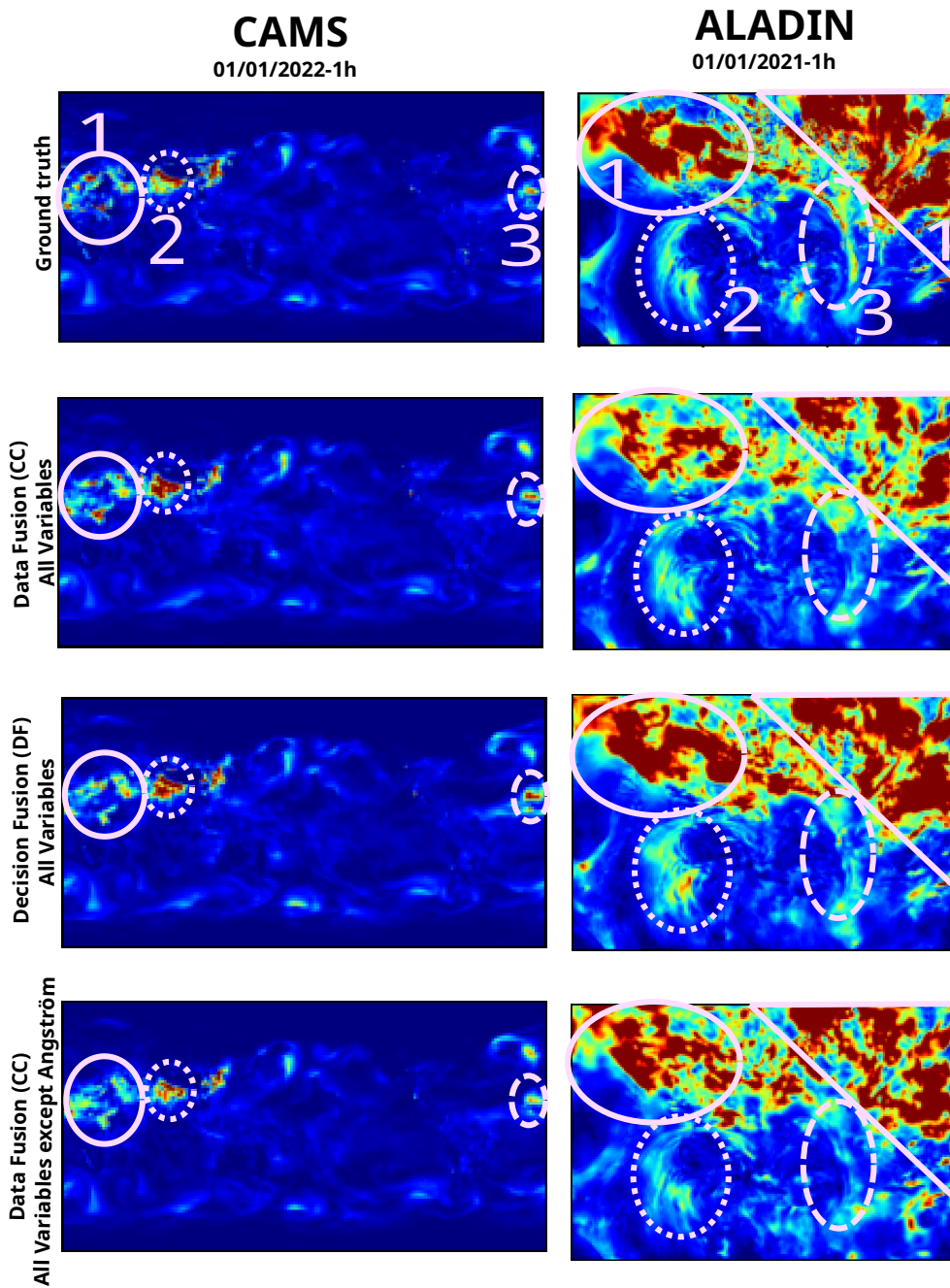


Figure 10. Outputs of the best models on both datasets for one given sample. The pink circles and numbers have been added afterwards to attract the reader's attention on some details of the image.

Looking at the left column Figure 10 gives us a bit more insight into the results. On this sample, it seems that the model using Data Fusion and all variables except the Ångström exponent is the one providing the best prediction for the area n°2 on the image. Other models overestimate the aerosol concentration in this area. The fact that this model has the worst *MBE*, and that it is negative could show a tendency to underestimation. The observations made on this sample are coherent with this assumption. The model using Data Fusion and all variables however provides good estimations for these two areas. Finally, the model using Decision Fusion (and all variables) underestimates the concentration in area n°1 and overestimates it in areas n°2 and 3. This model has the best *MBE*, but not the best *MAE*. Our hypothesis is that it overestimates certain areas and underestimates others, which compensates and leads to a small bias.

6.3 Comparison between the CAMS and ALADIN datasets

Table 3 shows the results obtained for the same models as in section 6.2 but on the ALADIN dataset, and the right column of Figure 10 shows their outputs for one given sample.

Table 3. Results on ALADIN Dataset. In the column "variables used", we put the initial of each used variable: W for Wind, H for Humidity, P for Pressure, T for Temperature, A for Ångström exponent. The AOD is always used as input. The column "fusion type" contains "Data" for Data Fusion, "Feature" for Feature Fusion, "Decision" for Decision Fusion.

Model Type	Fusion Type	Variables Used	Inference Time	MAE	QE	MBE	FSIM
UNet	Data	WHPTA	0.0064	8.37	0.35	-3.92	4.37%
UNet	Decision	WHPTA	0.0364	8.29	0.34	1.65	3.97%
UNet	Data	WHPT	0.0075	8.76	0.37	-1.04	4.64%

The table shows an important difference between the results obtained on the CAMS and the ALADIN dataset. However, even though the same metrics are used, these sets of results are not easily comparable to each other, as they are obtained on different data. Indeed, the ALADIN dataset contains images of much higher resolution than the CAMS dataset, these images do not represent the same geographical domain (Europe for ALADIN, the world for CAMS), and these dataset do not correspond to the same time period (July 2020 to June 2022 for CAMS, July 2019 to June 2021 for ALADIN). This explains why, in the CAMS dataset, the aerosol concentration values are comprised between 0 and 34 425 $\mu\text{g}/\text{m}^3$ with an average of 11.02 $\mu\text{g}/\text{m}^3$, while in the ALADIN dataset, they are comprised between 0 and 6 774 $\mu\text{g}/\text{m}^3$ with an average of 23.17 $\mu\text{g}/\text{m}^3$.

Figure 10 also shows a difference between results on the CAMS and ALADIN datasets. The model using Data Fusion with all variables underestimates the aerosol concentration in areas n°1 and 3, which is consistent with the fact that this model has the lowest *MBE* out of the three. The model using Data Fusion and all variables except the Ångström exponent also underestimates concentration in these areas, but less so. This is coherent with the fact that of all three models, this one has the second lowest *MBE*. The model using Decision Fusion does not underestimate concentration in areas n°1, underestimates the concentration in area n°3 as all other models, and overestimates the concentration in area n°2. It is also the model with the highest *MBE* and the best *MAE* out of all three.

Comparison between results on our two datasets does remain interesting, as the best performing methods on the CAMS dataset do not seem to correspond to the best performing ones on the ALADIN dataset. For example, let us look at the results from the table obtained with the Data Fusion strategy. One of these results is obtained while using all available variables as input, and the other is obtained using all variables except the Ångström exponent. Based solely on these two results, on the CAMS dataset it would seem using the Ångström exponent as part of the input variables leads to a smaller MBE, but we obtain higher values for all other metrics (except the inference time). On the ALADIN dataset the same situation and decision (of using the Ångström exponent) seem to lead to opposite results (higher MBE, smaller rest of the metrics). This also shows that the impact of the use of one specific input variable on the results of our models can not easily be interpreted. This is due to the interaction between the input variables themselves, and the very nature of the Neural Networks, which are often described (with reason) as black boxes.

6.4 Interpretation of the impact of the Ångström exponent

Let us look at Figure 11 to try and understand the impact of using the Ångström exponent on our results on the CAMS Dataset. This figure shows that the two metrics that are impacted the most by the use of the Ångström exponent are the *MAE* and the *MBE* (and their relative counterparts). Using the Ångström exponent seems to lead to a higher minimum value for the *MAE*. In other words, it helps avoiding our worst results (w.r.t. the *MAE*). The best *MBE* values are obtained when using the Ångström exponent. Using it therefore seems to lead to a lower bias.

Once again, these observations are valid for the CAMS Dataset, and for the chosen periods. We can not make a general conclusion on the use of the Ångström exponent as an input variable based on these observations alone. In particular, these observations are consistent with the results shown in table 2 (obtained with the CAMS Dataset), but not with those in table 3 (obtained with the ALADIN Dataset). This shows that our observations (about the Ångström exponent) on the CAMS Dataset can not automatically be assumed to be true for the ALADIN Dataset too.

6.5 Results of hybrid fusion method

Tables 4 and 5 shows the results of the two hybrid models described in section 4.3.4 on the CAMS and ALADIN datasets, respectively. Figure 12 shows the outputs of these models on one given sample.

Table 4. Results of hybrid models on the CAMS Dataset. The column "fusion type" contains "Hybrid1" for the model represented by Figure 5 and "Hybrid2" for the model represented by Figure 6.

Model Type	Fusion Type	Inference Time	MAE	QE	MBE	FSIM
UNet	Hybrid1	0.0098	4.39	0.28	0.2	2.96%
UNet	Hybrid2	0.0116	4.1	0.27	-0.86	2.9%

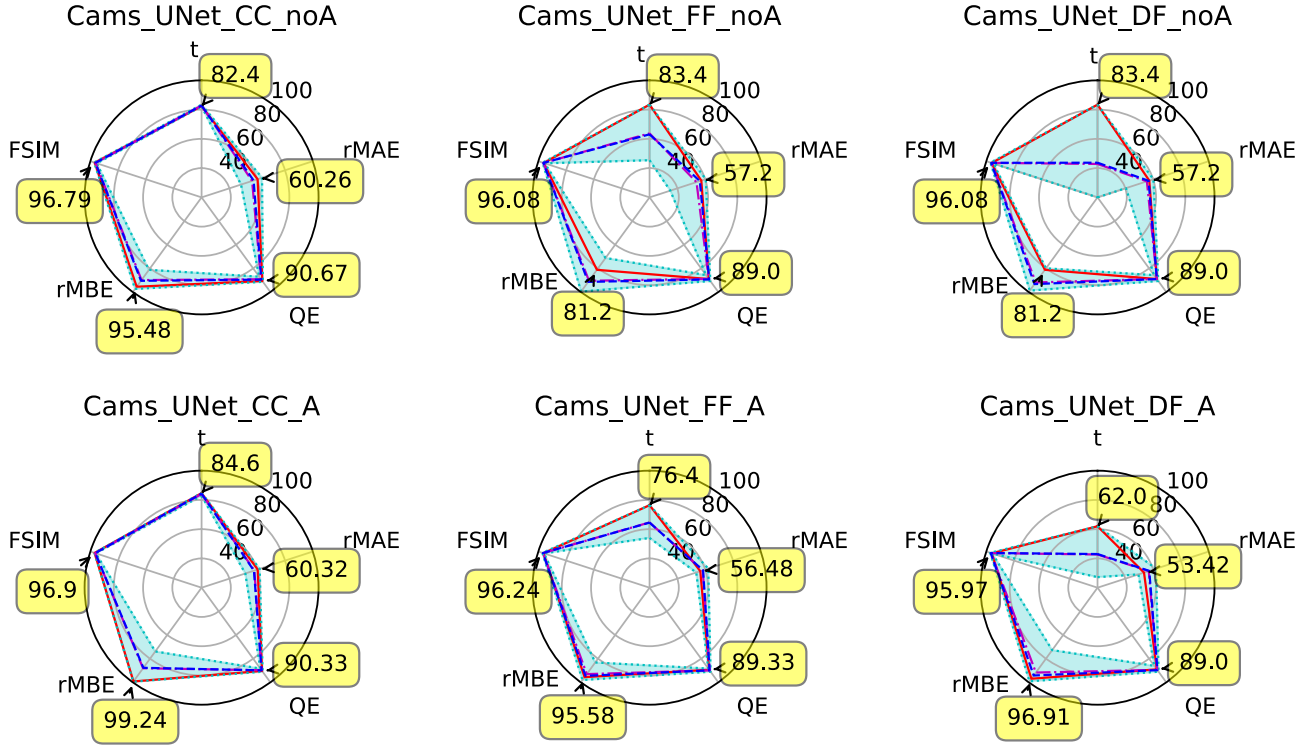


Figure 11. Overview of our experiments with (first line) and without (second line) using the Ångström exponent as an input variable. Regarding the fusion type, 'CC' means 'Channel Concatenation' or Data Fusion, 'FF' means Feature Fusion, and 'DF' means Decision Fusion. Cyan dots: max and min, with a cyan area in between. Blue dashes: median. Purple dashdots: average. Red plain line: best total score (corresponds to annotated values).

These results show that, from an Artificial Vision point of view, the second proposed hybrid model is better. However the first model appears to be more balanced, and is recommended in any situation where the *MBE* and Inference time are important metrics.

These models, while showing satisfying performance, show poorer performances than some of the results presented in table 2. Therefore we do not recommend the use of these hybrid models with the CAMS Dataset.

The left column of Figure 12 shows that both hybrid models produce a relatively adequate estimation for area n°3, underestimate concentration in area n°1, and overestimate it in area n°2. This is coherent with both models having relatively close metric values, and having *MBE* values close to 0.

These results show that, on the ALADIN dataset, the first proposed Hybrid model leads to better results than the second, on all metrics (except the Inference time).

The results obtained with this model are also better than all results presented in table 3. However, the model that were tested on the ALADIN dataset only correspond to the model that produced the best performances on the CAMS dataset. This means

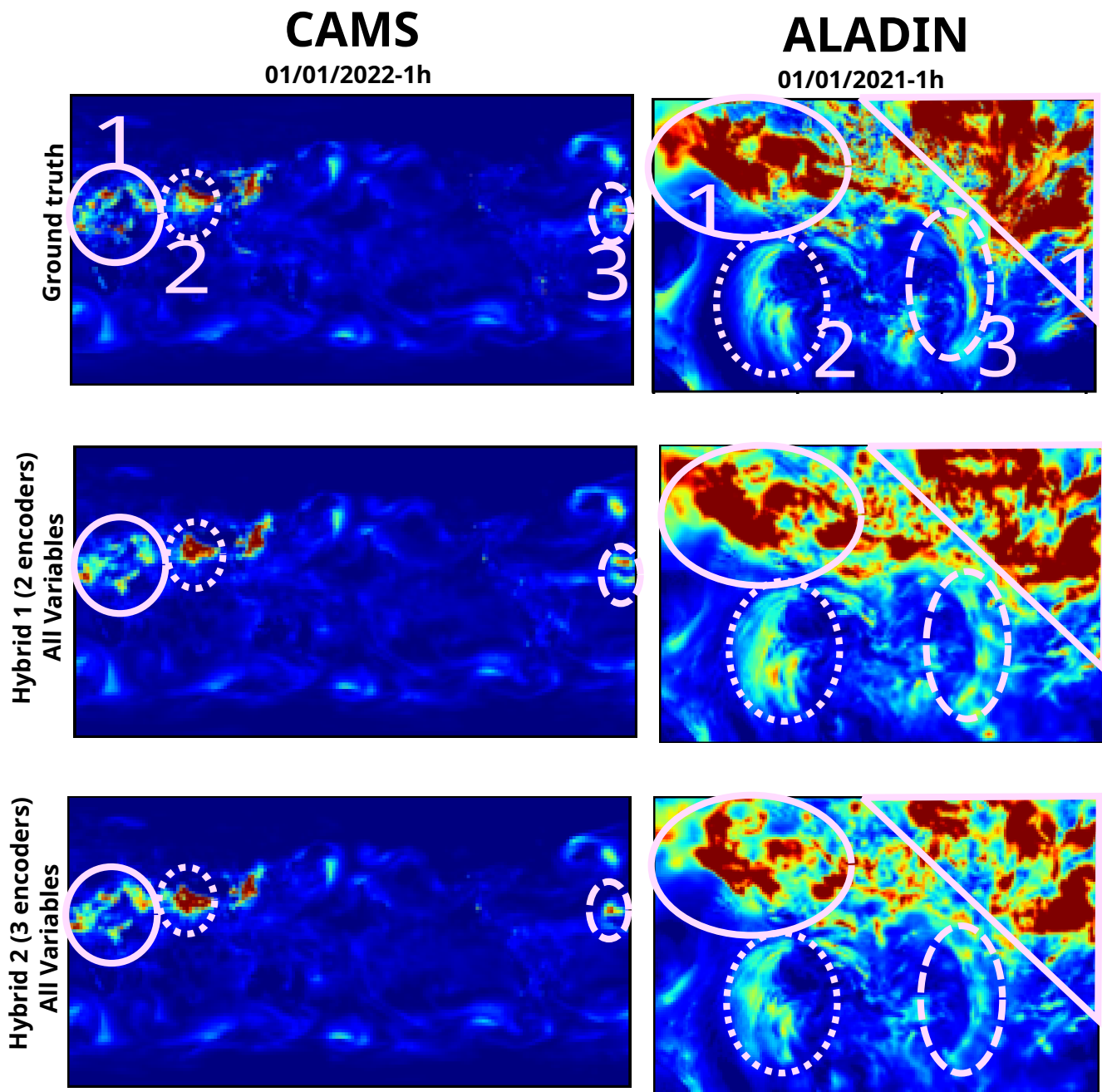


Figure 12. Outputs of hybrid models on both datasets for one given sample. The pink circles and numbers have been added afterwards to attract the reader's attention on some details of the image.

Table 5. Results of hybrid models on the ALADIN Dataset. The column "fusion type" contains "Hybrid1" for the model represented by Figure 5 and "Hybrid2" for the model represented by Figure 6.

Model Type	Fusion Type	Inference Time	MAE	QE	MBE	FSIM
UNet	Hybrid1	0.0108	7.69	0.33	-1.42	4%
UNet	Hybrid2	0.0103	7.9	0.33	-2.87	4.2%

545 that we can not conclude from these results that the Hybrid models work better than other models on the ALADIN dataset. To arrive to such a conclusion, we would need to realise an exhaustive study on our three fusion strategies, two models (GAN and UNet) and six input variables.

The right column of Figure 12 shows that the first hybrid model slightly underestimates concentration in areas n°1, and the second hybrid model underestimates it more. The first model slightly overestimates concentration in area n°2, while the second
550 model provides a more accurate estimation. These observation are coherent with both models having a low *MBE* and the second model having the lowest of the two. Interestingly enough, both these hybrid models seem to propose a better estimation of area n°3 than the models shown in Figure 10.

6.6 Comparison with SOTA

Table 6 shows a comparison of our best results with a few methods used as baseline on the CAMS dataset. It is important to
555 note that the Polynomial Interpolation and Random Forest Algorithm only use the AOD as input, while the kriging method only uses sparse values of the aerosol concentration (which represent our Boundary Conditions) as input.

Table 6. Comparison of baseline models with our best results on the CAMS dataset. Our best results on the CAMS dataset are those presented in table 2. "Poly. Interp." stands for Polynomial Interpolation, "Ord." for Ordinary, and "HE" for "hole-effect".

Model Type	Fusion Type	Variables Used	Inference Time	MAE	QE	MBE	FSIM
Poly. Interp. of Degree 3		AOD only	0.0007	6	0.41	-3	4.96%
Random Forest Algorithm		AOD only	1.1535	6.01	0.41	-2.63	4.95%
Ord. kriging with HE variogram		BC only	40.3732	6.03	0.38	-2.99	7.67%
UNet	Data	WHPTA	0.0077	4.38	0.29	-0.13	3.10%
UNet	Decision	WHPTA	0.0341	4.33	0.29	0.05	2.95%
UNet	Data	WHPT	0.0086	4.04	0.26	-2	2.8%

The Polynomial Interpolation method has a significantly smaller inference time than any other method discussed in this paper. However this is the only metric on which one of the baseline methods outperforms our best results. Indeed, our models outperform the chosen baselines by a large margin, in all metrics except this one.

560 Our hybrid models do not appear in table 6, as they are outperformed by the models presented in this table. However, as stated in section 6.5, their performances remain comparable. In other words, our hybrid models are outperformed by the models presented in table 6, but not by a large margin.

It has already been stated that, in a context where ground truths were less available, GANs would outperform UNets. Indeed their usefulness for this problem lies in their ability to realise semi-supervised learning. It is also interesting to note
565 that generally speaking, all models would probably benefit from a larger amount of data, as long as the training set remains representative of the actual data in a real-case scenario. The representativity of the dataset is paramount as it helps avoiding the overfitting problem often encountered in machine learning. More specifically, our deep learning models are the ones that would benefit the most from a larger amount of accessible data, as they contain more parameters.

7 Conclusion

570 In this paper we performed an extensive study on the use of several meteorological variables and column aerosol optical properties as inputs for a Deep Learning model to infer PM_{2.5} concentration from AOD using a scaling approach applicable globally. We tested different network architectures as well as the use of three different fusion strategies for the exploitation of these inputs, in order to investigate the optimal way of fusing those information for our specific application. Hybrid methods of fusion have been proposed, implemented and studied as well. Our experiments were conducted extensively on CAMS data in
575 order to assess model performances at global scale. We also performed limited experiment using the ALADIN dataset (instead of CAMS) over a large region covering Europe and the Mediterranean basin to study the impact of the datasets' characteristics on our results, especially its spatial resolution and geographic spatial coverage.

Based on five metrics used throughout to evaluate different models performances, our experiments have shown the superiority of UNets over BC-GANs in our context, as is shown by figure 8. However the sparse training set is, in our context,
580 significantly smaller than the complete set. We suggest in section 6.1 that this induces a reduced need for semi-supervised learning, and explains the difference in performance between UNets and GANs. The authors of (Dabrowski et al., 2023) show the superiority of their BC-GANs over UNets in their context, which includes sparse and complete training sets of more comparable size. This shows that the difference in performance between BC-GANs and UNets is not inherent to these models themselves. Therefore we recommend the use of a UNet in our context of semi-supervised learning, with our sparse training
585 set being significantly smaller than our complete training set. It remains difficult to deduce a superiority of one model over the other in the general sense from our experiments. The comparison between our results and the results of (Dabrowski et al., 2023) does show that the quantity of sparse data has a significant impact on the performances of these two models. Therefore this context parameter must be taken into account when recommending one of these models over the other.

Our results have also illustrated that increasing the number variables as input tends to augment model performances. This is
590 not surprising as the limited set of variables we used were selected for their known influence on PM_{2.5} surface concentration. This remains a tendency however, and not a guarantee as some exceptions have been observed where, depending on network architecture and fusion strategy, adding a variable may degrade performance. This is interesting because it is counter-intuitive

to the general belief that using more (relevant) data in deep learning yields better results and emphasises in particular the interest of studying the impact of network architecture for atmospheric applications.

595 Our experiments have also shown in section 6.3 the importance of dataset’s characteristics (here spatial resolution and coverage) and its impact not only on the results but on the conclusions that can be drawn from them as well. This is especially important in atmospheric sciences because geophysical variables have different scales of variability and network architecture should ideally be aligned with the spatial characteristics of input fields. Our work suggests that more work is needed to understand the impact of networks architecture on their ability to fully capture spatial features that are specific to atmospheric
600 sciences.

While identifying precisely the impact of each variable on the models’ performances would be useful, the observations made in section 6.4, and drawn from our results, highlight the difficulty of such a task.

The two fusion strategies that lead to our best results (shown in section 6.2) are the Data and Decision Fusion ones. According to our experiments, the Data Fusion strategy also seems to lead to more stable results. Moreover, it allows to build smaller
605 models, which in turn leads to shorter inference times and training times.

Our experiments on hybrid models did not show clear evidence of their advantage compared to other models, even though they do present comparable performances, as shown in section 6.5. Based on these conclusions, the Data Fusion strategy is the one we would recommend in a general case when all input variables are available at the same resolution and over the same area. Of course this recommendation depends on general context and more specifically on the definition of the desired outcome. For
610 example, using different metrics to measure performance might lead to a different recommendation.

Finally, our objective was not to develop a single and optimized model for PM2.5 inference from AOD but rather to study how multiple PM2.5 predictors could be used in order to best align the network architecture with the seek inference function. However while we did not try to conduct specific optimization and used only a limited set of predictors, we have proposed several architectures that yield PM2.5 inference performances comparable to other tailored models found in the literature (Ma
615 et al., 2022; Unik et al., 2023). The demonstrated performances obtained here should only be interpreted as baseline capabilities of the proposed models that could most likely be improved by extending further the time coverage of the learning database. As suggested by Zhou et al. (2024), we also strongly encourage a more systematic evaluation of models against a common test dataset and using standardized metrics. Since the code and data used for this article are both available, we suggest that our current results be seen as a benchmark for the task and context presented in this paper. Such a benchmark could be used
620 as a common ground for the evaluation of newly developed models of PM2.5 inference AOD data, therefore facilitating their comparison.

As stated before, the experiments realised in this paper have clearly illustrated the interest of using additional, carefully chosen, input variables in order to augment the performance of a scaling model to infer PM2.5 from AOD. Here we selected a limited number of meteorological variables and optical properties that are well known to drive surface level PM2.5 con-
625 centration. These variables are typically useful to establish the link between PM2.5 and AOD through a purely physics based model and not surprisingly our results demonstrate they are useful to establish this link through an Artificial Neural Network. Based on this insight, an interesting possibility of future work consists in applying the concept of Physics-Informed Neural

Networks (described in detail in section A) to this problem and study, depending of the fusion strategy used, at which level the incorporation of physics equations would be most relevant.

630 *Code and data availability.*

- The code used for these experiments is available on a Zenodo archive (Dabrowski, 2024a).

<https://doi.org/10.5281/zenodo.13947256>

The data from the CAMS model used during these same experiments is available on a different Zenodo archive (Dabrowski, 2024b).

<https://doi.org/10.5281/zenodo.13929498>

635 – The data from the ALADIN model was extracted from the dataset proposed by (Mallet and Nabat, 2024).

<https://doi.org/10.25326/703>

Appendix A: Related works

Generative Adversarial Networks (GANs). Since the authors of (Goodfellow et al., 2020) proposed this type of model, the popularity of GANs has increased consistently. They rely on the training of two networks : a generator and a discriminator.

640 The discriminator is presented with samples which can be either taken from the original data distribution, or generated (by the generator). Its main task is to differentiate these two kinds of inputs. On the other hand, if the discriminator makes an error and classifies a generated sample as real, then the generator is getting closer to its goal. The discriminator and generator’s losses are built in such a way that when one increases, the other decreases, and reversely. This why they are called adversarial networks.

Convolutional GANs are known for their ability to produce realistic images, which can fool both their discriminator but
645 also in some cases humans. They have also shown interesting performance in Image-to-Image translation tasks (Wang et al., 2019, 2020; Zhu et al., 2017b). This type of task is usually categorized as paired image translation such as in (Isola et al., 2017), or unpaired image translation such as in (Zhu et al., 2017a). In this article, our image translation task is a paired one.

Explainable Artificial Intelligence (XAI). Even though this work can not be classified as belonging to the field of XAI, the terminology this field proposes remains interesting in the context of this work. The general idea behind XAI is to build models
650 that can be understood by their users, or whose results can. While this field has existed for several decades (Confalonieri et al., 2021), its recent growth in popularity can be seen as a response to concerns about the black-box aspect of some neural networks models. This growth has been particularly remarkable in applications fields like finance, medicine, law, and even scientific production (Beckh et al., 2021; Murdoch et al., 2019; Belle and Papantonis, 2021; Roscher et al., 2020). In those fields, the ability to explain a model and its results can represent the ability to ensure safety, fairness or scientific rigour. In a
655 more general sense, it makes it easier for the user to trust the model.

According to (Roscher et al., 2020), in the context of XAI, there are three important elements to consider when evaluating the explainability of a model.

1. **Transparency:** an model is transparent if the processes that extract model parameters from training data and generate labels from testing data can be described and motivated by its designer.
- 660 2. **Interpretability:** its is the ability to generally understand what the model bases its decisions on. Some approaches for interpretable models are based on decision trees, as they can allow for an intuitive look on the decision-making process of a model.
3. **Explainability:** an explanation is the collection of features of the interpretable domain, that have contributed for a given example to produce a decision. For a model to be explainable, it generally needs to be possible to understand why the
665 model’s decision for a given datum A is different than for a given datum B.

It is interesting to note that domain knowledge can be used to enhance the explainability of a model (Beckh et al., 2021). In this sense, Physics-Informed Neural Networks can be seen as a type of XAI.

Physics-Informed Neural Networks (PINNs). Physics-Informed learning, introduced by the authors of (Raissi et al., 2019), can be considered today as its own research field. The term of Informed Networks suggest that the method makes use of prior

670 information about some specificity of the problem, for example its geometry. Physics-Informed Networks specifically make use of the physics of the problem to enhance their performances. This is usually done through the design of a physics-informed loss function, used during the training of the model. This loss function is often based off a differential equation that is verified by the data the model is using. As it can sometimes guide the training in a non-data oriented way, the use of this loss function reduces the need of these networks for labeled data, making them especially suitable for semi-supervised learning.

675 In physics, it is often necessary to use Initial and Boundary Conditions (BCs) to solve a given problem. In the literature around physics-based learning methods, two methods to take these BCs into account during training can be found. The soft constraint (or method) proposes to train the model to respect the BCs, through the use of an additional, tailored loss function. The hard constraint (or method) works through the transformation of the model outputs to enforce the respect of the BCs, and relies on pre-existing loss functions. When it comes to PINNs, the authors of (Sun et al., 2020) show that the hard constraint performs better than the soft.

Several authors have proposed to leverage the advantages shown by both adversarial and physics-informed approaches (Thanasutives et al., 2021; Nie et al., 2021), often calling these new models PI-GANs (Yang et al., 2019, 2020).

Kriging method. This spatial interpolation and extrapolation method was formalised by the author of (Matheron, 1963). In the statistical interpretation of the term, it is the optimal estimation method according to (Gratton, 2002). It is mathematically
685 described by equation A1.

$$F(x_p) = \sum_{i=1}^m W_i \cdot F(x_i) \quad (\text{A1})$$

$F(x_p)$, the value of function F at point x_p , can be estimated thanks to m surrounding points x_i , as the value of F at these points is known. However it remains necessary to determine the weights W_i of these points. The kriging method proposes to realise this through the estimation of what is called a variogram. To compute it, values of the variance of two points, and of the
690 distance between them, are needed.

This method has been described as performing better when provided with a significant volume of data, and when the values to estimate are following a normal distribution.

For each inference, new points x_i are used. As these points are the basis for the building of the kriging model, a new one is built for each inference. Because of this, kriging suffers from long inference time when compared to other methods presented
695 in this article.

Appendix B: Details of the approach

The purpose of this section is to show the architecture of the GAN's discriminator with our three main fusion strategies.

B1 Data Fusion / Channel Concatenation

Figure B1 shows the GAN's discriminator's architecture with the data fusion strategy. Related

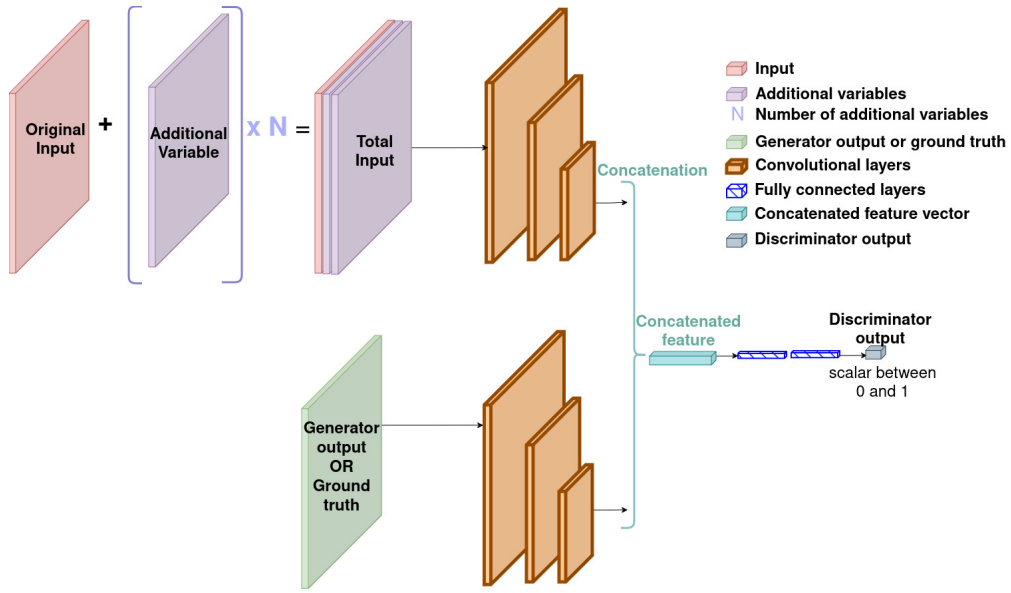


Figure B1. Architecture of the GAN’s discriminator with data fusion approach.

700 B2 Feature Fusion

Figure B2 shows the GAN’s discriminator’s architecture with the feature fusion strategy.

B3 Decision Fusion

Figure B3 shows the GAN’s discriminator’s architecture with the data fusion strategy.

Appendix C: Models complexity

705 Figure C1 shows the number of parameters of our models depending on the fusion method and the number of input images used. They correspond to the number of parameters for our UNets and BC-GANs, with both ALADIN and CAMS data.

Appendix D: Training loss, test loss and convergence

Figure D1 provides a graph of training loss values over iterations, showing clearly the convergence of the model. This corresponds to the training of a UNet model using exclusively the AOD as input. In this experiment as in all other experiments
710 presented in this article, the models are trained on 500 epochs.

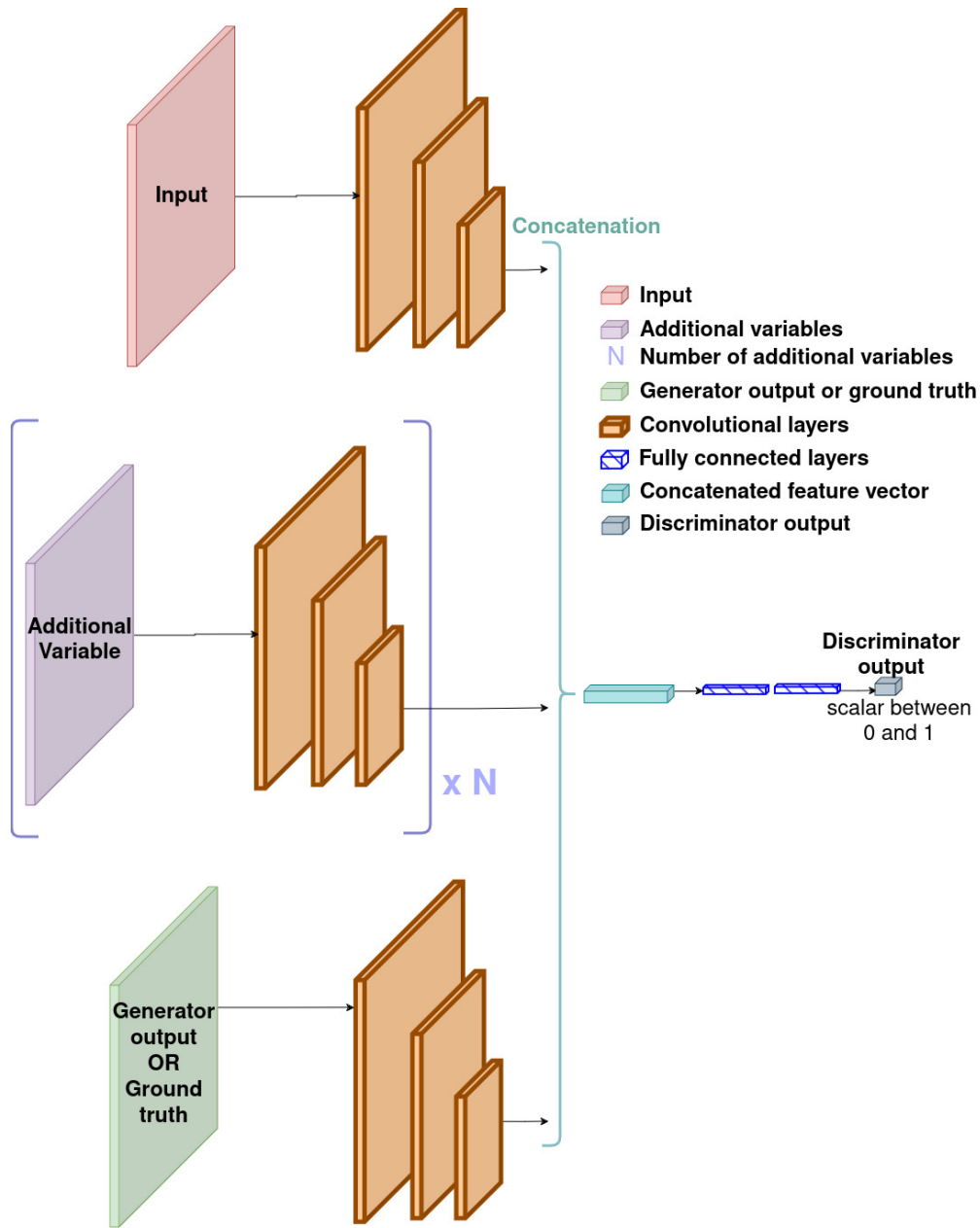


Figure B2. Architecture of the GAN's discriminator with feature fusion approach.

Figure D2 gives, for the same model, an overview of the MAE values for the different test samples. A few test samples stand out as having a significantly worse MAE than others, but the maximum MAE for these samples remains below $3\mu g/m^3$, which is satisfying.

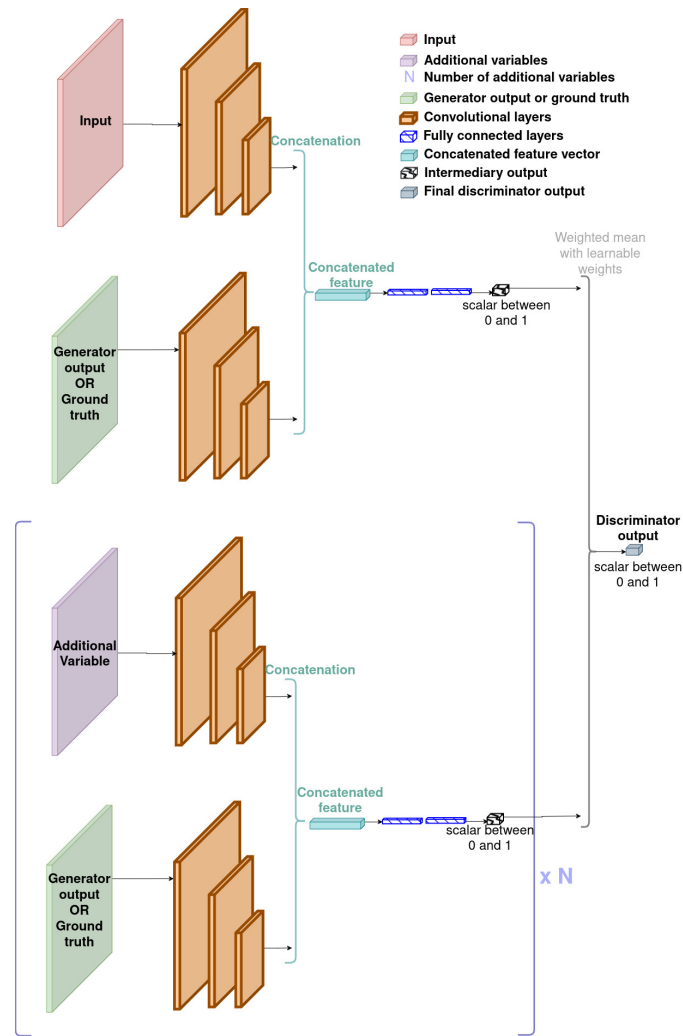


Figure B3. Architecture of the GAN's discriminator with decision fusion approach.

Author contributions.

715 **Matthieu Dabrowski** : main author, designed the models and experiments, performed the experiments, led the analysis and the writing of this article.

José Mennesson : provided insights on the models and metrics used, helped in analysing the results, and contributed to the writing of this article.

720 **Jérôme Riedi** : defined the atmospheric model datasets to be used in the experiments, provided insights about physics of input variables and PM2.5 values, helped in the analysis and the general writing of this article

Chaabane Djeraba : provided insight on the NN models and on the general research strategy, as well as useful and welcomed tips for the writing of this article.

Pierre Nabat : provided the ALADIN data and insights into specifics of atmospheric composition models.

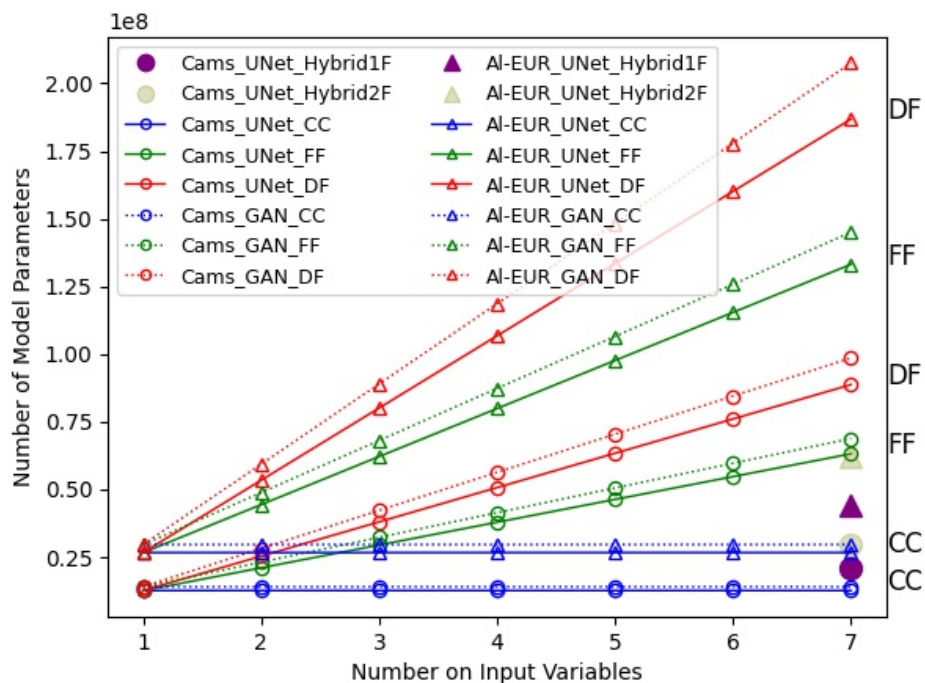


Figure C1. Number of parameters of each of our models depending on the number of input variables.

Competing interests. The authors declare an absence of competing interests.

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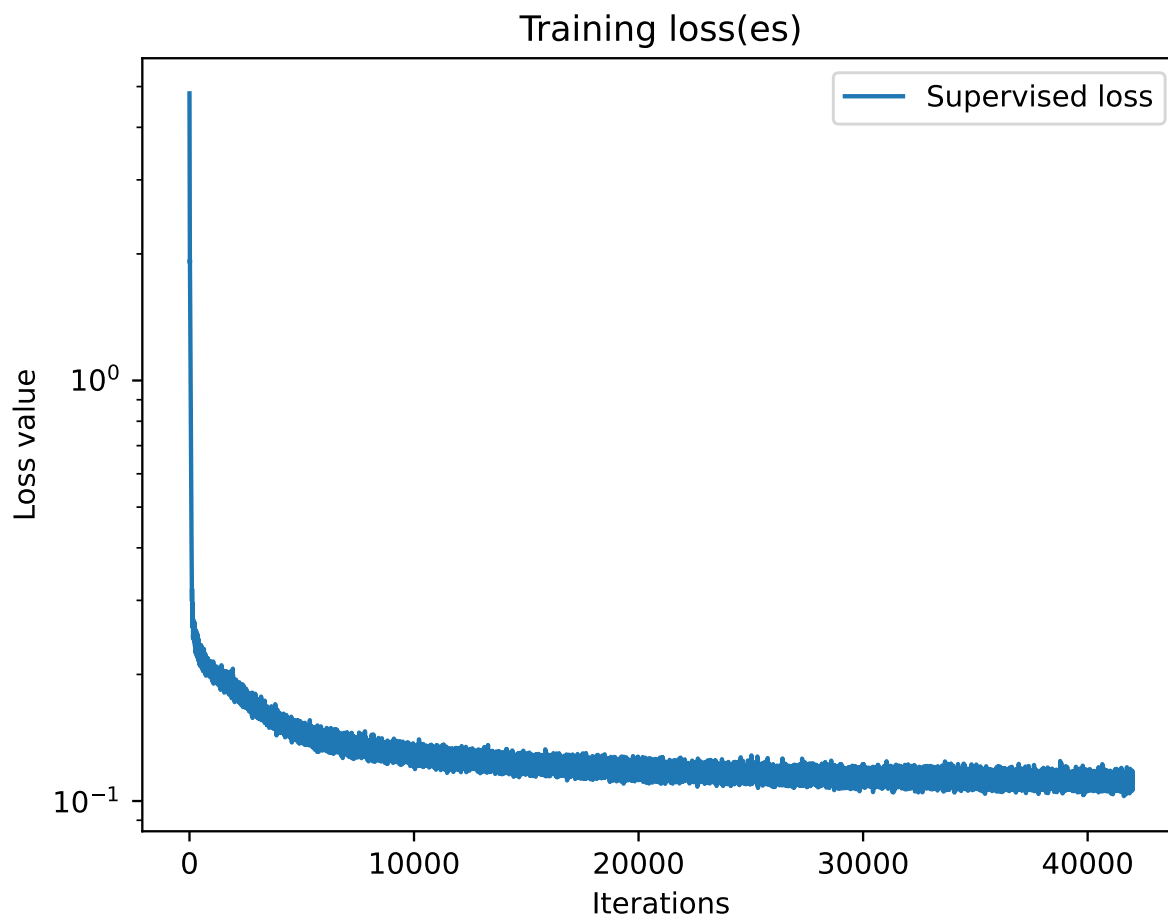


Figure D1. Graph of training loss during supervised learning over iterations

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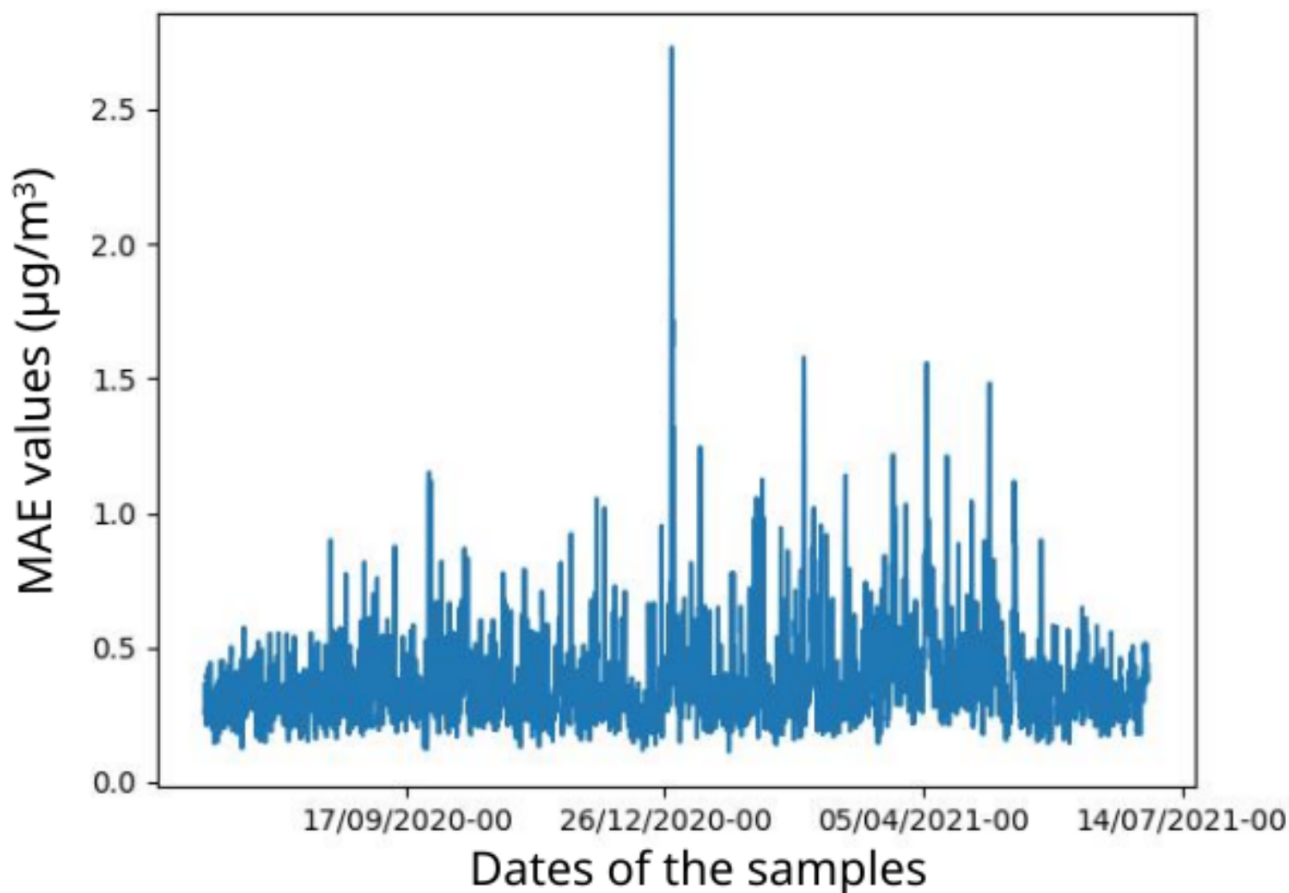


Figure D2. Graph of MAE values during testing over sample date

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