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OPERATORIAL TAU METHOD FOR FRACTIONAL DIFFERENTIAL EQUATIONS

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Abstract. The paper extends the applicability of our Matlab package *Chebpack* to find Chebyshev polynomial approximations of the solutions of some fractional differential equations. Numerical examples confirm the efficiency of this approach.

Keywords: fractional differential equations; Chebyshev polynomials; tau method.

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1. Introduction

The field of fractional calculus is almost as old as calculus itself. The non-integer derivatives are non-local operators which possess a memory effect explaining their significant uses in many applications such as the behavior of viscoelastic materials or polymers, for example.

Several methods to solve the fractional derivatives, fractional integrals and fractional differential equations have recently been proposed. Many interesting Matlab functions were already submitted to the *MathWorks*, *Inc. Matlab Central File Exchange*, where they are freely downloadable for sharing among the users.

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We refer here to the well-documented programs of Podlubny [24], [25] for fractional differential equations and [26] for distributed order ODEs, based on the Grünwald–Letnikov derivative. Another useful Matlab program is fde12.m of Garrappa [15] for (non-linear) differential equations of fractional order, based on the predictor-corrector method of Adams-Bashforth-Moulton described in Diethelm's paper [11].

We must remark that any algorithm using a discretization of a non-integer derivative has to take into account its non-local structure which imply, in general, high storage requirements and a great overall complexity of the algorithm.

There is a comprehensive literature on the theory, applications and numerical methods of the fractional calculus and the number of such textbooks and papers has grown rapidly in the recent years. The main sources used in this paper were [17], [9], [19], [23], [3], [14], [33], [22] and the references therein.

The aim of the presented paper is to extend the capabilities of our Matlab package *Chebpack* [32] for fractional calculus problems. *Chebpack* is based on the operational form of the Chebyshev spectral tau method [20], [8], [7]. The main advantage of *Chebpack* is a unified approach for initial value problems, boundary value problems, eigenproblems, nonlocal problems for ordinary, fractional or distributed order differential equations.

The Chebyshev polynomials were previously used for approximating the fractional derivatives in the Matlab package *Chebfun* of Trefethen [30]. Chebyshev and other kind of orthogonal polynomials were used for fractional differential equations in the papers [18], [16], [12], [13], [28], [4], [5], [6].

The paper is structured as follows: in section 2 we describe the tau method, in section 3 we give the discretization of the main fractional operators and in section 4 we give numerical examples to illustrate the facilities of *Chebpack* for fractional calculus. Finally, the paper ends with conclusions and references.

All the necessary Matlab code for reproducing the examples are now part of an updated version of *Chebpack* [32], in the folder *Examples*, subfolder *Fractional differential equations*.

2. The tau method

If we have to solve a differential or an integral problem, one of the most effective methods is to use the Chebyshev spectral methods [7], i.e. to approximate the solution *y* by a finite sum of a Chebyshev series

$$y(x) \approx \frac{1}{2}c_0T_0(x) + c_1T_1(x) + \dots + c_{n-1}T_{n-1}(x), x \in [-1,1].$$

Here $T_k(x) = \cos(k\cos^{-1}(x)), k = 0, 1, 2, ..., n - 1$ are the Chebyshev polynomials of the first kind and the coefficients $\mathbf{c} = \{c_k, k = 0, 1, ..., n - 1\}$ represented as a column vector are the unknowns. If $\mathscr{L} : C^{\infty}(-1, 1) \to C^{\infty}(-1, 1)$ is a linear operator then let \mathbf{u} be the corresponding coefficients of $\mathscr{L}(y)(x)$. The matrix $L : \mathbb{R}^n \to \mathbb{R}^n$ that maps \mathbf{c} into $\mathbf{u} = L\mathbf{c}$ is the Chebyshev approximation of \mathscr{L} .

Our Matlab package *Chebpack* [31], [32] implements the Chebyshev spectral method as a *Lanczos' tau method*, where we work in the spectral space of the coefficients. The linear operators of the differential or integral problem, such as differentiation, integration, product with the independent variable, modified argument, are discretized to their corresponding approximating matrices. The final form of the linear problem $\mathscr{L}(y)(x) = f(x)$ becomes, after the discretization, a pure algebraic linear problem in an *operational form* $L \cdot \mathbf{c} = \mathbf{f}$ with the supplementary conditions of the continuous problem included.

Consequently, *Chebpack* assumes the representation of the unknown functions in truncated Chebyshev polynomials series and the corresponding matrices L are given to act directly on the Chebyshev coefficients. Generally, this kind of discretization is performed by using the recurrence formula and the fast schemes of computation with Chebyshev polynomials, see [31] for more technical details on this topic.

Precisely, a sufficiently well-behaved function y(x) can be approximated by its physical representation $y(x_1)$, $y(x_2)$, ..., $y(x_n)$ of values of y at the given gridpoints **x**, for example at the *Chebyshev points of the first or of the second kind*

(2.1)
$$x_k^{(1)} = -\cos\frac{(2k-1)\pi}{2n}, x_k^{(2)} = -\cos\frac{(k-1)\pi}{n-1}, \ k = 1, ..., n$$

or by its spectral representation $\mathbf{c} = \{c_0, c_1, ..., c_{n-1}\}$ where

(2.2)
$$y_{n-1}(x) = \frac{c_0}{2}T_0(x) + c_1T_1(x) + \dots + c_{n-1}T_{n-1}(x)$$

is the unique polynomial obtained by interpolating y(x) through the points $x_1, ..., x_n$.

Of course, the Chebyshev polynomials are defined on dom = [-1, 1] but any interval dom = [a, b] can be shifted to [-1, 1] and we may use the shifted Chebyshev polynomials

(2.3)
$$T_k^*(x) = T_k(\alpha x + \beta), \ x \in [a,b], \ \alpha = \frac{2}{b-a}, \ \beta = -\frac{b+a}{b-a}, \ k = 0, 1, ..., n-1.$$

The code [x, w] = pd(n, dom, kind) of *Chebpack* calculates the *n* Chebyshev points **x** of the kind "kind" on dom as a column vector and the *n* quadrature weights **w** as a row vector for the Clenshaw-Curtis quadrature formula

$$\int_{a}^{b} y(x) dx \simeq \sum_{j=1}^{n} w_{j} y\left(x_{j}\right) \equiv \mathbf{w} \cdot y\left(\mathbf{x}\right).$$

The *fast* conversion between the spectral representation **c** of a function y and its physical values $\mathbf{v} = y(\mathbf{x})$ is performed by the functions v=t2x(c, kind) and c=x2t(v, kind) based on the Fast Chebyshev Transform.

If a function y is given by its Chebyshev coefficients c and we need its values at some points $\mathbf{x}_c \in dom$, a conversion matrix T_c is obtained from the code Tc=cpv(n, xc, dom), where

(2.4)
$$T_c = [T_0/2, T_1(\xi), ..., T_{n-1}(\xi)], \ \xi = \frac{2\mathbf{x}_c}{b-a} - \frac{b+a}{b-a} \in [-1, 1]$$

and ξ is transposed as a column vector. The code is based on the recurrence relations

(2.5)
$$T_1(x) = xT_0, \ T_{k+1}(x) = 2xT_k(x) - T_{k-1}(x), \ k = 1, 2, \dots$$

and we have $\mathbf{v}_c = T_c \mathbf{c}$. If T = cpv(n, x, dom) where **x** are the Chebyshev points, then the matrix *T* performs the (*non fast*, but useful if $n \le 256$) conversion between the coefficients **c** of the function *y* and the values $\mathbf{v} = y(\mathbf{x})$ through the formulas $\mathbf{v} = T\mathbf{c}$ and $\mathbf{c} = T^{-1}\mathbf{v}$.

It is important to remark that linear operators are better represented in the spectral space of the coefficients while the nonlinear operators are easily handled in the physical space of the values, see *Example 5*.

Another useful code is X=mult (n, dom) also based on the recurrence relations (2.5). If c is the column of Chebyshev coefficients of a function y(x), then $X \cdot c$ is the column of the Chebyshev coefficients of the multiplication by the independent variable xy(x) for $x \in dom$. If a is

the vector of the first *m* Chebyshev coefficients of a function a(x), then the vector of the coefficients of a(x)y(x) is $A \cdot \mathbf{c}$ where the multiplication matrix A is given by the code A=fact (a, m) based on the relations

(2.6)
$$2T_j(x)T_k(x) = T_{j+k}(x) + T_{|k-j|}(x), j, k = 0, 1, \dots$$

The differentiation matrix D is given by the code D=deriv(n, dom). If **c** is the column of Chebyshev coefficients of a function y(x), then $D \cdot \mathbf{c}$ is the column of the Chebyshev coefficients of the derivative $\frac{dy}{dx}$. The definition of D is based on the relations

(2.7)
$$T_0 = T'_1, \ T_1 = \frac{T'_2}{4}, \ T_k = \frac{T'_{k+1}}{2(k+1)} - \frac{T'_{k-1}}{2(k-1)}, \ k = 2, 3, ..., x \in [-1,1].$$

Similarly, the code [J, J0] = prim(n, dom) calculates the integration matrix J such that the coefficients of a primitive of y(x) are $J \cdot \mathbf{c}$. The coefficients of the particular primitive vanishing at a = dom(1) are obtained by using $J_0 \cdot \mathbf{c}$.

All the above codes take into account a general dom, see [31], [32] for more details.

3. The fractional operators

The operators of the fractional calculus can also be expressed by matrices. Let us start with *the Riemann–Liouville fractional integral operator of order q*

(3.1)
$$J^{q}y(x) = \frac{1}{\Gamma(q)} \int_{0}^{x} \frac{y(t)dt}{(x-t)^{1-q}},$$

where 0 < q < 1, $0 \le x \le b$. If $y \in L_1[0,b]$ is a sufficiently well-behaved function, see [9], then it has the spectral approximation

$$y(x) \simeq \sum_{k=0}^{n-1} c_k T_k \left(\alpha x - 1 \right),$$

where $\alpha = 2/b$ and the prime sign denotes the summation whose first term is halved. Then

$$J^{q}y(x_{j}) \simeq \frac{1}{\Gamma(q)} \sum_{k=0}^{n-1} c_{k} \int_{0}^{x_{j}} \frac{T_{k}(\alpha t-1) dt}{(x_{j}-t)^{1-q}}, \ j=1,...,n$$

approximates the physical values \mathbf{v} of $J^q y(\mathbf{x})$ at the Chebyshev points $\mathbf{x} = (x_j)_{j=1,...,n}$. Consequently, the spectral approximation of $J^q y(x)$ is given by the Chebyshev coefficients $T^{-1}\mathbf{v} \equiv I \cdot \mathbf{c}$, where *T* is given by (2.4), \mathbf{c} is the column vector of the coefficients $c_0, ..., c_{n-1}$ and *I* is the

basic integration matrix. If dom=[0,b] and x is given by x=pd(n, dom, kind), then the code

for 0 < q < 1 calculates the above matrix *I*.

The calculation starts with the elements I_{jk} , j = 1, 2, ..., n, k = 0, 1, ..., n - 1,

(3.2)
$$I_{jk} = \int_0^{x_j} \frac{T_k(\alpha t - 1) dt}{(x_j - t)^{1-q}}.$$

We have

$$\int_0^x \frac{dt}{(x-t)^{1-q}} = \frac{x^q}{q}, \quad \int_0^x \frac{tdt}{(x-t)^{1-q}} = \frac{x^{q+1}}{q(q+1)},$$
$$\int_0^x \frac{t^2dt}{(x-t)^{1-q}} = \frac{2x^{q+2}}{q(q+1)(q+2)}.$$

Consequently, the columns $I_{\cdot,k}$ for k = 0, 1, 2 are

$$I_{,0} = \frac{1}{2} \int_0^x \frac{dt}{(x-t)^{1-q}} = \frac{x^q}{2q}, I_{,1} = \int_0^x \frac{(\alpha t-1)dt}{(x-t)^{1-q}} = \alpha \frac{x^{q+1}}{q(q+1)} - \frac{x^q}{q},$$
$$I_{,2} = \int_0^x \frac{\left[2(\alpha t-1)^2 - 1\right]dt}{(x-t)^{1-q}} = 4\alpha^2 \frac{x^{q+2}}{q(q+1)(q+2)} - 4\alpha \frac{x^{q+1}}{q(q+1)} + \frac{x^q}{q}.$$

The columns for k > 2 are calculated iteratively, by using the recurrence formulas

$$T_{k+1}(x) = 2xT_k(x) - T_{k-1}(x), \ T_k(x) = \frac{T'_{k+1}(x)}{2(k+1)} - \frac{T'_{k-1}(x)}{2(k-1)}, \ -1 \le x \le 1.$$

shifted to $x \in [0, b]$. We have

$$\begin{split} I_{\cdot,k} &= \int_0^x \frac{T_k \left(\alpha t - 1\right)}{(x-t)^{1-q}} dt = \int_0^x \frac{2\left(\alpha t - 1\right) T_{k-1} \left(\alpha t - 1\right) - T_{k-2} \left(\alpha t - 1\right)}{(x-t)^{1-q}} dt \\ &= 2\alpha \int_0^x \frac{t T_{k-1} \left(\alpha t - 1\right)}{(x-t)^{1-q}} dt - 2I_{\cdot,k-1} - I_{\cdot,k-2} = \\ &= -2\alpha \int_0^x \frac{T_{k-1} \left(\alpha t - 1\right)}{(x-t)^{-q}} dt + 2\alpha x I_{\cdot,k-1} - 2I_{\cdot,k-1} - I_{\cdot,k-2}. \end{split}$$

But

$$\int_{0}^{x} \frac{T_{k-1}(\alpha t-1)}{(x-t)^{-q}} dt = \frac{1}{\alpha} \int_{0}^{x} \left(\frac{\frac{d}{dt}T_{k}(\alpha t-1)}{2k} - \frac{\frac{d}{dt}T_{k-2}(\alpha t-1)}{2(k-2)} \right) \frac{dt}{(x-t)^{-q}} = \\ = \frac{1}{2\alpha} \left\{ \left[\frac{T_{k}(\alpha t-1)}{k} - \frac{T_{k-2}(\alpha t-1)}{k-2} \right] \frac{1}{(x-t)^{-q}} \Big|_{t=0}^{t=x} \right\} + \\ + \frac{1}{2\alpha} \int_{0}^{x} \left[\frac{T_{k}(\alpha t-1)}{k} - \frac{T_{k-2}(\alpha t-1)}{k-2} \right] \frac{q}{(x-t)^{1-q}} dt = \\ = \frac{1}{2\alpha} \left\{ - \left[\frac{(-1)^{k}}{k} - \frac{(-1)^{k-2}}{k-2} \right] x^{q} + \frac{q}{k} I_{\cdot,k} - \frac{q}{k-2} I_{\cdot,k-2} \right\}$$

and, consequently,

(3.3)
$$I_{\cdot,k} = -\left\{\frac{2(-1)^{k}x^{q}}{k(k-2)} + \frac{q}{k}I_{\cdot,k} - \frac{q}{k-2}I_{\cdot,k-2}\right\} + 2\alpha xI_{\cdot,k-1} - 2I_{\cdot,k-1} - I_{\cdot,k-2},$$
$$I_{\cdot,k}\left(1 + \frac{q}{k}\right) = 2(\alpha x - 1)I_{\cdot,k-1} + \left(\frac{q}{k-2} - 1\right)I_{\cdot,k-2} - \frac{2(-1)^{k}x^{q}}{k(k-2)}.$$

Finally, the above calculated matrix I must be replaced by $\frac{1}{\Gamma(q)} \cdot T^{-1} \cdot I$ and, if the function y has the spectral representation given by its Chebyshev coefficients $\mathbf{c} = (c_j)_{j=0,...,n-1}$, then the vector $I \cdot \mathbf{c}$ will contain the spectral coefficients of $J^q y(x)$, for $q \in (0,1)$.

The function fracbas can be tested on the well known example

$$J^{1/2}e^{x} \equiv \frac{1}{\sqrt{\pi}} \int_{0}^{x} \frac{e^{t}}{\sqrt{x-t}} dt = e^{x} \operatorname{erf}\left(\sqrt{x}\right),$$

for x > 0, say for $x \in [0,2]$. The code test1.m

n=16;dom=[0,2];kind=2;x=pd(n,dom,kind);q=1/2;

I=fracbas (x, dom, q); c=x2t (exp(x), kind); reznum=t2x(I*c, kind); gives the numerical approximation reznum which reproduces the exact result at the points x with an error of 2.65×10^{-15} for n = 16.

Another example is the Abel integral equation of the second kind (compare with the results from [2], Example 7)

$$y(x) + \int_0^x \frac{y(t)dt}{\sqrt{x-t}} = 2\sqrt{x},$$

with exact solution $y_{ex}(x) = 1 - e^{\pi x} erfc(\sqrt{\pi x})$. This equation of the form $\left(\frac{1}{\sqrt{\pi}} + J^{0.5}\right) u(x) = 2\sqrt{\frac{x}{\pi}}$ is discretized in spectral form as

$$\left(\frac{E}{\sqrt{\pi}}+I\right)\mathbf{c}=\mathbf{b}$$

where *E* is the unit matrix, *I* is the discretization of $J^{\frac{1}{2}}$ given by fracbas and **b** is the column vector of the coefficients of the function $2\sqrt{\frac{x}{\pi}}$ at the Chebyshev points **x**. Finally, the coefficients **c** are converted to the numerical values **v** of the solution *y* at the points **x**. The code test2.m

gives for n = 64 the numerical values of the solution with an error of 1.7543×10^{-4} and needs a computing time of 6×10^{-3} seconds. For n = 512 the error becomes 2.7658×10^{-6} for an elapsed time of 3.3 seconds.

Of course, this method is not suitable for functions with singularities in [0,b], or singularities of lower-order derivatives, like \sqrt{x} or $erfc(\sqrt{x})$; in these cases, *n* must be excessively large. In practical applications, however, it is required to approximate the fractional integrals or derivatives of such badly-behaved functions. A nice procedure is given in [18] for the fractional derivatives of functions *f* such that $f(x) = x^{\alpha}g(x)$, $\alpha > -1$, where *g* is assumed to be a wellbehaved function. The idea is to use the above Chebyshev spectral approximation only for the function *g*. Recurrence relations for the Chebyshev coefficients of the fractional derivative are also given, but the algorithm is not put in an operational form. A future extension of *Chebpack* will take into account the approximation of such functions, too.

However, we may improve the above result by a small change in the Abel equation

$$y(x) - 2\sqrt{x} + \int_0^x \frac{y(t) - 2\sqrt{t}}{\sqrt{x-t}} dt = -\int_0^x \frac{2\sqrt{t}}{\sqrt{x-t}} dt.$$

The right-hand side is known to be the well-behaved function $-\pi x$ and for the new unknown function $v(x) = y(x) - 2\sqrt{x}$ the equation becomes

$$v(x) + \int_0^x \frac{v(t)dt}{\sqrt{x-t}} = -\pi x.$$

With the corresponding change in the above code (test2_improved.m) we obtain the physical values of v(x) and then $y(x) = v(x) + 2\sqrt{x}$. Now, for n = 64 the error becomes 6.4887×10^{-8} for almost the same computing time.

If q = 1 the code [~, J0]=prim(n, dom) gives the integration matrix of order q = 1 that vanishes at x = 0. For a general $q \ge 0$ the Riemann-Liouville integration matrix Jrl of order q is:

- the unit matrix for q = 0
- the matrix $I^{(q)}$ given by fracbas for 0 < q < 1
- the matrix J_0 given by prim for q = 1

- the matrix $I^{(q-[q])} \cdot J_0^{[q]}$ where [q] is the largest integer not larger than q (floor(q) in Matlab), J_0 is given by prim and I is given by fracbas for q - [q]. The above formulas are implemented in an extended code Jrl=primrl(x, dom, q).

Again, if *y* has the spectral representation given by its Chebyshev coefficients $\mathbf{c} = (c_j)_{j=0,...,n-1}$, the vector $Jrl \cdot \mathbf{c}$ will contain the spectral coefficients of $J^q y(x)$, for $q \ge 0$.

The *Riemann-Liouville derivative of order q* is

(3.4)
$${}^{RL}D^q y(x) = \frac{1}{\Gamma(m-q)} \left(\frac{d}{dx}\right)^m \int_0^x \frac{y(t)dt}{(x-t)^{m-q-1}}$$

where $q \ge 0$, $m = \lceil q \rceil$ is the smallest integer not smaller than q (ceil(q) in Matlab). In an operatorial (discrete) form, ${}^{RL}D^q y = D^m \cdot J^{m-q} y$ where D = deriv(n, dom) is the usual first order differentiation matrix and J^{m-q} is calculated as above. The fractional Riemann-Liouville matrix ${}^{RL}D^q$ is given by the code Drl=derivrl(x, dom, q).

The Caputo derivative of order q is

(3.5)
$${}^{C}D^{q}y(x) = \frac{1}{\Gamma(m-q)} \int_{0}^{x} \frac{y^{(m)}(t)dt}{(x-t)^{1+q-m}}$$

where q and m are as above. In the operatorial form, $D^q y = J^{m-q} \cdot D^m y$. The matrix $^C D^q$ is discretized by the code Dc=derivc (x, dom, q).

We remark that for 0 < q < 1, ${}^{RL}D^q y(x) = {}^{C}D^q y(x) + \frac{y(0)x^{-q}}{\Gamma(1-q)}$, see [9] for more details. All the above codes, fracbas, primrl, derivrl, derivc are now enclosed in an updated form of *Chebpack* in the folder level0. The same Chebyshev spectral method, *but not in operational form*, is used in Theorem 3.2 from [29] where the uniform approximation of the above fractional derivatives is proved.

Let us test the differentiation matrix for the function $y(x) = \sin x$, $x \in [0, 2\pi]$. The exact fractional derivative of order q = 0.5 is [9]

$${}^{RL}D^q \sin x = {}^C D^q \sin x = \frac{x^{-0.5}}{2i\Gamma(0.5)} \left[{}_1F_1(1, 0.5, ix) - {}_1F_1(1, 0.5, -ix) \right].$$

We approximate this derivative by using the above *Chebpack* matrix $^{C}D^{0.5}$, the *Chebfun* package [30] and the Podlubny's package [25]. The Matlab code is test3.m and the results are given in Table 1. For this example *Chebpack* has the most efficient running time and accuracy.

Method	Error	Time (sec)	Dimension
Exact	0	4.45	-
Chebpack	1.09e-14	0.005	64
Chebfun	3.58e-13	0.59	20
Podlubny	0.05	0.02	629

TABLE 1. Test of the fractional derivative $D^{0.5} \sin x$

Another test for a fractional differential problem is [24]

$$^{C}D^{q}y(t) + y(t) = 1, t \in [0, 5]$$

 $y(0) = 1, y'(0) = -1, q = 1.8$

with the exact solution

$$y(t) = E_{q,1}(-t^q) - tE_{q,2}(-t^q) + t^q E_{q,q+1}(-t^q).$$

The equation is discretized as a linear system $(D_c + E)\mathbf{c} = \mathbf{b}$, where \mathbf{c} are the Chebyshev coefficients of the solution y and \mathbf{b} are the Chebyshev coefficients of the constant function 1. The initial conditions are implemented in the last two rows of the above system by using the

matrix T given by T = cpv(n, 0, [0, 1]). Precisely, $T \cdot c = y(0) = 1$ and $T \cdot D \cdot c = y'(0) = -1$.

The Matlab code is test4.m

```
[x, solnum]=test4(n, dom, q)
x=pd(n, dom, 2); T=cpv(n, dom(1), dom);
Dc=derivc(x, dom, q); D=deriv(n, dom);
A=Dc+eye(n); b=x2t(ones(n, 1), 2); % the discrete form Ay=b
A(n-1,:)=T; b(n-1)=1; % y(0)=1
A(n,:)=T*D; b(n)=-1; % y'(0)=-1
sol=A\b; solnum=t2x(sol, 2);
```

The error of the numerical solution solnum for n = 64 is 1.63×10^{-7} and the computing time is 0.012 seconds. By using the Podlubny's package [24], [25] with h = 0.01 the error is 0.0114 and the computing time is 0.47 seconds while fde12 [15] with h = 0.01 gives an error of 9.24×10^{-6} with a computing time of 0.07 seconds.

4. Numerical examples

This section contains 5 examples for specific problems on fractional differential equations.

Example 1. Let us consider the initial-conditions problem for the multi-term Bagley-Torvik fractional differential equation [23]

$$y'' + {}^{C}D^{\frac{3}{2}}y + y = \begin{cases} 8, & x \le 1\\ 0, & x > 1 \end{cases}, x \in [0, 30]$$
$$y(0) = y'(0) = 0.$$

The "exact" solution of this problem is useless for comparison with the numerical solution. It is calculated in [23] and it uses the derivatives of the Mittag-Lefler function. The numerical solution on the grid x given by *Chebpack*, example1.m is compared with the numerical solution given by Podlubny's code and by fde12 code for some stepsizes h. The results of this comparison (the maximal absolute differences between *Chebpack* solution and other solutions) are given in Table 2

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Method	Difference	Time (sec)	Dimension
Chebpack	0	0.32	256
Podlubny h=0.075	0.48	0.27	400
Podlubny h=0.04	0.17	1.49	750
Podlubny h=0.02	0.07	11.16	1500
Podlubny h=0.01	0.04	86.38	3000
fde12 h=0.01	0.1202	0.49	3000
fde12 h=0.001	0.0957	5.27	30000

TABLE 2. Bagley-Torvik equation, discontinuous rhs

We remark that the code fde12 can be used for the Bagley-Torvik equation as a fractional differential system of order $\frac{1}{2}$ [10]

$$y_1^{\left(\frac{1}{2}\right)} = y_2, y_2^{\left(\frac{1}{2}\right)} = y_3, y_3^{\left(\frac{1}{2}\right)} = y_4, y_4^{\left(\frac{1}{2}\right)} = F - y_1 - y_4,$$

$$y_1(0) = 0, y_2(0) = 0, y_3(0) = 0, y_4(0) = 0,$$

where F is the function given in the right-hand side of the equation.

For a more relevant comparison let us consider the Bagley-Torvik equation

$$y'' + y^{\left(\frac{3}{2}\right)} + y = 2 + 4\sqrt{\frac{x}{\pi}} + x^2, x \in [0, 10]$$
$$y(0) = y'(0) = 0.$$

with the exact solution $y_{ex}(x) = x^2$. The results are given by example1bis.m in Table 3.

Method	Dimension	Time (sec)	Error
Chebpack	n=32	0.003	2.84e-14
Podlubny	h=0.01	3.55	0.01
fde12	h=0.001	5.23	1.08e-4

TABLE 3. Bagley-Torvik equation, continuous rhs

Example 2. Let us consider the test boundary value problem from [34]

$$y''(x) + 0.5^C D^q y(x) + y(x) = 3 + x^2 \left(\frac{x^{-q}}{\Gamma(3-q)} + 1\right), x \in [0,1], q = 0.5$$
$$y(0) = 1, y(1) = 2.$$

The exact solution is $y(x) = x^2 + 1$. The Matlab code in *Chebpack* is example2.m with n = 16, dom = [0, 1], q = 0.5. The maximal error is about 2×10^{-16} . For comparison, the error given in [34] for the same dimension of the problem (i.e. $h = \frac{1}{16}$) and by using a cubic spline approximation is 4.35×10^{-3} .

A more relevant example is the eigenvalue problem [1]

$$^{C}D^{1.5}y(x) + \left(\frac{1}{x} + \lambda\right)y(x) = 0, x \in (0,1),$$

 $y(0) = 0, y'(1) = 0.$

This problem is discretized as $A\mathbf{c} = -\lambda B\mathbf{c}$ where $A = D_c + X^{-1}$, *B* is the unit matrix of order *n* and **c** is the column vector of the coefficients of the eigenfunctions. Now the boundary conditions are implemented such that the two last rows of *B* are the two last rows of *A* divided by $\lambda^* = 1000$. Consequently, the boundary conditions are numerically imposed and two spurious eigenvalues $\lambda = -\lambda^*$ are introduced and can be easily eliminated while keeping the matrix *B* nonsingular. The approximating generalized eigenproblem $AY = -\lambda BY$ is then solved by using the function eig from Matlab.

The Matlab code is example2bis.m

A=Dc+inv(X);B=eye(n);

$$A(n-1,:) = T(1,:); B(n-1,:) = A(n-1,:) / 1000;$$

$$A(n, :) = T(n, :) * D; B(n, :) = A(n, :) / 1000;$$

[V,L]=eig(full(A),full(B));[LL,ind]=sort(diag(L));VV=V(:,ind); lam=-LL(1:3);

where n = 128, dom = [0, 1], q = 1.5. The numerical eigenvalues are compared with those from [1] in Table 4. The computing time was 0.13 seconds.

λ (Chebpack)	λ[1]
1.6610	1.6609
13.5509	13.5504
20.5147	20.5145

TABLE 4. Fractional eigenproblem

Example 3. Let us consider now a nonlocal boundary value problem adapted from [27]

$${}^{C}D^{\frac{3}{4}}y(x) + \frac{x^{2}e^{x}}{5}y(x) - e^{x}\int_{0}^{x}ty(t)dt = \frac{6x^{\frac{9}{4}}}{\Gamma\left(\frac{13}{4}\right)}, x \in [0,1],$$

$$y(0) + y(1) - \frac{e+1}{e-2}y'(0) + \frac{1}{6}y''(1) - 10\int_{0}^{1}ty(t)dt = 0.$$

with the exact solution $y(x) = x^3$. Here the nonlocal condition is implemented by using the matrix *T* and the integration matrix *J*₀. Precisely, if T=cpv(n, [0, 1], [0, 1]) and **c** are the Chebyshev coefficients of the solution *y*, then J0*X*c approximates the coefficients of the primitive $\int_0^x ty(t)dt$ and (T(2, :) -T(1, :))*J0*X*c implements the Newton-Leibniz formula to approximate $\int_0^1 ty(t)dt$. The Matlab code is example3.m and for n = 64 the computing time is 0.01 sec. while the error is about 2.45×10^{-12} .

Example 4. Let us consider a distributed order differential problem, precisely the distributed order oscillator [19]

$$y''(x) + D^{\varphi(\alpha)}y(x) + y(x) = \begin{cases} 8, & 0 \le x \le 1\\ 0, & x > 1 \end{cases}, x \in [0, 30],$$
$$y(0) = y'(0) = 0, \varphi(\alpha) = 6\alpha (1 - \alpha), \alpha \in [0, 1].$$

Here the distributed order derivative is defined as

$$D^{\varphi(\alpha)}y(x) = \int_0^1 \varphi(\alpha) \frac{d^\alpha}{dx^\alpha} y(x) d\alpha$$

where $\varphi(\alpha)$ is the weight function of distribution of order $\alpha \in [0, 1]$.

The distributed order derivative is discretized as

$$D^{\varphi(\alpha)} = \sum_{k=1}^{m} w_k \cdot \varphi(\alpha_k)^C D^{\alpha_k}$$

where the nodes α_k and the weights w_k are given by [alpha,w]=pd(m,[0,1],1) and it is coded in doderiv.m. This function is included in level0 of *Chebpack*.

The code for the above problem is example4.m, the dimension of the problem is n = 256 for dom = [0,30] and the discretization of the distributed order derivative is performed with m = 16. Chebpack needs 3.61 seconds computing time. Podlubny's code from [19] A.5 needs 1.28 seconds for h = 0.075 (dimension of the problem 400), 5.96 seconds for h = 0.0375 and 29.61 seconds for h = 0.01875. A magnified portion of the graphs of these numerical solutions for $x \in [5, 15]$ is shown in Figure 1.



FIGURE 1. The numerical solutions for Bagley-Torvik distributed order problem

Example 5. Finally, let us consider a nonlinear problem [21]

$$D^{1.2}y(x) = y^{4}(x) + \frac{\Gamma(2.8)}{\Gamma(1.6)}x^{0.6} - (x^{1.8} - x + 1)^{4}, x \in [0, 1]$$
$$y(0) = 1, y'(0) = -1.$$

This problem is of the form Ly = N(x, y). After discretization, the discrete problem becomes $\mathscr{L}\mathbf{c} = \mathscr{N}(\mathbf{c})$ where \mathbf{c} are the Chebyshev coefficients of y, \mathscr{L} is the Caputo differentiation

matrix of order 1.2 and \mathcal{N} is the nonlinear right-hand side function of the coefficients **c**. Here the coefficients **C** of the nonlinear part y^4 are calculated by using the scheme

$$\mathbf{c} \xrightarrow{\mathtt{t2x}} y(\mathbf{x}) \longrightarrow y^4(\mathbf{x}) \xrightarrow{\mathtt{x2t}} \mathbf{C}$$

Finally, after imposing the initial conditions in the last two rows of the matrix \mathscr{L} and of the vector $\mathscr{N}(\mathbf{c})$, we obtain a nonlinear system $\mathbf{c} = \mathscr{L}^{-1}\mathscr{N}(\mathbf{c})$. This system can be solved in this case by successive iterations started by the Chebyshev coefficients of the constant function yo(x) = 1, for all $x \in [0, 1]$. The code for this problem is example5.m and only 20 iterations are needed for a difference less than 1.e - 13 between two successive iterations. The exact solution is $y_{ex}(x) = x^{1.8} - x + 1$. The errors of the numerical solutions for different dimensions n are given in Table 5.

TABLE 5. Nonlinear problem

n	Error	Time (sec)	
32	1.1e-5	0.013	
64	1.1e-6	0.022	
128	1.2e-7	0.08	
256	1.3e-8	0.45	
512	1.4e-9	3.86	

We remark that the use of a grading exponent r in formula (3.1) from [21] makes the grid points more densely clustered near the left endpoint of the interval [0,1]. This distribution of the grid points is automatically performed in *Chebpack* by using the Chebyshev points (2.1) and consequently, the above results are consistent with the best results in [21].

5. Conclusion

The tau method implemented in our package *Chebpack* leads to very simple and efficient codes that can solve different kinds of problems for fractional differential equations in a unified approach. The Chebyshev grid points are automatically clustered near the left endpoint of the working interval but this is not enough for a good approximation if the solution has singularities at that point. This remaining problem to extend *Chebpack* to efficiently handle functions of algebraic singularity will be the aim of future work. All the necessary Matlab sources for reproducing the above tests and examples are now part of an updated version of *Chebpack* [32], in the folder *Examples*, subfolder *Fractional differential equations*.

Conflict of Interests

The author declares that there is no conflict of interests.

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