

FWAM Session B: Function Approximation and Differential Equations

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Wed, 10/30/19

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LECTURE 1: interpolation, integration, spectral methods

Motivations

exact func. $f(x)$ described by ∞ number of points

how handle approximately (but accurately) in computer, using least cost (bytes)?

- Interpolation: cheap but accurate look-up table for expensive $f(x)$
data fitting: given non-noisy data $f(x_i)$ at some x_i , model $f(x)$ at other points x ?

Contrast: fit noisy data = learning (pdf for) params in model, via likelihood/prior

- (Numerical) integration:
eg computing expectation values given a pdf

Contrast: Monte Carlo (random, high-dim.) integration, Thurs am

- Differentiation:
get gradient ∇f in order to optimize or
- Spectral (often Fourier) methods:

If $f(x)$ is smooth, handle very accurately without much extra cost

Deterministic (non-random) methods.

Integr/diff crucial for numerical ODEs and PDEs

topic of

Goals LECTURE I

TODO

teach range of practical methods focusing on 1D

pointers to dimensions $d > 1$

concepts:

convergence order

spectral methods

local vs global

adaptivity

rounding error

interpolation = func. representation, key to all else

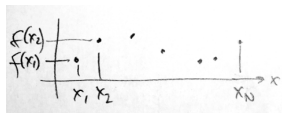
Interpolation in 1D ($d = 1$)

Say $y_j = f(x_j)$ known at nodes $\{x_j\}$

exact data, not noisy

want interpolant $\tilde{f}(x)$, s.t. $\tilde{f}(x_j) = y_j$

N -pt "grid"

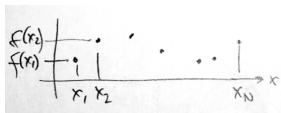


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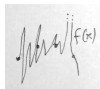
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hopeless w/o assumptions on f , eg smoothness, otherwise...

- extra info helps, eg f periodic, or $f(x) = \text{smooth} \cdot |x|^{-1/2}$

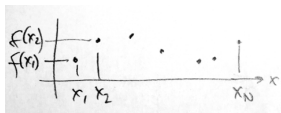


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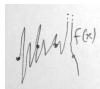
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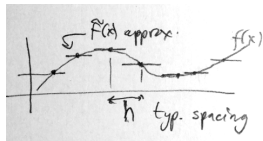


Simplest: use value at x_j nearest to x

"snap to grid"

Error $\max_x |\tilde{f}(x) - f(x)| = \mathcal{O}(h)$ as $h \rightarrow 0$

holds if f' bounded; can be nonsmooth but not crazy



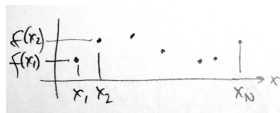
Recall notation " $\mathcal{O}(h)$ ": exists $C, h_0 > 0$ s.t. error $\leq Ch$ for all $h < h_0$

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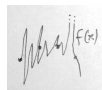
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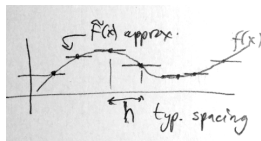


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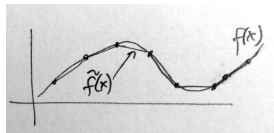
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Piecewise linear:

"connect the dots"

max error = $\mathcal{O}(h^2)$ as $h \rightarrow 0$

needs f'' bounded, ie smoother than before



Message: a higher order method is only higher order if f smooth enough

Interlude: convergence rates

Should know or measure convergence rate of any method you use

- “effort” parameter N eg # grid-points = $1/h^d$ where h = grid spacing, d = dim

We just saw algebraic conv. error = $\mathcal{O}(N^{-p})$, for order $p = 1, 2$

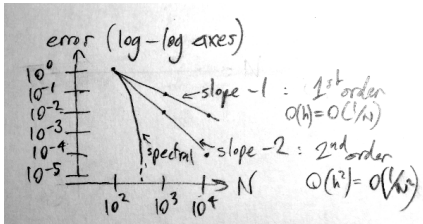
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Is only one graph in numerical analysis: “relative error vs effort”



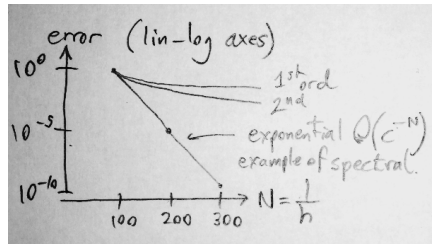
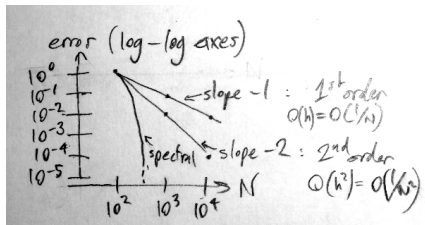
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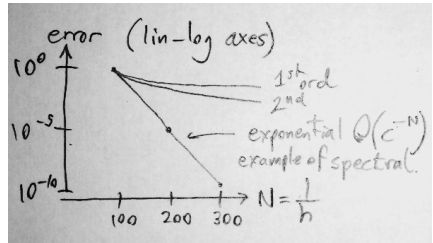
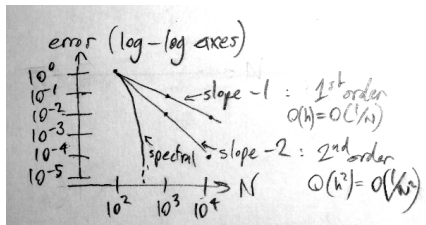
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Note how spectral gets many digits for small N

crucial for eg 3D prob.

“spectral” = “superalgebraic”, $\mathcal{O}(N^{-k})$ for any k

- how many digits to you want? for 1-digit (10% error), low order ok, easier to code

<rant> test your code w/ *known exact soln* to check error conv. <\rant>

What is the prefactor C in error $\leq Ch^k$? Has asymp. rate even kicked in yet? :)

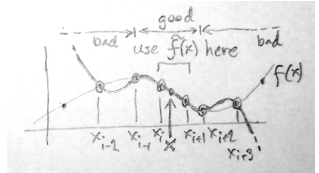
Higher-order interpolation for smooth f : the local idea

For any target x , use only set of nearest p nodes:

Exists unique degree- $(p-1)$ poly, $\sum_{k=0}^{p-1} c_k x^k$
which matches local data $(x_j, y_j)_{j=1}^p$

generalizes piecewise lin. idea

do **not** eval poly outside its central region!

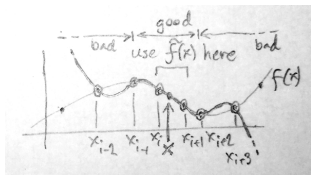


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if must have cont, recommend splines, eg cubic $p=3$: $\tilde{f} \in C^2$, meaning \tilde{f}'' is cont.

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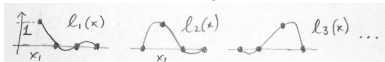
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How to find the degree- $(k-1)$ poly?

1) Crafty: solve square lin sys for coeffs $\sum_{k < p} x_j^k c_k = y_j$ $j = 1, \dots, p$
ie $V\mathbf{c} = \mathbf{y}$ $V = \text{"Vandermonde" matrix, is ill-cond. but works}$

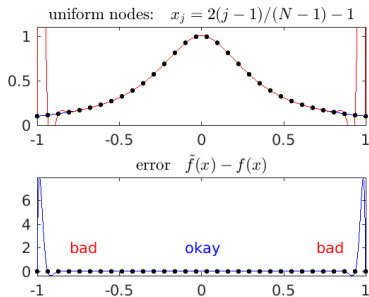
2) Traditional: barycentric formula $\tilde{f}(x) = \frac{\sum_{j=1}^p \frac{y_j}{x-x_j} w_j}{\sum_{j=1}^p \frac{1}{x-x_j} w_j}$ $w_j = \frac{1}{\prod_{i \neq j} (x_j - x_i)}$
[Tre13, Ch. 5]

Either way, $\tilde{f}(x) = \sum_{j=1}^p y_j \ell_j(x)$ where $\ell_j(x)$ is j th Lagrange basis func:



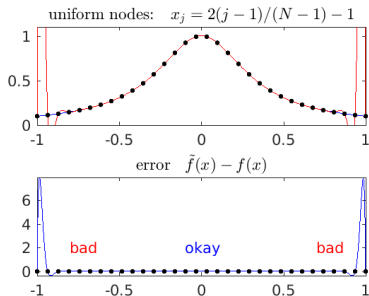
Global polynomial (Lagrange) interpolation?

Want increase order p . Use *all* data, get single $\tilde{f}(x)$, so $p = N$? “global”
 $p = N = 32$, smooth (analytic) $f(x) = \frac{1}{1+9x^2}$ on $[-1, 1]$: (Runge 1901)



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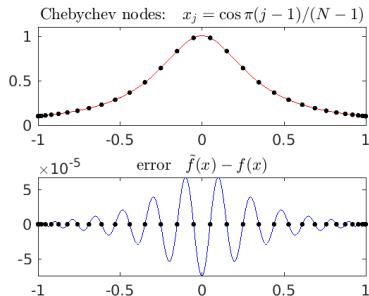
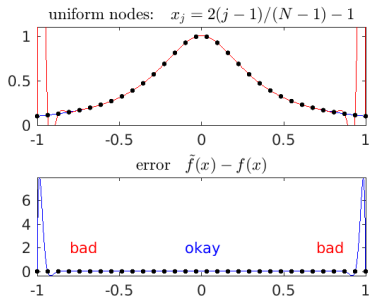
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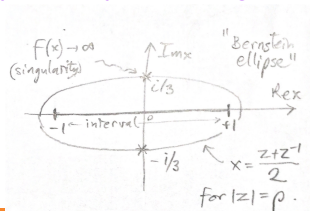
But exists good choice of nodes...

“Chebyshev”: means non-unif. grid density $\sim \frac{1}{\sqrt{1-x^2}}$

- our first spectral method

$\max \text{err} = \mathcal{O}(\rho^{-N})$ exponential conv!

$\rho > 1$ “radius” of largest ellipse in which f analytic



Node choice and adaptivity

Recap poly approx. $f(x)$ on $[a, b]$: are good & bad node sets $\{x_j\}_{j=1}^N$

Question: Do you get to *choose* the set of nodes at which f known?

- No: data fitting applications (or noisy variants: kriging, Gaussian processes, etc)
use local poly (central region only!), or something stable (splines)
- Yes: almost all else, interp., quadrature, PDE solvers so pick good nodes!

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Adaptivity idea global is inefficient if f smooth in most places, structured in a few

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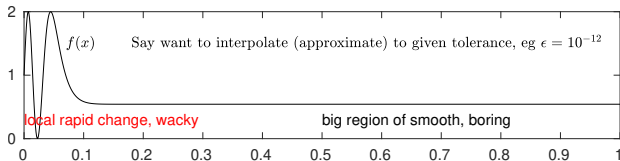
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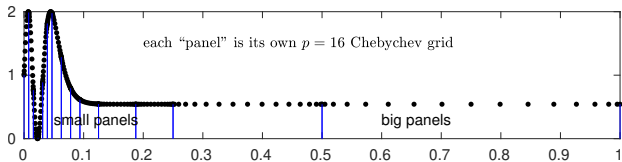
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automatically split
(recursively) panels
until $\max \text{err} \leq \epsilon$

via tests for local error

1D adaptive interpolator codes to try:

- [github/dbstein/function_generator](#) py+numba, fast (Stein '19)
- [chebfun for MATLAB](#) big- p Cheb. grids can exploit FFTs! (Trefethen et al.)

App.: replace nasty expensive $f(x)$ by cheap one!

optimal "look-up table"

Interpolation of periodic functions

Just did f on intervals $[a, b]$. Interpolation (& integration, etc.) of *periodic* f not same!

Periodic: $f(x + 2\pi) = f(x)$ for all x , $f(x) = \sum_{n \in \mathbb{Z}} \hat{f}_k e^{ikx}$ Fourier series

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instead of poly's, use **truncated** series $\tilde{f}(x) = \sum_{|k| < N/2} c_k e^{ikx}$ "trig. poly"

What's best you can do? if \hat{f}_n decays rapidly,

How get c_n ? unif. grid now good! Use FFT of vals $f(2\pi j/N)$,
 $j = 1, \dots, N$

$c_n =$

exp conv if f analytic in strip.

what if f not as smooth as this?

reapse to conv. order given by smoothness of f

why? \hat{f}_n decay

Messages: spectral conv usually requires understanding f off the real axis!

smooth = rapid Fourier series decay = rapid conv

Interpolation in $d > 1$

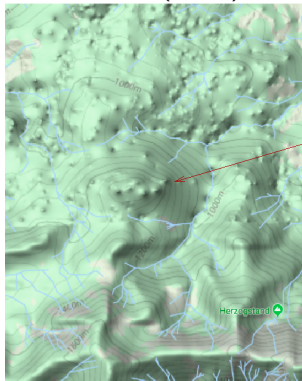
can you choose the points?

data fit: often no

PDE solve: usually yes - product grids, piecewise (FEM), etc

in $d > 1$ getting $f(\mathbf{x})$ from scattered data $\{\mathbf{x}_i\}$ is hard

Eg google terrain: $f(\mathbf{x})$ rough \rightarrow v low ord



height $f(\mathbf{x})$
interp from
unstructured
points in 2D,
kernel method

pock-marks!

interp from
Cartesian grid,
more accurate

unless you know f smooth, eg fit local multivariate polynomial
kriging

Numerical integration

Usually the user gets to choose the nodes x_j

Once have interpolant \tilde{f} from data $f(x_j)$, can *integrate it exactly*

“interpolatory quadrature”

Eg: piecewise linear gives composite trap rule $\mathcal{O}(N^{-2})$

periodic spectral gives periodic trap rule $\mathcal{O}(c^{-N})$ if analytic

TO DO

extrapolation

Rounding error [GC12, Ch. 5–6]

LECTURE II: numerical differential equations

For now we start with “elliptic”: time-independent problems

Motivations

eg steady-state (equilibrium) diffusion of a chemical

eg what electric potential caused by bunch of charges surrounded by H_2O ? (protein electrostatics)

Find u solving $\Delta u = f$, f = volume source term

Δ means Laplacian $\partial^2/\partial x^2 + \partial^2/\partial y^2 + \dots$ Δu is curvature of u

plus some BCs on u

eg viscous fluid flow: \mathbf{u} is velocity field, sat Stokes eqns

eg what is ground state of quantum system, solving $\Delta u = Eu$

Mike will in next talk overview this and 2 other flavors of PDE

References

- A Greenbaum and T P Chartier, *Numerical methods*, Princeton University Press, 2012.
- L. N. Trefethen, *Approximation theory and approximation practice*, SIAM, 2013, <http://www.maths.ox.ac.uk/chebfun/ATAP>.