

# A review of innovation-based methods to jointly estimate model and observation error covariance matrices in ensemble data assimilation

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2	<b>Observation Error Covariance Matrices in Ensemble Data Assimilation</b>
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## ABSTRACT

Data assimilation combines forecasts from a numerical model with observa-24 tions. Most of the current data assimilation algorithms consider the model and 25 observation error terms as additive Gaussian noise, specified by their covari-26 ance matrices Q and R, respectively. These error covariances, and specifically 27 their respective amplitudes, determine the weights given to the background 28 (i.e., the model forecasts) and to the observations in the solution of data as-29 similation algorithms (i.e., the analysis). Consequently, Q and R matrices sig-30 nificantly impact the accuracy of the analysis. This review aims to present and 3 to discuss, with a unified framework, different methods to jointly estimate the 32 Q and R matrices using ensemble-based data assimilation techniques. Most 33 of the methodologies developed to date use the innovations, defined as differ-34 ences between the observations and the projection of the forecasts onto the 35 observation space. These methodologies are based on two main statistical 36 criteria: (i) the method of moments, in which the theoretical and empirical 37 moments of the innovations are assumed to be equal, and (ii) methods that 38 use the likelihood of the observations, themselves contained in the innova-39 tions. The reviewed methods assume that innovations are Gaussian random 40 variables, although extension to other distributions is possible for likelihood-41 based methods. The methods also show some differences in terms of levels of 42 complexity and applicability to high-dimensional systems. The conclusion of 43 the review discusses the key challenges to further develop estimation meth-44 ods for Q and R. These challenges include taking into account time-varying 45 error covariances, using limited observational coverage, estimating additional 46 47

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#### 48 1. Introduction

In meteorology and other environmental sciences, an important challenge is to estimate the state of the system as accurately as possible. In meteorology, this state includes pressure, humidity, temperature and wind at different locations and elevations in the atmosphere. Data assimilation (hereinafter DA) refers to mathematical methods that use both model predictions (also called background information) and partial observations to retrieve the current state vector with its associated error. An accurate estimate of the current state is crucial to get good forecasts, and it is particularly so whenever the system dynamics is chaotic, such as it is the case for the atmosphere.

The performance of a DA system to estimate the state depends on the accuracy of the model 56 predictions, the observations, and their associated error terms. A simple, popular and mathemat-57 ically justifiable way of modeling these errors is to assume them to be independent and unbiased 58 Gaussian white noise, with covariance matrices  $\mathbf{Q}$  for the model and  $\mathbf{R}$  for the observations. Given 59 the aforementioned importance of  $\mathbf{Q}$  and  $\mathbf{R}$  in estimating the analysis state and error, a number of 60 studies dealing with this problem has arisen in the last decades. This review work presents and 61 summarizes the different techniques used to estimate simultaneously the Q and R covariances. 62 Before discussing the methods to achieve this goal, the mathematical formulation of DA is briefly 63 introduced. 64

#### 65 a. Problem statement

Hereinafter, the unified DA notation proposed in Ide et al. (1997) is used<sup>1</sup>. DA algorithms are used to estimate the state of a system,  $\mathbf{x}$ , conditionally on observations,  $\mathbf{y}$ . A classic strategy is to use sequential and ensemble DA frameworks, as illustrated in Fig. 1, and to combine two sources of information: model forecasts (in green) and observations (in blue). The ensemble framework

<sup>&</sup>lt;sup>1</sup>Other notations are also used in practice

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<sup>70</sup> uses different realizations, also called members, to track the state of the system at each assimilation
<sup>71</sup> time step.

The forecasts of the state are based on the usually incomplete and approximate knowledge of the system dynamics. The evolution of the state from time k - 1 to k is given by the model equation:

$$\mathbf{x}(k) = \mathscr{M}_k(\mathbf{x}(k-1)) + \boldsymbol{\eta}(k), \tag{1}$$

<sup>74</sup> where the model error  $\eta$  implies that the dynamic model operator  $\mathcal{M}_k$  is not perfectly known. <sup>75</sup> Model error is usually assumed to follow a Gaussian distribution with zero mean (i.e., the model <sup>76</sup> is unbiased) and covariance **Q**. The dynamic model operator  $\mathcal{M}_k$  in Eq. (1) has also an explicit <sup>77</sup> dependence on k, because it may depend on time-dependent external forcing terms. At time k, <sup>78</sup> the forecasted state is characterized by the mean of the forecasted states,  $\mathbf{x}^f$ , and its uncertainty <sup>79</sup> matrix, namely  $\mathbf{P}^f$ , which is also called the background error covariance matrix, and noted **B** in <sup>80</sup> DA.

The forecast covariance  $\mathbf{P}^{f}$  is determined by two processes. The first is the uncertainty propa-81 gated from k-1 to k by the model  $\mathcal{M}_k$  (the green shade within the dashed ellipse in Fig. 1, and 82 denoted by  $\mathbf{P}^{m}$ ). The second process is the model error covariance  $\mathbf{Q}$  accounted by the noise term 83 at time k in Eq. (1). Given that model error is largely unknown and originated by various and 84 diverse sources, the matrix  $\mathbf{Q}$  is also poorly known. Model error sources encompass the model  $\mathcal{M}$ 85 deficiencies to represent the underlying physics, including deficiencies in the numerical schemes, 86 the cumulative effects of errors in the parameters, and the lack of knowledge of the unresolved 87 scales. Its estimation is a challenge in general, but it is particularly so in geosciences because we 88 usually have far fewer observations than those needed to estimate the entries of  $\mathbf{Q}$  (Daley 1992; 89 Dee 1995). The sum of the two covariances  $\mathbf{P}^m$  and  $\mathbf{Q}$  gives the forecast covariance matrix,  $\mathbf{P}^f$ 90 (full green ellipse in Fig. 1). In the illustration given here, a large contribution of the forecast co-91

variance  $\mathbf{P}^{f}$  is due to  $\mathbf{Q}$ . This situation reflects what is common in ensemble DA, where  $\mathbf{P}^{m}$  can be too small, as a consequence of the ensemble undersampling of the initial condition error (i.e., the covariance estimated at the previous analysis). In that case, inflating  $\mathbf{Q}$  could partially compensate for the bad specification of  $\mathbf{P}^{m}$ .

<sup>96</sup> DA uses a second source of information, the observations **y**, which are assumed to be linked to <sup>97</sup> the true state **x** through the time-dependent operator  $\mathscr{H}_k$ . This step in DA algorithms is formalized <sup>98</sup> by the observation equation:

$$\mathbf{y}(k) = \mathscr{H}_k(\mathbf{x}(k)) + \boldsymbol{\epsilon}(k), \tag{2}$$

where the observation error  $\epsilon$  describes the discrepancy between what is observed and the truth. 99 In practice, it is important to remove as much as possible the large-scale bias in the observation 100 before DA. Then, it is common to state that the remaining error  $\epsilon$  follows a Gaussian and unbiased 101 distribution with a covariance  $\mathbf{R}$  (the blue ellipse in Fig. 1). This covariance takes into account er-102 rors in the observation operator  $\mathcal{H}$ , the instrumental noise and the representation error associated 103 with the observation, typically measuring a higher resolution state than the model represents. Op-104 erationally, a correct estimation of **R** that takes into account all these effects is often challenging 105 (Janjić et al. 2018). 106

DA algorithms combine forecasts with observations, based on the model and observation equa-107 tions, respectively given in Eq. (1) and Eq. (2). The corresponding system of equations is a non-108 linear state-space model. As illustrated in Fig. 1, this Gaussian DA process produces a posterior 109 Gaussian distribution with mean  $\mathbf{x}^a$  and covariance  $\mathbf{P}^a$  (red ellipse). The system given in Eqs. (1) 110 and (2) is representative of a broad range of DA problems, as described in seminal papers such 111 as Ghil and Malanotte-Rizzoli (1991), and still relevant today as referenced by Houtekamer and 112 Zhang (2016) and Carrassi et al. (2018). The assumptions made in Eqs. (1) and (2) about model 113 and observation errors (additive, Gaussian, unbiased, and mutually independent) are strong, yet 114

convenient from the mathematical and computational point of view. Nevertheless, these assumptions are not always realistic in real DA problems. For instance, in operational applications, systematic biases in the model and in the observations are recurring problems. Indeed, biases affect significantly the DA estimations and a specific treatment is required; see Dee (2005) for more details.

From Eqs. (1) and (2), noting that  $\mathcal{M}, \mathcal{H}$  and y are given, the only parameters that influence the 120 estimation of  $\mathbf{x}$  are the covariance matrices  $\mathbf{Q}$  and  $\mathbf{R}$ . These covariances play an important role in 121 DA algorithms. Their importance was early put forward in Hollingsworth and Lönnberg (1986), 122 in section 4.1 of Ghil and Malanotte-Rizzoli (1991) and Daley (1991) in section 4.9. The results 123 of DA algorithms highly depend on the two error covariance matrices  $\mathbf{Q}$  and  $\mathbf{R}$ , which have to be 124 specified by the users. But these covariances are not easy to tune. Indeed, their impact is hard to 125 grasp in real DA problems with high-dimensionality and nonlinear dynamics. We thus illustrate 126 the problem with a simple example first. 127

#### 128 b. Illustrative example

In either variational or ensemble-based DA methods, the quality of the reconstructed state (or hidden) vector **x** largely depends on the relative amplitudes between the assumed observation and model errors (Desroziers and Ivanov 2001). In Kalman filter based methods, the signal-to-noise ratio  $\|\mathbf{P}^{f}\| / \|\mathbf{R}\|$ , where  $\mathbf{P}^{f}$  depends on  $\mathbf{Q}$ , impacts the Kalman gain, which gives the relative weights of the observations against the model forecasts. Here, the  $\|.\|$  operator represents a matrix norm. For instance, Berry and Sauer (2013) used the Frobenius norm to study the effect of this ratio in the reconstruction of the state in toy models.

The importance of **Q**, **R** and  $\|\mathbf{P}^{f}\| / \|\mathbf{R}\|$  is illustrated with the aid of a toy example, using a scalar state *x* and simple linear dynamics. This simplified setup avoids several issues typical

<sup>138</sup> of realistic DA applications: the large dimension of the state, the strong nonlinearities and the <sup>139</sup> chaotic behavior. In this example, the dynamic model in Eq. (1) is a first-order autoregressive <sup>140</sup> model, denoted by AR(1) and defined by

$$x(k) = 0.95x(k-1) + \eta(k), \tag{3}$$

with  $\eta \sim \mathcal{N}(0, Q^t)$  where the superscript t means "true" and  $Q^t = 1$ . Furthermore, observations y 141 of the state are contaminated with an independent additive zero-mean and unit-variance Gaussian 142 noise, such that  $R^t = 1$  in Eq. (2) with  $\mathcal{H}(x) = x$ . The goal is to reconstruct x from the noisy ob-143 servations y at each time step. The AR(1) dynamic model defined by Eq. (3) has an autoregressive 144 coefficient close to one, representing a process which evolves slowly over time, and a stochastic 145 noise term  $\eta$  with variance Q<sup>t</sup>. Although the knowledge of these two sources of noise is crucial 146 for the estimation problem, identifying them is not an easy task. Given that the dynamic model is 147 linear and the error terms are additive and Gaussian in this simple example, the Kalman smoother 148 provides the best estimation of the state (see section 2 for more details). To evaluate the effect 149 of badly specified Q and R errors on the reconstructed state with the Kalman smoother, different 150 experiments were conducted with values of  $\{0.1, 1, 10\}$  for the ratio Q/R (in this toy example, we 151 use Q/R instead of  $\|\mathbf{P}^f\| / \|\mathbf{R}\|$  for simplicity). 152

Figure 2 shows, as a function of time, the true state (red line) and the smoothing Gaussian distributions represented by the 95% confidence intervals (gray shaded) and their means (black lines). We also report the Root Mean Squared Error (RMSE) of the reconstruction and the socalled "coverage probability", or percentage of *x* that falls in the 95% confidence intervals (defined as the mean  $\pm 1.96$  the standard deviation in the Gaussian case). In this synthetic experiment, the best RMSE and coverage probability obtained, applying the Kalman smoother with true  $Q^t = R^t = 1$ , are 0.71 and 95%, respectively. Using a small model error variance  $Q = 0.1Q^t$  in Fig. 2(a),

the filter gives a large weight to the forecasts given by the quasi-persistent autoregressive dynamic 160 model. On the other hand, with a small observation error variance  $R = 0.1R^t$  in Fig. 2(b), excessive 161 weight is given to the observation and the reconstructed state is close to the noisy measurements. 162 These results show the negative impact of independently badly scaled Q and R error variances. In 163 the case of overestimated model error variance as in Fig. 2(c), the mean reconstructed state vector 164 and thus its RMSE are identical to Fig. 2(b). In the same way, overestimated observation error 165 variance like in Fig. 2(d) gives similar mean reconstruction, as in Fig. 2(a). These last two results 166 are due to the fact that in both cases, the ratio Q/R are equal, respectively, to 10 and 0.1. Now, 167 we consider in Fig. 2(e) and Fig. 2(f) the case where the Q/R ratio is equal to 1, but, respectively, 168 using the simultaneous underestimation and overestimation of model and observation errors. In 169 both cases, the mean reconstructed state is equal to that obtained with the true error variances (i.e., 170 RMSE=0.71). The main difference is the gray confidence interval, which is supposed to contain 171 95% of the true trajectory: the spread is clearly underestimated in Fig. 2(e) and overestimated in 172 Fig. 2(f), with respective coverage probability of 36% and 100%. 173

We used a simple synthetic example, but for large dimensional and highly nonlinear dynamics, 174 such an underestimation or overestimation of uncertainty may have a strong effect and may cause 175 filters to collapse. The main issue in ensemble-based DA is an underdispersive spread, as in 176 Fig. 2(e). In that case, the initial condition spread is too narrow, and model forecasts (starting 177 from these conditions) would be similar and potentially out of the range of the observations. In 178 the case of an overdispersive spread, as in Fig. 2(f), the risk is that only a small portion of model 179 forecasts would be accurate enough to produce useful information on the true state of the system. 180 This illustrative example shows how important is the joint tuning of model and observation errors 181 in DA. Since the 1990s, a substantial number of studies have dealt with this topic. 182

In a seminal paper, Dee (1995) proposed an estimation method for parametric versions of  $\mathbf{Q}$ 184 and **R** matrices. The method, based on maximizing the likelihood of the observations, yields an 185 estimator which is a function of the innovation defined by  $\mathbf{v} - \mathcal{H}(\mathbf{x}^f)$ . Maximization is performed 186 at each assimilation step, with the current innovation computed from the available observations. 187 This technique was later extended to estimate the mean of the innovation, which depends on the 188 biases in the forecast and in the observations (Dee et al. 1999a). The methodology was then 189 applied to realistic cases in Dee et al. (1999b), making the maximization of innovation likelihood 190 a promising technique for the estimation of errors in operational forecasts. 191

<sup>192</sup> Following a distinct path, Desroziers and Ivanov (2001) proposed using the observation-minus-<sup>193</sup> analysis diagnostic. It is defined by  $\mathbf{y} - \mathscr{H}(\mathbf{x}^a)$  with  $\mathbf{x}^a$  the analysis (i.e., the output of DA algo-<sup>194</sup> rithms). The authors proposed an iterative optimization technique to estimate a scaling factor for <sup>195</sup> the background  $\mathbf{B} = \mathbf{P}^f$  and observation  $\mathbf{R}$  matrices. The procedure was shown to converge to a <sup>196</sup> proper fixed-point. As in Dee's work, the fixed-point method presented in Desroziers and Ivanov <sup>197</sup> (2001) is applied at each assimilation step, with the available observations at the current step.

Later, Chapnik et al. (2004) showed that the maximization of the innovation likelihood proposed by Dee (1995) makes the observation-minus-analysis diagnostic of Desroziers and Ivanov (2001) optimal. Moreover, the techniques of Dee (1995) and Desroziers and Ivanov (2001) have been further connected to the generalized cross-validation method previously developed by statisticians (Wahba and Wendelberger 1980).

These initial studies clearly nurtured the discussion of the estimation of observation **R**, model **Q**, or background  $\mathbf{B} = \mathbf{P}^{f}$  error covariance matrices in the modern DA literature. For demonstration purposes, the algorithms proposed in Dee (1995) and Desroziers and Ivanov (2001) were tested on

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realistic DA problems, using a shallow-water model on a plane with a simplified Kalman filter, and 206 using the French ARPEGE three-dimensional variational framework, respectively. In both cases, 207 although good performances have been obtained with a small number of iterations, the proposed 208 algorithms have shown some limits, in particular with regard to the simultaneous estimation of the 209 two sources of errors: observation and model (or background). In this context, Todling (2015) 210 pointed out that using only the current innovation is not enough to distinguish the impact of  $\mathbf{Q}$  and 211 **R**, which still makes their simultaneous estimation challenging. Given that our preliminary focus 212 here is to review methods for the joint estimate of Q and R, the work Dee (1995) and Desroziers 213 and Ivanov (2001) are not further detailed hereafter. After these two seminal studies, various 214 alternatives were proposed. They are based on the use of several types of innovations and are 215 discussed in this review. 216

#### 217 *d. Methods presented in this review*

The main topic of this review is the "joint estimation of **Q** and **R**". Thus, only methods based 218 on this specific goal are presented in detail. A history of what have been, in our opinion, the most 219 relevant contributions and the key milestones for Q and R covariance estimation in DA is sketched 220 in Fig. 3. The highlighted papers are discussed in this review, with a summary of the different 221 methodologies, given in Table 1. We distinguish four methods and we can classify them into 222 two categories: those which rely on moment-based methods, and those using likelihood-based 223 methods. Both methods make use of the innovations. The main concepts of the techniques are 224 briefly introduced below. 225

On the one hand, moment-based methods assume equality between theoretical and empirical statistical moments. A first approach is to study different type of innovations in the observation space (i.e., working in the space of the observations instead of the space of the state). It has

<sup>229</sup> been initiated in DA by Rutherford (1972) and Hollingsworth and Lönnberg (1986). A second
<sup>230</sup> approach extracts information from the correlation between lag innovations, namely innovations
<sup>231</sup> between consecutive times. On the other hand, likelihood-based methods aim to maximize likeli<sup>232</sup> hood functions with statistical algorithms. One option is to use a Bayesian framework, assuming
<sup>233</sup> prior distributions for the parameters of **Q** and **R** covariance matrices. Another option is to use the
<sup>234</sup> iterative expectation–maximization algorithm to maximize a likelihood function.

The four methodologies listed in Fig. 3 will be examined in this paper. Before doing that, it is worth mentioning existing review work that have attempted to summarize the methodologies in DA context and beyond.

#### 238 e. Other review papers

Other review papers on parameter estimation (including Q and R matrices) in state-space models 239 have appeared in the statistical and signal processing communities. The first one (Mehra 1972) 240 introduces moment- and likelihood-based methods in the linear and Gaussian case (i.e., when  $\eta$ 241 and  $\epsilon$  are Gaussians and  $\mathcal{M}$  is a linear operator in Eqs. (1) and (2)). Many extensions to nonlinear 242 state-space models have been proposed since the seminal work of Mehra, and these studies are 243 summarized in the recent review by Duník et al. (2017), with a focus on moment-based methods 244 and the extended Kalman filter (Jazwinski 1970). The book chapter by Buehner (2010) presents 245 another review of moment-based methods, with a focus on the modeling and estimation of spatial 246 covariance structures **Q** and **R** in DA with the ensemble Kalman filter algorithm (Evensen 2009). 247 In the statistical community, the recent development of powerful simulation techniques, known 248 as sequential Monte-Carlo algorithms or particle filters, has led to an extensive literature on the 249 statistical inference in nonlinear state-space models relying on likelihood-based approaches. A 250 recent and detailed presentation of this literature can be found in Kantas et al. (2015). However, 251

these methods typically require a large number of particles, which make them impractical for geophysical DA applications.

The review presented here focuses on methods proposed in DA, especially the moment- and likelihood-based techniques which are suitable for geophysical systems (i.e., with high dimensionality and strong nonlinearities).

#### 257 f. Structure of this review

The paper is organized as follows. Section 2 briefly presents the filtering and smoothing DA 258 algorithms used in this work. The main families of methods used in the literature to jointly 259 estimate error covariance matrices Q and R are then described. First, moment-based methods 260 are introduced in section 3. Then, we describe in section 4 the likelihood-based methods. We 261 also mention other alternatives in section 5, along with methods used in the past but not exactly 262 matching the scope of this review, and diagnostic tools to check the accuracy of  $\mathbf{Q}$  and  $\mathbf{R}$ . Finally, 263 in section 6, we provide a summary and discussion on what we consider to be the forthcoming 264 challenges in this area. 265

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#### 267 2. Filtering and smoothing algorithms

This review paper focuses on the estimation of **Q** and **R** in the context of ensemble-based DA methods. For the overall discussion of the methods and to set the notation, a short description of the ensemble version of the Kalman recursions is presented in this section: the ensemble Kalman filter (EnKF) and ensemble Kalman smoother (EnKS).

The EnKF and EnKS estimate various state vectors  $\mathbf{x}^{f}(k)$ ,  $\mathbf{x}^{a}(k)$ ,  $\mathbf{x}^{s}(k)$  and covariance matrices  $\mathbf{P}^{f}(k)$ ,  $\mathbf{P}^{a}(k)$ ,  $\mathbf{P}^{s}(k)$ , at each time step  $1 \le k \le K$ , where *K* represents the total number of assimila-

tion steps. Kalman-based algorithms assume a Gaussian prior distribution  $p(\mathbf{x}(k)|\mathbf{y}(1:k-1)) \sim \mathcal{N}(\mathbf{x}^{f}(k), \mathbf{P}^{f}(k))$ . Then, filtering and smoothing estimates correspond to the Gaussian posterior distributions  $p(\mathbf{x}(k)|\mathbf{y}(1:k)) \sim \mathcal{N}(\mathbf{x}^{a}(k), \mathbf{P}^{a}(k))$  and  $p(\mathbf{x}(k)|\mathbf{y}(1:K)) \sim \mathcal{N}(\mathbf{x}^{s}(k), \mathbf{P}^{s}(k))$  of the state conditionally to past/present observations and past/present/future observations respectively. The basic idea of the EnKF and EnKS is to use an ensemble  $\mathbf{x}_{1}, \dots, \mathbf{x}_{N_{e}}$  of size  $N_{e}$  to track Gaussian distributions over time with the empirical mean vector  $\bar{\mathbf{x}} = 1/N_{e}\sum_{i=1}^{N_{e}} \mathbf{x}_{i}$  and the empirical error covariance matrix  $1/(N_{e}-1)\sum_{i=1}^{N_{e}} (\mathbf{x}_{i}-\bar{\mathbf{x}}) (\mathbf{x}_{i}-\bar{\mathbf{x}})^{\mathrm{T}}$ .

Forecast step (forward in time):

$$\mathbf{x}_{i}^{f}(k) = \mathscr{M}_{k}\left(\mathbf{x}_{i}^{a}(k-1)\right) + \boldsymbol{\eta}_{i}(k)$$
(4a)

Analysis step (forward in time):

$$\mathbf{d}_{i}(k) = \mathbf{y}(k) - \mathscr{H}_{k}\left(\mathbf{x}_{i}^{f}(k)\right) + \varepsilon_{i}(k)$$
(4b)

$$\mathbf{K}^{f}(k) = \mathbf{P}^{f}(k) \mathscr{H}_{k}^{\mathrm{T}} \left( \mathscr{H}_{k} \mathbf{P}^{f}(k) \mathscr{H}_{k}^{\mathrm{T}} + \mathbf{R}(k) \right)^{-1}$$
(4c)

$$\mathbf{x}_{i}^{a}(k) = \mathbf{x}_{i}^{f}(k) + \mathbf{K}^{f}(k)\mathbf{d}_{i}(k)$$
(4d)

Reanalysis step (backward in time):

$$\mathbf{K}^{s}(k) = \mathbf{P}^{a}(k) \mathcal{M}_{k}^{\mathrm{T}} \left( \mathbf{P}^{f}(k+1) \right)^{-1}$$
(4e)

$$\mathbf{x}_{i}^{s}(k) = \mathbf{x}_{i}^{a}(k) + \mathbf{K}^{s}(k) \left(\mathbf{x}_{i}^{s}(k+1) - \mathbf{x}_{i}^{f}(k+1)\right)$$
(4f)

with  $\mathbf{K}^{f}(k)$  and  $\mathbf{K}^{s}(k)$  the filter and smoother Kalman gains, respectively. Here,  $\mathbf{P}^{f}(k)$  and  $\mathcal{H}_{k}\mathbf{P}^{f}(k)\mathcal{H}_{k}^{T}$  denote the empirical covariance matrices of  $\mathbf{x}_{i}^{f}(k)$  and  $\mathcal{H}_{k}(\mathbf{x}_{i}^{f}(k))$ , respectively. Then,  $\mathbf{P}^{f}(k)\mathcal{H}_{k}^{T}$  and  $\mathbf{P}^{a}(k)\mathcal{M}_{k}^{T}$  denote the empirical cross-covariance matrices between  $\mathbf{x}_{i}^{f}(k)$ and  $\mathcal{H}_{k}(\mathbf{x}_{i}^{f}(k))$  and between  $\mathbf{x}_{i}^{a}(k)$  and  $\mathcal{M}_{k}(\mathbf{x}_{i}^{a}(k))$ , respectively. These quantities are estimated using  $N_{e}$  ensemble members.

In some of the methods presented in this review, the ensembles are also used to approximate  $\mathcal{M}_k$ and  $\mathcal{H}_k$  by linear operators  $\mathbf{M}_k$  and  $\mathbf{H}_k$  such as

$$\mathbf{M}_{k} = \mathbf{E}_{k}^{\mathscr{M}(a)} (\mathbf{E}_{k-1}^{a})^{\dagger}$$
(5a)

$$\mathbf{H}_{k} = \mathbf{E}_{k}^{\mathcal{H}(f)} (\mathbf{E}_{k}^{f})^{\dagger}$$
(5b)

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with  $\dagger$  the pseudo-inverse,  $\mathbf{E}_{k}^{\mathscr{M}(a)}$ ,  $\mathbf{E}_{k-1}^{a}$ ,  $\mathbf{E}_{k}^{\mathscr{H}(f)}$  and  $\mathbf{E}_{k}^{f}$  the matrices containing along their columns the ensemble perturbation vectors (the centered ensemble vectors) of  $\mathscr{M}_{k}(\mathbf{x}_{i}^{a}(k-1))$ ,  $\mathbf{x}_{i}^{a}(k-1)$ ,  $\mathscr{H}_{k}(\mathbf{x}_{i}^{f}(k))$  and  $\mathbf{x}_{i}^{f}(k)$ , respectively.

In Eq. (4b), the innovation is denoted as **d** and tracked by  $\mathbf{d}_1(k), \dots, \mathbf{d}_{N_e}(k)$ . The innovation is the key ingredient of the methods presented in sections 3 and 4.

#### **3. Moment-based methods**

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In order to constrain the model and observational errors in DA systems, initial efforts were focused on the statistics of relevant variables which could contain information on covariances. The innovation, given in Eq. (4b), corresponds to the difference between the observations and the forecast in the observation space. This variable implicitly takes into account the **Q** and **R** covariances. Unfortunately, as explained in Blanchet et al. (1997), by using only current observations, their individual contributions cannot be easily disentangled. Thus, the techniques with only the classic innovation  $\mathbf{y}(k) - \mathscr{H}_k(\mathbf{x}^f(k))$  are not discussed further in this review.

Two main approaches have been proposed in the literature to address this issue. They are based on the idea of producing multiple equations involving  $\mathbf{Q}$  and  $\mathbf{R}$ . The first approach uses different type of innovation statistics (i.e., not only the classic one). The second approach is based on lag innovations, or differences between consecutive innovations. From a statistical point of view, they refer to the "methods of moments", where we construct a system of equations that links various moments of the innovations with the parameters and then replace the theoretical moments by the empirical ones in these equations.

#### *a. Innovation statistics in the observation space*

This first approach, based on the Desroziers diagnostic (Desroziers et al. 2005), is historical and now popular in the DA community. It does not exactly fit the topic of this review paper (i.e., estimating the model error **Q**), since it is based on the inflation of the background covariance matrix  $\mathbf{P}^{f}$ . However, this forecast error covariance is defined by  $\mathbf{P}^{f}(k) = \mathbf{M}_{k}\mathbf{P}^{a}(k-1)\mathbf{M}_{k}^{T} + \mathbf{Q}$  in the Kalman filter, considering a linear model operator  $\mathbf{M}_{k}$ . Thus, even if DA systems do not use an explicit model error perturbation controlled by  $\mathbf{Q}$ , the inflation of the background covariance matrix  $\mathbf{P}^{f}$  has similar effects, compensating for the lack of an explicit model uncertainty.

Desroziers et al. (2005) proposed examining various innovation statistics in the observation space. It is based on different type of innovation statistics between observations, forecasts and analysis, with all of them defined in the observation space: namely,  $\mathbf{d}^{o-f}(k) = \mathbf{y}(k) - \mathscr{H}_k(\mathbf{x}^f(k))$ as in Eq. (4b) and  $\mathbf{d}^{o-a}(k) = \mathbf{y}(k) - \mathscr{H}_k(\mathbf{x}^a(k))$ . In theory, in the linear and Gaussian case, for unbiased forecast and observation, and when  $\mathbf{P}^f(k)$  and  $\mathbf{R}(k)$  are correctly specified, the Desroziers innovation statistics should verify the equalities:

$$\int \mathbf{E} \left[ \mathbf{d}^{o-f}(k) \mathbf{d}^{o-f}(k)^{\mathrm{T}} \right] = \mathbf{H}_{k} \mathbf{P}^{f}(k) \mathbf{H}_{k}^{\mathrm{T}} + \mathbf{R}(k)$$
(6a)

$$\left( \mathbf{E} \left[ \mathbf{d}^{o-a}(k) \mathbf{d}^{o-f}(k)^{\mathrm{T}} \right] = \mathbf{R}(k)$$
(6b)

<sup>317</sup> with E the expectation operator. Equation (6a) is given by using Eq. (4b):

$$\mathbf{d}^{o-f}(k)\mathbf{d}^{o-f}(k)^{\mathrm{T}} = -\mathbf{y}(k)\mathbf{x}^{f}(k)^{\mathrm{T}}\mathbf{H}_{k}^{\mathrm{T}}$$
$$-\mathbf{H}_{k}\mathbf{x}^{f}(k)\mathbf{y}(k)^{\mathrm{T}}$$
$$+\mathbf{H}_{k}\mathbf{x}^{f}(k)\mathbf{x}^{f}(k)^{\mathrm{T}}\mathbf{H}_{k}^{\mathrm{T}}$$
$$+\mathbf{y}(k)\mathbf{y}(k)^{\mathrm{T}},$$
(7)

then applying the expectation operator and using the definition of  $\mathbf{P}^{f}$  and  $\mathbf{R}$ . The observationminus-forecast innovation statistics in Eq. (6a) is not useful to constrain model error  $\mathbf{Q}$ . Indeed,

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 $d^{o-f}$  does not depend explicitly on **Q**, but rather on the forecast error covariance matrix **P**<sup>*f*</sup>. Thus, the combination of Eq. (6a) and Eq. (6b) can be used as a diagnosis of the forecast and observational error covariances in the system. A mismatch between the Desroziers statistics and the actual covariances, namely the left- and right-hand side terms in Eq. (6a) and Eq. (6b), indicates inappropriate estimated covariances **P**<sup>*f*</sup>(*k*) and **R**(*k*).

The forecast covariance  $\mathbf{P}^{f}$  is sometimes badly estimated in ensemble-based assimilation sys-325 tems. The limitations may be attributed to a number of causes. The limited number of ensemble 326 members produces an over- or, most of the time, underestimation of the forecast variance. An-327 other limitation is the inaccuracies in methods used to sample initial condition or model error. The 328 underestimation of the forecast covariance produces negative feedback, and the estimated analysis 329 covariance  $\mathbf{P}^a$  is thus underestimated, which in turn produces a further underestimation of the fore-330 cast covariance in the next cycle. This feedback process leads to filter divergence, as was pointed 331 out by Pham et al. (1998), Anderson and Anderson (1999) or Anderson (2007). To avoid this 332 filter divergence, inflating the forecast covariance  $\mathbf{P}^{f}$  has been proposed. This covariance inflation 333 accounts for both sampling errors and the lack of representation of model errors, like a too small 334 amplitude for **Q** or the fact that a bias is omitted in  $\eta$  and  $\epsilon$ , Eqs. (1) and (2). In this context, the 335 diagnostics given by the Desroziers innovation statistics have been proposed as a tool to constrain 336 the required covariance inflation in the system. 337

<sup>338</sup> We distinguish three inflation methods: multiplicative, additive and relaxation-to-prior. In the <sup>339</sup> multiplicative case, the forecast error covariance matrix  $\mathbf{P}^{f}$  is usually multiplied by a scalar coeffi-<sup>340</sup> cient greater than 1 (Anderson and Anderson 1999). Using innovation statistics in the observation <sup>341</sup> space, adaptive procedures to estimate this coefficient have been proposed by Wang and Bishop <sup>342</sup> (2003), Anderson (2007), Anderson (2009) conditionally to the spatial location, Li et al. (2009), <sup>343</sup> Miyoshi (2011), Bocquet (2011), Bocquet and Sakov (2012), Miyoshi et al. (2013), Bocquet et al. (2015), El Gharamti (2018) and Raanes et al. (2019). In order to prevent excessive inflation or deflation, some authors have proposed assuming a priori distribution for the multiplicative inflation
factor. The most usual a priori distributions used by the authors are Gaussian in Anderson (2009),
inverse-gamma in El Gharamti (2018) or inverse chi-square in Raanes et al. (2019).

In practice, multiplicative inflation tends to excessively inflate in the data-sparse regions and 348 inflate too little in the densely observed regions. As a result, the spread looks like exaggeration of 349 data density (i.e., too much spread in sparsely observed regions, and vice versa). Additive inflation 350 solves this problem, but requires a lot of samples for additive noise; these drawbacks and benefits 351 are discussed in Miyoshi et al. (2010). In the additive inflation case, the diagonal terms of the 352 forecast and analysis empirical covariance matrices is increased (Mitchell and Houtekamer 2000; 353 Corazza et al. 2003; Whitaker et al. 2008; Houtekamer et al. 2009). This regularization also avoids 354 the problems corresponding to the inversion of the covariance matrices. 355

The last alternative is the relaxation-to-prior method. In application, this technique is more effi-356 cient than both additive and multiplicative inflations because it maintains a reasonable spread struc-357 ture. The idea is to relax the reduction of the spread at analysis. We distinguish the method pro-358 posed in Zhang et al. (2004), where the forecast and analysis ensemble perturbations are blended, 359 from the one given in Whitaker and Hamill (2012), which multiplies the analysis ensemble with-360 out blending perturbations. This last method is thus a multiplicative inflation, but applied after the 361 analysis, not the forecast. Finally, Ying and Zhang (2015) and Kotsuki et al. (2017b) proposed 362 methods to adaptively estimate the relaxation parameters using innovation statistics. Their con-363 clusions are that adaptive procedures for relaxation-to-prior methods are robust to sudden changes 364 in the observing networks and observation error settings. 365

<sup>366</sup> Closely connected to multiplicative inflation estimation is statistical modeling of the error vari-<sup>367</sup> ance terms proposed by Bishop and Satterfield (2013) and Bishop et al. (2013). From numerical

evidence based on the 10-dimensional Lorenz-96 model, the authors assume an inverse-gamma prior distribution for these variances. This distribution allows for an analytic Bayesian update of the variances using the innovations. Building on Bocquet (2011); Bocquet et al. (2015); Ménétrier and Auligné (2015), this technique was extended in Satterfield et al. (2018) to adaptively tune a mixing ratio between the true and sample variances. Adaptive covariance inflations are estimation methods directly attached to a traditional filtering

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373 method (such as the EnKF used here), with almost negligible overhead computational cost. In 374 practice, the use of this technique does not necessarily imply an additive error term  $\eta$  in Eq. (1). 375 Thus, it is not a direct estimation of Q but rather an inflation applied to  $P^{f}$  in order to compensate 376 for model uncertainties and sampling errors in the EnKFs, as explained in Raanes et al. (2019, 377 their section 4 and appendix C). Several DA systems work with an inflation method and use it for 378 its simplicity, low cost, and efficiency. As an example of inflation techniques, the most straight-379 forward inflation estimation is a multiplicative factor  $\lambda$  of the incorrectly scaled  $\tilde{\mathbf{P}}^{f}(k)$ , so that the 380 corrected forecast covariance is given by  $\mathbf{P}^{f}(k) = \lambda(k)\tilde{\mathbf{P}}^{f}(k)$ . The estimate of the inflation factor 381 is given by taking the trace of Eq. (6a): 382

$$\tilde{\lambda}(k) = \frac{\mathbf{d}^{o-f}(k)^{\mathrm{T}} \mathbf{d}^{o-f}(k) - \mathrm{Tr}\left(\mathbf{R}(k)\right)}{\mathrm{Tr}\left(\mathbf{H}_{k} \tilde{\mathbf{P}}^{f}(k) \mathbf{H}_{k}^{\mathrm{T}}\right)}.$$
(8)

The estimated inflation parameter  $\hat{\lambda}$  computed at each time *k* can be noisy. The use of temporal smoothing of the form  $\lambda(k+1) = \rho \tilde{\lambda}(k) + (1-\rho)\lambda(k)$  is crucial in operational procedures. Alternatively, Miyoshi (2011) proposed calculating the estimated variance of  $\lambda(k)$ , denoted as  $\sigma_{\lambda(k)}^2$ , using the central limit theorem. Then,  $\lambda(k+1)$  is updated using the previous estimate  $\lambda(k)$  and the Gaussian distribution with mean  $\tilde{\lambda}(k)$  and variance  $\sigma_{\lambda(k)}^2$ . From the Desroziers diagnostics, at each time step *k* and when sufficient observations are available, an estimate of **R**(k) is possible using Eq. (6b). For instance, Li et al. (2009) proposed estimating each component of a diagonal

and averaged **R** matrix. However, the diagonal terms cannot take into account spatial correlated error terms, and constant values for observation errors are not realistic. Then, Miyoshi et al. (2013) proposed additionally estimating the off-diagonal components of the time-dependent matrix  $\mathbf{R}(k)$ . The Miyoshi et al. (2013) implementation is summarized in the appendix, Algorithm 1.

The Desroziers diagnostic method has been applied widely to estimate the real observation error 394 covariance matrix **R** in Numerical Weather Prediction (NWP). The observations are coming from 395 different sources. In the case of satellite radiances, Bormann et al. (2010) applied three meth-396 ods, including the Desroziers diagnostic and the method detailed in Hollingsworth and Lönnberg 397 (1986) to estimate a constant diagonal term of **R** using the innovation  $\mathbf{d}^{o-f}$  and its correlations 398 in space, assuming that horizontal correlations in  $\mathbf{d}^{o-f}$  samples are purely due to  $\mathbf{P}^{f}$ . Weston 399 et al. (2014) and Campbell et al. (2017) then included the inter-channel observation error correla-400 tions of satellite radiances in DA and obtained improved results compared with the case using a 401 diagonal **R**. For spatial error correlations in **R**, Kotsuki et al. (2017a) estimated the horizontal ob-402 servation error correlations of satellite-derived precipitation data. Including horizontal observation 403 error correlations in DA for densely-observed data from satellites and radars is more challenging 404 than including inter-channel error correlations in DA. Indeed, the number of horizontally error-405 correlated observations is much larger, and some recent studies have been tackling this issue (e.g., 406 Guillet et al. (2019)). 407

To conclude, the Desroziers diagnostic is a consistency check and makes it possible to detect if the error covariances  $\mathbf{P}^{f}$  and  $\mathbf{R}$  are incorrect. When and how this method can result in accurate or inaccurate estimates, and convergence properties, have been studied in depth by Waller et al. (2016) and Ménard (2016). The Desroziers diagnostic is also useful to estimate off-diagonal terms of  $\mathbf{R}$ , for instance taking into account the spatial error correlations. However, covariance localization used in the ensemble Kalman filter might induce erroneous estimates of spatial correlations
(Waller et al. 2017).

#### 415 b. Lag innovation between consecutive times

Another way to estimate error covariances is to use multiple equations involving Q and R, 416 exploiting cross-correlations between lag innovations. More precisely, it involves the current in-417 novation  $\mathbf{d}(k) = \mathbf{d}^{o-f}(k)$  defined in Eq. (4b) and past innovations  $\mathbf{d}(k-1), \ldots, \mathbf{d}(k-l)$ . Lag 418 innovations were introduced by Mehra (1970) to recover Q and R simultaneously for Gaussian, 419 linear and stationary dynamic systems. In such a case,  $\{\mathbf{d}(k)\}_{k\geq 1}$  is completely characterized by 420 the lagged covariance matrix  $C_l = Cov(d(k), d(k-l))$ , which is independent of k. In other words, 421 the information encoded in  $\{\mathbf{d}(k)\}_{k\geq 1}$  is completely equivalent to the information provided by 422  $\{\mathbf{C}_l\}_{l>0}$ . Moreover, for linear systems in a steady state, analytic relations exist between  $\mathbf{Q}$ ,  $\mathbf{R}$  and 423  $E[\mathbf{d}(k)\mathbf{d}(k-l)^{T}]$ . However, these linear relations can be dependent and redundant for different 424 lags l. Therefore, as stated in Mehra (1970), only a limited number of Q components can be 425 recovered. 426

Bélanger (1974) extended these results to the case of time-varying linear stochastic processes, 427 taking  $\mathbf{d}(k)\mathbf{d}(k-l)^{\mathrm{T}}$  as "observations" of **Q** and **R** and using a secondary Kalman filter to update 428 them iteratively. On the one hand, considering the time-varying case may increase the number of 429 components in  $\mathbf{Q}$  that can be estimated. On the other hand, as pointed out in Bélanger (1974), 430 this method would no longer be analytically exact if  $\mathbf{Q}$  and  $\mathbf{R}$  were updated adaptively at each 431 time step. One numerical difficulty of Bélanger's method is that it needs to invert a matrix of size 432  $m^2 \times m^2$ , where m refers to the dimension of the observation vector. However, this difficulty has 433 been largely overcome by Dee et al. (1985) in which the matrix inversion is reduced to  $\mathcal{O}(m^3)$ , by 434 taking the advantage of the fact that the big matrix comes from some tensor product. 435

<sup>436</sup> More recent work have focused on high-dimensional and nonlinear systems using the extended <sup>437</sup> or ensemble Kalman filters. Berry and Sauer (2013) proposed a fast and adaptive algorithm in-<sup>438</sup> spired by the use of lag innovations proposed by Mehra. Harlim et al. (2014) applied the original <sup>439</sup> Bélanger algorithm empirically to a nonlinear system with sparse observations. Zhen and Harlim <sup>440</sup> (2015) proposed a modified version of Bélanger's method, by removing the secondary filter and <sup>441</sup> alternatively solving **Q** and **R** in a least-squares sense based on the averaged linear relation over a <sup>442</sup> long term.

Here, we briefly describe the algorithm of Berry and Sauer (2013), considering the lag-zero and lag-one innovations. The following equations are satisfied in the linear and Gaussian case, for unbiased forecast and observation when  $\mathbf{P}^{f}(k)$  and  $\mathbf{R}(k)$  are correctly specified:

$$\begin{cases} \mathbf{E} \left[ \mathbf{d}(k) \mathbf{d}(k)^{\mathrm{T}} \right] = \mathbf{H}_{k} \mathbf{P}^{f}(k) \mathbf{H}_{k}^{\mathrm{T}} + \mathbf{R}(k) = \boldsymbol{\Sigma}(k) \tag{9a} \\ \mathbf{E} \left[ \mathbf{d}(k) \mathbf{d}(k-1)^{\mathrm{T}} \right] = \mathbf{H}_{k} \mathbf{M}_{k} \mathbf{P}^{f}(k-1) \mathbf{H}_{k-1}^{\mathrm{T}} \\ -\mathbf{H}_{k} \mathbf{M}_{k} \mathbf{K}^{f}(k-1) \boldsymbol{\Sigma}(k-1). \tag{9b}$$

Equation (9a) is equivalent to Eq. (6a). Moreover, Eq. (9b) results from the fact that developing the expression of  $\mathbf{d}(k)$  using consecutively Eqs. (2), (1), (4a), and (4d), the innovation can be written as

$$\mathbf{d}(k) = \mathbf{y}(k) - \mathbf{H}_{k}\mathbf{x}^{f}(k)$$

$$= \mathbf{H}_{k}\left(\mathbf{x}(k) - \mathbf{x}^{f}(k)\right) + \boldsymbol{\epsilon}(k)$$

$$= \mathbf{H}_{k}\left(\mathbf{M}_{k}\mathbf{x}(k-1) - \mathbf{x}^{f}(k) + \boldsymbol{\eta}(k)\right) + \boldsymbol{\epsilon}(k)$$

$$= \mathbf{H}_{k}\left(\mathbf{M}_{k}\left(\mathbf{x}(k-1) - \mathbf{x}^{a}(k-1)\right) + \boldsymbol{\eta}(k)\right) + \boldsymbol{\epsilon}(k)$$

$$= \mathbf{H}_{k}\mathbf{M}_{k}\left(\mathbf{x}(k-1) - \mathbf{x}^{f}(k-1) - \mathbf{K}^{f}(k-1)\mathbf{d}(k-1)\right)$$

$$+ \mathbf{H}_{k}\boldsymbol{\eta}(k) + \boldsymbol{\epsilon}(k).$$
(10)

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$$\mathbf{H}_{k}\mathbf{M}_{k}\left(\mathbf{x}(k-1)-\mathbf{x}^{f}(k-1)\right)\mathbf{d}(k-1)^{\mathrm{T}}$$
$$-\mathbf{H}_{k}\mathbf{M}_{k}\left(\mathbf{K}^{f}(k-1)\mathbf{d}(k-1)\right)\mathbf{d}(k-1)^{\mathrm{T}}$$
$$+\mathbf{H}_{k}\boldsymbol{\eta}(k)\mathbf{d}(k-1)^{\mathrm{T}}+\boldsymbol{\epsilon}(k)\mathbf{d}(k-1)^{\mathrm{T}},$$
(11)

and assuming that the model  $\boldsymbol{\eta}$  and observation  $\boldsymbol{\epsilon}$  error noises are white and mutually uncorrelated, then  $\mathrm{E}\left[\boldsymbol{\eta}(k)\mathbf{d}(k-1)^{\mathrm{T}}\right] = 0$  and  $\mathrm{E}\left[\boldsymbol{\epsilon}(k)\mathbf{d}(k-1)^{\mathrm{T}}\right] = 0$ . Finally, developing  $\mathrm{E}\left[\mathbf{d}(k)\mathbf{d}(k-1)^{\mathrm{T}}\right]$ , Eq. (9b) is satisfied.

The algorithm in Berry and Sauer (2013) is summarized in the appendix, Algorithm 2. It is based on an adaptive estimation of  $\mathbf{Q}(k)$  and  $\mathbf{R}(k)$ , which satisfies the following relations in the linear and Gaussian case:

$$\tilde{\mathbf{P}}(k) = (\mathbf{H}_k \mathbf{M}_k)^{-1} \mathbf{d}(k) \mathbf{d}(k-1)^{\mathrm{T}} \mathbf{H}_{k-1}^{-\mathrm{T}},$$
  
+  $\mathbf{K}^f(k-1) \mathbf{d}(k-1) \mathbf{d}(k-1)^{\mathrm{T}} \mathbf{H}_{k-1}^{-\mathrm{T}}$  (12a)

$$\tilde{\mathbf{Q}}(k) = \tilde{\mathbf{P}}(k) - \mathbf{M}_{k-1} \mathbf{P}^a(k-2) \mathbf{M}_{k-1}^{\mathrm{T}},$$
(12b)

$$\tilde{\mathbf{R}}(k) = \mathbf{d}(k)\mathbf{d}(k)^{\mathrm{T}} - \mathbf{H}_{k}\mathbf{P}^{f}(k)\mathbf{H}_{k}^{\mathrm{T}}.$$
(12c)

In operational applications, when the number of observations is not equal to the number of components in state **x**, **H** is not a square matrix and Eq. (12a) is ill-defined. To avoid the inversion of **H**, Berry and Sauer (2013) proposed considering parametric models for **Q** and then solving a linear system associated with Eqs. (12a) and (12b). It is written as a least-squares problem such

457 that

$$\widetilde{\mathbf{Q}}(k) = \underset{\mathbf{Q}}{\operatorname{arg\,min}} ||\mathbf{d}(k)\mathbf{d}(k-1)^{\mathrm{T}} + \mathbf{H}_{k}\mathbf{M}_{k}\mathbf{K}^{f}(k-1)\mathbf{d}(k-1)\mathbf{d}(k-1)^{\mathrm{T}} - \mathbf{H}_{k}\mathbf{M}_{k}\mathbf{M}_{k-1}\mathbf{P}^{a}(k-2)\mathbf{M}_{k-1}^{\mathrm{T}}\mathbf{H}_{k-1}^{\mathrm{T}} - \mathbf{H}_{k}\mathbf{M}_{k}\mathbf{Q}\mathbf{H}_{k-1}^{\mathrm{T}}||.$$
(13)

In this adaptive procedure, joint estimations of  $\tilde{\mathbf{Q}}(k)$  and  $\tilde{\mathbf{R}}(k)$  can abruptly vary over time. Thus, the temporal smoothing of the covariances being estimated becomes crucial. As suggested by Berry and Sauer (2013), such temporal smoothing between current and past estimates is a reasonable choice:

$$\mathbf{Q}(k+1) = \boldsymbol{\rho} \tilde{\mathbf{Q}}(k) + (1-\boldsymbol{\rho})\mathbf{Q}(k), \qquad (14a)$$

$$\mathbf{R}(k+1) = \boldsymbol{\rho} \tilde{\mathbf{R}}(k) + (1-\boldsymbol{\rho})\mathbf{R}(k)$$
(14b)

with  $\mathbf{Q}(1)$  and  $\mathbf{R}(1)$  the initial conditions and  $\rho$  the smoothing parameter. When  $\rho$  is large (close to 1), weight is given to the current estimates  $\tilde{\mathbf{Q}}$  and  $\tilde{\mathbf{R}}$ , and when  $\rho$  is small (close to 0) it gives smoother  $\mathbf{Q}$  and  $\mathbf{R}$  sequences. The value of  $\rho$  is arbitrary and may depend on the system and how it is observed. For instance, in the case where the number of observations equals the size of the system, Berry and Sauer (2013) uses  $\rho = 5 \times 10^{-5}$  in order to estimate the full matrix  $\mathbf{Q}$  for the Lorenz-96 model.

The algorithm in Berry and Sauer (2013) only considers lag-zero and lag-one innovations. By incorporating more lags, Zhen and Harlim (2015) and Harlim (2018) showed that it makes it possible to deal with the case in which some components of Q are not identifiable from the method in Berry and Sauer (2013). For instance, let us consider the two-dimensional system with any stationary operator **M** and  $\mathbf{H} = [1,0]$ , meaning that only the first component of the system is

observed. This is a linear, Gaussian, stationary system, and Mehra's theory implies that two 473 parameters of **Q** are identifiable. However, using only lag-one innovations as in Berry and Sauer 474 (2013), Eq. (13) becomes a scalar equation and only one parameter of  $\mathbf{Q}$  can be determined. The 475 idea of considering more lag innovations to estimate more components of  $\mathbf{Q}$  was tested in Zhen 476 and Harlim (2015). Numerical results show that considering more than one lag can improve the 477 estimates of **Q** and **R**. For instance, Zhen and Harlim (2015) focused on the Lorenz-96 model. 478 Results show that when Q is stationary, the trace of Q and R are equal, and when observations are 479 taken at twenty fixed equally spaced grid points for every five integration time steps, the optimal 480 RMSE of the estimates of **Q** and **R** is achieved when four time lags are considered. But with more 481 lags, the performance is degraded. 482

To summarize, methods based on lag innovation between consecutive times have been studied for a long time in the signal processing community. The original methods (Mehra 1970; Bélanger 1974) were analytically established for linear systems with Gaussian noises. Inspired by these foundational ideas, empirical methods have been established for nonlinear systems in DA (Berry and Sauer 2013; Harlim et al. 2014; Zhen and Harlim 2015). Although these methods have not been tested in any operational experiment, the idea of using lagged innovations seems to have significant potential.

#### 490 4. Likelihood-based methods

This section focuses on methods based on the likelihood of the observations, given a set of statistical parameters. The conceptual idea behind what we refer to as likelihood-based methods is to determine the optimal statistical parameters (i.e.,  $\mathbf{Q}$  and  $\mathbf{R}$ ) that maximize the likelihood function for a given set of observations which may be distributed over time. In this way, the aim

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is to derive estimation methods that use the observations to find the most suitable, or most likely
 parameters.

Early studies in Dee (1995), Blanchet et al. (1997), Mitchell and Houtekamer (2000) and Liang et al. (2012) proposed finding the optimal  $\mathbf{Q}$  and  $\mathbf{R}$  that maximize the current innovation likelihood at time *k*. Unfortunately, if only the current observations are used, the joint estimation of  $\mathbf{Q}$  and  $\mathbf{R}$ is not well constrained (Todling 2015). To tackle this issue, several solutions have been recently proposed where the likelihood function considers observations distributed in time over several assimilation cycles.

The likelihood-based methods are broadly divided into two categories. One approach uses a Bayesian framework. It assumes a priori knowledge about the parameters and estimate jointly the posterior distribution of  $\mathbf{Q}$  and  $\mathbf{R}$  together with the state of the system, or alternatively to estimate them in a two-stage process<sup>2</sup>. The second one is based on the frequentist viewpoint and attempts a point estimate of the parameters by maximizing a total likelihood function.

#### 508 a. Bayesian inference

In the Bayesian framework, the elements of the covariance matrices  $\mathbf{Q}$  and  $\mathbf{R}$  are assumed to have a priori distributions which are controlled by hyperparameters. In practice, it is difficult to have prior distributions for each element of  $\mathbf{Q}$  and  $\mathbf{R}$ , especially for large DA systems. Instead, parametric forms are used for the matrices, typically describing the shape and level noise. We denote the corresponding parameters as  $\boldsymbol{\theta}$ .

<sup>&</sup>lt;sup>2</sup>Some of the methods presented in section 3 also use the Bayesian philosophy; for instance they assume a priori distribution for the multiplicative inflation parameter  $\lambda$  (Anderson 2009; El Gharamti 2018).

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The inference in the Bayesian framework aims to determine the posterior density  $p(\theta|\mathbf{y}(1:k))$ . Two techniques have appeared, the first based on a state augmentation and the second based on a rigorous Bayesian update of the posterior distribution.

#### 517 1) STATE AUGMENTATION

In the Bayesian framework,  $\theta$  is a random variable such that the state is augmented with these parameters by defining  $\mathbf{z}(k) = (\mathbf{x}(k), \theta)$ . To define an augmented state-space model, one has to define an evolution equation for the parameters. This leads to a new state-space model of the form of Eqs. (1) and (2) with  $\mathbf{x}$  replaced by  $\mathbf{z}$ . Therefore, the state and the parameters are estimated jointly using the DA algorithms.

State augmentation was first proposed in Schmidt (1966) and is known as the Schmidt-Kalman 523 filter. This technique was mainly used to estimate both the state of the system and additional pa-524 rameters, including bias, forcing terms and physical parameters. These kinds of parameters are 525 strongly related to the state of the system (Ruiz et al. 2013a). Therefore, they are identifiable 526 and suitable for an augmented state approach. However, Stroud and Bengtsson (2007) and later 527 Delsole and Yang (2010) formally demonstrated that augmentation methods fail for variance pa-528 rameters like Q and R. The explanation is that in the EnKF, the empirical forecast covariance  $\mathbf{P}^{f}$ 529 is computed using all the ensemble members, each one with a different realization of the random 530 variable  $\theta$ . Thus,  $\mathbf{P}^{f}$  and consequently the Kalman gain  $\mathbf{K}^{f}$ , are mixing the effects of  $\mathbf{Q}$  and  $\mathbf{R}$ 531 parameters contained in  $\theta$ . Therefore, after applying Eq. (4d), the update of z corresponding to 532 the  $\theta$  parameters is the same for all the parameters. To capture the impact of a single variance 533 parameter on the prediction covariance and circumvent the limitation of the state augmentation, 534 Scheffler et al. (2019) proposed to use an ensemble of states integrated with the same variance 535 parameter. The choice of an ensemble of states for each variance parameter leads to two nested 536

ensemble Kalman filters. The technique performs successfully under different model error covari ance structures but has an important computational cost.

Another critical aspect of state augmentation is that one needs to define an evolution model for 539 the augmented state  $\mathbf{z}(k) = (\mathbf{x}(k), \boldsymbol{\theta}(k))$ . If persistence is assumed in the parameters such that they 540 are constant in time, this leads to filter degeneracy, since the estimated variance of the error in  $\theta$ 541 is bound to decrease in time. To prevent or at least mitigate this issue, it was suggested to use an 542 independent inflation factor on the parameters (Ruiz et al. 2013b) or to impose artificial stochastic 543 dynamics for  $\theta$ , typically a random walk or AR(1) model, as introduced in Eq. (3) and proposed 544 in Liu and West (2001). The tuning of the parameters introduced in these artificial dynamics may 545 be difficult, and this introduces bias into the procedure, which is hard to quantify. 546

#### 547 2) BAYESIAN UPDATE OF THE POSTERIOR DISTRIBUTION

Instead of the inference of the joint posterior density using a state augmentation strategy, the state  $\mathbf{x}(k)$  and parameters  $\boldsymbol{\theta}$  can be divided into a two-step inference procedure using the following formula:

$$p(\mathbf{x}(k), \boldsymbol{\theta} | \mathbf{y}(1:k)) =$$

$$p(\mathbf{x}(k) | \mathbf{y}(1:k), \boldsymbol{\theta}) p(\boldsymbol{\theta} | \mathbf{y}(1:k)), \qquad (15)$$

which is a direct consequence of the conditional density definition. In Eq. (15),  $p(\mathbf{x}(k)|\mathbf{y}(1:k), \theta)$ represents the posterior distribution of the state, given the observations and the parameter  $\theta$ . It can be computed using a filtering DA algorithm. The second term on the right-hand side of Eq. (15) corresponds to the posterior distribution of the parameters, given the observations up to time *k*.

<sup>555</sup> The latter can be updated sequentially using the following Bayesian hierarchy:

$$p(\boldsymbol{\theta}|\mathbf{y}(1:k)) \propto$$

$$p(\mathbf{y}(k)|\mathbf{y}(1:k-1),\boldsymbol{\theta}) p(\boldsymbol{\theta}|\mathbf{y}(1:k-1)), \qquad (16)$$

where  $p(\mathbf{y}(k)|\mathbf{y}(1:k-1),\boldsymbol{\theta})$  is the likelihood of the innovations.

<sup>557</sup> Different approximations have been used for  $p(\theta|\mathbf{y}(1:k))$  in Eq. (16); these include parametric <sup>558</sup> models based on Gaussian (Stroud et al. 2018), inverse-gamma (Stroud and Bengtsson 2007) or <sup>559</sup> Wishart distributions (Ueno and Nakamura 2016), particle-based approximations (Frei and Künsch <sup>560</sup> 2012; Stroud et al. 2018) and grid-based approximation (Stroud et al. 2018).

The methods proposed in the literature also differ by the approximation used for the likelihood 561 of the innovations. We emphasize that  $p(\mathbf{y}(k)|\mathbf{y}(1:k-1),\boldsymbol{\theta})$  needs to be evaluated for different 562 values of  $\theta$  at each time step, and that this requires applying the filter from the initial time with 563 a single value of  $\theta$ , which is computationally impossible for applications in high dimensions. To 564 reduce computational time, it is generally assumed that  $\mathbf{x}^{f}$  and  $\mathbf{P}^{f}$  are independent of  $\boldsymbol{\theta}$ , and only 565 observations y(k-l:k-1) in a small time window from the current observation are used when 566 computing the likelihood of the innovations (see Ueno and Nakamura (2016); Stroud et al. (2018) 567 for a more detailed discussion). A summary of the Bayesian method from Stroud et al. (2018) is 568 given in the appendix, Algorithm 3. It was implemented within the EnKF framework and is one 569 of the most recent studies based on the Bayesian approach. 570

Applications of the Bayesian methodology in the DA context are now discussed. It has mainly been used to estimate shape and noise parameters of  $\mathbf{Q}$  and  $\mathbf{R}$  error covariance matrices. For instance, Purser and Parrish (2003) and Solonen et al. (2014) estimated statistical parameters controlling the magnitude of the variance and the spatial dependencies in the model error  $\mathbf{Q}$ , assuming that  $\mathbf{R}$  is known. There are also applications aimed at estimating parameters governing the shape

of the observation error covariance matrix **R** only: Frei and Künsch (2012) and Stroud et al. (2018) in the Lorenz-96 system, Winiarek et al. (2012, 2014) for the inversion of the source term of airborne radionuclides using a regional atmospheric model, and Ueno and Nakamura (2016) using a shallow-water model to assimilate satellite altimetry.

As pointed out in Stroud and Bengtsson (2007), Bayesian update algorithms work best when the 580 number of unknown parameters in  $\theta$  is small. This limitation may explain why the joint estimation 581 of parameters controlling both model and observation error covariances is not systematically ad-582 dressed. For instance, Stroud and Bengtsson (2007) used the EnKF with the Lorenz-96 model for 583 the estimation of a common multiplicative scalar parameter for predefined matrices  $\mathbf{Q}$  and  $\mathbf{R}$ . Al-584 ternatively, Stroud et al. (2018) tested the Bayesian method on different spatio-temporal systems 585 to estimate the signal-to-noise ratio between Q and R. Nevertheless, based on the experiments 586 about the importance of the signal-to-noise ratio  $\|\mathbf{P}^f\| / \|\mathbf{R}\|$  presented in Fig. 2, we know that this 587 estimation of the ratio is not optimal. 588

<sup>589</sup> Widely used in the statistical community, the Bayesian framework is useful incorporating phys-<sup>590</sup> ical knowledge about error covariance matrices and constraining their estimation process. In the <sup>591</sup> DA literature, authors have used a priori distributions for the shape and noise parameters of Q<sup>592</sup> or **R**, but rarely both. Operationally, only a limited number of parameters can be estimated. To <sup>593</sup> address this issue, Stroud and Bengtsson (2007) suggested combining Bayesian algorithms with <sup>594</sup> other techniques.

#### <sup>595</sup> *b. Maximization of the total likelihood.*

The innovation likelihood at time k,  $p(\mathbf{y}(k)|\mathbf{y}(1:k-1),\boldsymbol{\theta})$  in Eq. (16), can be maximized to find the optimal  $\boldsymbol{\theta}$  (i.e.,  $\mathbf{Q}$  and  $\mathbf{R}$  matrices or parameterizations of them). In practice, when this maximization is done at each time step, two issues arise. Firstly, the innovation covariance matrix

 $\Sigma(k) = \mathbf{H}_k \mathbf{P}^f(k) \mathbf{H}_k^{\mathrm{T}} + \mathbf{R}(k)$  combines the information about **R** and **Q**, the latter being contained 599 in  $\mathbf{P}^{f}$ . When using only time k, it is difficult to disentangle the model and observation error 600 covariances; in application, the aforementioned studies only estimated one of them. Secondly, 601 the number of observations at each time step is in general limited and, as pointed out by Dee 602 (1995), available observations should exceed "the number of tunable parameters by two or three 603 orders of magnitude". To overcome these limitations, a reasonable alternative is to use a batch of 604 observations within a time window and to assume  $\theta$  to be constant in time. The resulting total 605 likelihood expressed sequentially through conditioning is given by 606

$$p(\mathbf{y}(1:K)|\boldsymbol{\theta}) = \prod_{k=1}^{K} p(\mathbf{y}(k)|\mathbf{y}(1:k-1),\boldsymbol{\theta}).$$
(17)

Because it is an integration of innovation likelihoods over a long period of time from k = 1 to k =607 K, Eq. (17) provides more observational information to estimate Q and R. The maximization of 608 this total likelihood has been applied for the estimation of deterministic and stochastic parameters 609 (related to  $\mathbf{Q}$ ) using a direct sequential optimization procedure (Delsole and Yang 2010). Ueno 610 et al. (2010) used a grid-based procedure to estimate noise levels and spatial correlation lengths of 611 **Q** and a noise level for **R**. This grid-based method uses predefined sets of covariance parameters 612 and evaluates the different combinations to find the one that maximizes the likelihood criterion. 613 Brankart et al. (2010) also proposed a method using the same criterion but adding (at the initial 614 time) information on scale and correlation length parameters of  $\mathbf{Q}$  and  $\mathbf{R}$ . This information is only 615 given the first time, and is progressively forgotten over time, using a decreasing exponential factor. 616 The marginalization of the hidden state in Eq. (17) considers all the previous observations, and it 617 requires the use of a filter. The maximization of the total likelihood  $p(\mathbf{y}(1:K)|\boldsymbol{\theta})$  to estimate 618 model error covariance  $\mathbf{Q}$  was conducted in Pulido et al. (2018), where they used a gradient-based 619 optimization technique and the EnKF. 620

The likelihood function given in Eq. (17) only depends on the observations **y**. This likelihood can be written in a different way, taking into account both the observations and the hidden state **x**. Indeed, the marginalization of the hidden state to obtain the total likelihood can be produced using the whole trajectory of the state from k = 0 to the last time step *K* all at once. It is given by

$$p(\mathbf{y}(1:K)|\boldsymbol{\theta}) = \int p(\mathbf{x}(0:K), \mathbf{y}(1:K)|\boldsymbol{\theta}) d\mathbf{x}(0:K).$$
(18)

<sup>625</sup> The maximization of the total likelihood as a function of statistical parameters  $\theta$  is not possible, <sup>626</sup> since the total likelihood cannot be evaluated directly, nor its gradient with regard to the parameters <sup>627</sup> (Pulido et al. 2018). Shumway and Stoffer (1982) proposed using an iterative procedure based on <sup>628</sup> the expectation–maximization algorithm (hereinafter denoted as EM). They applied it to estimate <sup>629</sup> the parameters of a linear state-space model, with linear dynamics, and a linear observational <sup>630</sup> operator and Gaussian errors. The EM algorithm was introduced by Dempster et al. (1977).

Each iteration of the EM algorithm consists of two steps. In the expectation step (E-step), the posterior density  $p(\mathbf{x}(0:K)|\mathbf{y}(1:K), \boldsymbol{\theta}_{(n)})$  is determined conditioned on the batch of observations  $\mathbf{y}(1:K)$  and given the parameters  $\boldsymbol{\theta}_{(n)} = (\mathbf{Q}_{(n)}, \mathbf{R}_{(n)})$  from the previous iteration or initial guess. This is obtained through the application of a smoother like the EnKS. Then, the M-step relies on the maximization of an intermediate function, depending on the posterior density obtained in the E-step. The intermediate function is defined by the conditional expectation

$$\mathbf{E}\left[\log\left(p(\mathbf{x}(0:K),\mathbf{y}(1:K)|\boldsymbol{\theta})\right)|\mathbf{y}(1:K),\boldsymbol{\theta}_{(n)}\right].$$
(19)

<sup>637</sup> If as in Eqs. (1) and (2) the observational and model errors are assumed to be additive, unbiased <sup>638</sup> and Gaussian, the expression for the logarithm of the joint density in Eq. (19) is given by

$$-\frac{1}{2} \{ \sum_{k=1}^{K} \|\mathbf{x}(k) - \mathcal{M}(\mathbf{x}(k-1))\|_{\mathbf{Q}}^{2} + \log |\mathbf{Q}| + \|\mathbf{y}(k) - \mathcal{H}(\mathbf{x}(k))\|_{\mathbf{R}}^{2} + \log |\mathbf{R}| \} + c$$

$$(20)$$

(21a) (21b) r straightforward. is first guess and

where  $\|\mathbf{v}\|_{\mathbf{A}}^2$  is defined to be equal to  $\mathbf{v}^T \mathbf{A}^{-1} \mathbf{v}$  and *c* is a constant independent of  $\mathbf{Q}$  and  $\mathbf{R}$ . In this case, an analytic expression for the optimal error covariances at each iteration of the EM algorithm can be obtained. The estimators of the parameters that maximize Eq. (19) using Eq. (20) are

$$\mathbf{Q}_{(n+1)} = \frac{1}{K} \sum_{k=1}^{K} \mathrm{E}[(\mathbf{x}(k) - \mathscr{M}(\mathbf{x}(k-1))))$$
$$(\mathbf{x}(k) - \mathscr{M}(\mathbf{x}(k-1)))^{\mathrm{T}} | \mathbf{y}(1:K), \boldsymbol{\theta}_{(n)}]$$
(21a)

642 and

$$\mathbf{R}_{(n+1)} = \frac{1}{K} \sum_{k=1}^{K} \mathbf{E}[(\mathbf{y}(k) - \mathscr{H}(\mathbf{x}(k))) \\ (\mathbf{y}(k) - \mathscr{H}(\mathbf{x}(k)))^{\mathrm{T}} | \mathbf{y}(1:K), \boldsymbol{\theta}_{(n)}].$$
(21b)

The application of the EM algorithm for the estimation of  $\mathbf{Q}$  and  $\mathbf{R}$  is rather straightforward. 643 Starting from  $\mathbf{Q}_{(1)}$  and  $\mathbf{R}_{(1)}$ , an ensemble Kalman smoother is applied with this first guess and 644 the batch of observations  $\mathbf{y}(1:K)$  to obtain the posterior density  $p(\mathbf{x}(0:K)|\mathbf{y}(1:K),\boldsymbol{\theta}_{(1)})$ . Then 645 Eqs. (21a) and (21b) are used to update and obtain  $\mathbf{Q}_{(2)}$  and  $\mathbf{R}_{(2)}$ . Next, a new application of 646 the smoother is conducted using the parameters  $\mathbf{Q}_{(2)}$  and  $\mathbf{R}_{(2)}$  and the observations  $\mathbf{y}(1:K)$ , the 647 new resulting states are used in Eqs. (21a) and (21b) to estimate  $Q_{(3)}$  and  $R_{(3)}$ , and so on. As 648 a diagnostic of convergence or as a stop criterion, the product of innovation likelihood functions 649 given in Eq. (17) is evaluated using a filter. The EM algorithm guarantees that the total likelihood 650 increases in each iteration and that the sequence  $\theta_{(n)}$  converges to a local maximum (Wu 1983). 651 A summary of the EM method (using EnKF and EnKS) from Dreano et al. (2017) is given in the 652 appendix, Algorithm 4. 653

<sup>654</sup> EM is a well-known algorithm used in the statistical community. This procedure is parameter-<sup>655</sup> free and robust, due to the large number of observations used to approximate the likelihood when <sup>656</sup> using a long batch period (Shumway and Stoffer 1982). Although the use of the EM algorithm is

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still limited in DA, it is becoming more and more popular. Some studies have implemented the EM 657 algorithm for estimating only the observation error matrix **R**. For instance, Ueno and Nakamura 658 (2014) used the model proposed in Zebiak and Cane (1987) and satellite altimetry observations, 659 whereas Liu et al. (2017) used an air quality model for accidental pollutant source retrieval. But 660 the estimation of only the observation error covariance is limited, and other studies have tried 661 to jointly estimate model error  $\mathbf{Q}$  and  $\mathbf{R}$  matrices, for instance as in Tandeo et al. (2015) for an 662 orographic subgrid-scale nonlinear observation operator. Then, Dreano et al. (2017) and Pulido 663 et al. (2018) used the EM procedure to produce joint estimation of  $\mathbf{Q}$  and  $\mathbf{R}$  matrices in the Lorenz-664 63 and stochastic parameters of the Lorenz-96 systems, respectively. Recently, Yang and Mémin 665 (2019) extended the EM procedure for the estimation of physical parameters in a one-dimensional 666 shallow water model, more specifically for the identification of stochastic subgrid terms. Lastly, 667 an online adaptation of the EM algorithm for the estimation of  $\mathbf{Q}$  and  $\mathbf{R}$  at each time step, after the 668 filtering procedure, has been proposed in Cocucci et al. (2020). In this adaptive case, the likelihood 669 is averaged locally over time, see Cappé (2011) for more details. 670

<sup>671</sup> To our knowledge, EM has not been tested yet on operational systems with large observation-<sup>672</sup> and state-space. In that case, the use of parametric forms for the matrices **Q** and **R** is essential to <sup>673</sup> reduce the number of statistical parameters  $\theta$  to estimate. For instance, Dreano et al. (2017) and <sup>674</sup> Liu et al. (2017) showed that in the particular cases where covariances are diagonal or of the form <sup>675</sup>  $\alpha$ A with A a positive definite matrix, expressions in Eq. (21a) and Eq. (21b) are simplified, and a <sup>676</sup> suboptimal  $\theta$  in the space of the parametric covariance form can be obtained.

#### **5.** Other methods

In this section, we describe other methods that have been used to estimate Q and R, and that cannot be included in the categories presented in the previous sections. In particular, we report

here about methods that are applied either a posteriori, after DA cycles, or without applying any
 DA algorithms.

#### 682 a. Analysis (or reanalysis) increment approach

This first method is based on previous DA outputs. The key idea here is to use the analysis (or reanalysis) increments to provide a realistic sample of model errors from which statistical moments, such as the covariance matrix  $\mathbf{Q}$ , can be empirically estimated. This assumes that the sequence of reanalysis  $\mathbf{x}^{s}$  (or analysis  $\mathbf{x}^{a}$ ) is the best available representation of the true process  $\mathbf{x}$ . In that case, the following approximation in Eq. (1) is made:

$$\boldsymbol{\eta}(k) = \mathcal{M}\left(\mathbf{x}(k-1)\right) - \mathbf{x}(k)$$
$$\approx \mathcal{M}\left(\mathbf{x}^{s}(k-1)\right) - \mathbf{x}^{s}(k).$$
(22)

In this approximation, it is implicitly assumed that the estimated state is the truth, so that the initial condition at time k - 1 is neglected. A similar approximation of the true process by  $\mathbf{x}^a$  or  $\mathbf{x}^s$  in Eq. (2) can be used to estimate the observation error covariance matrix **R**.

<sup>691</sup> Operationally, the analysis (or reanalysis) increment method is applied after a DA filter (or <sup>692</sup> smoother) to estimate the **Q** matrix. This method was originally introduced by Leith (1978), and <sup>693</sup> later used to account for model error in the context of ensemble Kalman filters, using analysis and <sup>694</sup> reanalysis increments by Mitchell and Carrassi (2015), and in the context of weak-constraint vari-<sup>695</sup> ational assimilation by Bowler (2017). Along this line, Rodwell and Palmer (2007) also proposed <sup>696</sup> evaluating the average of instantaneous analysis increments to represent the systematic forecast <sup>697</sup> tendencies of a model.

#### <sup>698</sup> b. Covariance matching

The covariance matching method was introduced by Fu et al. (1993). It involves matching sample covariance matrices to their theoretical expectations. Thus, it is a method of moments, similar to the work in Desroziers et al. (2005), except that covariance matching is performed on a set of historical observations and numerical simulations (noted  $\mathbf{x}^{sim}$ ), without applying any DA algorithms. It has been extended by Menemenlis and Chechelnitsky (2000) to time-lagged innovations, as first considered in Bélanger (1974).

In the case of a constant and linear observation operator **H**, the basic idea in Fu et al. (1993) is to assume the following system

$$\mathbf{x}^{sim}(k) = \mathbf{x}(k) + \boldsymbol{\eta}^{sim}(k), \qquad (23a)$$

$$\boldsymbol{\eta}^{sim}(k) = \mathbf{A}\boldsymbol{\eta}^{sim}(k-1) + \boldsymbol{\eta}(k), \qquad (23b)$$

$$\mathbf{H}\mathbf{x}^{sim}(k) - \mathbf{y}(k) = \mathbf{H}\boldsymbol{\eta}^{sim}(k) + \boldsymbol{\epsilon}(k), \qquad (23c)$$

with **A** a transition matrix close to the identity matrix, assuming slow variations of the numerical simulation errors (noted  $\eta^{sim}$ ). In Eq. (23b) and Eq. (23c), the definitions of  $\eta$  and  $\epsilon$  errors remain similar, as in the general Eqs. (1) and (2).

Assuming that **Q** and **R** are constant over time,  $\epsilon$  is uncorrelated from **x** and from  $\eta^{sim}$ , then Eq. (23c) and Eq. (23a) yield to the following estimates of **R** and **P**<sup>sim</sup> (the latter represents the error covariance of the numerical simulations):

$$\widehat{\mathbf{R}} = \frac{1}{2} \{ \mathbf{E}[(\mathbf{y} - \mathbf{H}\mathbf{x}^{sim})(\mathbf{y} - \mathbf{H}\mathbf{x}^{sim})^{\mathrm{T}}] - \mathbf{E}[(\mathbf{H}\mathbf{x}^{sim})(\mathbf{H}\mathbf{x}^{sim})^{\mathrm{T}}] + \mathbf{E}[\mathbf{y}\mathbf{y}^{\mathrm{T}}] \}, \qquad (24a)$$
$$\mathbf{H}\widehat{\mathbf{P}}^{sim}\mathbf{H}^{\mathrm{T}} = \frac{1}{2} \{ \mathbf{E}[(\mathbf{y} - \mathbf{H}\mathbf{x}^{sim})(\mathbf{y} - \mathbf{H}\mathbf{x}^{sim})^{\mathrm{T}}] + \mathbf{E}[(\mathbf{H}\mathbf{x}^{sim})(\mathbf{H}\mathbf{x}^{sim})^{\mathrm{T}}] - \mathbf{E}[\mathbf{y}\mathbf{y}^{\mathrm{T}}] \}. \qquad (24b)$$

#### 713 c. Forecast sensitivity

In operational meteorology, it is critical to learn the sensitivity of the forecast accuracy to various 714 parameters of a DA system, in particular the error statistics of both the model and the observations. 715 This is why a significant portion of literature considers the tuning problem of **R** and **Q** through the 716 lens of the sensitivity of the forecast to these parameters. The computation of those sensitivities can 717 be seen as a first-order correction or diagnostic for such an estimation. The forecast sensitivities are 718 computed either using the adjoint model (Daescu and Todling 2010; Daescu and Langland 2013) 719 in the context of variational methods, or a forecast ensemble (Hotta et al. 2017) in the context of 720 the EnKF. 721

The basic idea is to compute at each assimilation cycle an innovation between forecast and analysis, noted  $\mathbf{d}^{f-a}(k) = \mathbf{x}^{f}(k) - \mathbf{x}^{a}(k)$ . Then, the forecast sensitivity is given by  $\mathbf{d}^{f-a}(k)^{\mathrm{T}}\mathbf{S}\mathbf{d}^{f-a}(k)$ with **S** a diagonal scaling matrix, to normalize the components of  $\mathbf{d}^{f-a}$ . **Q** and **R** estimates are the matrices that minimize  $\mathbf{d}^{f-a}(k)$ . The adjoint or the ensemble are thus used to compute the partial derivatives of this forecast sensitivity. w.r.t. **Q** and **R**.

#### 727 6. Conclusions and perspectives

As often considered in data assimilation, this review paper also deals with model and observation errors that are assumed additive and Gaussian with covariance matrices  $\mathbf{Q}$  and  $\mathbf{R}$ . The model error corresponds to the dynamic model deficiencies to represent the underlying physics, whereas the observation error corresponds to the instrumental noise and the representativity error. Model and

observation errors are assumed to be uncorrelated and white in time. The model and observations
 are also assumed unbiased, a strong assumption for real data assimilation applications.

The discussion starts with the aid of an illustration of the individual and joint impacts of im-734 properly calibrated covariances using a linear toy model. The experiments clearly showed that 735 to achieve reasonable filter accuracy (i.e., in terms of root mean squared error), it is crucial to 736 carefully define both Q and R. The effect on the coverage probability of a mis-specification of 737 **Q** or **R** is also highlighted. This coverage probability is related to the estimated covariance of 738 the reconstructed state, and thus to the uncertainty quantification in data assimilation. After the 739 one-dimensional illustration, the core of the paper gives an overview of various methods to jointly 740 estimate the **Q** and **R** error covariance matrices: they are summarized and compared below. 741

#### <sup>742</sup> a. Comparison of existing methods for estimating $\mathbf{Q}$ and $\mathbf{R}$

We mainly focused in this review on four methodologies for the joint estimation of the error co-743 variances Q and R. The methods are summarized in Table 1. They correspond to classic estimation 744 methods, based on statistical moments or likelihoods. The main difference between the four meth-745 ods comes from the innovations taken into account: the total innovation, as in the EM algorithm 746 proposed by Shumway and Stoffer (1982); lag innovations, following the idea given in Mehra 747 (1970); or different type of innovations in the observation space, as in Desroziers et al. (2005). 748 Additionally, to constrain the estimation, hierarchical Bayesian approaches use prior distributions 749 for the shape parameters of **Q** and **R**. 750

<sup>751</sup> Most of the methods estimate the model error  $\mathbf{Q}$ . The exception is the one using the Desroziers <sup>752</sup> diagnostic, dealing with different type of innovations in the observation space, which instead esti-<sup>753</sup> mates an inflation factor for  $\mathbf{P}^{f}$ . Moreover, the methods are mainly defined online, meaning that <sup>754</sup> they aim to estimate  $\mathbf{Q}$  and  $\mathbf{R}$  adaptively, together with the current state of the system. Conse-

quently, these methods require additional tunable parameters to smooth the estimated covariances over time. However, most of the methods presented in this review also have an offline variant. In that case, a batch of observations is used to estimate Q and R. In some methods, such as the EM algorithm, the parameters are determined iteratively. These offline approaches avoid the use of additional smoothing parameters.

Throughout this review paper, as usually stated in DA, it is assumed that model error  $\eta$  and 760 observation error  $\epsilon$ , defined in Eqs. (1) and (2), are Gaussian. Consequently, the distribution of the 761 innovations are also Gaussian. The four presented methods use this property to build estimates of 762 **Q** and **R** adequately. But, if  $\eta$  and  $\epsilon$  are non-Gaussian, Desroziers diagnostic and lag-innovation 763 methods are not suitable anymore. However, the EM procedures and Bayesian methods are still 764 relevant, although they must be used with an appropriate filter (e.g., particle filters), not Kalman-765 based algorithms (i.e., assuming a Gaussian distribution of the state). Recently, the treatment of 766 non-Gaussian error distributions in DA has been explored in Katzfuss et al. (2019), using hierarchi-767 cal state-space models. This Bayesian framework allows to handle unknown variables that cannot 768 be easily included in the state vector (e.g., parameters of  $\mathbf{Q}$  and  $\mathbf{R}$ ) and to model non-Gaussian 769 observations. 770

The four methods have been applied at different levels of complexity. For instance, Bayesian inference methods (due to their algorithm complexity) and the EM algorithm (due to its computational cost) have so far only been applied to small dynamic models. However, the online version of the EM algorithm is less consuming and opens new perspectives of applications on larger models. On the other hand, methods using innovation statistics in the observation space have already been applied to NWP models.

The four methods summarized in Table 1 show differences in maturity in terms of applications and methodological aspects. This review also shows that there are still remaining challenges and possible improvements for the four methods.

#### <sup>780</sup> b. Remaining challenges for each method

The first challenge concerns the improvements of adaptive techniques regarding additional pa-781 rameters that control the variations of  $\mathbf{Q}$  and  $\mathbf{R}$  estimates over time. Instead of using fixed values 782 for these parameters, for instance fixed  $\rho$  in the lag innovations or  $\sigma_{\lambda}^2$  in the inflation methods, 783 we suggest using time-dependent adaptations. This adaptive solution could avoid the problems 784 of instabilities close to the solution. Another option could be to adapt these procedures, working 785 with stable parameter values (small  $\rho$ , low  $\sigma_{\lambda}^2$ ) and iterating the procedures on a batch of obser-786 vations, as in the EM algorithm. This offline variant was suggested and tested in Desroziers et al. 787 (2005) with encouraging results. To the best of our knowledge, it has not yet been tested with 788 lag-innovation methods. 789

The second challenge concerns considering time-varying error covariance matrices. The adap-790 tive procedures, based on online estimations with temporal smoothing of Q and R, are supposed 791 to capture slowly evolving covariances. On the contrary, offline methods like the EM algorithm 792 are working on a batch of observations, assuming that **Q** and **R** are constant over the batch period. 793 Online solutions for the EM algorithm, with the likelihood averaged locally over time (Cocucci 794 et al. 2020), could also capture slow evolution of the covariances. Another simple solution could 795 be to work on small sets of observations, named as mini-batches, and to apply the EM algorithm 796 in each set using the previous estimates as an initial guess. These intermediate schemes are of 797 common use in machine learning. 798

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A third challenge has to do with the assumption, used by all of the methods described herein, that 799 observation and model errors are mutually independent. Nevertheless, as pointed out in Berry and 800 Sauer (2018), observation and model error are often correlated in real data assimilation problems 801 (e.g., for satellite retrieval of Earth observations that uses model outputs in the inversion process). 802 Methods based on Bayesian inference can, in principle, exploit existing model-to-observation cor-803 relations by using a prior joint distribution (i.e., not two individual ones). The explicit taking into 804 account of this correlation can then constrain the optimization procedure. This is not possible in 805 the other approaches described in this review, at least not in their standard known formulations, 806 and the presence of model-observation correlation can deteriorate their accuracy. 807

A fourth challenge is common to all the methods presented in this review. Iterative versions 808 of the presented algorithms need initial values or distributions for **R** and **Q** (or  $\mathbf{B} = \mathbf{P}^{f}$  in the 809 case of Desroziers). But, as mentioned in Waller et al. (2016) for the Desrorziers diagnostics, 810 there is no guarantee that the algorithms will converge to the optimal solution. Indeed, in such 811 an optimization problem, there are possibly several local and non-optimal solutions. Suboptimal 812 specifications of **R**, **Q**, or **B** in the initial DA cycle will affect the final estimation results. There 813 are several solutions to avoid this convergence problem: initialize the covariance matrices using 814 physical expertise, execute the iterative algorithms several times with different initial covariance 815 matrices, or use stochastic perturbations in the optimization algorithms to avoid to be trapped in 816 local solutions. These aspects of convergence and sensitivity to initial conditions have so far been 817 poorly addressed. It is therefore necessary to check which method is robust operationally. 818

The last remaining challenge concerns the estimation of other statistical parameters of the statespace model given in Eqs. (1) and (2) and associated filters. Indeed, the initial conditions  $\mathbf{x}(0)$  and  $\mathbf{P}(0)$  are crucial for certain satellite retrieval problems and have to be estimated. This is the case, for instance, when the time sequence of observations is short (i.e., shorter than the spinup time

of the filter with an uninformative prior) or when filtering and smoothing are repeated on various 823 iterations, as in the EM algorithm. Estimation methods should also consider the estimation of sys-824 tematic or time-varying biases, the deterministic part of  $\eta$  and  $\epsilon$ . This was initially proposed by 825 Dee et al. (1999a) and tested in Dee et al. (1999b) in the case of maximizing the innovation like-826 lihood, in Dee (2005) in a state augmentation formulation, and was adapted to a Bayesian update 827 formulation in Liu et al. (2017) and in Berry and Harlim (2017). Recently, the joint estimation of 828 bias and covariance error terms, for the treatment of brightness temperatures from the European 829 geostationary satellite, has been successfully applied in Merchant et al. (2020). 830

#### <sup>831</sup> c. Perspectives for geophysical DA

Beyond the aforementioned potential improvements in the existing techniques, specific research 832 directions need to be taken by the data assimilation community. The main one concerns the real-833 ization of a comprehensive numerical evaluation of the different methods for the estimation of Q 834 and **R**, built on an agreed experimental framework and a consensus model. Such an effort would 835 help to evaluate (i) the pros and cons of the different methods (including their capability to deal 836 with high dimensionality, localization in ensemble methods, and their practical feasibility), (ii) 837 their effects on different error statistics (RMSE, coverage probabilities, and other diagnostics), 838 (iii) the potential combination of the various methods (especially those considering constant or 839 adaptive covariances), and (iv) the capability to take into account other sources of error (due for 840 instance to improper parameterizations, multiplicative errors, or forcing terms). 841

The use of a realistic DA problem, with a high-dimensional state-space and a limited and heterogeneous observational coverage should be addressed in the future. In that realistic case, the observational information per degree of freedom will be significantly lower, and the estimates of  $\mathbf{Q}$  and  $\mathbf{R}$  will deteriorate. Parametric versions of these error covariance matrices will therefore be

44

<sup>846</sup> necessary. Among the parameters, some of them will control the variances, and will be different
<sup>847</sup> depending on the variable. Other parameters will control the spatial correlation lengths, that could
<sup>848</sup> be isotropic or anisotropic, depending on the region of interest and the considered variable. Cross<sup>849</sup> correlations between variables will also have to be considered. Consequently, **Q** and **R** will be
<sup>850</sup> block-matrices with as few parameters as possible.

A further challenge for future work is the evaluation of the feasibility of estimating non-additive, non-Gaussian, and time-correlated noises under the current estimation frameworks. In this way, the need for observational constraints for the stochastic perturbation methods in the NWP community could be considered within the estimation framework discussed in this review.

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### Four main algorithms to jointly estimate Q and R in data assimilation

```
- initialize inflation factor (for instance \lambda(1) = 1);
for k in 1:K do
      for \underline{i \text{ in } 1:N_e} do
            - compute forecast \mathbf{x}_i^f(k) using Eq. (4a);
            - compute innovation \mathbf{d}_i(k) using Eq. (4b);
      end
      - compute empirical covariance \tilde{\mathbf{P}}^{f}(k) of the \mathbf{x}_{i}^{f}(k);
      - compute \mathbf{K}^{f}(k) using Eq. (4c) where \tilde{\mathbf{P}}^{f}(k)\mathscr{H}_{k}^{\mathrm{T}} and \mathscr{H}_{k}\tilde{\mathbf{P}}^{f}(k)\mathscr{H}_{k}^{\mathrm{T}} are inflated by
        \lambda(k);
      for \underline{i \text{ in } 1:N_e} do
            - compute analysis \mathbf{x}_i^a(k) using Eq. (4d);
      end
      - compute mean innovations \mathbf{d}^{o-f}(k) and \mathbf{d}^{o-a}(k) with \mathbf{d}_i^{o-f}(k) = \mathbf{y}(k) - \mathscr{H}_k(\mathbf{x}_i^f(k))
        and \mathbf{d}_i^{o-a}(k) = \mathbf{y}(k) - \mathscr{H}_k(\mathbf{x}_i^a(k));
      - update \mathbf{R}(k) from Eq. (6b) using the cross-covariance between \mathbf{d}_i^{o-f}(k) and \mathbf{d}_i^{o-a}(k);
      - estimate \tilde{\lambda}(k) using Eq. (8) where \mathscr{H}_k \tilde{\mathbf{P}}^f(k) \mathscr{H}_k^{\mathrm{T}} is inflated by \lambda(k);
      - update \lambda(k+1) using temporal smoother;
end
```

Algorithm 1: Adaptive algorithm for the EnKF (Miyoshi et al. 2013)

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- initialize  $\mathbf{Q}(1)$  and  $\mathbf{R}(1)$ ;

for  $\underline{k \text{ in } 1:K}$  do

for  $\underline{i \text{ in } 1:N_e}$  do

- compute forecast  $\mathbf{x}_i^f(k)$  using Eq. (4a);

- compute innovation  $\mathbf{d}_i(k)$  using Eq. (4b);

end

```
- compute \mathbf{K}^{f}(k) using Eq. (4c);
```

for  $\underline{i \text{ in } 1:N_e}$  do

- compute analysis  $\mathbf{x}_i^a(k)$  using Eq. (4d);

## end

- apply Eq. (12a) to get  $\tilde{\mathbf{P}}(k)$  using linearizations of  $\mathbf{M}_k$  and  $\mathbf{H}_k$  given in Eqs. (5a) and

(5b);

- estimate  $\tilde{\mathbf{Q}}(k)$  using Eq. (12b);

- estimate  $\tilde{\mathbf{R}}(k)$  using Eq. (12c);

```
- update \mathbf{Q}(k+1) and \mathbf{R}(k+1) using temporal smoothers;
```

end

Algorithm 2: Adaptive algorithm for the EnKF (Berry and Sauer 2013)

- define a priori distributions for  $\theta$  (shape parameters of **Q** and **R**);

#### **for** <u>k in 1:K</u> **do**

for i in  $1:N_e$  do

- draw samples  $\boldsymbol{\theta}_i(k)$  from  $p(\boldsymbol{\theta}|\mathbf{y}(1:k-1));$ 

- compute forecast  $\mathbf{x}_i^f(k)$  using Eq. (4a) with  $\boldsymbol{\theta}_i(k)$ ;

- compute innovation  $\mathbf{d}_i(k)$  using Eq. (4b) with  $\boldsymbol{\theta}_i(k)$ ;

end

- compute  $\mathbf{K}^{f}(k)$  using Eq. (4c);

for  $\underline{i \text{ in } 1:N_e}$  do

- compute analysis  $\mathbf{x}_i^a(k)$  using Eq. (4d);

end

- approximate Gaussian likelihood of innovations  $p(\mathbf{y}(k)|\mathbf{y}(1:k-1), \boldsymbol{\theta}(k))$  using empirical mean  $\mathbf{\bar{d}}(k) = \frac{1}{N_e} \sum_{i=1}^{N_e} \mathbf{d}_i(k)$  and empirical covariance  $\Sigma(k) = \frac{1}{N_e-1} \sum_{i=1}^{N_e} (\mathbf{d}_i(k) - \mathbf{\bar{d}}(k)) (\mathbf{d}_i(k) - \mathbf{\bar{d}}(k))^{\mathrm{T}}$  with  $\mathbf{d}_i(k) = \mathbf{y}(k) - \mathscr{H}_k(\mathbf{x}_i^f(k))$ ; - update  $p(\boldsymbol{\theta}|\mathbf{y}(1:k))$  using Eq. (16); end

Algorithm 3: Adaptive algorithm for the EnKF (Stroud et al. 2018)

```
while p(\mathbf{y}(1:K)|\boldsymbol{\theta}_{(n)}) - p(\mathbf{y}(1:K)|\boldsymbol{\theta}_{(n-1)}) > \varepsilon do
```

for  $\underline{k \text{ in } 1:K}$  do

for  $\underline{i \text{ in } 1:N_e}$  do

- compute forecast  $\mathbf{x}_i^f(k)$  using Eq. (4a);

- compute innovation  $\mathbf{d}_i(k)$  using Eq. (4b);

end

```
- compute \mathbf{K}^{f}(k) using Eq. (4c);
```

for  $\underline{i \text{ in } 1:N_e}$  do

- compute analysis  $\mathbf{x}_i^a(k)$  using Eq. (4d);

end

end

```
for \underline{k \text{ in } K:1} do
```

- compute  $\mathbf{K}^{s}(k)$  using Eq. (4e);

for  $\underline{i \text{ in } 1:N_e}$  do

- compute reanalysis  $\mathbf{x}_i^s(k)$  using Eq. (4f);

end

end

```
- increment n \leftarrow n+1;
```

- estimate  $\mathbf{Q}_{(n)}$  using Eq. (21a);

```
- estimate \mathbf{R}_{(n)} using Eq. (21b);
```

end

Algorithm 4: EM algorithm for the EnKF/EnKS (Dreano et al. 2017)

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1145		in data assimilation.	•			•														63

Estimation method	Criteria	Estimation of covariance <b>Q</b>	Suitable for non-Gaussian errors	Application to the highest complexity model
Method of moments	Innovation statistics in the observation space	No (inflation of $\mathbf{P}^{f}$ instead)	No	NWP
Method of moments	Lag innovation between consecutive times	Yes	No	Lorenz-96
Likelihood methods	Bayesian update of the posterior distribution	No (or joint pa- rameter with <b>R</b> )	Yes (using particle filters, not EnKF)	Shallow water
Likelihood methods	Maximization of the to- tal likelihood	Yes	Yes (using particle filters, not EnKF)	Two-scale Lorenz-96

TABLE 1. Comparison of several methods to estimate error covariance matrices  $\mathbf{Q}$  and  $\mathbf{R}$  in data assimilation.

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FIG. 1. Sketch of sequential and ensemble data assimilation algorithms in the observation space (i.e., in the space of the observations **y**), where the observation operator  $\mathscr{H}$  is omitted for simplicity. The ellipses represent the forecast  $\mathbf{P}^{f}$  and analysis  $\mathbf{P}^{a}$  error covariances, while the model **Q** and observation **R** error covariances are the unknown entries of the state-space model in Eqs. (1) and (2). The forecast error covariance matrix is written  $\mathbf{P}^{f}$ and is the sum of  $\mathbf{P}^{m}$ , the forecasted state  $\mathbf{x}^{f}$  spread, and the model error **Q**. This scheme is a modified version based on Fig. 1 from Carrassi et al. (2018).



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