

Relations between Rayleigh Quotient Iteration and Classical Root Finding Algorithms

Jens-Peter M. Zemke
zemke@tu-harburg.de

Institut für Numerische Simulation
Technische Universität Hamburg-Harburg

Universiteit Utrecht, The Netherlands
2010/03/03



Outline

Classical Root Finding

- Newton's Method
- The Secant Method
- König's Method
- The Opitz-Larkin Method

Rayleigh Quotient Iteration

- John William Strutt's RQI
- Inverse Iteration
- Symmetric RQI
- Two-Sided RQI

The Hessenberg-Matrix Point Of View

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Newton's method

The best known method for the **computation of a root** of a rational function

$$f : \mathbb{C} \rightarrow \mathbb{C}, \quad f(z) := \frac{p(z)}{q(z)}, \quad p, q \in \mathbb{P}_m$$

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Newton's method costs **two function evaluations** per step, one evaluation of the function, one evaluation of its derivative.

The secant method

If the derivative of the function $f : \mathbb{C} \rightarrow \mathbb{C}$ is not at hand, we could use the first **divided difference** which gives the **secant method**:

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In general, **the secant method locally wins** (Raydan, 1993).

Schröder's and König's methods

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This family is nowadays known as **"König's method"**:

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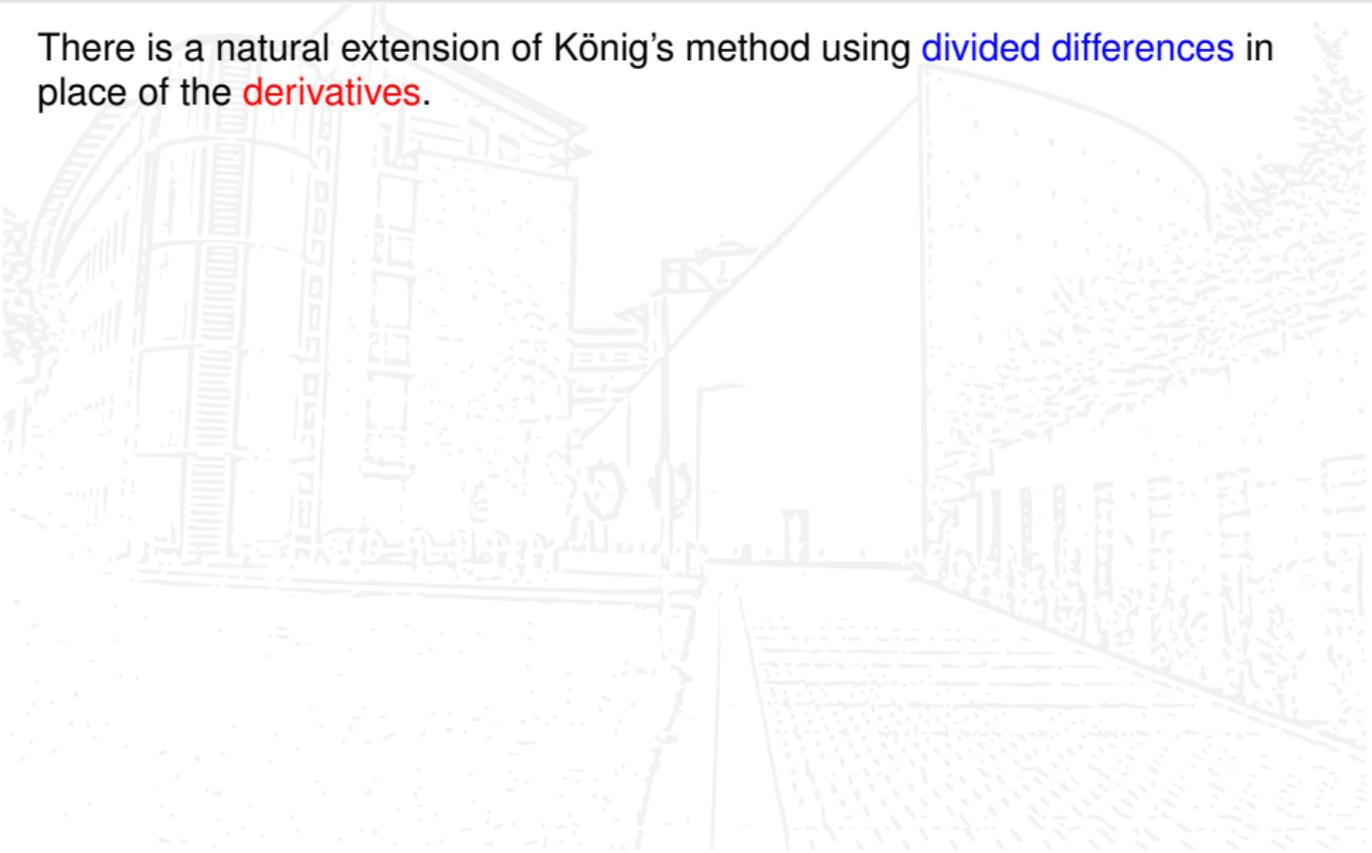
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König's method for $s = 1$ is **Newton's method**,

$$z_{k+1} = z_k + \frac{(1/f)(z_k)}{(1/f)'(z_k)} = z_k - \frac{1/f(z_k)}{f'(z_k)/(f(z_k))^2} = z_k - \frac{f(z_k)}{f'(z_k)}.$$

The Opitz-Larkin method

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Opitz **never published** a detailed version. Independently, **23 years later** F. M. Larkin re-developed Opitz' method and published parts of Opitz' results with proofs in (Larkin, 1981) and the predecessor (Larkin, 1980).

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We will refer to this method as **the Opitz-Larkin method**. The Opitz-Larkin method is **based on iterations** of the form

$$x_{k+1} = z_k + \frac{[z_1, z_2, \dots, z_{k-1}](1/f)}{[z_1, z_2, \dots, z_{k-1}, z_k](1/f)}.$$

The Opitz-Larkin method

Mostly, the z_i are all **distinct** and the next iterate is used as **new evaluation point** $z_{k+1} = x_{k+1}$,

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This variant of the Opitz-Larkin method converges with **R-order 2**.

Frequently, the Opitz-Larkin method is used with **truncation**:

$$z_{k+1} = z_k + \frac{[z_{k-p}, \dots, z_{k-1}](1/f)}{[z_{k-p}, \dots, z_{k-1}, z_k](1/f)},$$

see (Opitz, 1958, Seite 277, Gleichung (9)) and (Larkin, 1981, Section 4, pages 98–99).

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When we use **only confluent divided differences** in the truncated Opitz-Larkin method with truncation parameter $p = s$, we **recover** König's method:

$$\begin{aligned}
 z_{k+1} &= z_k + \frac{\overbrace{[z_k, \dots, z_k]}^s (1/f)}{\underbrace{[z_k, \dots, z_k, z_k]}_{s+1} (1/f)} \\
 &= z_k + \frac{(1/f)^{(s-1)}(z_k)/(s-1)!}{(1/f)^{(s)}(z_k)/s!} = z_k + s \frac{(1/f)^{(s-1)}(z_k)}{(1/f)^{(s)}(z_k)}.
 \end{aligned}$$

The Opitz-Larkin method

Truncated Opitz-Larkin with $p = 1$ is the secant method,

$$\begin{aligned}z_{k+1} &= z_k + \frac{[z_{k-1}](1/f)}{[z_{k-1}, z_k](1/f)} \\&= z_k + \frac{1}{f(z_{k-1})} \cdot \frac{z_{k-1} - z_k}{1/f(z_{k-1}) - 1/f(z_k)} \\&= z_k + \frac{f(z_k)f(z_{k-1})}{f(z_{k-1})} \cdot \frac{z_{k-1} - z_k}{f(z_k) - f(z_{k-1})} \\&= z_k - \frac{f(z_k)}{[z_{k-1}, z_k]f}.\end{aligned}$$

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Confluent truncated Opitz-Larkin with $p = 1$ is Newton's method.

The Opitz-Larkin method

In general, the Opitz-Larkin method is closely connected to **rational interpolation** of **the inverse function** (Larkin, 1981, Theorem 1, page 96):

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Theorem (Larkin 1981)

If, for any integer $k > 1$, there exists a rational function of the form

$$r_k(z) = \frac{q_d(z)}{z - \alpha}, \quad \forall z,$$

where q_d is a polynomial of degree $d \leq k - 2$, such that $q_d(\alpha) \neq 0$ and

$$r_k(z_j) = f(z_j)^{-1}, \quad j = 1, 2, \dots, k,$$

then

$$z_k + \frac{[z_1, z_2, \dots, z_{k-1}](1/f)}{[z_1, z_2, \dots, z_{k-1}, z_k](1/f)} = \alpha.$$

The Opitz-Larkin method

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In the earlier publication (Larkin, 1980) Larkin used **another approach** to obtain the rational interpolant and gave pointers to articles that investigated the **rate of convergence** of such (direct and inverse) rational interpolations.

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Most important are the articles (Tornheim, 1964) and (Jarratt and Nudds, 1965). We state the **main results** contained in these articles.

The Opitz-Larkin method

In the paper (Tornheim, 1964), Tornheim considered the case of **direct**

$$f^{(\ell_j)}(x_{i-j}) = \left(\frac{p}{q}\right)^{(\ell_j)}(x_{i-j})$$

and **inverse rational interpolation**

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where

$$\ell_j = 0, 1, \dots, m_j - 1, \quad m = \sum_{j=0}^k m_j = \deg(p) + \deg(q) + 1,$$

and k given distinct points $x_{i-j} \quad j = 1, \dots, k,$

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and gave its **rate of convergence** (Tornheim, 1964, Theorem 2).

The Opitz-Larkin method

Theorem (Tornheim 1964; conditions for the theorem)

Suppose an k -point iterative method is defined by the procedure to solve the equation $f(x) = 0$ by direct or inverse rational interpolation with m_j coincident interpolating points at x_{i-j} ($j = 1, \dots, k$) for the i -th iteration.

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$$\mathbf{M} = \begin{vmatrix} a_d & a_{d-1} & \cdots & a_{2d+2-m} \\ a_{d+1} & a_d & \cdots & \\ \vdots & \vdots & & \vdots \\ a_{m-2} & & \cdots & a_d \end{vmatrix} \neq 0.$$

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Here d is the degree of the numerator and e is the degree of the denominator of the rational function used; $d + e + 1 = m$; a_i is 0 if $i < 0$, otherwise it is the i th derivative of $f(x)$ (for direct interpolation) or of its inverse function (for inverse interpolation) at $x = x^*$.

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Theorem (Tornheim 1964; result of the theorem)

Then there is a neighborhood N^ of x^* such that if x_1, \dots, x_k are in N^* , the sequence $\{x_i\}$ converges to x^* .*

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Then there is a neighborhood N^* of x^* such that if x_1, \dots, x_k are in N^* , the sequence $\{x_i\}$ converges to x^* . Moreover, the **order of convergence** u , if it exists, is the **positive root of the equation**

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In the **context of the Opitz-Larkin method**, we have to consider the **limit of the positive root** for $k \rightarrow \infty$.

He also gave a **“comparison result”** that predicts faster convergence when the (inverse) function is evaluated to higher order at the **last iterates**.

The Opitz-Larkin method

Lemma (Tornheim 1964)

Suppose that the coefficients of

$$a(x) := x^n - a_1x^{n-1} - \dots - a_n$$

satisfy

$$a_1 + a_2 + \dots + a_n > 1, \quad a_1 \geq a_2 \geq \dots \geq a_n \geq 0.$$

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By *Descartes' rule of signs* the polynomial a has a **unique positive root** $u > 1$.
If the coefficients b_i of

$$b(x) := x^n - b_1x^{n-1} - \dots - b_n$$

are a *permutation* of the coefficients a_i , then the positive root v of b is **less than** u .

The Opitz-Larkin method

In the paper (Jarratt and Nudds, 1965), Jarratt and Nudds give a detailed treatment of the case of **rational interpolation** with

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As you might already have guessed: We are going to prove that **RQI is the Opitz-Larkin method**. One instance of RQI actually is the Opitz-Larkin method applied to the characteristic polynomial of the given matrix.

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In the **second edition** of the first volume of his book “The Theory of Sound” (Strutt, 1894), **John William Strutt**, 3rd Baron Rayleigh included on page 110 the following passage:

The stationary property of the roots of Lagrange's determinant (3) § 84, suggests a general method of approximating to their values. Beginning with assumed rough approximations to the ratios $A_1:A_2:A_3,\dots$ we may calculate a first approximation to p^2 from

$$p^2 = \frac{\frac{1}{2} c_{11}A_1^2 + \frac{1}{2} c_{22}A_2^2 + \dots + c_{12}A_1A_2 + \dots}{\frac{1}{2} a_{11}A_1^2 + \frac{1}{2} a_{22}A_2^2 + \dots + a_{12}A_1A_2 + \dots} \dots\dots (3).$$

With this value of p^2 we may recalculate the ratios $A_1:A_2,\dots$ from any $(m-1)$ of equations (5) § 84, then again by application of (3) determine an improved value of p^2 , and so on.]

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In **modern notation**, stated for the **Hermitean algebraic eigenvalue problem**

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and uses the linear system

$$(\mathbf{A} - \rho(\mathbf{v}_k)\mathbf{I}_n)\mathbf{v}_{k+1} = \mathbf{e}_j$$

for some **standard unit vector** \mathbf{e}_j to compute a new approximate eigenvector \mathbf{v}_{k+1} . He was, of course, only interested in **its direction**.

Original RQI

This is **repeated several times**, i.e.,

$$\mathbf{v}_{k+1} = (\mathbf{A} - \rho(\mathbf{v}_k)\mathbf{I}_n)^{-1}\mathbf{e}_j, \quad k = 0, 1, \dots$$

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Classical RQI can thus be stated in modern notation as

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for some suitably chosen $j \in \{1, 2, \dots, n\}$, which might vary, **depending on the computed approximate eigenvector**.

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for some suitably chosen $j \in \{1, 2, \dots, n\}$, which might vary, **depending on the computed approximate eigenvector**.

Classical Rayleigh quotient iteration mostly converges locally with **quadratic** order of convergence.

Inverse Iteration

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In the **most basic variant** of inverse iteration the **shift τ is never updated**, but the right-hand side is replaced by the latest approximate eigenvector:

$$\mathbf{v}_{k+1} = \frac{(\mathbf{A} - \tau \mathbf{I}_n)^{-1} \mathbf{v}_k}{\|(\mathbf{A} - \tau \mathbf{I}_n)^{-1} \mathbf{v}_k\|}, \quad k = 0, 1, \dots$$

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There exist **variants** which use **other scalings**, mostly using as left vector some standard unit vector $\ell = \mathbf{e}_j$:

$$\mathbf{v}_{k+1} = \frac{(\mathbf{A} - \tau \mathbf{I}_n)^{-1} \mathbf{v}_k}{\mathbf{e}_j^\top (\mathbf{A} - \tau \mathbf{I}_n)^{-1} \mathbf{v}_k}, \quad k = 0, 1, \dots$$

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In the latter context, $\mathbf{e}_j^T (\mathbf{A} - \tau \mathbf{I}_n)^{-1} \mathbf{v}_k \approx (\lambda - \tau)^{-1}$ gives an **eigenvalue approximation**.

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In both variants also the **Rayleigh quotient** can be used, the Rayleigh quotient uniquely solves the **least squares problem**

$$\rho(\mathbf{v}_k) = \operatorname{argmin}_{\rho \in \mathbb{C}} \|\mathbf{A}\mathbf{v}_k - \mathbf{v}_k\rho\|$$

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Both these methods typically exhibit a **quadratic convergence behavior**.

Symmetric RQI

When automatic computers became available, the **combination of inverse iteration with Rayleigh's original RQI** resulted in the locally **Q-cubically convergent** (symmetric) RQI

$$\mathbf{v}_{k+1} = \frac{(\mathbf{A} - \rho(\mathbf{v}_k)\mathbf{I}_n)^{-1}\mathbf{v}_k}{\|(\mathbf{A} - \rho(\mathbf{v}_k)\mathbf{I}_n)^{-1}\mathbf{v}_k\|}, \quad k = 0, 1, \dots$$

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Crandall was the **first** who investigated the three variants (the original Rayleigh quotient iteration; inverse iteration with fixed shift; symmetric RQI) and proved their convergence rates to be quadratic, linear, and cubic, respectively, see (Crandall, 1951).

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Ostrowski proved that unsymmetric RQI still has a **quadratic convergence rate**, (Ostrowski, 1959e). In (Ostrowski, 1959c), he also gave a variant that **recovers** the **cubic convergence rate** at the expense of the necessity to solve two linear systems every step instead of only one.

Two-Sided RQI

When $\mathbf{A} \in \mathbb{C}^{n \times n}$ is **no longer Hermitean**, the **cubic convergence is lost** and Ostrowski suggested in (Ostrowski, 1959c) the use of a **two-sided RQI**.

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Two-sided RQI is based on the **two-sided Rayleigh quotient**

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The iteration involves **two sequences of vectors**,

$$\begin{aligned} \mathbf{v}_{k+1} &= (\mathbf{A} - \rho(\mathbf{w}_k, \mathbf{v}_k) \mathbf{I}_n)^{-1} \mathbf{v}_k, \\ \mathbf{w}_{k+1} &= (\mathbf{A} - \rho(\mathbf{w}_k, \mathbf{v}_k) \mathbf{I}_n)^{-H} \mathbf{w}_k, \end{aligned} \quad k = 0, 1, \dots$$

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This trick **recovers the cubic convergence of RQI** at the price of the solution of an additional system.

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Ostrowski worked out a **more detailed analysis** than Crandall. He published a **series of six papers on RQI**, (Ostrowski, 1959b; Ostrowski, 1959c; Ostrowski, 1959d; Ostrowski, 1959e; Ostrowski, 1959a). He measured the rate of convergence with respect to the number of solutions of linear systems, which he called one **“Horner”**. He was a little unfair to the two-sided variant, as these two Horners are related to each other (one decomposition, two forward and backward substitutions with the same two triangular matrices).

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The proofs by Crandall and Ostrowski are [beautiful and worth reading](#).

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The proofs by Crandall and Ostrowski are **beautiful and worth reading**.

But we feel that **a more direct proof of convergence** for the different variants of RQI and related algorithms would be very helpful, especially when we want to investigate the **overall behavior**: the basins of attraction; global convergence; effects of perturbation and inexact methods, ...

Outline

Classical Root Finding

- Newton's Method
- The Secant Method
- König's Method
- The Opitz-Larkin Method

Rayleigh Quotient Iteration

- John William Strutt's RQI
- Inverse Iteration
- Symmetric RQI
- Two-Sided RQI

The Hessenberg-Matrix Point Of View

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The **implicit Q-Theorem** gives **uniqueness** of the upper part of the reduction to Hessenberg form in case of given first column \mathbf{q} , if we fix the **signs** of the elements in the lower diagonal, e.g., to be non-negative real.

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We use the implicit Q-Theorem to **unitarily transform** the pair (\mathbf{A}, \mathbf{q}) with $\|\mathbf{q}\|_2 = 1$ to the pair $(\mathbf{H}_n, \mathbf{e}_1)$, where \mathbf{H}_n is **upper Hessenberg** and \mathbf{e}_1 denotes the first standard unit vector.

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The following **Matlab-code** gives the transformed pair:

```
[Q,R] = qr(q);
[P,H] = hess(Q'*A*Q);
signs = sign(diag(H,-1));
S = diag(cumprod([1;signs]));
P = P*S;
H = S'*H*S;
```

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The **first resolvent identity** (Chatelin, 1993, Lemma 2.2.1, p. 63), valid for $z_1 \neq z_2$ from the resolvent set, gives

$$({}^{z_1}\mathbf{H}_n)^{-1}({}^{z_2}\mathbf{H}_n)^{-1} = (z_1\mathbf{I}_n - \mathbf{H}_n)^{-1}(z_2\mathbf{I}_n - \mathbf{H}_n)^{-1} \quad (1a)$$

$$= \frac{({}^{z_1}\mathbf{H}_n)^{-1} - ({}^{z_2}\mathbf{H}_n)^{-1}}{z_2 - z_1} = -[z_1, z_2]({}^z\mathbf{H}_n)^{-1}. \quad (1b)$$

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The first resolvent identity is based on the **trivial observation** that

$$(z_2\mathbf{I}_n - \mathbf{H}_n) - (z_1\mathbf{I}_n - \mathbf{H}_n) = (z_2 - z_1)\mathbf{I}_n.$$

Simplification

This identity can be **generalized** to k distinct points of evaluation:

$$\prod_{i=1}^k (z_i \mathbf{H}_n)^{-1} = (-1)^{k-1} [z_1, \dots, z_k] (z \mathbf{H}_n)^{-1}. \quad (2)$$

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The inverse of the characteristic matrix $z \mathbf{H}_n$ is the **rational function**

$$(z \mathbf{H}_n)^{-1} = \frac{\text{adj}(z \mathbf{H}_n)}{\chi(z)} =: \frac{\mathbf{P}_n(z)}{\chi(z)}, \quad \chi(z) := \det(z \mathbf{H}_n), \quad (3)$$

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Thus, **confluent divided differences** are **well-defined** and we do not need to restrict the points $\{z_i\}_{i=1}^k$ from the resolvent set.

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We only need here a **well-known result** on a recurrence for the determinants of unreduced Hessenberg matrices, see, e.g., (Franklin, 1968, Section 7.11, p. 252, Eqn. (8)), or, the **probably earliest reference** (Schweins, 1825, Erste Abtheilung, IV. Abschnitt, § 154, Seite 361, Gleichung 560)).

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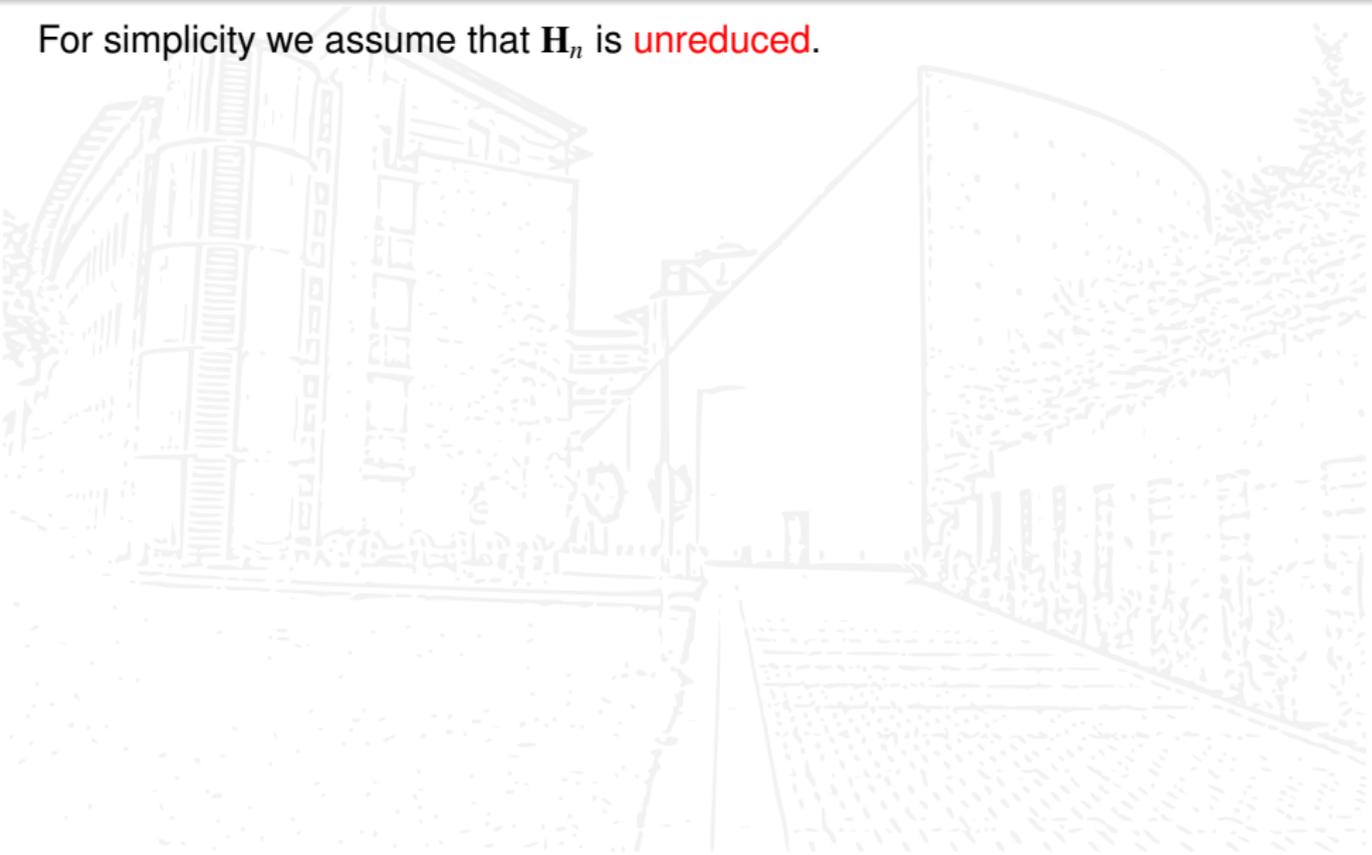
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There exist **short proofs** based on **Laplace expansion** and **Cramer's rule**.

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Polynomial vectors ν and $\check{\nu}$ are defined by

$$\nu(z) := \left(\frac{\chi_{j+1:n}(z)}{h_{j:n-1}} \right)_{j=1}^n \quad \text{and} \quad \check{\nu}(z) := \left(\frac{\chi_{1:j-1}(z)}{h_{1:j-1}} \right)_{j=1}^n. \quad (4)$$

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$$\nu(z) := \left(\frac{\chi_{j+1:n}(z)}{h_{j:n-1}} \right)_{j=1}^n \quad \text{and} \quad \check{\nu}(z) := \left(\frac{\chi_{1:j-1}(z)}{h_{1:j-1}} \right)_{j=1}^n. \quad (4)$$

The elements are denoted by $\nu_j(z)$ and $\check{\nu}_j(z)$, $j = 1, \dots, n$. We remark that $\nu_n \equiv 1 \equiv \check{\nu}_1$.

Simplification

For simplicity we assume that \mathbf{H}_n is **unreduced**.

We denote **products of sub-diagonal elements** of the unreduced Hessenberg matrices $\mathbf{H}_n \in \mathbb{C}^{n \times n}$ by

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The polynomials $\chi_{i:j}$ are the **characteristic polynomials** of **submatrices** of \mathbf{H}_n ,

$$\chi_{i:j}(z) := \det({}^z\mathbf{H}_{i:j}) = \det(z\mathbf{I}_{j-i+1} - \mathbf{H}_{i:j}).$$

Simplification

By (Z, 2006, Lemma 3.1, Eqn. (3.5)) for z in the **resolvent set**

$$({}^z\mathbf{H}_n)\nu(z) = \frac{\chi(z)}{h_{1:n-1}}\mathbf{e}_1 \Leftrightarrow \frac{\nu(z)h_{1:n-1}}{\chi(z)} = ({}^z\mathbf{H}_n)^{-1}\mathbf{e}_1, \quad (5a)$$

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The **repeated application of resolvents** to \mathbf{e}_1 results in

$$\left(\prod_{i=1}^k ({}^{z_i}\mathbf{H}_n)^{-1}\right)\mathbf{e}_1 = (-1)^{k-1}[z_1, \dots, z_k]({}^z\mathbf{H}_n)^{-1}\mathbf{e}_1 \quad (6)$$

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We note that $z\mathbf{I}_n - {}^z\mathbf{H}_n = z\mathbf{I}_n - (z\mathbf{I}_n - \mathbf{H}_n) = \mathbf{H}_n$.

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$$\mathbf{v}_{k+1} = \left(\prod_{i=1}^k (z_i \mathbf{H}_n)^{-1} \right) \mathbf{e}_1, \quad x_{k+1} = \frac{\mathbf{e}_n^\top \mathbf{H}_n \mathbf{v}_{k+1}}{\mathbf{e}_n^\top \mathbf{v}_{k+1}},$$

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and thus the approximate eigenvalues are given by the **Opitz-Larkin method**:

$$x_{k+1} = \frac{\mathbf{e}_n^\top \mathbf{H}_n \left(\prod_{i=1}^k (z_i \mathbf{H}_n)^{-1} \right) \mathbf{e}_1}{\mathbf{e}_n^\top \left(\prod_{i=1}^k (z_i \mathbf{H}_n)^{-1} \right) \mathbf{e}_1} = \frac{\mathbf{e}_n^\top (z_k \mathbf{I}_n - (z_k \mathbf{H}_n)) \left(\prod_{i=1}^k (z_i \mathbf{H}_n)^{-1} \right) \mathbf{e}_1}{\mathbf{e}_n^\top \left(\prod_{i=1}^k (z_i \mathbf{H}_n)^{-1} \right) \mathbf{e}_1} \quad (8a)$$

$$= z_k - \frac{\mathbf{e}_n^\top z_k \mathbf{H}_n \left(\prod_{i=1}^k (z_i \mathbf{H}_n)^{-1} \right) \mathbf{e}_1}{\mathbf{e}_n^\top \left(\prod_{i=1}^k (z_i \mathbf{H}_n)^{-1} \right) \mathbf{e}_1} = z_k - \frac{\mathbf{e}_n^\top \left(\prod_{i=1}^{k-1} (z_i \mathbf{H}_n)^{-1} \right) \mathbf{e}_1}{\mathbf{e}_n^\top \left(\prod_{i=1}^k (z_i \mathbf{H}_n)^{-1} \right) \mathbf{e}_1} \quad (8b)$$

$$= z_k + \frac{[z_1, \dots, z_{k-1}](1/\chi)}{[z_1, \dots, z_{k-1}, z_k](1/\chi)}. \quad (8c)$$

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When we **update the shifts** by choosing $z_{k+1} = x_{k+1}$ we obtain the **standard variant of the Opitz-Larkin method**. This method has asymptotically second order convergence against the roots of the characteristic polynomial χ .

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Inverse iteration with fixed shift $\tau = z_1 = z_2 = \dots = z_k$ results in the recurrence

$$x_{k+1} = \tau + \frac{[\tau, \dots, \tau](1/\chi)}{[\tau, \dots, \tau, \tau](1/\chi)} = \tau + k \frac{(1/\chi)^{(k-1)}(\tau)}{(1/\chi)^{(k)}(\tau)}. \quad (9)$$

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This knowledge together with an estimate for the **cost of preprocessing** (computing the LU decomposition; initializing a Krylov method using a seed system) and the **cost of the (approximate) solutions** of the systems enables to decide when to compute an update of the shift.

Simplification

The **original Rayleigh quotient iteration** (Strutt, 1894) with the symmetric Rayleigh quotient and, because of the symmetry, a **tridiagonal Hermitean Hessenberg matrix** \mathbf{H}_n , gives the update

$$z_{k+1} = \frac{\mathbf{e}_1^T (z_k \mathbf{H}_n)^{-H} \mathbf{H}_n (z_k \mathbf{H}_n)^{-1} \mathbf{e}_1}{\mathbf{e}_1^T (z_k \mathbf{H}_n)^{-H} (z_k \mathbf{H}_n)^{-1} \mathbf{e}_1} = \frac{\mathbf{e}_1^T \mathbf{H}_n (z_k \mathbf{H}_n)^{-2} \mathbf{e}_1}{\mathbf{e}_1^T (z_k \mathbf{H}_n)^{-2} \mathbf{e}_1} \quad (10a)$$

$$= \frac{\mathbf{e}_1^T (z_k \mathbf{I}_n - z_k \mathbf{H}_n) (z_k \mathbf{H}_n)^{-2} \mathbf{e}_1}{\mathbf{e}_1^T (z_k \mathbf{H}_n)^{-2} \mathbf{e}_1} \quad (10b)$$

$$= z_k - \frac{\mathbf{e}_1^T (z_k \mathbf{H}_n)^{-1} \mathbf{e}_1}{\mathbf{e}_1^T (z_k \mathbf{H}_n)^{-2} \mathbf{e}_1} = z_k + \frac{[z_k](\chi_{2:n}/\chi)}{[z_k, z_k](\chi_{2:n}/\chi)} \quad (10c)$$

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This is **Newton's method** on the **meromorphic function** r . As the poles of this meromorphic function are the eigenvalues of a submatrix, they interlace by Cauchy's interlace theorem the roots, which are the eigenvalues.

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Symmetric RQI for Hermitean matrices gives the update

$$z_{k+1} = z_k + \frac{[z_1, z_1, \dots, z_{k-1}, z_{k-1}, z_k](\chi_{2:n}/\chi)}{[z_1, z_1, \dots, z_{k-1}, z_{k-1}, z_k, z_k](\chi_{2:n}/\chi)}. \quad (11)$$

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This update has by the result of Tornheim asymptotically a **cubic convergence rate**, as we have to compute the limit of the real roots of the equations

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i.e., the maximal eigenvalue of a **Hessenberg matrix** with ones in the lower diagonal and twos in the last column. The **approximate eigenvector** of all ones to the approximate eigenvalue 3 gives the backward error $1/\sqrt{k}$ and the only real positive eigenvalue of the matrix is well separated, the other eigenvalues lie close to a circle of radius one around zero.

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If we take another standard unit vector \mathbf{e}_ℓ as left vector, we obtain the **Opitz-Larkin method applied to the meromorphic function**

$$m_\ell(z) = \frac{\chi(z)}{h_{1:\ell-1}\chi_{1+\ell:n}(z)}. \quad (12)$$

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The polynomials $\chi_{1+i:n}$ have degree $\deg(\chi_{1+i:n}) = n - i$ and leading coefficient one, thus they form a **basis of the space of polynomials** of degree less n .

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By luck or accident, we can construct a polynomial that is zero (of any order up to order $n - 1$) at one eigenvalue. This is of interest in case of (algebraically) multiple eigenvalues. In theory, there is always a left starting vector which ensures that the root is simple, as the multiple zero is reduced to a simple one.

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The **best choice** is the starting vector \mathbf{y} that represents the derivative of χ , i.e., the vector $\bar{\mathbf{y}}$ such that

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In this special case the rational function is the **Newton's update**

$$r(z; \bar{\mathbf{y}}) = \frac{\chi(z)}{\chi'(z)} \quad (15)$$

which has only **simple zeros** and poles between the eigenvalues.

Simplification

The Academic Example: The matrix $\mathbf{H}_4 = \text{triu}(\text{ones}(4), -1)$ has the eigenvalues 0 (double), 1, and 3, and the vector

$$\mathbf{y} = \begin{pmatrix} 4 \\ 0 \\ 2 \\ 2 \end{pmatrix} \quad (16)$$

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A variant of original RQI with starting vector \mathbf{e}_1 and test vector \mathbf{y} and updated shifts performs Newton's method on the Newton's update χ/χ' .

Simplification

The **two-sided RQI** method corresponds to a **confluent Opitz-Larkin** method with double nodes. In this method the left vector determines a polynomial, which is formed as a linear combination of characteristic polynomials of trailing submatrices.

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In single-sided RQI for non-Hermitian matrices, we change the vector \mathbf{y} that determines the denominator polynomial of the rational function

$$r(z; \mathbf{y}) = \frac{\chi(z)}{p(z; \mathbf{y})}$$

in every step and apply one step of the Opitz-Larkin method without confluent nodes. This gives **second order convergence**.

Conclusions and Outlook

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- ▶ **Much remains to be done . . .**

Thank you very much for your attention!

Hartelijk dank!

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