

# An introduction to fractional calculus

Fundamental ideas and numerics

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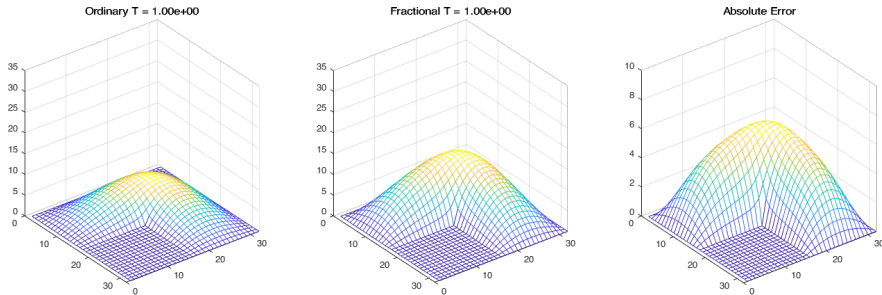
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May, 2022



# Subdiffusion equations

At the end of the last lecture we had observed the following behavior:



for the solution of:

$${}_C D_t^\alpha u = 0.05 \nabla^2 u, \quad \alpha = 0.3, 1.$$

The **visual effect** seemed to be a **slowing down of the diffusion**.

# Brownian motion (Metzler and Klafter 2000)

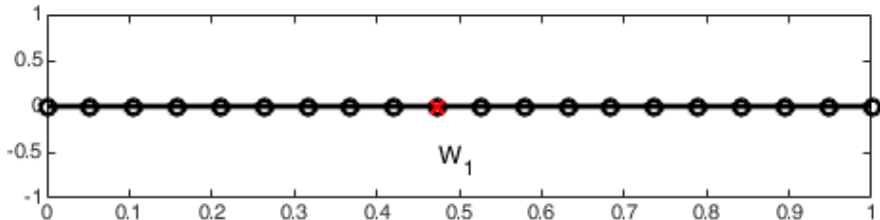
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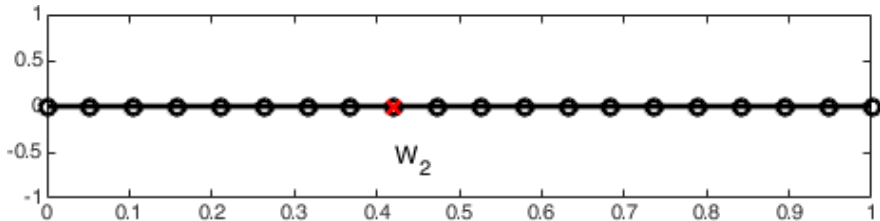
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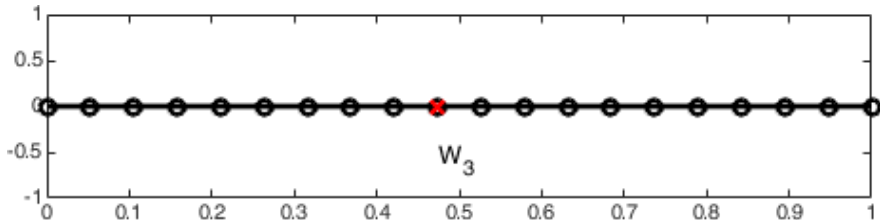
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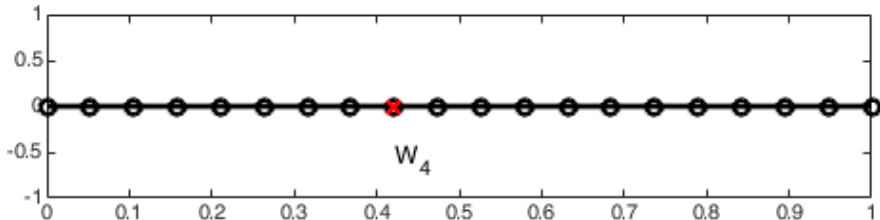
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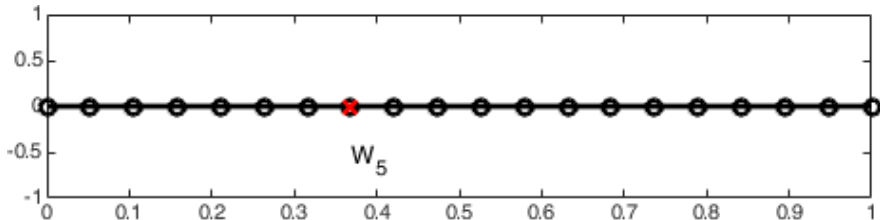
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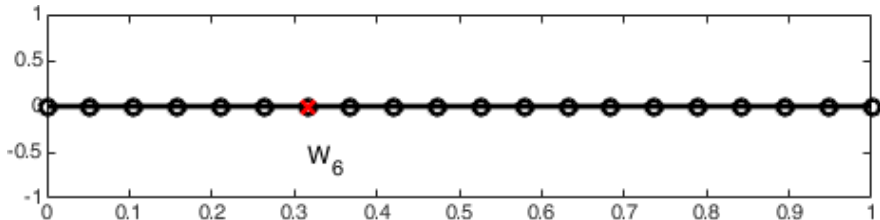




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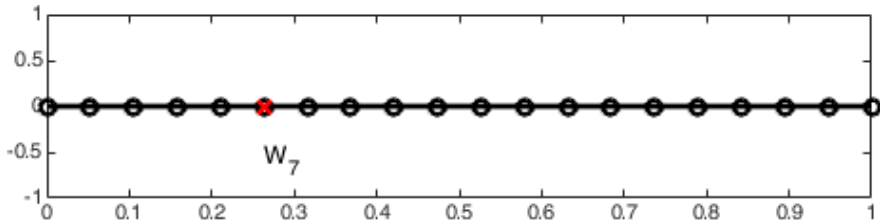
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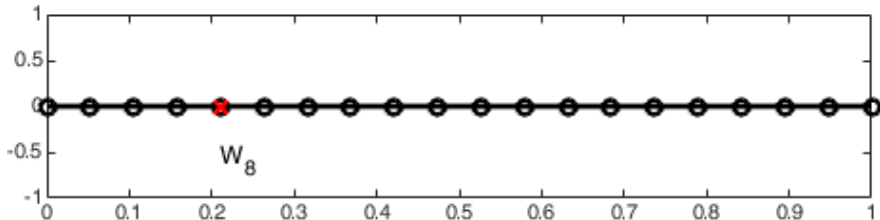
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- The prefactor  $1/2$  tells us that the **process is isotropic** with respect to the left/right direction.
- If we let  $\Delta t \rightarrow 0$ ,  $\Delta x \rightarrow 0$  and do a **Taylor expansion** in both  $\Delta$  and  $\Delta x$  we get

$$W_j(t + \Delta t) = W_j(t) + \Delta t \frac{\partial W_j}{\partial t} + O([\Delta t]^2), \quad \text{for } \Delta t \rightarrow 0,$$

$$W_{j\pm 1}(t) = W(x, t) \pm \Delta x \frac{\partial W}{\partial x} + \frac{(\Delta x)^2}{2} \frac{\partial^2 W}{\partial x^2} + O([\Delta x]^3), \quad \text{for } \Delta x \rightarrow 0,$$

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We now substitute the expansions

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$$W_j(t + \Delta t) = \frac{1}{2} W_{j-1}(t) + \frac{1}{2} W_{j+1}(t)$$

obtaining

$$W(x, t) + \Delta t \frac{\partial W}{\partial t} + O(\Delta t^2) = W(x, t) + \frac{1}{2} \Delta x^2 \frac{\partial^2 W}{\partial x^2} + O(\Delta x^3)$$



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$$\frac{\partial W}{\partial t} = \frac{\Delta x^2}{2\Delta t} \frac{\partial^2 W}{\partial x^2} + O(\Delta x^3 + \Delta t)$$

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obtaining

$$\frac{\partial W}{\partial t} = K_1 \frac{\partial^2 W}{\partial x^2}, \quad K_1 = \lim_{\substack{\Delta x \rightarrow 0 \\ \Delta t \rightarrow 0}} \frac{\Delta x^2}{2\Delta t} < \infty.$$

# Brownian motion

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$$\frac{\partial W}{\partial t} = K_1 \frac{\partial^2 W}{\partial x^2}$$

Let us call  $X$  the random variable measuring the distance covered in two consecutive jumps

- Assume that the *pdf* of  $X$  (appropriately normalised) has existing moments

$$\overline{X} = \sum_i X_i, \quad \overline{X^2},$$

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- Then the **central limit theorem** assures that exists

$$V = \frac{\overline{X}}{\Delta t} \text{ (Mean velocity)} \quad K = \frac{\overline{X^2} - \overline{X}^2}{2\Delta} \text{ (Diffusion coefficient)}$$

and that

$$W(x, t) = \frac{1}{2\sqrt{\pi K_1 t}} \exp(-x^2/4K_1 t).$$

# Brownian motion: the Fourier domain

---

We can rewrite

$$W(x, t) = \frac{1}{\sqrt{2\pi K_1 t}} \exp(-x^2/4K_1 t).$$

in the **Fourier domain** as

$$W(k, t) = \exp(-K_1 k^2 t), \quad W_0(x) = \lim_{t \rightarrow 0^+} W(x, t) = \delta(x),$$

that solve the **Fourier transformed diffusion equation**

$$\frac{\partial W}{\partial t} = -K_1 k^2 W(k, t),$$

that is a **relaxation equation**, for a fixed wavenumber  $k$ .

# From the discrete to the continuous

---

The **C**ontinuous **T**ime **R**andom **W**alk model (CTRW):

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$$\lambda(x) = \int_0^{+\infty} \psi(x, t) dt,$$

## Jump length

$\lambda(x)dx$  produces the probability for a jump length in the interval  $(x, x + dx)$ .

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🕒 The waiting time pdf

$$w(t) = \int_{-\infty}^{+\infty} \psi(x, t) dx$$

## Waiting time

$w(t)dt$  produces the probability for a waiting time in the interval  $(t, t + dt)$ .



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- If the jump length and waiting time are **independent random variables** then:

$$\psi(x, t) = w(t)\lambda(x)$$

# Characterization of CTRW

---

To categorise different CTRW one can look at the quantities

$$T = \int_0^{+\infty} tw(t) dt, \text{ (Characteristic waiting time),}$$

and

$$\Sigma^2 = \int_{-\infty}^{+\infty} x^2 \lambda(x) dx \text{ (Jump length variance),}$$

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The **master** (Langevin) **equation** for this process is then given by

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# Characterization of CTRW

Then if we use

$$\eta(x, t) = \int_{-\infty}^{+\infty} dx' \int_0^{+\infty} dt' \eta(x', t') \psi(x - x', t - t') + \delta(x) \delta(t),$$

we can write the pdf of being in  $x$  at time  $t$  as

$$W(x, t) = \int_0^t \eta(x, t') \Psi(t - t') dt', \quad \Psi(t) = 1 - \int_0^t w(t') dt',$$

where the latter is the cumulative probability assigned to the probability of **no jump event** during the time interval  $t - t'$ .

## Fact

If both  $T$  and  $\Sigma^2$  are finite the long-time limit corresponds to Brownian motion, e.g.,  $w(t) = \tau^{-1} \exp(-t/\tau)$ ,  $T = \tau$ ,  $\lambda(x) = (4\pi\sigma^2)^{-1/2} \exp(-x^2/4\sigma^2)$ ,  $\Sigma^2 = 2\sigma^2$ , we recover the standard diffusion equation.

# The CTRW in the Fourier-Laplace domain

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We take

$$W(x, t) = \int_0^t \eta(x, t') \Psi(t - t') dt', \quad \Psi(t) = 1 - \int_0^t w(t') dt',$$

and rewrite it again in the **Fourier-Laplace domain** (Fourier for the space variable, Laplace for the time one) as

$$W(k, u) = \frac{1 - w(u)}{u} \frac{W_0(k)}{1 - \psi(k, u)}, \quad W_0(k) = \int_{-\infty}^{+\infty} W_0(x) e^{-i2\pi kx} dx.$$

In the **Brownian case**

$$w(u) \sim 1 - u\tau + O(\tau^2), \quad \lambda(k) \sim 1 - \sigma^2 k^2 + O(k^4), \quad W_0(x) = \delta(x)$$

then

$$W(k, u) = \frac{1}{u + K_1 k^2}, \quad K_1 = \sigma^2/\tau.$$



# The case of long rests

## Long rests

The **characteristic waiting time**  $T = \int_0^{+\infty} tw(t) dt$  **diverges**, but the jump length variance  $\Sigma^2 = \int_{-\infty}^{+\infty} x^2 \lambda(x) dx$  is finite.

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- and then obtain the expression for  $W(k, u)$  in the Fourier-Laplace space

$$W(k, u) = w_0^{(k)}/u / (1 + K_\alpha u^{-\alpha} k^2).$$

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To get an expression of the equation we use the Laplace transform for fractional integrals:

$$\mathcal{L} \left\{ I_{[0,t]}^{-\alpha} W(x, t) \right\} = u^{-\alpha} W(x, u), \quad \alpha \geq 0,$$

and together with

$$W(k, u) = \frac{W_0(k)/u}{(1 + K_\alpha u^{-\alpha} k^2)}.$$

we infer the fractional integral equation

$$W(x, t) - W_0(x) = I_{[0,t]} K_\alpha \frac{\partial^2}{\partial x^2} W(x, t).$$

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we infer the fractional integral equation, and apply derivative w.r.t. to time

$$\frac{\partial}{\partial t} (W(x, t) - W_0(x)) = \frac{\partial}{\partial t} \left( I_{[0,t]} K_\alpha \frac{\partial^2}{\partial x^2} W(x, t) \right).$$

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We can compute also the mean squared displacement

$$\langle x^2(t) \rangle = \mathcal{L}^{-1} \left\{ \lim_{k \rightarrow 0} -\frac{d^2}{dk^2} W(k, u) \right\} = \frac{2K_\alpha}{\Gamma(1 + \alpha)} t^\alpha.$$

# The case of long rests

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We have obtained a Fractional Differential Equation:

$$\frac{\partial W}{\partial t} = {}_{RL}D_{[0,t]}^{\alpha} K_{\alpha} \frac{\partial^2}{\partial x^2} W(x, t), \quad 0 < \alpha < 1$$

but this is not the model we started looking at, that was

$${}_CD_{[0,t]}^{\alpha} \frac{\partial W}{\partial t} = K_{\alpha} \frac{\partial^2}{\partial x^2} W(x, t), \quad 0 < \alpha < 1$$

❓ Are they related?

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We have obtained a Fractional Differential Equation:

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We now have an *interpretation* of what a Fractional Derivative with respect to time is. We will come back to this when we will speak about fractional derivative with respect to space.

# “Exponential” Fractional Integrators

---

We start from the FDE

$$\begin{cases} {}_{CA}D_{[t_0, t]}^\alpha u(t) + \lambda y(t) = f(t), \\ u(0) = u_0, \end{cases} \quad \alpha \in \mathbb{R}_{>0}, \quad \lambda \in \mathbb{R}, \quad u(t) : [t_0, T] \rightarrow \mathbb{R}.$$

Then we rewrite the solution as

$$u(t) = e_{\alpha,1}(t - t_0; \lambda) u_0 + \int_{t_0}^t e_{\alpha,\alpha}(t - s; \lambda) f(s) \, ds, \quad e_{\alpha,\beta} = t^{\beta-1} E_{\alpha,\beta}(-\lambda t^\alpha),$$

for  $E_{\alpha,\beta}(z)$  the Mittag-Leffler (ML) function with two parameters.

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- 💡 We can use this formulation to build different PI rules,
- 💡 We can use it to address the problem

$${}_{CA}D_{[t_0, t]}^\alpha U(t) + Ay(t) = F(U(t)), \quad U(0) = U_0.$$

# Evaluation of the ML function

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For both the approaches we need reliable ways for **computing** the **ML function** on both the **real line** and with **matrix argument**.



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In general, we expect to mostly need matrix function–times–vector operations:

$$\mathbf{y} = E_{\alpha,\beta}(A)\mathbf{v}, \quad A \in \mathbb{R}^{n \times n}, \quad \mathbf{y}, \mathbf{v} \in \mathbb{R}^n.$$

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We postpone it to after we have discussed the actual necessities we have.

# PI - “Exponential” Fractional Integrators

---

We start from the formula

$$u(t) = e_{\alpha,1}(t - t_0; \lambda) u_0 + \int_{t_0}^t e_{\alpha,\alpha}(t - s; \lambda) f(s) \, ds, \quad e_{\alpha,\beta} = t^{\beta-1} E_{\alpha,\beta}(-\lambda t^\alpha),$$

and select a grid  $\{t_i\}_{i=0}^N$ , then

$$u(t_n) = e_{\alpha,1}(t_n - t_0; \lambda) u_0 + \sum_{j=0}^{n-1} \int_{t_j}^{t_{j+1}} e_{\alpha,\alpha}(t_n - s; \lambda) f(s) \, ds.$$

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$$u(t_n) = e_{\alpha,1}(t_n - t_0; \lambda) u_0 + \tau^\alpha \sum_{j=0}^{n-1} \int_0^1 e_{\alpha,\alpha}((t-t_j)/\tau - r; \tau^\alpha \lambda) f(t_j + r\tau) dr.$$

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# PI - “Exponential” Fractional Integrators

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Then a PI rule for

$$u(t_n) = e_{\alpha,1}(t_n - t_0; \lambda) u_0 + \tau^\alpha \sum_{j=0}^{n-1} \int_0^1 e_{\alpha,\alpha}((t-t_j)/\tau - r; \tau^\alpha \lambda) f(t_j + r\tau) dr.$$

is obtained by selecting  $q + 1$  *distinct* nodes  $0 \leq c_0 < c_1 < \dots < c_q \leq 1$  and replacing  $f(t_j + r\tau)$  with

$$p_j^{[q]}(t_j + r\tau) = \sum_{\ell=0}^q L_\ell^{[q]}(r) f(t_j + c_\ell \tau), \quad r \in [0, 1], \quad L_\ell^{[q]} \text{ Lagrange basis element of degree } q.$$



# PI - “Exponential” Fractional Integrators

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Then the PI rule is

$$u^{(n)} = e_{\alpha,1}(t_n - t_0; \lambda) y_0 + \tau^\alpha \sum_{j=0}^{n-1} \sum_{\ell=0}^q \omega_\ell^{[q;\alpha]}(n-j; \tau^\alpha \lambda) f(t_j + c_\ell \tau).$$

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And selecting the weights

$$\omega_\ell^{[q;\alpha]}(n, z) = \int_0^1 e_{\alpha,\alpha}(n-j-r; z) L_\ell^{[q]}(r) dr.$$

# PI - “Exponential” Fractional Integrators

Theorem (Garrappa and Popolizio 2011, Theorem 4.2)

Let  $\alpha > 0$  and  $f(t) \in \mathcal{C}^{q+2}([t_0, T])$ . The error of a  $q$ -step exponential PI rule is given by

$$u(t_n) - u^{(n)} = \tau^{q+1} \frac{C_0^{[q]}}{(q+1)!} \int_{t_0}^{t_n} e_{\alpha, \alpha}(t_n - s; \lambda) f^{(q+1)}(s) ds + O(\tau^{q+1+\alpha}),$$

where the constant  $C_0^{[q]}$  depends only on the nodes  $c_\ell$ .

- For  $q = 2$ ,  $c_0 = 0$ ,  $c_1 = 1/2$ ,  $c_2 = 1$ , one finds  $C_0^{[2]} = 0$ , thus an interpolatory formula of order  $O(\tau^{q+1+\alpha})$ .

💡 The **general idea** is to select nodes  $c_\ell$  in such way that

$$C_v^{[q]} = \int_0^1 \omega_q(r) \xi(1-v, 1-r) dr, \quad v \in \mathbb{R},$$

for  $\xi$ , the *Hurwitz zeta function*, are zeroed out in the error expansion for the method.

# The MOL/Matrix case

---

Let us go back to the case that sparked our interest in going “exponential”, that was the MOL problem

$$\begin{cases} {}_C D_{[0,t]}^\alpha \mathbf{u}(t) + A\mathbf{u}(t) = \mathbf{g}(t), & t > 0, \\ \mathbf{u}(0) = \mathbf{u}_0. \end{cases}$$

By the variation of constant formula, we have seen that we can express the solution as

$$\mathbf{u}(t) = E_{\alpha,1}(-t^\alpha A)\mathbf{u}_0 + \int_0^t (t-s)^{\alpha-1} E_{\alpha,\alpha}(-A(t-s)^\alpha) \mathbf{g}(s) \, ds.$$

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- In the general case we then have to apply one of the PI rules to compute the integral term,
- If  $\mathbf{g}(s) = \sum_{k=0}^q s^k \mathbf{v}_k$  for some vectors, we can compute the integral on the right-hand side in *closed form* and obtain

$$\mathbf{u}(t) = E_{\alpha,1}(-t^\alpha A)\mathbf{y}_0 + \sum_{k=0}^q \Gamma(k+1) t^{\alpha+k} E_{\alpha,\alpha+k+1}(-t^\alpha A) \mathbf{v}_k, \quad t > 0.$$

# Matrix functions: the normal case

---

If  $A$  is a normal matrix, and  $f$  is a function existing on the spectrum of  $A$ , then

$$f(A) = Uf(\Lambda)U^H, \quad U^H U = I, \quad \Lambda = \text{diag}(\lambda_1, \dots, \lambda_n), \quad A\mathbf{u}_i = \lambda_i\mathbf{u}_i, \quad U = [\mathbf{u}_1, \dots, \mathbf{u}_n].$$

This is, e.g., sufficient for the cases in which

- $A$  is the discretization of a self-adjoint operator,
- $A$  is symmetric.

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What about the *non-normal* and *nond-diagonalizable* case? For diagonalizable matrices, we can use the eigendecomposition at the same way.

# Matrix functions: the Jordan Canonical Form

## Jordan Canonical Form

We recall that any matrix  $A \in \mathbb{C}^{n \times n}$  can be expressed in Jordan canonical form as

$$Z^{-1}AZ = J = \text{diag}(J_1, \dots, J_p), \quad \text{for } J_k = J_k(\lambda_k) = \begin{bmatrix} \lambda_k & 1 & & \\ & \lambda_k & \ddots & \\ & & \ddots & 1 \\ & & & \lambda_k \end{bmatrix} \in \mathbb{C}^{m_k \times m_k},$$

where  $Z$  is nonsingular and  $m_1 + m_2 + \dots + m_p = n$ . If each block in which the eigenvalue  $\lambda_k$  appears is of size 1 then  $\lambda_k$  is said to be a *semisimple* eigenvalue.

- This is a *theoretical object*, it is useful to prove and define *things*, **not to implement things**.



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- This is a *theoretical object*, it is useful to prove and define *things*, **not to implement things**.
- Now that we have a decomposition of the matrix, we need to introduce a suitable definition of **being defined on the spectrum**.

# Matrix functions: the general case

---

Let us denote by  $\lambda_1, \dots, \lambda_s$  the distinct eigenvalues of  $A$ , and by  $n_i$  the order of the largest Jordan block in which the  $\lambda_i$  appears, i.e., the *index* of the eigenvalue  $\lambda_i$ .

## Defined on the spectrum

The function  $f$  is *defined on the spectrum of  $A$*  if the values

$$f^{(j)}(\lambda_i), \quad j = 0, 1, \dots, n_i - 1, \quad i = 1, \dots, s,$$

exist, where  $f^{(j)}$  denotes the  $j$ th derivative of  $f$ , with  $f^{(0)} = f$ .

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⚠ Again for the ML function and  $\alpha > 0$  we have no problem with this.

# Matrix functions: the general case

## Matrix function

Let  $f$  be defined on the spectrum of  $A \in \mathbb{C}^{n \times n}$ , which is represented in Jordan canonical form as  $Z^{-1}AZ = J$ ,

$$f(A) = Zf(J)Z^{-1} = Z \operatorname{diag}(f(J_1), \dots, f(J_p))Z^{-1},$$

where

$$f(J_k) = \begin{bmatrix} f(\lambda_k) & f'(\lambda_k) & \dots & \frac{f^{(m_k-1)}(\lambda_k)}{(m_k-1)!} \\ & f(\lambda_k) & \ddots & \vdots \\ & & \ddots & f'(\lambda_k) \\ & & & f(\lambda_k) \end{bmatrix}.$$

Moreover, let  $f$  be a multivalued function and suppose some eigenvalues occur in more than one Jordan block. If the same choice of branch of  $f$  is made in each block, then we say that  $f(A)$  is a *primary matrix function*.

# Matrix functions: computing $f(A)$ and $f(A)\mathbf{v}$

---

To march our scheme for

$$\mathbf{u}(t) = E_{\alpha,1}(-t^\alpha A)\mathbf{u}_0 + \int_0^t (t-s)^{\alpha-1} E_{\alpha,\alpha}(-A(t-s)^\alpha) \mathbf{g}(s) \, ds.$$

we need to compute operations of the form  $f(A)\mathbf{v}$ , *nevertheless*, we will have to compute  $f(\cdot)$  at least on some **small matrix**.

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## Schur decomposition and matrix functions

Given a matrix  $A$  there exist always a matrix  $Q$  such that  $Q^*Q = I$ , and a upper triangular matrix  $T$  such that  $A = QTQ^*$ . Then, if  $f$  is **defined on the spectrum** of  $A$  we can compute  $f(A)$  as  $f(A) = Qf(T)Q^*$ .

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But how do we compute the matrix function of an upper triangular matrix?

# Matrix functions: the upper triangular case

---

**Assumption** we assume that  $T$  is such that each block  $T_{i,j}$  has **clustered eigenvalues**, and distinct diagonal blocks have *far enough* eigenvalues.

❗ If the **assumption** doesn't hold we look for a block permutation.

$$\left[ \begin{array}{cc|c} (T_{1,1})_{1,1} & (T_{1,1})_{1,2} & T_{1,2} \\ 0 & (T_{1,1})_{2,2} & \\ \hline \mathbf{0} & & (T_{2,2})_{1,1} & (T_{2,2})_{1,2} \\ & & 0 & (T_{2,2})_{2,2} \end{array} \right]$$

⚠ Close eigenvalues may lead to severe *accuracy loss*, even far apart eigenvalues can produce more inaccurate answers than expected, see (Davies and Higham 2003).



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- To evaluate  $f(T_{ii})$  we use the Taylor series in  $\sigma$

$$f(T_{i,i}) = \sum_{k=0}^{+\infty} \frac{f^{(k)}}{k!} M^k,$$

for  $\sigma = \text{trace}(T_{i,i})/m$ ,  $m = \dim(T_{i,i})$ , and  $M = T_{i,i} - \sigma I$ .

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For the **off-diagonal blocks** we apply the block-Parlett recurrence

$$F_{i,j} = f(T_{i,j}), \quad i = 1, \dots, n;$$

**for**  $j = 2, \dots, n$  **do**

**for**  $i = j - 1, j - 2, \dots, 1$  **do**

        Solve Sylvester equation for  $F_{i,j}$ :

$$T_{i,i}F_{j,j} - F_{i,j}T_{j,j} = F_{i,i}T_{i,j} - T_{i,j}F_{j,j} \\ + \sum_{k=0}^{j-1} (F_{i,k} - T_{k,j} - T_{i,k}F_{k,j}).$$

**end**

**end**

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## What we need

To use the algorithm we have sketched out, we need to be able to compute the derivatives of the ML function sufficiently accurately.

# Derivatives of the ML function

---

The key observation for this task is

$$\frac{d^k}{dz^k} E_{\alpha, \beta}(z) = \sum_{j=0}^{+\infty} \frac{(j+k)_k z^j}{\Gamma(\alpha j + \alpha k + \beta)} = \frac{k!}{\Gamma(k+1)} \sum_{j=0}^{+\infty} \frac{\Gamma(j+k+1) z^j}{j! \Gamma(\alpha j + \alpha k + \beta)} = k! E_{\alpha, \alpha k + \beta}^{k+1}(z),$$

where

$$E_{\alpha, \beta}^{\gamma}(z) = \frac{1}{\Gamma(\gamma)} \sum_{j=0}^{+\infty} \frac{\Gamma(1+\gamma) z^j}{j! \Gamma(\alpha j + \beta)},$$

is called the **Prabhakar function**.

# Derivatives of the ML function

---

The key observation for this task is

$$\frac{d^k}{dz^k} E_{\alpha,\beta}(z) = \sum_{j=0}^{+\infty} \frac{(j+k)_k z^j}{\Gamma(\alpha j + \alpha k + \beta)} = \frac{k!}{\Gamma(k+1)} \sum_{j=0}^{+\infty} \frac{\Gamma(j+k+1) z^j}{j! \Gamma(\alpha j + \alpha k + \beta)} = k! E_{\alpha,\alpha k + \beta}^{k+1}(z),$$

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is called the **Prabhakar function**.

Its **efficient computation** can be obtained, similarly to the ML function, by means of a *Laplace transform inversion*

$$\mathcal{L} \left\{ t^{\beta-1} E_{\alpha,\beta}^{\gamma}(t^{\alpha} z) \right\} (s) = \frac{s^{\alpha\gamma-\beta}}{(s^{\alpha} - t^{\alpha} z)^{\gamma}}, \quad \Re(s) > 0, \quad |t^{\alpha} z s^{-\alpha}| < 1.$$

# Computing the Prabhakar function (Garrappa 2015)

---

We select  $t = 1$  in

$$\mathcal{L} \left\{ t^{\beta-1} E_{\alpha,\beta}^{\gamma}(t^{\alpha}z) \right\} (s) = \frac{s^{\alpha\gamma-\beta}}{(s^{\alpha} - t^{\alpha}z)^{\gamma}}, \quad \Re(s) > 0, \quad |t^{\alpha}zs^{-\alpha}| < 1.$$

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- Since

$$\frac{d^k}{dz^k} E_{\alpha, \beta}(z) = k! E_{\alpha, \alpha k + \beta}^{k+1}(z) = \frac{k!}{2\pi i} \int_{\mathcal{C}} e^s H_k(s; z) ds \equiv I_k(z),$$



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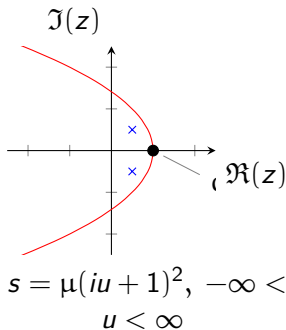
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- we use the *Optimal Parabolic Contour* we have already discussed in **Lecture 2** to determine the deformation of the Bromwich line to evaluate

$$I_k^{[N]} = \frac{k!h}{2\pi i} \sum_{j=-N}^N e^{\sigma(u_j)} H_k(\sigma(u_j); z) \sigma'(u_j).$$



## An alternative option (Higham and Liu 2021)

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We needed the ML derivatives to apply Schur-Parlett to non-diagonalizable matrices.

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### Diagonalization by perturbation

Let  $A$  be nonnormal

$$\tilde{A} = A + E$$

for  $E$  a suitable perturbation is *likely to be diagonalizable*. **Diagonalizable matrices are dense in  $\mathbb{C}^{n \times n}$** , for a given  $A$  and machine precision  $\epsilon$  then the best approximate diagonalization can be measured in terms of

$$\sigma(A, \epsilon) = \inf_{E, V} \sigma(A, V, E, \epsilon) = \inf_{E, V} \{\kappa_2(V)\epsilon + \|E\|_2\}.$$

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We can expect to measure on  $f(A)$  by estimating

$$\|f(A + E) - f(A)\| \lesssim \|L_f(A, E)\| \leq \|L_f(A)\| \|E\|,$$

for  $L_f(A, E)$  the **Fréchet derivative** of  $f$  at  $A$  in direction  $E$ ,  $\|L_f(A)\| = \max_{\|E\|=1} \{\|L_f(A, E)\|\}.$

## An alternative option (Higham and Liu 2021)

### Fréchet derivative

The **Fréchet derivative** of a matrix function  $f : \mathbb{C}^{n \times n} \rightarrow \mathbb{C}^{n \times n}$  at a point  $X \in \mathbb{C}^{n \times n}$  is a linear mapping  $L : \mathbb{C}^{n \times n} \rightarrow \mathbb{C}^{n \times n}$   $E \mapsto L_f(X, E)$  such that for all  $E \in \mathbb{C}^{n \times n}$  we find

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Thus, in our estimate we have

$$\|f(A + E) - f(A)\| \lesssim \|L_f(A, E)\| \leq \|L_f(A)\| \|E\|,$$

and therefore the change in  $f$  induced by  $E$  grows as  $\|L_f(A)\|_2 \|E\|_2$  and there are many cases in which  $\|L_f(A)\|_2 \gg 1$ .

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💡 The idea from (Higham and Liu 2021) is to use a *structured perturbation*:  
“take  $E$  to be upper triangular standard Gaussian matrix.”

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The idea in few steps

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What precision do we need?

To have  $\kappa_1(V)u_h \lesssim u$  we select for  $c_mu \approx \min_i |\text{diag}(\tilde{t}_{1,1}I - \tilde{T}_{2,2})|$

$$u_h \lesssim \frac{c_mu^2}{\max_{i < j} |\tilde{t}_{i,j}| (\max_{i < j} |\tilde{t}_{i,j}|/c_mu + 1)^{k-2}}, \quad k = \text{"size of the Jordan block"} \geq 2.$$

# From small to large matrices

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We now know how to compute  $E_{\alpha,\beta}(A)$  for a *small matrix*  $A$ , either with

- Classical Schur-Parlett algorithm with Laplace inversion technique for the needed derivative of the ML function (Garrappa and Popolizio 2018),  
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What about *large matrices*?

## 💡 Projection methods for matrix functions

We can exploit the *subspace projection* idea, take  $V \in \mathbb{R}^{n \times k}$  spanning a given subspace  $\mathcal{W}_k$

$$f(A)\mathbf{v} \approx Vf(V^T AV)V^T \mathbf{v} \quad V^T AV \in \mathbb{R}^{k \times k}, \quad k \ll n.$$



# Krylov Projection Methods

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## A general framework

Given a set of scalars  $\{\sigma_1, \dots, \sigma_{k-1}\} \subset \overline{\mathbb{C}}$  (the extended complex plane), that are not eigenvalues of  $A$ , let

$$q_{k-1}(z) = \prod_{j=1}^{k-1} (\sigma_j - z).$$

The **rational Krylov** subspace of order  $k$  associated with  $A$ ,  $\mathbf{v}$  and  $q_{k-1}$  is defined by

$$\mathcal{Q}_k(A, \mathbf{v}) = [q_{k-1}(A)]^{-1} \mathcal{K}_k(A, \mathbf{v}), \quad \mathcal{K}_k(A, \mathbf{v}) = \text{Span}\{\mathbf{v}, A\mathbf{v}, \dots, A^{k-1}\mathbf{v}\}.$$

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Given  $\{\mu_1, \dots, \mu_{k-1}\} \subset \overline{\mathbb{C}}$  such that  $\sigma_j \neq \mu_j^{-2}$ , we define the matrices

$$C_j = (\mu_j \sigma_j A - I) (\sigma_j I - A)^{-1}, \text{ and } \mathcal{Q}_k(A, \mathbf{v}) = \text{Span}\{\mathbf{v}, C_1 \mathbf{v}, \dots, C_{k-1} \cdots C_2 C_1 \mathbf{v}\}.$$

# Krylov Projection Methods: special cases

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$$(\mu_j, \sigma_j) = \begin{cases} (1, \infty), & \text{for } j \text{ even,} \\ (0, 0), & \text{for } j \text{ odd.} \end{cases}$$

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Shift-And-Invert  $\mathcal{W}_k(A, \mathbf{v}) = \text{Span}\{\mathbf{v}, (\sigma I - A)^{-1}\mathbf{v}, \dots, (\sigma I - A)^{-(k-1)}\mathbf{v}\}$ , take  $\mu_j = 0$  and  $\sigma_j = \sigma$  for each  $j$ ,

# The ML function (Moret and Novati 2011)

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To estimate the convergence behavior of general projection methods in the non-normal we need the concept of **field of values** (or *numerical range*.)

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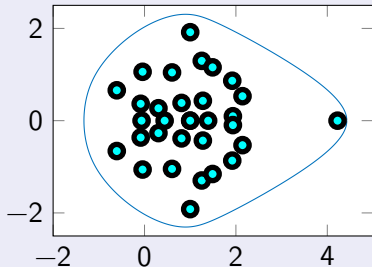
To estimate the convergence behavior of general projection methods in the non-normal we need the concept of **field of values** (or *numerical range*.)

## Field of Values/Numerical Range

Given  $A \in \mathbb{C}^{N \times N}$  we denote its **field of values** as

$$W(A) = \left\{ \frac{\langle \mathbf{x}, A\mathbf{x} \rangle}{\langle \mathbf{x}, \mathbf{x} \rangle}, \quad \mathbf{0} \neq \mathbf{x} \in \mathbb{C}^N \right\},$$

where  $\langle \cdot, \cdot \rangle$  represents the Euclidean inner product.





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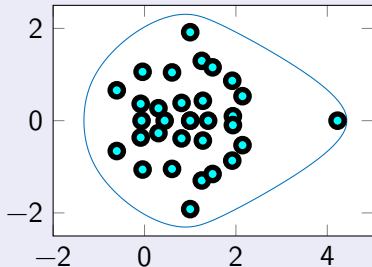
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It has many **properties**, e.g.,  $W(A) \subseteq D(0, \|A\|)$  (disk centered on 0 with radius  $\|A\|$ ), is *compact*, sub-additive  $W(A + B) \subseteq W(A) + W(B)$ , unitarily invariant  $W(UAU^H) = UW(A)U^H$ , etc. see (Benzi 2021).

# The ML function (Moret and Novati 2011)

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## Assumptions:

(A1) We assume that  $\exists a > 0$ ,  $\theta \in [0, \pi/2)$  such that

$$W(A) \subset \Sigma_{\theta, a} = \{\lambda \in \mathbb{C} : |\arg(\lambda) - a| \leq \theta\}.$$

(A2)  $\beta > 0$ ,  $\alpha \in (0, 2)$  be such that  $\alpha\pi/2 < \pi - \theta$ ,  $\varepsilon > 0$  and

$$\frac{\alpha\pi}{2} < \mu \leq \min\{\pi, \alpha\pi\}, \quad \mu < \pi - \theta.$$

**Method of choice:** we use **polynomial Krylov method**  $\mathcal{K}_m(A, \mathbf{v})$ :

$$AV_m = V_m H_m + h_{m+1, m} \mathbf{v}_{m+1} \mathbf{e}_m^T, \quad \text{Span } V_m = \text{Span}\{\mathbf{v}_i\}_{i=1}^m = \mathcal{K}_m(A, \mathbf{v}), \quad H_m = V_m^H A V_m.$$

**We want to bound:**

$$R_m = E_{\alpha, \beta}(-A)\mathbf{v} - V_m E_{\alpha, \beta}(-H_m)\mathbf{e}_1, \quad m \geq 1.$$

# The ML function (Moret and Novati 2011)

We first express the error in *integral form*, starting from (Podlubny 1999, Theorem 1.1)

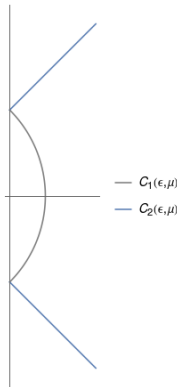
$$E_{\alpha,\beta}(z) = \frac{1}{2\alpha\pi i} \int_{C(\varepsilon,\mu)} \frac{\exp(\lambda^{1/\alpha})\lambda^{1-\beta/\alpha}}{\lambda - z} d\lambda, \quad z \in G^-(\varepsilon,\mu),$$

where

- $\forall \varepsilon > 0, 0 < \mu < \pi$

$$C(\varepsilon,\mu) = \bigcup \begin{cases} C_1(\varepsilon,\mu) = \{\lambda : \lambda = \varepsilon \exp(i\varphi), & -\mu \leq \varphi \leq \mu\}, \\ C_2(\varepsilon,\mu) = \{\lambda : \lambda = r \exp(\pm i\mu), & r \geq \varepsilon\}. \end{cases}$$

- The contour  $C(\varepsilon,\mu)$  divides the complex plane into two domains,  $G^-(\varepsilon,\mu)$  and  $G^+(\varepsilon,\mu)$  lying respectively on the left and on the right of  $C(\varepsilon,\mu)$ .



# An Expression for the Error

---

From the previous we find

$$E_{\alpha,\beta}(-A) = \frac{1}{2\alpha\pi i} \int_{C(\varepsilon,\mu)} \exp(\lambda^{1/\alpha}) \lambda^{1-\beta/\alpha} (\lambda I + A)^{-1} d\lambda, \quad \sigma(-A) \in G^-(\varepsilon, \mu),$$

and together with

$$R_m = E_{\alpha,\beta}(-A)\mathbf{v} - V_m E_{\alpha,\beta}(-H_m)\mathbf{e}_1, \quad m \geq 1,$$

we write

$$R_m = \frac{1}{2\alpha\pi i} \int_{C(\varepsilon,\mu)} \exp(\lambda^{1/\alpha}) \lambda^{1-\beta/\alpha} \delta_m(\lambda) d\lambda,$$

for

$$\begin{aligned} \delta_m(\lambda) &= (\lambda I + A)^{-1} \mathbf{v} - V_m (\lambda I + H_m)^{-1} \mathbf{e}_1 \\ &= (\lambda I + A)^{-1} \mathbf{v} - V_m (\lambda I + H_m)^{-1} V_m^H \mathbf{v}. \end{aligned}$$

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Observe now that

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Therefore we have

$$\Delta_m (\lambda I + A) V_m = 0.$$

# An Expression for the Error

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Observe now that

$$\delta_m(\lambda) = (\lambda I + A)^{-1} \mathbf{v} - V_m(\lambda I + H_m)^{-1} V_m^H \mathbf{v} = \Delta_m \mathbf{v},$$

By using the **Arnoldi relation**, since  $\mathbf{v}_{m+1} \perp V_m$ :

$$V_m^H (\lambda I + A) V_m = \lambda I + H_m,$$

Therefore we have

$$\Delta_m (\lambda I + A) V_m = 0.$$

For an arbitrary  $\mathbf{y} \in \mathbb{C}^m$  we have then

$$(\lambda I + A)^{-1} \mathbf{v} - V_m(\lambda I + H_m)^{-1} V_m^H \mathbf{v} = \Delta_m (\mathbf{v} - (\lambda I + A) V_m \mathbf{y}) = \Delta_m p_m(A) \mathbf{v},$$

where  $p_m(z)$  is a **polynomial of degree  $\leq m$**  with  $p_m(-\lambda) = 1$ .



# An Expression for the Error

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We have therefore proved that

$$\|\delta_m(A)\| \leq \|(\lambda I + A)^{-1} - V_m(\lambda I + H_m)^{-1}V_m^H\| \|p_m(A)\mathbf{v}\|, \forall p_m \in \mathbb{P}_{\leq m}[z] \text{ with } p_m(-\lambda) = 1.$$

By using (Diele, Moret, and Ragni [2008/09](#), Lemma 2) we also have the following expression

$$\|\delta_m(\lambda)\| = \frac{\prod_{j=1}^m h_{j+1,j}}{|\det(\lambda I + H_m)|} \|(\lambda I + A)^{-1}\mathbf{v}_{m+1}\|.$$

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To obtain the first bound we call then

$$D(\lambda) = \text{dist}(\lambda, W(-A)) \quad \forall \lambda \in C(\varepsilon, \mu).$$

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## Representation function

Using (A1) and (A2) we can find a function  $\nu(\varphi)$  such that

$$\forall \lambda = |\lambda| \exp(\pm i\varphi) \in C(\varepsilon, \mu) \quad D(\lambda) \geq \nu(\varphi)|\lambda|, \quad \nu(\varphi) \geq \nu > 0.$$

# A First Error Bound

---

Theorem (Moret and Novati 2011, Theorem 3.2)

Let assumptions (A1) and (A2) hold, then for  $m \geq 1$  and for every  $M > 0$  we have

$$\|R_m\| \leq \frac{\exp(M) \prod_{j=1}^m h_{j+1,j}}{\pi \nu^{m+1} M^{m\alpha+\beta-1}} \left( \frac{\mu}{\alpha} + \frac{\exp(-M(|\cos(\mu/\alpha)| + 1))}{m\alpha - 1 + \beta} \right).$$

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**Proof.** We use  $\|(\lambda I + A)^{-1}\| \leq D(\lambda)^{-1}$  and  $W(H_m) \subseteq W(A)$  in the error expression  $R_m$

$$\begin{aligned} \|R_m\| &= \left\| \frac{1}{2\alpha\pi i} \int_{C(\varepsilon, \mu)} \exp(\lambda^{1/\alpha}) \lambda^{1-\beta/\alpha} \delta_m(\lambda), d\lambda \right\| \\ &\leq \frac{\prod_{j=1}^m h_{j+1,j}}{2\pi\alpha} \int_{C(\varepsilon, \mu)} \frac{|\exp(\lambda^{1/\alpha}) \lambda^{1-\beta/\alpha}|}{D(\lambda)^{m+1}} |d\lambda|. \end{aligned}$$

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$$\|R_m\| \leq \frac{\prod_{j=1}^m h_{j+1,j}}{2\pi\alpha} (I_1 + I_2),$$

with

$$I_1 = \int_{C_1(\varepsilon, \mu)} \frac{|\exp(\lambda^{1/\alpha}) \lambda^{1-\beta/\alpha}|}{D(\lambda)^{m+1}} |\mathrm{d}\lambda| \leq 2\varepsilon^{\frac{1-\beta}{\alpha}-m} \int_0^\mu \frac{\exp(\varepsilon^{1/\alpha} \cos(\varphi/\alpha))}{\nu(\varphi)^{m+1}} \mathrm{d}\varphi,$$

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$$\|R_m\| \leq \frac{\prod_{j=1}^m h_{j+1,j}}{2\pi\alpha} \left( 2\varepsilon^{\frac{1-\beta}{\alpha}-m} \int_0^\mu \frac{\exp(\varepsilon^{1/\alpha} \cos(\varphi/\alpha))}{\nu(\varphi)^{m+1}} d\varphi + I_2 \right),$$

with

$$\begin{aligned} I_2 &= \int_{C_2(\varepsilon, \mu)} \frac{|\exp(\lambda^{1/\alpha}) \lambda^{1-\beta/\alpha}|}{D(\lambda)^{m+1}} |d\lambda| \leq \frac{2}{\nu^{m+1}} \int_\varepsilon^{+\infty} \frac{r^{\frac{1-\beta}{\alpha}} \exp(-r^{\frac{1}{\alpha}} |\cos(\mu/\alpha)|)}{r^{m+1}} dr \\ &= \frac{2}{\nu^{m+1}} \int_{\varepsilon^{1/\alpha}}^{+\infty} \frac{\exp(-s |\cos(\mu/\alpha)|)}{s^{m\alpha+\beta}} ds \leq \frac{2\alpha \exp(-\varepsilon^{1/\alpha} |\cos(\mu/\alpha)|)}{(m\alpha + \beta - 1) \nu^{m+1} \varepsilon^{\frac{m\alpha+\beta-1}{\alpha}}}. \end{aligned}$$

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The result follows then by setting  $\varepsilon = M^\alpha$  and simplifying the expression.  $\square$

 With the same proof another bound for the case of small  $\alpha$  can be obtained.

# A First Error Bound: small $\alpha$ s

Theorem (Moret and Novati 2011, Theorem 3.2)

Let assumptions (A1) and (A2) hold, then for  $m \geq 1$  and for every  $M > 0$  we have

$$\|R_m\| \leq \frac{\exp(M) \prod_{j=1}^m h_{j+1,j}}{\pi \nu^{m+1} M^{m\alpha+\beta-1}} \left( \frac{\mu}{\alpha} + \frac{\exp(-M(|\cos(\mu/\alpha)| + 1))}{m\alpha - 1 + \beta} \right).$$

Corollary (Moret and Novati 2011, Corollary 3.3)

Let assumptions (A1) and (A2) hold. Let  $m \geq 1$  be such that  $m\alpha + \beta > 0$ , then for every  $M > 0$ , we have

$$\|R_m\| \leq \frac{\exp(M) \prod_{j=1}^m h_{j+1,j}}{4 \nu^{m+1} M^{m\alpha}} \frac{4M^{1-\beta}}{\pi} \left( \frac{\mu}{\alpha} + \frac{\exp(-M(1 + |\cos(\mu/\alpha)|))}{M|\cos(\mu/\alpha)|} \right).$$

## A First Error Bound: some observations

---

⚙ The ML function is entire for  $\alpha > 0 \Rightarrow$  superlinear convergence for large enough  $m$ :

$$M = m\alpha + \beta - 1 \Rightarrow \|R_m\| \propto \left(\frac{\exp(1)}{M}\right)^M \nu^{-(m+1)} \prod_{j=1}^m h_{j+1,j}.$$

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- ⚙ To better understand this, we use that for every monic polynomial of degree  $m$  we find

$$\prod_{j=1}^m h_{j+1,j} \leq \|q_m(A)\nu\|,$$

Therefore, if we take  $q_m$  as the **monic Faber polynomial** associated to a closed convex subset  $\Omega \supset W(-A)$  we get the bound in terms of the **logarithmic capacity**  $\gamma$  of  $\Omega$ .

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- ⚙ To better understand this, we use that for every monic polynomial of degree  $m$  we find

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Therefore, if we take  $q_m$  as the **monic Faber polynomial** associated to a closed convex subset  $\Omega \supset W(-A)$  we get the bound in terms of the **logarithmic capacity**  $\gamma$  of  $\Omega$ .

$\Rightarrow$  we have discovered:

$$\|R_m\| \propto \left(\frac{\exp(1)}{m\alpha}\right)^{m\alpha} \left(\frac{\gamma}{\nu}\right)^m.$$

# Specialized bounds

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The bound can be refined under stricter hypotheses.

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Theorem (Moret and Novati 2011, Theorem 3.5)

Assume that  $A$  is Hermitian with  $\sigma(A) \subseteq [a, b] \subset [0, +\infty)$ . Assume that  $0 < \alpha < 1$ ,  $\beta \geq \alpha$ . Let  $\mu \leq \pi/2$ ,  $\alpha\pi/2 < \mu < \alpha\pi$ . Then for every index  $m \geq 1$  and for every  $M > 0$  we have

$$\|R_m\| \leq \frac{4M^{1-\beta}}{\pi} \left( \frac{\mu}{\alpha} + \frac{\exp(-M(1 + |\cos(\mu/\alpha)|))}{M|\cos(\mu/\alpha)|} \right) \exp(M) \Phi(u(M^\alpha \exp(i\mu)))^{-m}.$$

for  $\Phi(u) = u + \sqrt{u^2 - 1}$ ,  $u(z) = (|b+z| + |a+z|)/b-a$ .

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## Limiting relation

If  $\alpha \rightarrow 0$ ,  $\beta = 1$ , we have  $E_{0,1}(-z) = (1+z)^{-1}$ ,  $|z| < 1$ . Then setting  $\mu = \alpha\pi$  and letting  $M = 1$ , we find

$$\|R_m\| \leq \frac{4(\pi \exp(1) - \exp(-1))}{\pi \Phi(u(1))^m}$$



# The Shift-and-Invert Method (Moret and Novati 2011)

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We remain under the assumptions (A1) and (A2) and consider the matrix

$$Z = (I + hA)^{-1}, \quad h > 0,$$

together with the space  $\mathcal{K}_m(Z, \mathbf{v})$ .

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We can write the **analogous Arnoldi relation** for  $U_m = [\mathbf{u}_1, \dots, \mathbf{u}_m]$  spanning  $\mathcal{K}_m(Z, \mathbf{v})$ :

$$ZU_m = U_m S_m + s_{m+1,m} u_{m+1} \mathbf{e}_m^T, \quad S_m = U_m^H Z U_m.$$

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The **approximation** is then given by

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We can repeat the general error analysis using

$$R_m = E_{\alpha,\beta}(-A)\mathbf{v} - U_m E_{\alpha,\beta}(-B_m) \mathbf{e}_1 = \frac{1}{2\pi\alpha i} \int_{C(\varepsilon,\mu)} \exp(\lambda^{1/\alpha}) \lambda^{(1-\beta)/\alpha} b_m(\lambda) d\lambda,$$

for  $b_m(\lambda) = (\lambda I + A)^{-1} \mathbf{v} - U_m (\lambda I + B_m)^{-1} \mathbf{e}_1$ .

# Error bound (Moret and Novati 2011)

Theorem (Moret and Novati 2011, Theorem 4.3)

For every matrix  $A$  satisfying (A1) and (A2), assume  $0 < \alpha < 1$  and  $\beta \geq \alpha$ . Then, there exists a function  $g(h)$ , continuous in any bounded interval  $0 < h_1 \leq h \leq h_2$ , such that for  $m \geq 2$ ,

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## Theorem (Moret and Novati 2011, Theorem 4.5)

Assume that  $A$  is Hermitian with  $\sigma(A) \subseteq [a, +\infty)$ ,  $a \geq 0$ . Assume  $0 < \alpha \leq 2/3$  and  $\beta \geq \alpha$ . Then, for every  $m \geq 1$  we have

$$\|R_m\| \leq \frac{K_1 Q_m h^{\frac{\beta-1}{\alpha}}}{(1+\sqrt{2})^{m-1}} + \frac{K_2 h^{\beta/\alpha}}{(m-1)^2} \exp\left(-\frac{h^{-1/\alpha}}{\sqrt{2}}\right),$$

where  $Q_m = \max_{0 \leq |\varphi| \leq 3\alpha\pi/4} \exp\left(h^{-1/\alpha} \cos \varphi / \alpha\right) (1 - \cos \varphi)^{\frac{m-1}{2}}$ , with  $K_1, K_2$  constants.

# ML function, what have we found?

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- 🔧 To obtain a complete method one still has to find a way to repeatedly compute the matrix functions in

$$\mathbf{u}(t) = E_{\alpha,1}(-t^\alpha A)\mathbf{u}_0 + \int_0^t (t-s)^{\alpha-1} E_{\alpha,\alpha}(-A(t-s)^\alpha) \mathbf{g}(s) \, ds.$$

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## Other extensions

A variant with *restart* is discussed in (Moret and Popolizio [2014](#)), the combination with other matrix-functions in (Moret and Novati [2019](#)).





# Bibliography I

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-  Benzi, M. (2021). "Some uses of the field of values in numerical analysis". In: *Boll. Unione Mat. Ital.* 14.1, pp. 159–177. ISSN: 1972-6724. DOI: [10.1007/s40574-020-00249-2](https://doi.org/10.1007/s40574-020-00249-2). URL: <https://doi.org/10.1007/s40574-020-00249-2>.
-  Davies, P. I. and N. J. Higham (2003). "A Schur-Parlett algorithm for computing matrix functions". In: *SIAM J. Matrix Anal. Appl.* 25.2, pp. 464–485. ISSN: 0895-4798. DOI: [10.1137/S0895479802410815](https://doi.org/10.1137/S0895479802410815). URL: <https://doi.org/10.1137/S0895479802410815>.
-  Diele, F., I. Moret, and S. Ragni (2008/09). "Error estimates for polynomial Krylov approximations to matrix functions". In: *SIAM J. Matrix Anal. Appl.* 30.4, pp. 1546–1565. ISSN: 0895-4798. DOI: [10.1137/070688924](https://doi.org/10.1137/070688924). URL: <https://doi.org/10.1137/070688924>.
-  Garrappa, R. (2015). "Numerical evaluation of two and three parameter Mittag-Leffler functions". In: *SIAM J. Numer. Anal.* 53.3, pp. 1350–1369. ISSN: 0036-1429. DOI: [10.1137/140971191](https://doi.org/10.1137/140971191). URL: <https://doi.org/10.1137/140971191>.





# Bibliography II

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-  Garrappa, R. and M. Popolizio (2011). “On accurate product integration rules for linear fractional differential equations”. In: *J. Comput. Appl. Math.* 235.5, pp. 1085–1097. ISSN: 0377-0427. DOI: [10.1016/j.cam.2010.07.008](https://doi.org/10.1016/j.cam.2010.07.008). URL: <https://doi.org/10.1016/j.cam.2010.07.008>.
-  — (2018). “Computing the matrix Mittag-Leffler function with applications to fractional calculus”. In: *J. Sci. Comput.* 77.1, pp. 129–153. ISSN: 0885-7474. DOI: [10.1007/s10915-018-0699-5](https://doi.org/10.1007/s10915-018-0699-5). URL: <https://doi.org/10.1007/s10915-018-0699-5>.
-  Higham, N. J. and X. Liu (2021). “A multiprecision derivative-free Schur-Parlett algorithm for computing matrix functions”. In: *SIAM J. Matrix Anal. Appl.* 42.3, pp. 1401–1422. ISSN: 0895-4798. DOI: [10.1137/20M1365326](https://doi.org/10.1137/20M1365326). URL: <https://doi.org/10.1137/20M1365326>.
-  Metzler, R. and J. Klafter (2000). “The random walk’s guide to anomalous diffusion: a fractional dynamics approach”. In: *Phys. Rep.* 339.1, p. 77. ISSN: 0370-1573. DOI: [10.1016/S0370-1573\(00\)00070-3](https://doi.org/10.1016/S0370-1573(00)00070-3). URL: [https://doi.org/10.1016/S0370-1573\(00\)00070-3](https://doi.org/10.1016/S0370-1573(00)00070-3).

# Bibliography III

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-  Moret, I. and P. Novati (2011). “On the convergence of Krylov subspace methods for matrix Mittag-Leffler functions”. In: *SIAM J. Numer. Anal.* 49.5, pp. 2144–2164. ISSN: 0036-1429. DOI: 10.1137/080738374. URL: <https://doi.org/10.1137/080738374>.
-  — (2019). “Krylov subspace methods for functions of fractional differential operators”. In: *Math. Comp.* 88.315, pp. 293–312. ISSN: 0025-5718. DOI: 10.1090/mcom/3332. URL: <https://doi.org/10.1090/mcom/3332>.
-  Moret, I. and M. Popolizio (2014). “The restarted shift-and-invert Krylov method for matrix functions”. In: *Numer. Linear Algebra Appl.* 21.1, pp. 68–80. ISSN: 1070-5325. DOI: 10.1002/nla.1862. URL: <https://doi.org/10.1002/nla.1862>.
-  Podlubny, I. (1999). *Fractional differential equations*. Vol. 198. Mathematics in Science and Engineering. An introduction to fractional derivatives, fractional differential equations, to methods of their solution and some of their applications. Academic Press, Inc., San Diego, CA, pp. xxiv+340. ISBN: 0-12-558840-2.

# Bibliography IV

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Sokolov, I. M. and J. Klafter (2005). “From diffusion to anomalous diffusion: a century after Einstein’s Brownian motion”. In: *Chaos* 15.2, pp. 026103, 7. ISSN: 1054-1500. DOI: 10.1063/1.1860472. URL: <https://doi.org/10.1063/1.1860472>.