

Computation of eigenvalues using multiscale techniques

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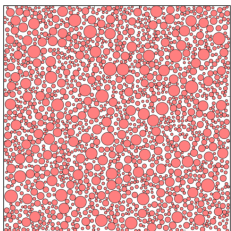
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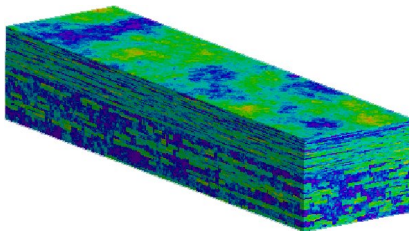
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Multiscale problems

Applications such as



▷ composite materials



▷ flow in a porous medium

require numerical solution of partial differential equations with rough data (module of elasticity, conductivity, or permeability).

Major challenge: Features on **multiple non-separated scales**.

Multiscale methods

Let A be rapidly varying data and consider a differential equation and its corresponding numerical approximation,

$$\mathcal{L}(A)u = f \qquad \mathcal{L}_h(A)u_h = f_h.$$

Classical finite element methods typically give

$$\|u - u_h\| \leq C(A, A')h^\gamma.$$

Multiscale methods seek an upscaled representation

$$\mathcal{L}_H(A)u_H = f_H$$

fulfilling $\|u_h - u_H\| \leq C(A)H^\gamma$ with C independent of A' .

How well is the spectrum of \mathcal{L} preserved?

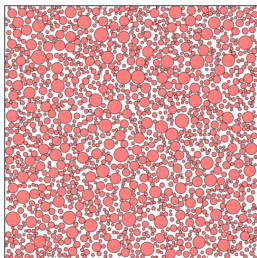
- 1 **Model problem**
- 2 Upscaling technique and error analysis
- 3 Numerical experiments
- 4 Applications to non-linear eigenvalue problems
- 5 Conclusions

Model multiscale eigenvalue problem

Prototypical self-adjoint eigenvalue problem

$$-\nabla \cdot A \nabla u = \lambda u \quad \text{in } \Omega \quad u = 0 \quad \text{on } \partial\Omega$$

with data $0 < \alpha \leq A \leq \beta < \infty$

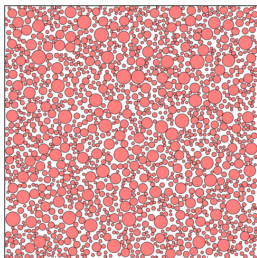


Model multiscale eigenvalue problem

Prototypical self-adjoint eigenvalue problem (variational form): find $u \in V := H_0^1(\Omega)$ and $\lambda \in \mathbb{R}$ such that

$$a(u, v) := \int_{\Omega} (A \nabla u) \cdot \nabla v \, dx = \lambda \int_{\Omega} u \cdot v \, dx \quad \text{for all } v \in V$$

with data $0 < \alpha \leq A \leq \beta < \infty$

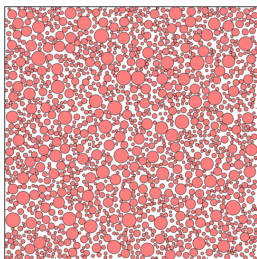


Model multiscale eigenvalue problem

Prototypical self-adjoint eigenvalue problem (FE approximation):
 $u_h \in V_h \subset V$ and $\lambda_h \in \mathbb{R}$ such that

$$a(u_h, v) := \int_{\Omega} (A \nabla u_h) \cdot \nabla v \, dx = \lambda_h \int_{\Omega} u_h \cdot v \, dx \quad \text{for all } v \in V_h$$

with data $0 < \alpha \leq A \leq \beta < \infty$



Model multiscale eigenvalue problem

Prototypical self-adjoint eigenvalue problem (FE approximation):
 $u_h \in V_h \subset V$ and $\lambda_h \in \mathbb{R}$ such that

$$a(u_h, v) := \int_{\Omega} (\mathbf{A} \nabla u_h) \cdot \nabla v \, dx = \lambda_h \int_{\Omega} u_h \cdot v \, dx \quad \text{for all } v \in V_h$$

with data $0 < \alpha \leq A \leq \beta < \infty$

Numerical error (piecewise linear continuous FE approximation)

- For an eigenpair $(u^{(k)}, \lambda^{(k)})$ with $u^{(k)} \in H^2(\Omega)$ it holds

$$\lambda^{(k)} \leq \lambda_h^{(k)} \leq \lambda^{(k)} + C(\mathbf{A}, \mathbf{A}', k) h^2,$$
$$\| \| u^{(k)} - u_h^{(k)} \| \| := \| \mathbf{A}^{1/2} \nabla (u^{(k)} - u_h^{(k)}) \|_{L^2(\Omega)} \leq C(\mathbf{A}, \mathbf{A}', k) h.$$

- The mesh size h has to resolve the variations in \mathbf{A} , e.g. $h < \epsilon$ if \mathbf{A} is periodic.

Investigate how the Localized Orth. Decomposition (LOD) in



A. Målqvist and D. Peterseim.

Localization of Elliptic Multiscale Problems.
Mathematics of Computation 2014

preserves the (low part of the) spectrum of $-\nabla \cdot A \nabla$.

Without assumptions on scales (A') or regularity (u):

$$\lambda_h \leq \lambda_H^{\text{ms}} \leq \lambda_h + CH^4,$$
$$\|u_h - u_H^{\text{ms}}\| \leq CH^2.$$



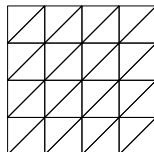
A. Målqvist and D. Peterseim.

Computation of eigenvalues by numerical upscaling.
accepted for publication in Numerische Mathematik

- 1 Model problem
- 2 **Upscaling technique and error analysis**
- 3 Numerical experiments
- 4 Applications to non-linear eigenvalue problems
- 5 Conclusions

Multiscale decomposition

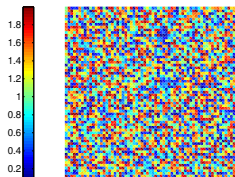
- (coarse) FE mesh \mathcal{T} with parameter H
- P1-FE space $V_H := \{v \in V \mid \forall T \in \mathcal{T}, v|_T \in P_1(T)\}$
- $\mathfrak{I}_{\mathcal{T}} : V \rightarrow V_H$ a Clément interpolation operator



Decomposition

$$V = V_H \oplus V^f \quad \text{with } V^f := \text{kernel } \mathfrak{I}_{\mathcal{T}} = \{v \in V \mid \mathfrak{I}_{\mathcal{T}} v = 0\}$$

Example:



rough coefficient

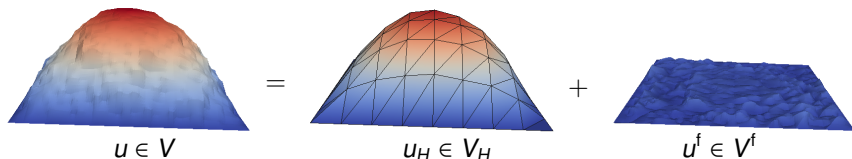
Multiscale decomposition

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- P1-FE space $V_H := \{v \in V \mid \forall T \in \mathcal{T}, v|_T \in P_1(T)\}$
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Example:



Orthogonalization

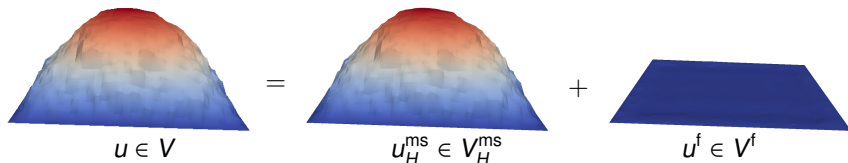
- For each $v \in V_H$ define finescale projection $\mathfrak{F}v \in V^f$ by

$$a(\mathfrak{F}v, w) = a(v, w) \quad \text{for all } w \in V^f$$

Orthogonal Decomposition

$$V = V_H^{\text{ms}} \oplus V^f \quad \text{with } V_H^{\text{ms}} := (V_H - \mathfrak{F}V_H)$$

Example:



Ideal multiscale representation

Given the space V_H^{ms} we construct a Galerkin approximation:

Ideal method

Find $u_H^{\text{ms}} \in V_H^{\text{ms}}$, $\lambda_H^{\text{ms}} \in \mathbb{R}$ such that

$$a(u_H^{\text{ms}}, v) = \lambda_H^{\text{ms}}(u_H^{\text{ms}}, v), \quad \forall v \in V_H^{\text{ms}}.$$

- We note that $\dim(V_H^{\text{ms}}) = \dim(V_H)$.
- For V_H^{ms} to be useful we need a discrete localized basis.
- But first of all we need to show that λ_H^{ms} is a good approximation of λ .

A priori error bound (ideal case)

For the k :th eigenvalue it holds

Theorem

$$\lambda^{(k)} \leq \lambda_H^{ms,(k)} \leq \lambda^{(k)} + CH^4,$$

C independent of A' and only H^1 -regularity of the eigenfunctions.

Sketch of proof for the *lowest* eigenvalue:

- Let $u^{(1)} := u = u_c + u_f$ with $u_c \in V_H^{ms}$ and $u_f \in V^f$, such that $\|u\|_{L^2(\Omega)} = 1$. Then

$$\begin{aligned} \lambda_H^{ms,(1)} &\leq \frac{a(u_c, u_c)}{(u_c, u_c)} \leq \frac{a(u, u)}{(u_c, u_c)} = \frac{a(u, u)}{(u - u_f, u - u_f)} \\ &= \frac{\lambda^{(1)}}{(u, u) - 2(u, u_f) + (u_f, u_f)} \leq \frac{\lambda^{(1)}}{1 - 2(u, u_f)}. \end{aligned}$$

A priori error bound (ideal case)

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$$\lambda^{(k)} \leq \lambda_H^{ms,(k)} \leq \lambda^{(k)} + CH^4,$$

C independent of A' and only H^1 -regularity of the eigenfunctions.

Sketch of proof for the *lowest* eigenvalue:

- Since $\mathfrak{I}_{\mathcal{T}} u_f = 0$, $(\mathfrak{I}_{\mathcal{T}} u, u_f) = 0$ (weighted Clement, CV99), $a(u_c, u_f) = 0$, and $\|u\|^2 = \lambda^{(1)}$, we have

$$\begin{aligned} (u, u_f) &= (u - \mathfrak{I}_{\mathcal{T}} u, u_f - \mathfrak{I}_{\mathcal{T}} u_f) \leq CH^2 \|u\| \cdot \|u_f\| \leq C' H^2 \|u_f\|, \\ \|u_f\|^2 &= a(u, u_f) = \lambda^{(1)} (u - \mathfrak{I}_{\mathcal{T}} u, u_f - \mathfrak{I}_{\mathcal{T}} u_f) \leq CH^2 \|u_f\|. \end{aligned}$$

- We conclude $\lambda_H^{ms,(1)} \leq \frac{\lambda^{(1)}}{1-CH^4} \leq \lambda^{(1)} + 2CH^4$.

A priori error bound (ideal case)

For the k :th eigenfunction it holds

Theorem

$$\| \| u^{(k)} - u_H^{ms,(k)} \| \| \leq CH^2,$$

C independent of A' and only H^1 -regularity of the eigenfunctions.

- Similar arguments using $\mathfrak{S}_{\mathcal{T}} u_f = 0$ and $(\mathfrak{S}_{\mathcal{T}} u, u_f) = 0$.
- Only $H^1(\Omega)$ regularity is assumed.

Can we find a localized discrete basis that approximates V_H^{ms} ?

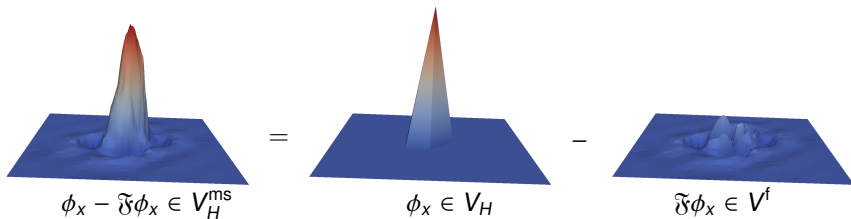
Modified nodal basis

- \mathcal{N} denotes set of interior vertices of \mathcal{T}
- $\phi_x \in V_H$ denotes classical nodal basis function ($x \in \mathcal{N}$)
- $\mathfrak{F}\phi_x \in V^f$ denotes finescale correction of ϕ_x ($x \in \mathcal{N}$)

Ideal multiscale FE space

$$V_H^{\text{ms}} = \text{span} \{ \phi_x - \mathfrak{F}\phi_x \mid x \in \mathcal{N} \}$$

Example



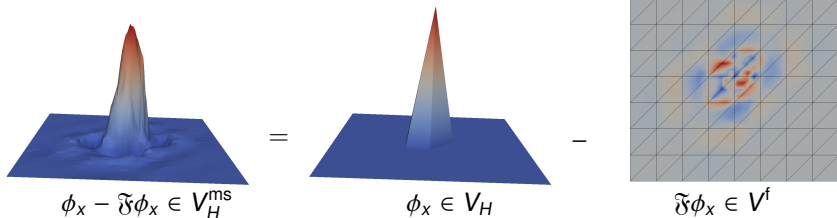
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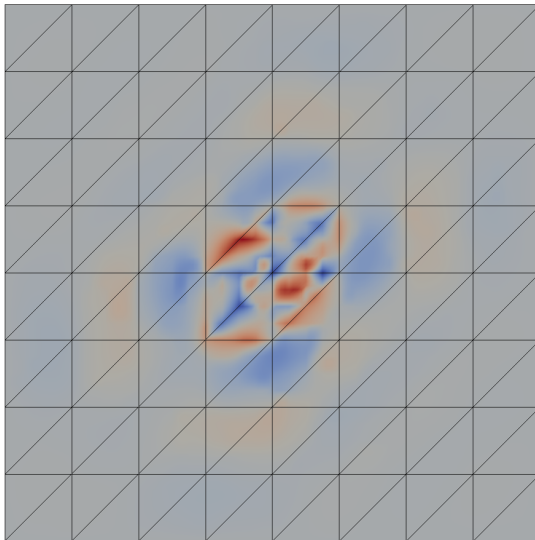
Ideal multiscale FE space

$$V_H^{\text{ms}} = \text{span} \{ \phi_x - \mathfrak{F}\phi_x \mid x \in \mathcal{N} \}$$

Example

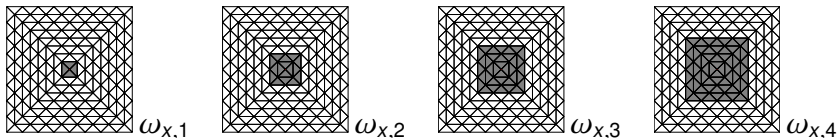


Modified nodal basis



Localization

- Define nodal patches of ℓ -th order $\omega_{x,\ell}$ about $x \in \mathcal{N}$



- Localized corrections $\mathfrak{F}\phi_{x,\ell} \in V^f(\omega_{x,\ell}) := \{v \in V^f \mid v|_{\Omega \setminus \omega_{x,\ell}} = 0\}$
solve

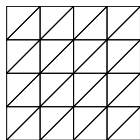
$$a(\mathfrak{F}\phi_{x,\ell}, w) = a(\phi_x, w) \quad \text{for all } w \in V^f(\omega_{x,\ell})$$

Localized multiscale FE spaces

$$V_{H,\ell}^{\text{ms}} = \text{span}\{\phi_x - \mathfrak{F}\phi_{x,\ell} \mid x \in \mathcal{N}\}$$

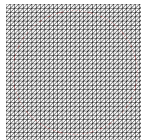
Fine scale discretization

- Finescale mesh



\mathcal{T}

mesh refinement



\mathcal{T}_h with $h \leq H$

- Reference FE space

$$V_h := \{v \in V \mid \forall T \in \mathcal{T}(\Omega), v|_T \in P_1(T)\}$$

- Reference FE solution $u_h \in V_h$ and $\lambda_h \in \mathbb{R}$ solves

$$a(u_h, v) = \lambda_h(u_h, v) \quad \text{for all } v \in V_h$$

- Fully discrete corrections $\mathfrak{F}\phi_{x,\ell}^h \in V_h^f(\omega_{x,\ell}) := V^f(\omega_{x,\ell}) \cap V_h$:

$$a(\mathfrak{F}\phi_{x,\ell}^h, w) = a(\phi_x, w) \quad \text{for all } w \in V_h^f(\omega_{x,\ell})$$

Localized Orthogonal Decomposition (LOD)

Fully discrete multiscale FE spaces

$$V_{H,\ell}^{\text{ms},h} = \text{span}\{\phi_x - \mathfrak{F}\phi_{x,\ell}^h \mid x \in \mathcal{N}\}$$

Fully discrete multiscale approximation $u_{H,\ell}^{\text{ms},h} \in V_{H,\ell}^{\text{ms},h}$, $\lambda_{H,\ell}^{\text{ms},h} \in \mathbb{R}$

$$a(u_{H,\ell}^{\text{ms},h}, v) = \lambda_{H,\ell}^{\text{ms},h} (u_{H,\ell}^{\text{ms},h}, v) \quad \text{for all } v \in V_{H,\ell}^{\text{ms},h}$$

Remarks:

- $\dim V_{H,\ell}^{\text{ms},h} = |\mathcal{N}| = \dim V_H$
- The basis functions have local support, with overlap depending on ℓ , and are independent.

A priori error analysis (discrete case)

Lemma (Truncation error)

$$\| \mathfrak{F}\phi_x^h - \mathfrak{F}\phi_{x,\ell}^h \| \leq C_1 \gamma^\ell \| \mathfrak{F}\phi_x^h \|.$$

$C_1 < \infty$ and $\gamma < 1$ depends on β/α , not A' .

By choosing $\ell = C_2 \log(H^{-1})$ with appropriate C_2 we guarantee that the truncation leads to a higher order perturbation:

Theorem

$$\begin{aligned} \lambda_h^{(k)} &\leq \lambda_{H,\ell}^{ms,h,(k)} \leq \lambda_h^{(k)} + CH^4, \\ \| u_h^{(k)} - u_{H,\ell}^{ms,h,(k)} \| &\leq CH^2, \end{aligned}$$

with C independent of A' and the regularity of the eigenfunctions.

A priori error analysis (discrete case)

The result can be improved using a postprocessing technique:



J. Xu and A. Zhou.

A two-grid discretization scheme for eigenvalue problems.
Mathematics of Computation 2001.

Find $u_h^p \in V_h$ s.t.

$$a(u_h^p, v) = \lambda_{H,\ell}^{\text{ms},h}(u_{H,\ell}^{\text{ms},h}, v), \quad v \in V_h,$$

and letting $\lambda_h^p = a(u_h^p, u_h^p)/(u_h^p, u_h^p)$.

Theorem

$$\begin{aligned} \lambda_h^{(k)} &\leq \lambda_h^{p,(k)} \leq \lambda_h^{(k)} + CH^6, \\ \|\| u_h^{(k)} - u_h^{p,(k)} \|\| &\leq CH^4. \end{aligned}$$

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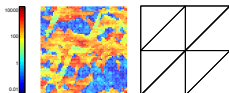
Eigenvalue Problem



k	$\lambda_h^{(k)}$	$e^{(k)}(1/2 \sqrt{2})$	$e^{(k)}(1/4 \sqrt{2})$	$e^{(k)}(1/8 \sqrt{2})$	$e^{(k)}(1/16 \sqrt{2})$
1	9.6436869	0.003494567	0.000034466	0.000000546	0.000000010
2	15.1989274	0.009621397	0.000079887	0.000000845	0.000000010
3	19.7421815	0.023813222	0.000213097	0.000002073	0.000000023
4	29.5281571	0.096910157	0.000724615	0.000006574	0.000000076
5	31.9265496	0.094454625	0.000874659	0.000009627	0.000000138
6	41.4922250	-	0.002395227	0.000019934	0.000000254
7	44.9604884	-	0.002443271	0.000019683	0.000000223
8	49.3631826	-	0.003651870	0.000028869	0.000000308
9	49.3655623	-	0.004266472	0.000032835	0.000000355
10	56.7389993	-	0.006863742	0.000055219	0.000000618
11	65.4085991	-	0.011534878	0.000082414	0.000000856
12	71.0947630	-	0.012596114	0.000090083	0.000001002
13	71.6064671	-	0.014249938	0.000098736	0.000001006
14	79.0043994	-	0.021801461	0.000164436	0.000001605
15	89.3706421	-	0.033550079	0.000211985	0.000002296
16	92.3648207	-	0.040060692	0.000239441	0.000002295
17	97.4459210	-	0.037438984	0.000284936	0.000002724
18	98.7545147	-	0.044544409	0.000269854	0.000002559
19	98.7545639	-	0.047835987	0.000276139	0.000002539
20	101.6755971	-	0.038203654	0.000297356	0.000002909

Table : Errors $e^{(k)}(H) =: \frac{\lambda_H^{\text{ms},(k)} - \lambda_h^{(k)}}{\lambda_h^{(k)}}$ and $h = 2^{-7} \sqrt{2}$.

Eigenvalue Problem

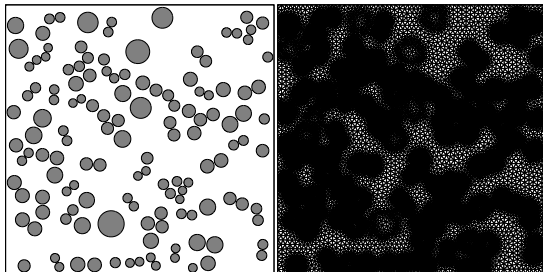


k	$\lambda_h^{(k)}$	$e^{(k)}(1/2\sqrt{2})$	$e^{(k)}(1/4\sqrt{2})$	$e^{(k)}(1/8\sqrt{2})$	$e^{(k)}(1/16\sqrt{2})$
1	21.4144522	5.472755371	0.237181706	0.010328293	0.000781683
2	40.9134676	-	0.649080539	0.032761482	0.002447049
3	44.1561133	-	1.687388874	0.097540102	0.004131422
4	60.8278691	-	1.648439518	0.028076168	0.002079812
5	65.6962136	-	2.071005692	0.247424446	0.006569640
6	70.1273082	-	4.265936007	0.232458016	0.016551520
7	82.2960238	-	3.632888104	0.355050163	0.013987920
8	92.8677605	-	6.850048057	0.377881216	0.049841235
9	99.6061234	-	10.305084010	0.469770376	0.026027378
10	109.1543283	-	-	0.476741452	0.005606426
11	129.3741945	-	-	0.505888044	0.062382302
12	138.2164330	-	-	0.554736550	0.039487317
13	141.5464639	-	-	0.540480876	0.043935515
14	145.7469718	-	-	0.765411709	0.034249528
15	152.6283573	-	-	0.712383825	0.024716759
16	155.2965039	-	-	0.761104705	0.026228034
17	158.2610708	-	-	0.749058367	0.091826207
18	164.1452194	-	-	0.840736127	0.118353184
19	171.1756923	-	-	0.946719951	0.111314058
20	179.3917590	-	-	0.928617606	0.119627862

Table : Errors $e^{(k)}(H) =: \frac{\lambda_H^{\text{ms},(k)} - \lambda_h^{(k)}}{\lambda_h^{(k)}}$ and $h = 2^{-7} \sqrt{2}$.

Eigenvalue problem, non-nested meshes

In many applications $V_H \not\subset V_h$, e.g., a domain with inclusions,

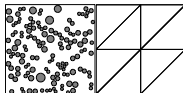


where A varies between 100 (gray) and 1 (white).

Using the nodal interpolant $I_h : C(\Omega) \rightarrow V_h$ we can construct a space $\tilde{V}_H = \{I_h \phi_x, x \in \mathcal{N}\} \subset V_h$ with $V_H = \text{span}(\phi_x) \not\subset V_h$.

Then $\mathfrak{I}_{\mathcal{T}} : V \rightarrow \tilde{V}_H$ is used to get V^f and V_H^{ms} .

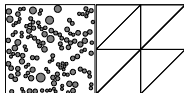
Eigenvalue Problem, non-nested meshes



ℓ	$\lambda_h^{(\ell)}$	$e^{(\ell)}(1/2\sqrt{2})$	$e^{(\ell)}(1/4\sqrt{2})$	$e^{(\ell)}(1/8\sqrt{2})$	$e^{(\ell)}(1/16\sqrt{2})$
1	25.6109462	0.025518831	0.000572341	0.000017083	0.000000700
2	58.9623566	-	0.005235813	0.000090490	0.000002710
3	67.5344854	-	0.006997582	0.000154850	0.000006488
4	98.2808694	-	0.023497502	0.000358178	0.000011675
5	121.2290664	-	0.052366141	0.000563438	0.000016994
6	125.2014779	-	0.066627585	0.000747688	0.000019934
7	156.0597873	-	0.145676350	0.001579177	0.000034329
8	168.2376096	-	0.095360287	0.001320185	0.000043781
9	197.4467434	-	0.343991317	0.002888471	0.000049479
10	209.4657306	-	-	0.003223901	0.000056318
11	222.4472476	-	-	0.003431462	0.000080284
12	245.5656759	-	-	0.005906282	0.000102243
13	253.7074603	-	-	0.006215809	0.000121646
14	288.0756442	-	-	0.013859535	0.000180899
15	298.8903269	-	-	0.010587124	0.000138404
16	311.4410556	-	-	0.012159268	0.000161510
17	324.6865434	-	-	0.012143676	0.000176624
18	336.7931865	-	-	0.016554437	0.000233067
19	379.5697606	-	-	0.023254268	0.000325324
20	386.9938901	-	-	0.028772395	0.000383532

Table : Errors $e^{(k)}(H) =: \frac{\lambda_H^{\text{ms},(k)} - \lambda_h^{(k)}}{\lambda_h^{(k)}}$ and $h \approx 2^{-9} - 2^{-7}$.

Eigenvalue Problem, non-nested, postpr.



ℓ	$\lambda_h^{(\ell)}$	$e^{(\ell)}(1/2\sqrt{2})$	$e^{(\ell)}(1/4\sqrt{2})$	$e^{(\ell)}(1/8\sqrt{2})$	$e^{(\ell)}(1/16\sqrt{2})$
1	25.6109462	0.001559704	0.000003765	0.000000008	3.5e-10
2	58.9623566	-	0.000191532	0.000000213	1.9e-08
3	67.5344854	-	0.000284980	0.000000474	0.000000001
4	98.2808694	-	0.002239689	0.000002253	0.000000004
5	121.2290664	-	0.007461217	0.000005065	0.000000008
6	125.2014779	-	0.011284614	0.000006826	0.000000008
7	156.0597873	-	0.042466017	0.000023867	0.000000024
8	168.2376096	-	0.025093182	0.000027547	0.000000042
9	197.4467434	-	0.186960343	0.000072471	0.000000051
10	209.4657306	-	-	0.000105777	0.000000079
11	222.4472476	-	-	0.000131569	0.000000129
12	245.5656759	-	-	0.000286351	0.000000213
13	253.7074603	-	-	0.000268463	0.000000255
14	288.0756442	-	-	0.000915102	0.000000473
15	298.8903269	-	-	0.000762135	0.000000403
16	311.4410556	-	-	0.000873769	0.000000504
17	324.6865434	-	-	0.000955392	0.000000642
18	336.7931865	-	-	0.001335246	0.000000977
19	379.5697606	-	-	0.002896202	0.000001886
20	386.9938901	-	-	0.007202657	0.000001908

Table : Errors $e^{(k)}(H) =: \frac{\lambda_H^{\text{ms},(k)} - \lambda_h^{(k)}}{\lambda_h^{(k)}}$ and $h \approx 2^{-9} - 2^{-7}$.

- 1 Model problem
- 2 Upscaling technique and error analysis
- 3 Numerical experiments
- 4 **Applications to non-linear eigenvalue problems**
- 5 Conclusions

The quadratic eigenvalue problem

Consider the quadratic eigenvalue problem (QEP): find $u \in V$, $\|u\|_{L^2(\Omega)} = 1$, and $\lambda \in \mathbb{C}$ such that

$$(A\nabla u, \nabla v) + \lambda c(u, v) + \lambda^2(u, v) = 0, \quad \forall v \in V.$$

This equation appears in structural mechanics and describes damped vibrations. Discretization gives an algebraic QEP,

$$Kx + \lambda Cx + \lambda^2 Mx = 0,$$

where K is stiffness, C is damping, and M is mass matrix.

If C is symmetric, real, and positive: the eigenvalues are complex conjugate with negative real part.



F. Tisseur and K. Meerbergen.

The quadratic eigenvalue problem SIAM Review 2001.

The quadratic eigenvalue problem

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- Rayleigh damping $C = \alpha_0 M + \alpha_1 K$ leads to unchanged eigenmodes.
- Systems of Rayleigh damped components, α_0, α_1 are functions.

QEP: Linearization and approximation

Linearization: ($y = \lambda x$)

$$\begin{bmatrix} K & 0 \\ 0 & M \end{bmatrix} \begin{bmatrix} x \\ y \end{bmatrix} = \lambda \begin{bmatrix} -C & -M \\ M & 0 \end{bmatrix} \begin{bmatrix} x \\ y \end{bmatrix}$$

On weak form: find (u_1, u_2) and $\lambda \in \mathbb{C}$ such that,

$$a(u, v) = \lambda b(u, v),$$

$$a(u, v) = (A \nabla u_1, \nabla v_1) + (u_2, v_2),$$

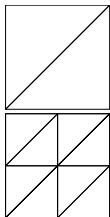
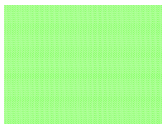
$$b(u, v) = -c(u_1, v_1) - (u_2, v_1) + (u_1, v_2).$$

Multiscale formulation:

Find $u_H^{\text{ms}} \in V_H^{\text{ms}} \times V_H^{\text{ms}}$ and $\lambda_H^{\text{ms}} \in \mathbb{C}$ such that,

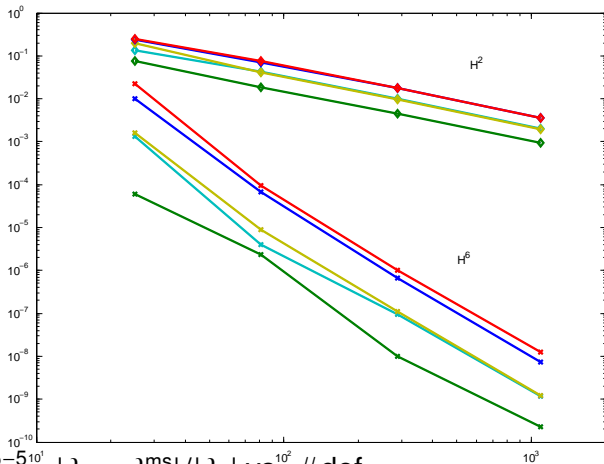
$$a(u_H^{\text{ms}}, v) = \lambda_H^{\text{ms}} b(u_H^{\text{ms}}, v), \quad \forall v \in V_H^{\text{ms}} \times V_H^{\text{ms}}.$$

Numerical experiment



$$H = 2^{-1}, 2^{-2}, \dots, 2^{-5 \cdot 10^1}$$

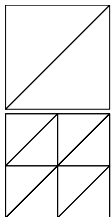
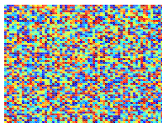
$$h = 2^{-6}, k = \infty$$



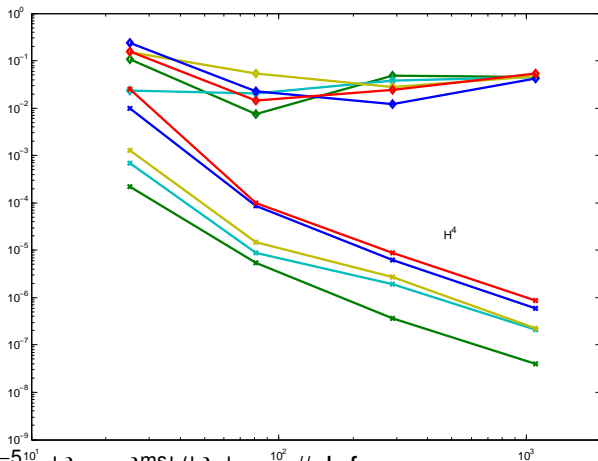
$$|\lambda_h - \lambda_H^{\text{ms}}| / |\lambda_h| \text{ vs. } \# \text{dof}$$

$$A = 1, c(u, v) = \int_{\Omega} (1 + \sin(10x)) u v \, dx$$

Numerical experiment



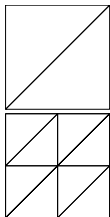
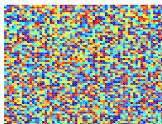
$$H = 2^{-1}, 2^{-2}, \dots, 2^{-5 \cdot 10^1}$$
$$h = 2^{-6}, k = \infty$$



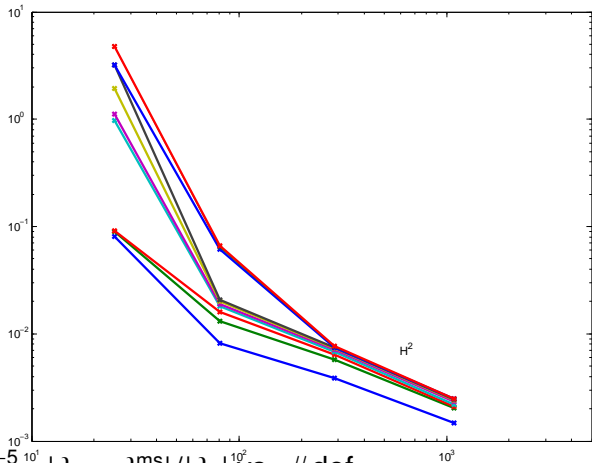
$|\lambda_h - \lambda_H^{\text{ms}}| / |\lambda_h|$ vs. 10^2 #dof

$$A \text{ (pic)}, c(u, v) = \int_{\Omega} (1 + \sin(10x)) u v \, dx$$

Numerical experiment



$H = 2^{-1}, 2^{-2}, \dots, 2^{-5}$
 $h = 2^{-6}, k = \infty$



$$\pi c(u, v) = \int_{\Omega} \left(\text{atan}(10x - 5) + \frac{\pi}{2} \right) A \nabla u \nabla v \, dx$$

QEP: Analysis

We note that the operator $B : V \times V \rightarrow V \times V$ defined by,

$$a(Bu, v) = b(u, v),$$

has eigenvalues $Bu = \mu u = \lambda^{-1} u$ since,

$$b(u, v) = a(Bu, v) = a(\mu u, v) = \mu a(u, v).$$

Furthermore, a is coercive, bounded, symmetric and b is bounded if $|c(u, v)| \leq C\|u\|_V\|v\|_V$.

 J. Descloux, N. Nassif, and J. Rappaz.

On spectral approximation I & II. RAIRO 1978. (non compact)

Error in eigenvalues

Theorem (Descloux, Nassif, and Rappaz)

Let λ be an eigenvalue of B with algebraic multiplicity m . Let $\{\tilde{\lambda}_j\}_{j=1}^m$ be Galerkin approximations using a discretized function space $\tilde{V} \subset V$. Then,

$$\left| \lambda - \left(\frac{1}{m} \sum_{j=1}^m (\tilde{\lambda}_j)^{-1} \right)^{-1} \right| \leq C \sup_{u \in M(\lambda)} \inf_{\chi \in \tilde{V}} \|u - \chi\|_V \cdot \sup_{v \in M^*(\lambda)} \inf_{\chi \in \tilde{V}} \|v - \chi\|_V,$$

where $M(\lambda)$ is the generalized eigenvectors and $M^*(\lambda)$ is the generalized adjoint eigenvectors associated with λ .

The choice $\tilde{V} = V_H^{\text{ms}}$ fulfills the requirements for this theorem.

Error in eigenvalues

Conjecture

Let $c(u, v) = \int_{\Omega} \alpha_0 A \nabla u \cdot \nabla v \, dx + \int_{\Omega} \alpha_1 uv \, dx$, with $\|\alpha_i - \mathfrak{I}_{\mathcal{T}} \alpha_i\| \leq CH$ and let λ be an eigenvalue with algebraic multiplicity m . Let $\{\lambda_{h,j}^{ms}\}_{j=1}^m$ be LOD approximations using a $V_H^{ms} \subset V$. Then,

$$\sup_{u \in M(\lambda)} \inf_{\chi \in V_H^{ms}} \|u - \chi\|_V \cdot \sup_{v \in M^*(\lambda)} \inf_{\chi \in V_H^{ms}} \|v - \chi\|_V \leq CH^2,$$

and therefore,

$$\left| \lambda - \left(\frac{1}{m} \sum_{j=1}^m (\lambda_{H,j}^{ms})^{-1} \right)^{-1} \right| \leq CH^2.$$

This holds without assuming more regularity than $u \in H^1(\Omega)$. Localization and fine grid discretization is not considered.

The Gross-Pitaevskii equation

Consider the Gross-Pitaevskii equation: find $u \in V$, $\|u\|_{L^2(\Omega)} = 1$, and $\lambda \in \mathbb{R}$ such that

$$(A\nabla u, \nabla v) + (bu, v) + (u^3, v) = \lambda(u, v), \quad \forall v \in V.$$

The equation describes the quantum states of a boson gas cooled down to an ultra-low temperature.

- We reuse the same discrete space $V_{H,\ell}^{ms,h}$ i.e. we ignore the low order non-linearity on the fine scale.
- We then solve the upscaled non-linear eigenvalue problem on the coarse scale.

 P. Henning, A. Målqvist, and D. Peterseim.

Two-level discretization techniques for **ground state** computations of Bose-Einstein condensates.

SIAM Journal on Numerical Analysis 2014.

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Theorem

$$\begin{aligned} \lambda &\leq \lambda_h^p \leq \lambda + CH^2 \|u - u_h\|_{H^1(\Omega)} + CH^4, \\ \|u - u_h^p\|_{H^1(\Omega)} &\leq C \|u - u_h\|_{H^1(\Omega)} + CH^3. \end{aligned}$$

for the ground state, with C independent on the regularity of u and variations in A .

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Conclusion

- The Localized Orthogonal Decomposition (LOD) technique preserves the low spectrum of the operator. In particular the eigenvalue error is proportional to H^4 after postprocessing H^6 .
- Numerical experiments indicates even higher rates possibly due to additional regularity in the solution that is not taken advantage of in the analysis.
- The technique is applicable also for non-linear eigenvalue problems without pre-asymptotic effects in the convergence.
- More work is needed in the analysis for the quadratic eigenvalue problem.
- Numerical tests with more complicated damping is needed.

Thank you for your attention!