

"Optimal" Interpolation (Loren 1981)

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In multi-dimension. n = size of state x
 p = number of observations y^o

$$x = \begin{pmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{pmatrix} \text{ state vector} \quad \varepsilon = \begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \vdots \\ \varepsilon_n \end{pmatrix} \text{ error}$$

$$\Sigma = \overline{\varepsilon \varepsilon^T} = \begin{pmatrix} \overline{\varepsilon_1^2} & \overline{\varepsilon_1 \varepsilon_2} & \cdots & \overline{\varepsilon_1 \varepsilon_n} \\ \overline{\varepsilon_2 \varepsilon_1} & \overline{\varepsilon_2^2} & \cdots & \overline{\varepsilon_2 \varepsilon_n} \\ \vdots & \vdots & \ddots & \vdots \\ \overline{\varepsilon_n \varepsilon_1} & \overline{\varepsilon_n \varepsilon_2} & \cdots & \overline{\varepsilon_n^2} \end{pmatrix} \text{ error covariance}$$

$$y^o = \begin{pmatrix} y_1^o \\ \vdots \\ y_p^o \end{pmatrix} \text{ observations}$$

$$H = \begin{pmatrix} H_{11} & H_{12} & \cdots & H_{1n} \\ \vdots & \vdots & \ddots & \vdots \\ H_{p1} & H_{p2} & \cdots & H_{pn} \end{pmatrix} = \begin{pmatrix} h_1 \\ \vdots \\ h_p \end{pmatrix} \text{ observation operator}$$

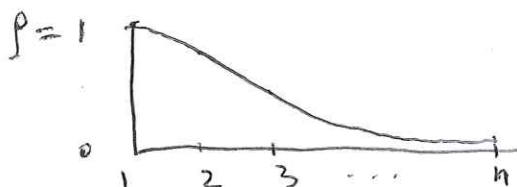
True covariance Σ is not available, use isotropic, stationary covariance B with a given correlation length scale as an assumption.

For a 1D domain case, a row in B encode the correlation function in space, the first row for example:

$$(\overline{\varepsilon_1^2} \quad \overline{\varepsilon_1 \varepsilon_2} \quad \overline{\varepsilon_1 \varepsilon_3} \quad \cdots \quad \overline{\varepsilon_1 \varepsilon_n}) = \varepsilon_1 (\varepsilon_1 \varepsilon_2 \cdots \varepsilon_n)$$

$$= \sigma_1 (\sigma_1 \quad \sigma_2 \rho_{12} \quad \sigma_3 \rho_{13} \quad \cdots \quad \sigma_n \rho_{1n})$$

$(1 \quad \rho_{12} \quad \rho_{13} \quad \cdots \quad \rho_{1n})$ is the Auto correlation Function (ACF)



- To estimate B , one can use the "NMC" method (explained later)

- Using B in place of Σ introduces suboptimality.

Update equation for "optimal" Interpolation (OI):

$$\underset{n \times 1}{X_a} = \underset{n \times 1}{X_b} + \underset{n \times p}{W} (\underset{p \times 1}{y^o} - \underset{p \times 1}{h(X_b)})$$

$$\underset{n \times p}{W} = \underset{n \times n}{B} \underset{n \times p}{H}^T (\underset{p \times n}{H} \underset{n \times n}{B} \underset{n \times p}{H}^T + \underset{p \times p}{R})^{-1}$$

analysis increment: $\Delta X = X_a - X_b$

covariance between observation and state:

$$BH^T = \overline{\varepsilon_b (H \varepsilon_b)^T} = \overline{\begin{pmatrix} \varepsilon_1 \\ \varepsilon_2 \\ \vdots \\ \varepsilon_n \end{pmatrix}_b \begin{pmatrix} h_1 \varepsilon_b & h_2 \varepsilon_b & \dots & h_p \varepsilon_b \end{pmatrix}}$$

variance of observation prior:

$$HBH^T = \overline{(H \varepsilon_b)(H \varepsilon_b)^T} = \overline{\begin{pmatrix} h_1 \varepsilon_b \\ h_2 \varepsilon_b \\ \vdots \\ h_p \varepsilon_b \end{pmatrix} \begin{pmatrix} h_1 \varepsilon_b & h_2 \varepsilon_b & \dots & h_p \varepsilon_b \end{pmatrix}}$$

$$= \begin{pmatrix} \sigma_{y_1}^2 & \dots & \sigma_{y_1}^b \sigma_{y_p}^b \rho_{1p} \\ \vdots & \ddots & \vdots \\ 0 & \dots & \sigma_{y_p}^2 \end{pmatrix}$$

where

$$\sigma_{y_i}^b \sigma_{y_j}^b \rho_{ij} = \overline{(h_i \varepsilon_b)(h_j \varepsilon_b)}$$

variance of observation: $R = \begin{pmatrix} \sigma_{y_1^o}^2 & \dots & 0 \\ \vdots & \ddots & \vdots \\ 0 & \dots & \sigma_{y_p^o}^2 \end{pmatrix}$

analysis error covariance:

$$A = (I - WH)B$$

recall in one-variable case:

$$\sigma_a^2 = \frac{\sigma_b^2 \sigma_o^2}{\sigma_b^2 + \sigma_o^2} = \left(1 - \underbrace{\frac{\sigma_b^2}{\sigma_b^2 + \sigma_o^2}}_w\right) \sigma_b^2$$

Problem: It is often not feasible to calculate W due to large n and p .

Solutions:

1. Assimilate observations sequentially, assuming their errors are uncorrelated (R is diagonal)
2. Perform analysis on a subdomain of x where observations have impact:

$$(BH^T)_{ij} = \overline{\varepsilon_{b_i}(H_j \varepsilon_b)} = 0 \quad \text{if } \sqrt{i^2 + j^2} > \text{Radius of Influence (ROI)}$$

Sequential Algorithm:

For $i = 1, 2, \dots, p$

$$\sigma_{yb}^2 = \overline{(h_i \varepsilon_b)^2} \quad \sigma_{y_o}^2 = R_{ii}$$

$$W = Bh_i^T (\sigma_{yb}^2 + \sigma_{y_o}^2)^{-1}$$

$$x_a = x_b + W(y_i^o - h_i x_b)$$

$$A = (I - Wh_i)B$$

let $x_b = x_a$, $B = A$

end

"NMC" method: (Parrish and Derber, 1992)

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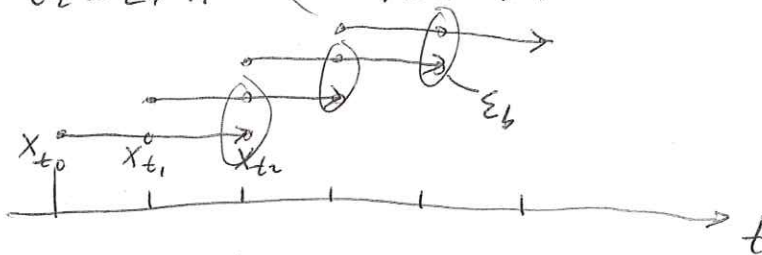
$$B \approx \alpha \overline{(x_{t_1} - x_{t_2})(x_{t_1} - x_{t_2})^T} = \alpha \cdot \overline{\epsilon_b \epsilon_b^T}$$

Scaling factor

Expected value calculated from averaging many realizations

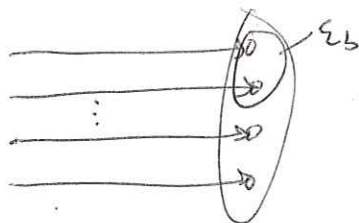
$$t_1 = 48 \text{ h} \quad \left(\text{or } 24 \text{ h} \right)$$

$$t_2 = 24 \text{ h} \quad \left(\text{or } 12 \text{ h} \right)$$



48-h forecast - 24-h forecast
is a "sample" of model forecast error

other option: using ensemble forecast



forecasts using different IC/physics/
models, that characterize errors in models.

$$B = \alpha \overline{\epsilon_b \epsilon_b^T}$$