# Optimization of Polynomial Roots, Eigenvalues and Pseudospectra 

Michael L. Overton<br>Courant Institute of Mathematical Sciences NYU<br>Banff Stability Workshop<br>Nov 5, 2012

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Globally Optimizing
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Affine Constraint
with
V. Blondel (Louvain)
M. Gürbüzbalaban
(NYU)
A. Megretski (MIT)
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and the Root
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\section*{The Root Radius and the Root Abscissa}

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Let \(\rho\) denote the root radius of a polynomial:
\[
\rho(p)=\max \{|z|: p(z)=0, z \in \mathbf{C}\} .
\]

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Stabilization

Let \(\rho\) denote the root radius of a polynomial:
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We say \(p\) is Schur stable if \(\rho(p)<1\).

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As functions of the polynomial coefficients, the radius \(\rho\) and abscissa \(\alpha\) are

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As functions of the polynomial coefficients, the radius \(\rho\) and abscissa \(\alpha\) are
not convex

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As functions of the polynomial coefficients, the radius \(\rho\) and abscissa \(\alpha\) are

■ not convex
■ not Lipschitz near polynomials with a multiple root

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As functions of the polynomial coefficients, the radius \(\rho\) and abscissa \(\alpha\) are
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■ not Lipschitz near polynomials with a multiple root So, in general, global minimization of the radius or abscissa over an affine family of monic polynomials, pushing the roots as far as possible towards the origin or left in the complex plane, seems hard.

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Indeed, variations on the question of whether a polynomial family contains one that is stable (has roots inside the unit circle or in the left-half plane) have been studied for decades.

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Indeed, variations on the question of whether a polynomial family contains one that is stable (has roots inside the unit circle or in the left-half plane) have been studied for decades.
But if an affine family of monic polynomials of degree \(n\) has \(n-1\) free parameters, this question can be answered efficiently.

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\(\square\) not convex
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Indeed, variations on the question of whether a polynomial family contains one that is stable (has roots inside the unit circle or in the left-half plane) have been studied for decades.
But if an affine family of monic polynomials of degree \(n\) has \(n-1\) free parameters, this question can be answered efficiently.
Equivalently, there is just one affine constraint on the coefficients.

\section*{Optimizing the Root Radius, Real Case}

Theorem RRR. Let \(B_{0}, B_{1}, \ldots, B_{n}\) be real scalars (with

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Theorem RRR. Let \(B_{0}, B_{1}, \ldots, B_{n}\) be real scalars (with \(B_{1}, \ldots, B_{n}\) not all zero) and consider the affine family
\(P=\left\{z^{n}+a_{1} z^{n-1}+\ldots+a_{n-1} z+a_{n}: B_{0}+\sum_{j=1}^{n} B_{j} a_{j}=0, a_{i} \in \mathbf{R}\right\}\).
The optimization problem
\[
\rho^{*}:=\inf _{p \in P} \rho(p)
\]
has a globally optimal solution of the form
\[
p^{*}(z)=(z-\gamma)^{n-k}(z+\gamma)^{k} \in P
\]
for some integer \(k\) with \(0 \leq k \leq n\), where \(\gamma=\rho^{*}\).
Proof: uses implicit function theorem.
Algorithm: for each \(k=0, \ldots, n\), substitute coefficients of \((z-\gamma)^{n-k}(z+\gamma)^{k}\) into the constraint to give a polynomial with \(n\) roots that are candidates for \(\gamma\). Choose smallest such \(|\gamma|\).

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\section*{Case}

Theorem RRC. Let \(B_{0}, B_{1}, \ldots, B_{n}\) be complex scalars (with
\(B_{1}, \ldots, B_{n}\) not all zero) and consider the affine family
\(P=\left\{z^{n}+a_{1} z^{n-1}+\ldots+a_{n-1} z+a_{n}: B_{0}+\sum_{j=1}^{n} B_{j} a_{j}=0, a_{i} \in \mathbf{C}\right\}\).
The optimization problem
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\rho^{*}:=\inf _{p \in P} \rho(p)
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has an optimal solution of the form
\[
p^{*}(z)=(z-\gamma)^{n} \in P
\]
with \(-\gamma\) given by a root of smallest magnitude of the polynomial
\[
h(z)=B_{n} z^{n}+B_{n-1}\binom{n}{n-1} z^{n-1}+\ldots+B_{1}\binom{n}{1} z+B_{0}
\]

\section*{Optimizing the Root Radius: Complex Case}

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Theorem RRC. Let \(B_{0}, B_{1}, \ldots, B_{n}\) be complex scalars (with
\(B_{1}, \ldots, B_{n}\) not all zero) and consider the affine family
\(P=\left\{z^{n}+a_{1} z^{n-1}+\ldots+a_{n-1} z+a_{n}: B_{0}+\sum_{j=1}^{n} B_{j} a_{j}=0, a_{i} \in \mathbf{C}\right\}\).
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h(z)=B_{n} z^{n}+B_{n-1}\binom{n}{n-1} z^{n-1}+\ldots+B_{1}\binom{n}{1} z+B_{0}
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Proof: A complicated inductive argument.

\section*{Optimizing the Root Abscissa: Real Case}

Theorem RAR. Let \(B_{0}, B_{1}, \ldots, B_{n}\) be real scalars (with

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\section*{Root Abscissa, Real Case, Continued}

Furthermore, the optimal value \(\alpha^{*}\) is attained by a minimizing
polynomial \(p^{*}\) if and only if \(-\alpha^{*}\) is a root of \(h\) (as opposed to one of its derivatives), and in this case we can take
\[
p^{*}(z)=(z-\gamma)^{n} \in P
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with \(\gamma=\alpha^{*}\).
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Furthermore, the optimal value \(\alpha^{*}\) is attained by a minimizing polynomial \(p^{*}\) if and only if \(-\alpha^{*}\) is a root of \(h\) (as opposed to one of its derivatives), and in this case we can take
\[
p^{*}(z)=(z-\gamma)^{n} \in P
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with \(\gamma=\alpha^{*}\).
When the optimal abscissa is not attained, for all \(\epsilon>0\) can find an approximately optimal polynomial
\[
p_{\epsilon}(z):=\left(z-M_{\epsilon}\right)^{\ell}\left(z-\left(\alpha^{*}+\epsilon\right)\right)^{n-\ell} \in P
\]
with \(0<\ell \leq n\) and \(M_{\epsilon} \rightarrow-\infty\) as \(\epsilon \rightarrow 0\).

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Thus, as in the real radius case, two roots play a role, but only one is finite.

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with \(0<\ell \leq n\) and \(M_{\epsilon} \rightarrow-\infty\) as \(\epsilon \rightarrow 0\).
Thus, as in the real radius case, two roots play a role, but only one is finite.

In practice, bad idea to make \(\epsilon\) too small: then \(\left|M_{\epsilon}\right|\) becomes large.

\section*{Optimizing the Abscissa: Real vs. Complex Case}

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We observed that, in the real case, the optimal value is not attained when one of the derivatives of \(h\) has a real root to the right of the rightmost real root of \(h\).

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We observed that, in the real case, the optimal value is not attained when one of the derivatives of \(h\) has a real root to the right of the rightmost real root of \(h\).

However, it is not possible that a derivative of \(h\) has a complex root to the right of the rightmost complex root of \(h\). This follows immediately from the Gauss-Lucas theorem.

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\section*{Optimizing the}

Abscissa: Real vs. Complex Case

We observed that, in the real case, the optimal value is not attained when one of the derivatives of \(h\) has a real root to the right of the rightmost real root of \(h\).

However, it is not possible that a derivative of \(h\) has a complex root to the right of the rightmost complex root of \(h\). This follows immediately from the Gauss-Lucas theorem.

This suggests the optimal abscissa value might always be attained in the complex case.

\section*{Optimizing the Abscissa: Real vs. Complex Case}

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Indeed, this is the case...

\section*{Optimizing the Root Abscissa: Complex Case}

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\author{
Optimizing the Root
}

Abscissa: Complex

\section*{Case}

Theorem RAC. Let \(B_{0}, B_{1}, \ldots, B_{n}\) be complex scalars (with \(B_{1}, \ldots, B_{n}\) not all zero) and consider the affine family
\[
P=\left\{z^{n}+a_{1} z^{n-1}+\ldots+a_{n-1} z+a_{n}: B_{0}+\sum_{j=1}^{n} B_{j} a_{j}=0, a_{i} \in \mathbf{C}\right\}
\]

\section*{Optimizing the Root Abscissa: Complex Case}

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The optimization problem
\[
\alpha^{*}:=\inf _{p \in P} \alpha(p)
\]
has an optimal solution of the form
\[
p^{*}(z)=(z-\gamma)^{n} \in P
\]
with \(-\gamma\) given by a root with largest real part of the polynomial
\[
h(z)=B_{n} z^{n}+B_{n-1}\binom{n}{n-1} z^{n-1}+\ldots+B_{1}\binom{n}{1} z+B_{0}
\]

\section*{Example: Static Output Feedback Stabilization}

Given the dynamical system with input and output:

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\section*{Case}

Given the dynamical system with input and output:
\[
\dot{x}=F x+G u, \quad y=H x
\]
where \(F \in \mathbf{R}^{n \times n}, G \in \mathbf{R}^{n \times \ell}, H \in \mathbf{R}^{m \times n}\).
SOF: find a controller \(K \in \mathbf{R}^{\ell \times m}\) so that, setting \(u=K y\)
\[
\dot{x}=(F+G K H) x
\]
is stable, that is all eigenvalues of \(F+G K H\) are in the left half-plane, or prove that this is not possible.

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A major open problem in control.

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is stable, that is all eigenvalues of \(F+G K H\) are in the left half-plane, or prove that this is not possible.
A major open problem in control.
But, if \(p=1\) and \(m=n-1\) (one input and \(n-1\) outputs)
\[
\operatorname{det}(\lambda I-F-G K H)=\operatorname{det}(\lambda I-F)+K H \operatorname{adj}(\lambda I-F) G .
\]

This is a monic polynomial with affine dependence on the \(n-1\) entries of \(K \in \mathbf{R}^{1 \times(n-1)}\) so the SOF problem can be solved explicitly using Theorem RAR.

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Stabilization

Another set of classical problems in control that, in a certain case, can be solved using the theorems given above.

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Another set of classical problems in control that, in a certain case, can be solved using the theorems given above.

An example: stabilizing the two-mass-spring dynamical system by a second-order controller.

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\section*{Case} case, can be solved using the theorems given above.

An example: stabilizing the two-mass-spring dynamical system by a second-order controller.

Then, maximizing the closed-loop asymptotic decay rate is equivalent to solving the optimization problem
\[
\min _{n \subset D} \max _{\sim \subset C}\{\operatorname{Re} z: p(z)=0\}
\]
where
\(P=\left\{\left(z^{4}+2 z^{2}\right)\left(x_{0}+x_{1} z+z^{2}\right)+y_{0}+y_{1} z+y_{2} z^{2}: x_{0}, x_{1}, y_{0}, y_{1}, y_{2} \in \mathbf{R}\right\}\)

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We can minimize the root abscissa explicitly using Theorem RAR as \(P\) is a set of monic polynomials with degree 6 whose coefficients depend affinely on 5 real parameters.

\section*{Caveats}

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Multiple roots are very sensitive to perturbation!

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A random perturbation of size \(\epsilon\) to the coefficients of a polynomial with a root that has multiplicity \(k\) moves the roots by \(O\left(\epsilon^{1 / k}\right)\).

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In practice, might want to locally optimize a more robust measure of stability: see Part III.

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Independently of this, the monomial basis is a poor choice numerically unless the polynomial has very small degree.

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Nonetheless, the optimal value can be computed accurately even if \(n\) is fairly large.

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Independently of this, the monomial basis is a poor choice numerically unless the polynomial has very small degree.

Nonetheless, the optimal value can be computed accurately even if \(n\) is fairly large.

Affpolymin: publicly available Matlab code implementing the algorithms implicit in Theorems RRR, RRC, RAR, RAC.

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The control applications are due to Chen, Rantzer and Henrion.
A publicly available Matlab code implementing the constructive algorithms implicit in the theorems:
www.cs.nyu.edu/overton/software/affpoly/

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\section*{The Spectral Radius and the Spectral Abscissa}

Now let \(\rho: \mathbf{C}^{n \times n} \rightarrow \mathbf{R}\) denote spectral radius:
\[
\rho(A)=\max \{|z|: \operatorname{det}(A-z I)=0, z \in \mathbf{C}\} .
\]

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An eigenvalue is active if its modulus (real part) equals the spectral radius (abscissa).

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An eigenvalue is active if its modulus (real part) equals the spectral radius (abscissa).
\(\rho(A)<1\) is the stability condition for the discrete-time dynamical system \(\xi_{k+1}=A \xi_{k}\).
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Part II
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The Spectral Radius
and the Spectral
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The Reduced
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Eigenvalues of the
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\section*{The Spectral Radius and the Spectral Abscissa}

Now let \(\rho: \mathbf{C}^{n \times n} \rightarrow \mathbf{R}\) denote spectral radius:
\[
\rho(A)=\max \{|z|: \operatorname{det}(A-z I)=0, z \in \mathbf{C}\} .
\]

Let \(\alpha: \mathbf{C}^{n \times n} \rightarrow \mathbf{R}\) denote spectral abscissa:
\[
\alpha(A)=\max \{\operatorname{Re}(z): \operatorname{det}(A-z I)=0, z \in \mathbf{C}\} .
\]

An eigenvalue is active if its modulus (real part) equals the spectral radius (abscissa).
\(\rho(A)<1\) is the stability condition for the discrete-time dynamical system \(\xi_{k+1}=A \xi_{k}\).
\(\alpha(A)<0\) is the stability condition for the continuous-time dynamical system \(\dot{\xi}=A \xi\).

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The spectral functions \(\rho\) and \(\alpha\) are not convex and are not Lipschitz near a matrix with an active multiple eigenvalue.

\section*{No Extension of Part I}

\section*{Part I}

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For example, consider the matrix family
\[
A(x)=\left[\begin{array}{cc}
x & 1 \\
-1 & x
\end{array}\right] .
\]

This matrix depends affinely on a single parameter \(x\), but its characteristic polynomial, a monic polynomial of degree 2 , does not, so the results of Part I do not apply.

The minimal spectral radius of \(A(x)\) is attained by \(x=0\), for which the eigenvalues are \(\pm \mathbf{i}\).

\section*{The Diaconis-Holmes-Neal Sampler}

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A nonreversable Markov chain for Monte Carlo simulation. For \(x \in(0,1)\), the transition matrix is \(A(x) \in \mathbf{R}^{2 n \times 2 n}\) is


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Diaconis et. al. showed that for \(x=1 / n\), the corresponding nonreversible chain reaches a stationary state in \(O(n)\) steps, compared to \(O\left(n^{2}\right)\) steps for a similar reversible chain.

\section*{The Reduced Spectral Radius}

\section*{Part I}

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Spectral Radius
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Transition Matrix,
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The rate of convergence is determined by
\[
\tilde{\rho}(A(x))=\max \{|z|: \operatorname{det}(A(x)-z I)=0, z \in \mathbf{C}, z \neq 1\} .
\]

It is easy to prove that this is minimized over \(x \in[0,1]\) by
\[
x_{\mathrm{opt}}=\frac{\sin (\pi / n)}{1+\sin (\pi / n)}>\frac{1}{n}
\]

\section*{The Reduced Spectral Radius}

\section*{Part I}

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\[
x_{\mathrm{opt}}=\frac{\sin (\pi / n)}{1+\sin (\pi / n)}>\frac{1}{n}
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For \(x<x_{\mathrm{opt}}\), the active eigenvalues (the ones with largest modulus excluding 1) all occur in conjugate pairs.

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\]

For \(x<x_{\mathrm{opt}}\), the active eigenvalues (the ones with largest modulus excluding 1) all occur in conjugate pairs.

For \(x=x_{\mathrm{opt}}\), one conjugate pair has coalesced to a double real eigenvalue (corresponding to a \(2 \times 2\) Jordan block).

\section*{The Reduced Spectral Radius}

\section*{Part I}

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■ Blue: eigenvalues when \(x=1 / n\) (all complex)

Eigenvalues of the Transition Matrix, \(n=10\)

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■ Blue: eigenvalues when \(x=1 / n\) (all complex)
- Red: eigenvalues when \(x=x_{\mathrm{opt}}\) (one double real eigenvalue)

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■ Blue: eigenvalues when \(x=1 / n\) (all complex)
■ Red: eigenvalues when \(x=x_{\mathrm{opt}}\) (one double real eigenvalue)
■ Black: eigenvalues when \(x>x_{\text {opt }}\) ( \(\tilde{\rho}\) increases rapidly)

\section*{Reduced Spectral Radius as a Function of \(x\)}

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Note the big improvement changing \(x\) from \(1 / n\) to \(x_{\text {opt }}\).

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Note the big improvement changing \(x\) from \(1 / n\) to \(x_{\text {opt }}\).
Much better to underestimate \(x_{\mathrm{opt}}\) than overestimate. Similar plots apply to optimal damping for one-dimensional wave equation, optimal choice of parameter for SOR (successive over-relaxation), etc etc.

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Much better to underestimate \(x_{\mathrm{opt}}\) than overestimate. Similar plots apply to optimal damping for one-dimensional wave equation, optimal choice of parameter for SOR (successive over-relaxation), etc etc.
Convergence rate deteriorates as \(n\) increases.

\section*{Adding More Parameters}

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Not surprising that with one free parameter, we can only make one pair of eigenvalues coalesce.
Let's change \(A(x)\) to have multiple parameters:

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\section*{Still doubly stochastic. Can we now reduce \(\tilde{\rho}\) further?}

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Let's change \(A(x)\) to have multiple parameters:


Still doubly stochastic. Can we now reduce \(\tilde{\rho}\) further?
No! It appears that \(\mathbf{x}_{\mathrm{opt}}=\left[x_{\mathrm{opt}}, \ldots, x_{\mathrm{opt}}\right]^{T}\) is locally optimal.

\section*{Checking Local Optimality}

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Numerically: by running an optimization method suitable for nonsmooth objectives at randomly generated points near \(\mathbf{x}_{\text {opt }}\). We repeatedly obtained convergence to \(\mathbf{x}_{\mathrm{opt}}\).

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■ \(\mathrm{x}_{\text {opt }}\) satisfies a necessary condition for local optimality
■ if we remove some redundancy by setting \(x_{j}=x_{n-1-j}\) for \(j=1,2, \ldots,\left\lfloor\frac{n-1}{2}\right\rfloor\) and \(x_{n-1}=x_{n}\), we find \(\mathbf{x}_{\text {opt }}\) satisfies a sufficient condition for local optimality.

\section*{Checking Local Optimality}

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Special analysis needed because optimization objective is not Lipschitz at a matrix with an active multiple eigenvalue.

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■ if we remove some redundancy by setting \(x_{j}=x_{n-1-j}\) for \(j=1,2, \ldots,\left\lfloor\frac{n-1}{2}\right\rfloor\) and \(x_{n-1}=x_{n}\), we find \(\mathbf{x}_{\text {opt }}\) satisfies a sufficient condition for local optimality.

Special analysis needed because optimization objective is not Lipschitz at a matrix with an active multiple eigenvalue.
Essential to the analysis: each active eigenvalue corresponds to a single Jordan block, in this case with sizes \(2,1, \ldots, 1\).

\section*{Checking Local Optimality}

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\section*{Transition Matrix,}

Numerically: by running an optimization method suitable for nonsmooth objectives at randomly generated points near \(\mathbf{x}_{\text {opt }}\). We repeatedly obtained convergence to \(\mathrm{x}_{\mathrm{opt}}\).

Theoretically: by variational analysis. We found that
■ \(\mathrm{x}_{\text {opt }}\) satisfies a necessary condition for local optimality
■ if we remove some redundancy by setting \(x_{j}=x_{n-1-j}\) for \(j=1,2, \ldots,\left\lfloor\frac{n-1}{2}\right\rfloor\) and \(x_{n-1}=x_{n}\), we find \(\mathbf{x}_{\text {opt }}\) satisfies a sufficient condition for local optimality.

Special analysis needed because optimization objective is not Lipschitz at a matrix with an active multiple eigenvalue.
Essential to the analysis: each active eigenvalue corresponds to a single Jordan block, in this case with sizes \(2,1, \ldots, 1\).
Too complicated to explain in talk, but see references for more information.

\section*{Surface Approximation By Subdivision}

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An example from surface approximation by subdivision: several fixed eigenvalues, want to reduce modulus of others to optimize the smoothness of the surface: after much numerical computation, found that all can be reduced nearly to zero

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■ triangular mesh case: optimal multiple zero eigenvalue verified analytically, with multiple Jordan blocks of order 2, 1, 1, 1.

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■ triangular mesh case: optimal multiple zero eigenvalue verified analytically, with multiple Jordan blocks of order 2, 1, 1, 1.
- quadrilateral mesh case: numerically reduced moduli of eigenvalues to about \(10^{-4}\) and estimated that the apparently optimal multiple zero eigenvalue has multiple Jordan blocks of order 5, 3, 2, 2.

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An example from surface approximation by subdivision: several fixed eigenvalues, want to reduce modulus of others to optimize the smoothness of the surface: after much numerical computation, found that all can be reduced nearly to zero

■ triangular mesh case: optimal multiple zero eigenvalue verified analytically, with multiple Jordan blocks of order 2, 1, 1, 1.
- quadrilateral mesh case: numerically reduced moduli of eigenvalues to about \(10^{-4}\) and estimated that the apparently optimal multiple zero eigenvalue has multiple Jordan blocks of order 5, 3, 2, 2.
In both cases, the active eigenvalue zero has not only algebraic multiplicity \(>1\) but also geometric multiplicity \(>1\). The latter results from special structure and will not occur generically.

\section*{Numerical Optimization of Nonsmooth, Nonconvex \(f\)}

Ordinary gradient method with line search: fails, typically

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\section*{Numerical Optimization of Nonsmooth, Nonconvex \(f\)}

Ordinary gradient method with line search: fails, typically
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BFGS quasi-Newton method with line search: empirically same property with much less computation.
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When using these methods to minimize the nonsmooth, nonconvex, non-Lipschitz functions \(\rho(A(x))\) or \(\alpha(A(x))\), make no attempt to predict active eigenvalues or estimate their multiplicities; just use gradients which exist at almost every \(x\)
\[
\frac{\partial}{\partial x_{k}} \alpha(A(x))=\left\langle\frac{\partial A}{\partial x_{k}}(x), \frac{1}{v^{*} u} v u^{*}\right\rangle=\operatorname{Re} \frac{u^{*} \frac{\partial A}{\partial x_{k}}(x) v}{u^{*} v}
\]
where \(v\) and \(u\) are right and left eigenvectors for the rightmost eigenvalue \(\lambda\).

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\]
where \(v\) and \(u\) are right and left eigenvectors for the rightmost eigenvalue \(\lambda\).
Hanso (Hybrid Algorithm for Nonsmooth Optimization):
publicly available Matlab software.

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Variational Analysis of Non-Lipschitz Spectral Functions J.V. Burke and M.L. Overton, Math. Programming (2001)

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Orr-Sommerfeld
Matrix ( \(n=99\),

\section*{Pseudospectra}

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The area swept out in the complex plane by the eigenvalues under perturbation.
\[
\sigma_{\epsilon}(A)=\{z \in \mathbf{C}: \operatorname{det}(A+E-z I)=0 \text { for some } E \text { with }\|E\| \leq \epsilon\}
\]

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A more robust measure of system behaviour than eigenvalues.

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A more robust measure of system behaviour than eigenvalues.
\[
\begin{aligned}
& \text { For }\|\cdot\|=\|\cdot\|_{2}, \\
& \qquad \sigma_{\epsilon}(A)=\left\{z \in \mathbf{C}:\left\|(A-z I)^{-1}\right\| \geq \epsilon^{-1}\right\}
\end{aligned}
\]

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A more robust measure of system behaviour than eigenvalues.
For \(\|\cdot\|=\|\cdot\|_{2}\),
\[
\begin{aligned}
\sigma_{\epsilon}(A) & =\left\{z \in \mathbf{C}:\left\|(A-z I)^{-1}\right\| \geq \epsilon^{-1}\right\} \\
& =\left\{z \in \mathbf{C}: s_{n}(A-z I) \leq \epsilon\right\}
\end{aligned}
\]
where \(s_{n}\) denotes smallest singular value:
\[
A-z I=U \operatorname{diag}(s) V^{*}
\]
with \(U^{*} U=V^{*} V=I\).

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\end{aligned}
\]
where \(s_{n}\) denotes smallest singular value:
\[
A-z I=U \operatorname{diag}(s) V^{*}
\]
with \(U^{*} U=V^{*} V=I\).
Let \(f(x, y)=s_{n}(A-(x+i y) I)\). Then pseudospectra are lower level sets of \(f\).

\section*{Pseudospectra}

\section*{Orr-Sommerfeld Matrix ( \(n=99, \epsilon=10^{-4}, 10^{-3}, 10^{-2}\) )}

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Pseudospectra
Orr-Sommerfeld
Matrix ( \(n=99\),


Black dots are eigenvalues and colored curves are pseudospectral boundaries. Note the pseudospectra are not convex.

\section*{Constructing \(E\) given \(z \in \partial \sigma_{\epsilon}(A)\)}

\section*{Let}

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Orr-Sommerfeld
Matrix ( \(n=99\),
\[
A-z I=U \operatorname{diag}(s) V^{*}=\sum_{j=1}^{n} s_{j} u_{j} v_{j}^{*}, \quad s_{n}=\epsilon
\]
\[
\text { with } U^{*} U=V^{*} V=I
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\section*{Constructing \(E\) given \(z \in \partial \sigma_{\epsilon}(A)\)}

Let
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\]
with \(U^{*} U=V^{*} V=I\).
Then if we set \(u=u_{n}, v=v_{n}, E=-\epsilon u v^{*}\) we have
\[
\operatorname{det}(A-z I+E)=0
\]
so \(z\) is an eigenvalue of \(A+E\).

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\operatorname{det}(A-z I+E)=0
\]
so \(z\) is an eigenvalue of \(A+E\).
Key point: can choose \(E\) to have rank one.
```

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\section*{Pseudospectra}

Constructing \(E\) given \(z \in \partial \sigma_{\epsilon}(A)\)
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\operatorname{det}(A-z I+E)=0
\]
so \(z\) is an eigenvalue of \(A+E\).
Key point: can choose \(E\) to have rank one. Furthermore
\[
(A-z I) v=\epsilon u, \quad u^{*}(A-z I)=\epsilon v^{*}
\]

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Constructing \(E\) given \(z \in \partial \sigma_{\epsilon}(A)\)
Let

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\[
A-z I=U \operatorname{diag}(s) V^{*}=\sum_{j=1}^{n} s_{j} u_{j} v_{j}^{*}, \quad s_{n}=\epsilon
\]
with \(U^{*} U=V^{*} V=I\).
Then if we set \(u=u_{n}, v=v_{n}, E=-\epsilon u v^{*}\) we have
\[
\operatorname{det}(A-z I+E)=0
\]
so \(z\) is an eigenvalue of \(A+E\).
Key point: can choose \(E\) to have rank one. Furthermore
\[
(A-z I) v=\epsilon u, \quad u^{*}(A-z I)=\epsilon v^{*}
\]
so
\[
(A-z I+E) v=0, \quad u^{*}(A-z I+E)=0 .
\]

Thus the right and left singular vectors of \(A-z I\) for the singular value \(\epsilon\) are also right and left eigenvectors of \(A+E\) for the eigenvalue \(z\).

\section*{Pseudospectral Radius and Abscissa}

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Pseudospectral radius: modulus of outermost point in \(\sigma_{\epsilon}(A)\)
\[
\rho_{\epsilon}(A)=\max \left\{|z|: z \in \sigma_{\epsilon}(A)\right\}
\]

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Computing these quantities: nontrivial because \(\sigma_{\epsilon}(A)\) is not convex.

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Computing these quantities: nontrivial because \(\sigma_{\epsilon}(A)\) is not convex.

Criss-cross algorithm for computing the pseudospectral abscissa \(\alpha_{\epsilon}(A)\) : based on repeatedly computing eigenvalues of \(2 n \times 2 n\) Hamiltonian matrices and checking whether any are imaginary, and computing SVDs for each imaginary eigenvalue.

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Criss-cross algorithm for computing the pseudospectral abscissa \(\alpha_{\epsilon}(A)\) : based on repeatedly computing eigenvalues of \(2 n \times 2 n\) Hamiltonian matrices and checking whether any are imaginary, and computing SVDs for each imaginary eigenvalue.
Too expensive if \(n\) large.

\section*{Approximating the Pseudospectral Abscissa if \(n\) is Big}

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We want a rightmost point \(z\) of \(\sigma_{\epsilon}(A)\), so \(s_{n}(A-z I)=\epsilon\). Let \(v\) and \(u\) be corresponding right and left singular vectors. We know that \(z\) is an eigenvalue of \(B=A-\epsilon u v^{*}\) with right and left eigenvectors \(v\) and \(u\).

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Let us generate a sequence
\[
B^{(k)}=A-\epsilon u^{(k)}\left(v^{(k)}\right)^{*}
\]
with \(\left\|u^{(k)}\right\|=\left\|v^{(k)}\right\|=1\). We want \(u^{(k)} \rightarrow u, v^{(k)} \rightarrow v\).

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\[
B^{(k)}=A-\epsilon u^{(k)}\left(v^{(k)}\right)^{*}
\]
with \(\left\|u^{(k)}\right\|=\left\|v^{(k)}\right\|=1\). We want \(u^{(k)} \rightarrow u, v^{(k)} \rightarrow v\).
No Hamiltonian eigenvalue decompositions or SVDs allowed. The only matrix operations are the computation of eigenvalues with largest real part and their corresponding right and left eigenvectors, which can be done efficiently using the implicitly restarted Arnoldi method (ARPACK).

\section*{RP-Compatible Right and Left Eigenvectors}

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A pair of right and left eigenvectors \(p\) and \(q\) for a simple eigenvalue \(\lambda\) is called \(R P\)-compatible if \(\|p\|=\|q\|=1\) and \(p^{*} q\) is real and positive, and therefore in the interval \((0,1]\).

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This defines right and left eigenvectors uniquely up to \(p \leftarrow e^{i \theta} p\), \(q \leftarrow e^{i \theta} q\).

\section*{New Algorithm to Approximate \(\alpha_{\epsilon}(A)\)}

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1. Let \(z^{(0)}\) be a rightmost eigenvalue of \(A\), with RP-compatible right and left eigenvectors \(v^{(0)}\) and \(u^{(0)}\). Set \(B^{(0)}=A-\epsilon u^{(0)}\left(v^{(0)}\right)^{*}\).

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2. For \(k=1,2, \ldots\) : let \(z^{(k)}\) be a rightmost eigenvalue of \(B^{(k)}\) with RP-compatible right and left eigenvectors \(v^{(k)}\) and \(u^{(k)}\). Set \(B^{(k+1)}=A-\epsilon u^{(k)}\left(v^{(k)}\right)^{*}\).

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Clearly, \(\operatorname{Re} z^{(k)} \leq \alpha_{\epsilon}(A)\) for all \(k\).

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Clearly, \(\operatorname{Re} z^{(k)} \leq \alpha_{\epsilon}(A)\) for all \(k\).
Almost always: \(z^{(k)} \rightarrow z\), a locally rightmost point of \(\sigma_{\epsilon}(A)\), and \(v^{(k)}\) and \(u^{(k)}\) converge to right and left singular vectors \(v\) and \(u\) corresponding to smallest singular value of \(A-z I\).

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Clearly, \(\operatorname{Re} z^{(k)} \leq \alpha_{\epsilon}(A)\) for all \(k\).
Almost always: \(z^{(k)} \rightarrow z\), a locally rightmost point of \(\sigma_{\epsilon}(A)\), and \(v^{(k)}\) and \(u^{(k)}\) converge to right and left singular vectors \(v\) and \(u\) corresponding to smallest singular value of \(A-z I\).
Often, but not always, \(z\) is a globally rightmost point so
\(\operatorname{Re} z=\alpha_{\epsilon}(A)\).

\section*{New Algorithm to Approximate \(\alpha_{\epsilon}(A)\)}

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Almost always: \(z^{(k)} \rightarrow z\), a locally rightmost point of \(\sigma_{\epsilon}(A)\), and \(v^{(k)}\) and \(u^{(k)}\) converge to right and left singular vectors \(v\) and \(u\) corresponding to smallest singular value of \(A-z I\).
Often, but not always, \(z\) is a globally rightmost point so \(\operatorname{Re} z=\alpha_{\epsilon}(A)\).
We have theorems characterizing fixed points of the algorithm and proving local convergence at a geometric rate for \(\epsilon\) small.

\section*{Orr-Sommerfeld Matrix ( \(n=99, \epsilon=10^{-4}, 10^{-2}\) )}

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Orr-Sommerfeld


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\section*{Pseudospectra}

Orr-Sommerfeld


\section*{Minimizing \(\alpha_{\epsilon}(A(x))\) over Parametrized Matrix \(A(x)\)}

For given \(x\) in parameter space \(\mathbf{R}^{p}\), compute \(\alpha_{\epsilon}(A(x))\) by criss-cross algorithm if \(n\) small and otherwise by the new algorithm.

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\section*{Minimizing \(\alpha_{\epsilon}(A(x))\) over Parametrized Matrix \(A(x)\)}

For given \(x\) in parameter space \(\mathbf{R}^{p}\), compute \(\alpha_{\epsilon}(A(x))\) by criss-cross algorithm if \(n\) small and otherwise by the new algorithm.
Like \(\alpha, \alpha_{\epsilon}\) is nonsmooth and nonconvex, but unlike \(\alpha\), it is locally Lipschitz for \(\epsilon>0\) (although \(\sigma_{\epsilon}\) is not).

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For given \(x\) in parameter space \(\mathbf{R}^{p}\), compute \(\alpha_{\epsilon}(A(x))\) by criss-cross algorithm if \(n\) small and otherwise by the new algorithm.
Like \(\alpha, \alpha_{\epsilon}\) is nonsmooth and nonconvex, but unlike \(\alpha\), it is locally Lipschitz for \(\epsilon>0\) (although \(\sigma_{\epsilon}\) is not).
Derivatives:
\[
\frac{\partial}{\partial x_{k}} \alpha_{\epsilon}(A(x))=\left\langle\frac{\partial A}{\partial x_{k}}(x), \frac{1}{v^{*} u} v u^{*}\right\rangle=\operatorname{Re} \frac{u^{*} \frac{\partial A}{\partial x_{k}}(x) v}{u^{*} v}
\]
where \(v\) and \(u\) are right and left singular vectors for the singular value \(\epsilon\) of \(A-z I\) with \(z\) the rightmost point of \(\sigma_{\epsilon}(A)\), equivalently RP-compatible right and left eigenvectors for the eigenvalue \(z\) of \(A-\epsilon u v^{*}\).

\section*{Minimizing \(\alpha_{\epsilon}(A(x))\) over Parametrized Matrix \(A(x)\)}

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As earlier, use Gradient Sampling or BFGS.

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As earlier, use Gradient Sampling or BFGS.
Example: \(A(x)=F+G K H\) with \(x=\operatorname{vec}(K)\), a static output feedback control design problem for a turbo generator with \(n=10, \ell=m=2\), so controller \(K \in \mathbf{R}^{2 \times 2}\).

\section*{A Turbo Generator Control Problem}

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Pseudospectra for Turbo-Generator with No Feedback



Pseudospectra for open-loop turbo generator plant with no feedback. Matrix ( \(n=99\),

\section*{Turbo Generator with Optimized Eigenvalues}

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Pseudospectra for turbo generator plant with feedback computed by minimizing the spectral abscissa \(\alpha\)

\author{
Pseudospectra
}

\section*{Turbo Generator with Optimized \(\epsilon\)-Pseudospectrum}

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Pseudospectra for Turbo-Generator when \(1 \sigma^{-1.5}\)-Pseudospectrum is Optimized



Pseudospectra for turbo generator plant with feedback computed by minimizing the pseudospectral abscissa \(\alpha_{\epsilon}\) with \(\epsilon=10^{-1.5}\)

\section*{Turbo Generator with Optimized Dist. to Instability}

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Pseudospectra for Turbo-Generator when Complex Stability Radius is Optimized



Pseudospectra for turbo generator plant with feedback computed by maximizing the distance to instability: largest \(\epsilon\) so that \(\alpha_{\epsilon}(A(x)) \leq 0\).

\section*{References for Part III}

Origins of Pseudospectra in 1980s:

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Landau, Varah, Godunov, Demmel, Wilkinson, Trefethen, ?.
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## Thanks a lot for your attention!

