

# A Semi-incremental Scheme for Fatigue Damage Computations

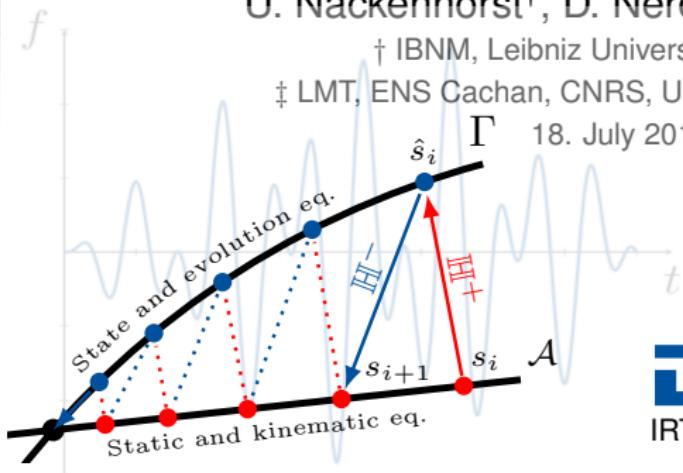
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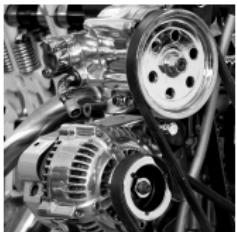


**DFG**  
IRTG-1627

Deutsche  
Forschungsgemeinschaft

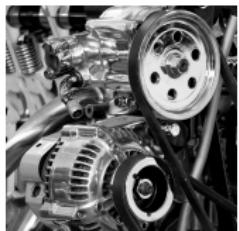
# Fatigue damage

## Fluctuating loads



# Fatigue damage

Fluctuating loads



Material degradation  
lower load carrying capacity

Large number of load cycles  
thousands or millions

Continuum damage model  
 $\dot{D} = dD/dt$

Computational expense  
memory and time

# Fatigue damage

## Fluctuating loads



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lower load carrying capacity

Large number of load cycles  
thousands or millions

Continuum damage model  
 $\dot{D} = dD/dt$



## Computational expense memory and time

## Model order reduction (MOR) techniques

- 1 State of the art**
- 2 Reduced order model for cyclic loading**
- 3 Challenges and workarounds**
- 4 Numerical examples**
- 5 Conclusions and current research**

# Outline

- 1 State of the art**
- 2 Reduced order model for cyclic loading**
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## Short-term plastic/damage computations

Proper Orthogonal Decomposition (POD)

[Kerfriden, 2012; Ryckelynck, 2011].

Proper Generalised Decomposition (PGD)

[Vitse, 2016; Bhattacharyya, 2017].

## Long-term fatigue computations

Temporal homogenisation

[Fish and Yu, 2002; Devulder, 2010]

(Modified) jump cycle approach

[Desmorat, 2005; Bhattacharyya, 2018]

Space-time finite element method

[Bhamare, 2014; Fritzen, 2018]

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PGD in a modified LATIN framework

## 1 State of the art

## 2 Reduced order model for cyclic loading

## 3 Challenges and workarounds

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# Mechanical problem

Infinitesimal strain and isothermal quasi-static settings

## Admissibility equations (**AD.**)

Static admissibility

$$\nabla \cdot \boldsymbol{\sigma} + \boldsymbol{b} = \mathbf{0} \quad \text{in } \Omega \times [0, T]$$

$$\boldsymbol{\sigma} \cdot \boldsymbol{n} = \bar{t} \quad \text{on } \partial\Omega_N \times \mathcal{I}$$

Kinematic admissibility

$$\boldsymbol{\varepsilon} = \nabla^s \boldsymbol{u} \quad \text{in } \Omega \times [0, T]$$

$$\boldsymbol{u} = \bar{\boldsymbol{u}} \quad \text{on } \partial\Omega_D \times \mathcal{I}$$

## Nonlinear constitutive model (**CM.**)

State equations

$$\boldsymbol{\sigma} = f(\psi(\boldsymbol{\varepsilon}, \boldsymbol{\varepsilon}^P, \boldsymbol{q}))$$

$$\boldsymbol{Q} = g(\psi(\boldsymbol{\varepsilon}, \boldsymbol{\varepsilon}^P, \boldsymbol{q}))$$

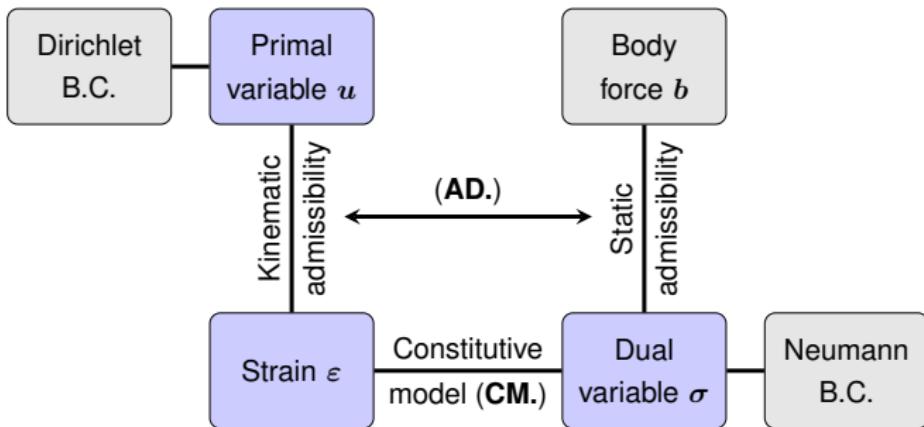
Evolution equations

$$\dot{\boldsymbol{\varepsilon}}^P = \hat{f}(\phi(\boldsymbol{\sigma}, \boldsymbol{Q}))$$

$$\dot{\boldsymbol{q}} = \hat{g}(\phi(\boldsymbol{\sigma}, \boldsymbol{Q}))$$

An **incremental** scheme:

Solve **iteratively**

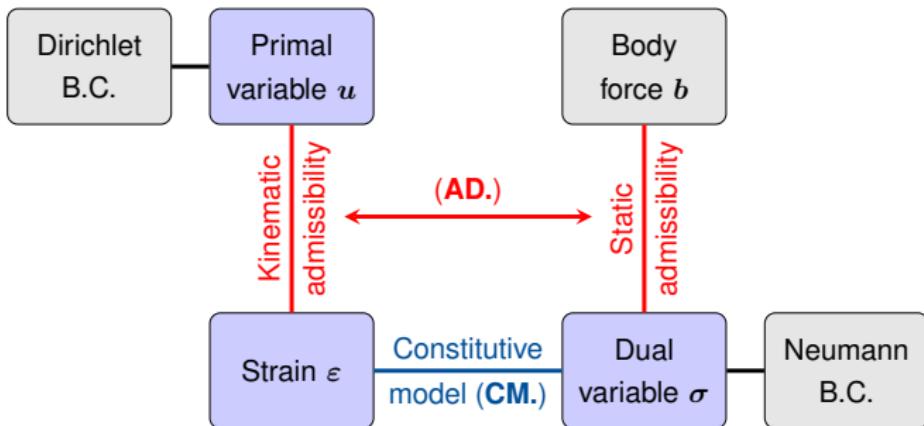


for each time step **consecutively**.

# LATIN vs. incremental schemes

A **LATIN** scheme:

Solve iteratively **(AD.)** and **(CM.)**



for all time steps **simultaneously**.

# What does simultaneously mean?

école  
normale  
supérieure  
paris-saclay

1 1 1  
1 0 2  
1 0 0 4

Leibniz  
Universität  
Hannover

# LATIN solution algorithm

Start with an elastic initialisation  $s_0$

Evaluate (CM.) to get  $s_{\text{local}}$  (local stage)

Solve (AD.) to get  $s_{\text{global}}$  (global stage)

Transfer data using affine relations (search direction eq.)

$$(\boldsymbol{\sigma}_{\text{global}} - \boldsymbol{\sigma}_{\text{local}}) - \mathbb{H}^\top : (\boldsymbol{\varepsilon}_{\text{global}} - \boldsymbol{\varepsilon}_{\text{local}}) = \mathbf{0}$$

Iterate until convergence with an energy error indicator

$$\xi = \frac{\|\boldsymbol{s}_{\text{global}} - \boldsymbol{s}_{\text{local}}\|}{\frac{1}{2} \|\boldsymbol{s}_{\text{global}} + \boldsymbol{s}_{\text{local}}\|}, \quad \|\boldsymbol{s}\|^2 = \frac{1}{2T} \int_{\Omega \times \mathcal{I}} (\boldsymbol{\sigma} : \mathbb{C}^{-1} : \boldsymbol{\sigma} + \boldsymbol{\varepsilon} : \mathbb{C} : \boldsymbol{\varepsilon}) \, d\Omega \, dt.$$

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# Static admissibility

Weak form at iteration  $i + 1$

$$\int_{\Omega \times \mathcal{I}} \nabla^s \mathbf{w}^* : \boldsymbol{\sigma}_{\text{global}}^{(i+1)} \, d\Omega \, dt = \int_{\Omega \times \mathcal{I}} \mathbf{w}^* \cdot \mathbf{b} \, d\Omega \, dt + \int_{\partial\Omega_N \times \mathcal{I}} \mathbf{w}^* \cdot \bar{\mathbf{t}} \, dS \, dt$$

Correction  $\Delta \bullet_{\text{global}}^{(i+1)} = \bullet^{(i+1)} - \bullet^{(i)}$

$$\int_{\Omega \times \mathcal{I}} \nabla^s \mathbf{w}^* : \Delta \boldsymbol{\sigma}_{\text{global}}^{(i+1)} \, d\Omega \, dt = 0$$

Static admissibility and global search direction

$$\int_{\Omega \times \mathcal{I}} \nabla^s \mathbf{w}^* : \mathbb{H}^- : \Delta \boldsymbol{\varepsilon}_{\text{global}}^{(i+1)} \, d\Omega \, dt = - \int_{\Omega \times \mathcal{I}} \nabla^s \mathbf{w}^* : \underbrace{\hat{\mathbf{f}}_{\text{local}}}_{\text{known}} \, d\Omega \, dt$$

# Proper Generalised Decomposition

Low-rank approximation

$$\boldsymbol{u}(\boldsymbol{x}, t) = \sum_{j=1}^N \boldsymbol{v}_j(\boldsymbol{x}) \circ \boldsymbol{\lambda}_j(t)$$

Enrichment to  $(\mu)$  previously generated modes

$$\Delta \boldsymbol{u}_{i+1}(\boldsymbol{x}, t) = \boldsymbol{v}_{\mu+1}(\boldsymbol{x}) \circ \boldsymbol{\lambda}_{\mu+1}(t)$$

POD-like update of  $(\mu)$  previously generated modes

$$\Delta \boldsymbol{u}_{i+1}(\boldsymbol{x}, t) = \sum_{j=1}^{\mu} \underbrace{\boldsymbol{v}_j(\boldsymbol{x})}_{\text{known}} \circ \Delta \boldsymbol{\lambda}_j(t)$$

# Alternated directions algorithm

Galerkin finite element discretisation

Spatial problem, with homogeneous boundary conditions:

$$\int_{\Omega} \nabla^s \mathbf{v}^* : \left( \int_{\mathcal{I}} \lambda \mathbb{H}^- \lambda dt \right) : \nabla^s \mathbf{v} d\Omega = - \int_{\Omega} \nabla^s \mathbf{v}^* : \left( \int_{\mathcal{I}} \lambda \hat{\mathbf{f}} dt \right) d\Omega$$

Temporal problem, with zero initial conditions:

$$\int_{\mathcal{I}} \lambda^* \left( \int_{\Omega} \nabla^s \mathbf{v} : \mathbb{H}^- : \nabla^s \mathbf{v} d\Omega \right) \lambda dt = - \int_{\mathcal{I}} \lambda^* \left( \int_{\Omega} \nabla^s \mathbf{v} : \hat{\mathbf{f}} d\Omega \right) dt$$

# What is actually solved?

Choosing the search directions  $\mathbb{H}^-$  as

$$\mathbb{H}^- = \alpha \mathbb{C} \quad \alpha \in ]0, 1]$$

Enrichment step

$$\begin{aligned} \gamma \underline{\underline{K}} \underline{v} &= \underline{f} & \gamma \in \mathbb{R} & \underline{\underline{K}} \in \mathbb{R}^{n \times n} & \underline{v}, \underline{f} \in \mathbb{R}^n & \text{with B.C.} \\ a \underline{\lambda} &= \underline{b} & a \in \mathbb{R} & \underline{\lambda}, \underline{b} \in \mathbb{R}^{n_t} & & \text{with I.C.} \end{aligned}$$

Temporal update step

$$\tilde{\underline{\underline{A}}} \tilde{\underline{\lambda}}^\top = \tilde{\underline{\underline{B}}} \quad \tilde{\underline{\underline{A}}} \in \mathbb{R}^{\mu \times \mu} \quad \tilde{\underline{\underline{B}}} \in \mathbb{R}^{\mu \times n_t} \quad \tilde{\underline{\lambda}} = [\Delta \underline{\lambda}_1, \dots, \Delta \underline{\lambda}_\mu]$$

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# Challenges in a PGD framework

The cost of integration over all generalised coordinates

$$\int_{\mathcal{I}} \int_{\Omega} \bullet \, d\Omega \, dt \quad \text{:(}$$

The fast increase in the number of modes

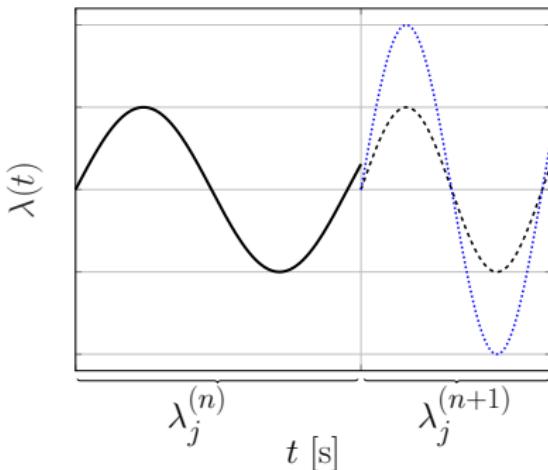
$$\boldsymbol{u}(\boldsymbol{x}, t) = \sum_{j=1}^{\mu} \boldsymbol{v}_j(\boldsymbol{x}) \circ \boldsymbol{\lambda}_j(t) \text{ with large } \mu \quad \text{:(}$$

Efficient variable amplitude and frequency simulations

# Semi-incremental extension

The temporal domain is divided into cycles

Cycles are simulated consecutively



Already generated  $\{\lambda_j(t)\}_{j=1}^\mu$  are scaled to  $\tilde{t} \in [0, 1]$

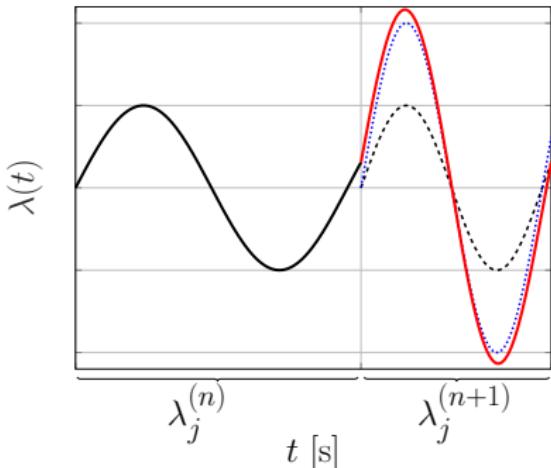
Scaled back to the temporal coordinate of the current cycle

# Semi-incremental extension

Continuity of the temporal modes is an issue

Temporal modes are vertically scaled and shifted via

$$\tilde{\lambda}_j^{(n+1)}(\tilde{t}) = m \tilde{\lambda}_j^{(n)}(\tilde{t}) + g \tilde{t} + h \quad \text{with I.C. and B.C.}$$

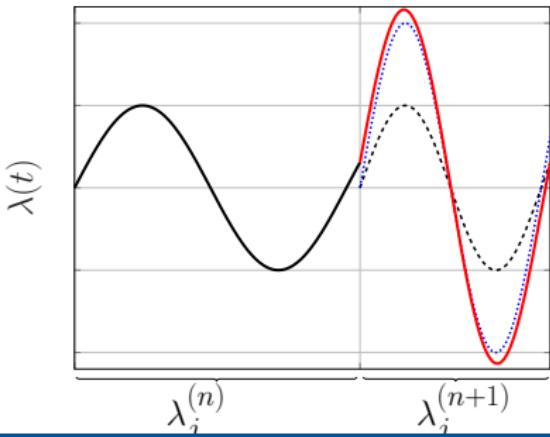


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Variable loading / integration over confined domains 😊

# SVD Compression of PGD

Large number of modes

Slow temporal update step

$$\tilde{\underline{A}} \tilde{\underline{\Lambda}}^T = \tilde{\underline{B}} \quad \tilde{\underline{A}} \in \mathbb{R}^{\mu \times \mu} \quad \tilde{\underline{B}} \in \mathbb{R}^{\mu \times n_t} \quad \tilde{\underline{\Lambda}} = [\Delta \underline{\lambda}_1, \dots, \Delta \underline{\lambda}_\mu]$$

PGD enrichment with a Gram-Schmidt scheme is not optimal

SVD is optimal but expensive to compute

# SVD Compression of PGD

Given a solution  $\underline{\underline{U}} = \underline{\underline{V}} \underline{\underline{\Lambda}}^T \in \mathbb{R}^{n \times n_t}$  with

$$\underline{\underline{V}} = [\underline{v}_1, \dots, \underline{v}_\mu] \in \mathbb{R}^{n \times \mu} \quad \underline{\underline{\Lambda}} = [\underline{\lambda}_1, \dots, \underline{\lambda}_\mu] \in \mathbb{R}^{n_t \times \mu}$$

Exploit the outer product format of PGD

$$\underline{\underline{V}} = \underline{\underline{Q}}_v \underline{\underline{R}}_v \quad \underline{\underline{\Lambda}} = \underline{\underline{Q}}_\lambda \underline{\underline{R}}_\lambda$$

Compute SVD of a small matrix

$$\underline{\underline{T}} = \underline{\underline{R}}_v \underline{\underline{R}}_\lambda^T \in \mathbb{R}^{\mu \times \mu} \quad \underline{\underline{T}} \approx \hat{\underline{\underline{V}}} \hat{\underline{\underline{S}}} \hat{\underline{\underline{\Lambda}}}^T$$

Solution with a compressed basis

$$\underline{\underline{U}} \approx \underline{\underline{Q}}_v \hat{\underline{\underline{V}}} \quad \hat{\underline{\underline{S}}} \quad \hat{\underline{\underline{\Lambda}}}^T \underline{\underline{Q}}_\lambda^T$$

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Solution with a compressed basis

$$U \sim \mathcal{O}(\hat{\underline{\underline{V}}} \hat{\underline{\underline{S}}} \hat{\underline{\underline{\Lambda}}}^T)$$

Non-demanding optimal decomposition 😊

# Is this efficient?

A low-rank approximation of  $\varepsilon_{\text{global}}$

What about  $\sigma_{\text{global}}$ , the local stage and the error indicator?

$$(\sigma_{\text{global}} - \sigma_{\text{local}}) - \mathbb{H}^\top : (\varepsilon_{\text{global}} - \varepsilon_{\text{local}}) = \mathbf{0}$$

What is missing is PGD of  $\varepsilon_{\text{local}}$

Randomised SVD (RSVD) algorithm

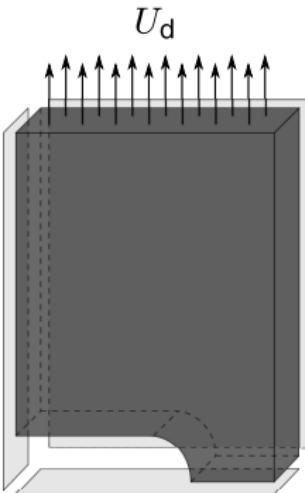
A dedicated arithmetic toolbox to handle operations on PGD

# Outline

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- 2 Reduced order model for cyclic loading**
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# Numerical results

A plate subjected to cyclic loading (Cr-Mo steel at 580°C)



# Viscoplastic viscodamage material

Stress-strain

$$\sigma = (1 - D) \mathbb{C} : \varepsilon^e$$

$$\dot{\varepsilon}^p = \dot{\lambda} N$$

Kinematic hardening

$$\beta = \frac{2}{3} c \alpha$$

$$\dot{\alpha} = \dot{\lambda} \left( \tilde{N} - \frac{3}{2} \frac{a}{c} \beta \right)$$

Isotropic hardening

$$R = R_\infty (1 - e^{-b r})$$

$$\dot{r} = \dot{\lambda}$$

Damage

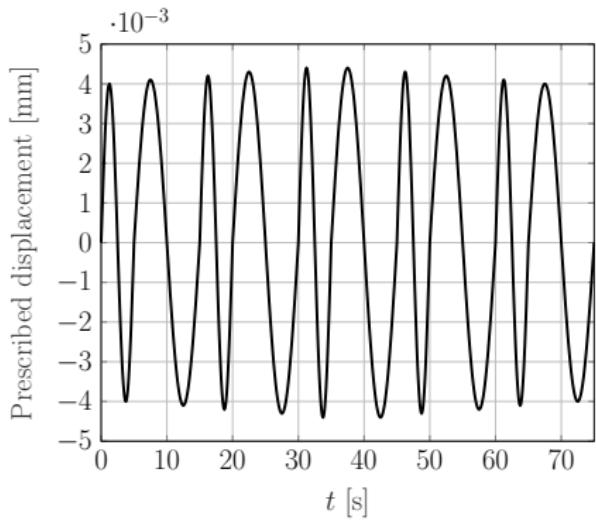
$$Y = \frac{1}{2} \varepsilon^e : \mathbb{C} : \varepsilon^e$$

$$\dot{D} = \frac{\dot{\lambda}}{1 - D} \left( \frac{Y}{S} \right)^s$$

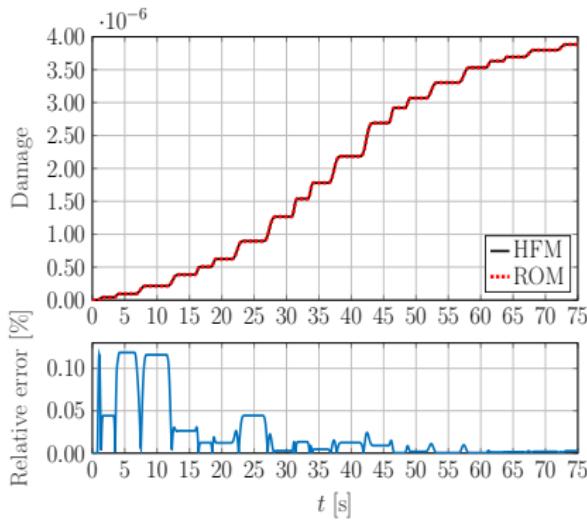
# Model verification

$1884 \cdot 41 \cdot 10$  DOF

With respect to a modified Newton-Raphson scheme



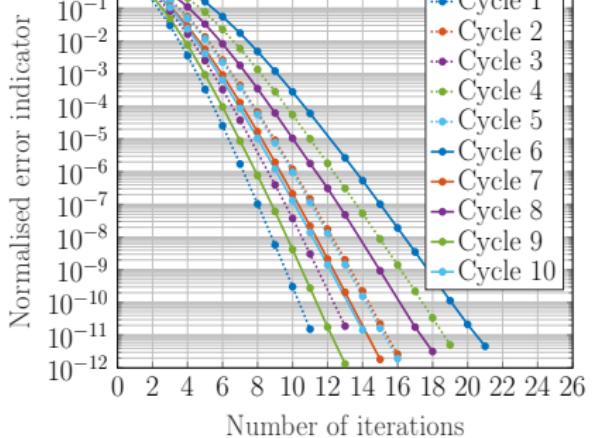
Prescribed displacement



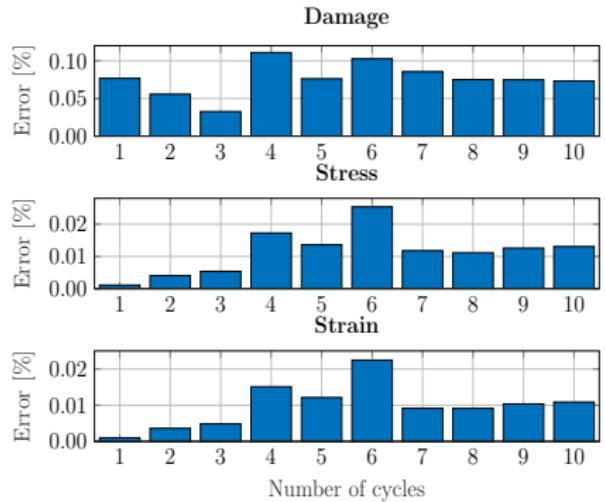
Damage evolution

# Model verification

$$\|e\|_{\Omega \times \mathcal{I}}^2 = \frac{1}{T |\Omega|} \int_{\mathcal{I}} \int_{\Omega} e : e \, d\Omega \, dt$$



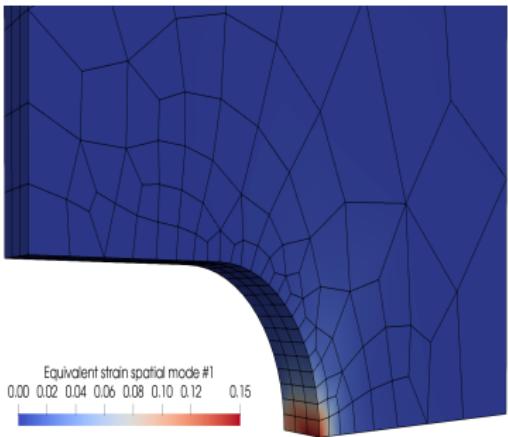
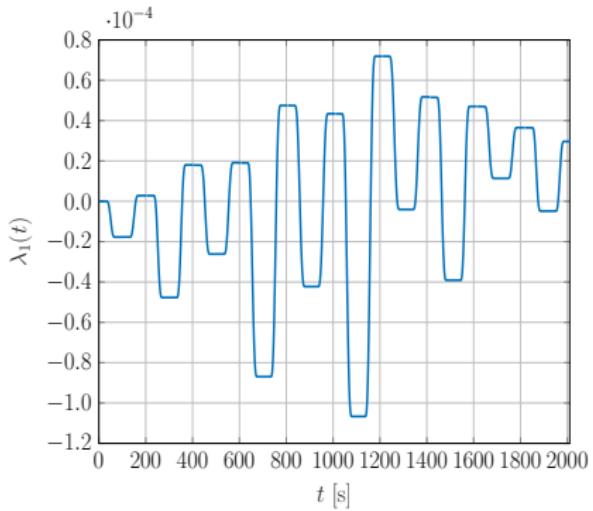
Error indicator



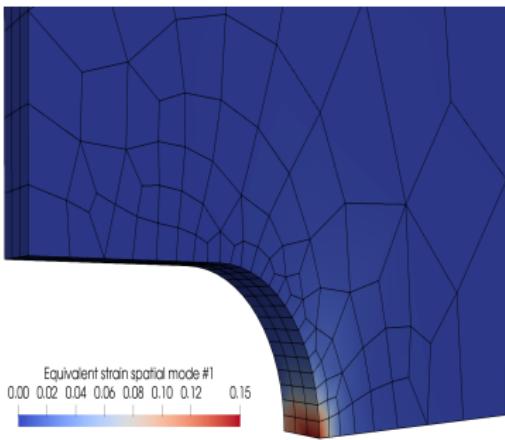
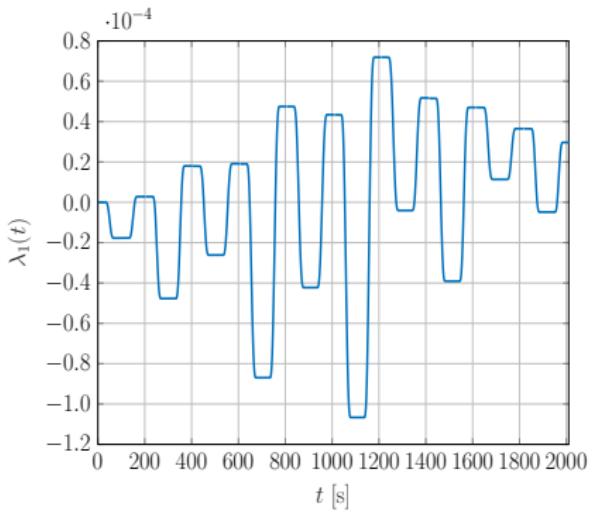
Space-time average relative error

# Model verification

The first temporal and spatial modes



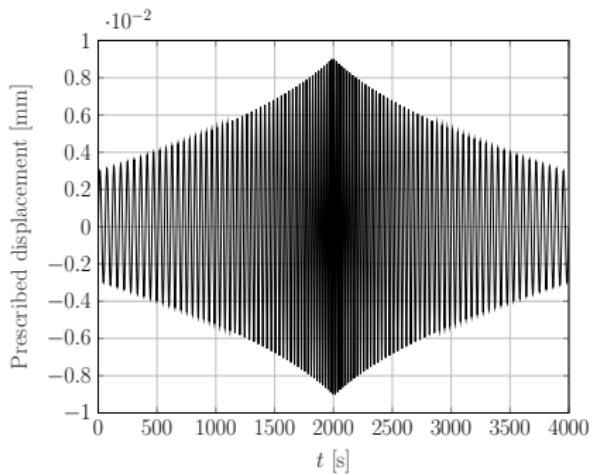
## The first temporal and spatial modes



12 modes / Speedup factor  $\sim 25$

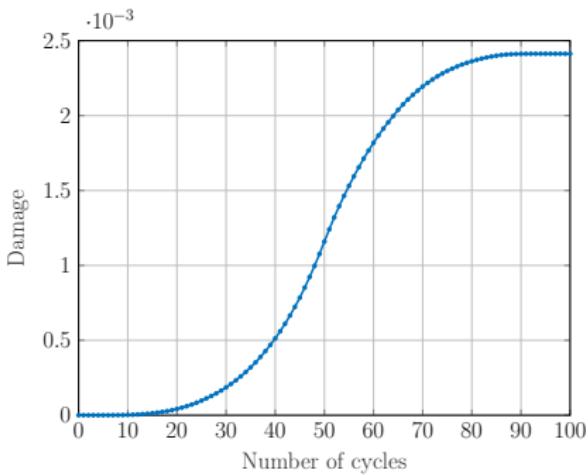
# Variable amplitudes & frequencies

Amplitudes:  $[30, 90] \cdot 10^{-4}$  mm



Prescribed displacement

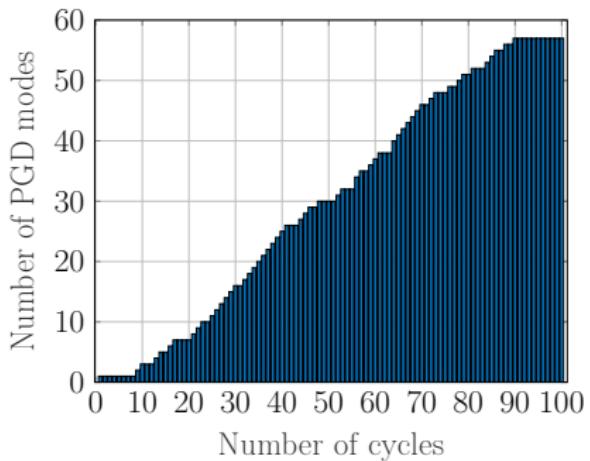
Time periods: [20, 60] sec



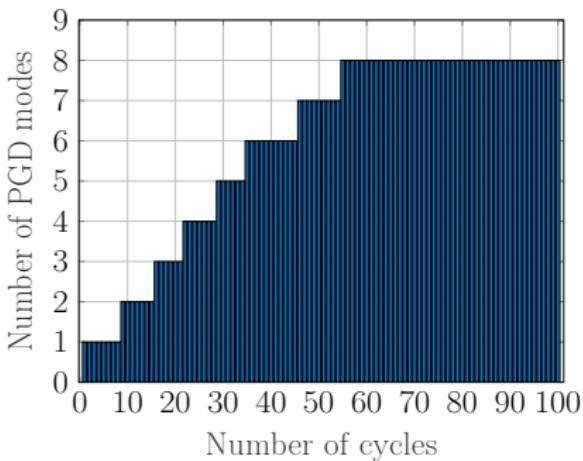
Damage evolution

# Variable amplitudes & frequencies

The growth of the ROB using an SVD scheme



Gram-Schmidt

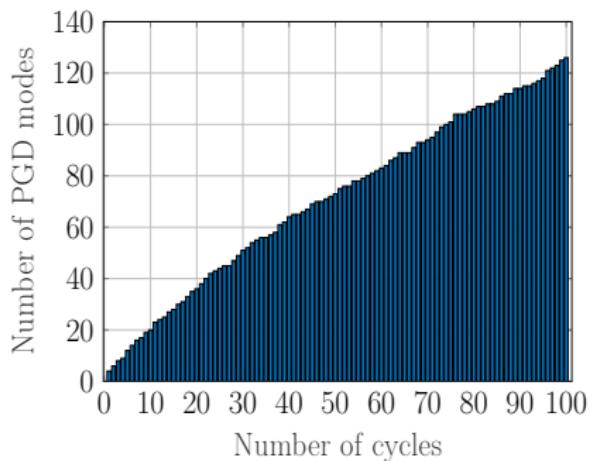


SVD scheme

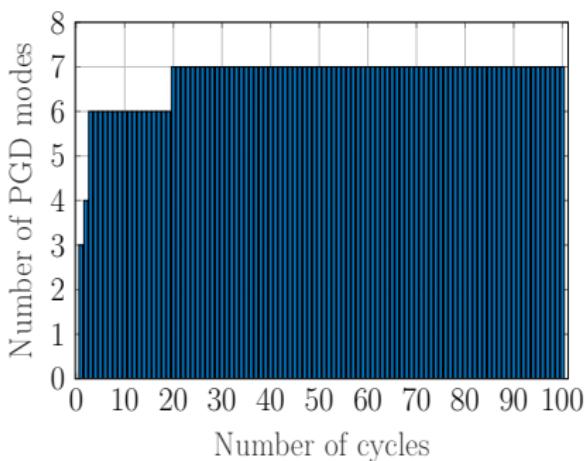
# Different ortho. schemes

$50,547 \cdot 33 \cdot 100$  DOF

In case of random loading and fine discretisation



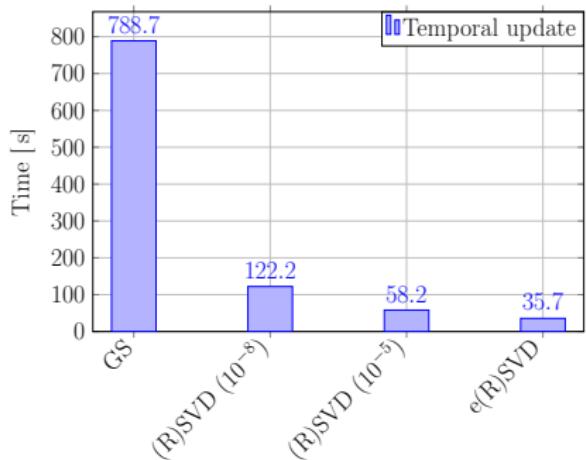
Gram-Schmidt



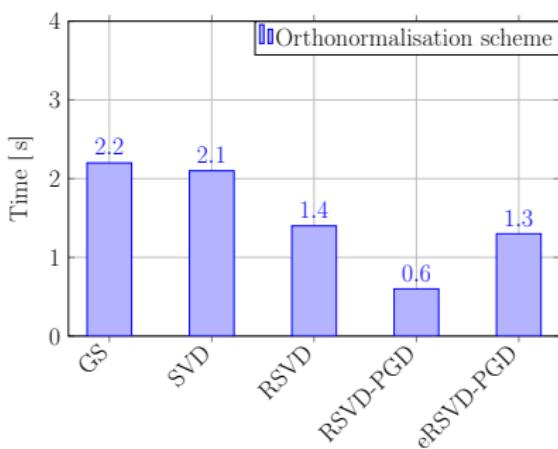
SVD scheme

# Different ortho. schemes

The required time to update and orthonormalise the modes



Temporal update

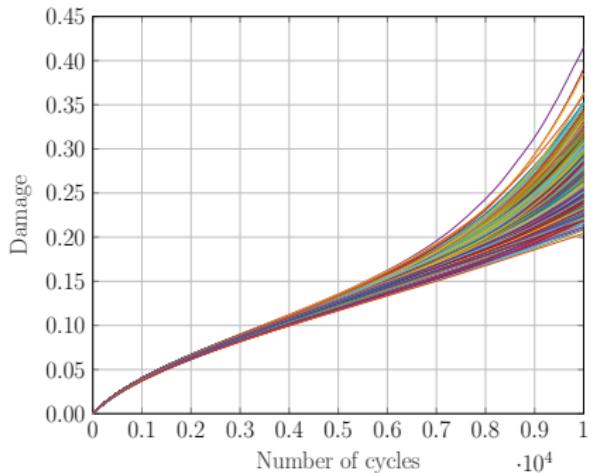


Orthonormalisation step

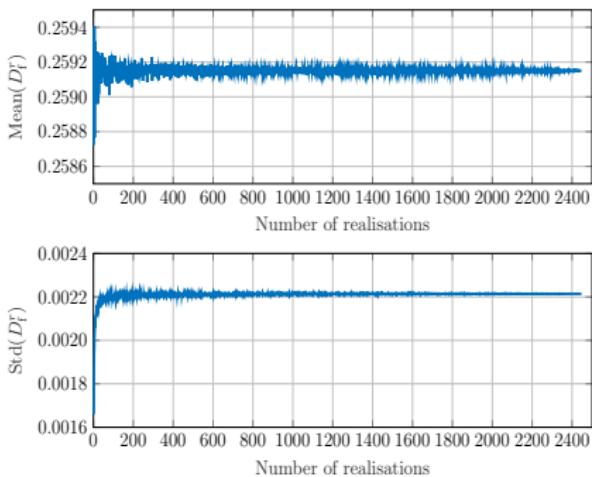
# Random amplitude loading

$1884 \cdot 41 \cdot 10^4$  DOF

$10^4$  cycles, uniform distribution in  $[53, 56] \cdot 10^{-2}$  mm



Different damage realisations

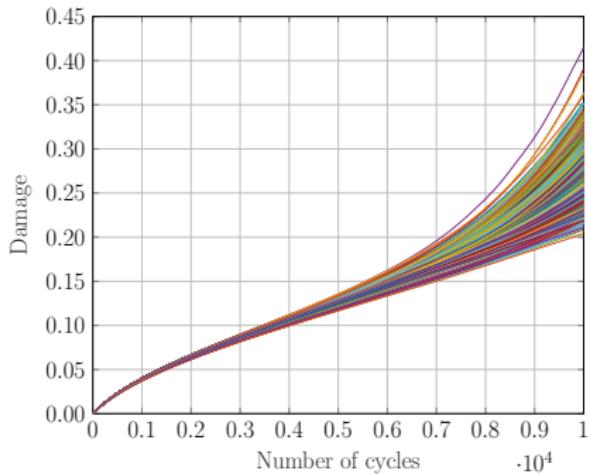


The mean and STD of  $D_f$

# Random amplitude loading

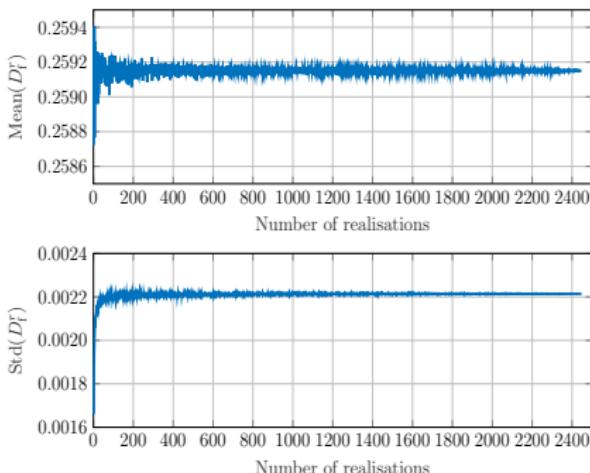
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$10^4$  cycles, uniform distribution in  $[53, 56] \cdot 10^{-2}$  mm



Critical damage value  $D_c = 0.3$

Probability of failure  $P_f = 5.4\%$



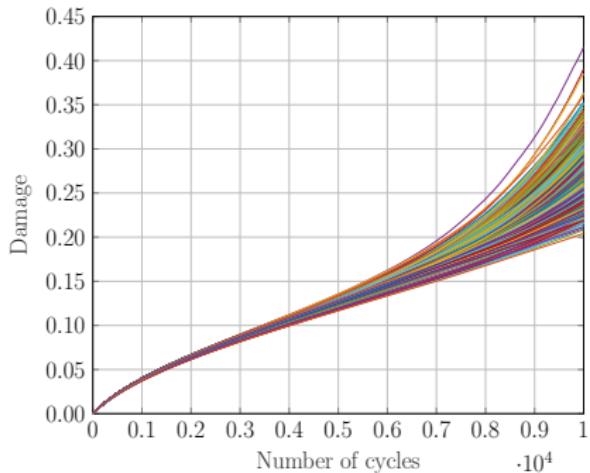
Requirements:

$[15 - 35]$  min and  $[1 - 1.5]$  GB

# Random amplitude loading

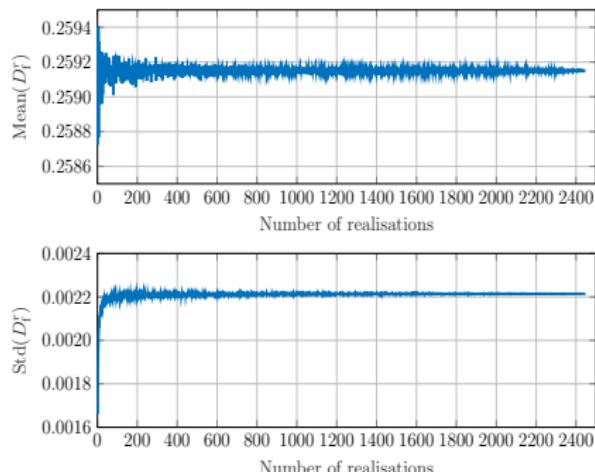
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Requirements:

$[15 - 35]$  min and  $[1 - 1.5]$  GB

Approx. 10 modes / Time-saving factors  $50 \sim 100$

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# Conclusion and future research

Efficient semi-incremental scheme for damage problems

Can handle variable amplitude and frequency loadings

Provides a minimal expansion of PGD

Most operations are done over decomposed QoI

Open-source code: [gitlab.com/shadialameddin/romfem](https://gitlab.com/shadialameddin/romfem)

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## Current research

Reducing the cost of evaluating the constitutive model

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Reducing the cost of evaluating the constitutive model

Thank you for your attention