

# Reduced order modeling of parameter dependent nonlinear eigenvalue bifurcation problems 

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## Outline

(1) Sparsity in PDE solution

Industrial application
Model reduction/sparsification Eigenvalue Methods

## Sparsity in PDE solutions

$\triangleright$ Numerical solution of PDE $L u=f$, with differential operator $L$ in a domain $\Omega \subset \mathbb{R}^{d}$ with boundary $\Gamma$ and appropriate boundary conditions given on $\Gamma$.
$\triangleright$ Let $\mathcal{V}$ be an ansatz function space in which we know or expect the solution to be.
$\triangleright$ Choose another (or the same) space $\mathcal{W}$ as test space.
$\triangleright$ Classical Galerkin or Petrov-Galerkin approach: Seek solution $u$ in some finite dimensional ansatz space $\mathcal{V}_{n} \subset \mathcal{V}$ (spanned by a basis or frame) $\mathcal{B}=\left\{\phi_{1}, \ldots, \phi_{n}\right\}$, i.e. the solution is represented as $u=\sum_{i=1}^{n} u_{i} \phi_{i}$ and $(L u-f, w)=0$ of $|(L u-f, w)|<\epsilon$ for all $w \in W$, where $\epsilon$ is a given tolerance.

## Questions

$\triangleright$ Can $u$ be sparsely represented in $\mathcal{V}$ ? Sure if the solution lies in $\mathcal{V}$, just take $u \in \mathcal{B}$.
$\triangleright$ Can $u$ be sparsely represented in $\mathcal{V}_{n} \subset \mathcal{V}$.
$\triangleright$ What is a good basis/frame of $\mathcal{V}_{n}$ so that $u$ can be sparsely represented.
$\triangleright$ What are conditions for the basis/frame so that the finite dimensional version $L_{n} u_{n}=f_{n}$ has a sparse $L_{h}$, or a sparse inverse $L_{h}^{-1}$.
$\triangleright$ Is there a cheap $(O(n))$ method to get a sparse solution?
$\triangleright$ Can we have all this together?
$\triangleright$ Is there a 'eierlegende Wollmilchsau', a swiss army knife for PDE solution?

## Outline

# (1) Sparsity in PDE solution <br> (2) Industrial application <br> Model reduction/sparsification Eigenvalue Methods 

## Disk brake squeal

Current project with Audi and Opel and several SMEs (2012-14) Joint with N. Gräbner, U. von Wagner, TU Berlin, Mechanics and N. Hoffmann, TU Hamburg-Harburg, Mechanics,
S. Quraishi, C. Schröder, TU Berlin Mathematics. Goals:
$\triangleright$ Develop mechanics based discrete FE model of disk brake.
$\triangleright$ Identification of energy dissipation effects.
$\triangleright$ Model and simulate nonlinear effects in brake squeaking.
$\triangleright$ Reduced order (compressed) model for a given range of disk speeds.
$\triangleright$ Sparse representation of operator and solution.
$\triangleright$ Finally, passive and active remedies to avoid squeaking.

## Disk brake



View of the brake model

## Finite element model



## Dynamics of disc brake

Differential-algebraic equation (DAE)

$$
M \ddot{q}+\left(C_{1}+\frac{\omega_{\text {ref }}}{\omega} C_{r}+\frac{\omega}{\omega_{\text {ref }}} C_{g}\right) \dot{q}+\left(K_{1}+K_{r}+\left(\frac{\omega}{\omega_{\text {ref }}}\right)^{2} K_{g}\right) q=f .
$$

$\triangleright M$ is symmetric semi-definite mass matrix,
$\triangleright C=C_{1}+C_{g}+C_{r}$ is a 'damping matrix'

- $C_{1}$ is symmetric,
- $C_{g}$ (due to gyroscopic effects) is skew-symmetric,
- $C_{r}$ is friction induced damping (symmetric),
- $\omega$ is the angular velocity, $\omega_{\text {ref }}$ reference.
$\triangleright K=K_{1}+K_{r}+K_{g}$ is a 'stiffness matrix'
- $K_{1}$ is symmetric, dominant part,
- $K_{r}$ describes circulatory effects (non symmetric),
- $K_{g}$ is geometric stiffness matrix.


## Complex eigenvalue analysis

$\triangleright$ Setting $q(t)=e^{\lambda t} u$, we get a quadratic eigenvalue problem (QEP):

$$
P_{\omega}(\lambda) u=\left(\lambda^{2} M+\lambda C(\omega)+K(\omega)\right) u=0 .
$$

$\triangleright$ Likelihood of brake to squeal is correlated with magnitude of positive real part of eigenvalue.
$\triangleright$ Compute eigenvalues in right half plane for lots of parameter values e.g. $\omega \in(2 \pi, 2 \pi \times 20)$.

## Nature of FE matrices

| $C=C_{1}+\frac{\omega_{\text {ref }}}{\omega} C_{r}+\frac{\omega}{\omega_{r e t}} C_{g}$, |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: |
| $\begin{aligned} & K=K_{1}+K_{r}+\left(\frac{\omega}{\omega_{\text {ref }}}\right)^{2} K_{g} \\ & n=842,638, \omega=17 \times 2 \pi \end{aligned}$ |  |  |  |  | $\begin{gathered} C_{1} \\ \square \\ \substack{n z=3 e+02 \\ C_{R}} \end{gathered}$ |
| matrix | pattern | 2-norm | structural rank |  |  |
| M | symm | 5e-2 | 842,623 | $\stackrel{\circ}{\circ}$ |  |
| $D_{1}$ | symm | 1e-19 | 160 | $n z=3 e+0$ | $n z=4 e+0$ |
| $\mathrm{D}_{\mathrm{G}}$ | skew | 1.5e-1 | 217500 | $\mathrm{K}_{1}$ | $\mathrm{K}_{\mathrm{R}}$ |
| $\mathrm{D}_{\mathrm{R}}$ | symm | 7e-2 | 2120 |  |  |
| $\mathrm{K}_{1}$ | symm | 2e13 | full | $n z=4 e+07$ | $n z=1 e+05$ |
| $K_{R}$ | - | 3 C 4 | 2110 | ${ }_{6}$ |  |
| $K_{G E O}$ | symm | 40 | 842,623 |  |  |

## Challenges

$\triangleright$ The discrete modeling is done directly with matrices, so space discretization cannot easily be done in an adaptive FEM way.
$\triangleright$ The set of ansatz functions (dictionary) is fixed, not a choice.
$\triangleright$ It is difficult to enrich the space with 'better functions'.
$\triangleright$ We have to work in an algebraic framework.
$\triangleright$ How to bring in sparsity?

## Outline

(1) Sparsity in PDE solution Eigenvalue Methods

## Projection approach

$\triangleright$ Project QEP: $P_{\omega}(\lambda) x=\left(\lambda^{2} M+\lambda C(\omega)+K(\omega)\right) x=0$ into a small $d$-dimension subspace $Q$ independent of $\omega$.
$\triangleright$ Projected QEP

- $\tilde{P}_{\omega}(\lambda)=Q^{T} P_{\omega}(\lambda) Q=\left(\lambda^{2} Q^{T} M Q+\lambda Q^{T} C(\omega) Q+Q^{T} K(\omega) Q\right)$
$\triangleright$ How to choose $Q$ to get sufficiently accurate approximation of eigenvalues with positive real part;
$\triangleright$ How to choose $Q$ to capture the important (to analyze and modify the squeaking) dynamics of the system;
$\triangleright$ Ideally $Q$ should contain good approximations to the wanted eigenvectors for all parameter values;
$\triangleright$ We should be able to construct $Q$ in a reasonable amount of time.


## Model reduction approaches

1. Traditional approach, often with Algebraic Multi Level Sub-structuring (AMLS).
2. New proper orthogonal decomposition (POD) based approach.

## Traditional approach

$\triangleright$ Traditional approach to get a subspace $Q$ :

- $Q_{T R A D}$ matrix of dominant eigenvectors.
- Select dominant eigenvectors by solving the GEVP $K_{1} v=-\lambda^{2} M v$
$\triangleright$ Advantages:
- Only need to solve a large sparse, symmetric and definite GEVP.
$\triangleright$ Disadvantages:
- Subspace does not take into account damping and parameter dependence.
- The reduced model often does not capture the important dynamics.
- Poor approximation of true eigenpairs.


## Real eigenforms

Undamped model without circulatory and gyroscopic forces: $\left(\lambda^{2} M+K+K_{g}\right) x=0$.


Figure: Trad. eigenmode at 1859 Hz

## Complex eigenforms



Figure: Eigenform at 1873 Hz with positive real part and a phase of 0, 45, 90, and 135.

## Measurement of brake vibrations



Gitter der Messpunkte


Betriebsschwingform $(1750 \mathrm{~Hz})$

Measurements indicate subcritical Hopf bifurcations, i.e. eigenvalues crossing imaginary axis for certain disk frequencies. Traditional approach deals only with purely imaginary evs.

## Stability regions, linear vs. nonlinear




Bifurcation diagram linear analysis (blue), nonlinear analysis (red). Coefficient of friction $\mu$ via disk frequency $\Omega$.

## Outline

# (1) Sparsity in PDE solution Industrial application Model reduction/sparsification 

(4) Eigenvalue Methods

## Linearization

We use the classical companion linearization to turn the quadratic into a linear generalized eigenvalue problem

$$
\boldsymbol{A}_{\tau}(\omega) x(\omega)=\lambda_{\tau} B_{\tau}(\omega) x(\omega)
$$

with

$$
\left[\begin{array}{cc}
K_{\tau}(\omega) & 0 \\
0 & I_{n}
\end{array}\right]\left[\begin{array}{c}
x(\omega) \\
\lambda_{\tau}(\omega) x(\omega)
\end{array}\right]=\lambda_{\tau}(\omega)\left[\begin{array}{cc}
-C_{\tau}(\omega) & -M_{\tau} \\
I_{n} & 0
\end{array}\right]\left[\begin{array}{c}
x(\omega) \\
\lambda_{\tau} x(\omega)
\end{array}\right] .
$$

## Shift and invert Arnoldi

$\triangleright$ Compute eigenvalue and eigenvector approximations near a given shift point $\tau$ via the Shift-and-invert Arnoldi method.
$\triangleright$ Given $v_{0} \in \mathbb{C}^{n}$ and $A \in \mathbb{C}^{n \times n}$, the Krylov subspace of $\mathbb{C}^{n}$ of order $k$ associated with $A$ is

$$
\mathcal{K}_{k}\left(A, v_{0}\right)=\operatorname{span}\left\{v_{0}, A v_{0}, A^{2} v_{0} \ldots, A^{k-1} v_{0}\right\} .
$$

$\triangleright$ Arnoldi obtains an orthonormal basis of this space and an Arnoldi relation

$$
A V_{k}=V_{k} H_{k}+f e_{k}^{*},
$$

$\triangleright$ The columns of $V_{k}$ are approximation of $k$-dimensional invariant subspace of $A$.
$\triangleright H_{k}$ is upper Hessenberg, its evs are Ritz approximations to evs of $A$ associated to $V_{k}$.
$\triangleright$ We apply Arnoldi with shift $\tau$ and frequency $\omega$ to the matrix $A=B_{\tau}(\omega)^{-1} A_{\tau}(\omega)$. In every step we have to multiply with $A_{\tau}(\omega)$ and to solve a linear system with the matrix $B_{\tau}(\omega)$.
$\triangleright$ We construct a measurement matrix $X \in \mathbb{R}^{n, k m}$ containing the 'unstable' eigenvectors for a sequence of angular velocities,

$$
X=\left[X\left(\omega_{1}\right), X\left(\omega_{2}\right), X\left(\omega_{3}\right), \ldots X\left(\omega_{k}\right)\right]
$$

$\triangleright$ Perform a singular value decomposition (SVD) of $X$

$$
X=\left[u_{1}, u_{2}, u_{3}, \ldots u_{k m}\right]\left[\begin{array}{ccccc}
\sigma_{1} & & & & \\
& \sigma_{2} & & & \\
& & \sigma_{3} & & \\
& & & \ddots & \\
& & & & \sigma_{k m}
\end{array}\right]\left[v_{1}, v_{2}, v_{3}, \ldots v_{k m}\right]^{T}
$$

## Compression

$\triangleright$ We use approximation

provided $\sigma_{d+1}, \sigma_{d+2}, \ldots \sigma_{k m}$ are small.
$\triangleright$ We choose $Q=\left[u_{1}, u_{2}, u_{3}, \ldots u_{d}\right]$ to project $P_{\omega}(\lambda)$.

## Some results on small $n \approx 5000$ matrices

$\triangleright$ POD for uniformly spaced $p$ parameters

$$
p=2^{j}+1, j=0,1,2,3
$$


$\triangleright$ Increasing dimension does not improve TRAD approach

## Realistic $n \approx 800,000$ matrices

$\triangleright$ The evp is completely singular $M, D$ have a 12 dimensional common nullspace and $K$ has relative size $10^{-14}$ in that nullspace.
$\triangleright$ Shifted matrix $\widetilde{K_{\tau}}=\tau^{2} \widetilde{M}+\tau \widetilde{C}+\widetilde{K}$ has condition number $\sim 10^{14}$ for a range of target points. Most likely due to bad FEM model.
$\triangleright$ Need to solve many large scale evps to get measurement matrix $\widetilde{X}=\left[X\left(\omega_{1}\right), X\left(\omega_{2}\right), X\left(\omega_{3}\right) \cdots X\left(\omega_{p}\right)\right]$.
$\triangleright$ It is not clear which parameter values $\omega_{i}$ are important.
$\triangleright$ Where to look for eigenvalues in the right half plane.
$\triangleright$ Scaling of matrices with scalar parameters to make them comparable in norm.
$\triangleright$ Diagonal scaling of matrices to improve conditioning.

## Which shifts to trust?



$\triangleright$ Different shift gives different evs in the overlapping region.

## Some CPU timings

$\triangleright$ Construction of subspace (One time investment)

- Each shift of eigs (Arnoldi method) ~ 20 min
- Eigenpairs for each parameter value $\sim 3$ targets $\sim 1$ hour
- POD measurement vectors for 2 parameters $\sim 2$ hours (or just 20 min on 6 processors)
- Constructing POD subspace (SVD) ~ 1 min
- Constructing 300 dimensional TRAD subspace $\sim 45 \mathrm{~min}$
$\triangleright$ Solution for every $\omega$
- Solution with 300 dimensional TRAD subspace $\sim 30$ sec
- Solution with 100 dimensional POD subspace $\sim 10 \mathrm{sec}$


## Evs for $\omega=17 \times 2 \pi$

$\triangleright$ POD model for $\omega=[1,20] \times 2 \pi$
$\triangleright$ Color coded with residual $\mathcal{R}=\frac{\left\|\left(\lambda_{i}^{2} M+\lambda_{i} C+K\right) u_{i}\right\|_{\infty}}{\left\|\left(1 \lambda_{i}{ }^{2}|M|+\left|\lambda_{i} \| C\right|+|K|\right) \mid u_{i}\right\|_{\infty}}$
$\triangleright U_{P O D}$ : 100, $U_{\text {TRAD }}: 300$ (Industry Recommendation)

$\triangleright$ all +'s are red (TRAD approach has very high residual)

## TRAD misses important eigenvalue



## TRAD misses important eigenvalue

$\triangleright$ Place shift point

$\tau=7.5+16500 i$ near an eigenvalue found from POD
Running eigs with this shift result in an exact eigenvalue
$\lambda=7.5414+16508 i$ very close to POD result

## TRAD misses important eigenvalue

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Running eigs with this shift result in an exact eigenvalue
$\lambda=7.5414+16508 i$
${ }^{11}$ very close to POD result
111.5
$\triangleright$ TRAD misses it

## What did we learn?

$\triangleright$ POD is better than traditional approach but not satisfactory.
$\triangleright$ Discrete FE and quasi-uniform grids followed by expensive model reduction is really a waste.
$\triangleright$ Numerical linear algebra methods that we currently use are not efficient (in particular those in commercially available codes).
$\triangleright$ For evp everything is partially heuristic.
$\triangleright$ Can we get error estimates? Can we bring in adaptivity? Dictionary learning?
$\triangleright$ Can we disprove the engineers that say that uniform mesh and brute force linear algebra is best.

## A compressed sensing point of view

$\triangleright$ The vectors $q(t)$ represent coefficient vectors for the infinite dimensional solution represented in an FEM basis $\phi_{1}(x, t), \ldots, \phi_{N}(x, t)$ in space-time.
$\triangleright$ The eigenvectors $x_{i}(\omega)$ also represent coefficient vectors in this FEM basis to synthesize the fundamental solution matrix of the DAE.
$\triangleright$ Every eigenvector $x_{i}(\omega)$ is the coefficient vector of a non-sparse function $\xi_{i}(x, \omega, t)$, because it typically linearly combines many FEM basis functions.
$\triangleright$ The POD basis represents a small set of linear combinations of the $\xi_{i}(\omega)$, given by functions $\psi_{j}(x, t) j=1, \ldots, d$ which are independent of $\omega$.
$\triangleright$ Consider the dictionary
$\mathcal{D}=\left\{\phi_{1}(x, \omega, t), \ldots, \phi_{N}(x, \omega, t)\right\} \cup\left\{\xi_{1}(x, \omega, t), \ldots, \xi_{\ell}(x, \omega, t)\right\}$.
$\triangleright$ Choosing the POD basis is selecting a small 'sparse' set of linear combinations from $\mathcal{D}$.

## Conclusions and Questions.

$\triangleright$ Real world industrial problems as motivation for studying, functions spaces, dictionaries, ...
$\triangleright$ Can we use this analogy to get convergence proofs, error bounds, complexity analysis?
$\triangleright$ What kind of sparsity should we go for?
$\triangleright$ How should we construct FE dictionaries?
$\triangleright$ Can we convince the engineers?
$\triangleright$ Can we make this practical?
$\triangleright$ Can we remove the brake squeal?

Thank you very much for your attention.

