An Algorithm for the Generalized Symmetric Tridiagonal Eigenvalue Problem

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Generalized Eigenvalue Problem

Def

 $T, S \in \mathbb{R}^{n \times n}$. We call (T, S) pencil.

We consider *only*

symmetric and

tridiagonal

pencils.

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Def (Problem)

Find $\lambda \in \mathbb{R}$ such that $Tv = \lambda Sv$, with $v \in \mathbb{C}^n$. T, S symmetric implies $\lambda \in \mathbb{R}$.

Algorithm philosophy

We find zeros of the polynomial equation

$$\mathcal{F}_{(T,S)}(\lambda) = \det(T - \lambda S) = 0$$

using an iterative method, living on real line.

Brainstorming

We want:

Fast and secure iterative method.

Starting points for our method.

Scalability.

We have: Lague

Laguerre's method.

Cuppen's divide and conquer method.

Symmetric tridiagonal matrices.

We add:

Unreducible condition.

Dynamic programming (Bottom-up).

Efficient matrix storing.

Rapid tour

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Unreducible pencil

Def (as in [?]) $(T,S) \text{ is an } \textit{unreducible pencil if } t_{i,i+1}^2 + s_{i,i+1}^2 \neq 0$ for $i=1,2,\ldots,n-1.$

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(T,S) is an unreducible pencil if $t_{i,i+1}^2 + s_{i,i+1}^2 \neq 0$ for i = 1, 2, ..., n - 1.

exempli gratia:

Bad
$$T = I, S = 0$$

 $T = I, S = I$

Good
$$T = I, S = trid(-1, 2, -1)$$

 $T = trid(-1, 2, -1), S = trid(-1, 2, -1)$
 $T = trid(rnd_{sub}, rnd_{diag}, rnd_{sub}), S = I,$
with rnd_{sub} random number $\neq 0$.

Matrix storing

$$T = trid(sub, diag, super)$$

But T is symmetric, so sub = super. We define and use

```
integer, parameter :: dp = kind(1.d0)
real(dp), dimension(1:n,0:1) :: T, S
```

Listing 1: T, S as couple of array

```
with T(:,0) = diag and T(:,1) = super.
```

Remark

We don't use T(1,1) and S(1,1).

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$$g_u(X) = (x - X)^2 \sum_{i=1}^n \frac{(u - \lambda_i)^2}{(x - \lambda_i)^2} - (u - X)^2$$

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Bolzano's Theorem tell us that there are two zeros of g_u between λ_m and λ_{m+1} . We call them X_-, X_+ .

$$\lambda_m < X_- < x < X_+ < \lambda_{m+1}$$

We have one freedom: the u parameter.

Calling $\hat{X}_- = min_u X_-$ and $\hat{X}_+ = max_u X_+$ we can obtain

$$\lambda_m \approx \hat{X}_- < x < \hat{X}_+ \approx \lambda_{m+1}$$

and with algebraic manipulations:

$$\hat{X}_{-}, \hat{X}_{+} = L_{\pm}(x) = x + \frac{n}{-\frac{f'}{f} \pm \sqrt{(n-1)[(n-1)(-\frac{f'}{f}) - n\frac{f''}{f}]}}$$

with
$$\frac{f'}{f} = \frac{(\mathcal{F}_{\mathcal{T},S}(\lambda))'}{\mathcal{F}_{\mathcal{T},S}(\lambda)}$$
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$$x_-^{(1)} \qquad x_0 \qquad x_+^{(1)} \qquad \cdots$$

It's clear how we use $\lambda_m \approx \hat{X}_- < x < \hat{X}_+ \approx \lambda_{m+1}$:

$$\begin{array}{cccc}
\lambda_m & \chi_0 & \lambda_{m+1} \\
\hline
 & \bigcirc & \bigcirc
\end{array}$$

Def (Laguerre's iteration)

If
$$mlt(\lambda_m) = mlt(\lambda_{m+1}) = 1$$
 then

 $x_+^{(k)} = L_+^k(x) = L_+(L_+(\dots(x_0)))$
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Proof We can prove that

$$\lambda_m \leftarrow \dots x_-^{(2)} < x_-^{(1)} < x_0 < x_+^{(1)} < x_+^{(2)} \cdots \rightarrow \lambda_{m+1}$$



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So Laguerre's method is secure.

We also have un important property:

Teo

If we choose* $\lambda_m < x_0$ s.t. $sign\Big(\frac{f'(x_0)}{f(x_0)}\Big) = sign(\lambda_m - x_0)$ then $\{x_-^{(k)}\}_{k=1,\dots}$ converges monotonically in asymptotically cubic rate to λ_m .

mon. cubic to λ_{m+1}



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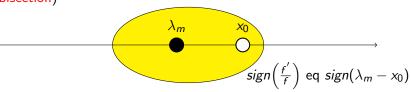
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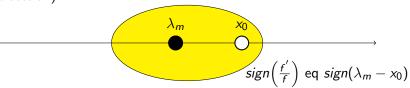
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The Laguerre's method is fast.

*for $x_0 < \lambda_{m+1}$ s.t. $sign\left(\frac{f'(x_0)}{f(x_0)}\right) = sign(\lambda_m - x_0)$ we have $\{x_+^{(k)}\}_{k=1,\dots}$ conv. mon. cubic to λ_{m+1}

It's clear that we need a powerfull method to obtain x_0 and an algorithm to estimate $mlt(\lambda_m)$.

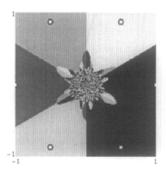
Overstimate $mlt(\lambda_m)$ (as we can read in [?]) causes no trouble, so the most important aspects of our calculation are:

Good x_0 .

Good evaluation of $L_{\pm}(x)$.

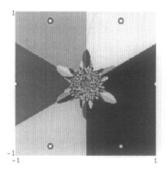
Complex case (little digression)

According to [?], we observe that with Laguerre's method searching $z \in \mathbb{C}$ such that $z^n - 1 = 0$ for n > 4 it's difficult because near the origin there is a Julia fractal set for starting point z_0 . (figure: n = 6)



Complex case (little digression)

According to [?], we observe that with Laguerre's method searching $z \in \mathbb{C}$ such that $z^n - 1 = 0$ for n > 4 it's difficult because near the origin there is a Julia fractal set for starting point z_0 . (figure: n = 6)



So if we want to solve the Generalized Eigenvalues Problem with $T, S \in \mathbb{C}$ we have great problems to place the starting point if n > 4.

Three-term recurrence

For a generic $x \in \mathbb{R}$ we call $\rho_n(x) = \det(T_n - xS_n)$, and $\rho_{n-1}(x) = \det(T_{n-1} - xS_{n-1})$, with T_{n-1} leading principal submatrix.

We have

$$\begin{split} \rho_0 &:= 1, \; \rho_1 := t_{1,1} - x s_{1,1} \\ \rho_i &:= (t_{i,i} - x s_{i,i}) \rho_{i-1} - (t_{i-1,i} - x s_{i-1,i})^2 \rho_{i-2}, \; i = 2, 3, \dots, n \end{split}$$

Three-term recurrence

We can proove it with the Laplace expansion (e.g. n = 4)

$$T_{4} - xS_{4} = \begin{pmatrix} a & b & 0 & 0 \\ b & c & d & 0 \\ 0 & d & e & f \\ 0 & 0 & f & g \end{pmatrix}$$

$$det (T_{4} - \lambda S_{4}) = g \begin{bmatrix} a & b & 0 \\ b & c & d \\ 0 & d & e \end{bmatrix} - f \begin{bmatrix} a & b & 0 \\ b & c & d \\ 0 & 0 & f \end{bmatrix}$$

$$= g \begin{bmatrix} a & b & 0 \\ b & c & d \\ 0 & d & e \end{bmatrix} - f^{2} \begin{bmatrix} a & b \\ b & c \end{bmatrix}$$

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Remark

If
$$x = \lambda$$
 then $\rho_n(\lambda) = \mathcal{F}_{(T_n, S_n)}(\lambda) = f(\lambda)$

Remark (important results with ${\color{red}\bullet}$ in Appendix)

 $\rho_0, \rho_1, \dots, \rho_n$ is a *Sturm sequence* of polynomials so, $\forall x \in \mathbb{R}$, we have

$$\kappa(x) := \#$$
 eigenvalues less then x
 $\kappa(x) = \#$ consecutive sign changes in $\{\rho_i\}_{i=0,\dots,n}$

Obviusly
$$f' = \rho'_n, f'' = \rho''_n$$
.



Obviusly
$$f^{'}=\rho_{n}^{'}, f^{''}=\rho_{n}^{''}.$$
 So*

IF
$$\left(\kappa(x_0) \geq m \text{ AND } -\frac{f'}{f} < 0\right)$$
 OR $\left(\kappa(x_0) < m \text{ AND } -\frac{f'}{f} \geq 0\right)$ THEN

Obviusly
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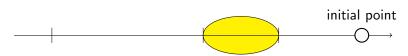
$$\mathsf{IF}\,\left(\kappa(\mathsf{x}_0) \geq m\;\mathsf{AND}\; -\frac{f'}{f} < 0\right)\;\mathsf{OR}\;\left(\kappa(\mathsf{x}_0) < m\;\mathsf{AND}\; -\frac{f'}{f} \geq 0\right)\;\mathsf{THEN}$$

We have cubic convergence in Laguerre's method with x_0 as starting point.



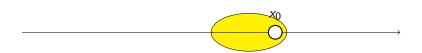






with bisection we can find the neighbourhood of λ_m

we can now use the Laguerre's iteration





We define
$$\xi_i = \frac{\rho_i}{\rho_{i-1}}$$
 for $i = 2, \dots, n$ and we have $\rho_i = \prod_{k=1}^i \xi_k$.

We define $\xi_i = \frac{\rho_i}{\rho_{i-1}}$ for $i=2,\ldots,n$ and we have $\rho_i = \prod_{k=1}^i \xi_k$. We also define $\eta_i = \frac{\rho_i'}{\rho_i}, \zeta_i = \frac{\rho_i''}{\rho_i}$ for $i=0,1,\ldots,n$ and finally we have

$$\kappa(x) = \text{ number of negative terms in } \{\xi_i\}_{i=1}^n$$

$$-\frac{f'(x)}{f(x)} = \eta_n$$

$$\frac{f''(x)}{f(x)} = \zeta_n$$

So we have to calculate three three-term recurrences.

^{*}we consider all elements of diag and super of (T,S) as non zero.

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* Total $\leq 2 + 38n$ multiplications.

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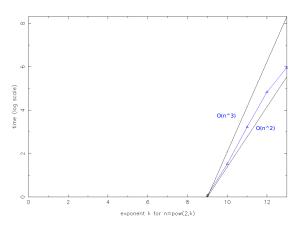
* Total $\leq 2 + 38n$ multiplications.

For every step in Laguerre's iteration we have to do 2+7+38n (at most) multiplications and 1 square root extraction. Because the convergence is cubic we hope in a small number of iteration.

^{*}we consider all elements of diag and super of (T,S) as non zero.

Complexity

*Author's prevision: $O(n^2)$. With (I, S) and eigenvalues between 0.5 and 1.5 we have

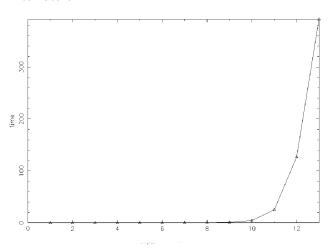




^{*}They use n < 100, my experiment ends up with $n = 2^{13} = 8192$.

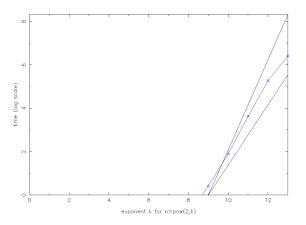
Complexity

And in linear scale:



Complexity

With $T = (a_i, 1, a_i)$, $S = (b_i, c_i, b_i)$ and a_i, b_i random numbers such that $|a_i| \le 10^{-3}$, $|b_i| \le 10^{-1}$, $|c_i| = i10^{-1} + |2b_i|$ we have



We can appreciate little differences between T = (0, 1, 0) and $T = (a_i, 1, a_i)$



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$$\xi_i = \frac{\rho_i}{\rho_{i-1}}$$
, it's self-scaled.

Teo

$$f[f(x)] = f[det[T - xS]] = (1 + \gamma)det[(T + \delta T) - x(S + \delta S)]$$

where $|\gamma| \leq n\epsilon$, with ϵ machine precision, and both δT and δS are symmetric tridiagonal matrices satisfying entrywise inequalities $|\delta T|_{\infty} \leq 2.51\epsilon |T|_{\infty} + \sqrt{\epsilon_u}, |\delta S|_{\infty} \leq 3.51\epsilon |S|_{\infty}$, where ϵ_u is the underflow threshold (in double precision is 10^{-308}).

We are interested not only in $\mathrm{fl}[f(x)]=\mathrm{fl}\Big[\prod_{i=1}^n\mathrm{fl}[\xi_i]\Big]$, but also in $\mathrm{fl}[\eta_n]$ and $\mathrm{fl}[\zeta_n]$, because we use these three value to calculate the Laguerre's iteration.

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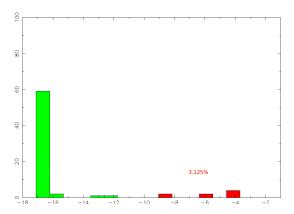
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The authors don't report this important aspect of analysis. (fl[η_n], fl[ζ_n] suffer from similar problems and have similar backward analysis, but they are -in partucular ζ_{n^-} enourmous in magnitude; sometimes this can cause troubles)

Exponent of absolute error (n = 256)

With (I, S) and eigenvalue between 0.5 and 1.5 we have



Sometimes (about 3 over 100 eigenvalues with $n \ge 256$) Laguerre's iteration breaks the monotonic convergence!

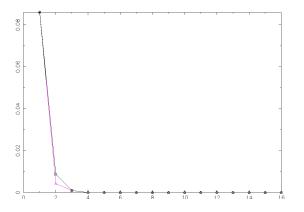
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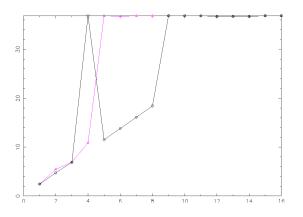
 $|\lambda_i - x_i| \leq 10^{-i}$ and $i = 1, \dots, 16$ we have

n = 256, dim = 64 and pencil=(T(193 : 4dim), S(193 : 4dim))



For $i=1,\ldots,16$, absolute errors between λ_{32} and x_i with κ , absolute errors between λ_{32} and x_i with inertia.

n = 256, dim = 64 and pencil=(T(193 : 4dim), S(193 : 4dim))



For $i=1,\ldots,16$, absolute errors between λ_{32} and x_i with κ , absolute errors between λ_{32} and x_i with inertia.

I use a different scale: $log(\frac{1}{|error|})$



When x_0 for Laguerre's iteration is into our good neighbourhood but $10^{-9} \le |x_0 - \lambda_m| \le 10^{-1}$, we can have troubles...

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Searching initial points

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Scalability.

We have: Laguerre's method.

Cuppen's divide and conquer method.

Symmetric tridiagonal matrices.

We add: Unreducible condition.

Dynamic programming (Bottom-up).

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Ideas:

If n = 1 we have $t \cdot v = \lambda \cdot s \cdot v$ with $s \neq 0$, so $\lambda = \frac{t}{s}$.

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 we have $t\cdot v=\lambda\cdot s\cdot v$ with $s\neq 0$, so $\lambda=\frac{t}{s}$. If $n=2$ we have
$$\begin{bmatrix} t_{1,1} & t_{1,2} \\ t_{2,1} & t_{2,2} \end{bmatrix} \begin{bmatrix} v_1 \\ v_2 \end{bmatrix} = \lambda \begin{bmatrix} s_{1,1} & s_{1,2} \\ s_{2,1} & s_{2,2} \end{bmatrix} \begin{bmatrix} v_1 \\ v_2 \end{bmatrix}$$

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$$S^{-1}T = \delta^{-1} \cdot \begin{pmatrix} \alpha & \gamma \\ \gamma & \beta \end{pmatrix}$$

we resolve $det[S^{-1}T - \lambda I] = 0$ with

$$\lambda_{1,2} = \frac{1}{2\delta} \left(\alpha + \beta \pm \sqrt{\alpha^2 + \beta^2 + \gamma^2 - \alpha\beta} \right)$$

 $^*\alpha = s_{2,2}t_{1,1} - s_{1,2}t_{1,2}, \beta = -s_{1,2}t_{1,2} + s_{1,1}t_{2,2}$ $\gamma = -s_{1,2}t_{1,2} + s_{1,1}t_{1,2}, \delta = s_{1,1}s_{2,2} - s_{1,2}^2.$



And if n = 4?

$$\begin{pmatrix} t_{1,1} & t_{1,2} & 0 & 0 \\ t_{2,1} & t_{2,2} & t_{2,3} & 0 \\ 0 & t_{3,2} & t_{3,3} & t_{3,4} \\ 0 & 0 & t_{4,3} & t_{4,4} \end{pmatrix} \begin{pmatrix} v_1 \\ v_2 \\ v_3 \\ v_4 \end{pmatrix} = \lambda \begin{pmatrix} s_{1,1} & s_{1,2} & 0 & 0 \\ s_{2,1} & s_{2,2} & s_{2,3} & 0 \\ 0 & s_{3,2} & s_{3,3} & s_{3,4} \\ 0 & 0 & s_{4,3} & s_{4,4} \end{pmatrix} \begin{pmatrix} v_1 \\ v_2 \\ v_3 \\ v_4 \end{pmatrix}$$

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We can solve the blue part and obtain μ_1, μ_2 , the cyan part and obtain μ_3, μ_4 .

From μ_j we can enter in our neighbourhood of λ_j using simple bisection:

^{*}This is an important part of our method.

From μ_i we can enter in our neighbourhood of λ_i using simple bisection:

Pseudocode:

```
set a_i = 0 and b_i = \mu_i
set x = \mu_i
# IF \kappa(x) < i
THEN set a_i = x
ELSE set b_i = x
*IF -\frac{f'(x)}{f(x)} = sign(\lambda_j - x)
THEN stop
ELSE set x = \frac{b_j - a_j}{2} and go to #
set x_0 = x
```

Authors of [?, ?] doesn't explain this point.

Loode

Loode



^{*}This is an important part of our method.

$$x_0^1, L(x_0^1), L(L(x_0^1)), \cdots \rightsquigarrow \lambda_1$$

$$x_0^2, L(x_0^2), L(L(x_0^2)), \cdots \rightsquigarrow \lambda_2$$
Hopefully we have
$$x_0^3, L(x_0^3), L(L(x_0^3)), \cdots \rightsquigarrow \lambda_3$$

$$x_0^4, L(x_0^4), L(L(x_0^4)), \cdots \rightsquigarrow \lambda_4$$

Consider (\hat{T}, \hat{S}) with

$$\hat{T} = \begin{pmatrix} T_0 & 0 \\ 0 & T_1 \end{pmatrix}$$

$$\hat{S} = \begin{pmatrix} S_0 & 0 \\ 0 & S_1 \end{pmatrix}$$

with T_0, T_1, S_0, S_1 symmetric tridiagonal, and let be

$$\hat{\lambda}_1 \leq \hat{\lambda}_2 \leq \dots \leq \hat{\lambda}_n$$

eigenvalues of (\hat{T}, \hat{S}) , they are what we call μ_1, \ldots, μ_n .

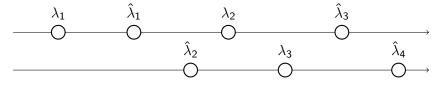
Teo (A sort of "interlacing" with ▶ proof)

$$-\infty < \lambda_1 \le \hat{\lambda}_1$$
$$\hat{\lambda}_{i-1} \le \lambda_i \le \hat{\lambda}_{i+1}$$
$$\hat{\lambda}_n \le \lambda_n < \infty$$

with i = 2, 3, ..., n - 1.

Remark

It's possible that



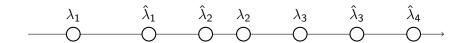
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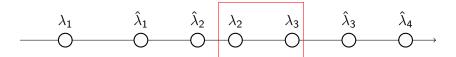
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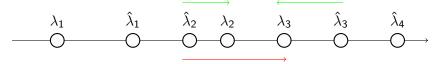
Remark

It's possible that



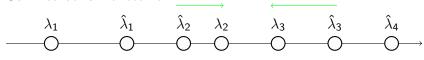
Remark

Our method is monotonic



Remark

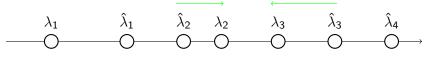
Our method is monotonic



We can reach λ_2 only from $\hat{\lambda}_2$, moving from left to right (and similar λ_3 only from $\hat{\lambda}_3$, moving from right to left).

Remark

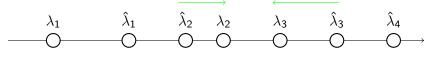
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We can reach λ_2 only from $\hat{\lambda}_2$, moving from left to right (and similar λ_3 only from $\hat{\lambda}_3$, moving from right to left). If $|\lambda_2 - \lambda_3| \leq 10^{-14}$ we can have troubles.

Remark

Our method is monotonic



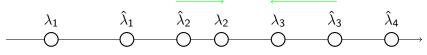
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If $|\lambda_2 - \lambda_3| \le 10^{-14}$ we can have troubles.

Luckily we also have a *multiplicity estimator* (called EstMlt) that works only with $\hat{\lambda}_1, \dots, \hat{\lambda}_n$.

Remark

Our method is monotonic



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If $|\lambda_2 - \lambda_3| \le 10^{-14}$ we can have troubles.

Luckily we also have a *multiplicity estimator* (called EstMlt) that works only with $\hat{\lambda}_1, \dots, \hat{\lambda}_n$.

If mlt=2 then we consider $\lambda_2=\lambda_3$, i.e. we said that λ_2 has multiplicity 2.

EstMIt

We have j, x, $\mathrm{sgn}\Big(-\frac{f'(x)}{f(x)}\Big)$ and $\hat{\lambda}_1, \dots, \hat{\lambda}_n$ as *INPUT*.

Pseudocode:

```
mlt = 1
 do k = 1, ...

m = j + k \operatorname{sgn}\left(-\frac{f'(x)}{f(x)}\right)
 if m \le 0 then
 something goes wrong
  set mlt = 1
  go to #
  end if
 if |\hat{\lambda}_j - \hat{\lambda}_m| \leq 0.01 |\hat{\lambda}_j - x| then mlt = mlt + 1
 else go to #
  end if
```

We want: Fast and secure iterative method.

Starting points for our method.

Scalability.

We have: Laguerre's method.

Cuppen's divide and conquer method.

Symmetric tridiagonal matrices.

We add: Unreducible condition.

Dynamic programming (Bottom-up).

Efficient matrix storing.

The author's point of view is parallel computating.



My point of view is sequential computating.





Instead of recurtion I chose dynamic programming, *i.e.* "solving complex problems by breaking them down into simpler subproblems. [...] In general, to solve a given problem, we need to solve different parts of the problem (subproblems), then combine the solutions of the subproblems to reach an overall solution." (cit. Wikipedia)

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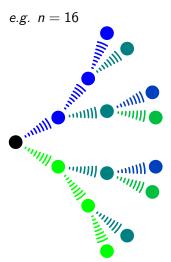
We can

- ▶ compute eigenvalues for all the pencils 2×2 (they are 2^{k-1})
- ▶ compute eigenvalues for all the pencils 4×4 (they are 2^{k-2})
- •
- ▶ compute eigenvalues for all the pencils $\frac{n}{2} \times \frac{n}{2}$ (they are 2)
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- ▶ compute eigenvalues for all the pencils $\frac{n}{2} \times \frac{n}{2}$ (they are 2)
- \triangleright compute eigenvalues of (T, S)

This is a "Bottom-up approach" because we start with dimension 2 and finish with dimension n.



We start from leaves (dim=2) and finish with the root (dim=16)

Code:

```
WHILE dim < n
DO k = 0, \frac{n}{dim} - 1
set Tstart = 1 + k\dim \text{ and } Tend = (1 + k)\dim
set Sstart = 1 + k\dim \text{ and } Send = (1 + k)\dim
IF dim = 2
THEN
explicit calculation
ELSE
search eigenvalues of
(T(Tstart...Tend, 0...1), S(Sstart...Send, 0...1))
knowing eigenvalues of
(T(Tstart \dots \frac{Tend}{2}, 0 \dots 1), S(Sstart \dots \frac{Send}{2}, 0 \dots 1))
and
\left(T\left(\frac{Tend}{2}+1\dots Tend,0\dots 1\right),S\left(\frac{Send}{2}+1,0\dots 1\right)\right)
END DO
```

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...We can search only eigenvalues in one particular interval of real line.

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e.g. we are interested in $[0,10^{-16}]$ or $[10^3,10^8]$ or we want to use Gershgorin theorems to isolate particolar eigenvalues.

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e.g. we are interested in $[0,10^{-16}]$ or $[10^3,10^8]$ or we want to use Gershgorin theorems to isolate particolar eigenvalues.

...If also T is *definite* we can search eigenvalues in $[1,\infty]$ using our method with (S,T) and $\lambda_j=\frac{1}{\mu_i}$.

"Real world" matrices

Consider *piecewise linear finite element discretization of* the Sturm-Liouville problem

$$-\frac{d}{dx}\left(p(x)\frac{du}{dx}\right)+q(x)u=\lambda u$$

where $u=u(x), 0< x<\pi$ and $u(0)=u^{'}(\pi)=0$ and p(x)=1, q(x)=6 ($h=\frac{1}{n+1}$). We obtain*

$$t_{k,k} = -6h(2k^2 - 2k - 1)$$
 $t_{k,k+1} = -\frac{1}{h} + h$
 $s_{k,k} = \frac{4h}{6}$
 $s_{k,k+1} = \frac{h}{6}$



^{*}authors doesn't report (T, S)

"Real world" matrices

Thank you.

Appendix A: proof of Laguerre's convergence

According to [?, p.444] we can write

$$x_{\pm}^{(k+1)} = x_{\pm}^{(k)} - \frac{nf}{f' \pm H^{\frac{1}{2}}}$$

$$H = (n-1)^{2} (f')^{2} - n(n-1)ff''$$

If we choose the sign so that the $|f'\pm H^{\frac{1}{2}}|$ has the larger absolute value then we can approx $x_+^{(k+1)}-\lambda_m$ as

$$x_{\pm}^{(k+1)} - \lambda_m \approx \frac{1}{2} (x_{\pm}^{(k)} - \lambda_m)^3 \frac{(n-1)\Sigma_2' - (\Sigma_1')^2}{n-1}$$

$$\Sigma_2' = \sum_{i \neq n} \frac{1}{(\lambda_m - \lambda_i)^2}$$

$$\Sigma_1' = \sum_{i \neq n} \frac{1}{\lambda_m - \lambda_i}$$

Appendix A: proof of Laguerre's convergence

So we have convergence and $x_{\pm}^{(k+1)} - \lambda_m \approx \text{number } (x_{\pm}^{(k)} - \lambda_m)^3$ tell us that the convergence is cubic.



Appendix B: proof of property about Sturm sequence



Def

For $\alpha \in [0,1]$ we define the pencil $(T(\alpha), S(\alpha)) := ((1-\alpha)\hat{T} + \alpha T, (1-\alpha)\hat{S} + \alpha S).$

Lemm

 $(T(\alpha), S(\alpha))$ is a symmetric definite pencil for each $\alpha \in [0, 1]$. Calling $\lambda_1(\alpha) \leq \lambda_2(\alpha) \leq \cdots \leq \lambda_n(\alpha)$ the n real eigenvalues of the pencil $(T(\alpha), S(\alpha))$ we have

Lemm

Each $\lambda_i(\alpha)$ is a continuous function of $\alpha \in [0,1]$.

Teo (A sort of "interlacing")

$$-\infty < \lambda_1 \le \hat{\lambda}_1$$
$$\hat{\lambda}_{i-1} \le \lambda_i \le \hat{\lambda}_{i+1}$$
$$\hat{\lambda}_n \le \lambda_n < \infty$$

with
$$i = 2, 3, ..., n - 1$$
.

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Classic interlacing still works for $-\infty < \lambda_1 \le \hat{\lambda}_1$ and for $\hat{\lambda}_n \le \lambda_n < \infty$.

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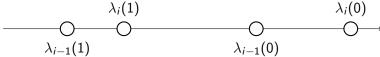
with i = 2, 3, ..., n - 1.

Classic interlacing still works for $-\infty < \lambda_1 \le \hat{\lambda}_1$ and for $\hat{\lambda}_n < \lambda_n < \infty$.

We have to proove that $\lambda_i \geq \hat{\lambda}_{i-1}$ (and similar $\lambda_i \leq \hat{\lambda}_{i+1}$).

```
(Proof by contradiction) if we consider \lambda_i < \hat{\lambda}_{i-1} for some i \in \{2, 3, ..., n-1\} (that, in our new writing, is \lambda_i(1) < \lambda_{i-1}(0) because all \lambda_j(1) are eigenvalues of (T, S) and all \lambda_j(0) are eigenvalues of (\hat{T}, \hat{S})) then
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$$\begin{array}{c|c} \lambda_i(1) & \lambda_i(0) \\ \hline \lambda_{i-1}(1) & \lambda_{i-1}(0) \end{array}$$

in symbols: $\lambda_{i-1}(1) \leq \lambda_i(1) < \lambda_{i-1}(0) \leq \lambda_i(0)$.

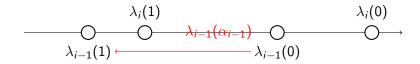
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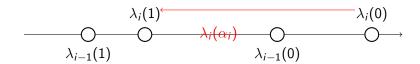
$$\begin{array}{c|c} \lambda_i(1) & \lambda_i(0) \\ \hline \lambda_{i-1}(1) & \lambda_{i-1}(0) \end{array}$$

in symbols: $\lambda_{i-1}(1) \leq \lambda_i(1) < \lambda_{i-1}(0) \leq \lambda_i(0)$. For all $\tilde{\lambda} \in [\lambda_i(1), \lambda_{i-1}(0)]$ we can find α_i, α_{i-1} such as $\tilde{\lambda} = \lambda_i(\alpha_i) = \lambda_{i-1}(\alpha_{i-1})$.









$$H(\alpha, \lambda) := det[T(\alpha) - \lambda S(\alpha)] =$$

$$\begin{bmatrix} \ddots & \ddots & & \\ \ddots & t_{k,k} - \lambda s_{k,k} & \alpha(t_{k,k+1} - s_{k,k+1}) \\ & \alpha(t_{k,k+1} - s_{k,k+1}) & t_{k+1,k+1} - \lambda s_{k+1,k+1} & \ddots \\ & & \ddots & & \ddots \end{bmatrix}$$

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Sc

$$H(\alpha, \lambda) = p(\lambda) + \alpha^2 q(\lambda)$$

with p, q polynomials.

If $q(\tilde{\lambda}) \neq 0$ then $\tilde{\alpha} = +\sqrt{H(\tilde{\alpha}, \tilde{\lambda}) - \frac{p(\tilde{\lambda})}{q(\tilde{\lambda})}} \in [0, 1]$. only one solution in α .

If $q(\tilde{\lambda}) \neq 0$ then $\tilde{\alpha} = +\sqrt{H(\tilde{\alpha}, \tilde{\lambda}) - \frac{p(\tilde{\lambda})}{q(\tilde{\lambda})}} \in [0, 1]$. only one solution in α . But we have, for all $\tilde{\lambda}$, two solution in α , named α_i and α_{i-1} .

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Then $q(\lambda) = 0$ for all $\tilde{\lambda}$.

We have $\tilde{\lambda}$ eigenvalue of (\hat{T}, \hat{S}) , for all $\tilde{\lambda}$ and (\hat{T}, \hat{S}) have exactly n eigenvalues.

If
$$q(\tilde{\lambda}) \neq 0$$
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conctraddiction

▶ back

Appendix D: code

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Grazie (ancora) per l'attenzione.