Numerical Analysis and Computing

Lecture Notes #4 — Solutions of Equations in One Variable, Interpolation and Polynomial Approximation — Accelerating Convergence; Zeros of Polynomials; Deflation; Müller's Method; Lagrange Polynomials; Neville's Method

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Introduction

"[...] it is rare to have the luxury of quadratic convergence."

(Burden-Faires, p.86 9th)

There are a number of methods for squeezing faster convergence out of an **already computed sequence** of numbers.

We here explore one method which seems the have been around since the beginning of numerical analysis... **Aitken's \Delta^2 method**. It can be used to accelerate convergence of a sequence that is linearly convergent, regardless of its origin or application.

A review of extrapolation methods can be found in:

"Practical Extrapolation Methods: Theory and Applications," Avram Sidi, Number 10 in Cambridge Monographs on Applied and Computational Mathematics, Cambridge University Press, June 2003. ISBN: 0-521-66159-5

Recall: Convergence of a Sequence

Definition

Suppose the sequence $\{p_n\}_{n=0}^{\infty}$ converges to p, with $p_n \neq p$ for all n. If positive constants λ and α exists with

$$\lim_{n\to\infty}\frac{|p_{n+1}-p|}{|p_n-p|^{\alpha}}=\lambda,$$

then $\{p_n\}_{n=0}^{\infty}$ converges to p of order α , with asymptotic error constant λ .

Linear convergence means that $\alpha = 1$, and $\lambda \in (0, 1)$.

Assume $\{p_n\}_{n=0}^{\infty}$ is a **linearly convergent sequence** with limit p.

Further, assume we are far out into the tail of the sequence (n large), and the signs of the successive errors agree, i.e.

$$\operatorname{sign}(p_n-p)=\operatorname{sign}(p_{n+1}-p)=\operatorname{sign}(p_{n+2}-p)=\ldots$$

so that

$$\frac{p_{n+2}-p}{p_{n+1}-p}\approx\frac{p_{n+1}-p}{p_n-p}\approx\lambda\quad\text{(the asymptotic limit)}.$$

This would indicate

$$(p_{n+1}-p)^2 \approx (p_{n+2}-p)(p_n-p),$$

 $p_{n+1}^2 - 2p_{n+1}\mathbf{p} + \mathbf{p^2} \approx p_{n+2}p_n - (p_{n+2}+p_n)\mathbf{p} + \mathbf{p^2}.$

We solve for p and get...

 Π/Π

We solve for p and get...

$$p \approx \frac{p_{n+2}p_n - p_{n+1}^2}{p_{n+2} - 2p_{n+1} + p_n}.$$

A little bit of algebraic manipulation put this into the equivalent "classical" Aitken form:

$$\hat{p}_n = p = p_n - \frac{(p_{n+1} - p_n)^2}{p_{n+2} - 2p_{n+1} + p_n}.$$

Aitken's Δ^2 Method is based on the assumption that the \hat{p}_n we compute from p_{n+2} , p_{n+1} and p_n is a better approximation to the actual limit p.

The analysis needed to rigorously **prove** this is beyond the scope of this class, see *e.g.* Sidi's book.



Given a sequence finite $\{p_n\}_{n=0}^N$ or infinite $\{q_n\}_{n=0}^\infty$ sequence which converges linearly to some limit.

Define the new sequences

$$\hat{p}_n = p_n - \frac{(p_{n+1} - p_n)^2}{p_{n+2} - 2p_{n+1} + p_n}, \quad n = 0, 1, \dots, N - 2,$$

or

$$\hat{q}_n = q_n - \frac{(q_{n+1} - q_n)^2}{q_{n+2} - 2q_{n+1} + q_n}, \quad n = 0, 1, \dots, \infty.$$

Aitken's Δ^2 Method

Example

Consider the sequence $\{p_n\}_{n=0}^{\infty}$, where the sequence is generated by the fixed point iteration $p_{n+1} = \cos(p_n)$, $p_0 = 0$.

Iteration	p _n	ρ̂ _n
0	0.000000000000000	0.685073357326045
1	1.0000000000000000	0.7 28010361467617
2	0 .540302305868140	0.73 3665164585231
3	0.857553215846393	0.73 6906294340474
4	0 .654289790497779	0.73 8050421371664
5	0.7 93480358742566	0.73 8636096881655
6	0.7 01368773622757	0.73 8876582817136
7	0.7 63959682900654	0.73 8992243027034
8	0.7 22102425026708	0.7390 42511328159
9	0.7 50417761763761	0.7390 65949599941
10	0.73 1404042422510	0.7390 76383318956
11	0.7 44237354900557	0.73908 1177259563*
12	0.73 5604740436347	0.73908 3333909684*

Note: Bold digits are correct; \hat{p}_{11} needs p_{13} , and \hat{p}_{12} additionally needs p_{14} .



Faster Convergence for "Aitken-Sequences"

Theorem (Convergence of Aitken- Δ^2 -Sequences)

Suppose $\{p_n\}_{n=0}^{\infty}$ is a sequence that converges linearly to the limit p, and for n large enough we have $(p_n-p)(p_{n+1}-p)>0$. Then the Aitken-accelerated sequence $\{\hat{p}_n\}_{n=0}^{\infty}$ converges fast to p in the sense that

$$\lim_{n\to\infty}\left[\frac{\hat{p}_n-p}{p_n-p}\right]=0.$$

We can combine Aitken's method with fixed-point iteration in order to get a "fixed-point iteration on steroids." (or should that be Erythropoietin (EPO), or possibly Clenbuterol?!)



Steffensen's Method: Fixed-Point Iteration on Steroids

Suppose we have a (linearly converging) fixed point iteration:

$$p_0, p_1 = g(p_0), p_2 = g(p_1), \dots$$

Once we have p_0 , p_1 and p_2 , we can compute

$$\hat{p}_0 = p_0 - \frac{(p_1 - p_0)^2}{p_2 - 2p_1 + p_0}.$$

At this point we "restart" the fixed point iteration with $p_0 = \hat{p}_0$, e.g.

$$p_3 = \hat{p}_0, \quad p_4 = g(p_3), \quad p_5 = g(p_4),$$

and compute

$$\hat{p}_3 = p_3 - \frac{(p_4 - p_3)^2}{p_5 - 2p_4 + p_3}.$$

Steffensen's Method: The Quadratic, g-g-A, Waltz! Quadratic Convergence

Algorithm: Steffensen's Method

Input: Initial approximation p_0 ; tolerance TOL; maximum number of iterations N_0 .

Output: Approximate solution p, or failure message.

- 1. Set i = 1
- 2. While $i \le N_0$ do 3--6
- 3* Set $p_1 = g(p_0)$, $p_2 = g(p_1)$, $p = p_0 (p_1 p_0)^2/(p_2 2p_1 + p_0)$
- 4. If $|p-p_0| < TOL$ then
- 4a. output *p*
- 4b. stop program
- 5. Set i = i + 1
- 6. Set $p_0 = p$
- 7. Output: "Failure after N_0 iterations."

Steffensen's Method: Potential Breakage

3* If at some point $p_2-2p_1+p_0=0$ (which appears in the denominator), then we stop and select the current value of p_2 as our approximate answer.

Both Newton's and Steffensen's methods give quadratic convergence. In Newton's method we compute one function value and one derivative in each iteration. In Steffensen's method we have two function evaluations and a more complicated algebraic expression in each iteration, **but no derivative**. It looks like we got something for (almost) nothing.

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However, in order the guarantee quadratic convergence for Steffensen's method, the fixed point function g must be 3 times continuously differentiable, e.g. $f \in C^3[a,b]$, (see theorem-2.15 in Burden-Faires $^{9\text{th}}$). Newton's method "only" requires $f \in C^2[a,b]$ (BF $^{9\text{th}}$ theorem-2.6).

Aitken's Δ^2 and Steffensen's Methods

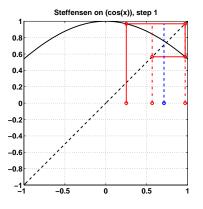
Example

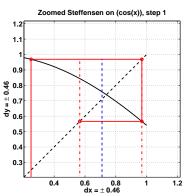
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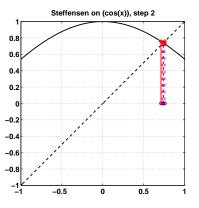
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9	0.7 50417761763761	0.7390 65949599941	0.7390851 06356719 ^s
10	0.73 1404042422510	0.7390 76383318956	0.7390851 51307330
11	0.7 44237354900557	0.73908 1177259563*	0.7390851 21028058
12	0.73 5604740436347	0.73908 3333909684*	0.739085133215161 s

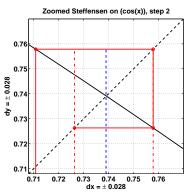
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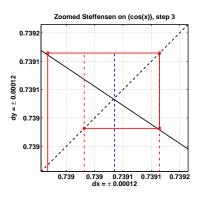


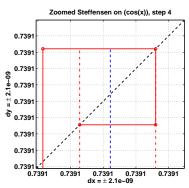












Zeros of Polynomials

Definition: Degree of a Polynomial

A **polynomial of degree** *n* has the form

$$P(x) = a_n x^n + a_{n-1} x^{n-1} + \dots + a_1 x + a_0, \quad a_n \neq 0$$

where the a_i 's are constants (either real, or complex) called the **coefficients** of P.

Why look at polynomials? — We'll be looking at the problem P(x) = 0 (i.e. f(x) = 0 for a special class of functions.)

Polynomials are the basis for many approximation methods, hence being able to solve polynomial equations fast is valuable.

We'd like to use Newton's method, so we need to compute P(x) and P'(x) as efficiently as possible.

Fundamentals

Theorem (The Fundamental Theorem of Algebra)

If P(x) is a polynomial of degree $n \ge 1$ with real or complex coefficients, then P(x) = 0 has at least one (possibly complex) root.

The proof is surprisingly(?) difficult and requires understanding of complex analysis... We leave it as an exercise for the motivated student!

Key Consequences of the Fundamental Theorem of Algebra

Corollary

If P(x) is a polynomial of degree $n \ge 1$ with real or complex coefficients then there exists unique constants x_1, x_2, \ldots, x_k (possibly complex) and unique positive integers m_1, m_2, \ldots, m_k such that $\sum_{i=1}^k m_i = n$ and

$$P(x) = a_n(x - x_1)^{m_1}(x - x_2)^{m_2} \cdots (x - x_k)^{m_k}$$

- The collection of zeros is unique.
- m_i are the multiplicities of the individual zeros.
- A polynomial of degree n has exactly n zeros, counting multiplicity.



Corollary

Let P(x) and Q(x) be polynomials of degree at most n. If x_1, x_2, \ldots, x_k , with k > n are **distinct** numbers with $P(x_i) = Q(x_i)$ for $i = 1, 2, \ldots, k$, then P(x) = Q(x) for all values of x.

— If two polynomials of degree n agree at at least (n+1) points, then they must be the same.

Let

$$P(x) = a_n x^n + a_{n-1} x^{n-1} + \cdots + a_1 x + a_0.$$

If we are looking to evaluate $P(x_0)$ for any x_0 , let

$$b_n = a_n$$
, $b_k = a_k + b_{k+1}x_0$, $k = (n-1), (n-2), \dots, 1, 0$,

then $b_0 = P(x_0)$. We have only used n multiplications and n additions.

At the same time we have computed the decomposition

$$P(x) = (x - x_0)Q(x) + b_0,$$

where

$$Q(x) = \sum_{k=0}^{n-1} b_{k+1} x^k.$$

Horner's Method: Evaluating Polynomials Quickly

Huh?!? Where did the expression come from? — Consider

$$P(x) = a_n x^n + a_{n-1} x^{n-1} + \dots + a_1 x + a_0$$

$$= (a_n x^{n-1} + a_{n-1} x^{n-2} + \dots + a_1) x + a_0$$

$$= ((a_n x^{n-2} + a_{n-1} x^{n-3} + \dots) x + a_1) x + a_0$$

$$= \underbrace{(\dots ((a_n x^{n-2} + a_{n-1}) x + \dots) x + a_1) x + a_0}_{b_{n-1}}$$

Horner's method (first published by Theophilus Holdred(!) in 1820) is "simply" the computation of this parenthesized expression from the inside-out...

Now, if we need to compute $P'(x_0)$ we have

$$P'(x)\Big|_{x=x_0}=(x-x_0)Q'(x)+Q(x)\Big|_{x=x_0}=Q(x_0)$$

Which we can compute (again using Horner's method) in (n-1) multiplications and (n-1) additions.

Proof? We really ought to prove that Horner's method works. It basically boils down to lots of algebra which shows that the coefficients of P(x) and $(x - x_0)Q(x) + b_0$ are the same...

A couple of examples may be more instructive...

For
$$P(x) = x^4 - x^3 + x^2 + x - 1$$
, compute $P(5)$:
 $x_0 = 5$ $\begin{vmatrix} a_4 = 1 & a_3 = -1 & a_2 = 1 & a_1 = 1 & a_0 = -1 \\ b_4 x_0 = 5 & b_3 x_0 = 20 & b_2 x_0 = 105 & \mathbf{b_1 x_0} = \mathbf{530} \end{vmatrix}$
 $b_4 = 1$ $b_3 = 4$ $b_2 = 21$ $b_1 = 106$ $\mathbf{b_0} = \mathbf{529}$

Hence, P(5) = 529, and

$$P(x) = (x - 5)(x^3 + 4x^2 + 21x + 106) + 529$$

Similarly we get P'(5) = Q(5) = 436



Algorithm: Horner's Method

Algorithm: Horner's Method

Input: Degree n; coefficients a_0, a_1, \ldots, a_n ; x_0

Output:
$$y = P(x_0), z = P'(x_0).$$

- 1. Set $y = a_n$, $z = a_n$
- 2. For $j = (n-1), (n-2), \dots, 1$ Set $y = x_0y + a_j, z = x_0z + y$
- 3. Set $y = x_0y + a_0$
- 4. Output (y, z)
- 5. End program



Deflation — Finding All the Zeros of a Polynomial

If we are solving our current favorite problem

$$P(x) = 0$$
, $P(x)$ a polynomial of degree n ,

and we are using Horner's method of computing $P(x_i)$ and $P'(x_i)$, then after N iterations, x_N is an approximation to one of the roots of P(x) = 0.

We have

$$P(x) = (x - x_N)Q(x) + b_0, \quad b_0 \approx 0.$$

At this point, let $\hat{r}_1 = x_N$ be the first root, and $Q_1(x) = Q(x)$.

We can now find a second root by applying Newton's method to $Q_1(x)$.



Deflation — Finding All the Zeros of a Polynomial

After some number of iterations of Newton's method we have

$$Q_1(x) = (x - \hat{r}_2)Q_2(x) + b_0^{(2)}, \quad b_0^{(2)} \approx 0$$

If P(x) is an n^{th} -degree polynomial with n real roots, we can apply this procedure (n-2) times to find (n-2) approximate zeros of P(x): \hat{r}_1 , \hat{r}_2 , ..., \hat{r}_{n-2} , and a quadratic factor $Q_{n-2}(x)$.

At this point we can solve $Q_{n-2}(x) = 0$ using the quadratic formula, and we have n roots of P(x) = 0.

This procedure is called **Deflation**.

Quality of Deflation

Now, the big question is "are the approximate roots \hat{r}_1 , \hat{r}_2 , ..., \hat{r}_n good approximations of the roots of P(x)??"

Unfortunately, sometimes, no.

In each step we solve the equation to some tolerance, i.e.

$$|b_0^{(k)}| < tol$$

Even though we may solve to a tight tolerance (10^{-8}) , the errors accumulate and the inaccuracies increase iteration-by-iteration...

Question: Is deflation therefore useless???



Improving the Accuracy of Deflation

The problem with deflation is that the zeros of $Q_k(x)$ are not good representatives of the zeros of P(x), especially for high k's.

As k increases, the quality of the root \hat{r}_k decreases.

Maybe there is a way to get all the zeros with the same quality?

Improving the Accuracy of Deflation

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As k increases, the quality of the root \hat{r}_k decreases.

Maybe there is a way to get all the zeros with the same quality?

The idea is quite simple... in each step of deflation, instead of just accepting \hat{r}_k as a root of P(x), we re-run Newton's method on the **full polynomial** P(x), with \hat{r}_k as the starting point — a couple of Newton iterations should quickly converge to the root of the full polynomial.

Improved Deflation — Algorithm Outline

Algorithm Outline: Improved Deflation

- 1. Apply Newton's method to $P(x) \rightarrow \hat{\mathbf{r}}_1, \mathbf{Q}_1(\mathbf{x})$.
- 2. For k = 2, 3, ..., (n-2) do 3--4
- 3. Apply Newton's method to $\mathbf{Q}_{\mathbf{k}-\mathbf{1}} \ \to \ \hat{\mathbf{r}}_{\mathbf{k}}^*, \ \mathbf{Q}_{\mathbf{k}}^*(\mathbf{x}).$
- 4. Apply Newton's method to P(x) with \hat{r}_k^* as the initial point $\to \hat{r}_k$ Apply Horner's method to $Q_{k-1}(x)$ with $x=\hat{r}_k \to Q_k(x)$
- 5. Use the quadratic formula on $\mathbf{Q}_{n-2}(\mathbf{x})$ to get the two remaining roots.

Note: "Inside" Newton's method, the evaluations of polynomials and their derivatives are also performed using Horner's method.



Deflation & Improvement

Wilkinson Polynomials

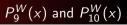
The Wilkinson Polynomials

$$P_n^{\mathsf{W}}(x) = \prod_{k=1}^n (x-k)$$

have the roots $\{1, 2, ..., n\}$, but provide surprisingly tough numerical root-finding problems. (Additional details in Math 543.)

In the next few slides we show the results of Deflation and Improved Deflation applied to Wilkinson polynomials of degree 9, 10, 12, and 13.

Deflation & Improvement



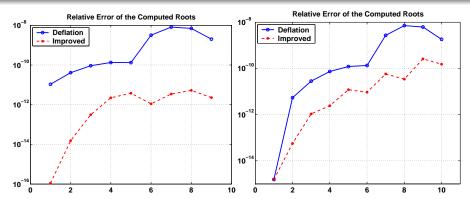
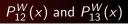


Figure: [LEFT] The result of the two algorithms on the Wilkinson polynomial of degree 9; in this case the roots are computed so that $|b_0^{(k)}| < 10^{-6}$. [RIGHT] The result of the two algorithms on the Wilkinson polynomial of degree 10; in this case the roots are computed so that $|b_0^{(k)}| < 10^{-6}$. In both cases the lower line corresponds to improved deflation and we see that we get an improvement in the relative error of several orders of magnitude.

Deflation & Improvement



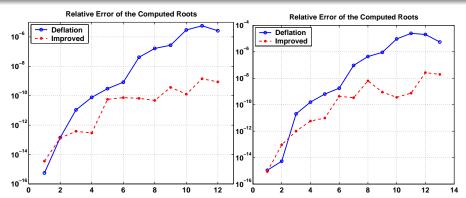


Figure: [LEFT] The result of the two algorithms on the Wilkinson polynomial of degree 12; in this case the roots are computed so that $|b_0^{(k)}| < 10^{-4}$. [RIGHT] The result of the two algorithms on the Wilkinson polynomial of degree 13; in this case the roots are computed so that $|b_0^{(k)}| < 10^{-3}$. In both cases the lower line corresponds to improved deflation and we see that we get an improvement in the relative error of several orders of magnitude.

What About Complex Roots???

One interesting / annoying feature of polynomials with real coefficients is that they may have complex roots, *e.g.* $P(x) = x^2 + 1$ has the roots $\{-i, i\}$. Where by definition $i = \sqrt{-1}$.

If the initial approximation given to Newton's method is real, all the successive iterates will be real... which means we will not find complex roots.

One way to overcome this is to start with a complex initial approximation and do all the computations in complex arithmetic.

Another solution is Müller's Method...



Müller's Method

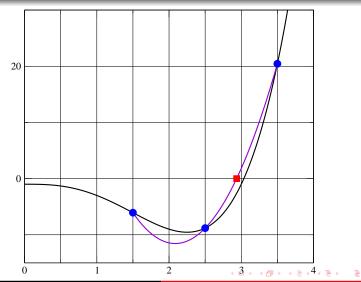
Müller's method is an extension of the Secant method...

Recall that the secant method uses two points x_k and x_{k-1} and the function values in those two points $f(x_k)$ and $f(x_{k-1})$. The zero-crossing of the linear interpolant (the secant line) is used as the next iterate x_{k+1} .

Müller's method takes the next logical step: it uses **three points**: x_k , x_{k-1} and x_{k-2} , the function values in those points $f(x_k)$, $f(x_{k-1})$ and $f(x_{k-2})$; a second degree polynomial fitting these three points is found, and its zero-crossing is the next iterate x_{k+1} .

Next slide: $f(x) = x^4 - 3x^3 - 1$, $x_{k-2} = 1.5$, $x_{k-1} = 2.5$, $x_k = 3.5$.

Müller's Method — Illustration $f(x) = x^4 - 3x^3 - 1$



Müller's Method — Fitting the Quadratic Polynomial

We consider the quadratic polynomial

$$m(x) = a(x - x_k)^2 + b(x - x_k) + c$$

at the three fitting points we get

$$f(x_{k-2}) = a(x_{k-2} - x_k)^2 + b(x_{k-2} - x_k) + c$$

$$f(x_{k-1}) = a(x_{k-1} - x_k)^2 + b(x_{k-1} - x_k) + c$$

$$f(x_k) = c$$

We can solve for a, b, and c:

$$a = \frac{(x_{k-1} - x_k)(f(x_{k-2}) - f(x_k)) - (x_{k-2} - x_k)(f(x_{k-1}) - f(x_k))}{(x_{k-2} - x_k)(x_{k-1} - x_k)(x_{k-2} - x_{k-1})}$$

$$b = \frac{(x_{k-2} - x_k)^2(f(x_{k-1}) - f(x_k)) - (x_{k-1} - x_k)^2(f(x_{k-2}) - f(x_k))}{(x_{k-2} - x_k)(x_{k-1} - x_k)(x_{k-2} - x_{k-1})}$$

$$c = f(x_k)$$

Müller's Method — Identifying the Zero

We now have a quadratic equation for $(x - x_k)$ which gives us two possibilities for x_{k+1} :

$$x_{k+1} - x_k = \frac{-2c}{b \pm \sqrt{b^2 - 4ac}}$$

In Müller's method we select

$$x_{k+1} = x_k - \frac{2c}{b + \text{sign}(b)\sqrt{b^2 - 4ac}}$$

we are maximizing the (absolute) size of the denominator, hence we select the root closest to x_k .

Note that if $b^2 - 4ac < 0$ then we automatically get complex roots.

Müller's Method — Algorithm

Algorithm: Müller's Method

Input: x_0 , x_1 , x_2 ; tolerance tol; max iterations N_0

Output: Approximate solution p, or failure message.

- 1. Set $h_1 = (x_1 x_0)$, $h_2 = (x_2 x_1)$, $\delta_1 = [f(x_1) f(x_0)]/h_1$, $\delta_2 = [f(x_2) f(x_1)]/h_2$, $d = (\delta_2 \delta_1)/(h_2 + h_1)$, j = 3.
- 2. While $j \le N_0$ do 3--7
- 3. $b = \delta_2 + h_2 d$, $D = \sqrt{b^2 4f(x_2)d}$ complex?
- 4. If |b-D| < |b+D| then set E = b+D else set E = b-D
- 5. Set $h = -2f(x_2)/E$, $p = x_2 + h$
- 6. If |h| < tol then output p; stop program
- 7. Set $x_0 = x_1$, $x_1 = x_2$, $x_2 = p$, $h_1 = (x_1 x_0)$, $h_2 = (x_2 x_1)$, $\delta_1 = [f(x_1) f(x_0)]/h_1$, $\delta_2 = [f(x_2) f(x_1)]/h_2$, $d = (\delta_2 \delta_1)/(h_2 + h_1)$, j = j + 1
- 8. output "Müller's Method failed after N_0 iterations."

Now We Know "Everything" About Solving f(x) = 0!?

Let's recap... Things to remember...

The relation between **root finding** (f(x) = 0) and **fixed point** (g(x) = x).

Key algorithms for root finding: Bisection, Secant Method, and **Newton's Method**. — Know what they are (the updates), how to start (one or two points? bracketing or not bracketing the root?), can the method break, can breakage be fixed? Convergence properties.

Also, know the mechanics of the Regula Falsi method, and understand why it can run into trouble.

Fixed point iteration: Under what conditions do FP-iteration converge for all starting values in the interval?

Recap, continued...

Basic error analysis: order α , asymptotic error constant λ . — Which one has the most impact on convergence? Convergence rate for general fixed point iterations?

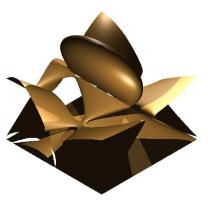
Multiplicity of zeros: What does it mean? How do we use this knowledge to "help" Newton's method when we're looking for a zero of high multiplicity?

Convergence acceleration: Aitken's Δ^2 -method. Steffensen's Method.

Zeros of polynomials: Horner's method, Deflation (with improvement), Müller's method.

New Favorite Problem:

Interpolation and Polynomial Approximation



Weierstrass Approximation Theorem

The following theorem is the basis for polynomial approximation:

Theorem (Weierstrass Approximation Theorem)

Suppose
$$f \in C[a, b]$$
. Then $\forall \epsilon > 0 \exists a \text{ polynomial } P(x) : |f(x) - P(x)| < \epsilon, \forall x \in [a, b].$

Note: The bound is *uniform*, *i.e.* valid for all x in the interval.

Note: The theorem says nothing about how to find the polynomial, or about its order.

Illustrated: Weierstrass Approximation Theorem

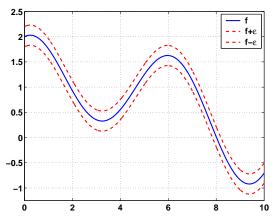


Figure: Weierstrass approximation Theorem guarantees that we (maybe with substantial work) can find a polynomial which fits into the "tube" around the function f, no matter how thin we make the tube.

Candidates: the Taylor Polynomials???

Natural Question:

Are our old friends, the Taylor Polynomials, good candidates for polynomial interpolation?

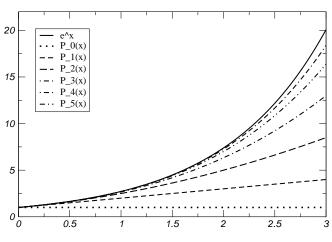
Answer:

No. The Taylor expansion works very hard to be accurate in the neighborhood of *one point*. But we want to fit data at many points (in an extended interval).

[Next slide: The approximation is great near the expansion point $x_0 = 0$, but get progressively worse at we get further away from the point, even for the higher degree approximations.]



Taylor Approximation of e^x on the Interval [0, 3]



Lookahead: Polynomial Approximation

Clearly, Taylor polynomials are not well suited for approximating a function over an **extended** interval.

We are going to look at the following:

- Lagrange polynomials Neville's Method. [This Lecture]
- Newton's divided differences.
- Hermite interpolation.
- Cubic splines Piecewise polynomial approximation.
- (Parametric curves)
- (Bézier curves)



Interpolation: Lagrange Polynomials

Idea: Instead of working hard at *one point*, we will prescribe a number of points through which the polynomial must pass.

As warm-up we will define a function that passes through the points $(x_0, f(x_0))$ and $(x_1, f(x_1))$. First, lets define

$$L_0(x) = \frac{x - x_1}{x_0 - x_1}, \quad L_1(x) = \frac{x - x_0}{x_1 - x_0},$$

and then define the interpolating polynomial

$$P(x) = L_0(x)f(x_0) + L_1(x)f(x_1),$$

then
$$P(x_0) = f(x_0)$$
, and $P(x_1) = f(x_1)$.

- P(x) is the unique linear polynomial passing through $(x_0, f(x_0))$ and $(x_1, f(x_1))$.

An *n*-degree polynomial passing through n+1 points

We are going to construct a polynomial passing through the points $(x_0, f(x_0)), (x_1, f(x_1)), (x_2, f(x_2)), \ldots, (x_N, f(x_n)).$

We define $L_{n,k}(x)$, the **Lagrange coefficients:**

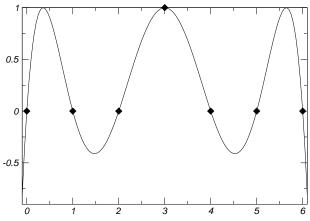
$$L_{n,k}(x) = \prod_{i=0, i \neq k}^{n} \frac{x - x_i}{x_k - x_i} = \frac{x - x_0}{x_k - x_0} \cdots \frac{x - x_{k-1}}{x_k - x_{k-1}} \cdot \frac{x - x_{k+1}}{x_k - x_{k+1}} \cdots \frac{x - x_n}{x_k - x_n},$$

which have the properties

$$L_{n,k}(x_k) = 1; \quad L_{n,k}(x_i) = 0, \ \forall i \neq k.$$



Example of $L_{n,k}(x)$



This is $L_{6,3}(x)$, for the points $x_i = i, i = 0, ..., 6$.

The nth Lagrange Interpolating Polynomial

We use $L_{n,k}(x)$, k = 0, ..., n as building blocks for the Lagrange interpolating polynomial:

$$P(x) = \sum_{k=0}^{n} f(x_k) L_{n,k}(x),$$

which has the property

$$P(x_i) = f(x_i), \quad \forall i = 0, \ldots, n.$$

This is the unique polynomial passing through the points $(x_i, f(x_i)), i = 0, ..., n$.



Error bound for the Lagrange interpolating polynomial

Suppose x_i , i = 0, ..., n are distinct numbers in the interval [a, b], and $f \in C^{n+1}[a, b]$. Then $\forall x \in [a, b] \exists \xi(x) \in (a, b)$ so that:

$$f(x) = P_{Lagrange}(x) + \frac{f^{(n+1)}(\xi(x))}{(n+1)!} \prod_{i=0}^{n} (x - x_i),$$

where $P_{Lagrange}(x)$ is the n^{th} Lagrange interpolating polynomial.

Compare with the error formula for Taylor polynomials

$$f(x) = P_{\mathsf{Taylor}}(x) + \frac{f^{(n+1)}(\xi(x))}{(n+1)!} (x - x_0)^{n+1},$$

Problem: Applying the error term may be difficult...



The Lagrange and Taylor Error Terms

Just to get a feeling for the non-constant part of the error terms in the Lagrange and Taylor approximations, we plot those parts on the interval [0,4] with interpolation points $x_i = i, i = 0, 1, ..., 4$:

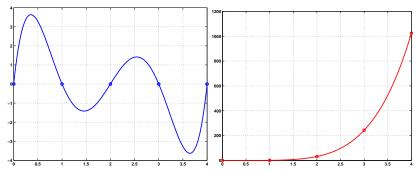


Figure: [LEFT] The non-constant error terms for the Lagrange interpolation oscillates in the interval [-4, 4] (and takes the value zero at the node point x_k), and [RIGHT] the non-constant error term for the Taylor extrapolation grows in the interval [0, 1024].

Practical Problems

Applying (estimating) the error term is difficult.

The degree of the polynomial needed for some desired accuracy is not known until after cumbersome calculations — checking the error term.

If we want to increase the degree of the polynomial (to e.g. n+1) the previous calculations are not of any help...

Building block for a fix: Let f be a function defined at x_0, \ldots, x_n , and suppose that m_1, m_2, \ldots, m_k are k (< n) distinct integers, with $0 \le m_i \le n \ \forall i$. The Lagrange polynomial that agrees with f(x) the k points $x_{m_1}, x_{m_2}, \ldots, x_{m_k}$, is denoted $P_{m_1, m_2, \ldots, m_k}(x)$.

Note: $\{m_1, m_2, \dots, m_k\} \subset \{0, 1, \dots, n\}.$

Increasing the degree of the Lagrange Interpolation

Theorem

Let f be defined at x_0, x_1, \ldots, x_k , and x_i and x_j be two distinct points in this set, then

$$P(x) = \frac{(x - x_j)P_{0,\dots,j-1,j+1,\dots,k}(x) - (x - x_i)P_{0,\dots,i-1,i+1,\dots,k}(x)}{x_i - x_j}$$

is the k^{th} Lagrange polynomial that interpolates f at the k+1 points x_0, \ldots, x_k .

Recursive Generation of Higher Degree Lagrange Interpolating Polynomials

-X ₀	P_0				
x_1	P_1	$P_{0,1}$			
<i>X</i> ₂	P_2	$P_{1,2}$	$P_{0,1,2}$		
<i>X</i> 3	P_3	$P_{2,3}$	$P_{1,2,3}$	$P_{0,1,2,3}$	
<i>X</i> 4	P_4	$P_{3,4}$	$P_{2,3,4}$	$P_{1,2,3,4}$	P _{0,1,2,3,4}

Neville's Method

The notation in the previous table gets cumbersome... We introduce the notation $Q_{Last,Degree} = P_{Last-Degree,...,Last}$, the table becomes:

Compare with the old notation:

Algorithm: Neville's Method — Iterated Interpolation

Algorithm: Neville's Method

To evaluate the polynomial that interpolates the n+1 points $(x_i, f(x_i))$, i = 0, ..., n at the point x:

- 1. Initialize $Q_{i,0} = f(x_i)$.
- 2.

FOR
$$i=1:n$$
FOR $j=1:i$

$$Q_{i,j}=\frac{(x-x_{i-j})Q_{i,j-1}-(x-x_i)Q_{i-1,j-1}}{x_i-x_{i-j}}$$
END

3. Output the Q-table.

END

Homework #3

http://webwork.sdsu.edu

- Will open on 09/12/2014 at 09:30am PDT.
- Will close no earlier than 09/24/2014 at 09:00pm PDT.