Conditioning of the 3DVAR Data Assimilation Problem

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Abstract

Variational data assimilation schemes are commonly used in major numerical weather prediction (NWP) centres around the world. The convergence of the variational scheme and the sensitivity of the analysis to perturbations are dependent on the conditioning of the Hessian of the linearized least-squares variational equation. The problem is ill-conditioned and hence is difficult to solve quickly and accurately. To make the scheme operationally feasible, NWP centres perform a control variable transform with the aim of preconditioning the problem to reduce the condition number of the Hessian. In this paper we investigate the conditioning of the 3DVar problem for a single periodic system parameter. We give bounds on the condition number of both the original and preconditioned 3DVar problems and demonstrate the reasons for the superior performance of the preconditioned system. We also exhibit the effect of the observation error variances and the positions of the observations on the conditioning of the system.

1 Introduction

Variational data assimilation is popularly used in numerical weather and ocean forecasting to combine observations with a model forecast in order to produce a best estimate of the current state of the system and enable accurate prediction of future states. The estimate minimizes a weighted nonlinear least-squares measure of the error between the model forecast and the available observations and is found using an iterative optimization algorithm. Under certain statistical assumptions the solution to the variational data assimilation problem, known as the *analysis*, yields the *maximum a posteriori* Bayesian estimate of the state of the system [15].

In practice an incremental version of the variational scheme is implemented in many operational centres, including the Met Office [19] and the European Centre for Medium-Range Weather Forecasting (ECMWF) [18]. This method solves a sequence of linear approximations to the nonlinear least-squares problem and is equivalent to an approximate Gauss-Newton method for determining the analysis [14]. Each approximate linearised least-squares problem is solved using an inner gradient iteration method, such as the conjugate gradient method, and the linearization state is then updated in an outer iteration loop. Generally only a very few outer loops are performed. For use in operational forecasting the complete iteration scheme must produce an accurate solution to the variational problem rapidly, in *real* time.

The rate of convergence of the inner loop of the variational scheme and the sensitivity of the solution to perturbations are largely determined by the condition number, that is, the ratio of the largest and smallest eigenvalues, of the Hessian of the linear least-squares objective function [8]. Experimental results indicate that in operational systems the Hessian is ill-conditioned and that this is a result of the ill-conditioning of the background error covariance matrix [16]. In practice the system is preconditioned by transforming the state variables to new variables where the errors are assumed to be approximately uncorrelated [2]. Experimental comparisons have demonstrated that the preconditioning significantly improves the speed and accuracy of the assimilation scheme [7], [16].

Explanations are offered in the literature for the ill-conditioning of the variational assimilation problem and for the benefits of preconditioning in the operational system [1],[20],[16]. In [1], an analysis of the preconditioned system in a simplified 3DVar system with only 2 grid points shows that the conditioning of the preconditioned Hessian is dependent on the accuracy and density of observations. In their paper, Andersson *et al.* take *p* observations at each grid point with error variance σ_o^2 and a background with error variance σ_b^2 and find an approximation to the condition number given by

$$\kappa \approx 2p \left(\frac{\sigma_b^2}{\sigma_o^2}\right) + 1. \tag{1}$$

This approximation is supported experimentally in the ECMWF operational 4DVar system in [20], where it is shown that for dense surface observations, the conditioning of the problem improves as the observations become less accurate. The causes for poor conditioning for dense observations are thus attributed to accurate observations, increasing number of observations (larger p) and large background error variances.

In this paper we examine the conditioning and preconditioning of a more general 3DVar problem theoretically. We derive expressions for the eigenvalues and hence bounds on the conditioning of the Hessian of the problem in the case of a single, periodic, spatially-distributed system parameter. We consider three questions: how does the condition number of the Hessian depend on the length-scale in the correlation structures; how does preconditioning compare with the conditioning of the original Hessian; and how do the error variances of the observations and the distances between observations affect the conditioning of the Hessian.

In the next section we introduce the incremental variational assimilation method. In Section 3 we look at the conditioning of two particular background error covariance matrices. We consider the conditioning of the Hessian and the preconditioned Hessian in Sections 4 and 5. In Section 6 we investigate how the position of observations affects the conditioning and in Section 7 we summarize the conclusions.

2 Variational Data Assimilation

The aim of the variational assimilation scheme is to find an optimal estimate for the initial state of the system \mathbf{x}_0 (the *analysis*) at time t_0 given a *prior* estimate \mathbf{x}_0^b (the *background*) and observations \mathbf{y}_i , $i = 0, \ldots, n$, subject to the nonlinear forecast model given by

$$\mathbf{x}_i = \mathcal{M}(t_i, t_{i-1}, \mathbf{x}_{i-1}), \tag{2}$$

$$\mathbf{y}_i = \mathcal{H}_i(\mathbf{x}_i) + \boldsymbol{\delta}_i,\tag{3}$$

for i = 0, ..., n. Here \mathcal{M} and \mathcal{H}_i denote the evolution and observation operators of the system. The errors $(\mathbf{x}_0 - \mathbf{x}_0^b)$ in the background and the errors δ_i in the observations are assumed to be random with mean zero and covariance matrices **B** and **R**_i, respectively. The assimilation problem is then to minimize, with respect to \mathbf{x}_0 , the objective function

$$J(\mathbf{x}_{0}) = \frac{1}{2} (\mathbf{x}_{0} - \mathbf{x}_{0}^{b})^{T} \mathbf{B}^{-1} (\mathbf{x}_{0} - \mathbf{x}_{0}^{b}) + \frac{1}{2} \sum_{i=0}^{n} (\mathcal{H}_{i}(\mathbf{x}_{i}) - \mathbf{y}_{i})^{T} \mathbf{R}_{i}^{-1} (\mathcal{H}_{i}(\mathbf{x}_{i}) - \mathbf{y}_{i}),$$
(4)

subject to the model forecast equations (2)–(3). If observations are given at several points t_i , i = 0, 1, ..., n over a time window $[t_0, t_n]$ with n > 0, the assimilation scheme is known as the four-dimensional variational method (4DVar). If observations are given only at the initial time with n = 0, then the optimization problem reduces to the three-dimensional data assimilation problem (3DVar).

2.1 Incremental variational assimilation

In operational NWP centres, to reduce computational cost, a sequence of linear approximations to the nonlinear least-squares problem (4) is solved. Given the current estimate of the analysis \mathbf{x}_0 , the nonlinear objective function is linearized about the corresponding model trajectory \mathbf{x}_i , $i = 1, \ldots, n$, satisfying the nonlinear forecast model. An increment $\delta \mathbf{x}_0$ to the current estimate of the analysis is then calculated by minimizing the linearized objective function subject to the linearized model equations. The linear minimization problem is solved in an inner loop by a gradient iteration method. The current estimate of the analysis is then updated with the computed increment and the process is repeated in the outer loop of the algorithm. This data assimilation scheme is known as *incremental variational assimilation* [5], [14].

The linearised objective function, which is minimized with respect to $\delta \mathbf{x}_0$ in each outer loop, can be written

$$\tilde{J}[\delta \mathbf{x}_0] = \frac{1}{2} [\delta \mathbf{x}_0 - (\mathbf{x}_0^b - \mathbf{x}_0)]^T \mathbf{B}^{-1} [\delta \mathbf{x}_0 - (\mathbf{x}_0^b - \mathbf{x}_0)] + \frac{1}{2} (\hat{\mathbf{H}} \delta \mathbf{x}_0 - \hat{\mathbf{d}})^T \hat{\mathbf{R}}^{-1} (\hat{\mathbf{H}} \delta \mathbf{x}_0 - \hat{\mathbf{d}}), \quad (5)$$

subject to the linearized model equations

$$\delta \mathbf{x}_i = \mathbf{M}(t_i, t_{i-1}) \,\delta \mathbf{x}_{i-1},\tag{6}$$

where

$$\hat{\mathbf{H}} = \begin{bmatrix} \mathbf{H}_0^T, (\mathbf{H}_1 \mathbf{M}(t_1, t_0))^T, \dots, (\mathbf{H}_n \mathbf{M}(t_n, t_0))^T \end{bmatrix}^T, \\ \hat{\mathbf{d}}^T = \begin{bmatrix} \mathbf{d}_0^T, \mathbf{d}_1^T, \dots, \mathbf{d}_n^T \end{bmatrix}, \text{ with } \mathbf{d}_i = \mathbf{y}_i - \mathcal{H}_i(\mathbf{x}_i).$$

The matrices $\mathbf{M}(t_i, t_0)$ and \mathbf{H}_i are linearizations of the evolution and observation operators $\mathcal{M}(t_i, t_0, \mathbf{x}_0)$ and $\mathcal{H}_i(\mathbf{x}_i)$ about the current estimated state trajectory \mathbf{x}_i , $i = 0, \ldots, n$ and $\hat{\mathbf{R}}$ is a block diagonal matrix with diagonal blocks equal to \mathbf{R}_i .

The minimizer of (5) is also the solution to $\nabla \tilde{J} = 0$, which may be written explicitly as the linear system

$$(\mathbf{B}^{-1} + \hat{\mathbf{H}}^T \hat{\mathbf{R}}^{-1} \hat{\mathbf{H}}) \delta \mathbf{x}_0 = \mathbf{B}^{-1} (\mathbf{x}_0^b - \mathbf{x}_0) + \hat{\mathbf{H}}^T \hat{\mathbf{R}}^{-1} \hat{\mathbf{d}}.$$
 (7)

Iterative gradient methods are used to solve the inner loop (5), or equivalently (7). The gradients are found by an adjoint procedure.

2.2 Condition number

A measure of the accuracy and efficiency with which the data assimilation problem can be solved is given by the *condition number* of the Hessian matrix

$$\mathbf{A} = (\mathbf{B}^{-1} + \hat{\mathbf{H}}^T \hat{\mathbf{R}}^{-1} \hat{\mathbf{H}})$$
(8)

of the linearized objective function (5) [8]. Our aim here is to establish explicit bounds on the condition number of **A** and to investigate its properties in terms of the background and observation error covariance matrices **B** and $\hat{\mathbf{R}}$.

The condition number of the Hessian, which is a square, symmetric, positive definite matrix, is defined in the 2-norm by

$$\kappa(\mathbf{A}) = ||\mathbf{A}||_2 ||\mathbf{A}^{-1}||_2 \equiv \frac{\lambda_{\max}(\mathbf{A})}{\lambda_{\min}(\mathbf{A})},\tag{9}$$

where $\lambda_{\max}(\mathbf{A})$ and $\lambda_{\min}(\mathbf{A})$ denote the maximum and minimum eigenvalues of the matrix respectively. The condition number measures the sensitivity of the solution to the linearized least-squares problem (5), or equivalently the solution to the gradient equation (7), to perturbations in the data of the problem. If the condition number of the Hessian, $\kappa(\mathbf{A})$, is very large, the problem is 'ill-conditioned' and, even for small perturbations to the system, the relative error in the solution may be extremely large. For the gradient methods that are commonly used to solve the problem, such as the conjugate gradient method, the rate of convergence then may also be very slow. For example, for the conjugate gradient method the error in the computed solution after k iterations is bounded in proportion to $\left(\left(\sqrt{\kappa(\mathbf{A})} - 1\right)/(\sqrt{\kappa(\mathbf{A})} + 1)\right)^k$, which shows that slow convergence can be expected if the Hessian is 'ill-conditioned', with a large condition number.

Here we consider the conditioning of the 3DVar linearized least-squares problem in the case of a single periodic system parameter spatially distributed.

3 Conditioning of the Background Matrix

In [16] it has been proposed that an ill-conditioned background matrix contributes to an illconditioned Hessian. Here we consider the conditioning of two background error covariance matrices commonly used in data assimilation in the case of a single periodic system parameter on a one-dimensional uniform grid of N points.

We write the background error covariance matrix as $\mathbf{B} = \sigma_b^2 \mathbf{C}$, where \mathbf{C} denotes the correlation structure of the background errors and σ_b^2 is the background error variance. The condition number $\kappa(\mathbf{B})$ then equals the condition number $\kappa(\mathbf{C})$. We assume that the correlation structure is homogeneous where the correlations depend only on the distance between states and not position. Under these conditions the correlation matrices are symmetric and have a circulant form [9], which we exploit to derive our theoretical bounds. For example, the Gaussian, Markov and SOAR correlation matrices have this structure, as do those based on Laplacian smoothing. A circulant matrix is a special form of Toeplitz matrix where each row is a cyclic permutation of the previous row

$$\mathbf{C} = \begin{pmatrix} c_0 & c_1 & c_2 & c_3 & \dots & c_{N-2} & c_{N-1} \\ c_{N-1} & c_0 & c_1 & c_2 & \dots & c_{N-3} & c_{N-2} \\ c_{N-2} & c_{N-1} & c_0 & & & \vdots \\ & & \ddots & \ddots & & & \\ \vdots & & & \ddots & \ddots & & \\ \vdots & & & \ddots & & c_2 \\ c_2 & & & \ddots & c_0 & c_1 \\ c_1 & c_2 & \dots & c_{N-2} & c_{N-1} & c_0 \end{pmatrix}$$

The eigenvalues of such a matrix are the discrete Fourier transform of the coefficients of the first row of the matrix [9] and are given by

$$\nu_m = \sum_{k=0}^{N-1} c_k e^{-2\pi i m k/N}.$$
(10)

Similarly the corresponding eigenvectors are given by the discrete exponential function,

$$\mathbf{v}_m = \frac{1}{\sqrt{N}} (1, e^{-2\pi i m/N}, \dots, e^{-2\pi i m(N-1)/N})^T.$$
(11)

3.1 Conditioning of the Gaussian background error covariance matrix

We first consider the Gaussian correlation matrix \mathbf{C} ([6], [11]) with entries given by

$$c_{i,j} = \rho^{|i-j|^2}$$
(12)

for |i - j| < N/2, where $\rho = \exp\left(\frac{-\Delta x^2}{2L^2}\right)$, and by periodicity for the remaining entries. The coefficient $c_{i,j}$ denotes the correlation between background errors at positions i and j, L is the correlation length-scale and determines the strength of the spatial error correlations, Δx is the grid spacing and N is the number of grid points. A large length-scale means that the errors are strongly correlated over the whole grid. The maximum eigenvalue of this correlation matrix is

$$\lambda_{\max}(\mathbf{C}) = \sigma_b^2 \sum_{k=0}^{N-1} \rho^{k^2},\tag{13}$$

with corresponding eigenvector $\mathbf{v}_{\max} = \frac{1}{\sqrt{N}} (1, \dots, 1)^T$. Similarly the minimum eigenvector is

$$\lambda_{\min}(\mathbf{C}) = \sigma_b^2 \sum_{k=0}^{N-1} (-1)^k \rho^{k^2}, \tag{14}$$

with corresponding eigenvector $\mathbf{v}_{\min} = \frac{1}{\sqrt{N}} (1, -1, 1, \dots, -1)^T$.

The condition number is given by the ratio of the maximum to minimum eigenvalues and is highly sensitive to changes in length-scale, as shown in Figure 1 for a grid spacing of $\Delta x = 0.1$ and N = 500 grid points. The matrix becomes very ill-conditioned as the length-scale increases,



Figure 1: Condition number of the periodic Gaussian background error covariance matrix **B** against length-scale.

primarily due to a rapid reduction in its smallest eigenvalue, seen in Figure 3, compared with the modest growth of the largest eigenvalue, seen in Figure 2, as we increase the length-scale. We note that the Gaussian covariance matrix is known to be inherently ill-conditioned. Kostinski [12] shows that the ill-conditioned behaviour of the Gaussian correlation function, which [13] calls the *Gaussian Anomaly*, is due to the fact that the function lies on the boundary of a parametric family of auto-correlation functions with positive Fourier transforms.

3.2 Conditioning of the Laplacian background error covariance matrix

Another commonly used background matrix (see [10]) applies a second derivative smoothing derived from the discrete Laplacian operator. The correlation matrix is defined in terms of its inverse as

$$\mathbf{C}^{-1} = \gamma^{-1} \left(\mathbf{I} + \frac{L^4}{2\Delta x^4} (\mathbf{L})^2 \right), \tag{15}$$



Figure 2: Change in largest eigenvalue of the periodic Gaussian background error covariance matrix \mathbf{B} with length-scale.



Figure 3: Change in smallest eigenvalue of the periodic Gaussian background error covariance matrix ${f B}$ with length-scale.

where the matrix \mathbf{L} is given by

$$\mathbf{L} = \begin{pmatrix} -2 & 1 & 0 & 0 & \dots & 0 & 1 \\ 1 & -2 & 1 & 0 & \dots & 0 & 0 \\ & \ddots & \ddots & & & & \\ \vdots & & \ddots & \ddots & & & & \\ 0 & & & \ddots & \ddots & & & 1 \\ 1 & 0 & \dots & & & 1 & -2 \end{pmatrix},$$

and we define γ so that the maximum value of an element of **C** is unity. Using circulant theory we find the maximum eigenvalue to be

$$\lambda_{\max}(\mathbf{C}^{-1}) = \frac{1}{\gamma} \left(1 + 16 \left(\frac{L^4}{2\Delta x^4} \right) \right), \tag{16}$$

with corresponding eigenvector $\mathbf{w}_{\max} = \frac{1}{\sqrt{N}} (1, -1, 1, \dots, -1)^T$. The smallest eigenvalue is

$$\lambda_{\min}(\mathbf{C}^{-1}) = \frac{1}{\gamma},\tag{17}$$

with corresponding eigenvector $\mathbf{w}_{\min} = \frac{1}{\sqrt{N}} (1, 1, 1, \dots, 1)^T$. The conditioning of the Laplacian correlation matrix is therefore

$$\kappa(\mathbf{C}) = \left(1 + 16\frac{L^4}{2\Delta x^4}\right). \tag{18}$$

The conditioning grows in proportion to L^4 and hence is also quite poorly conditioned. However, as Figure 4 shows, the condition number is many orders of magnitude smaller than that of the Gaussian error covariance matrix at all length-scales.

4 Conditioning of the Hessian

In this section we consider the conditioning of the Hessian of the 3DVar linearized least-squares problem

$$\mathbf{A} = (\mathbf{B}^{-1} + \mathbf{H}^T \mathbf{R}^{-1} \mathbf{H}) \tag{19}$$

in the case of a single periodic system parameter with background error variance σ_b^2 . We examine the Hessian for each of the two background error correlation matrices defined in Section 3. We write the observational error covariance matrix in the form $\mathbf{R} = \sigma_o^2 \mathbf{I}_p$, where p is the number of observations. We assume that the observations are direct measurements of the state variables. Then $\mathbf{H}^T \mathbf{H}$ is a diagonal matrix, where the k^{th} diagonal element is unity if the k^{th} state variable is observed and is zero otherwise. Under these conditions we can prove the following bounds on the condition number of the Hessian matrix for the 3DVar problem

$$\left(\frac{1+\frac{p}{N}\frac{\sigma_b^2}{\sigma_o^2}\lambda_{\min}(\mathbf{C})}{1+\frac{p}{N}\frac{\sigma_b^2}{\sigma_o^2}\lambda_{\max}(\mathbf{C})}\right)\kappa(\mathbf{C}) \le \kappa(\mathbf{B}^{-1} + \mathbf{H}^T\mathbf{R}^{-1}\mathbf{H}) \le \left(1+\left(\frac{\sigma_b^2}{\sigma_o^2}\right)\lambda_{\min}(\mathbf{C})\right)\kappa(\mathbf{C}), \quad (20)$$

where $\lambda_{\max}(\mathbf{C})$ and $\lambda_{\min}(\mathbf{C})$ are the largest and smallest eigenvalues of \mathbf{C} respectively. A proof of this result is given in Appendix A.



Figure 4: Condition number of Laplacian matrix \mathbf{B} against lengthscale.



Figure 5: Condition number of the Hessian (red) and bounds (blue) against length-scale for Gaussian error covariance matrix **B**.

We see that with σ_b fixed, as σ_o increases and the observations become less accurate, the upper bound on the condition number of the Hessian decreases and both the upper and lower bounds converge to $\kappa(\mathbf{C}) = \kappa(\mathbf{B})$. As σ_o decreases, the lower bound goes to unity and, unless σ_o is much smaller than $\lambda_{min}(\mathbf{C})$, the upper bound remains of order $\kappa(\mathbf{C})$. We expect, therefore, that the conditioning of the Hessian will be dominated by the condition number of \mathbf{C} as the correlation length-scales change in the background errors. We demonstrate this in Figure 5 for the Gaussian background covariance matrix with $\sigma_o^2 = \sigma_b^2 = 0.1$, N = 500 grid points and p = 250 observations. Similarly Figure 6 shows the conditioning of the Hessian for the same configuration but using the Laplacian background matrix. (Since the conditioning of the Laplacian is better than that of the Gaussian, a wider range of length-scales is shown in Figure 6.) In these cases including observations has little effect on the conditioning of the assimilation problem.



Figure 6: Condition number of the Hessian (red) and bounds (blue) against length-scale for Laplacian error covariance matrix \mathbf{B} .

5 Preconditioned variational assimilation

A well-known technique for improving the convergence of an iterative method for solving a linear least-squares problem is to apply a linear transformation to 'precondition' the system and thus reduce the condition number of the Hessian [8]. The strategy used in many forecasting centres is to precondition the Hessian symmetrically using the square root of the background error covariance matrix $\mathbf{B}^{1/2}$ [4],[17],[19],[2]. The preconditioning is implemented using a control variable transform to new variables $\delta \mathbf{z} = \mathbf{B}^{-1/2} \delta \mathbf{x}_0$, which are thus uncorrelated. In terms of the new control variable, the 3DVar problem is to minimize with respect to $\delta \mathbf{z}$ the transformed objective function

$$\hat{J}[\delta \mathbf{z}] = \frac{1}{2} [\delta \mathbf{z} - (\mathbf{z}_0^b - \mathbf{z}_0)]^T [\delta \mathbf{z} - (\mathbf{z}_0^b - \mathbf{z}_0)] + \frac{1}{2} (\mathbf{H} \mathbf{B}^{1/2} \delta \mathbf{z} - \mathbf{d})^T \mathbf{R}^{-1} (\mathbf{H} \mathbf{B}^{1/2} \delta \mathbf{z} - \mathbf{d}), \quad (21)$$

where $\mathbf{z}_0^b = \mathbf{B}^{-1/2} \mathbf{x}_0^b$ and $\mathbf{z}_0 = \mathbf{B}^{-1/2} \mathbf{x}_0$. The Hessian of the preconditioned objective function is now given by

$$\mathbf{I}_n + \mathbf{B}^{1/2} \mathbf{H}^T \mathbf{R}^{-1} \mathbf{H} \mathbf{B}^{1/2}.$$
 (22)

In general there are fewer observations than states of the system and therefore the matrix $\mathbf{B}^{1/2}\mathbf{H}^T\mathbf{R}^{-1}\mathbf{H}\mathbf{B}^{1/2}$ is not of full rank, but is positive semi-definite. It follows that the smallest eigenvalue of (22) is unity and the condition number of the preconditioned Hessian is equal to its largest eigenvalue. We can then establish that the condition number satisfies

$$1 + \frac{\sigma_b^2}{\sigma_o^2} \gamma \le \kappa (\mathbf{I}_n + \mathbf{B}^{1/2} \mathbf{H}^T \mathbf{R}^{-1} \mathbf{H} \mathbf{B}^{1/2}) \equiv \kappa (\mathbf{I}_p + \frac{\sigma_b^2}{\sigma_o^2} \mathbf{H} \mathbf{C} \mathbf{H}^T) \le 1 + \frac{\sigma_b^2}{\sigma_o^2} ||\mathbf{H} \mathbf{C} \mathbf{H}^T||_{\infty}, \qquad (23)$$

where $\gamma = \frac{1}{p} \sum_{i,j \in J} c_{i,j}$ and J is the set of indices of the variables that are observed. A proof of this result is given in Appendix B.

We see that the upper bound on the condition number is significantly reduced by preconditioning. For the Gaussian background error covariance matrix we show this in Figure 7 for the case with the same data as in Figure 5. The condition number of the preconditioned problem is shown to be of order unity and to increase roughly linearly. In comparison to the case without preconditioning, there is a dramatic reduction in the condition number from order 10^7 to size $\cong 3.5$ at length-scale L = 0.2. Similar, although less dramatic results, are obtained with the Laplacian correlation matrix. Figure 8 shows the condition number of the preconditioned Hessian with the Laplacian background error covariance matrix **B** using the same configuration as for Figure 6.

If we fix σ_b , but increase σ_o , then the bounds on the conditioning tend towards one. If, however, we decrease σ_o , then both bounds will increase. Hence an increase in the accuracy of the observations results in poorer conditioning of the preconditioned Hessian.

6 Spacing of the Observations

We now consider the condition number of the preconditioned Hessian as a function of the density or separation of the observations. From the definition of the correlation matrix \mathbf{C} the coefficients $c_{i,j}$ are expected to decrease as the distance |i-j| increases. Both upper and lower bounds in



Figure 7: Condition number (red) and bounds (blue) against length-scale for the preconditioned Hessian with the Gaussian background error covariance matrix.



Figure 8: Condition number (red) and bounds (blue)against length-scale for the preconditioned Hessian with the Laplacian background error covariance matrix.

(23) depend on sums of the elements of the matrix \mathbf{HCH}^T , which can be viewed as a 'reduced' correlation matrix. The reduced matrix is simply the correlation matrix \mathbf{C} with all non-observed rows and columns deleted. As the separation of the observations increases, the elements of the reduced matrix decrease, due to the decrease in the coefficients with distance. Similarly, as the observations become more dense and the number of observations increases, the bounds in (23) become larger as more elements are summed. We therefore expect the conditioning of the problem to decrease as the separation of the observations increases or the density decreases.

The effect of spacing of observations on the condition number can be further illustrated if we consider a few special cases. If we observe every q^{th} grid point, where q divides N exactly, then \mathbf{HCH}^T is a circulant matrix. For the Gaussian background error covariance matrix, whose enteries are all positive, the upper and lower bounds are equal and hence we have an exact expression for the condition number

$$1 + \frac{\sigma_b^2}{\sigma_o^2} \sum_{j=1}^p (\mathbf{H}\mathbf{C}\mathbf{H}^T)_{1,j}.$$
 (24)

The larger the spacing between observations, the smaller the elements in the sum (24) are. Similarly the fewer the number of observations, the smaller the total sum is. Hence the conditioning is improved by separating and thinning the observations.

If we have only two observations, at the grid point positions k and m, then we can also write an exact expression for the conditioning. We have

$$\kappa(\mathbf{I} + \mathbf{B}^{1/2}\mathbf{H}^T\mathbf{R}^{-1}\mathbf{H}\mathbf{B}^{1/2}) \equiv \lambda_{max}(\mathbf{I} + \mathbf{B}^{1/2}\mathbf{H}^T\mathbf{R}^{-1}\mathbf{H}\mathbf{B}^{1/2}) = 1 + \frac{\sigma_b^2}{\sigma_o^2}(1 + |c_{k,m}|).$$
(25)

As a function of observation separation, the conditioning changes in proportion to the corresponding background error correlation. In the Gaussian case, as the grid points become farther apart the condition number decays exponentially, as shown in Figure 9.

7 Conclusions

We have examined the conditioning of the 3DVar incremental variational data assimilation problem in the case of a single periodic system parameter. We have derived an expression for the conditioning of two classes of background error correlation structures and demonstrated that the correlation matrices become ill-conditioned as the length-scale increases. In particular the Gaussian correlation matrix has been shown to be especially sensitive to changes in lengthscale. We have also established bounds on the condition number of the Hessian in the 3DVAR case and shown that this is dominated by the background error correlations as the lengthscales increase. The theory shows also that for a fixed background variance, as the variance of the observation errors increases, the conditioning of the Hessian decreases. We have obtained theoretical bounds on the preconditioned Hessian (preconditioned by the symmetric square root of the background covariance matrix) and demonstrated that the preconditioning provides a dramatic reduction in the condition number of the problem. For the preconditioned system we have also shown theoretically that the conditioning of the problem improves as the separation between observations is increased and the density is reduced. We also confirmed the result



Figure 9: Condition number of the preconditioned Hessian for two observations as the grid-point separation is increased. The background error covariance matrix is Gaussian with a length-scale of 0.2.

found in [1] for a two grid-point observing system, which shows that the conditioning of the preconditioned system reduces as the accuracy of the observations is decreased.

Experiments in the Met Office operational variational assimilation system support the theoretical results presented here and confirm that they hold in a more general case. These results will be published in a forthcoming report.

There are two natural extensions to the work presented here. The first is to extend the results for the preconditioned system to encompass more general systems. One approach analogous to our treatment here is to use the dual form of the Hessian [3]

$$\mathbf{I}_p + \mathbf{R}^{-1/2} \mathbf{H} \mathbf{B} \mathbf{H}^T \mathbf{R}^{-1/2}.$$
 (26)

The largest eigenvalue of this matrix is the condition number of the preconditioned Hessian and a general upper bound can be found, as in Appendix B, to be

$$\kappa(\mathbf{I}_N + \mathbf{B}^{1/2}\mathbf{H}^T\mathbf{R}^{-1}\mathbf{H}\mathbf{B}^{1/2}) \le 1 + ||\mathbf{R}^{-1/2}\mathbf{H}\mathbf{B}\mathbf{H}^T\mathbf{R}^{-1/2}||_{\infty}.$$
(27)

For more general systems the conditioning of the assimilation problem may be easier to analyse using (27). The other major extension is to incorporate a forecast model and examine the conditioning of the 4DVar problem. This will be the considered in a future report.

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Appendix A: Bounds on the Condition Number of the Hessian

To bound the condition number of the Hessian

$$\mathbf{A} = \mathbf{B}^{-1} + \mathbf{H}^T \mathbf{R}^{-1} \mathbf{H},\tag{28}$$

we bound the maximum and minimum eigenvalues of the Hessian.

If \mathbf{A}_1 and \mathbf{A}_2 are $n \times n$ symmetric matrices, then

$$\lambda_k(\mathbf{A}_1) + \lambda_n(\mathbf{A}_2) \le \lambda_k(\mathbf{A}_1 + \mathbf{A}_2) \le \lambda_k(\mathbf{A}_1) + \lambda_1(\mathbf{A}_2)$$
(29)

where $\lambda_k(\mathbf{A})$ is the *k*th largest eigenvalue of a matrix \mathbf{A} [8]. Then, noting that $\mathbf{R} = \sigma_o^2 \mathbf{I}_p$, $\mathbf{H}\mathbf{H}^T = \mathbf{I}_p$ and therefore

$$\lambda_1(\mathbf{H}^T \mathbf{R}^{-1} \mathbf{H}) = \sigma_o^{-2} \quad \text{and} \quad \lambda_n(\mathbf{H}^T \mathbf{R}^{-1} \mathbf{H}) = 0,$$
(30)

we find that the maximum and minimum eigenvalues of the Hessian have the following bounds

$$\lambda_{\max}(\mathbf{B}^{-1}) \le \lambda_{\max}(\mathbf{B}^{-1} + \sigma_o^{-2}\mathbf{H}^T\mathbf{H}) \le \lambda_{\max}(\mathbf{B}^{-1}) + \sigma_o^{-2},\tag{31}$$

$$\lambda_{\min}(\mathbf{B}^{-1}) \le \lambda_{\min}(\mathbf{B}^{-1} + \sigma_o^{-2}\mathbf{H}^T\mathbf{H}) \le \lambda_{\min}(\mathbf{B}^{-1}) + \sigma_o^{-2}.$$
(32)

An improvement in the bounds can be achieved using the *Rayleigh quotient*, $R_{\mathbf{A}}(\mathbf{v})$, which, for a Hermitian matrix \mathbf{A} and non-zero vector \mathbf{v} , is defined to be

$$R_{\mathbf{A}}(\mathbf{v}) = \frac{\mathbf{v}^T \mathbf{A} \mathbf{v}}{\mathbf{v}^T \mathbf{v}}.$$
(33)

The maximum and minimum eigenvalues of \mathbf{A} are the maximum and minimum values of $R_{\mathbf{A}}(\mathbf{v})$ respectively, where \mathbf{v} is the corresponding eigenvector. We first consider the eigenvector corresponding to the largest eigenvalue of \mathbf{B}^{-1} . For the the Gaussian covariance matrix this is \mathbf{v}_{\min} and for the Laplacian covariance this is \mathbf{w}_{\max} (see Section 3). In both cases the Rayleigh quotient with respect to $\mathbf{H}^T \mathbf{R}^{-1} \mathbf{H} = \sigma_o^{-2} \mathbf{H}^T \mathbf{H}$ is simply $\sigma_o^{-2} p/N$, where p is the number of observations and N is the number of grid points. Then an improved lower bound on the maximum eigenvalue of the Hessian is given by

$$\lambda_{\max}(\mathbf{A}) = \max_{\mathbf{v} \in \mathbb{R}^n} \left(\frac{\mathbf{v}^T \mathbf{A} \mathbf{v}}{\mathbf{v}^T \mathbf{v}} \right) \ge \mathbf{z}_{\max}^T \mathbf{A} \mathbf{z}_{\max} = \lambda_{\max}(\mathbf{B}^{-1}) + \sigma_o^{-2} \frac{p}{N}, \tag{34}$$

where \mathbf{z}_{max} is either \mathbf{v}_{min} or \mathbf{w}_{max} . Similarly we consider \mathbf{v}_{max} and \mathbf{w}_{min} , which are the eigenvectors corresponding to the smallest eigenvalues of the Gaussian and Laplacian covariance matrices respectively. The same Rayleigh quotient, $\sigma_o^{-2}p/N$, is found with these vectors with respect to $\sigma_o^{-2}\mathbf{H}^T\mathbf{H}$. We find an improved upper bound on the smallest eigenvalue of the Hessian as

$$\lambda_{\min}(\mathbf{A}) = \min_{\mathbf{v} \in \mathbb{R}^n} \left(\frac{\mathbf{v}^T \mathbf{A} \mathbf{v}}{\mathbf{v}^T \mathbf{v}} \right) \le \mathbf{z}_{\min}^T(\mathbf{A}) \mathbf{z}_{\min} = \lambda_{\min}(\mathbf{B}^{-1}) + \sigma_o^{-2} \frac{p}{N}, \tag{35}$$

where \mathbf{z}_{max} is either \mathbf{v}_{max} or \mathbf{w}_{min} . The following bounds are found for the maximum and minimum eigenvalues

$$\lambda_{\max}(\mathbf{B}^{-1}) + \sigma_o^{-2} \frac{p}{n} \le \lambda_{\max}(\mathbf{A}) \le \lambda_{\max}(\mathbf{B}^{-1}) + \sigma_o^{-2}.$$
(36)

and

$$\lambda_{\min}(\mathbf{B}^{-1}) \le \lambda_{\min}(\mathbf{A}) \le \lambda_{\min}(\mathbf{B}^{-1}) + \sigma_o^{-2} \frac{p}{N}.$$
(37)

Noting that, $\lambda_{\max}(\mathbf{B}) = \lambda_{\min}(\mathbf{B}^{-1})$, $\kappa(\mathbf{B}^{-1}) = \kappa(\mathbf{B})$ and using the bounds (37) and (36), the bounds on the condition number of (8) are produced after factoring out $\kappa(\mathbf{B})$ giving the result (20).

Appendix B: Bounds on Condition Number of the Preconditioned Hessian

Using results from [3] we know that the preconditioned Hessian has exactly the same eigenvalues (with an additional n - p unit eigenvalues) as the Hessian of the dual problem $\mathbf{I}_p + \mathbf{R}^{-1/2}\mathbf{HBHR}^{-1/2}$, which for our configuration is equal to

$$\mathbf{I}_{p} + \frac{1}{\sigma_{o}^{2}} \mathbf{H} \mathbf{B} \mathbf{H}^{T} = \mathbf{I}_{p} + \frac{\sigma_{b}^{2}}{\sigma_{o}^{2}} \mathbf{H} \mathbf{C} \mathbf{H}^{T}.$$
(38)

A lower bound can be achieved by considering the Rayleigh quotient on (38). We define a unit vector $\mathbf{v} \in \mathbb{R}^p$ which is equal to $\mathbf{v} = \frac{1}{\sqrt{p}}(1, 1, \dots, 1)^T$. Hence

$$\mathbf{v}^T \mathbf{H} \mathbf{B} \mathbf{H}^T \mathbf{v} = \frac{\sigma_b^2}{p} \sum_{i=1}^p \sum_{j=1}^p (\mathbf{H} \mathbf{C} \mathbf{H}^T)_{i,j}.$$
 (39)

Thus we find that

$$\lambda_{\max}(\mathbf{I}_N + \sigma_o^{-2}\mathbf{B}^{1/2}\mathbf{H}^T\mathbf{H}\mathbf{B}^{1/2}) = \max_{||\mathbf{x}||=1}(\mathbf{x}^T(\mathbf{I}_p + \sigma_o^{-2}\mathbf{H}\mathbf{B}\mathbf{H}^T)\mathbf{x}) \ge 1 + \frac{\sigma_b^2}{p\sigma_o^2}\sum_{i=1}^p\sum_{j=1}^p(\mathbf{H}\mathbf{C}\mathbf{H}^T)_{i,j},$$
(40)

which establishes the lower bound.

The upper bound can be found simply by noting that $\lambda(\mathbf{A}) \leq ||\mathbf{A}||_p$ for any p. Hence

$$\lambda_{\max}(\mathbf{H}\mathbf{C}\mathbf{H}^T) \le ||\mathbf{H}\mathbf{C}\mathbf{H}^T||_{\infty}.$$
(41)

Therefore we have

$$\kappa \left(\mathbf{I}_p + \frac{\sigma_b^2}{\sigma_o^2} (\mathbf{H}\mathbf{C}\mathbf{H}^T) \right) \le 1 + \frac{\sigma_b^2}{\sigma_o^2} ||\mathbf{H}\mathbf{C}\mathbf{H}^T||_{\infty},$$
(42)

which establishes the bounds given in (23).

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