Assimilation Algorithms Lecture 1: Basic Concepts

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Outline

- History and Terminology
- Elementary Statistics The Scalar Analysis Problem
- Extension to Multiple Dimensions
- Optimal Interpolation
- Summary



Outline

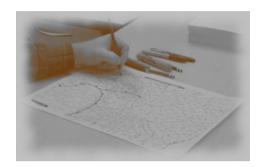
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Terminology

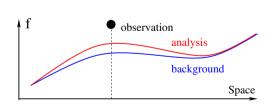
Analysis

- Analysis: The process of approximating the true state of a (geo)physical system at a given time using the available knowledge.
- **X** For example:
 - Hand analysis of synoptic observations (1850 LeVerrier, Fitzroy).
 - Polynomial Interpolation (1950s Panofsky)



Background

★ An important step forward was made by Gilchrist and Cressman (1954), who introduced the idea of using a previous numerical forecast to provide a preliminary estimate of the analysis.



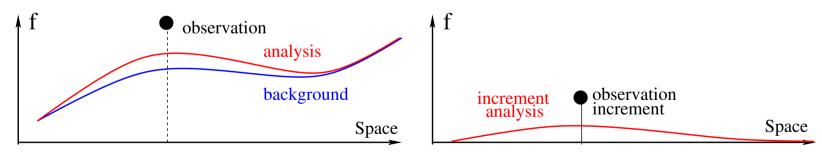
This prior estimate was called the background.



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Optimal interpolation

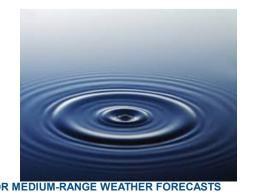
- ➤ Bergthorsson and Döös (1955) took the idea of using a background field a step further by casting the analysis problem in terms of increments which were added to the background.
- ✗ The increments were weighted linear combinations of nearby observation increments (observation minus background), with the weights determined statistically.
- ✗ This idea of statistical combination of background and synoptic observations led ultimately to Optimal Interpolation.
- The use of statistics to merge model fields with observations is fundamental to all current methods of analysis.





Data Assimilation

- X An important change of emphasis happened in the early 1970s with the introduction of primitive-equation models.
- Primitive equation models support inertia-gravity waves. This makes them much more fussy about their initial conditions than the filtered models that had been used hitherto.
- X The analysis procedure became much more intimately linked with the model. The analysis had to produce an initial state that respected the model's dynamical balances.
- Unbalanced increments from the analysis procedure would be rejected as a result of geostrophic adjustment.
- Initialisation techniques (which suppress inertia-gravity waves) became important.





Data Assimilation

The idea that the analysis procedure must present observational information to the model in a way in which it can be absorbed (i.e. not rejected by geostrophic adjustment) led to the coining of the term data assimilation.

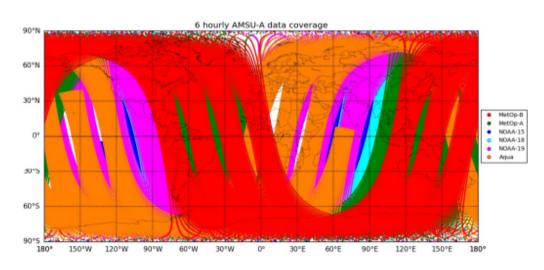
Google define: Assimilate

- X To incorporate nutrients into the body after digestion
- X To incorporate or absorb knowledge into the mind
- The social process of absorbing one cultural group into harmony with another
- ✗ The process by which the Borg integrate beings and cultures into their collective.
- X The process of objectively adapting the model state to observations in a statistically optimal way taking into account model and observation errors



Data Assimilation

- ✗ A final impetus towards the modern concept of data assimilation came from the increasing availability of asynoptic observations from satellite instruments.
- X It was no longer sufficient to think of the analysis purely in terms of spatial interpolation of contemporaneous observations.
- X The time dimension became important, and the model dynamics assumed the role of propagating observational information in time to allow a synoptic view of the state of the system to be generated from asynoptic data.



Example of satellite data coverage in 6 hours (AMSU-A data).



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Problem

Suppose we want to estimate the temperature of this room, given:

- \times A prior estimate: T_b .
 - E.g., we measured the temperature an hour ago, and we have some idea (i.e. a model) of how the temperature varies as a function of time, the number of people in the room, whether the windows are open, etc.
- \times A thermometer: T_o .

Errors

- \times Denote the true temperature of the room by T_t .
- **X** The errors in T_b and T_a are:

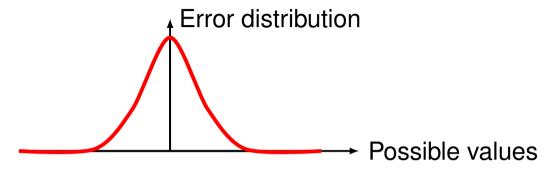
$$\varepsilon_b = T_b - T_t$$

$$\varepsilon_o = T_o - T_t$$

 \times ε_b and ε_o are random variables (or stochastic variables)

Hypotheses

X We will assume that the error statistics of T_b and T_o are known.



We will assume that T_b and T_o have been adjusted (bias corrected) so that their mean errors are zero:

$$\overline{\varepsilon_b} = \overline{\varepsilon_o} = 0$$
.

X There is usually no reason for ε_b and ε_o to be connected in any way:

$$\overline{\varepsilon_o \varepsilon_b} = 0$$
.

X The quantity $\overline{\varepsilon_o \varepsilon_b}$ represents the covariance between the error of our prior estimate and the error of our thermometer measurement.

X We estimate the temperature of the room as a linear combination of T_b and T_o :

$$T_a = \alpha T_o + \beta T_b + \gamma$$

- \blacksquare Denote the error of our estimate as $\varepsilon_a = T_a T_t$.
- **X** We want the estimate to be unbiased: $\overline{\varepsilon_a} = 0$.
- **X** We have:

$$T_a = T_t + \varepsilon_a = \alpha (T_t + \varepsilon_o) + \beta (T_t + \varepsilon_b) + \gamma$$

Taking the mean and rearranging gives:

$$\overline{\varepsilon_a} = (\alpha + \beta - 1) T_t + \gamma$$

- \times Since this holds for any T_t , we must have
 - $\Rightarrow \gamma = 0$, and
 - $\Rightarrow \alpha + \beta 1 = 0.$
- X I.e. $T_a = \alpha T_o + (1 \alpha) T_b$

The general Linear Unbiased Estimate is:

$$T_a = \alpha T_o + (1 - \alpha) T_b$$

- X Now consider the error of this estimate.
- \times Subtracting T_t from both sides of the equation gives

$$\varepsilon_a = \alpha \varepsilon_o + (1 - \alpha) \varepsilon_b$$

X The variance of the estimate is:

$$\overline{\varepsilon_a^2} = \alpha^2 \overline{\varepsilon_o^2} + 2\alpha (1 - \alpha) \overline{\varepsilon_o \varepsilon_b} + (1 - \alpha)^2 \overline{\varepsilon_b^2}$$

X With the previous hypothesis $\overline{\varepsilon_o \varepsilon_b} = 0$:

$$\overline{\varepsilon_a^2} = \alpha^2 \overline{\varepsilon_o^2} + (1 - \alpha)^2 \overline{\varepsilon_b^2}$$

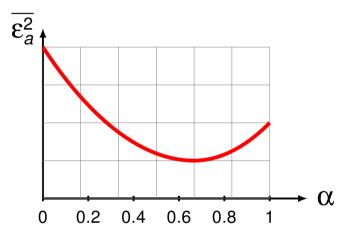
$$\overline{\epsilon_a^2} = \alpha^2 \overline{\epsilon_o^2} + (1 - \alpha)^2 \overline{\epsilon_b^2}$$

We can easily derive some properties of our estimate:

$$\frac{d\overline{\varepsilon_a^2}}{d\alpha} = 2\alpha\overline{\varepsilon_o^2} - 2(1-\alpha)\overline{\varepsilon_b^2}$$

$$imes$$
 For $\alpha=0$, $\overline{\epsilon_a^2}=\overline{\epsilon_b^2}$ and $\frac{d\overline{\epsilon_a^2}}{d\alpha}=-2\overline{\epsilon_b^2}<0$

$$imes$$
 For $\alpha=$ 1, $\overline{\epsilon_a^2}=\overline{\epsilon_o^2}$ and $\frac{d\overline{\epsilon_a^2}}{d\alpha}=2\overline{\epsilon_o^2}>0$



From this we can deduce:

- **x** For $0 \le \alpha \le 1$, $\overline{\varepsilon_a^2} \le \max(\overline{\varepsilon_b^2}, \overline{\varepsilon_o^2})$
- **X** The minimum-variance estimate occurs for $\alpha \in (0,1)$.
- **X** The minimum-variance estimate satisfies $\overline{\varepsilon_a^2} < \min(\overline{\varepsilon_b^2}, \overline{\varepsilon_o^2})$, which means it is lower than the variance of each piece of information.

The minimum-variance estimate occurs when

$$\frac{d\overline{\varepsilon_a^2}}{d\alpha} = 2\alpha\overline{\varepsilon_o^2} - 2(1-\alpha)\overline{\varepsilon_b^2} = 0$$

$$\Rightarrow \quad \alpha = \frac{\overline{\varepsilon_b^2}}{\overline{\varepsilon_b^2} + \overline{\varepsilon_o^2}}.$$

It is not difficult to show that the error variance of this minimum-variance estimate is:

$$\frac{1}{\overline{\varepsilon_a^2}} = \frac{1}{\overline{\varepsilon_b^2}} + \frac{1}{\overline{\varepsilon_o^2}},$$

and the analysis is:

$$\frac{1}{\overline{\varepsilon_a^2}}T_a = \frac{1}{\overline{\varepsilon_b^2}}T_b + \frac{1}{\overline{\varepsilon_o^2}}T_o.$$

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- Now, let's turn our attention to the multi-dimensional case.
- x Instead of a scalar prior estimate T_b , we now consider a vector x_b .
- We can think of \mathbf{x}_b as representing the entire state of a numerical model at some time.
- \mathbf{x} The elements of \mathbf{x}_b might be grid-point values, spherical harmonic coefficients, etc., and some elements may represent temperatures, humidity, others wind components, etc.
- \times We refer to \mathbf{x}_b as the background
- \times Similarly, we generalise the observation to a vector y.
- **y** can contain a disparate collection of observations at different locations, and of different variables.



✗ The major difference between the simple scalar example and the multi-dimensional case is that there is no longer a one-to-one correspondence between the elements of the observation vector and those of the background vector.



- X It is no longer trivial to compare observations and background.
- ✗ Observations are not necessarily located at model gridpoints
- ✗ The observed variables (e.g. radiances) may not correspond directly with any of the variables of the model.
- ✗ To overcome this problem, we must assume that our model is a more-or-less complete representation of reality, so that we can always determine "model equivalents" of the observations.

- **X** We formalise this by assuming the existence of an observation operator, \mathcal{H} .
- \mathbf{x} Given a model-space vector, \mathbf{x} , the vector $\mathcal{H}(\mathbf{x})$ can be compared directly with \mathbf{y} , and represents the "model equivalent" of \mathbf{y} .

$$\mathbf{x} \xrightarrow{\mathcal{H}(\cdot)} \mathcal{H}(\mathbf{x}) o \underbrace{\qquad \qquad} \leftarrow \mathbf{y}$$

For now, we will assume that \mathcal{H} is perfect. I.e. it does not introduce any error, so that:

$$\mathcal{H}(\mathbf{x}_t) = \mathbf{y}_t$$

where \mathbf{x}_t is the true state, and \mathbf{y}_t contains the true values of the observed quantities.

As we did in the scalar case, we will look for an analysis that is a linear combination of the available information:

$$\mathbf{x}_a = \mathbf{F}\mathbf{x}_b + \mathbf{K}\mathbf{y} + \mathbf{c}$$

where **F** and **K** are matrices, and where **c** is a vector.

- $m{x}$ If \mathcal{H} is linear, we can proceed as in the scalar case and look for a linear unbiased estimate.
- \times In the more general case of nonlinear \mathcal{H} , we will require that error-free inputs ($\mathbf{x}_b = \mathbf{x}_t$ and $\mathbf{y} = \mathbf{y}_t$) produce an error-free analysis ($\mathbf{x}_a = \mathbf{x}_t$):

$$\mathbf{x}_t = \mathbf{F}\mathbf{x}_t + \mathbf{K}\mathcal{H}(\mathbf{x}_t) + \mathbf{c}$$

Since this applies for any \mathbf{x}_t , we must have $\mathbf{c} = 0$ and

$$\mathbf{F} \equiv \mathbf{I} - \mathbf{K} \mathcal{H}(\cdot)$$

Our analysis equation is thus:

$$\mathbf{x}_a = \mathbf{x}_b + \mathbf{K} (\mathbf{y} - \mathcal{H} (\mathbf{x}_b))$$



$$\mathbf{x}_a = \mathbf{x}_b + \mathbf{K} (\mathbf{y} - \mathcal{H}(\mathbf{x}_b))$$

Remember that in the scalar case, we had

$$T_a = \alpha T_o + (1 - \alpha) T_b$$

= $T_b + \alpha (T_o - T_b)$

- imes We see that the matrix **K** plays a role equivalent to that of the coefficient α .
- **X** K is called the gain matrix.
- X It determines the weight given to the observation increment
- ✗ It handles the transformation of information defined in "observation space" to the space of model variables.

- X The next step in deriving the analysis equation is to describe the statistical properties of the analysis errors.
- **X** We define

✗ We will assume that the errors are small, so that

$$\mathcal{H}(\mathbf{x}_b) = \mathcal{H}(\mathbf{x}_t) + \mathbf{H}\varepsilon_b + O(\varepsilon_b^2)$$

where **H** is the Jacobian of \mathcal{H} (if **H** is nonlinear).

Substituting the expressions for the errors into our analysis equation, and using $\mathcal{H}(\mathbf{x}_t) = \mathbf{y}_t$, gives (to first order):

$$\mathbf{\epsilon}_{a} = \mathbf{\epsilon}_{b} + \mathbf{K} \left(\mathbf{\epsilon}_{o} - \mathbf{H} \mathbf{\epsilon}_{b} \right)$$

- **X** As in the scalar example, we will assume that the mean errors have been removed, so that $\overline{\varepsilon_b} = \overline{\varepsilon_o} = 0$. We see that this implies that $\overline{\varepsilon_a} = 0$.
- ✗ In the scalar example, we derived the variance of the analysis error, and defined our optimal analysis to minimise this variance.
- X In the multi-dimensional case, we must deal with covariances.



Covariance

 \times The covariance between two variables x_i and x_i is defined as

$$\operatorname{cov}(x_i, x_j) = \overline{(x_i - \overline{x_i})(x_j - \overline{x_j})}$$

- **X** Given a vector $\mathbf{x} = (x_1, x_2, \dots, x_N)^T$, we can arrange the covariances into a covariance matrix, \mathbf{C} , such that $C_{ij} = \text{cov}(x_i, x_j)$.
- X Equivalently:

$$\mathbf{C} = \overline{(\mathbf{x} - \overline{\mathbf{x}})(\mathbf{x} - \overline{\mathbf{x}})^{\mathrm{T}}}$$

- Covariance matrices are symmetric and positive definite
 - \Rightarrow symmetric: $\mathbf{C}^T = \mathbf{C}$
 - \Rightarrow positive definite: $\mathbf{z}^T \mathbf{C} \mathbf{z}$ is positive for every non-zero vector \mathbf{z}

The analysis error is:

$$\epsilon_a = \epsilon_b + \mathbf{K}(\epsilon_o - \mathbf{H}\epsilon_b)$$

$$= (\mathbf{I} - \mathbf{K}\mathbf{H})\epsilon_b + \mathbf{K}\epsilon_o$$

Forming the analysis error covariance matrix gives:

$$\begin{split} \overline{\epsilon_{a}\epsilon_{a}^{\mathrm{T}}} &= \overline{\left[(\mathbf{I} - \mathbf{K} \mathbf{H})\epsilon_{b} + \mathbf{K}\epsilon_{o} \right] \left[(\mathbf{I} - \mathbf{K} \mathbf{H})\epsilon_{b} + \mathbf{K}\epsilon_{o} \right]^{\mathrm{T}}} \\ &= (\mathbf{I} - \mathbf{K} \mathbf{H}) \overline{\epsilon_{b}\epsilon_{b}^{\mathrm{T}}} (\mathbf{I} - \mathbf{K} \mathbf{H})^{\mathrm{T}} + (\mathbf{I} - \mathbf{K} \mathbf{H}) \overline{\epsilon_{b}\epsilon_{o}^{\mathrm{T}}} \mathbf{K}^{\mathrm{T}} \\ &+ \mathbf{K} \overline{\epsilon_{o}\epsilon_{b}^{\mathrm{T}}} (\mathbf{I} - \mathbf{K} \mathbf{H})^{\mathrm{T}} + \mathbf{K} \overline{\epsilon_{o}\epsilon_{o}^{\mathrm{T}}} \mathbf{K}^{\mathrm{T}} \end{split}$$

Assuming that the background and observation errors are uncorrelated (i.e. $\overline{\varepsilon_o \varepsilon_b^{\mathrm{T}}} = \overline{\varepsilon_b \varepsilon_o^{\mathrm{T}}} = 0$), we find:

$$\overline{\epsilon_{\textit{a}}\epsilon_{\textit{a}}^{\mathrm{T}}} = (\mathbf{I} - \mathbf{K}\mathbf{H}) \overline{\epsilon_{\textit{b}}\epsilon_{\textit{b}}^{\mathrm{T}}} (\mathbf{I} - \mathbf{K}\mathbf{H})^{\mathrm{T}} + \mathbf{K} \, \overline{\epsilon_{\textit{o}}\epsilon_{\textit{o}}^{\mathrm{T}}} \mathbf{K}^{\mathrm{T}}$$

$$\overline{\boldsymbol{\varepsilon}_{a}\boldsymbol{\varepsilon}_{a}^{\mathrm{T}}} = \left(\mathbf{I} - \mathbf{K}\mathbf{H}\right)\overline{\boldsymbol{\varepsilon}_{b}\boldsymbol{\varepsilon}_{b}^{\mathrm{T}}}\left(\mathbf{I} - \mathbf{K}\mathbf{H}\right)^{\mathrm{T}} + \mathbf{K}\,\overline{\boldsymbol{\varepsilon}_{o}\boldsymbol{\varepsilon}_{o}^{\mathrm{T}}}\,\mathbf{K}^{\mathrm{T}}$$

X This expression is the equivalent of the expression we obtained for the error of the scalar analysis:

$$\overline{\varepsilon_a^2} = (1 - \alpha)^2 \overline{\varepsilon_b^2} + \alpha^2 \overline{\varepsilon_o^2}$$

- **X** Again, we see that **K** plays essentially the same role in the multi-dimensional analysis as α plays in the scalar case.
- $f{x}$ In the scalar case, we chose α to minimise the variance of the analysis error.
- What do we mean by the minimum-variance analysis in the multi-dimensional case?

- Note that the diagonal elements of a covariance matrix are variances $C_{ii} = \text{cov}(x_i, x_i) = \overline{(x_i \overline{x_i})^2}$.
- ✗ Hence, we can define the minimum-variance analysis as the analysis that minimises the sum of the diagonal elements of the analysis error covariance matrix.
- X The sum of the diagonal elements of a matrix is called the trace.
- In the scalar case, we found the minimum-variance analysis by setting $\frac{d\varepsilon_a^2}{d\alpha}$ to zero.
- X In the multidimensional case, we are going to set

$$\frac{\partial trace(\overline{\epsilon_a \epsilon_a^T})}{\partial \mathbf{K}} = \mathbf{0}$$

× Note: $\frac{\partial \operatorname{trace}(\overline{\varepsilon_a \varepsilon_a^{\mathrm{T}}})}{\partial \mathbf{K}}$ is the matrix whose ij^{th} element is $\frac{\partial \operatorname{trace}(\overline{\varepsilon_a \varepsilon_a^{\mathrm{T}}})}{\partial \mathcal{K}_{ij}}$.

- **X** We have: $\overline{\varepsilon_a \varepsilon_a^{\mathrm{T}}} = (\mathbf{I} \mathbf{K} \mathbf{H}) \overline{\varepsilon_b \varepsilon_b^{\mathrm{T}}} (\mathbf{I} \mathbf{K} \mathbf{H})^{\mathrm{T}} + \mathbf{K} \overline{\varepsilon_o \varepsilon_o^{\mathrm{T}}} \mathbf{K}^{\mathrm{T}}$.
- X The following matrix identities come to our rescue:

$$\frac{\partial trace(\mathbf{K}\mathbf{A}\mathbf{K}^{T})}{\partial \mathbf{K}} = \mathbf{K}(\mathbf{A} + \mathbf{A}^{T})$$

$$\frac{\partial trace(\mathbf{K}\mathbf{A})}{\partial \mathbf{K}} = \mathbf{A}^{T}$$

$$\frac{\partial trace(\mathbf{A}\mathbf{K}^{T})}{\partial \mathbf{K}} = \mathbf{A}$$

X Applying these to $\partial \operatorname{trace}(\overline{\varepsilon_a \varepsilon_a^T})/\partial \mathbf{K}$ gives:

$$\frac{\partial \text{trace}(\overline{\epsilon_a \epsilon_a^{\text{T}}})}{\partial \mathbf{K}} = 2\mathbf{K} \left[\mathbf{H} \overline{\epsilon_b \epsilon_b^{\text{T}}} \mathbf{H}^{\text{T}} + \overline{\epsilon_o \epsilon_o^{\text{T}}} \right] - 2\overline{\epsilon_b \epsilon_b^{\text{T}}} \mathbf{H}^{\text{T}} = \mathbf{0}$$

$$m{\mathsf{X}}$$
 Hence: $m{\mathsf{K}} = \overline{\epsilon_b \epsilon_b^\mathrm{T}} \, m{\mathsf{H}}^\mathrm{T} \left[m{\mathsf{H}} \, \overline{\epsilon_b \epsilon_b^\mathrm{T}} \, m{\mathsf{H}}^\mathrm{T} + \overline{\epsilon_o \epsilon_o^\mathrm{T}} \right]^{-1}$.

$$\mathbf{K} = \overline{\epsilon_b \epsilon_b^{\mathrm{T}}} \mathbf{H}^{\mathrm{T}} \left[\mathbf{H} \, \overline{\epsilon_b \epsilon_b^{\mathrm{T}}} \, \mathbf{H}^{\mathrm{T}} + \overline{\epsilon_o \epsilon_o^{\mathrm{T}}} \right]^{-1}$$

- This optimal gain matrix is called the Kalman Gain Matrix.
- Note the similarity with the optimal gain we derived for the scalar analysis: $\alpha = \overline{\varepsilon_b^2}/(\overline{\varepsilon_b^2} + \overline{\varepsilon_o^2})$.
- The variance of analysis error for the optimal scalar problem was:

$$\frac{1}{\overline{\varepsilon_a^2}} = \frac{1}{\overline{\varepsilon_b^2}} + \frac{1}{\overline{\varepsilon_o^2}}$$

The equivalent expression for the multi-dimensional case is:

$$\left[\overline{\boldsymbol{\epsilon}_{a}\boldsymbol{\epsilon}_{a}^{\mathrm{T}}}\right]^{-1} = \left[\overline{\boldsymbol{\epsilon}_{b}\boldsymbol{\epsilon}_{b}^{\mathrm{T}}}\right]^{-1} + \mathbf{H}^{\mathrm{T}}\left[\overline{\boldsymbol{\epsilon}_{o}\boldsymbol{\epsilon}_{o}^{\mathrm{T}}}\right]^{-1}\mathbf{H}$$

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Notation

- ✗ The notation we have used for covariance matrices can get a bit cumbersome.
- X The standard notation is:

$$egin{aligned} \mathbf{P}^a &\equiv \overline{\mathbf{\epsilon}_a \mathbf{\epsilon}_a^{\mathrm{T}}} \ \mathbf{P}^b &\equiv \overline{\mathbf{\epsilon}_b \mathbf{\epsilon}_b^{\mathrm{T}}} \ \mathbf{R} &\equiv \overline{\mathbf{\epsilon}_o \mathbf{\epsilon}_o^{\mathrm{T}}} \end{aligned}$$

- \times In many analysis schemes, the true covariance matrix of background error, \mathbf{P}^b , is not known, or is too large to be used.
- ✗ In this case, we use an approximate background error covariance matrix.
 This approximate matrix is denoted by **B**.

Alternative Expression for the Kalman Gain

Finally, we derive an alternative expression for the Kalman gain:

$$\mathbf{K} = \mathbf{P}^b \mathbf{H}^{\mathrm{T}} \left[\mathbf{H} \mathbf{P}^b \mathbf{H}^{\mathrm{T}} + \mathbf{R}
ight]^{-1}$$

Multiplying both sides by $\left[\mathbf{P}^{b^{-1}} + \mathbf{H}^{\mathrm{T}}\mathbf{R}^{-1}\mathbf{H}\right]$ gives:

$$\begin{split} \left[\mathbf{P}^{b^{-1}} + \mathbf{H}^{\mathrm{T}}\mathbf{R}^{-1}\mathbf{H}\right]\mathbf{K} &= \left[\mathbf{H}^{\mathrm{T}} + \mathbf{H}^{\mathrm{T}}\mathbf{R}^{-1}\mathbf{H}\mathbf{P}^{b}\mathbf{H}^{\mathrm{T}}\right]\left[\mathbf{H}\mathbf{P}^{b}\mathbf{H}^{\mathrm{T}} + \mathbf{R}\right]^{-1} \\ &= \mathbf{H}^{\mathrm{T}}\mathbf{R}^{-1}\left[\mathbf{R} + \mathbf{H}\mathbf{P}^{b}\mathbf{H}^{\mathrm{T}}\right]\left[\mathbf{H}\mathbf{P}^{b}\mathbf{H}^{\mathrm{T}} + \mathbf{R}\right]^{-1} \\ &= \mathbf{H}^{\mathrm{T}}\mathbf{R}^{-1} \end{split}$$

Hence:

$$\mathbf{K} = \left[\mathbf{P}^{b^{-1}} + \mathbf{H}^{\mathrm{T}}\mathbf{R}^{-1}\mathbf{H}\right]^{-1}\mathbf{H}^{\mathrm{T}}\mathbf{R}^{-1}$$

- \times Expression 1: need the inverse of a matrix of dimension size(\mathbf{R})
- \times Expression 2: need the inverse of a matrix of dimension size(\mathbf{P}^b)

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Optimal Interpolation

- Optimal Interpolation is a statistical data assimilation method based on the multi-dimensional analysis equations we have just derived.
- The method was used operationally at ECMWF from 1979 until 1996, when it was replaced by 3D-Var.
- X The basic idea is to split the global analysis into a number of boxes which can be analysed independently:

$$\mathbf{x}_a^{(i)} = \mathbf{x}_b^{(i)} + \mathbf{K}^{(i)} \left[\mathbf{y}^{(i)} - \mathcal{H}^{(i)}(\mathbf{x}_b) \right]$$

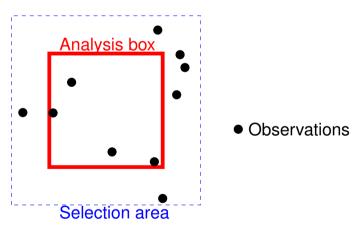
where

$$\mathbf{x}_a = egin{pmatrix} \mathbf{x}_a^{(1)} \\ \mathbf{x}_a^{(2)} \\ \vdots \\ \mathbf{x}_a^{(M)} \end{pmatrix} \qquad \mathbf{x}_b = egin{pmatrix} \mathbf{x}_b^{(1)} \\ \mathbf{x}_b^{(2)} \\ \vdots \\ \mathbf{x}_b^{(M)} \end{pmatrix} \qquad \mathbf{K} = egin{pmatrix} \mathbf{K}^{(1)} \\ \mathbf{K}^{(2)} \\ \vdots \\ \mathbf{K}^{(M)} \end{pmatrix}$$

Optimal Interpolation

$$\mathbf{x}_a^{(i)} = \mathbf{x}_b^{(i)} + \mathbf{K}^{(i)} \left(\mathbf{y}^{(i)} - \mathcal{H}^{(i)}(\mathbf{x}_b) \right)$$

- In principle, we should use *all* available observations to calculate the analysis for each box. However, this is too expensive.
- To produce a computationally-feasible algorithm, Optimal Interpolation (OI) restricts the observations used for each box to those observations which lie in a surrounding selection area:





Optimal Interpolation

The gain matrix used for each box is:

$$\mathbf{K}^{(i)} = \left(\mathbf{P}^b \mathbf{H}^{\mathrm{T}}\right)^{(i)} \left[\left(\mathbf{H} \mathbf{P}^b \mathbf{H}^{\mathrm{T}}\right)^{(i)} + \mathbf{R}^{(i)} \right]^{-1}$$

- **X** Now, the dimension of the matrix $\left[\left(\mathbf{H} \mathbf{P}^b \mathbf{H}^T \right)^{(i)} + \mathbf{R}^{(i)} \right]$ is equal to the number of observations in the selection box.
- ✗ Selecting observations reduces the size of this matrix, making it feasible to use direct solution methods to invert it.
- Note that to implement Optimal Interpolation, we have to specify $(\mathbf{P}^b\mathbf{H}^T)^{(i)}$ and $(\mathbf{H}\mathbf{P}^b\mathbf{H}^T)^{(i)}$. This effectively limits us to very simple observation operators, corresponding to simple interpolations.
- X This, together with the artifacts introduced by observation selection, was one of the main reasons for abandoning Optimal Interpolation in favour of 3D-Var.

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Summary

- We derived the linear analysis equation for a simple scalar example.
- X We showed that a particular choice of the weight α given to the observation resulted in an optimal minimum-variance analysis.
- ✗ We repeated the derivation for the multi-dimensional case. This required the introduction of the observation operator.
- ✗ The derivation for the multi-dimensional case closely parallelled the scalar derivation.
- ✗ The expressions for the gain matrix and analysis error covariance matrix were recognisably similar to the corresponding scalar expressions.
- Finally, we considered the practical implementation of the analysis equation, in an Optimal Interpolation data assimilation scheme.

