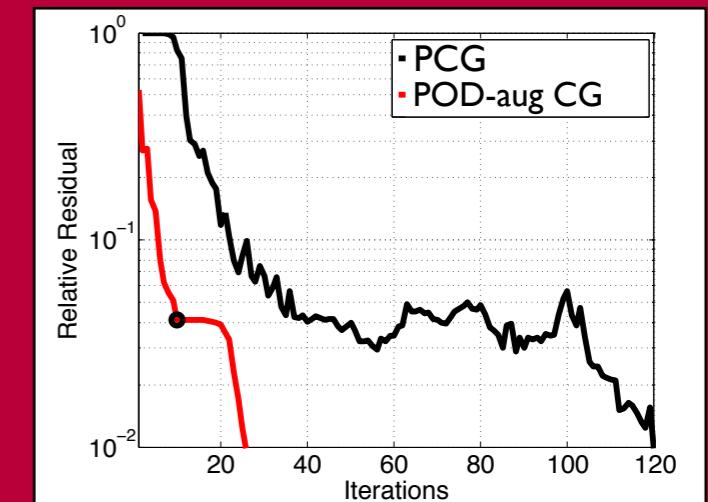
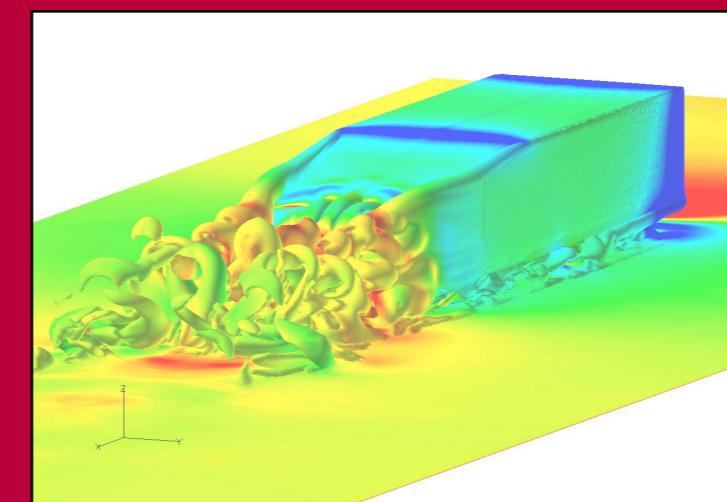
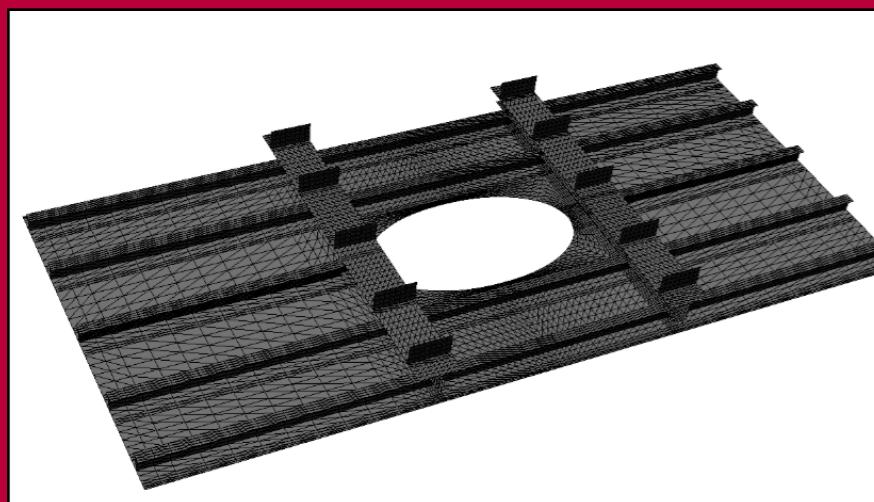


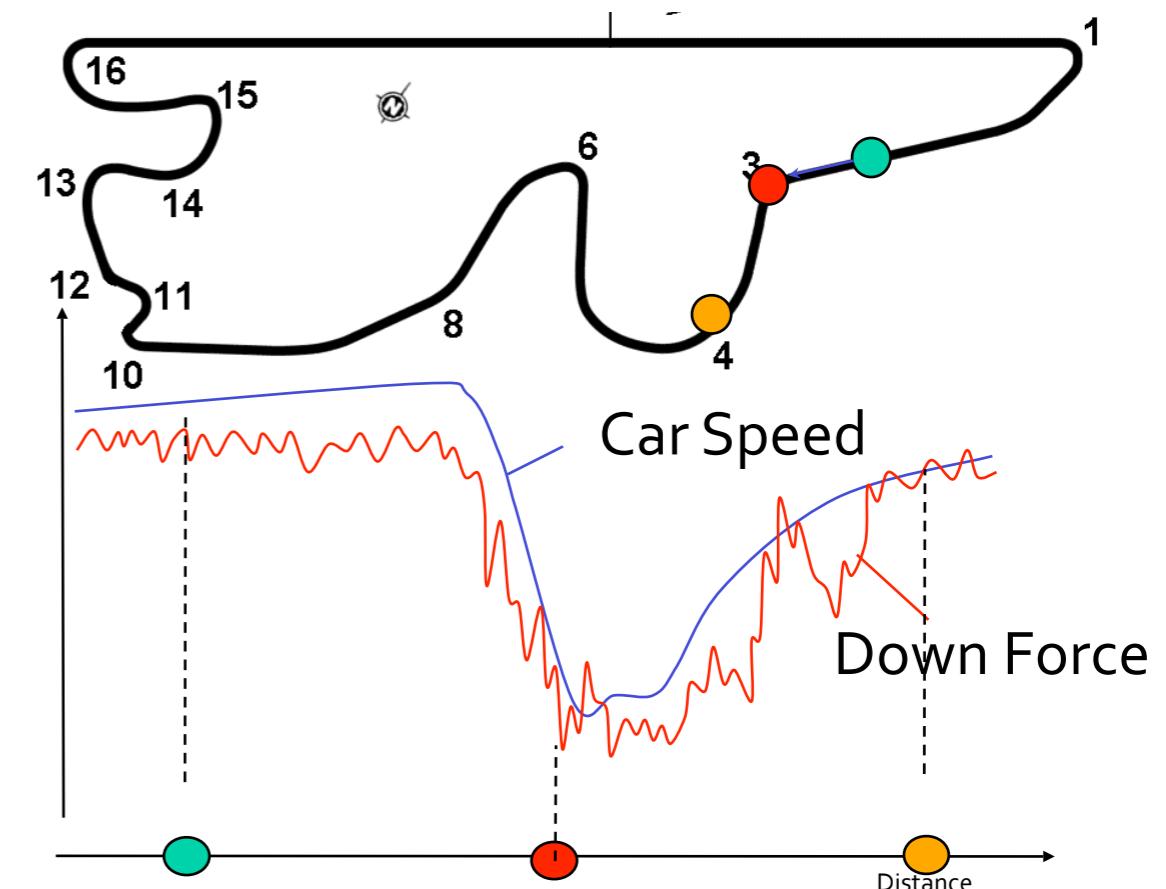


# Model reduction-based iterative methods for real-time simulation and repeated analyses of mathematical models



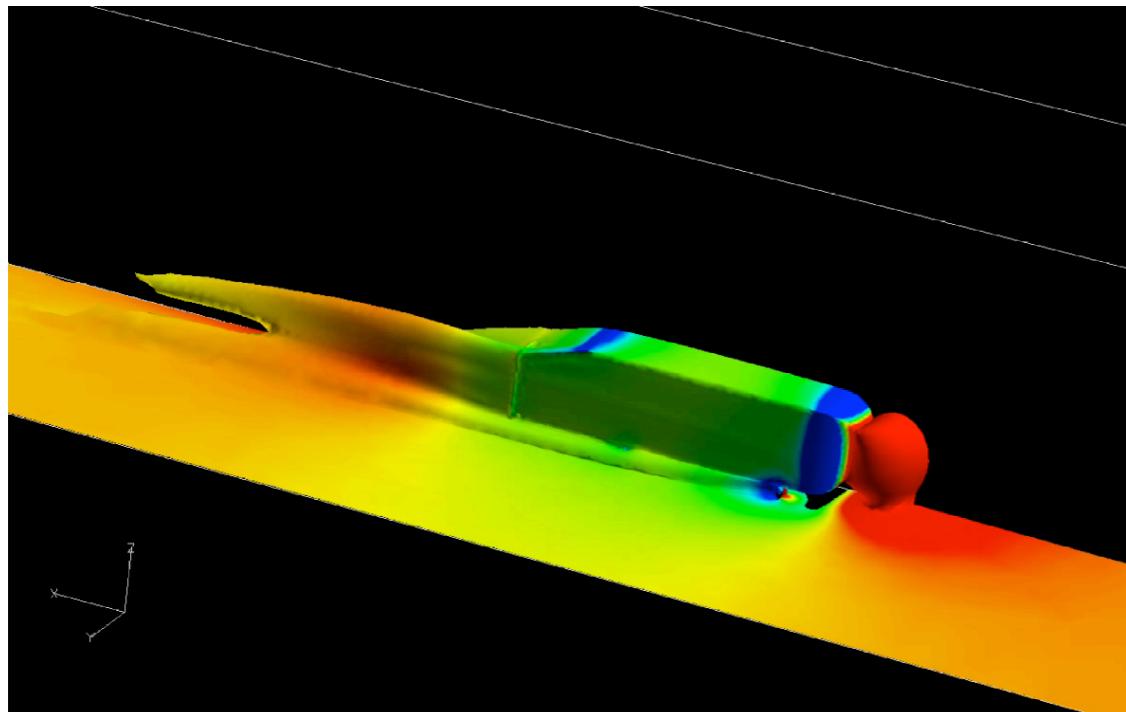
Kevin Carlberg and Charbel Farhat  
Stanford University  
Department of Aeronautics & Astronautics

Linear Algebra and Optimization Seminar  
Stanford, CA  
October 27, 2010

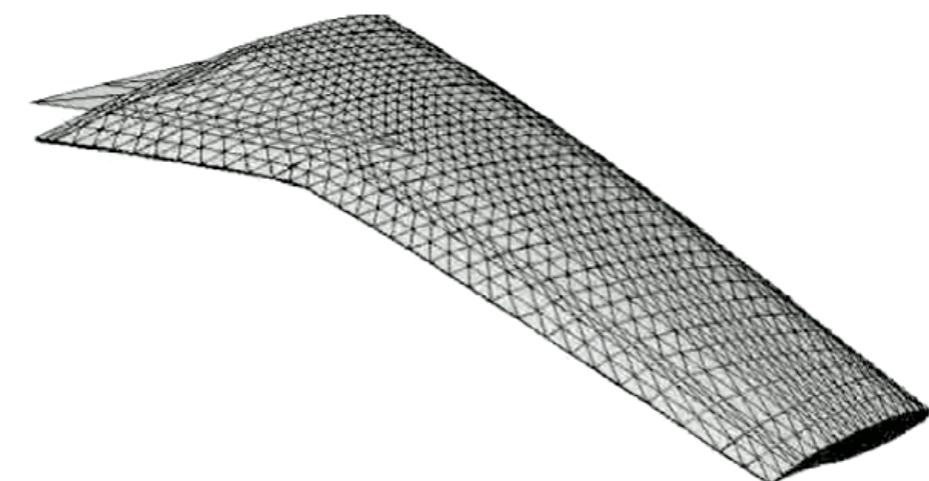


## Requirements

- Fast analysis during the season ("online")
- Can sacrifice some accuracy
- Can run expensive analyses before the season ("offline")



Nonlinear analysis



Design optimization

## Requirements

- Stringent accuracy requirement
- Want to minimize the overall computational cost of analyses



# Model reduction problem types



## Real-time analysis

- Fast-turnaround design
- “In-field” analysis
- Model predictive control

## Repeated analyses

- Nonlinear analysis
- Design optimization
- Parameter space sampling

Multiple objectives: **error** and **cost**

## Real-time analysis

Online cost more important:

minimize error  
subject to **online cost**  $\leq \tau$

## Repeated analyses

Error more important:

minimize **total** cost  
subject to error  $\leq \epsilon$



# Model reduction problem types



## Real-time analysis

- Fast-turnaround design
- “In-field” analysis
- Model predictive control

## Repeated analyses

- Nonlinear analysis
- Design optimization
- Parameter space sampling

Multiple objectives: **error** and **cost**

## Real-time analysis

Online cost more important:

minimize error  
subject to **online cost**  $\leq \tau$

## Repeated analyses

Error more important:

minimize **total** cost  
subject to error  $\leq \epsilon$

## 1) Offline

- Sample input space
- Build surrogate model

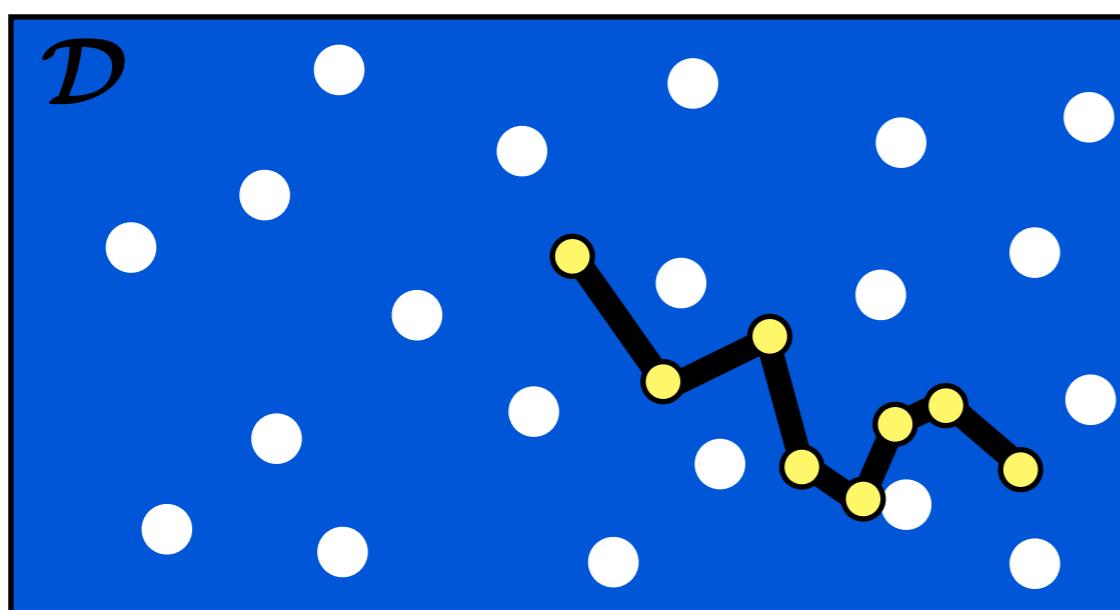
## 2) Online

- Analysis with surrogate model

### Real-time analysis

✓ High offline cost okay

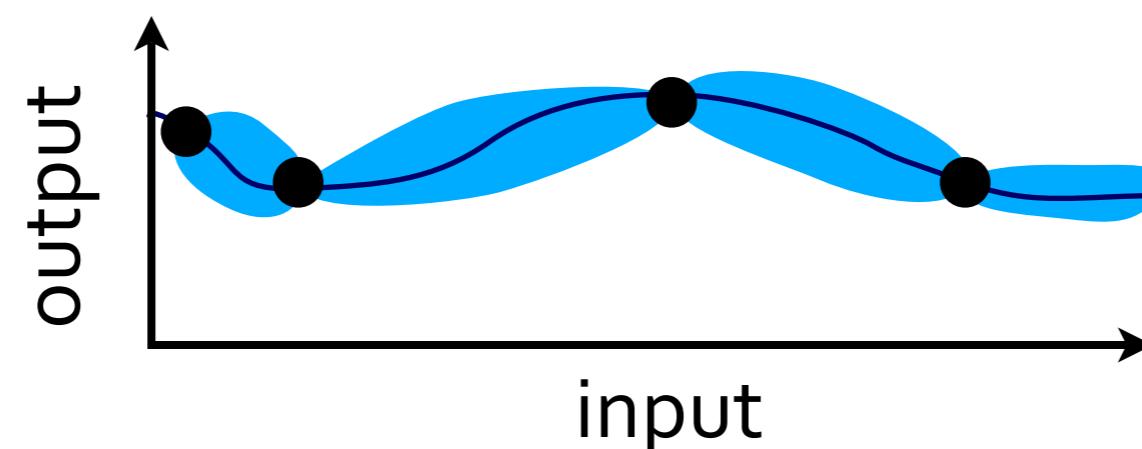
✓ Very low online cost



- $\mathcal{D}$  Input space
- Full-order model evaluation
- Surrogate model evaluation
- Online trajectory

## 1) Data fitting (response surfaces, Kriging)

- Directly model input-output map
  - ✗ “Blind” to problem physics
  - ✗ Impractical for large-dimensional input spaces



## 2) Reduced-order modeling

- Approximately solve state equations, then compute outputs
  - ✓ Robust: queries the physics online
  - ✓ Can apply arbitrary excitations/boundary conditions



# Reduced-order modeling



- Linear dynamical systems: “mature”

$$x^{n+1} = Ax^n + Bu^n$$

$$y^n = Cx^n + Du^n$$

SVD-based

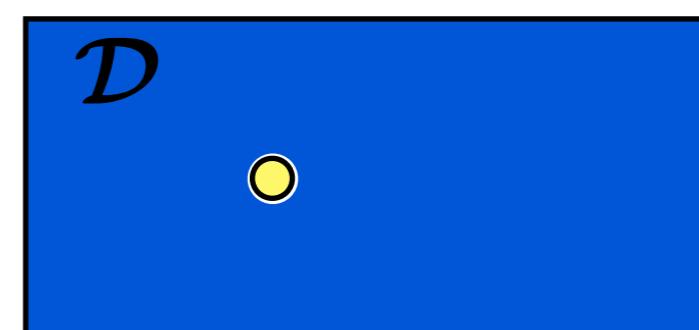
Krylov-based

- Balanced truncation
- Hankel norm approximation
- Balanced POD
- Partial realization
- Padé approximation
- Rational interpolation

- Nonlinear systems: major problems remain

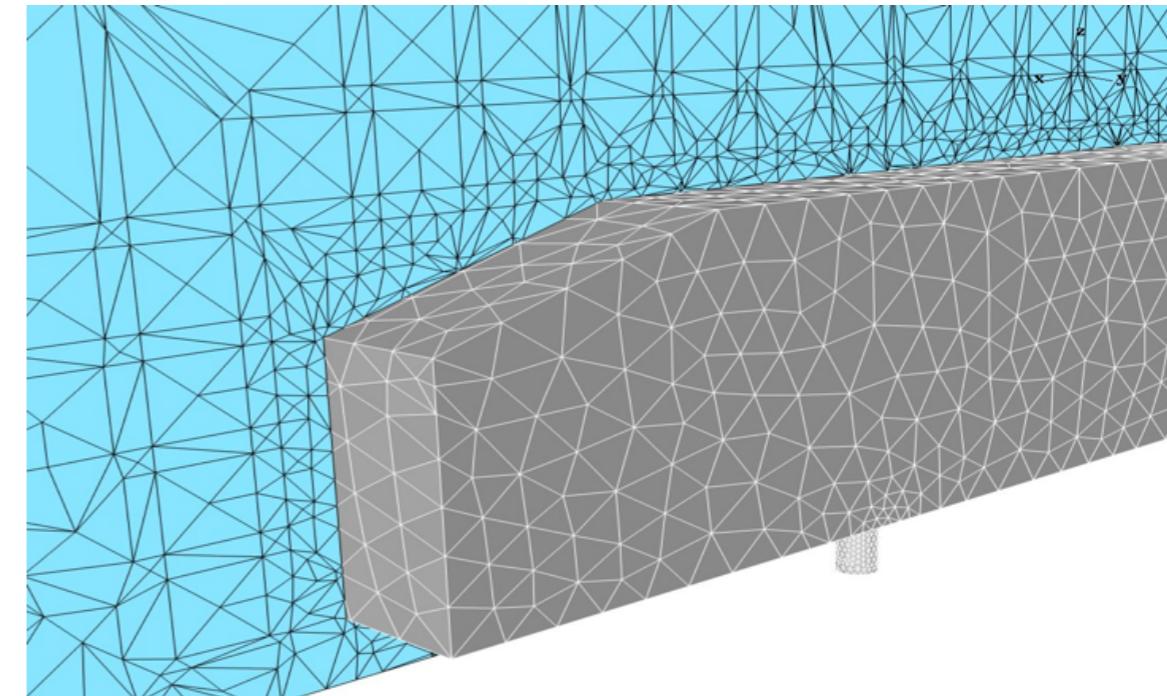
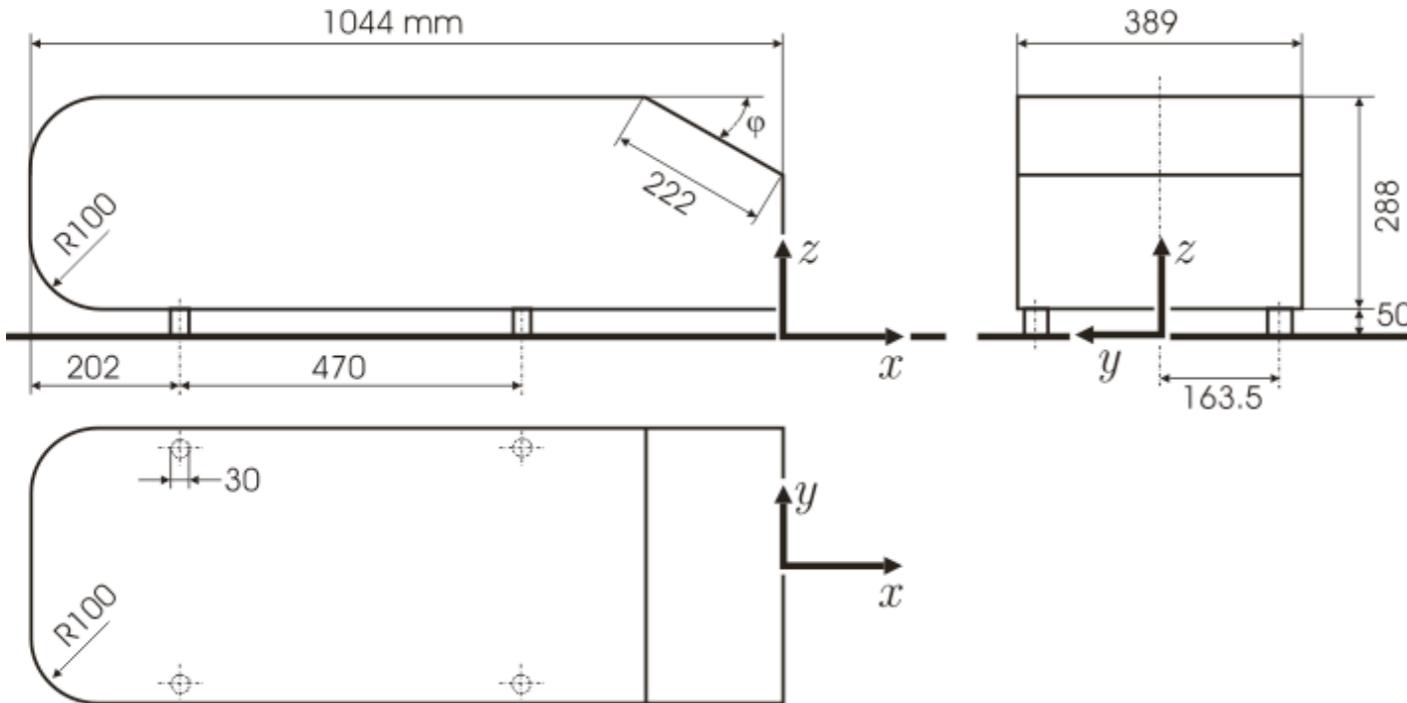
1. Accuracy

2. Computational cost

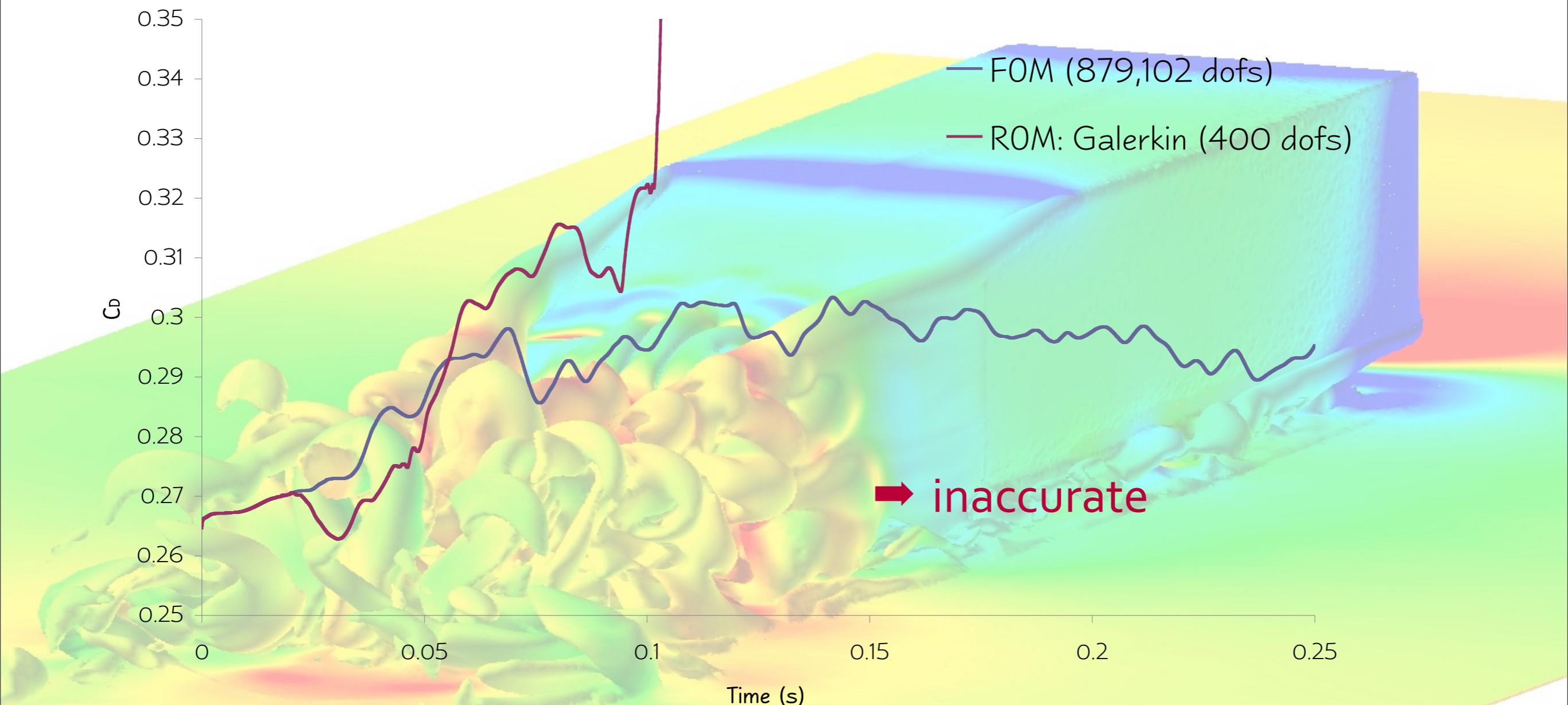




# Nonlinear example: Ahmed body



- Navier-Stokes simulation
  - $\phi = 20 \text{ deg}$
  - $V = 60 \text{ m/s}$
  - $\text{Re} = 4.29 \times 10^6$
  - DES turbulence model
  - 146,517 nodes
  - 837,894 tetrahedra
- Full-order model (FOM): 879,102 degrees of freedom
- POD/Galerkin reduced-order model (ROM): 400 degrees of freedom



- CPU time (first 0.1 s of simulated time)
- FOM: 9,292 s
- ROM POD/Galerkin: 24,984 s

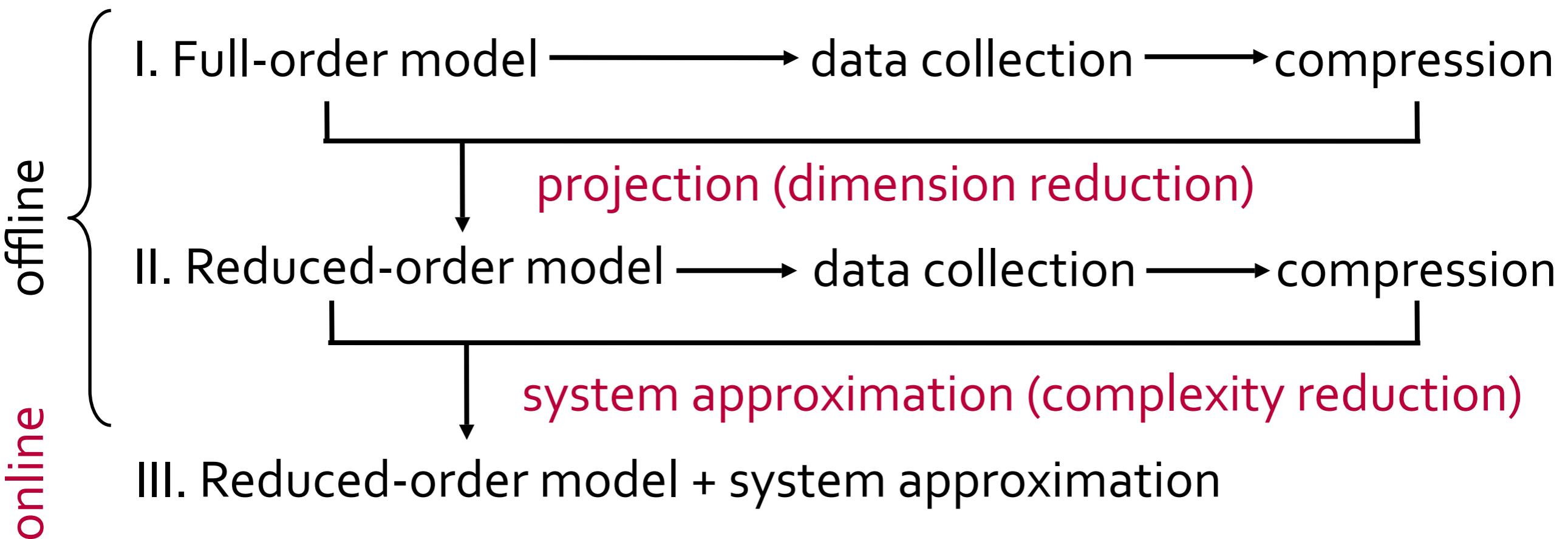
→ computationally expensive



# Methodology



- Consecutively introduce approximations that satisfy 2 properties



1. **Consistency:** In the limit of **no compression**, reproduce exactly the solution of the previous model for sampled problems
2. **Optimality:** Minimize an error measure with respect to the previous model (*a priori* convergence)



# Mathematical framework



- Nonlinear PDE

$$\mathcal{L}(u; x, t) = 0$$

- Semi-discretized PDE in space (ODE)

$$\mathcal{L}^d(u; t) = 0$$

- Fully-discretized PDE with implicit time integration

$$R(y^{n+1}; y^n, \dots, y^0; t^n) = 0$$

- FOM: sequence of nonlinear systems of equations of general form

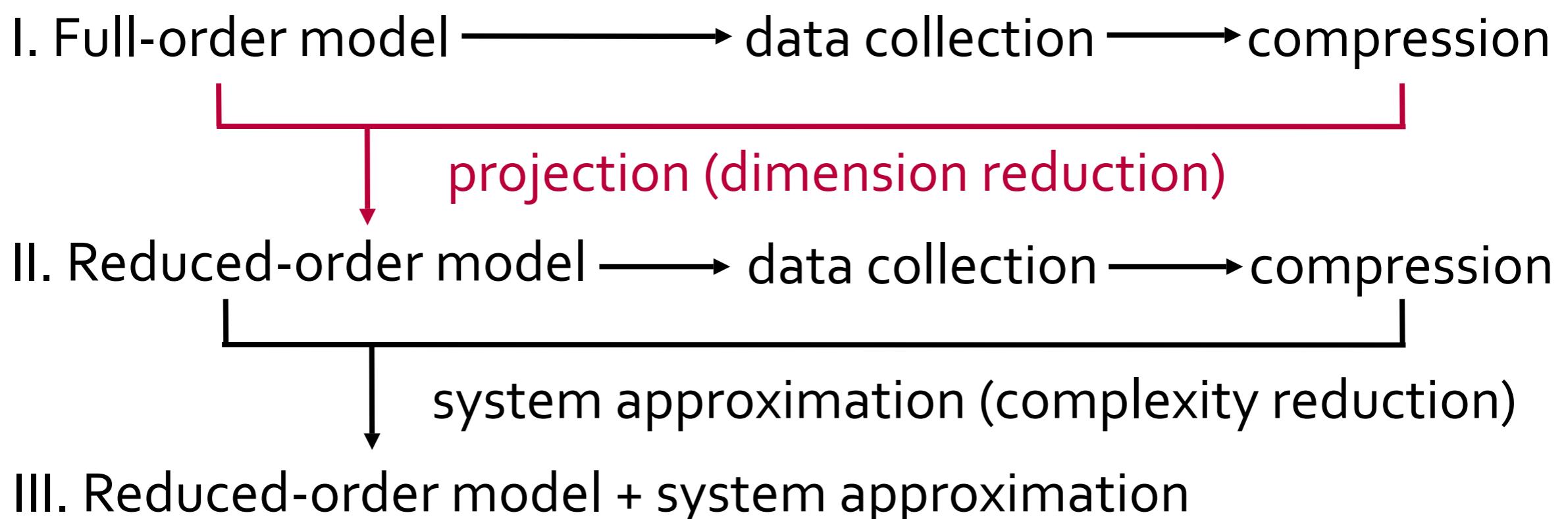
$$R(y) = 0$$

- $R : \mathbb{R}^N \rightarrow \mathbb{R}^N, (w) \mapsto R(w)$

- dimension N “large”



# Projection





# Petrov-Galerkin Projection



- To decrease the dimension, search for a solution in the affine subspace:  $y^{(0)} + \text{range}(\Phi_y)$  of dimension  $n_y \ll N$

$$y \approx y^{(0)} + \Phi_y y_r$$

- Enforce residual orthogonal to  $n_y$ -dimensional subspace  $\text{range}(\Psi)$

$$\Psi^T R(y^{(0)} + \Phi_y y_r) = 0$$

- Solve reduced system by Newton's method for  $k=1, \dots, K$  (converged)

$$\Psi^T J^{(k)} \Phi_y p^{(k)} = -\Psi^T R^{(k)}$$

$$y_r^{(k+1)} = y_r^{(k)} + \alpha^{(k)} p^{(k)}$$

$$\triangleright R^{(k)} \equiv R(y^{(0)} + \Phi_y y_r^{(k)}), \quad J^{(k)} \equiv \frac{dR}{dw} \left( y^{(0)} + \Phi_y y_r^{(k)} \right)$$



## Proposition

The projection approximation is **consistent** if:

1.  $\Phi_y$  is a **POD basis** computed with snapshots ( $y - y^{(0)}$ ) collected during the evaluation of Model I
2.  $y$  is sufficiently close to  $y^{(0)}$
3. The reduced Jacobian  $\Psi^T \frac{dR}{dw}(\cdot) \Phi_y : \Omega \rightarrow \mathbb{R}^{n_y \times n_y}$  is Lipschitz continuous, with  $\Omega \subset \mathbb{R}^N$  an open set containing  $y - y^{(0)}$
4. The reduced Jacobian at the solution  $\Psi^T \frac{dR}{dw}(y) \Phi_y$  is nonsingular

- 1: Defines data and procedure for computing  $\Phi_y$
- 2–4: Newton's method convergence conditions



# Proper orthogonal decomposition



- Given  $n_x$  “snapshots”  $x^i$ , the first  $k$  POD vectors satisfy:

$$\text{span}\{\phi^i\}_{i=1}^k = \arg \min_{\mathcal{S} \in \mathcal{G}(k, N)} \sum_{j=1}^{n_x} \| (I - \Pi_{\mathcal{S}}) x_j \|^2$$

- $\Pi_{\mathcal{S}}$ : orthogonal projection onto  $\mathcal{S}$
- $\mathcal{G}(k, N)$ : set of  $n$ -dimensional subspaces of  $\mathbb{R}^N$
- $\Phi = [\phi^1 \dots \phi^k]$ : POD basis in matrix form
- Properties
  - “No compression”:  $k = n_x$  and  $\text{range}(\Phi) = \text{range}([x^1 \dots x^{n_x}])$
  - Efficient computation by singular value decomposition (SVD)



# Projection optimality



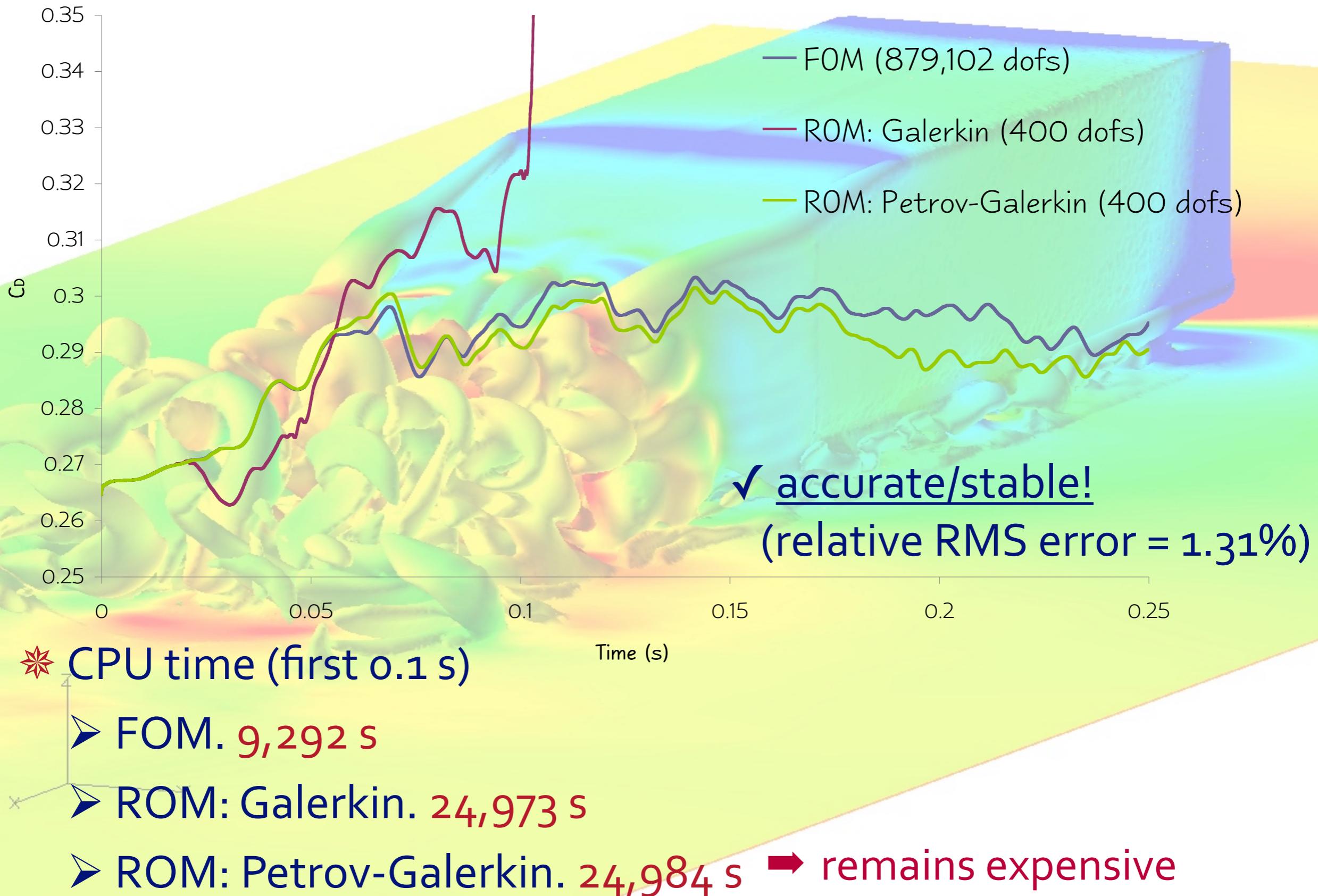
- Want solution  $p^{(k)}$  to satisfy for some norm  $\Theta$

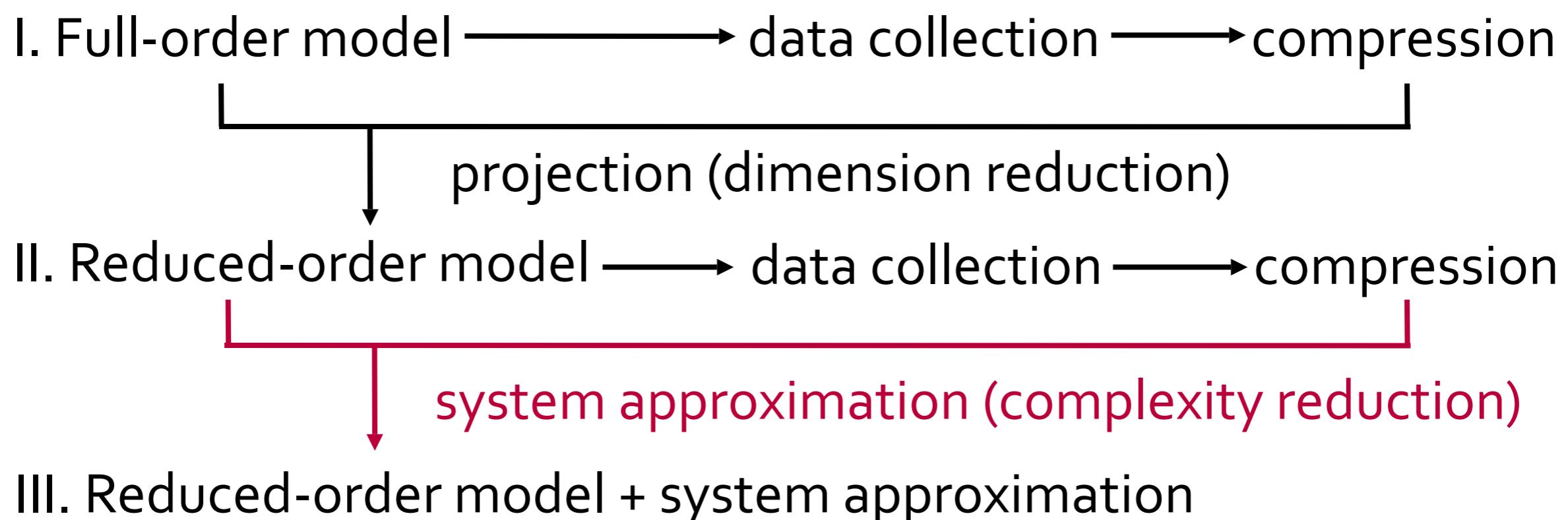
$$p^{(k)} = \arg \min_{p \in \mathbb{R}^{n_y}} \|\Phi_y p - (J^{(k)})^{-1} R^{(k)}\|_\Theta \quad (1)$$

- Galerkin:  $\Psi = \Phi_y$  satisfies (1) with  $\Theta = J^{(k)}$  only if  $J^{(k)}$  SPD
- Petrov-Galerkin:  $\Psi = J^{(k)} \Phi_y$  satisfies (1) with  $\Theta = J^{(k)\top} J^{(k)}$  if  $J^{(k)}$  nonsingular
- Least-squares Petrov-Galerkin projection
  - Equivalent to **globally convergent Gauss-Newton** method for
$$\min \|R(y + \Phi_y y_r)\|_2$$
  - $\Psi$  is state-dependent, changes each Newton iteration
  - Can only be introduced on fully discrete system



# Ahmed body: ROM results





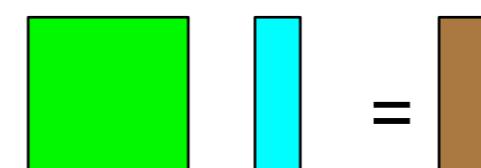


# Computational bottleneck



- Reduced-order Newton iterations are

$$[J^{(k)}\Phi_y]^T [J^{(k)}\Phi_y] p^{(k)} = -[J^{(k)}\Phi_y]^T R^{(k)}$$



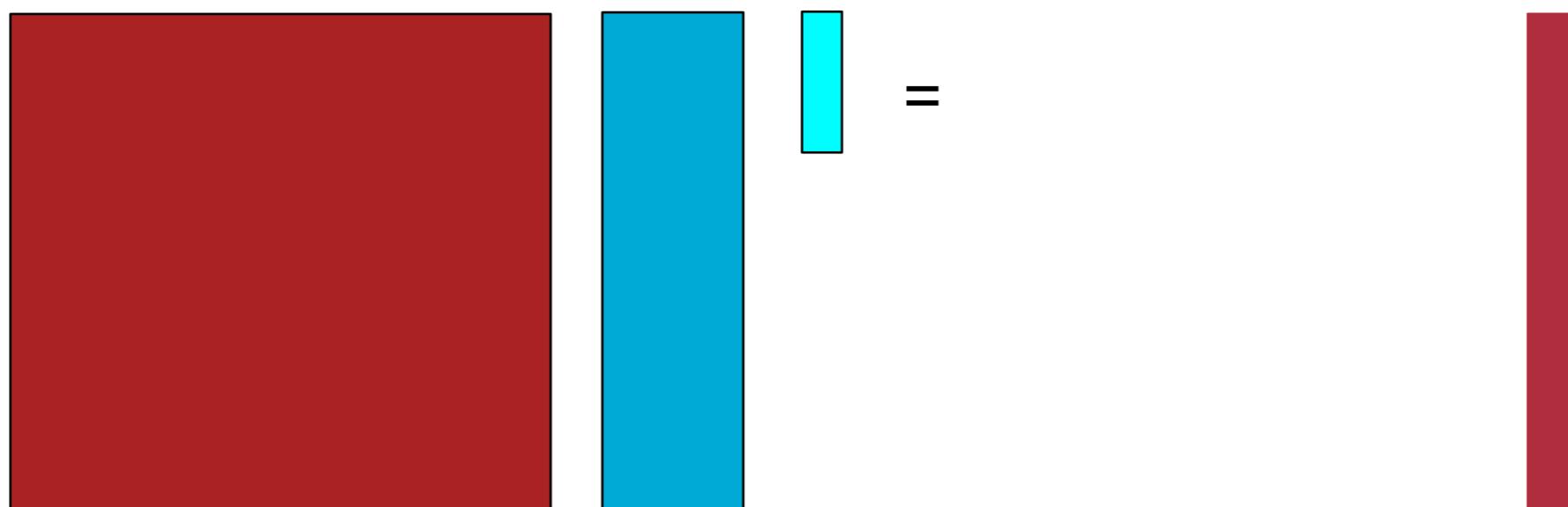


# Computational bottleneck



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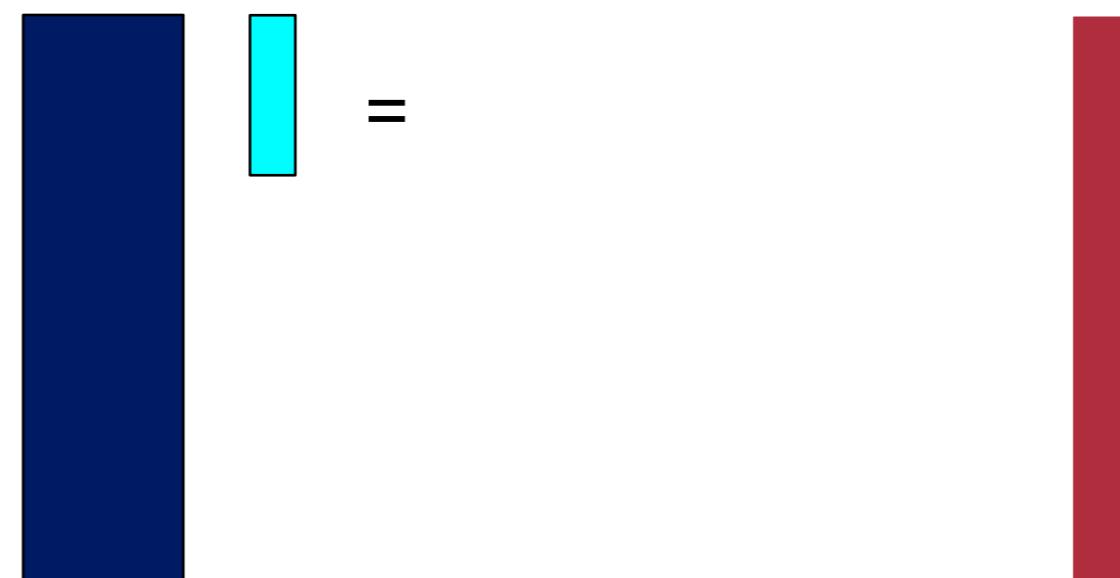


# Computational bottleneck



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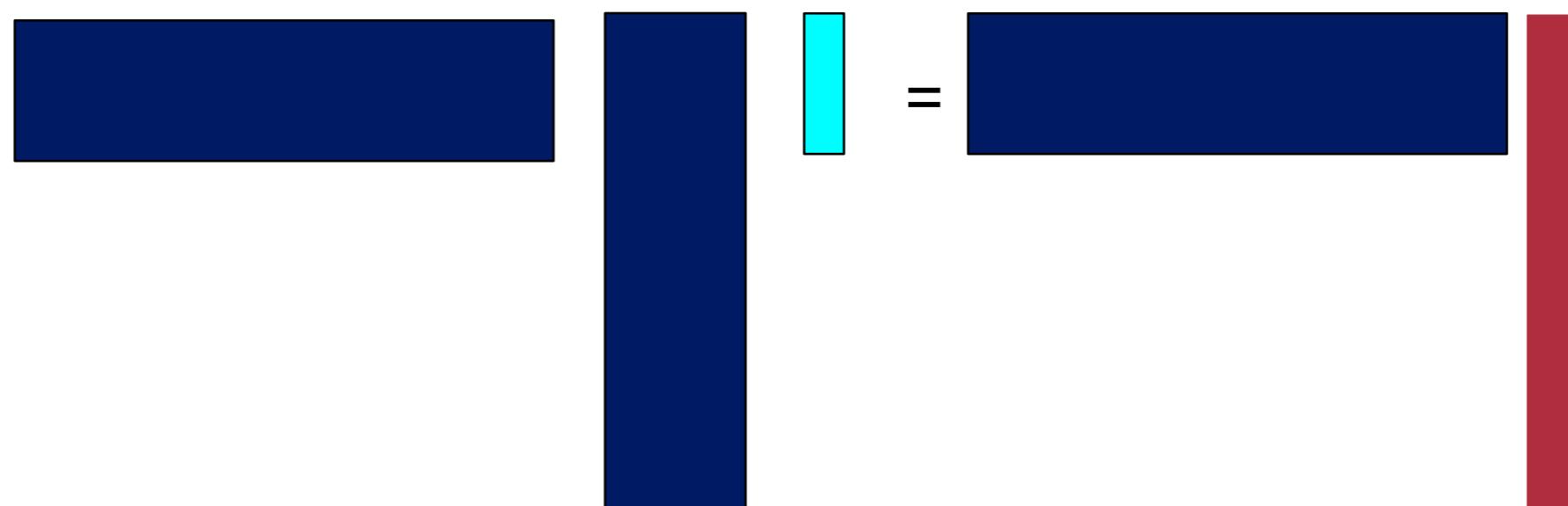


# Computational bottleneck



- Reduced-order Newton iterations are

$$[J^{(k)}\Phi_y]^T [J^{(k)}\Phi_y] p^{(k)} = -[J^{(k)}\Phi_y]^T R^{(k)}$$



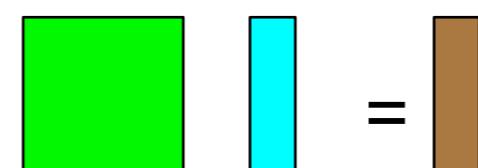


# Computational bottleneck



- Reduced-order Newton iterations are

$$[J^{(k)}\Phi_y]^T [J^{(k)}\Phi_y] p^{(k)} = -[J^{(k)}\Phi_y]^T R^{(k)}$$



- Operation count at each iteration scales with large dimension N
- Expensive even though reduced equations are of small dimension

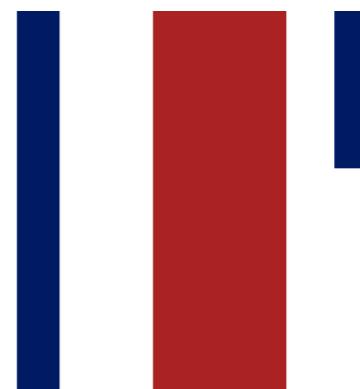


# System approximation



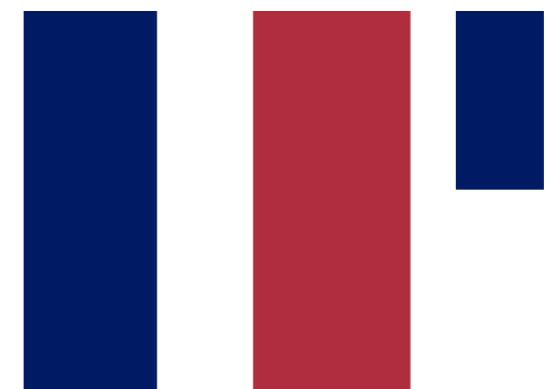
- Tensor approximations:  $R^{(k)} \in \text{range}(\Phi_R)$ ,  $J^{(k)} \Phi_y \in \text{range}(\Phi_J)$

$$R^{(k)} \approx \Phi_R R_r^{(k)}$$



invariant  
iteration-dependent

$$J^{(k)} \Phi_y \approx \Phi_J J_r^{(k)}$$

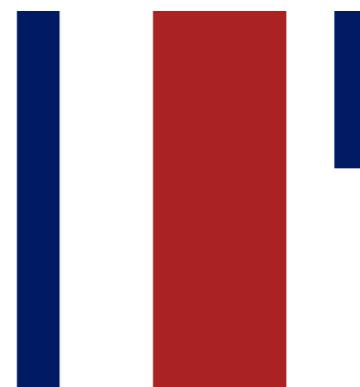


- Newton iterations become

$$[J^{(k)} \Phi_y]^T [J^{(k)} \Phi_y] p^{(k)} = -[J^{(k)} \Phi_y]^T R^{(k)}$$

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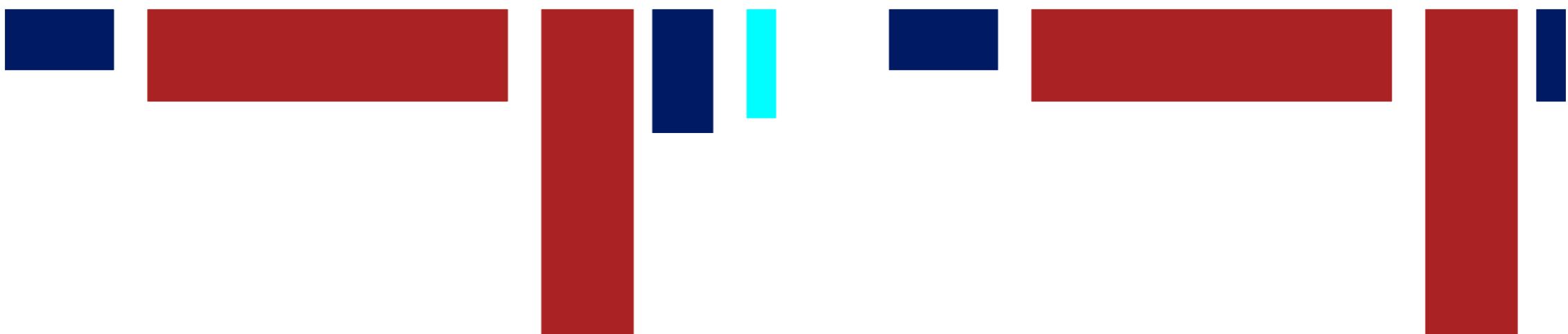
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invariant  
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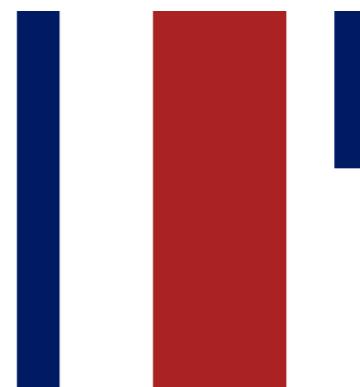
- Newton iterations become

$$(J_r^{(k)})^T \Phi_J^T \Phi_J J_r^{(k)} p^{(k)} = -(J_r^{(k)})^T \Phi_J^T \Phi_R R_r^{(k)}$$



- Tensor approximations:  $R^{(k)} \in \text{range}(\Phi_R)$ ,  $J^{(k)} \Phi_y \in \text{range}(\Phi_J)$

$$R^{(k)} \approx \Phi_R R_r^{(k)}$$



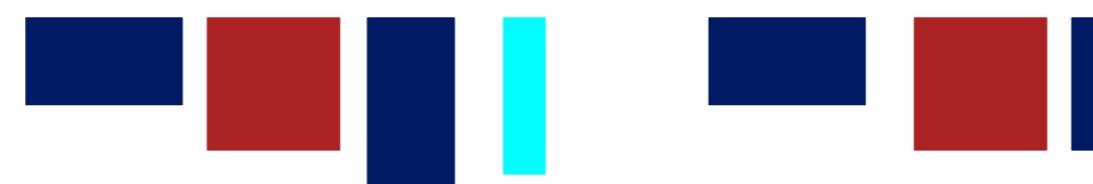
$$J^{(k)} \Phi_y \approx \Phi_J J_r^{(k)}$$



invariant  
iteration-dependent

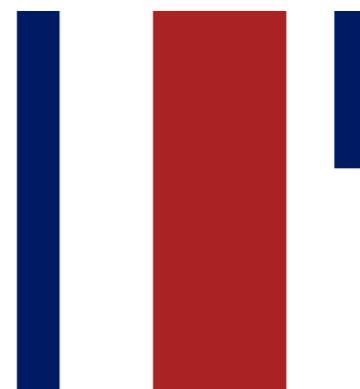
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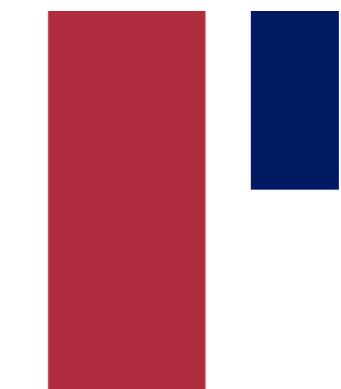


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$$R^{(k)} \approx \Phi_R R_r^{(k)}$$



$$J^{(k)} \Phi_y \approx \Phi_J J_r^{(k)}$$



invariant  
iteration-dependent

- Newton iterations become

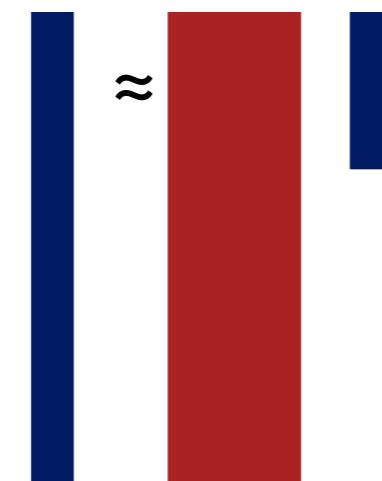
$$(J_r^{(k)})^T \Phi_J^T \Phi_J J_r^{(k)} p^{(k)} = -(J_r^{(k)})^T \Phi_J^T \Phi_R R_r^{(k)}$$



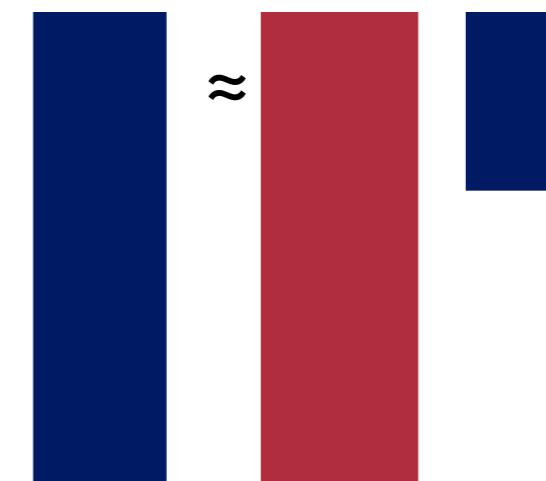
- Form  $\Phi_J^T \Phi_J$ ,  $\Phi_J^T \Phi_R$  offline: online operations independent of  $N$ !
- Similar: Barrault et al., 2004; Grepl et al., 2007; Nguyen & Peraire, 2008; Chatarantabut & Sorensen, 2009; Galbally et al., 2010

- Tensor approximations

$$R^{(k)} \approx \Phi_R R_r^{(k)}$$



$$J^{(k)} \Phi_y \approx \Phi_J J_r^{(k)}$$



- Property 1: Optimality

- $R_r^{(k)}, J_r^{(k)}$  computed online by gappy data reconstruction

- Property 2: Consistency

- $\Phi_R, \Phi_J$  computed offline by POD with specific snapshots



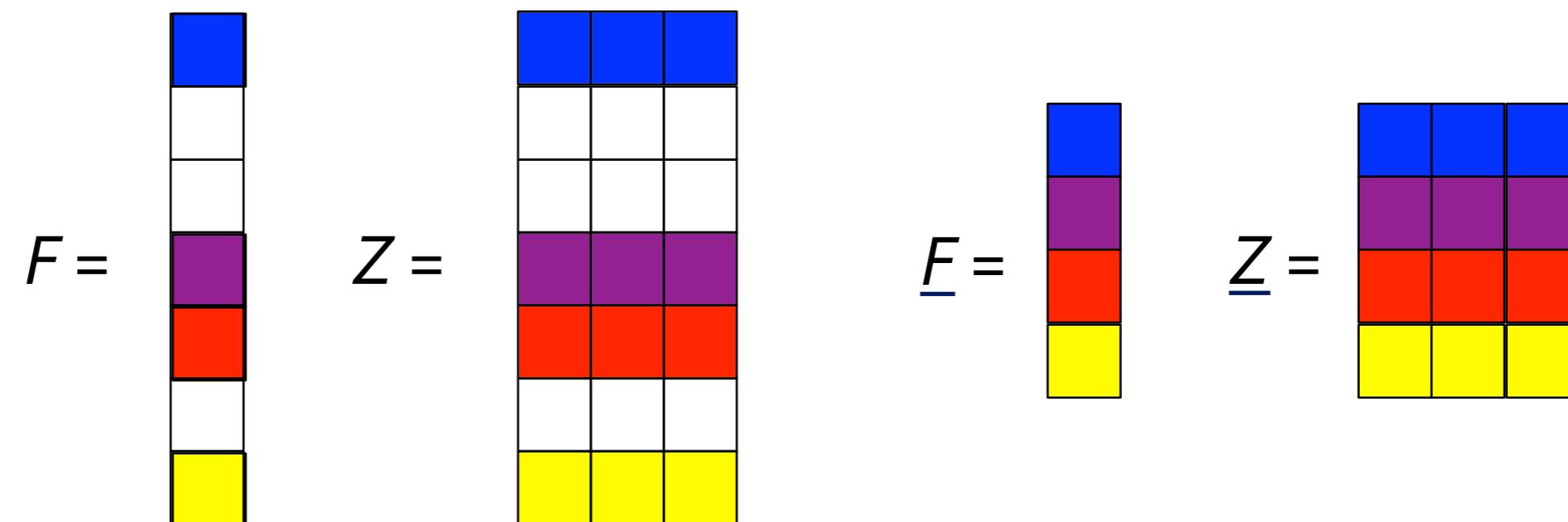
# Gappy data reconstruction

- Goal: accurately reconstruct vector  $F \in \mathbb{R}^N$  [Everson & Sirovich, 1995]
- Given: 1) a basis  $Z = [z^1 \dots z^p]$   
2) some sampled entries of the vector  $F_i, i \in I$  with  $|I| = q \geq p$  and  $q \ll N$
- Define restriction:  $\underline{F} \equiv [F]_i, i \in I$   
 $\underline{Z} \equiv [z^1 \dots z^p]$

---

Example

$N=7$   
 $p=3$   
 $q=4$   
 $I=\{1,4,5,7\}$





# Gappy data reconstruction



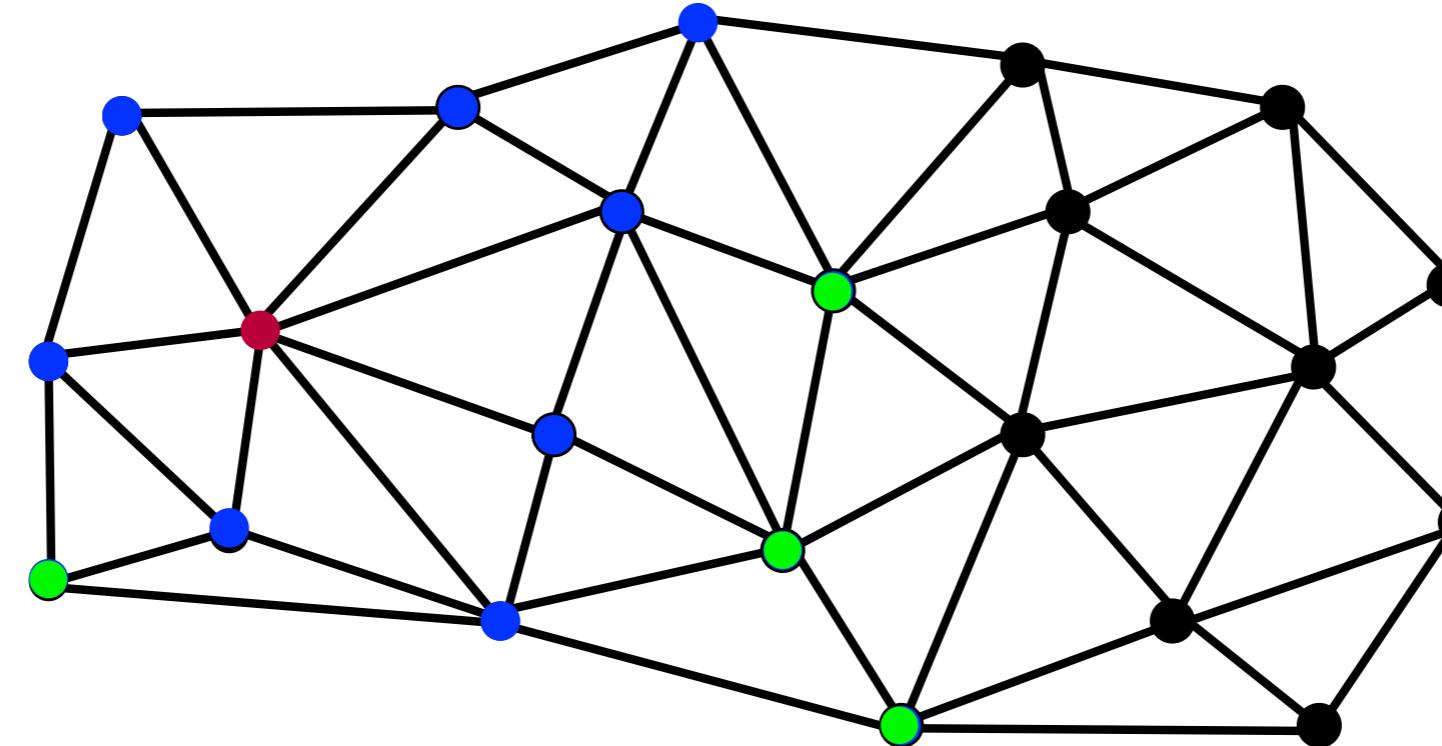
- Least squares minimization on sampled entries

$$\underline{F}_r = \arg \min_{x \in \mathbb{R}^p} \|\underline{Z}\underline{x} - \underline{E}\|_2$$

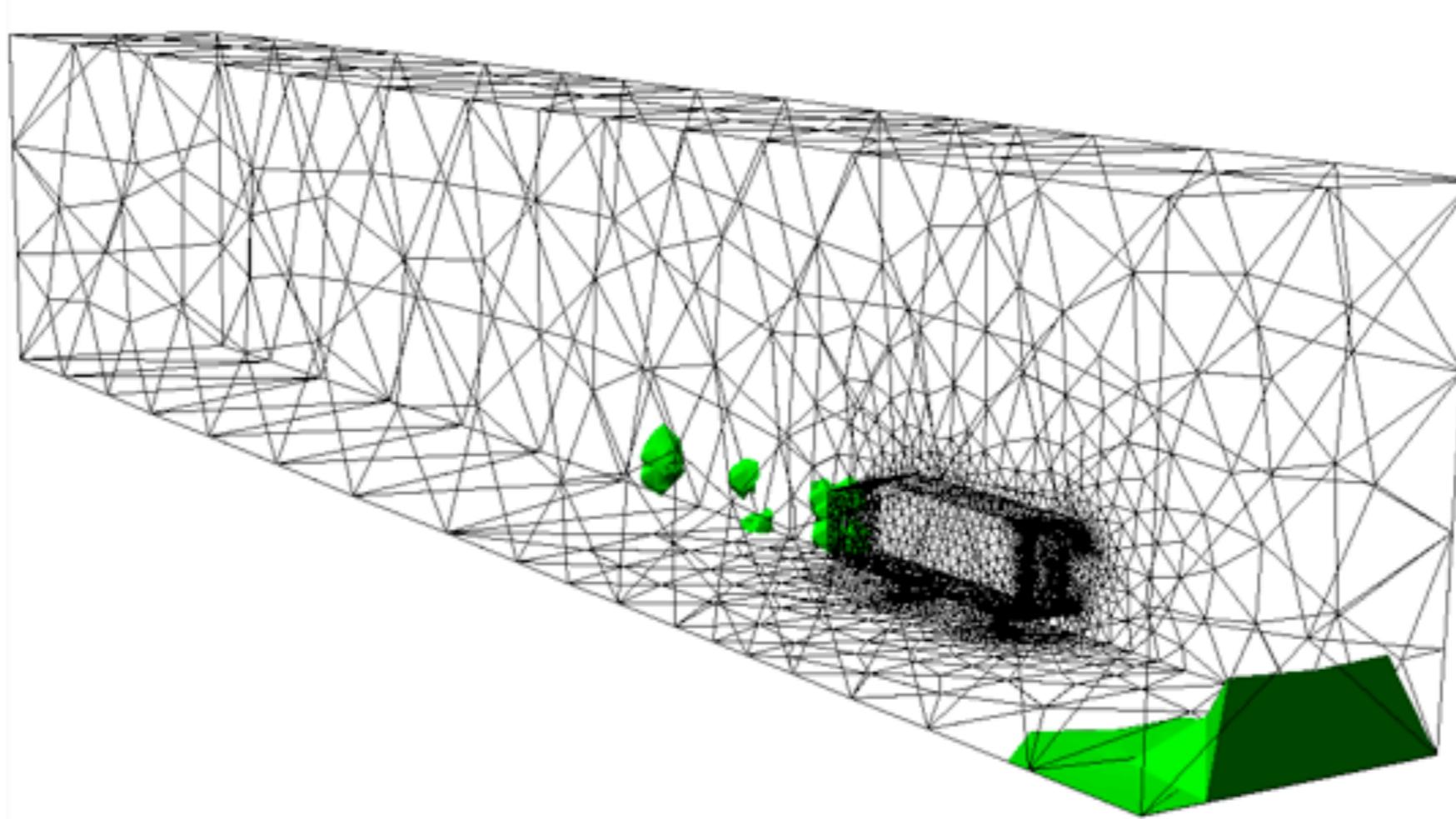
- Reconstruct all entries of the vector

$$\underline{F} \approx \underline{Z}\underline{F}_r$$

- Optimality
  - error  $\|\underline{Z}\underline{F}_r - \underline{E}\|_2$  monotonically decreases as p increases
- Apply at each Newton iteration to reconstruct
  1.  $R^{(k)}$ , with  $F_r = R_r^{(k)}$  and  $Z = \Phi_R$
  2.  $J^{(k)} \Phi_y$ , with  $F_r = J_r^{(k)}$  and  $Z = \Phi_J$



- $R^{(k)}$  and  $J^{(k)} \Phi_y$  computed at **sample nodes**
- To do this, must compute the state at:
  1. **Sample nodes**
  2. **Neighbors** of sample nodes
  3. **Neighbors of neighbors** of sample nodes
- In general, this depends on the Jacobian's sparsity pattern



- Only compute the physics in isolated **mesh clusters**
- Gappy POD “fills in” the rest of the mesh



# System approximation consistency



## Proposition

The system approximation is **consistent** if  $\Phi_R$  and  $\Phi_J$  are POD bases computed with snapshots that satisfy:

1.  $R^{(k)}$  from the Model II simulation is a snapshot for  $\Phi_R$
2.  $J^{(k)} \Phi_y p^{(k)}$  from the Model II simulation is a snapshot for  $\Phi_J$ ,
3. Each column of  $J^{(k)} \Phi_y$  from the Model II simulation is a snapshot for  $\Phi_J$

- Leads to hierarchy of snapshot collection procedures characterized by tradeoffs between consistency and offline cost/storage



# Snapshot collection hierarchy



ID	0	1	2	3
Snapshots for $y$	$y_I - y_I^{(o)}$	$y_I - y_I^{(o)}$	$y_I - y_I^{(o)}$	$y_I - y_I^{(o)}$
Snapshots for $R^{(k)}$	$R_I^{(k)}$	$R_{II}^{(k)}$	$R_{II}^{(k)}$	$R_{II}^{(k)}$
Snapshots for $J^{(k)}\Phi_y$	$R_I^{(k)}$	$R_{II}^{(k)}$	$[J^{(k)}\Phi_y p^{(k)}]_{II}$	$[J^{(k)}\Phi_y]_{II}$
# simulations	1	2	2	2
# snapshots per Newton iteration	1	1	2	$1 + n_y$
consistency conditions satisfied	none	1	1, 2	1, 2, 3

- $(\cdot)_I, (\cdot)_{II}$ : snapshot saved during Model I, II
- $(\cdot)^{(k)}$ : snapshot saved at each Newton iteration



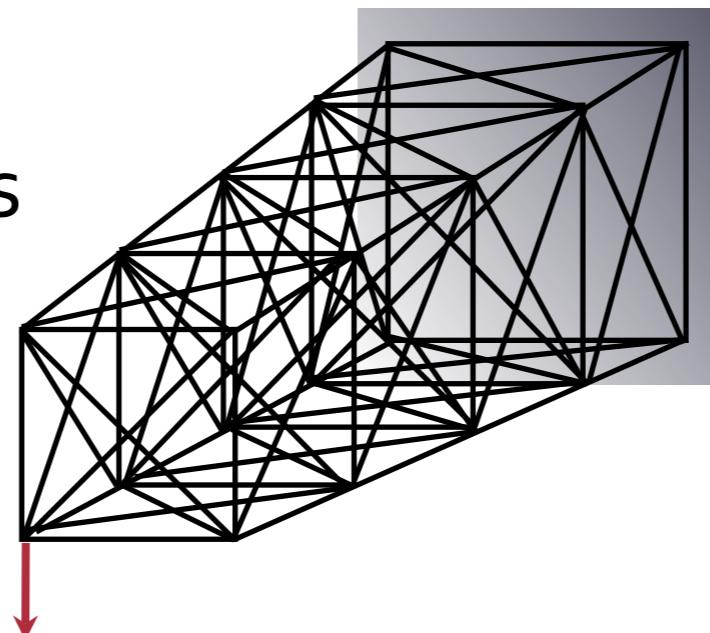
# Snapshot collection hierarchy

ID	0	1	2	3
Snapshots for $y$	$y_I - y_I^{(o)}$	$y_I - y_I^{(o)}$	$y_I - y_I^{(o)}$	$y_I - y_I^{(o)}$
Snapshots for $R^{(k)}$	$R_I^{(k)}$	$R_{II}^{(k)}$	$R_{II}^{(k)}$	$R_{II}^{(k)}$
Snapshots for $J^{(k)}\Phi_y$	$R_I^{(k)}$	$R_{II}^{(k)}$	$[J^{(k)}\Phi_y p^{(k)}]_{II}$	$[J^{(k)}\Phi_y]_{II}$
# simulations	1	2	2	2
# snapshots per Newton iteration	1	1	2	$1 + n_y$
consistency conditions satisfied	none	1	1, 2	1, 2, 3

- ✗ Procedure 0: satisfies no consistency conditions
- ✗ Procedure 3: consistent, but prohibitive cost
- ✓ Procedure 2: similar cost as 1, more consistency conditions

- Geometrically nonlinear structural dynamics

16 nonlinear 3D bars  
per bay



- Full-order model: 12,000 degrees of freedom (1000 bays)

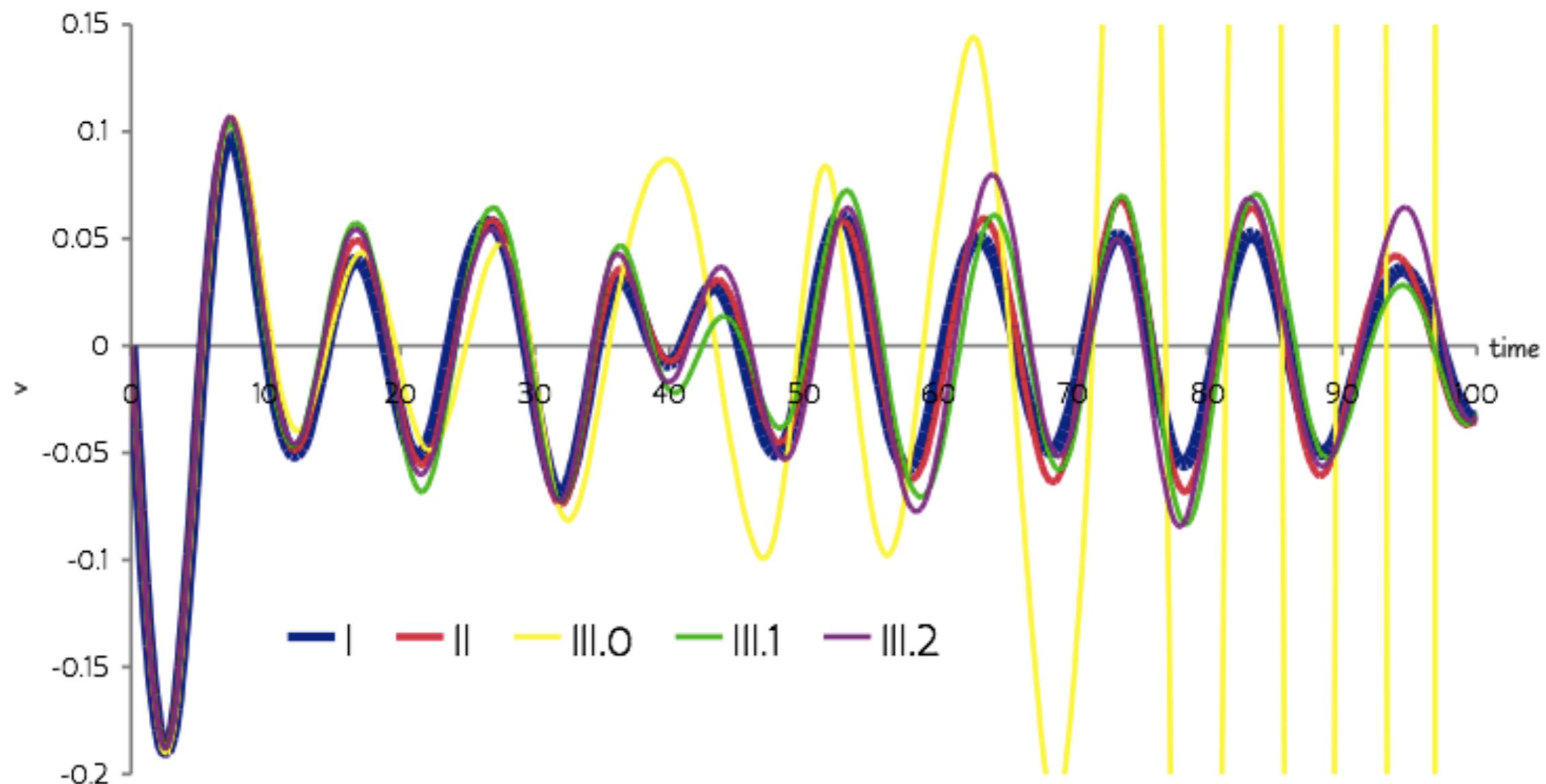
- Output of interest: downward velocity at tip

- Truncation of POD bases

➤  $\Phi_y$ :  $n_y = 26$  (99% of  $y_I - y_I^{(o)}$  snapshot energy)

➤  $\Phi_R, \Phi_y$ :  $n_R = n_J = 28$  (99% of  $R_I^{(k)}$  snapshot energy)

- $|I|=28$  (interpolation)



- III.0, which satisfies no consistency conditions, is unstable



# Consistency importance

# consistency  
conditions  
satisfied

Model	$e_1$ (%)	Total Newton it	Speed-up over I
II	3.44	601	1.03
III.o	unstable	—	—
III.1	6.31	603	158
III.2	6.01	891	122

- Model II: Good accuracy, but insufficient speed-up
- Model III.o: inaccurate, perhaps because not consistent
- Models III.1, III.2: accurate, promising speed-ups
- Meeting more consistency conditions improves accuracy/stability



# Least-squares v. interpolation

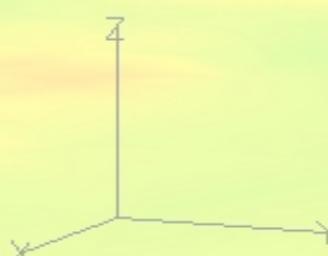
- $n_y=26, n_R=n_J=28$

		Relative error $e_l$ (%)			Speed-up over Model I		
		III.0	III.1	III.2	III.0	III.1	III.2
		28	unstable	6.31	6.01	120	158
least squares	42	4.95	3.77	3.63	137	152	119
	56	4.03	3.50	3.44	129	136	116

- Increasing ||| improves accuracy, even stabilizing III.0
  - Least-squares reconstruction performs better than interpolation



- Conclusions
  - Carefully build approximations to be consistent & optimal
  - Least-squares Petrov-Galerkin led to **accuracy**
  - System approximation led to **low computational complexity**
- Future work
  - Currently implementing in AERO-F flow solver





# Repeated analyses



## Real-time analysis

- Fast-turnaround design
- “In-field” analysis
- Model predictive control

## Repeated analyses

- Nonlinear analysis
- Design optimization
- Parameter space sampling

Multiple objectives: **error** and **cost**

## Real-time analysis

Online cost more important:

minimize error  
subject to **online cost**  $\leq \tau$

## Repeated analyses

Error more important:

minimize **total** cost  
subject to error  $\leq \epsilon$

## 1) Offline

- Sample input space
- Build surrogate model

Real-time analysis

✓ High offline cost  
okay

Repeated analyses

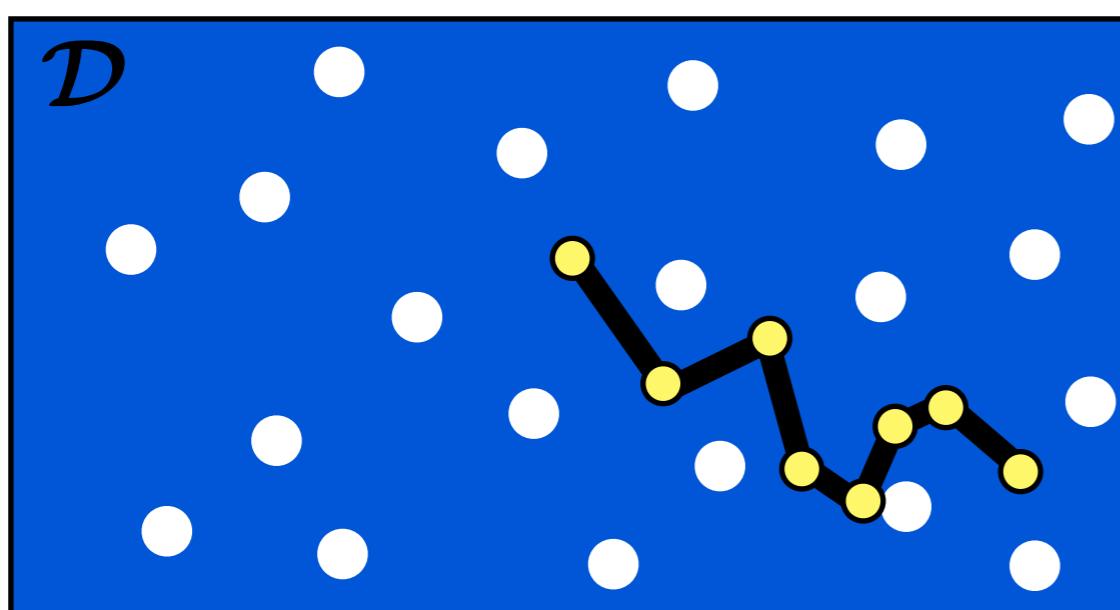
✗ May preclude  
total cost savings

## 2) Online

- Analysis with surrogate

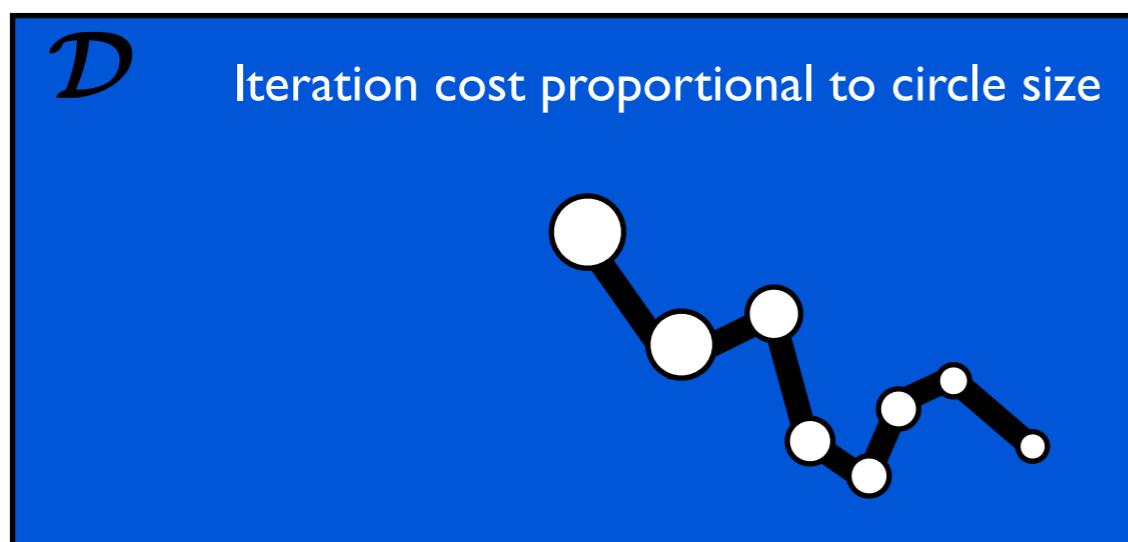
✓ Very low online cost

✗ May not be  
accurate enough



- $\mathcal{D}$  Input space
- Full-order model evaluation
- Surrogate model evaluation
- Online trajectory

- Guiding philosophy
  1. Avoid extra computations
  2. Fully exploit data generated from previous analyses
  3. Use data to accelerate solver convergence to required tolerance
- Procedure
  1. Execute only required analyses (no offline sampling)
  2. Build a POD basis on-the-fly
  3. Use the POD basis within an iterative method to accelerate analyses



$\mathcal{D}$  Input space  
○ Accelerated analyses  
— Trajectory



# Repeated analyses



- For  $i=1, \dots, n_{mat}$  and  $j=1, \dots, n_{RHS}$ , solve

$$A^{(i)} x^{(i,j)} = b^{(i,j)}$$

- $A^{(i)} \in \mathbb{R}^{N \times N}$  sparse, **SPD**, and “nearby”
- Solve each system by preconditioned conjugate gradient (PCG)

$$\tilde{x}^{(i,j)} = \arg \min_{v \in \mathcal{S}} \|x^{(i,j)} - v\|_{A^{(i)}}$$

- $\mathcal{S} = \mathcal{K}_k$ :  $k$ -dimensional Krylov subspace of  $\mathbb{R}^N$
- Current PCG acceleration approach: **augmented CG methods**



# Augmented CG methods



$$\mathcal{S} = \mathcal{K}_k + \mathcal{Y}$$

- Choices of  $\mathcal{Y}$

1. All previous search directions

- Multi-RHS [O'Leary, 1980; Saad, 1987; Farhat *et al.*, 1994; Erhel, *et al.*, 2000]
- Multi-matrices [Rey, 1994; Roux, 1995; Farhat, 1995; Farhat *et al.*, 2000, Risler *et al.*, 2000]
  - ✗ Large cost/storage for many linear systems

2. Approximated extreme eigenvectors (deflation)

- Multi-RHS [Chapman *et al.*, 1997; Saad *et al.*, 2000]
- Multi-matrices [Rey *et al.*, 1998; Parks *et al.*, 2007]
  - ✗ Only effective if a few eigenvalues hamper convergence

- Goals: 1) Feasibility

- 2) Effective in the general case



# Proper orthogonal decomposition



- Given 1)  $n_w$  snapshots  $w_j$ , 2)  $n_w$  weights  $\gamma_j$ , 3) a  $\Theta$ -norm
- The first  $n \leq n_w$  POD vectors satisfy:

$$\text{span}\{\phi_i\}_{i=1}^n = \arg \min_{S \in \mathcal{G}(n, N)} \sum_{j=1}^{n_w} \| (I - \Pi_S^\Theta) \gamma_j w_j \|_\theta^2$$

- $\Pi_S^\Theta$ :  $\Theta$ -orthogonal projection onto  $S$
- $\mathcal{G}(n, N)$ : set of  $n$ -dimensional subspaces of  $\mathbb{R}^N$
- $\Phi(n) \equiv [\phi_1 \ \cdots \ \phi_n]$ :  $n$ -dimensional POD basis in matrix form
- Key properties
  1. Optimal ordering
    - First  $n$  POD basis vectors span optimal  $n$ -dimensional space
  2.  $\Theta$ -orthonormality:  $\Phi(n)^T \Theta \Phi(n) = I$



# POD strategy



- Align POD with minimizing the error for a nearby “target system”

$$\bar{A}\bar{x}^{(j)} = \bar{b}^{(j)}$$

- To do this, choose three ingredients to be:
  1. Snapshots  $w_j$ : all previous search directions
  2. Weights  $\gamma_j$ : estimate the solution for the target problem

$$\bar{x}^{(j)} \approx \bar{x}_{\text{est}}^{(j)} = \sum_{j=1}^{n_w} \gamma_j w_j$$

- 3.  $\Theta$  : matrix for the target problem  $\bar{A}$



- Compute one POD basis for each RHS

$$\Phi_j(n) \equiv [\phi_1^{(j)} \ \cdots \ \phi_n^{(j)}]$$

- Key properties

1. Optimal ordering

→ POD vectors optimally ordered to minimize **error at target**

2.  $\bar{A}$  -orthonormality

$$\Phi_j(n)^T \bar{A} \Phi_j(n) = I$$

→  $\Phi_j(n)^T A^{(i)} \Phi_j(n) \approx I$  for  $A^{(i)}$  near  $\bar{A}$



# POD-augmented CG algorithm



$$\mathcal{S} = \mathcal{K}_k + \text{range}(\Phi_j(n))$$

1. Directly solve  $n_1$ -dimensional reduced equations ( $n_1$  small)

$$\Phi_j(n_1)^T A^{(i)} \Phi_j(n_1) \hat{x} = \Phi_j(n_1)^T b^{(i,j)}$$

$$\tilde{x}_1^{(i,j)} = \Phi_j(n_1) \hat{x}$$

‣ Accurate (Property 1) and low cost ( $n_1$  small)

2. Iteratively solve  $n$ -dimensional reduced equations ( $n > n_1$ )

$$\boxed{\Phi_j(n)^T A^{(i)} \Phi_j(n) \hat{x} = \Phi_j(n)^T \left( b^{(i,j)} - A^{(i)} \tilde{x}_1^{(i,j)} \right)}$$

$$\tilde{x}_2^{(i,j)} = \tilde{x}_1^{(i,j)} + \Phi_j(n_1) \hat{x}$$

- Use augmented CG without forming reduced matrix
- More accurate (Property 1) and low cost (Property 2)



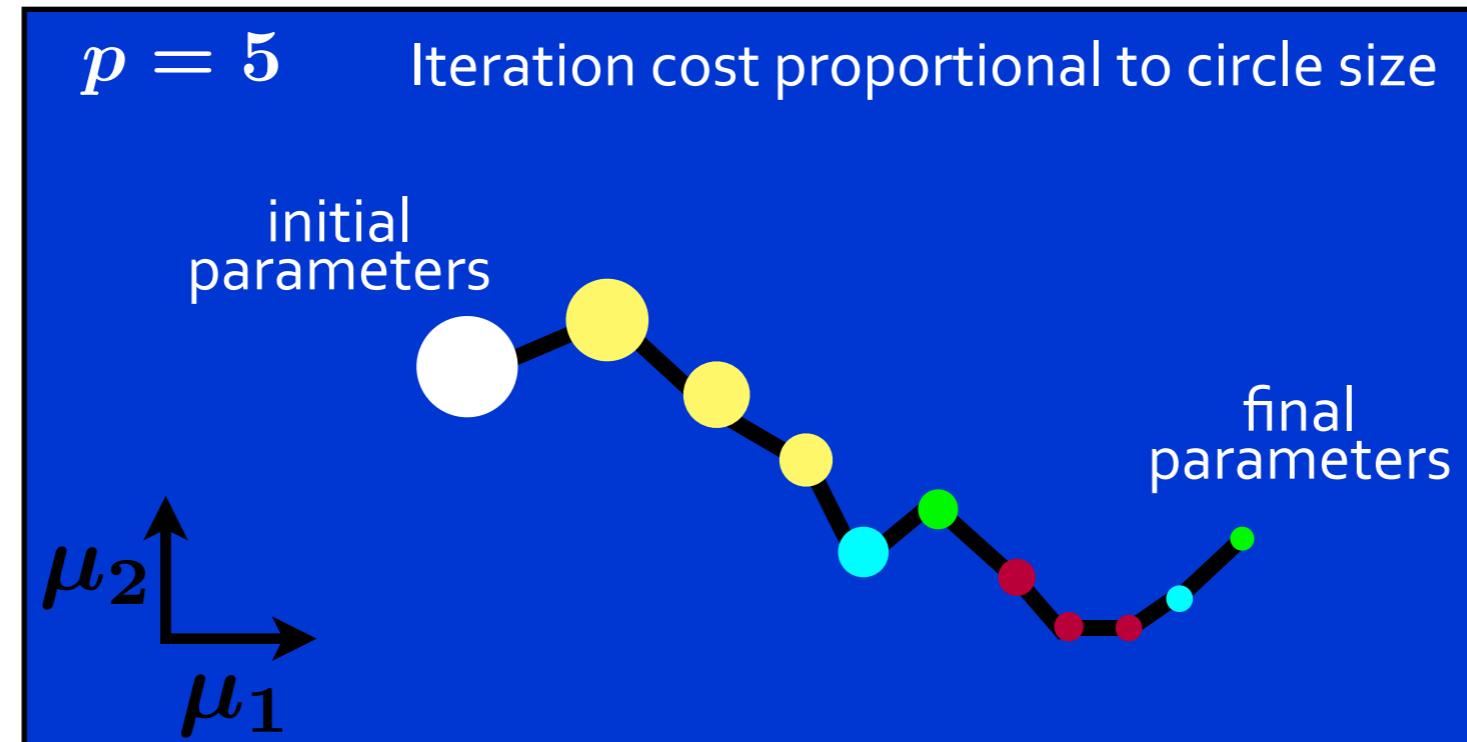
### 3. Iteratively solve full state equations to specified tolerance

$$A^{(i)} \hat{x} = b^{(i,j)} - A^{(i)} \tilde{x}_2^{(i,j)}$$

$$\tilde{x}^{(i,j)} = \tilde{x}_2^{(i,j)} + \hat{x}$$

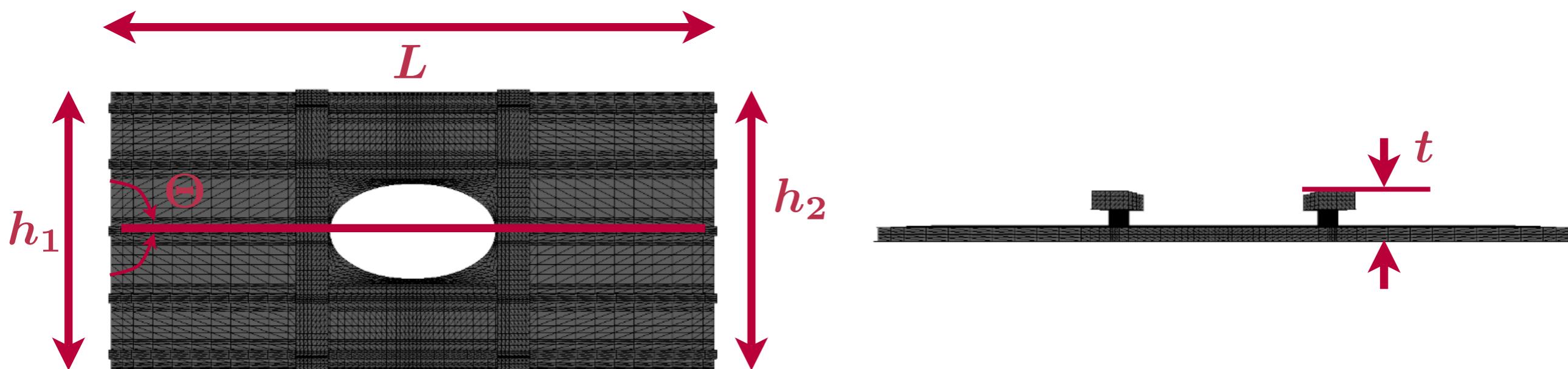
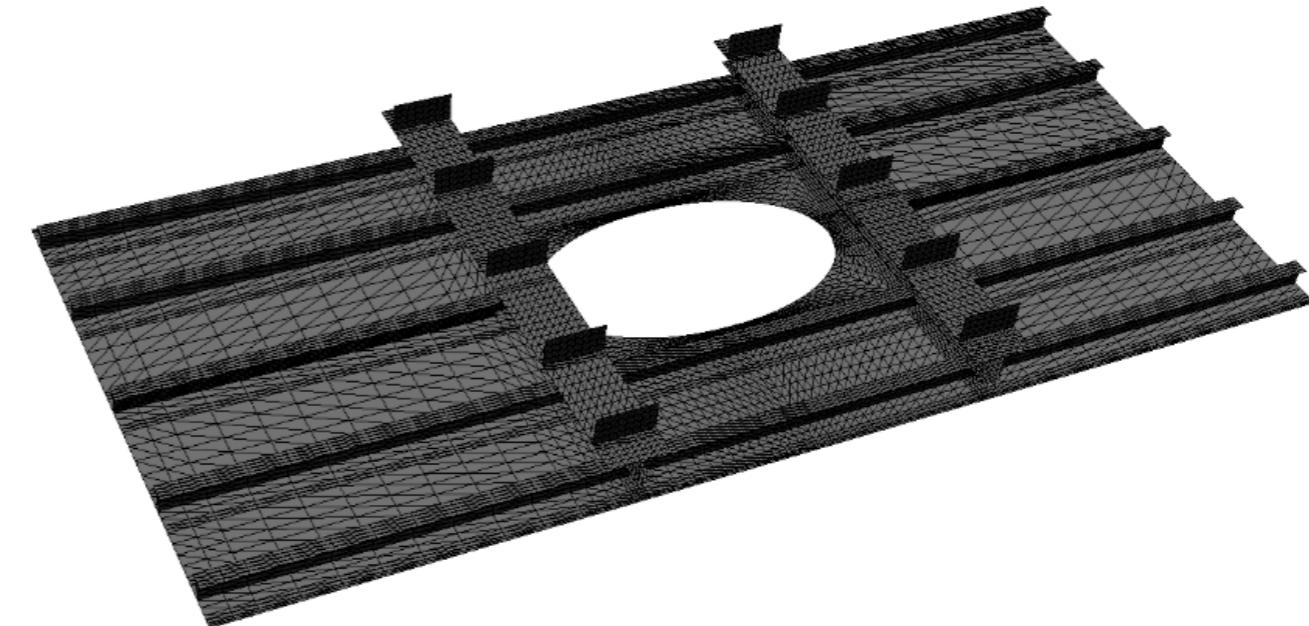
- Maintain  $A^{(i)}$ -conjugacy to old search directions using augmented PCG [Farhat *et al.*, 1994]
- Multiple-RHS
  - Sequentially execute Stages 1-3 for  $j=1, \dots, n_{RHS}$
  - Stage 1 approximation space includes search directions from all previous RHS

- $p = \#$  mat before recomputing POD basis (data compression)



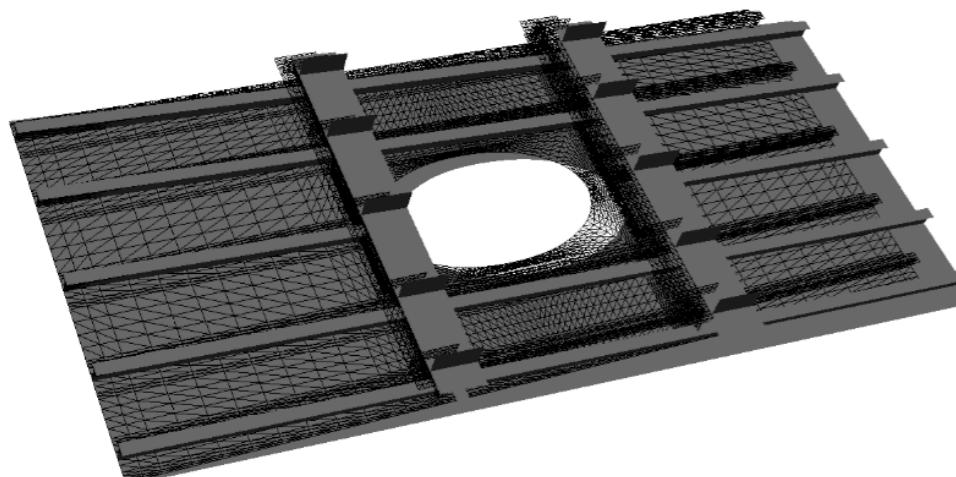
	$i$	Stage 1 basis	Stage 2 basis	Compute POD?
○	1			
●	$2 : p - 1$	$W$		
○	$np$	$W$		✓
●	$np + 1$	$\Phi(n_1)$	$\Phi(n)$	
●	$np + 2 : (n + 1)p - 1$	$[W, \Phi(n_1)]$	$[W, \Phi(n)]$	

# Example: V-22 Osprey wing panel

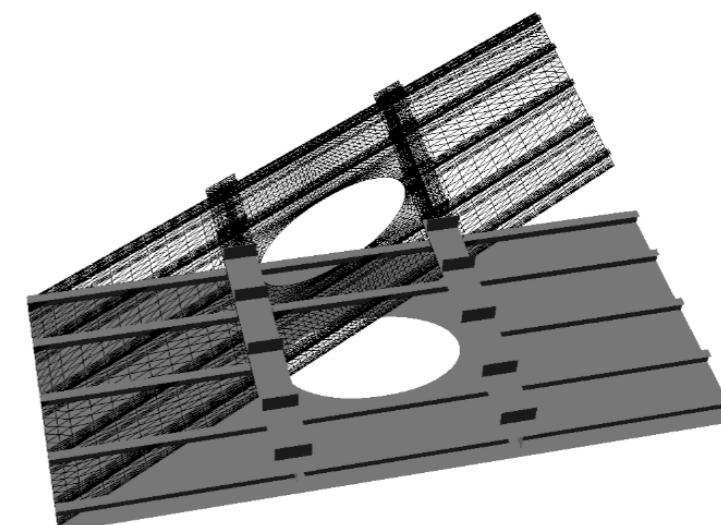


- Finite element model with 56,916 degrees of freedom
- 13 design variables (5 shape, 8 material)

- Problem Statement:  $n_{mat}=11$ ,  $n_{RHS}=14$ 
  - Given: 10 randomly-chosen (not necessarily nearby) previously-queried designs and designs for  $i=11$ :



Design A



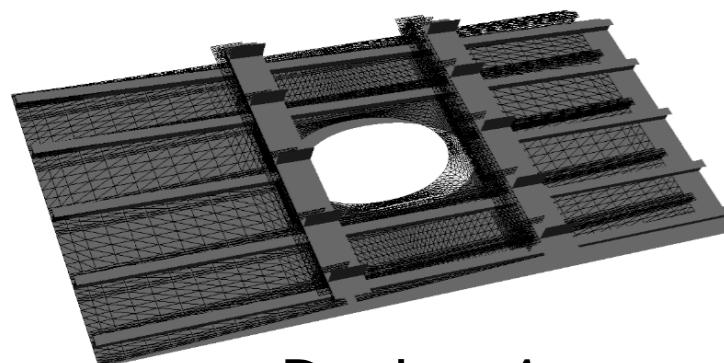
Design B

- Compute:  $\tilde{x}^{(11,j)}$ ,  $j = 1, \dots, n_{RHS}$  satisfying

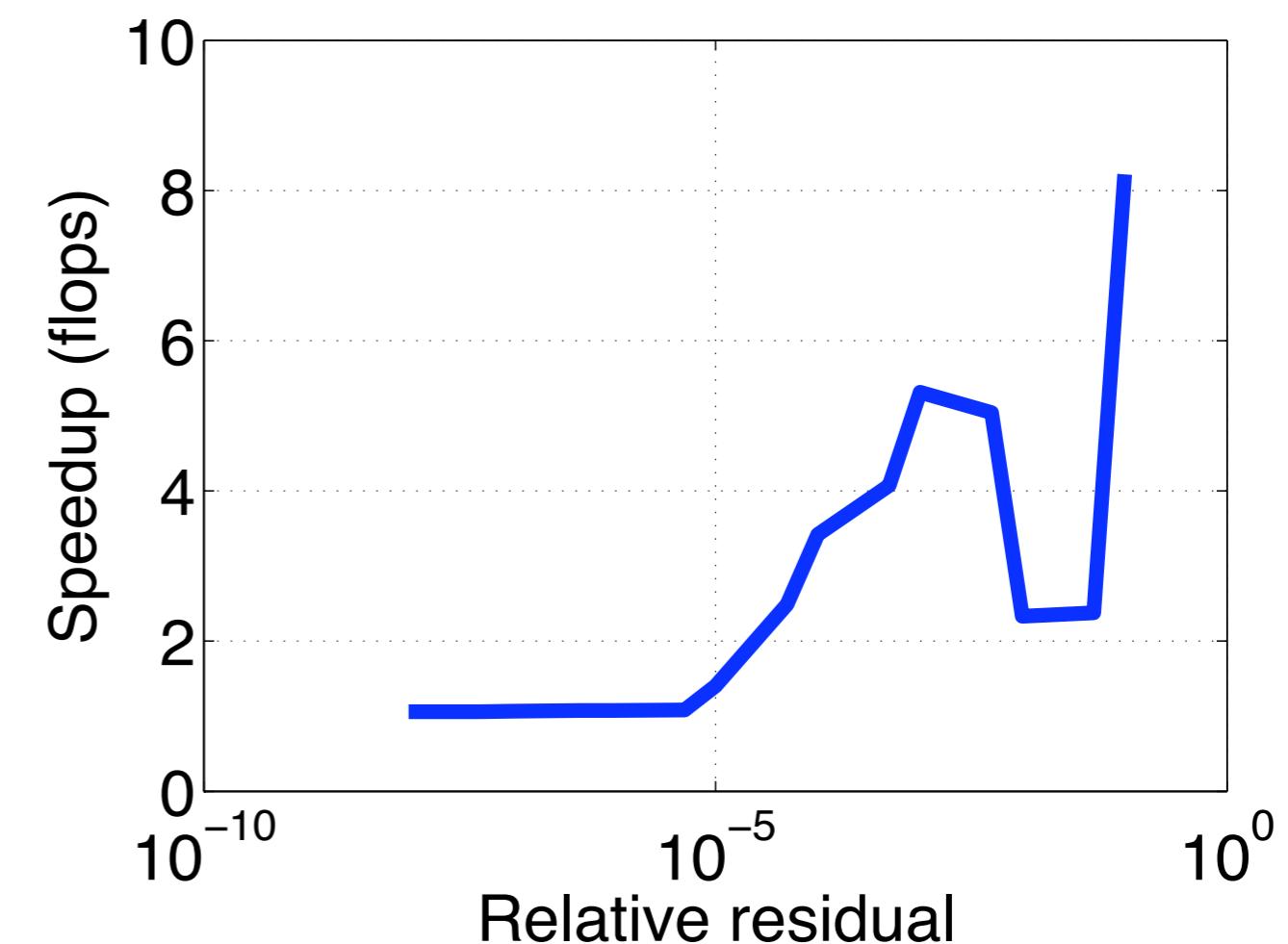
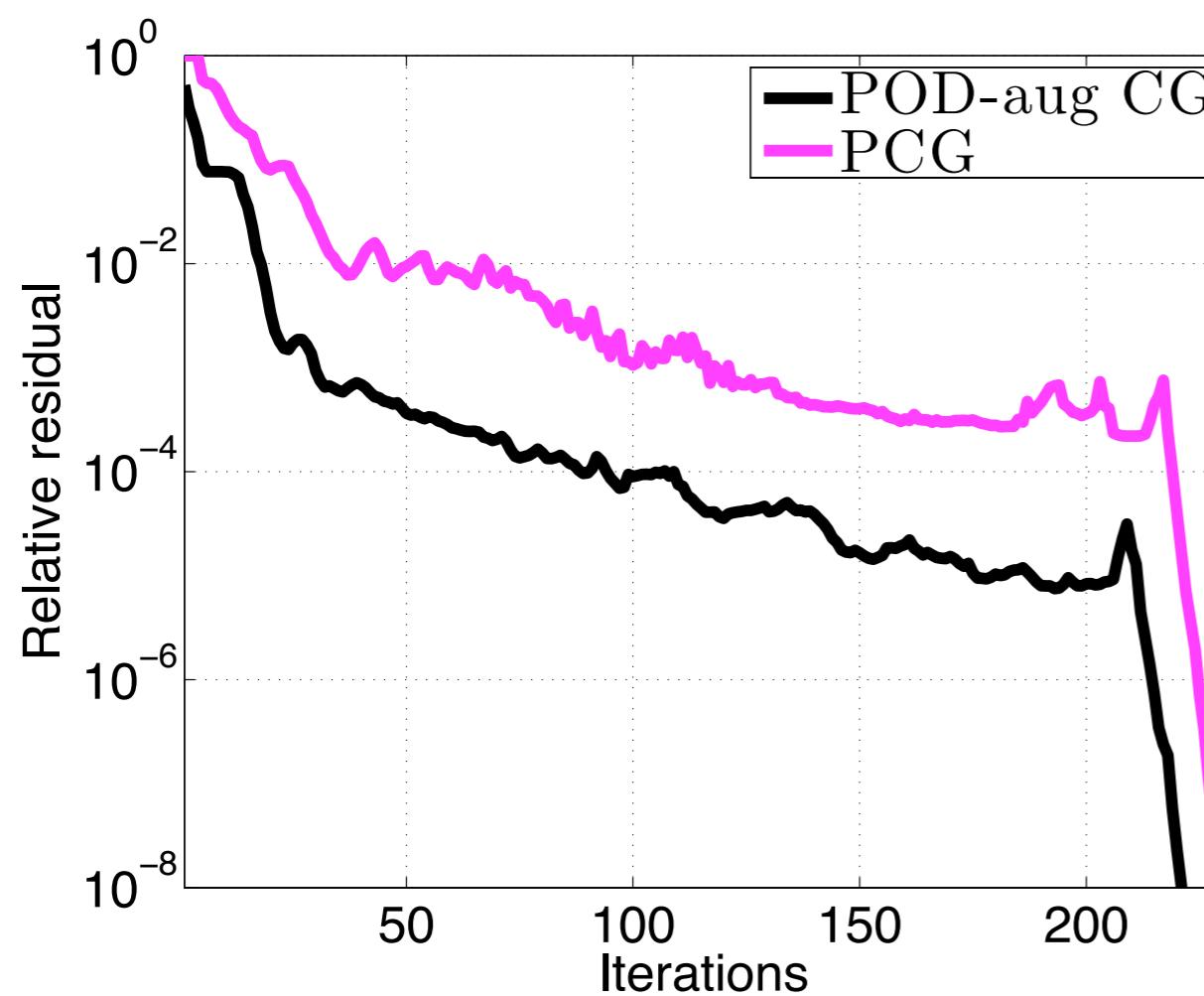
$$\frac{\|b^{(11,j)} - A^{(11)}\tilde{x}^{(11,j)}\|_2}{\|b^{(11,j)}\|_2} \leq \epsilon$$

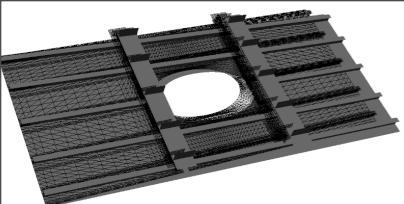


# Results: Design B, $n_{RHS} = 1$



Design A

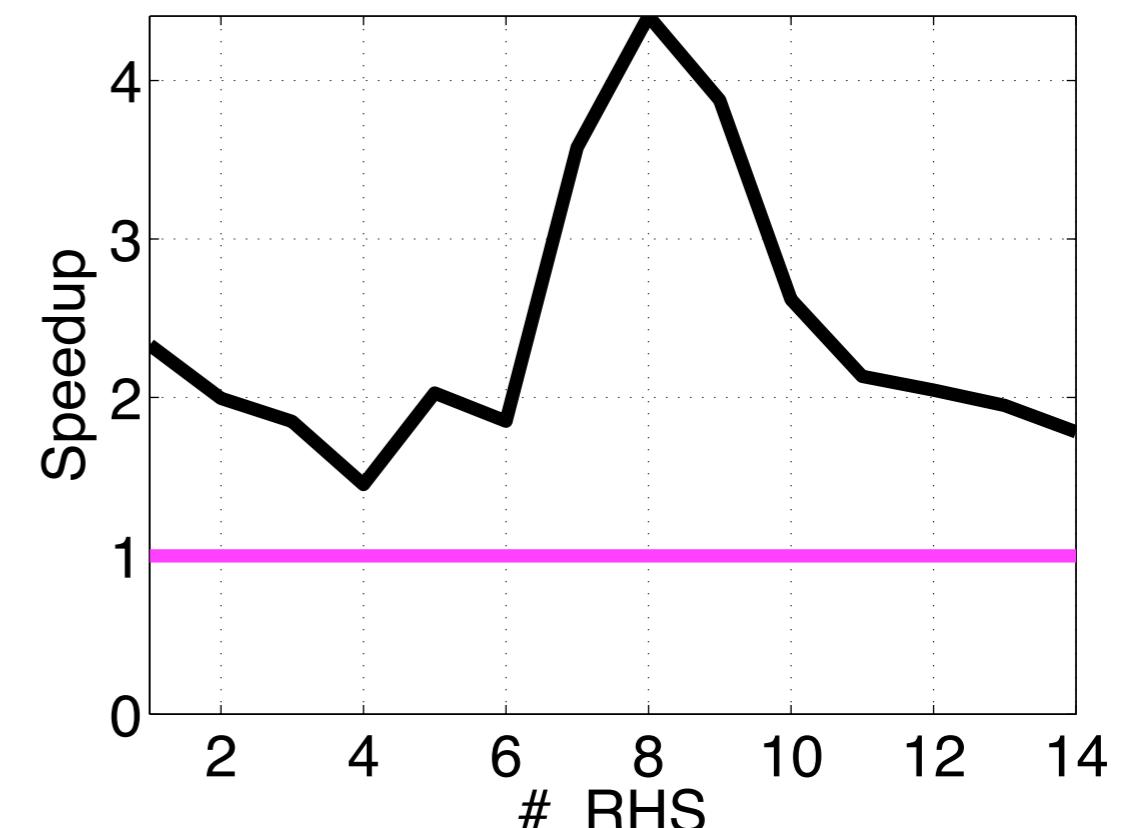
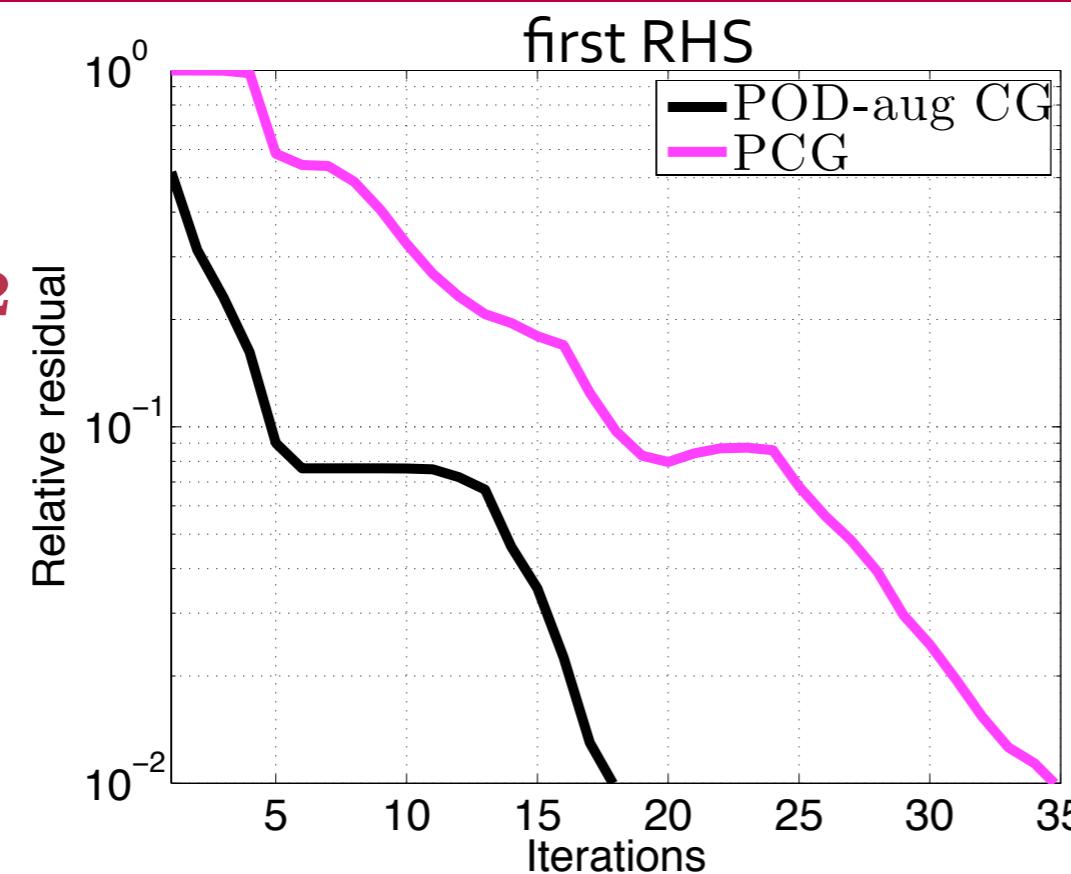




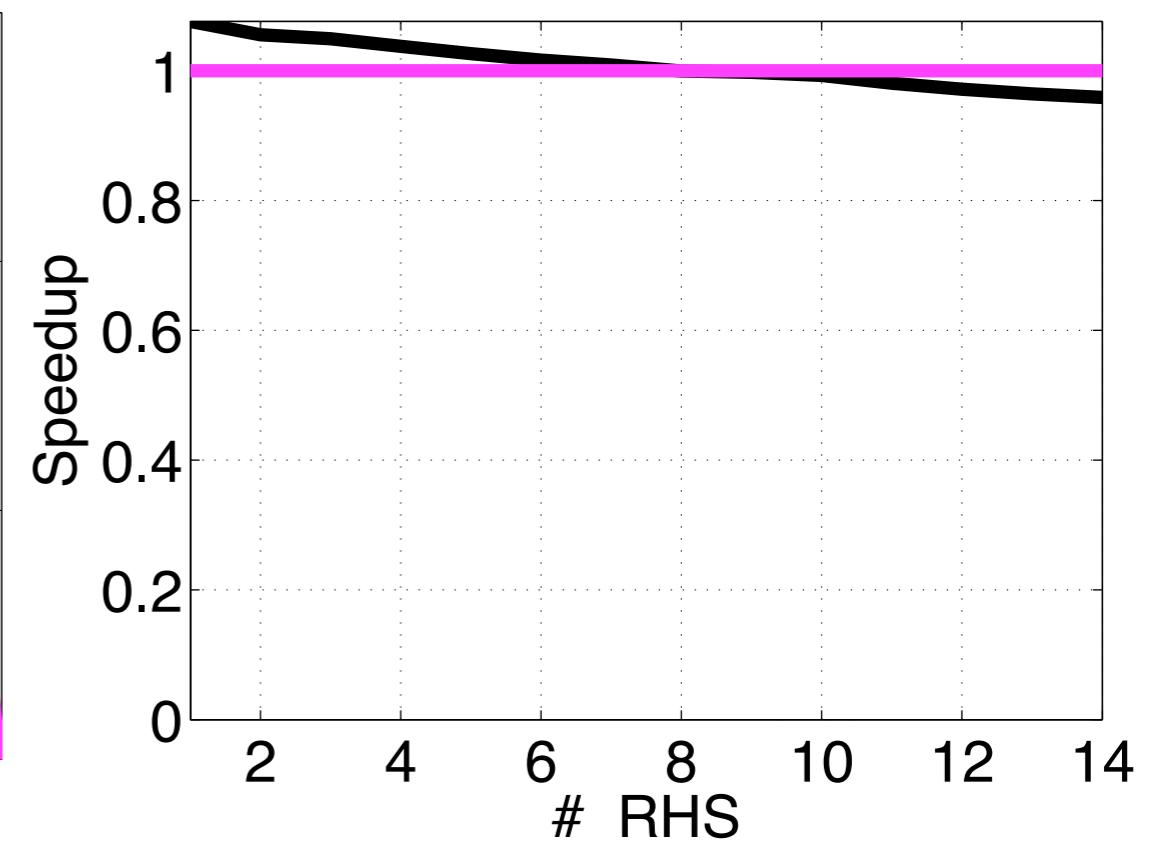
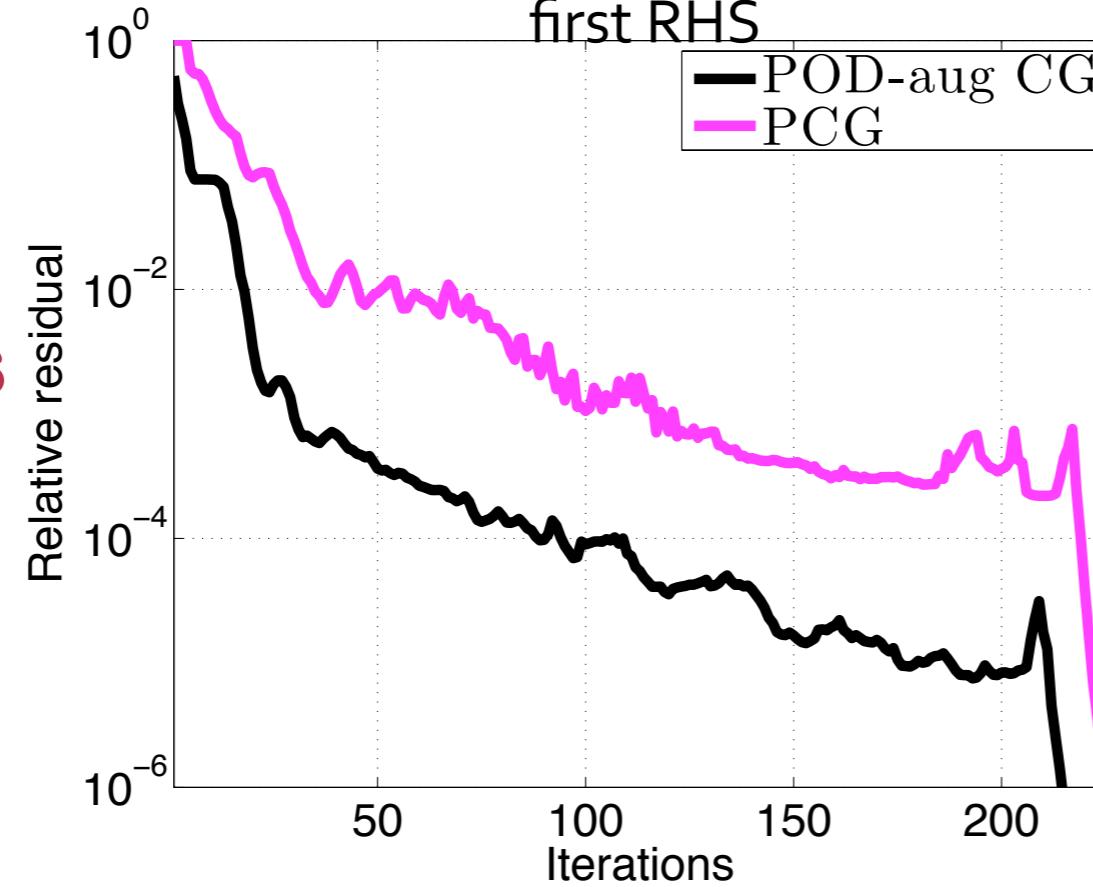
# Results: Design A, $n_{RHS} = 14$

Design A

$$\epsilon = 10^{-2}$$

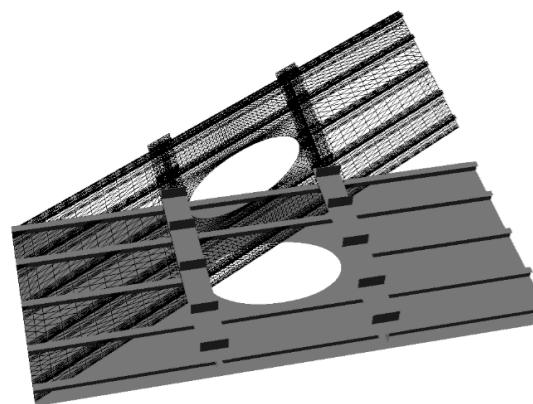


$$\epsilon = 10^{-6}$$

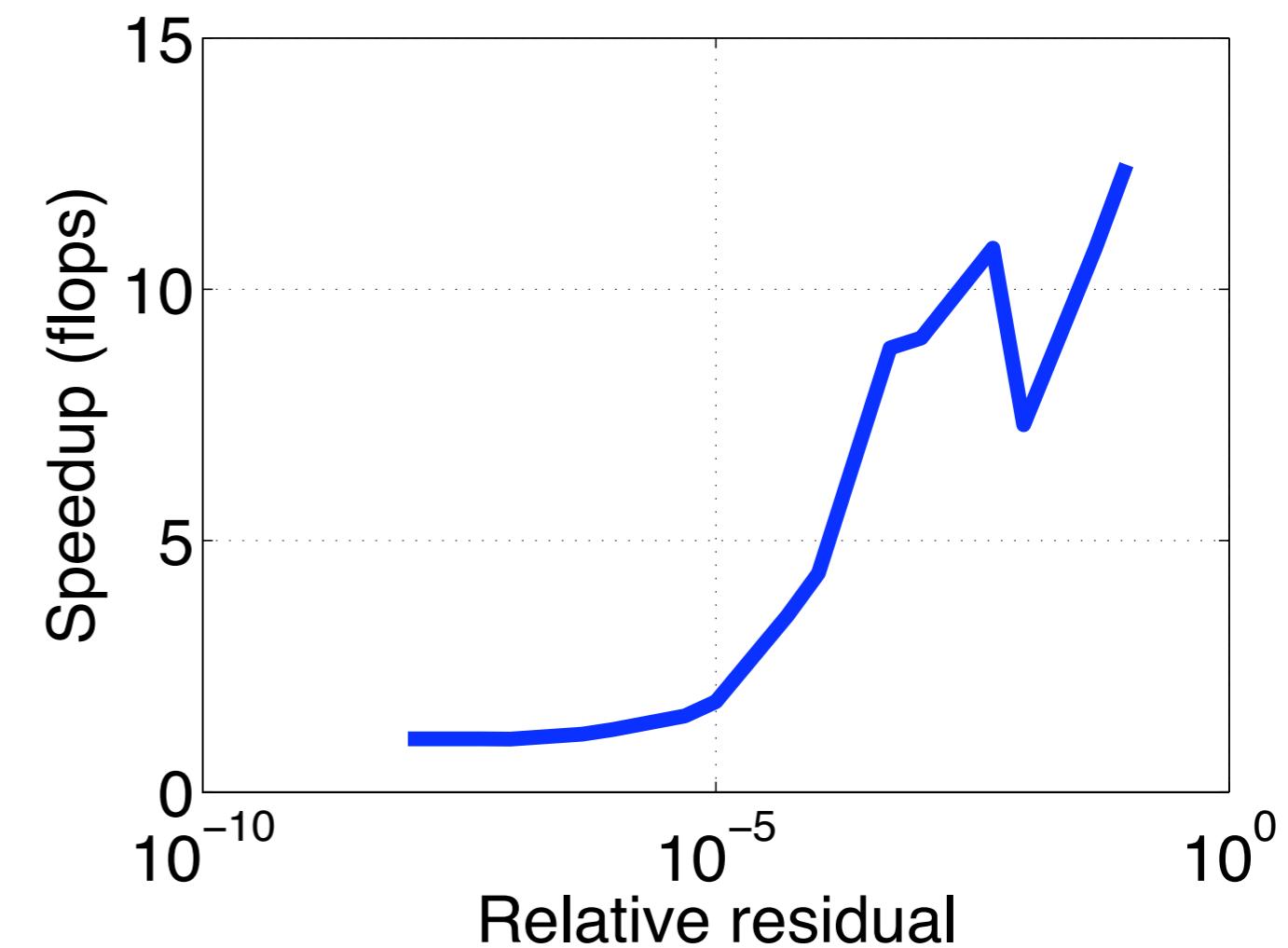
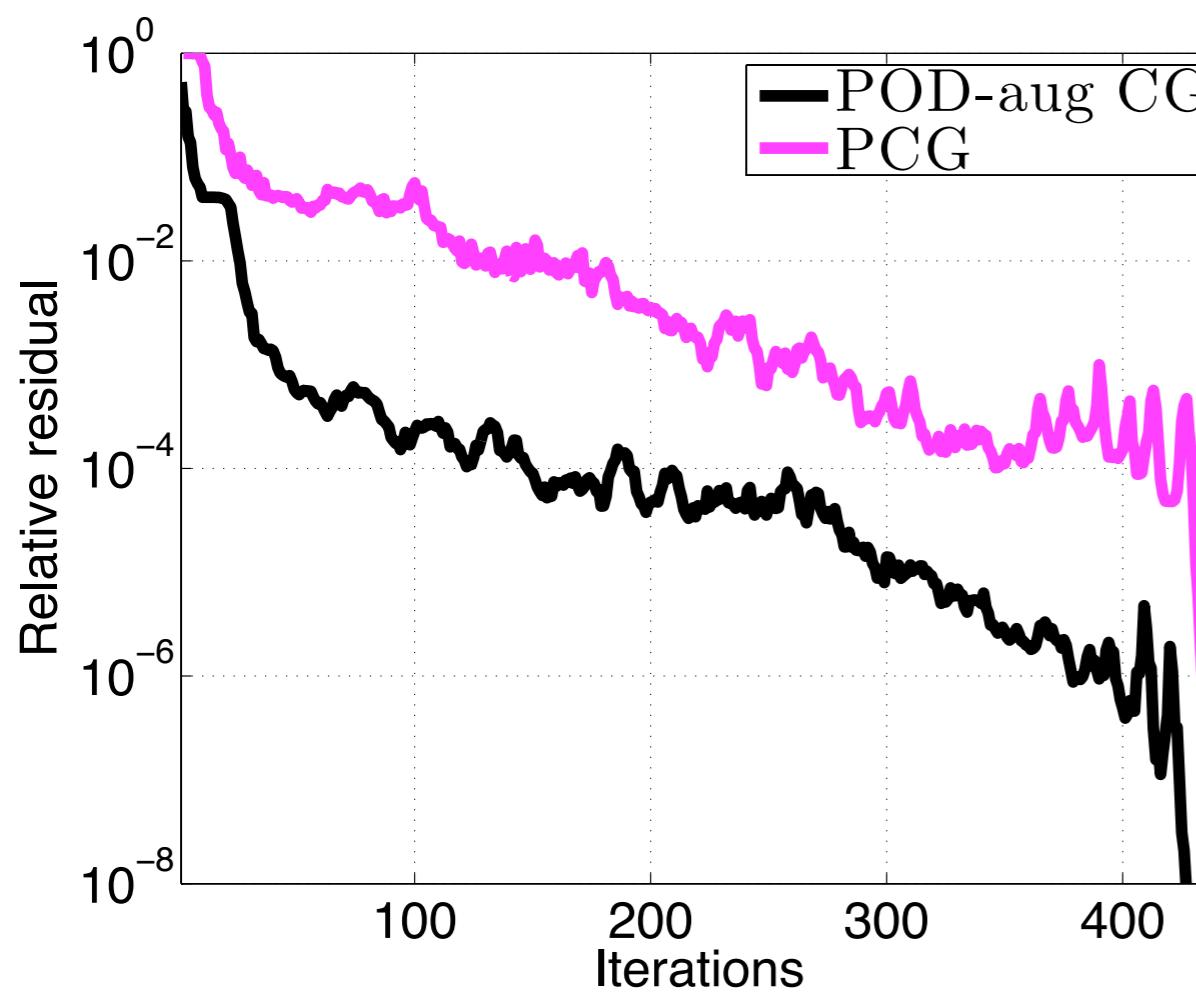


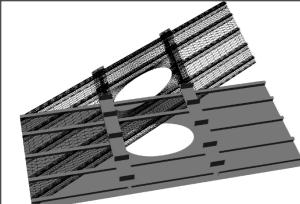


# Results: Design B, $n_{RHS} = 1$



Design B

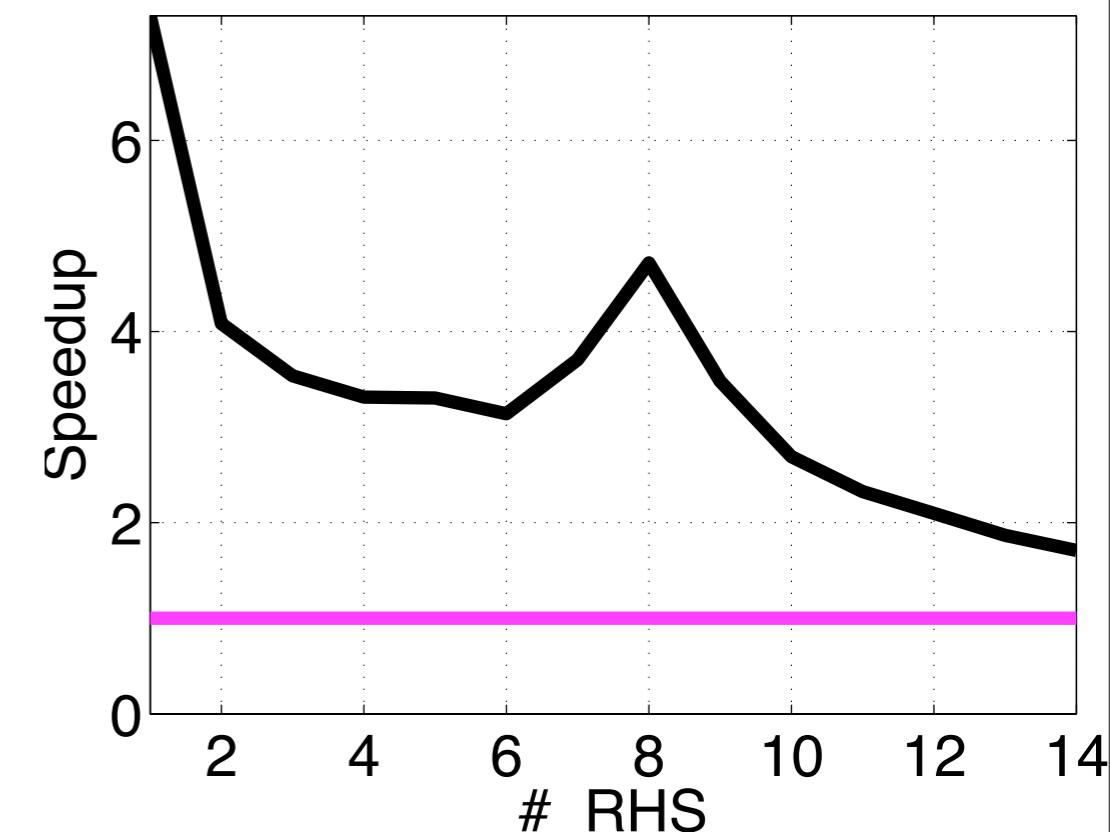
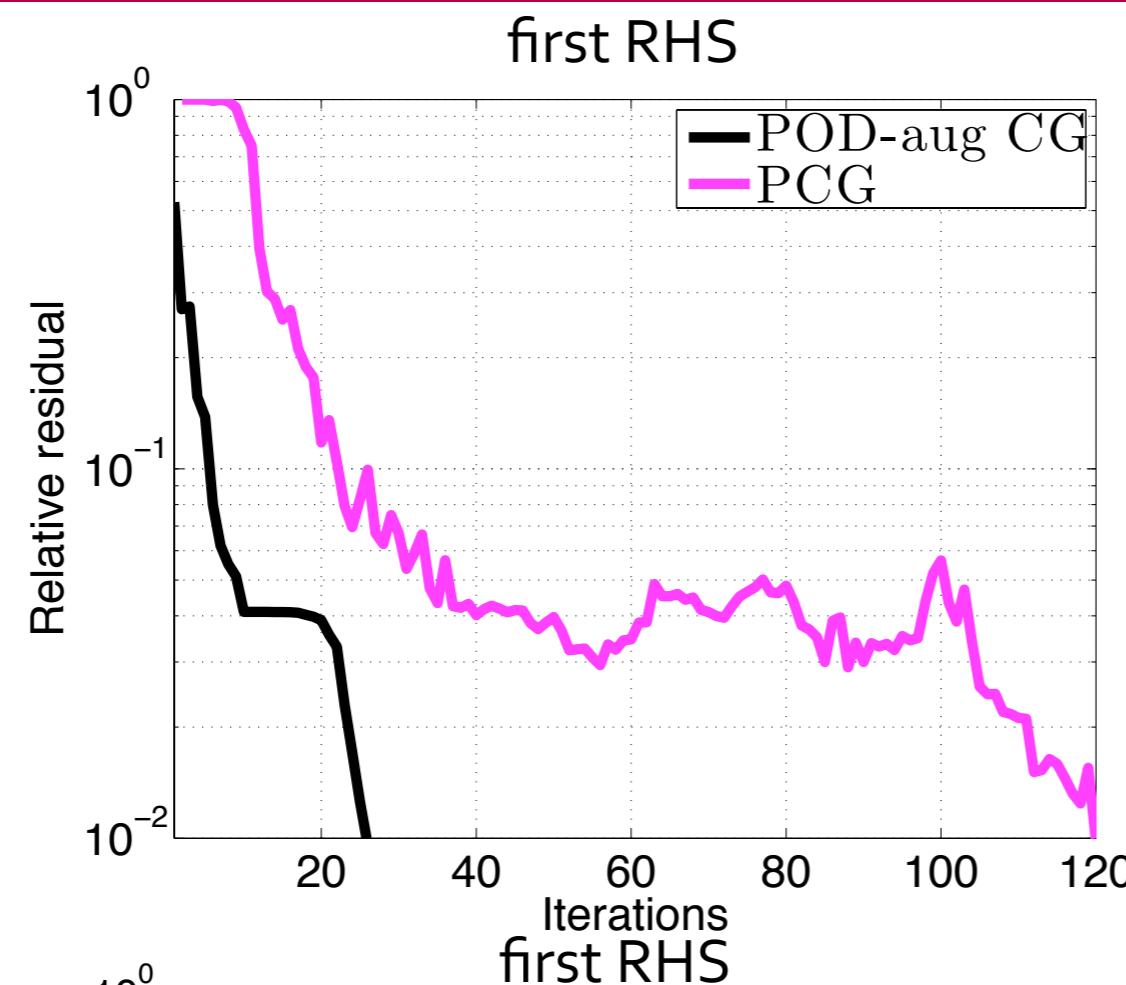




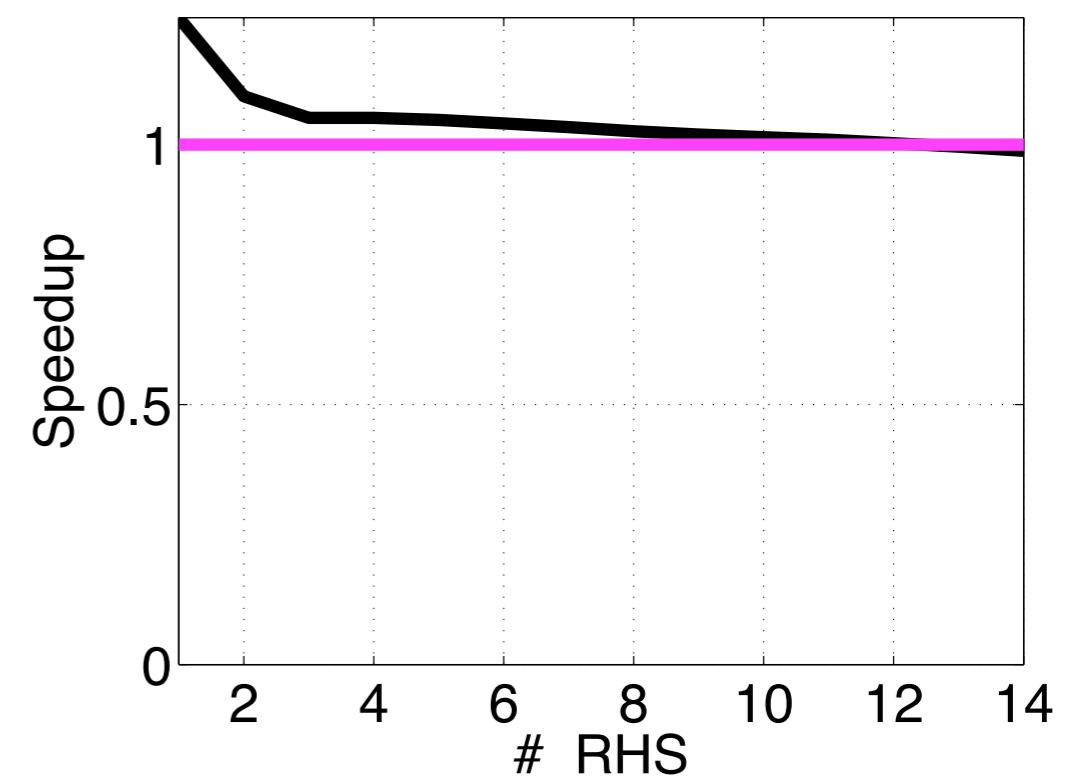
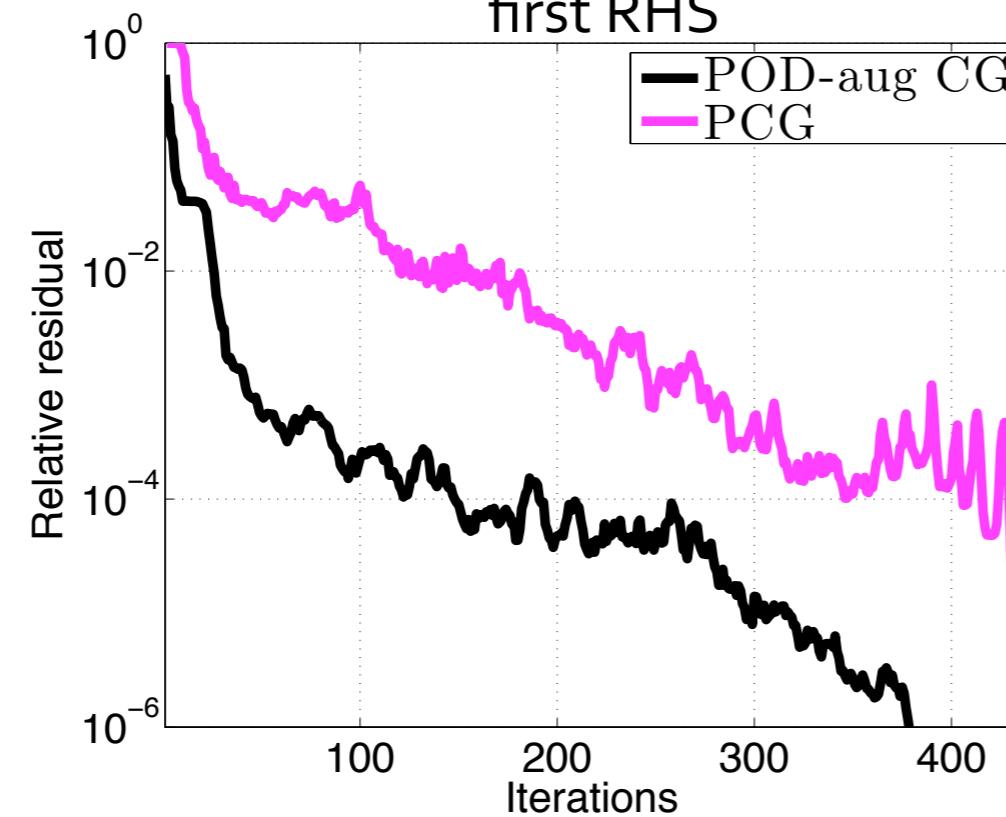
# Results: Design B, $n_{RHS} = 14$

Design B

$$\epsilon = 10^{-2}$$



$$\epsilon = 10^{-6}$$





# Conclusions: POD-augmented CG



- A POD-based augmented conjugate gradient method
  - Accelerates convergence to inexact tolerances
  - Efficiency due to choice of POD snapshots, weights, and norm
  - Highest speedups for **modest tolerances, few right-hand sides**
- Future work
  - Combine with other augmented CG approaches
    - Deflation: include approximated eigenvectors in stage 1
  - Extend to systems with non-SPD matrices



# Questions?



- Thanks to Charbel Bou-Mosleh, David Amsallem, and Julien Cortial for their contributions to this research.
- References
  - K. Carlberg, C. Bou-Mosleh, and C. Farhat, “Efficient nonlinear model reduction via a least-squares Petrov-Galerkin projection and compressive tensor approximations,” International Journal of Numerical Methods in Engineering, to appear, 2010.
  - K. Carlberg and C. Farhat, “An Adaptive POD-Krylov Reduced-Order Model for Structural Optimization,” Proceedings of 8th World Congress on Structural and Multidisciplinary Optimization, Lisbon, Portugal, June 1–5, 2009.