Parallel Algorithms for Four-Dimensional Variational Data Assimilation

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- Four-Dimensional Variational Data Assimilation 4D-Var is the method used by most national and international Numerical Weather Forecasting Centres to provide initial conditions for their forecast models.
- 4D-Var combines observations with a prior estimate of the state, provided by an earlier forecast.
- The method is described as Four-Dimensional because it takes into account observations that are distributed in space and over an interval of time (typically 6 or 12 hours), often called the analysis window.
- It does this by using a complex and computationally expensive numerical model to propagate information in time.
- In many applications of 4D-Var, the model is assumed to be perfect.
- In this talk, I will concentrate on so-called weak-constraint 4D-Var, which takes into account imperfections in the model.

 Weak-Constraint 4D-Var represents the data-assimilation problem as a very large least-squares problem.

$$\begin{split} J(x_0, x_1, \dots, x_N) &= \frac{1}{2} \left(x_0 - x_b \right)^{\mathrm{T}} B^{-1} \left(x_0 - x_b \right) \\ &+ \frac{1}{2} \sum_{k=0}^{N} \left(y_k - \mathcal{H}_k(x_k) \right)^{\mathrm{T}} R_k^{-1} \left(y_k - \mathcal{H}_k(x_k) \right) \\ &+ \frac{1}{2} \sum_{k=1}^{N} \left(q_k - \bar{q} \right)^{\mathrm{T}} Q_k^{-1} \left(q_k - \bar{q} \right) \end{split}$$

where $q_k = x_k - \mathcal{M}_k(x_{k-1})$.

Here, the cost function J is a function of the states x₀, x₁,..., x_N defined at the start of each of a set of sub-windows that span the analysis window.

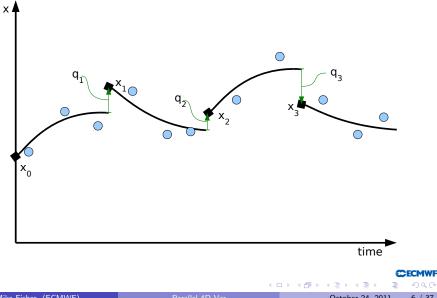
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$$\begin{aligned} J(x_0, x_1, \dots, x_N) &= & \frac{1}{2} (x_0 - x_b)^{\mathrm{T}} B^{-1} (x_0 - x_b) \\ &+ \frac{1}{2} \sum_{k=0}^{N} (y_k - \mathcal{H}_k(x_k))^{\mathrm{T}} R_k^{-1} (y_k - \mathcal{H}_k(x_k)) \\ &+ \frac{1}{2} \sum_{k=1}^{N} (q_k - \bar{q})^{\mathrm{T}} Q_k^{-1} (q_k - \bar{q}) \end{aligned}$$

- Each x_i contains $\approx 10^7$ elements.
- Each y_i contains $\approx 10^5$ elements.
- The operators \mathcal{H}_k and \mathcal{M}_k , and the matrices B, R_k and Q_k are represented by codes that apply them to vectors.
- We do not have access to their elements.

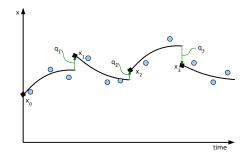
$$\begin{split} J(x_0, x_1, \dots, x_N) &= \frac{1}{2} (x_0 - x_b)^{\mathrm{T}} B^{-1} (x_0 - x_b) \\ &+ \frac{1}{2} \sum_{k=0}^{N} (y_k - \mathcal{H}_k(x_k))^{\mathrm{T}} R_k^{-1} (y_k - \mathcal{H}_k(x_k)) \\ &+ \frac{1}{2} \sum_{k=1}^{N} (q_k - \bar{q})^{\mathrm{T}} Q_k^{-1} (q_k - \bar{q}) \end{split}$$

- \mathcal{H}_k and \mathcal{M}_k involve integrations of the numerical model, and are computationally expensive.
- The covariance matrices B, R_k and Q_k are less expensive to apply.



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The cost function contains 3 terms:

- $\frac{1}{2}(x_0 x_b)^T B^{-1}(x_0 x_b)$ ensures that x_0 stays close to the prior estimate.
- $\frac{1}{2} \sum_{k=0}^{N} (y_k \mathcal{H}_k(x_k))^T R_k^{-1} (y_k \mathcal{H}_k(x_k))$ keeps the estimate close to the observations (blue circles).
- $\frac{1}{2} \sum_{k=1}^{N} (q_k \bar{q})^{\mathrm{T}} Q_k^{-1} (q_k \bar{q})$ keeps the jumps between sub-windows small.

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Gauss-Newton (Incremental) Algorithm

- It is usual to minimize the cost function using a modified Gauss-Newton algorithm.
- Nearly all the computational cost is in the inner loop, which minimizes the quadratic cost function:

$$\begin{aligned} \hat{J}(\delta x_{0}^{(n)}, \dots, \delta x_{N}^{(n)}) &= \frac{1}{2} \left(\delta x_{0} - b^{(n)} \right)^{\mathrm{T}} B^{-1} \left(\delta x_{0} - b^{(n)} \right) \\ &+ \frac{1}{2} \sum_{k=0}^{N} \left(H_{k}^{(n)} \delta x_{k} - d_{k}^{(n)} \right)^{\mathrm{T}} R_{k}^{-1} \left(H_{k}^{(n)} \delta x_{k} - d_{k}^{(n)} \right) \\ &+ \frac{1}{2} \sum_{k=1}^{N} \left(\delta q_{k} - c_{k}^{(n)} \right)^{\mathrm{T}} Q_{k}^{-1} \left(\delta q_{k} - c_{k}^{(n)} \right) \end{aligned}$$

 $\delta q_k = \delta x_k - M_k^{(n)} \delta x_{k-1}$, and where $b^{(n)}$, $c_k^{(n)}$ and $d_k^{(n)}$ come from the outer loop:

$$b^{(n)} = x_b - x_0^{(n)}, \qquad c_k^{(n)} = \bar{q} - q_k^{(n)}, \qquad d_k^{(n)} = y_k - \mathcal{H}_k(x_k^{(n)})$$

Gauss-Newton (Incremental) Algorithm

$$\begin{split} \hat{J}(\delta x_{0}^{(n)}, \dots, \delta x_{N}^{(n)}) &= \frac{1}{2} \left(\delta x_{0} - b^{(n)} \right)^{\mathrm{T}} B^{-1} \left(\delta x_{0} - b^{(n)} \right) \\ &+ \frac{1}{2} \sum_{k=0}^{N} \left(H_{k}^{(n)} \delta x_{k} - d_{k}^{(n)} \right)^{\mathrm{T}} R_{k}^{-1} \left(H_{k}^{(n)} \delta x_{k} - d_{k}^{(n)} \right) \\ &+ \frac{1}{2} \sum_{k=1}^{N} \left(\delta q_{k} - c_{k}^{(n)} \right)^{\mathrm{T}} Q_{k}^{-1} \left(\delta q_{k} - c_{k}^{(n)} \right) \end{split}$$

- Note that 4D-Var requires tangent linear versions of M_k and H_k:
 M⁽ⁿ⁾_k and H⁽ⁿ⁾_k, respectively
- It also requires the transposes (adjoints) of these operators:
 - $(M_k^{(n)})^{\mathrm{T}}$ and $(H_k^{(n)})^{\mathrm{T}}$, respectively

Why do we need more parallel algorithms?

- In its usual implementation, 4D-Var is solved by applying a conjugate-gradient solver.
- This is highly sequential:
 - Iterations of CG.
 - Tangent Linear and Adjoint integrations run one after the other.
 - Model timesteps follow each other.
- Computers are becoming ever more parallel, but processors are not getting faster.
- Unless we do something to make 4D-Var more parallel, we will soon find that 4D-Var becomes un-affordable (even with a 12-hour window).
- We cannot make the model more parallel.
 - The inner loops of 4D-Var run with a few 10's of grid columns per processor.
 - This is barely enough to mask inter-processor communication costs.
- We have to use more parallel algorithms.

Parallelising within an Iteration

- The model is already parallel in both horizontal directions.
- The modellers tell us that it will be hard to parallelise in the vertical (and we already have too little work per processor).
- We are left with parallelising in the time direction.

Dropping the outer loop index (n), the inner loop of weak-constraints 4D-Var minimises:

$$\begin{split} \hat{J}(\delta x_0, \dots, \delta x_N) &= \frac{1}{2} \left(\delta x_0 - b \right)^{\mathrm{T}} B^{-1} \left(\delta x_0 - b \right) \\ &+ \frac{1}{2} \sum_{k=0}^{N} \left(H_k \delta x_k - d_k \right)^{\mathrm{T}} R_k^{-1} \left(H_k \delta x_k - d_k \right) \\ &+ \frac{1}{2} \sum_{k=1}^{N} \left(\delta q_k - c_k \right)^{\mathrm{T}} Q_k^{-1} \left(\delta q_k - c_k \right) \end{split}$$

where $\delta q_k = \delta x_k - M_k \delta x_{k-1}$, and where *b*, c_k and d_k come from the outer loop:

$$b = x_b - x_0$$

$$c_k = \bar{q} - q_k$$

$$d_k = y_k - \mathcal{H}_k(x_k)$$
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We can simplify this further by defining some 4D vectors and matrices:

$$\delta \mathbf{x} = \begin{pmatrix} \delta x_0 \\ \delta x_1 \\ \vdots \\ \delta x_N \end{pmatrix} \qquad \qquad \delta \mathbf{p} = \begin{pmatrix} \delta x_0 \\ \delta q_1 \\ \vdots \\ \delta q_N \end{pmatrix}$$

These vectors are related through $\delta q_k = \delta x_k - M_k \delta x_{k-1}$. We can write this relationship in matrix form as:

$$\delta \mathbf{p} = \mathbf{L} \delta \mathbf{x}$$

where:

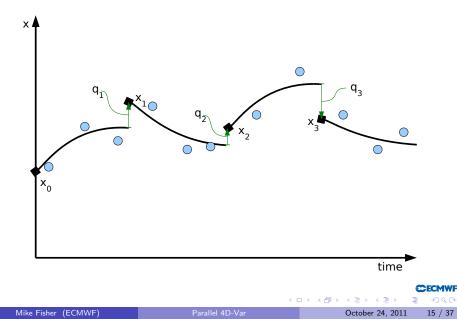
$$\mathbf{L} = \begin{pmatrix} I & & & \\ -M_1 & I & & \\ & -M_2 & I & \\ & & \ddots & \ddots & \\ & & & -M_N & I \\ & & & -M_N & I \end{pmatrix}$$

$$\mathbf{L} = \begin{pmatrix} I & & & \\ -M_1 & I & & & \\ & -M_2 & I & & \\ & & \ddots & \ddots & \\ & & & -M_N & I \end{pmatrix}$$

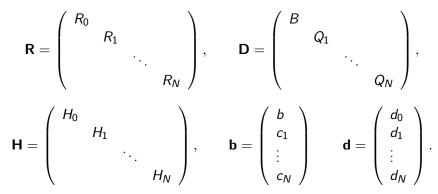
 $\delta \mathbf{p} = \mathbf{L} \delta \mathbf{x}$ can be done in parallel: $\delta q_k = \delta x_k - M_k \delta x_{k-1}$. We know all the $\delta x_{k-1}'s$. We can apply all the $M_k's$ simultaneously.

 $\delta \mathbf{x} = \mathbf{L}^{-1} \delta \mathbf{p}$ is sequential: $\delta x_k = M_k \delta x_{k-1} + \delta q_k$. We have to generate each δx_{k-1} in turn before we can apply the next M_k .

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We will also define:



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With these definitions, we can write the inner-loop cost function either as a function of $\delta \mathbf{x}$:

$$J(\delta \mathbf{x}) = (\mathbf{L} \delta \mathbf{x} - \mathbf{b})^{\mathrm{T}} \mathbf{D}^{-1} (\mathbf{L} \delta \mathbf{x} - \mathbf{b}) + (\mathbf{H} \delta \mathbf{x} - \mathbf{d})^{\mathrm{T}} \mathbf{R}^{-1} (\mathbf{H} \delta \mathbf{x} - \mathbf{d})$$

Or as a function of $\delta \mathbf{p}$:

$$J(\delta \mathbf{p}) = (\delta \mathbf{p} - \mathbf{b})^{\mathrm{T}} \mathbf{D}^{-1} (\delta \mathbf{p} - \mathbf{b}) + (\mathbf{H} \mathbf{L}^{-1} \delta \mathbf{p} - \mathbf{d})^{\mathrm{T}} \mathbf{R}^{-1} (\mathbf{H} \mathbf{L}^{-1} \delta \mathbf{p} - \mathbf{d})$$

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Forcing Formulation

$$J(\delta \mathbf{p}) = (\delta \mathbf{p} - \mathbf{b})^{\mathrm{T}} \mathbf{D}^{-1} (\delta \mathbf{p} - \mathbf{b}) + (\mathbf{H} \mathbf{L}^{-1} \delta \mathbf{p} - \mathbf{d})^{\mathrm{T}} \mathbf{R}^{-1} (\mathbf{H} \mathbf{L}^{-1} \delta \mathbf{p} - \mathbf{d})$$

- This version of the cost function is sequential.
 - It contains L⁻¹.
- It closely resembles strong-constraint 4D-Var.
- We can precondition it using $D^{1/2}$:

$$J(\chi) = \chi^{\mathrm{T}} \chi + (\mathbf{H} \mathbf{L}^{-1} \delta \mathbf{p} - \mathbf{d})^{\mathrm{T}} \mathbf{R}^{-1} (\mathbf{H} \mathbf{L}^{-1} \delta \mathbf{p} - \mathbf{d})$$

where $\delta \mathbf{p} = \mathbf{D}^{1/2} \chi + \mathbf{b}$.

- This guarantees that the eigenvalues of J" are bounded away from zero.
- We understand how to minimise this.

4D State Formulation

$J(\delta \mathbf{x}) = (\mathbf{L}\delta \mathbf{x} - \mathbf{b})^{\mathrm{T}} \mathbf{D}^{-1} (\mathbf{L}\delta \mathbf{x} - \mathbf{b}) + (\mathbf{H}\delta \mathbf{x} - \mathbf{d})^{\mathrm{T}} \mathbf{R}^{-1} (\mathbf{H}\delta \mathbf{x} - \mathbf{d})$

• This version of the cost function is parallel.

- It does not contain L⁻¹.
- We could precondition it using $\delta \mathbf{x} = \mathbf{L}^{-1} (\mathbf{D}^{1/2} \chi + \mathbf{b})$.
- This would give exactly the same $J(\chi)$ as before.
- But, we have introduced a sequential model integration (i.e. L⁻¹) into the preconditioner.

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Plan A: State Formulation, Approximate Preconditioner

- In the forcing $(\delta \mathbf{p})$ formulation, and in its Lagrangian dual formulation (4D-PSAS) \mathbf{L}^{-1} appears in the cost function.
 - These formulations are inherently sequential.
 - We cannot modify the cost function without changing the problem.
- In the 4D-state (δx) formulation, L^{-1} appears in the preconditioner.
 - We are free to modify the preconditioner as we wish.
- This suggests we replace L^{-1} by a cheap approximation:

$$\delta \mathbf{x} = \mathbf{\tilde{L}}^{-1} (\mathbf{D}^{1/2} \chi + \mathbf{b})$$

- If we do this, we can no longer write the first term as: $\chi^{\rm T}\chi$.
- We have to calculate $\delta \mathbf{x}$, and explicity evaluate it as

$$(\mathbf{L}\delta \mathbf{x} - \mathbf{b})^{\mathrm{T}}\mathbf{D}^{-1}(\mathbf{L}\delta \mathbf{x} - \mathbf{b})$$

• This is where we run into problems: **D** is very ill-conditioned.

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Plan A: State Formulation, Approximate Preconditioner

• When we approximate L^{-1} in the preconditioner, the Hessian of the first term of the cost function (with respect to χ) is no longer the identity matrix, but:

$\boldsymbol{\mathsf{D}}^{\mathrm{T}/2}\boldsymbol{\tilde{\mathsf{L}}}^{-\mathrm{T}}\boldsymbol{\mathsf{L}}^{\mathrm{T}}\boldsymbol{\mathsf{D}}^{-1}\boldsymbol{\mathsf{L}}\boldsymbol{\tilde{\mathsf{L}}}^{-1}\boldsymbol{\mathsf{D}}^{1/2}$

- Because **D** is ill-conditioned, This is likely to have some very small (and some very large) eigenvalues, unless \tilde{L} is a very good approximation for **L**.
- So far, I have not found a preconditioner that gives condition numbers for the minimisation less than $O(10^9)$.
- We need a Plan B!

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Plan B: Saddle Point Formulation

$$J(\delta \mathbf{x}) = (\mathbf{L} \delta \mathbf{x} - \mathbf{b})^{\mathrm{T}} \mathbf{D}^{-1} (\mathbf{L} \delta \mathbf{x} - \mathbf{b}) + (\mathbf{H} \delta \mathbf{x} - \mathbf{d})^{\mathrm{T}} \mathbf{R}^{-1} (\mathbf{H} \delta \mathbf{x} - \mathbf{d})$$

At the minimum:

$$\nabla J = \mathbf{L}^{\mathrm{T}} \mathbf{D}^{-1} (\mathbf{L} \delta \mathbf{x} - \mathbf{b}) + \mathbf{H}^{\mathrm{T}} \mathbf{R}^{-1} (\mathbf{H} \delta \mathbf{x} - \mathbf{d}) = \mathbf{0}$$

Define:

$$\lambda = \mathbf{D}^{-1}(\mathbf{b} - \mathbf{L}\delta \mathbf{x}), \qquad \mu = \mathbf{R}^{-1}(\mathbf{d} - \mathbf{H}\delta \mathbf{x})$$

Then:

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$$\left(\begin{array}{ccc} \mathbf{D} & \mathbf{0} & \mathbf{L} \\ \mathbf{0} & \mathbf{R} & \mathbf{H} \\ \mathbf{L}^{\mathrm{T}} & \mathbf{H}^{\mathrm{T}} & \mathbf{0} \end{array}\right) \left(\begin{array}{c} \lambda \\ \mu \\ \delta \mathbf{x} \end{array}\right) = \left(\begin{array}{c} \mathbf{b} \\ \mathbf{d} \\ \mathbf{0} \end{array}\right)$$

- This is called the saddle point formulation of 4D-Var.
- The matrix is a saddle point matrix.
- The matrix is real, symmetric, indefinite.
- Note that the matrix contains no inverse matrices.
- We can apply the matrix without requiring a sequential model integration (i.e. we can parallelise over sub-windows).
- We can hope that the problem is well conditioned (since we don't multiply by D⁻¹).

Alternative derivation:

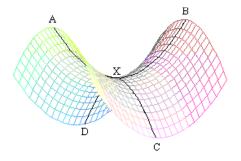
$$\min_{\delta \mathbf{p}, \delta \mathbf{w}} J(\delta \mathbf{p}, \delta \mathbf{w}) = (\delta \mathbf{p} - \mathbf{b})^{\mathrm{T}} \mathbf{D}^{-1} (\delta \mathbf{p} - \mathbf{b}) + (\delta \mathbf{w} - \mathbf{d})^{\mathrm{T}} \mathbf{R}^{-1} (\delta \mathbf{w} - \mathbf{d})$$

subject to $\delta \mathbf{p} = \mathbf{L} \delta \mathbf{x}$ and $\delta \mathbf{w} = \mathbf{H} \delta \mathbf{x}$.

$$\mathcal{L}(\delta \mathbf{x}, \delta \mathbf{p}, \delta \mathbf{w}, \lambda, \mu) = (\delta \mathbf{p} - \mathbf{b})^{\mathrm{T}} \mathbf{D}^{-1} (\delta \mathbf{p} - \mathbf{b}) + (\delta \mathbf{w} - \mathbf{d})^{\mathrm{T}} \mathbf{R}^{-1} (\delta \mathbf{w} - \mathbf{d}) + \lambda^{\mathrm{T}} (\delta \mathbf{p} - \mathbf{L} \delta \mathbf{x}) + \mu^{\mathrm{T}} (\delta \mathbf{w} - \mathbf{H} \delta \mathbf{x})$$

•
$$\frac{\partial \mathcal{L}}{\partial \lambda} = \mathbf{0} \Rightarrow \qquad \delta \mathbf{p} = \mathbf{L} \delta \mathbf{x}$$

• $\frac{\partial \mathcal{L}}{\partial \mu} = \mathbf{0} \Rightarrow \qquad \delta \mathbf{w} = \mathbf{H} \delta \mathbf{x}$
• $\frac{\partial \mathcal{L}}{\partial \delta \mathbf{p}} = \mathbf{0} \Rightarrow \qquad \mathbf{D}^{-1} (\delta \mathbf{p} - \mathbf{b}) + \lambda = \mathbf{0}$
• $\frac{\partial \mathcal{L}}{\partial \delta \mathbf{w}} = \mathbf{0} \Rightarrow \qquad \mathbf{R}^{-1} (\delta \mathbf{w} - \mathbf{d}) + \mu = \mathbf{0}$
• $\frac{\partial \mathcal{L}}{\partial \delta \mathbf{x}} = \mathbf{0} \Rightarrow \qquad \mathbf{L}^{\mathrm{T}} \lambda + \mathbf{H}^{\mathrm{T}} \mu = \mathbf{0}$
• $\frac{\partial \mathcal{L}}{\partial \delta \mathbf{x}} = \mathbf{0} \Rightarrow \qquad \mathbf{L}^{\mathrm{T}} \lambda + \mathbf{H}^{\mathrm{T}} \mu = \mathbf{0}$
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Lagrangian: $\mathcal{L}(\delta \mathbf{x}, \delta \mathbf{p}, \delta \mathbf{w}, \lambda, \mu)$

- 4D-Var solves the primal problem: minimise along AXB.
- 4D-PSAS solves the Lagrangian dual problem: maximise along CXD.
- The saddle point formulation finds the saddle point of \mathcal{L} .
- The saddle point formulation is neither 4D-Var nor 4D-PSAS.

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- To solve the saddle point system, we have to precondition it.
- Preconditioning saddle point systems is the subject of much current research. It is something of a dark art!
 - See e.g. Benzi and Wathen (2008), Benzi, Golub and Liesen (2005).
- Most preconditioners in the literature assume that **D** and **R** are expensive, and **L** and **H** are cheap.
- The opposite is true in 4D-Var!

Example: Diagonal Preconditioner:

$$\mathcal{P}_D = \left(egin{array}{ccc} \hat{\mathbf{D}} & \mathbf{0} & \mathbf{0} \ \mathbf{0} & \hat{\mathbf{R}} & \mathbf{0} \ \mathbf{0} & \mathbf{0} & -\mathbf{S} \end{array}
ight)$$

where $\boldsymbol{\mathsf{S}}\approx-\boldsymbol{\mathsf{L}}^{\mathrm{T}}\boldsymbol{\mathsf{D}}^{-1}\boldsymbol{\mathsf{L}}-\boldsymbol{\mathsf{H}}^{\mathrm{T}}\boldsymbol{\mathsf{R}}^{-1}\boldsymbol{\mathsf{H}}$

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• One possibility is an approximate constraint preconditioner (Bergamaschi, *et al.*, 2007 & 2011):

$$\tilde{\mathcal{P}} = \left(\begin{array}{ccc} \mathbf{D} & \mathbf{0} & \mathbf{\tilde{L}} \\ \mathbf{0} & \mathbf{R} & \mathbf{0} \\ \mathbf{\tilde{L}}^{\mathrm{T}} & \mathbf{0} & \mathbf{0} \end{array} \right)$$

$$\Rightarrow \tilde{\mathcal{P}}^{-1} = \left(\begin{array}{ccc} \mathbf{0} & \mathbf{0} & \mathbf{\tilde{L}}^{-\mathrm{T}} \\ \mathbf{0} & \mathbf{R}^{-1} & \mathbf{0} \\ \mathbf{\tilde{L}}^{-1} & \mathbf{0} & -\mathbf{\tilde{L}}^{-1}\mathbf{D}\mathbf{\tilde{L}}^{-\mathrm{T}} \end{array} \right)$$

• Note that $\tilde{\mathcal{P}}^{-1}$ does not contain \mathbf{D}^{-1} .

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 $\bullet\,$ With this preconditioner, we can prove some nice results for the case $\tilde{L}=L$:

$$\begin{pmatrix} \mathbf{0} & \mathbf{0} & \mathbf{L}^{-\mathrm{T}} \\ \mathbf{0} & \mathbf{R}^{-1} & \mathbf{0} \\ \mathbf{L}^{-1} & \mathbf{0} & -\mathbf{L}^{-1}\mathbf{D}\mathbf{L}^{-\mathrm{T}} \end{pmatrix} \begin{pmatrix} \mathbf{D} & \mathbf{0} & \mathbf{L} \\ \mathbf{0} & \mathbf{R} & \mathbf{H} \\ \mathbf{L}^{\mathrm{T}} & \mathbf{H}^{\mathrm{T}} & \mathbf{0} \end{pmatrix} \begin{pmatrix} \lambda \\ \mu \\ \delta \mathbf{x} \end{pmatrix} = \tau \begin{pmatrix} \lambda \\ \mu \\ \delta \mathbf{x} \end{pmatrix}$$

$$\Rightarrow \mathbf{I} + \begin{pmatrix} \mathbf{0} & \mathbf{L}^{-\mathrm{T}}\mathbf{H}^{\mathrm{T}} & \mathbf{0} \\ \mathbf{0} & \mathbf{0} & \mathbf{R}^{-1}\mathbf{H} \\ \mathbf{0} & -\mathbf{L}^{-1}\mathbf{D}\mathbf{L}^{-\mathrm{T}}\mathbf{H}^{\mathrm{T}} & \mathbf{0} \end{pmatrix} \begin{pmatrix} \lambda \\ \mu \\ \delta \mathbf{x} \end{pmatrix} = \tau \begin{pmatrix} \lambda \\ \mu \\ \delta \mathbf{x} \end{pmatrix}$$

$$\Rightarrow \mathbf{H}\mathbf{L}^{-1}\mathbf{D}\mathbf{L}^{-\mathrm{T}}\mathbf{H}^{\mathrm{T}}\boldsymbol{\mu} + (\tau - 1)^{2}\mathbf{R}\boldsymbol{\mu} = \mathbf{0}$$

 \Rightarrow ($\tau - 1$) is imaginary (or zero) since $\mathbf{H}\mathbf{L}^{-1}\mathbf{D}\mathbf{L}^{-T}\mathbf{H}^{T}$ is positive semi-definite and **R** is positive definite.

- The eigenvalues τ of $\tilde{\mathcal{P}}^{-1}\mathcal{A}$ lie on the line $\Re(\tau) = 1$ in the complex plane.
- Their distance above/below the real axis is:

$$\pm \sqrt{\frac{\mu_i^{\mathrm{T}} \mathbf{H} \mathbf{L}^{-1} \mathbf{D} \mathbf{L}^{-\mathrm{T}} \mathbf{H}^{\mathrm{T}} \mu_i}{\mu_i^{\mathrm{T}} \mathbf{R} \mu_i}}$$

where μ_i is the μ component of the *i*th eigenvector.

- The fraction under the square root is, roughly speaking, the ratio of background+model error variance to observation error variance associated with the pattern μ_i.
- In the usual implementation of 4D-Var, the condition number is given by the ratio of these variances, not the square-root.

- For the preconditioner, we need an approximate inverse of L.
- One approach is to use the following identity (exercise for the reader!):

$$\mathbf{L}^{-1} = \mathbf{I} + (\mathbf{I} - \mathbf{L}) + (\mathbf{I} - \mathbf{L})^2 + \ldots + (\mathbf{I} - \mathbf{L})^{N-1}$$

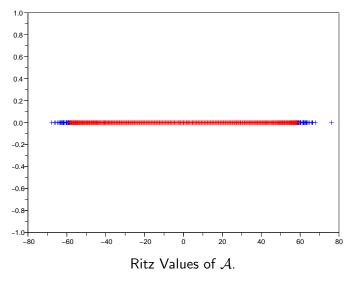
- Since this is a power series expansion, it suggests truncating the series at some order < N 1.
- (A very few iterations of a Krylov solver may be a better idea. I've not tried this yet.)

Results

- The practical reaults shown in the next few slides are for a simplified (toy) analogue of a real system.
- The model is a two-level quasi-geostrophic channel model with 1600 gridpoints.
- The model has realistic error-growth and time-to-nonlinearity
- There are 100 observations of streamfunction every 3 hours, and 100 wind observations every 6 hours.
- The error covariances are assumed to be horizontally isotropic and homogeneous, with a Gaussian spatial structure.

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OOPS QG model. 24-hour window with 8 sub-windows.

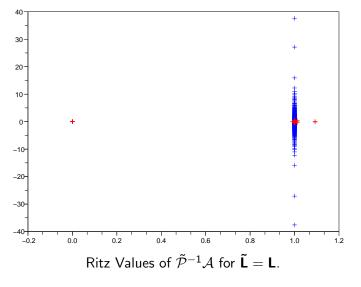


Converged Ritz values after 500 Arnoldi iterations are shown in blue. Unconverged values in red. >

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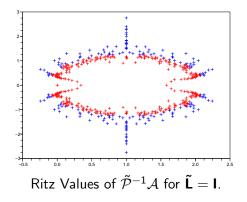
OOPS QG model. 24-hour window with 8 sub-windows.



Converged Ritz values after 500 Arnoldi iterations are shown in blue. Unconverged values in red. >

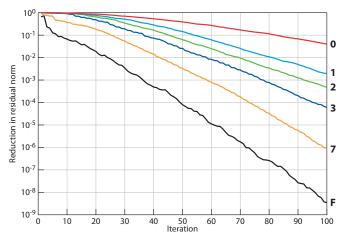
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- \bullet It is much harder to prove results for the case $\tilde{L}\neq L.$
- Experimentally, it seems that many eigenvalues continue to lie on $\Re(\tau) = 1$, with the remainder forming a cloud around $\tau = 1$.



Converged Ritz values after 500 Arnoldi iterations are shown in blue. Unconverged values in red

OOPS, QG model, 24-hour window with 8 sub-windows.

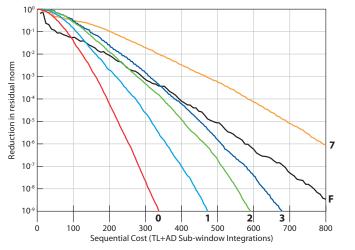


Convergence as a function of iteration for different truncations of the series expansion for L. ("F" = Forcing formulation.)

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OOPS, QG model, 24-hour window with 8 sub-windows.



Convergence as a function of sequential sub-window integrations for different truncations of the series expansion for L.

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Parallel 4D-Va

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Conclusions

- 4D-Var was analysed from the point of view of parallelization.
- 4D-PSAS and the forcing formulation are inherently sequential.
- The 4D-state problem is parallel, but ill-conditioning of **D** makes it difficult to precondition.
- The saddle point formulation is parallel, and seems easier to precondition.
 - A saddle point method for strong-constraint 4D-Var was proposed by Thierry Lagarde in his PhD thesis (2000: Univ Paul Sabatier, Toulouse). It didn't catch on:
 - Parallelization was not so important 10 year ago.
 - In strong constraint 4D-Var, we only get a factor-of-two speed-up. This is not enough to overcome the slower convergence due to the fact that the system is indefinite.
- The ability to also parallelize over sub-windows allows a much bigger speed-up in weak-constraint 4D-Var.
- The saddle point formulation is already fast enough to be useful.
- Better solvers and preconditioners can only make faster.