

A Spectral Element Based Parareal Method for the Long Time Integration of Time-Fractional Differential Equations

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- 1 Introduction
- 2 Spectral element method for TFDE
- 3 Parareal method for TFDE
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Introduction

We consider the long time integration of the time fractional differential equation (TFDE),

$$\begin{cases} {}_0D_t^\alpha u(t) = \lambda u(t) + f(t), & t \in \Omega, \\ u(0) = u_0, \end{cases} \quad (1)$$

where $\lambda \in \mathbb{C}$. ${}_0D_t^\alpha u(t)$ is the Caputo derivative of order α of function u ,

$${}_0D_t^\alpha u(t) \triangleq {}_0D_t^{\alpha-n} \frac{d^n u(t)}{dt^n} = \frac{1}{\Gamma(n-\alpha)} \int_0^t (t-s)^{n-1-\alpha} \frac{d^n u(s)}{ds^n} ds,$$

where n is an integer and $\alpha \in (n-1, n]$.

Challenges:

- 1 insure high accuracy
- 2 reduce computational cost

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- 1 insure high accuracy
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Solutions:

- 1 high order spectral element method
- 2 parallel-in-time (parareal) method

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Spectral collocation method based on Jacobian polynomial

Denote $J_j^{a,b}(t)$ the j -th order Jacobi polynomial with index (a, b) defined on $[-1, 1]$.

$\{J_j^{a,b}(t)\}_{j=0}^N$ satisfies the three terms recurrence relation:

$$J_0^{a,b}(t) = 1,$$

$$J_1^{a,b}(t) = (a + b + 2)t + (a - b),$$

$$J_{j+1}^{a,b}(t) = (A_j^{a,b}t - B_j^{a,b})J_j^{a,b}(t) - C_j^{a,b}J_{j-1}^{a,b}(t), \quad 1 \leq j \leq N-1,$$

where $A_j^{a,b}, B_j^{a,b}, C_j^{a,b}$ are known coefficients.

Fractional derivative of Jacobian polynomial

Consider the fractional integral of Jacobi polynomials,

$$\hat{J}_j^{a,b,\alpha}(t) \triangleq {}_{-1}D_t^{-\alpha} J_j^{a,b}(t) = \frac{1}{\Gamma(\alpha)} \int_{-1}^t (t-s)^{\alpha-1} J_j^{a,b}(s) ds$$

Following [C. Li, F. Zeng, and F. Liu, 2012, FCAA], $\{\hat{J}_j^{a,b,\alpha}(t)\}_{j=0}^N$ satisfies the three term relation:

$$\begin{aligned} \hat{J}_0^{a,b,\alpha}(t) &= \frac{(t+1)^\alpha}{\Gamma(\alpha+1)}; \\ \hat{J}_1^{a,b,\alpha}(t) &= \frac{a+b+2}{2} \left(\frac{t(t+1)^\alpha}{\Gamma(\alpha+1)} - \alpha \frac{(t+1)^{\alpha+1}}{\Gamma(\alpha+2)} \right) + \frac{a-b}{2} \hat{J}_0^{a,b,\alpha}(t); \\ \hat{A}_j^{a,b} \hat{J}_{j+1}^{a,b,\alpha}(t) &= \hat{B}_j^{a,b} \hat{J}_j^{a,b,\alpha}(t) + \hat{C}_j^{a,b} \hat{J}_{j-1}^{a,b,\alpha}(t) + \hat{D}_j^{a,b} (t+1)^\alpha, \end{aligned}$$

Fractional derivative of Jacobian polynomial

Using the relation

$$\frac{d^m}{dx^m} J_j^{a,b}(t) = d_{j,m}^{a,b} J_{j-m}^{a+m,b+m}(t), \quad m \geq j, \quad m \in \mathbb{N},$$

the fractional derivative of a Jacobi polynomial can be expressed as

$$(J_j^{a,b})^{(\alpha)}(t) \triangleq {}_{-1}D_t^\alpha J_j^{a,b}(t) = d_{j,n}^{a,b} \hat{J}_{j-n}^{a+n,b+n,n-\alpha}(t), \quad n-1 < \alpha < n. \quad (2)$$

Spectral collocation method

Assume $u(t)$ is a polynomial of degree no greater than N .
Then we have

$$u(t) = \sum_{i=0}^N \hat{u}_i J_i^{a,b}(t),$$

Hence, the fractional derivative of $u(t)$ is expressed as

$${}_{-1}D_t^\alpha u(t) = \sum_{j=0}^N \hat{u}_j (J_j^{a,b})^{(\alpha)}(t).$$

Matrix form

Let $\{t_i\}_{i=0}^N$ be a set of Gauss quadrature points, e.g., the L-G-L points. Define the fractional differentiation matrix $\mathcal{D}^\alpha$ as

$$\mathcal{D}_{ij}^\alpha = (J_j^{a,b})^{(\alpha)}(t_i),$$

Denote

$$\mathbf{U} = (u(t_0), u(t_1), \dots, u(t_N))^T, \hat{\mathbf{U}} = (\hat{u}(t_0), \hat{u}(t_1), \dots, \hat{u}(t_N))^T,$$

Then, equation (1) can be written as matrix form,

$$\mathcal{D}^\alpha \hat{\mathbf{U}} = \lambda \mathbf{U} + F$$

Matrix form

Suppose $\mathcal{J}$ is a $(N+1) \times (N+1)$ matrix defined by $\mathcal{J}_{ij} = J_j^{a,b}(t_i)$, then $\mathbf{U} = \mathcal{J}\hat{\mathbf{U}}$.

The matrix equation can be expressed as modal form

$$\mathcal{D}^\alpha \hat{\mathbf{U}} = \lambda \mathcal{J} \hat{\mathbf{U}} + F$$

and a nodal form,

$$\mathcal{D}^\alpha \mathcal{J}^{-1} \mathbf{U} = \lambda \mathbf{U} + F$$

Combined with the initial condition, both matrix equations can be solved uniquely.

Numerical example 1

Consider (1) with $\lambda = 0$,

$${}_{-1}D_t^\alpha u(t) = f(t), \quad t \in [-1, 1].$$

We choose $u(t) = \sin_M(t+1)$ as our test example, where $\sin_M(x)$ is Taylor of expansion of $\sin(x)$ up to degree M .

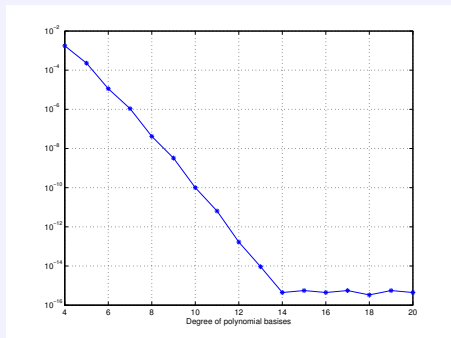


Figure: Maximum errors of the spectral collocation method with $\alpha = 0.5$.

Spectral Element Method for the TFDE

Let $\Omega = [t_l, t_r]$, which is divided into K elements:

$$t_l = t_0 < t_1 < \cdots < t_K = t_r, \Omega_i = [t_{i-1}, t_i], h_i = t_i - t_{i-1}, 1 \leq i \leq K.$$

For $t \in [t_i, t_{i+1}]$ and $n - 1 < \alpha \leq n$, we have

$$\begin{aligned} {}_0D_t^\alpha u(t) = & \frac{1}{\Gamma(n - \alpha)} \left(\sum_{k=1}^i \int_{\Omega_k} (t - s)^{n-1-\alpha} \frac{\partial^n u(s)}{\partial s^n} ds \right) \\ & + \frac{1}{\Gamma(n - \alpha)} \left(\int_{t_i}^t (t - s)^{n-1-\alpha} \frac{\partial^n u(s)}{\partial s^n} ds \right). \end{aligned}$$

The last term can be evaluated using (2). This induces a stiffness matrix, M_0 , which comes from $\mathcal{D}^\alpha$.

For each j , we define the following integral,

$$\tilde{J}_j^{a,b,\alpha}(t) \triangleq \frac{1}{\Gamma(\alpha)} \int_{-1}^1 (t-s)^{\alpha-1} J_j^{a,b}(s) ds, \quad t > 1.$$

How to compute $\tilde{J}_j^{a,b,\alpha}(t)$? Exact integral or numerical integral?

- 1 If $t \gg s$, Gauss quadrature;
- 2 If $t = 1$, formula (2) and Jacobian-Gauss quadrature;
- 3 If t close to 1, accuracy of both Gauss and Jacobian-Gauss quadrature would be reduced.

To address this, we return to the three terms recursive relation for Jacobian polynomials. We obtain a new three terms recursive relation for $\tilde{J}_j^{a,b,\alpha}(t)$:

$$\tilde{J}_0^{a,b,\alpha}(t) = \frac{(t+1)^\alpha - (t-1)^\alpha}{\Gamma(\alpha+1)},$$

$$\tilde{J}_1^{a,b,\alpha}(t) = \frac{a+b+2}{2\Gamma(2+\alpha)} ((t-\alpha)(t+1)^\alpha - (t+\alpha)(t-1)^\alpha) + \frac{a-b}{2} \tilde{J}_0^{a,b,\alpha}(t),$$

$$\tilde{A}_j^{a,b} \tilde{J}_{j+1}^{a,b,\alpha}(t) = \tilde{B}_j^{a,b} \tilde{J}_j^{a,b,\alpha}(t) + \tilde{C}_j^{a,b} \tilde{J}_{j-1}^{a,b,\alpha}(t) + \tilde{D}_j^{a,b} (t+1)^\alpha + \tilde{E}_j^{a,b} (t-1)^\alpha,$$

where $\tilde{A}_j^{a,b}, \tilde{B}_j^{a,b}, \tilde{C}_j^{a,b}, \tilde{D}_j^{a,b}, \tilde{E}_j^{a,b}$ are the coefficients.

Then, the integral over element of $(J_j^{a,b})^{(n)}(s)$ can be computed using

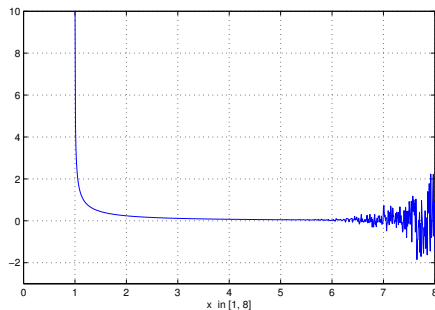
$$\begin{aligned} \tilde{D}^\alpha J_j^{a,b}(t) &\triangleq \frac{1}{\Gamma(n-\alpha)} \int_{-1}^1 (t-s)^{n-\alpha-1} (J_j^{a,b})^{(n)}(s) ds \\ &= d_{j,n}^{a,b} \tilde{J}_{j-n}^{a+n,b+n,n-\alpha}(t), \quad t > 1. \end{aligned} \quad (3)$$

How well does this work?

We compute

$$\tilde{J}_{14}^{a,b,\alpha}(t) = \frac{1}{\Gamma(\alpha)} \int_{-1}^1 (t-s)^{\alpha-1} J_{14}^{a,b}(s) ds,$$

for $t \in [1, 8]$, $\alpha = 0.5$. The value of $\tilde{J}_{14}^{a,b,\alpha}(t)$ is shown in the figure.



Hybrid strategy

- smooth part : Gauss quadrature;
- near singular part : three term recurrence relation.

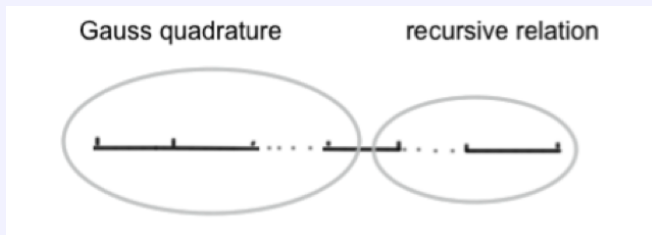


Figure: Hybrid strategy

Denote $\mathbf{U}_k$ as the vector of function values at all Gauss points in element k ,

$$\mathbf{U}_k = (u_{k,0}, u_{k,1}, \dots, u_{k,N})^T$$

Then,

$$M_0 \mathbf{U}_k + M_1 \mathbf{U}_{k-1} + \dots + M_{k-1} \mathbf{U}_1 = \lambda \mathbf{U}_k + \mathbf{f}_k, \quad k = 1, 2, \dots, K.$$

$$(\lambda \mathcal{I} - M_0) \mathbf{U}_k = M_1 \mathbf{U}_{k-1} + \dots + M_{k-1} \mathbf{U}_1 - \mathbf{f}_k, \quad k = 1, 2, \dots, K.$$

Combined with the initial condition for each element, the equation can be solved uniquely.

Analysis

For the approximation of fractional derivative, we have the following lemma,

Lemma 2.1

(Q. Xu, J. S. Hesthaven, 2012) Suppose $u(x)$ is a smooth function defined in $\Omega = [a, b]$. Ω_h is a discretization of the domain with interval weight h , $u_h(x)$ is a piecewise interpolation of u in $P_h^N = \text{span}\{1, x, \dots, x^N\}$. $\forall x \in I_i$, u_h interpolate u at all Gauss points. Then we have

$$\| {}_a D_x^\alpha u - {}_a D_x^\alpha u_h \|_{L_2} \leq \begin{cases} C, & n > N \\ Ch^{N+1-n}, & n \leq N, N-n \text{ odd} \\ Ch^{N+1-\alpha}, & n \leq N, N-n \text{ even} \end{cases}$$

where $n-1 < \alpha < n$, C is a constant independent of x .

Convergence

Since u is solved element by element, for $0 < \alpha < 1$ the following result is obtained.

Proposition 2.1

Suppose u is a smooth function, u_h is the numerical solution of equation (1) using spectral elements method based on Gauss points, $0 < \alpha < 1$, $e = u - u_h$,

$$\|e\|_2 \leq \begin{cases} Ch^N, & N \text{ is even} \\ Ch^{N+1-\alpha}, & N \text{ is odd} \end{cases}$$

Numerical examples

Consider the following equation,

$${}_0D_t^\alpha u(t) = -u(t) + f(t), \quad t \in \Omega = [0, 3\pi],$$

We choose the exact solution to be $u(t) = \sin_M(t)$.

$N = 1$						
K	$\alpha = 0.1$		$\alpha = 0.4$		$\alpha = 0.8$	
	$\ e_h\ _2$	Order	$\ e_h\ _2$	Order	$\ e_h\ _2$	Order
20	6.70e-04	1.68	5.15e-03	1.50	2.74e-02	1.18
40	2.09e-04	1.69	1.82e-03	1.51	1.21e-02	1.18
60	1.05e-04	1.66	9.87e-04	1.50	7.51e-03	1.19
80	6.54e-05	-	6.42e-04	-	5.34e-03	-
$N = 2$						
K	$\alpha = 0.1$		$\alpha = 0.4$		$\alpha = 0.8$	
	$\ e_h\ _2$	Order	$\ e_h\ _2$	Order	$\ e_h\ _2$	Order
20	3.33e-06	2.94	2.42e-05	2.89	1.88e-04	2.27
40	4.33e-07	2.85	3.27e-06	2.90	3.90e-05	2.25
60	1.36e-07	2.75	1.01e-06	2.84	1.56e-05	2.26
80	6.18e-08	-	4.45e-07	-	8.18e-06	-

Numerical examples

$N = 3$						
K	$\alpha = 0.1$		$\alpha = 0.4$		$\alpha = 0.8$	
	$\ e_h\ _2$	Order	$\ e_h\ _2$	Order	$\ e_h\ _2$	Order
20	3.10e-08	3.94	2.82e-07	3.81	2.40e-06	3.33
40	2.01e-09	3.95	2.02e-08	3.81	2.39e-07	3.29
60	4.06e-10	3.94	4.31e-09	3.80	6.30e-08	3.27
80	1.31e-10	-	1.45e-09	-	2.46e-08	-
$N = 4$						
K	$\alpha = 0.1$		$\alpha = 0.4$		$\alpha = 0.8$	
	$\ e_h\ _2$	Order	$\ e_h\ _2$	Order	$\ e_h\ _2$	Order
20	2.31e-10	4.96	1.73e-09	4.83	1.54e-08	4.41
40	7.42e-12	4.96	6.08e-11	4.74	7.28e-10	4.35
60	9.93e-13	4.96	8.91e-12	4.45	1.25e-10	4.31
80	2.38e-13	-	2.48e-12	-	3.61e-11	-
$N = 5$						
K	$\alpha = 0.1$		$\alpha = 0.4$		$\alpha = 0.8$	
	$\ e_h\ _2$	Order	$\ e_h\ _2$	Order	$\ e_h\ _2$	Order
20	1.51e-12	5.82	1.26e-11	5.85	1.22e-10	5.40
40	2.66e-14	-	2.20e-13	5.72	2.89e-12	5.33
60	2.89e-15	-	2.16e-14	-	3.34e-13	5.12
80	1.23e-14	-	5.88e-15	-	7.66e-14	-

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Parareal algorithm for ODEs

The parallel-in-time approach, called parareal method, was proposed in (Lions, Maday, and Turinici, 2001).

Consider the following systems of ODEs:

$$\begin{aligned}u_t &= f(u, t), \quad t \in \Theta := (0, T], \\u(0) &= u_0,\end{aligned}$$

The parareal algorithm is defined using two propagation operators.

- A coarse solver $\mathcal{G}(t_2, t_1, u_1)$, provides a cheap but inaccurate approximation to $u(t_n)$ with $u(t_{n-1}) = u_{n-1}$,
- A fine solver $\mathcal{F}(t_2, t_1, u_1)$, provides an expensive but accurate approximation of $u(t_n)$ from $u(t_{n-1}) = u_{n-1}$.

First, the time domain is decomposed into K elements:

$$0 = T_0 < \cdots < T_n < \cdots < T_K = T, \quad T_n = n\Delta T, \quad \Delta T = \frac{T}{K}.$$

Assume that u_n^k is the k 'th iterated approximation to u_n . The process of parareal algorithm is shown as,

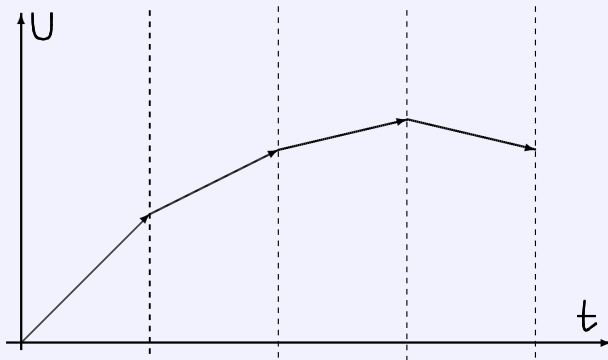
- 1 Initial guess : compute u_n^0 using coarse solver,
- 2 Assuming that u_n^k is known, compute the solution in parallel using the fine solver for each element,
- 3 The parareal iteration

$$u_{n+1}^{k+1} = \mathcal{G}u_n^{k+1} + \mathcal{F}u_n^k - \mathcal{G}u_n^k.$$

- 4 The algorithm is terminated if $|u_n^k - u_n^{k-1}| \leq \tau$.

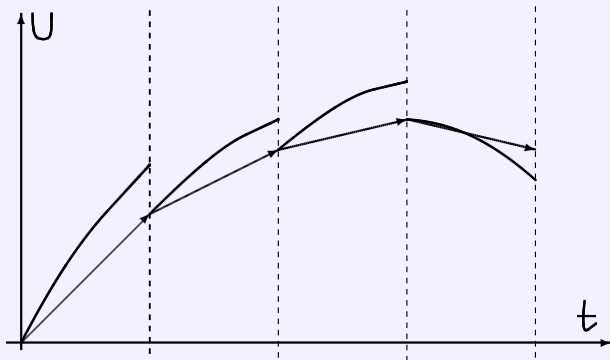
Parareal method for TFDE

1. Initial guess using coarse solver, (serial)



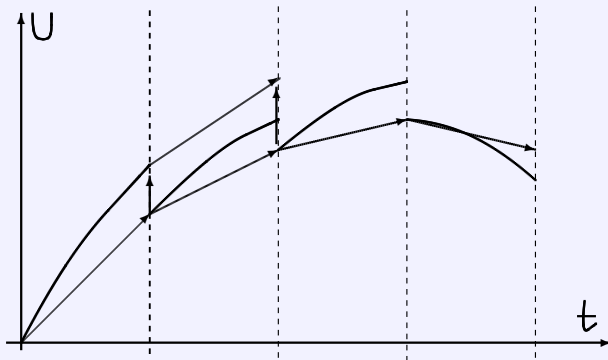
Parareal method for TFDE

2. Using fine solver in each element, (parallel)



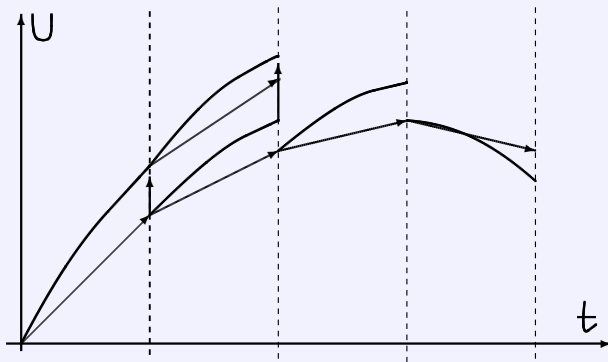
Parareal method for TFDE

3. Compute the new solution using coarse solver



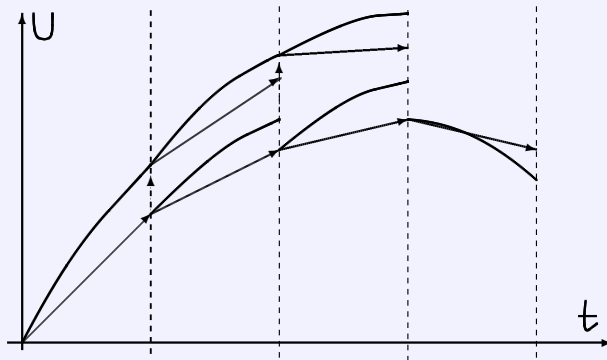
Parareal method for TFDE

Correct the solution



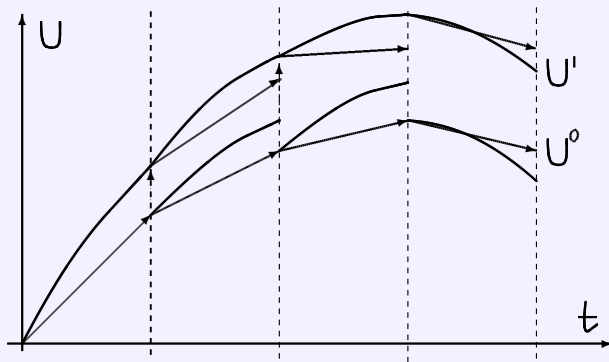
Parareal method for TFDE

Advance iteration



Parareal method for TFDE

Advance iteration



Parareal method for TFDE

Consider the fractional ordinary differential equation with $\alpha < 1$

$$\begin{cases} {}_0D_t^\alpha u(t) = \lambda u(t), & t \in \Omega, \\ u(0) = u_0, \end{cases}$$

Both the coarse solver and the fine solver are defined based on our spectral element method in this talk.

- Coarse grid : $N_c + 1$ collocation points.
- Fine grid: $N_f + 1$ collocation points.
- Memory part : $U_{0:n} = \{U_0, U_1, \dots, U_n\}$.
- Denote v_n as the value of $u(t)$ at right interface of element n , that's $v_n = U_{n,N} = U_{n+1,0}$.

Challenge 1: Dealing with the memory part.

We propose the following parareal method for TFDE,

$$U_{n+1}^{k+1} = \mathcal{G}U_{0:n}^{k+1} + \mathcal{F}U_{0:n}^k - \mathcal{G}U_{0:n}^k, \quad 0 \leq n \leq N-1.$$

Challenge 1: Dealing with the memory part.

We propose the following parareal method for TFDE,

$$U_{n+1}^{k+1} = \mathcal{G}U_{0:n}^{k+1} + \mathcal{F}U_{0:n}^k - \mathcal{G}U_{0:n}^k, \quad 0 \leq n \leq N-1.$$

Challenge 2: Convergence.

The spectral element method above,

- fine solution converges to exact solution as $\Delta t \rightarrow 0$ or $K \rightarrow \infty$.
- Will this parareal solution converges to fine solution as iteration number k increase?

Now we consider the values at each interface of elements, v_n .

We have

$$U_n = \sum_{j=1}^{n-1} (\lambda \mathcal{I} - M_0)^{-1} M_j U_{n-j}, \quad n = 1, 2, \dots, K.$$

For the coarse grid, we define an operator $\mathcal{R}$, s.t.

$$v_n^k = \sum_{j=1}^n \mathcal{R}_j v_{n-j}^k$$

where k denotes k' th iteration, v_0 is provided by the initial condition.

The fine integration operator $\mathcal{R}_j^{\mathcal{F}}$ can be defined accordingly.

Introduce the parareal method for v_n^{k+1} ,

$$v_n^{k+1} = \sum_{j=1}^n \mathcal{R}_j v_{n-j}^{k+1} + \sum_{j=1}^n \mathcal{R}_j^{\mathcal{F}} v_{n-j}^k - \sum_{j=1}^n \mathcal{R}_j v_{n-j}^k, \quad 0 \leq n \leq K-1.$$

Denote $\mathbf{v}^k = (v_0^k, v_1^k, \dots, v_K^k)^T$

$$\mathcal{M} = \begin{pmatrix} 0 & 0 & \cdots & 0 \\ \mathcal{R}_1 & 0 & \cdots & 0 \\ \vdots & \cdots & \ddots & 0 \\ \mathcal{R}_{K+1} & \cdots & \mathcal{R}_1 & 0 \end{pmatrix}$$

$$\mathcal{N} = \begin{pmatrix} 0 & 0 & \cdots & 0 \\ \mathcal{R}_1^{\mathcal{F}} - \mathcal{R}_1 & 0 & \cdots & 0 \\ \vdots & \cdots & \ddots & 0 \\ \mathcal{R}_{K+1}^{\mathcal{F}} - \mathcal{R}_{K+1} & \cdots & \mathcal{R}_1^{\mathcal{F}} - \mathcal{R}_1 & 0 \end{pmatrix}$$

then we get the matrix form,

$$\mathbf{v}^{k+1} = \mathcal{M}\mathbf{v}^{k+1} + \mathcal{N}\mathbf{v}^k, \quad 0 \leq n \leq K-1.$$

Assume $u(t_n)$ = fine solution at t_n , then $u(t_n) = \sum_{j=1}^n \mathcal{R}_j^{\mathcal{F}} u(t_{n-j})$.

Denote $e_n^k = u(t_n) - v_n^k$ and $e^k = (e_0^k, e_1^k, \dots, e_K^k)^T$, then

$$\begin{aligned} e_n^{k+1} &= u(t_n) - v_n^{k+1} \\ &= \sum_{j=1}^n \mathcal{R}_j^{\mathcal{F}} u(t_{n-j}) - \sum_{j=1}^n \mathcal{R}_j v_{n-j}^{k+1} - \sum_{j=1}^n \mathcal{R}_j^{\mathcal{F}} v_{n-j}^k + \sum_{j=1}^n \mathcal{R}_j v_{n-j}^k \\ &= \sum_{j=1}^n (\mathcal{R}_j^{\mathcal{F}} - \mathcal{R}_j) e_{n-j}^k + \sum_{j=1}^n \mathcal{R}_j e_{n-j}^{k+1} \end{aligned}$$

In matrix form, we have

$$e^{k+1} = \mathcal{M}e^{k+1} + \mathcal{N}e^k$$

Since $\mathcal{M}, \mathcal{N}$ are strictly lower triangular matrix,

$$\begin{aligned} e^{k+1} &= \mathcal{N} \left(\sum_{j=0}^K \mathcal{M}^j \right) e^k \\ &= \mathcal{N}^{k+1} \left(\sum_{j=0}^K \mathcal{M}^j \right)^{k+1} e^0 \end{aligned}$$

Assume the convergence order of coarse solver is $p + 1$, then $(\mathcal{R}^{\mathcal{F}} - \mathcal{R}) \sim (\Delta t)^{p+1}$, $\|\mathcal{N}^k e^0\| \sim (K(\Delta t)^{p+1})^k \sim (\Delta t)^{kp}$.

If $|(\sum_{j=0}^K \mathcal{M}^j)^k|$ is bounded, the parareal solution converges to the fine solution with convergence order $(\Delta t)^{kp}$, (k is number of iterations).

For the case $N_c = 1$. Our method reduce to L_1 scheme. The distribution of $\mathcal{R}_j$ with $K = 20, \alpha = 0.5$ is shown in the following figure.

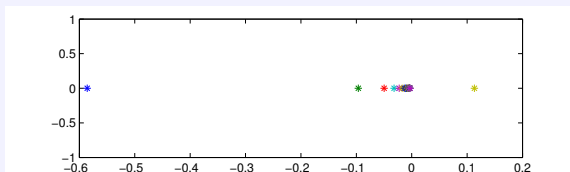


Figure: The distribution of $\mathcal{R}_j$

- 1 For $N_c = 1, \lambda < 0$, $|(\sum_{j=0}^K \mathcal{M}^j)^k|$ is bounded.
- 2 For a high order coarse solver ($N_c \geq 2$), convergence of parareal method can be verified numerically.

Assume $u_e(t)$ is the exact solution. Let $\eta_n = u_e(t_n) - u(t_n)$, $\epsilon_n^k = u_e(t_n) - v_n^k$, then $\|\epsilon^k\| \leq C(\|e^k\| + \|\eta\|)$.

We consider

$${}_0D_t^\alpha u(t) = -u(t) + f(t), \quad t \in [0, T].$$

To test the numerical accuracy, we choose the exact solution $u_e(t) = \sin_{60}(t)$. $f(t)$ can be calculated according to this setting.

Example 3.1

Let $T = 2\pi$. The number of processors $N_p = 10$, and $N_F = 12$. $N_C = 2, 4, 6$. The absolute error $\max |v^k - u_e|$ is shown as,

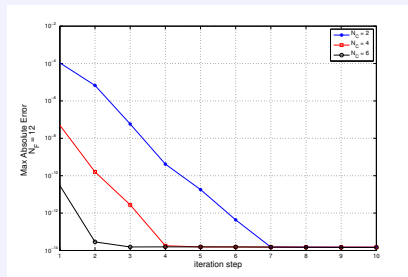


Figure: The accuracy of the parareal method in Example 3.1.

Example 3.2

Let $f(t) = \sin(t)$, $T = 15\pi$, $\alpha = 0.5$. The number of processors $N_p = 20$, and $N_F = 22$. $N_C = 1$. $e^k = \max |v^k - u|$.

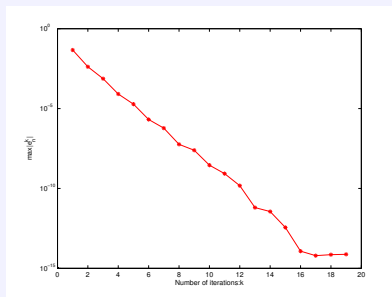


Figure: Convergence

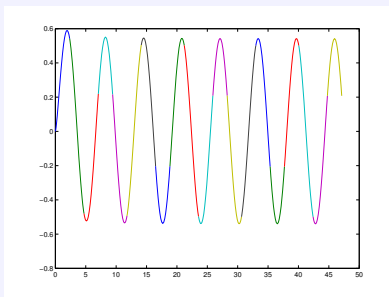


Figure: The parareal solution

Example 3.3

Let $f(t) = \sin(t)$, $T = 15\pi$, $\alpha = 0.1, 0.5, 0.9$. The number of processors $N_p = 20$, and $N_F = 16$. $N_C = 1$. The error $e^k = \max |v^k - u|$.

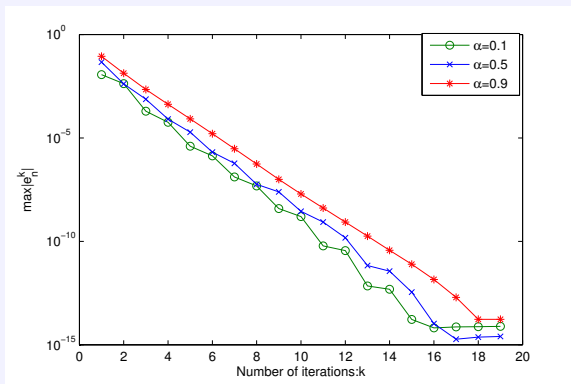


Figure: The error for different α

Example 3.4

Let $f(t) = \sin(t)$, $T = 15\pi$, $\alpha = 0.6$. The number of processors $N_p = 20$, and $N_F = 16$. $N_C = 1, 2, 3, 4$. The error $e^k = \max |v^k - u|$.

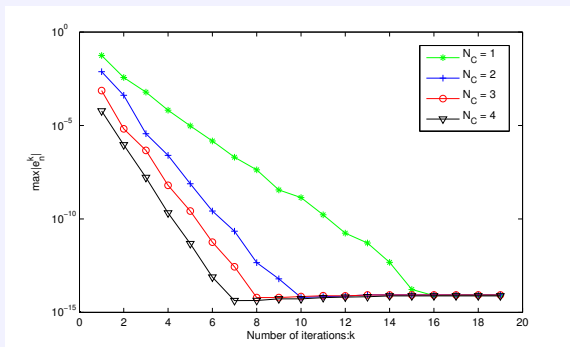


Figure: The error for different coarse solvers

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Parareal method for subdiffusion equation

We consider the following equation with periodic boundary condition,

$${}_0D_t^\alpha u(x, t) = \frac{\partial^2 u(x, t)}{\partial x^2}, \quad x \in [0, 2\pi], t \in [0, T].$$

with initial condition $u(x, 0) = \sin(x)$. The exact solution is $u(x, t) = E_\alpha(-t^\alpha) \sin(x)$, where $E_\alpha(t)$ is Mittag-Leffler function with one parameter.

Using Fourier transform in space, this equation becomes a system of decoupled ODEs for each Fourier mode ω .

$${}_0D_t^\alpha \hat{u} = -\omega^2 \hat{u}.$$

Parareal method for subdiffusion equation

Example 4.1

Let $T = 2, \alpha = 0.2$. The number of processors $N_p = 30$, and $N_F = 16$. $N_C = 2$. Number of Fourier modes $m = 40$. The error $e^k = \max |v^k - u|$.

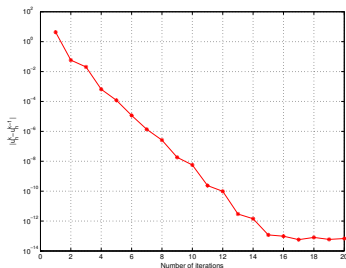


Figure: Convergence

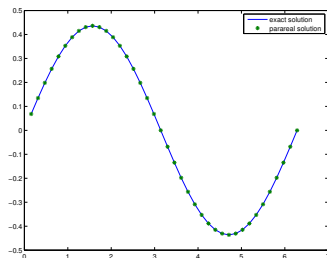


Figure: The parareal solution

Parareal method for subdiffusion equation

We consider the parareal method for long time integration of subdiffusion equation.

Example 4.2

Let $T = 100$, $\alpha = 0.2, 0.5, 0.8$. The number of processors $N_p = 30$, and $N_F = 16$. $N_C = 2$, Number of Fourier modes $m = 60$.

We consider the error between parareal solution and fine solution.

Parareal method for subdiffusion equation

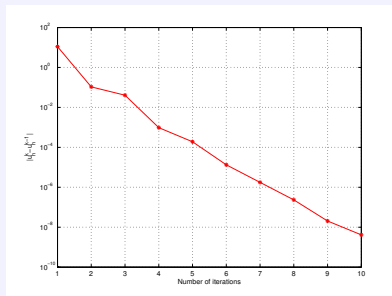


Figure: Convergence for $\alpha = 0.2$

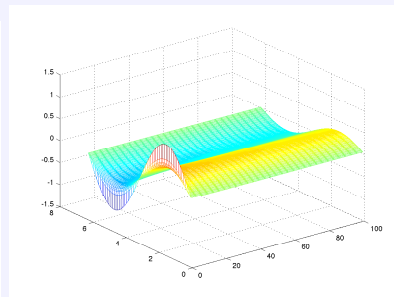


Figure: The parareal solution for $\alpha = 0.2$

Parareal method for subdiffusion equation

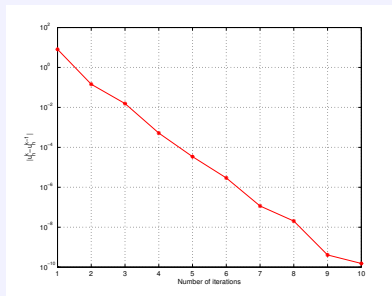


Figure: Convergence for $\alpha = 0.5$

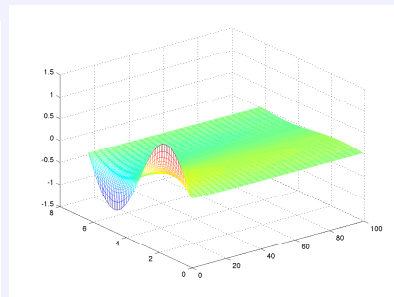


Figure: The parareal solution for $\alpha = 0.5$

Parareal method for subdiffusion equation

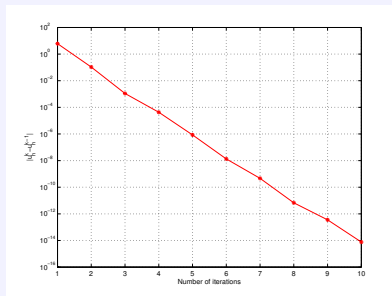


Figure: Convergence for $\alpha = 0.8$

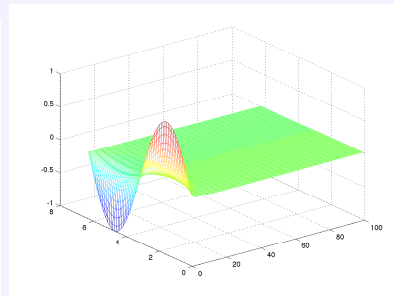


Figure: The parareal solution for $\alpha = 0.8$

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Summary

- 1 hybrid strategy for near singular integration of memory part;
- 2 spectral element method for TFDE and the convergence;
- 3 parareal method for time fractional differential equation;
- 4 convergence of parareal method for TFDE.

Future work

- 1 Analysis of high order coarse solver;
- 2 Parareal method for fractional diffusion-wave equation.

Thanks for your attention!