A PRACTICAL GUIDE TO PRABHAKAR FRACTIONAL CALCULUS

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ABSTRACT. The Mittag–Leffler function is universally acclaimed as the Queen function of fractional calculus. The aim of this work is to survey the key results and applications emerging from the three-parameter generalization of this function, known as the Prabhakar function. Specifically, after reviewing key historical events that led to the discovery and modern development of this peculiar function, we discuss how the latter allows one to introduce an enhanced scheme for fractional calculus. Then, we summarize the progress in the application of this new general framework to physics and renewal processes. We also provide a collection of results on the numerical evaluation of the Prabhakar function.

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1. INTRODUCTION AND OUTLINE

Fractional calculus and the theory of special functions have managed to attract an increasing attention from the mathematical and physical community, especially in the last fifty years. In particular, the strict connection between these two research topics has been acting as the driving force for the latest developments in the literature on these subjects. The aim of this work is to build on these premises, collecting and summarizing the main works and results on the emerging Prabhakar’s approach to fractional calculus. In this spirit, in the following we provide the reader, who already has a general understanding of the basics of fractional calculus, with a comprehensive practical guide to the main aspects of Prabhakar calculus. In other words, the target audience for this review consists of researchers with an elementary background in fractional calculus who need a quick and hands-on training in order to start working on this compelling topic.

The review is therefore organized as follows: First, in Section 2 we summarize the key historical events that have led to the discovery and modern developments on the Mittag–Leffler function, its generalizations, and the extension of fractional calculus based on these exotic objects. In Section 3 we provide some physical motivations for the need of this new modeling scheme. In Section 4 we then recap the main mathematical features and properties of the Prabhakar function. In Section 5 we use the discussion in Section 4 as a springboard for the rigorous description of the theory of fractional calculus based on the Prabhakar function. In Section 6 we profit of the preliminary mathematical sections in order to frame some anomalous physical phenomena, mentioned in Section 3, within this generalized theory of calculus. In Section 7 we analyze the implications that these operators bring onto the theory of probability, with particular attention for renewal processes. In Section 8 we provide a bird’s-eye view on the numerical methods for evaluating the Prabhakar function. Finally, in Section 9 we provide some concluding remarks and an overview on some open problems.

2. History of the Mittag–Leffler function and its generalizations

In the beginning of the twentieth century the Swedish mathematician Magnus Gösta Mittag–Leffler, while working on summation methods for divergent series, defined a new special function which soon became known as the Mittag–Leffler (ML) function. Originally proposed in [107, 108], the ML function was then analyzed in a systematic fashion by A. Buhl in [8] around twenty years from its original formulation. Most notably, this function turns out to be a one-parameter version of a more general (two-parameter) function promoted by A. Wiman in [156], just a few years after the seminal works of M. G. Mittag–Leffler. Later on, in 1948, Pollard [125] investigated the complete monotonicity of the (one-parameter) ML function, while other efforts aimed at a deeper understanding of the main features of this function and its two-parameter generalization were made by P. Humbert, et al. in 1953 [73–75]. Then, the first extensive collection of results on ML functions, with one and two parameters, appeared in 1955 in the third volume of the Bateman Manuscript Project [30]. This project, aimed at redacting an encyclopedia on special functions, was started by Harry Bateman. However, the English mathematician passed away before the completion of this endeavor, which was ultimately brought to its end by A. Erdélyi, W. Magnus, F. Oberhettinger, and F. G. Tricomi. At that time the role of the ML function and its generalizations was yet to be recognized in the literature. For this reason these functions only appear in the Bateman series as a miscellaneous topic. The current notoriety of the ML function and its generalization can however be traced back
to the 1930s, specifically to the work of E. Hille and J. D. Tamarkin [72], when the connection between these objects and the solution of weakly singular linear Volterra integral equations with power law kernels was drawn for the first time. Along these lines, in 1954 J. H. Barrett employed the ML function to express the solution of some linear differential equations of non-integer order [2], thus identifying the fundamental role of this function in fractional calculus.

The increasing interest in the theory of differential equations of non-integer order thus motivated the mathematical community to investigate at a deeper level the properties and features of the ML function and its generalization, as it is clearly highlighted by several era-defining monographs (see, e.g. [28, 81, 96, 106, 120]) and fundamental works (see, e.g. [62, 82, 89, 98]). Nonetheless, the first and only consistent and comprehensive monograph [47] dedicated solely to this important topic was published in 2014 as the result of a joint endeavor of R. Gorenflo, A. A. Kilbas, F. Mainardi, and S. Rogosin.

Aside from the (mostly) mathematical interest in the ML function, in the early days of fractional calculus, it is worth remarking that this subject had also a significant impact in the physics literature. This is particularly highlighted by the early works of K. S. Cole on the electrical conductance of biological systems [17] (see also [25]) and by the fundamental contributions of B. Gross to the study of dielectric and mechanical relaxation, see e.g. [54, 55, 59]. Afterwards, M. Caputo and F. Mainardi gave their crucial contributions to fractional viscoelasticity [13, 14] by showing that the ML function always appears in the material functions of a system described by a fractional-order constitutive equation. It is also worth noting that [14] contains the very first plot of the ML function appeared in the literature. These are just a few early studies on the implications of the ML function in applied sciences that served as the backbone for all the subsequent developments in fractional calculus. Indeed, in light of the striking relevance of this function for the mathematical modeling of several phenomena, in 2007 the ML function was named the Queen function of fractional calculus [99] by F. Mainardi and R. Gorenflo. Since we do not have the pretense of providing a full and complete picture of the history and implications of this topic, we refer the interested reader to [47] by R. Gorenflo, et al.

The clear impact of the ML function and its two-parameter version on both the physical and mathematical community has then stimulated the quest for the generalization of these objects beyond two parameters. Among the many attempts we wish to commend the early contributions of V. Kiryakova and Y. Luchko on the so-called multi-index Mittag–Leffler functions [83], [95], [84], [85]. If not for a few exceptions, among which we find the works of V. Kiryakova mentioned above, most of the proposed generalizations were just curiosity-driven and in general do not carry any relevance for physics or real-world applications. Contrarily, another special praise is due to the three-parameter extension of the ML function introduced by the Indian mathematician Tilak Raj Prabhakar in [126], dating back to 1971. Although this function was originally analyzed within the framework of weakly singular Volterra integral equations from a pure mathematical perspective, it turns out that it plays an important role in the description of anomalous dielectric relaxation. An example of the effectiveness of the Prabhakar function in this context is provided by the Havriljak–Negami empirical law [64–66]. Notably, despite the short time-span between the original proposal of T. R. Prabhakar and the empirical studies of S. Havriliak and S. Negami, the connection between these works was drawn not longer than a decade ago, with the works of A. Hanyga and M. Seredyńśka [61] together with the study of E. Capelas De Oliveira, et al. [12]. Nonetheless, it is worth mentioning that several authors had previously investigated the time-domain representation of the
Havriliak–Negami model without highlighting (explicitly) its relation with the Prabhakar function (see, e.g. [68,69,111–113,147,150,155]).

The first contribution aimed at providing a fractional interpretation for weakly singular Volterra operators involving the Prabhakar function was first developed in 2002 by A. A. Kilbas, M. Saigo, and R. K. Saxena in [79]. Specifically, in this paper the notion of the Prabhakar fractional integral was first identified as an independent entity worth of investigation. Besides, in [80] one of the left-inverse operators of the Prabhakar fractional integral, which will later become known as Prabhakar fractional derivative (of the Riemann-Liouville type), was found. Over a decade later, in 2013, M. D’Ovidio and F. Polito [29] provided a regularization of the Prabhakar fractional derivative and named it after the Indian mathematician, thus originating the current terminology on the subject. These preliminary analyses then inspired the seminal work by R. Garra, et al. [33] where the features of the Prabhakar derivatives and some of their applications are discussed in depth. This work, in particular, contributed to kick-start a series of studies concerning applications of these operators to different fields of research (see, e.g. [10,11,16,138,142,148]). A particular praise is due to the work of L. Beghin and E. Orsingher who first derived the connection between the Prabhakar and Wright functions [5]. Furthermore, it is of paramount importance to mention the contribution of R. B. Paris [117] in the study of the asymptotic behavior of the Fox–Wright functions, of which the Prabhakar function represents a special case, which has then led to specific results on the Prabhakar function [32,118].

Finally, we conclude this bird’s-eye view on the history of this subject by mentioning some recent results carrying some more practical implications. First, the conditions for the complete monotonicity of the Prabhakar function have been extensively analyzed throughout the last decade in [12,97,149] by several authors. Using these studies A. Giusti [42] has then laid down a connection between the so called Prabhakar fractional calculus and Kochubei’s general fractional calculus [87]. Lastly, it is worth recalling that a discussion of the numerical aspects of the Prabhakar function has been carried out by R. Garrappa in [35], who also designed a Matlab code for computing the Prabhakar function [34]. Note that, as of today, this is the sole code (freely available on the Mathwork website) for the computation of the Prabhakar function.

3. ANOMALOUS PHYSICS: A CRY FOR HELP

In this section we provide some physical examples showing the need for an extension of ordinary calculus based on the Prabhakar function. In this regard, the most telling framework is provided by the so called anomalous phenomena emerging in several physical settings.

Let us start off with the classical theory of dielectrics [76]. A dielectric material is a poor conductor of electricity whose main feature is that its internal electric charges are not allowed to flow freely under the effect of an external electric field. On the contrary, they will mostly tend to shift a little from their equilibrium position. This will result in a very low electric current, since some charges will still be able to move through the material, and a dielectric polarization of the system. Experimentally it is observed that once the external stress is turned off a dielectric material will tend to lose its acquired polarization over time. This time lag, known in physics as dielectric relaxation, is due to the fact that the atomic and/or molecular polarization in response to a changing electric field is not an instantaneous process (see e.g. [78]). A key quantity in the study of electromagnetism in matter is given by the polarization density $\mathbf{P}$, or simply polarization, which is defined
as the average dipole moment of the material per unit volume and it accounts for the contribution of the “bound charges” that are responsible for the dielectric polarization. This picture allows us to split the electric charge density appearing in Gauss’ law into two contributions, namely the one given by the charges which are free to move through the material and a second piece due to bound charges. This suggests the introduction of a new entity

$$D = \varepsilon_0 \mathbf{E} + \mathbf{P},$$

known as the electric displacement, where $\mathbf{E}$ is the electric field and $\varepsilon_0$ denotes the vacuum permittivity, whose divergence depends solely on the free charge.

For a linear dielectric one has that the polarization density is proportional to the electric field, namely

(3.1)$$P = \varepsilon_0 \chi \mathbf{E},$$

which leads to

(3.2)$$D = \varepsilon \mathbf{E} = \varepsilon_0 \varepsilon_r \mathbf{E} = \varepsilon_0 \left(1 + \chi\right) \mathbf{E},$$

with $\chi$ the electric susceptibility of the material, $\varepsilon_r = 1 + \chi$ the relative permittivity, and $\varepsilon = \varepsilon_0 \varepsilon_r$ the permittivity of the system. On the other hand, it is well known that not all dielectrics display this very simple behavior, hence the polarization density can be, in general, a complicated function of the electric field. Nonetheless, if one wishes to keep the relation between $P$ and $\mathbf{E}$ linear while introducing a non-instantaneous response to a time-varying electric field, then a straightforward generalization of (3.1) is obtained via a Fourier convolution. In detail, given two sufficiently regular functions $f(t)$ and $g(t)$ the convolution of these two functions, in the Fourier sense, is defined as

(3.3)$$(f * g)(t) := \int_{-\infty}^{\infty} f(t-\tau) g(\tau) \, d\tau.$$

Hence, if we now denote by $\chi(t)$ a time dependent electric susceptibility such that $\chi(t) = 0$ for $t < 0$, then (3.1) can be generalized as

(3.4)$$P(t) = \varepsilon_0 \left(\chi * \mathbf{E}\right)(t) \equiv \varepsilon_0 \int_{-\infty}^{t} \chi(t-\tau) \mathbf{E}(\tau) \, d\tau,$$

or, in the frequency domain

(3.5)$$\hat{P}(\omega) = \varepsilon_0 \hat{\chi}(\omega) \hat{\mathbf{E}}(\omega),$$

where the hat denotes the Fourier transform, i.e.

$$\hat{f}(\omega) \equiv \mathcal{F}[f(t);\omega] := \int_{-\infty}^{\infty} \exp(-i\omega t) \, f(t) \, dt.$$

Eq. (3.4) is particularly instructive since it highlights the role of $\chi(t)$ as a measure of the response of the material to a sudden variation in the electric field $\mathbf{E}(t)$. In other words, the time dependent electric susceptibility acts as a “response function” of the dielectric which, in this scenario, is pictured as a passive linear system (see e.g. [37, 151], and references therein). More precisely, the response function of the dielectric can be defined, in the frequency domain, as

(3.6)$$\hat{\phi}(\omega) := \frac{\hat{\chi}(\omega)}{\hat{\chi}(0)}.$$

Now, since we are considering a causal linear system and we are only interested in the history of the system from a certain time $t = 0$ (at which we have applied the external perturbation) up to the time $t > 0$, Eq. (3.6) can be recast as [37]
(3.7) \[ \tilde{\phi}(s) = \hat{\phi}(\omega)\big|_{s=i\omega}, \]
where the tilde denotes the Laplace transform, i.e.
\[ \tilde{f}(s) \equiv \mathcal{L}\{f(t) ; s\} := \int_0^\infty \exp(-s t) f(t) \, dt. \]

Alongside with the response function one can also introduce another quantity, known as relaxation function, which is defined as
\[ \Psi(t) := 1 - \int_0^t \phi(\tau) \, d\tau, \quad t \geq 0, \]
from which one can also infer that \( \phi(t) = -\Psi'(t) \), where the prime denotes the derivative with respect to time.

The standard lore on dielectrics is described by the Debye model \([26, 71]\) according to which an ideal dielectric is schematically represented as a series of non-interacting dipoles. Such a model then leads to a (normalized) complex susceptibility given by
\[ \hat{\chi}_D(\omega) = \frac{1}{1 + i\omega \tau_D}, \]
which yields \([37,71]\)
\[ \phi_D(t) = \exp(-t/\tau_D), \quad \Psi_D(t) = \exp(-t/\tau_D), \]
with \( \tau_D \) denoting the relaxation time.

While this very simple and powerful model provides a satisfactory description of the features of a wide class of standard dielectrics, there are several experimental evidences \([77,78,110]\) that show large deviations from the exponential relaxation for many materials. An example of these anomalous dielectrics is provided by the Cole–Cole relaxation \([18,19]\) displayed by certain biological tissues \([91, 129]\). The key feature of these systems is a fractional-power law fall-off shown by the complex susceptibility, at high frequencies, which leads to a stretched relaxation over a wider range of frequencies. In detail, the Cole–Cole relaxation is modeled in terms of a complex susceptibility that reads
\[ \hat{\chi}_{CC}(\omega) = \frac{1}{1 + (i\omega \tau_*)^\alpha}, \quad 0 < \alpha \leq 1 \]
that yields
\[ \phi_{CC}(t) = \frac{\exp(-t/\tau_*)}{\tau_*} \left[ -\left( \frac{t}{\tau_*} \right)^\alpha \right], \]
and
\[ \Psi_{CC}(t) = E_{\alpha,1} \left[ -\left( \frac{t}{\tau_*} \right)^\alpha \right], \]
where \( \tau_* \) represents a typical relaxation timescale for the system.

Note that, although the Cole–Cole model breaks the standard exponential relaxation replacing it with fractional-power law tails, it can still be framed within the conventional picture of fractional calculus based on Caputo derivatives. This is, however, not the case for several other empirical laws for anomalous dielectrics such as the Davidson–Cole \([24]\) and Havriliak–Negami \([63]\) models, described by \([37]\)
\[ \hat{\chi}_{DC}(\omega) = \frac{1}{(1 + i\omega \tau_*)^\gamma}, \quad 0 < \gamma \leq 1 \]
and \([37,67,97]\)
\[
\hat{\chi}_{HN}(\omega) = \frac{1}{\left[1 + (i\omega\tau_*)^\alpha\right]^\gamma}, \quad 0 < \alpha \leq 1, \quad 0 < \alpha\gamma \leq 1
\]
respectively, for which the dielectric decay is inherently related to the Prabhakar function [12, 36, 37, 49, 50, 97], as we shall discuss in detail in Section 6.

Although anomalous dielectric relaxation is probably the most apparent exemplification of the need for a fractional theory of calculus based on the Prabhakar function, there exist also several other different physical systems in which this formalism naturally emerges. For instance, this scenario comes up when dealing with certain kind of anomalous diffusion processes [131–136, 139, 141], and when employing the formal duality between dielectric materials and viscoelastic systems [23, 41, 43]. More on these physical cases can be found in Section 6.

4. Mathematical preliminaries on the Prabhakar function

The Prabhakar function (also known as the three-parameter Mittag-Leffler function, see e.g. [47]), introduced by T. R. Prabhakar in 1971 [126], is defined as

\[
E_{\alpha,\beta,\gamma}(z) = \sum_{k=0}^{\infty} \frac{(\gamma)_k z^k}{k! \Gamma(\alpha k + \beta)}, \quad \alpha, \beta, \gamma \in \mathbb{C}, \quad \text{Re}(\alpha) > 0, \quad z \in \mathbb{C},
\]

where \((\gamma)_k \equiv \Gamma(\gamma + k)/\Gamma(\gamma)\) is the Pochhammer symbol and \(\Gamma(\cdot)\) denotes the Euler gamma function. Note that \(E_{\gamma,1,1}(z)\) is an entire function of order \(\rho = 1/\text{Re}(\alpha)\) and type \(\sigma = 1\) [47].

Although the three parameters \(\alpha, \beta\) and \(\gamma\) are allowed to assume values in \(\mathbb{C}\), and several of the properties we are going to present hold with complex parameters, in this work we will restrict our attention just to real parameters. The reason for this choice is that real parameters turn out to be of particular interest in physics, as we shall discuss later on. Therefore, for the remainder of this work we will implicitly assume

\[\alpha, \beta, \gamma \in \mathbb{R}, \quad \alpha > 0.\]

In view of the increasing relevance of the Prabhakar function in current literature, in the following we summarize the most relevant results and open problems on this compelling research topic.

4.1. Main properties and relations with other functions. The Prabhakar function has important connections with the standard ML function, its two-parameter generalization, and other special functions. Since for \(k \geq 1\) one has

\[
(\gamma)_k = \frac{\Gamma(\gamma + k)}{\Gamma(\gamma)} = \gamma(\gamma + 1) \cdots (\gamma + k - 1),
\]

then for \(\gamma = 0\) we clearly have

\[E_{\alpha,\beta,0}(z) = \frac{1}{\Gamma(\beta)},\]

whereas, for \(\gamma = 1\) we recover the widely known two-parameter ML function, i.e.

\[E_{\alpha,\beta,1}(z) = E_{\alpha,\beta}(z) = \sum_{k=0}^{\infty} \frac{z^k}{\Gamma(\alpha k + \beta)}.\]

Hence, if one sets \(\alpha = \beta = \gamma = 1\) the classical exponential function \(E_{1,1,1}(z) = e^z\) is retrieved.

Being the Prabhakar function distinguished from the two-parameter ML function just for the presence of a third parameter \(\gamma\), one is compelled to focus on properties related to this additional parameter.
A first formula, already present in the original work by Prabhakar [126], enables one to operate the reduction of the third parameter, i.e.

\[ E_{\alpha,\beta}^{\gamma+1}(z) = \frac{E_{\alpha,\beta-1}^{\gamma}(z) + (1 - \beta + \alpha)E_{\alpha,\beta}^{\gamma}(z)}{\alpha \gamma}, \]

and a further reduction formula was later derived in [32] and reads

\[ E_{\alpha,\beta}^{\gamma+1}(z) = \frac{E_{\alpha,\beta-\alpha-1}^{\gamma}(z) + (1 - \beta + \alpha)E_{\alpha,\beta-\alpha}^{\gamma}(z)}{\alpha \gamma z}, \quad z \neq 0. \]

Both these formulas can be used to lower the value of the third (additional) parameter, nonetheless these results can probably be better appreciated when \( \gamma \) is an integer. Specifically, let \( \gamma = k \in \mathbb{N} \), then

\[ E_{\alpha,\beta}^{k+1}(z) = \frac{1}{\alpha^k k!} \sum_{j=0}^{k} d_j^{(k)} E_{\alpha,\beta-j}(z), \]

and

\[ E_{\alpha,\beta}^{k+1}(z) = \frac{1}{\alpha^k z^k k!} \sum_{j=0}^{k} d_j^{(k)} E_{\alpha,\beta-ak-j}(z), \quad z \neq 0, \]

which means that one can explicitly express \( E_{\alpha,\beta}^{k+1} \) as a combination of two-parameter ML functions. The coefficients \( d_j^{(k)} \) appearing in both (4.4) and (4.5) are given by the recursive expression

\[ d_j^{(k)} = \begin{cases} (1 - \beta + \alpha) d_0^{(k-1)} & j = 0, \\ d_{j-1}^{(k-1)} + (1 - \beta + \alpha + j) d_j^{(k-1)} & j = 1, \ldots, k - 1, \\ 1 & j = k. \end{cases} \]

originally derived by R. Garrappa and M. Popolizio in [40].

The Prabhakar function is also directly related to the Fox–Wright functions [31, 157]. Specifically, the latter are defined as

\[ p\Psi_q(z) \equiv p\Psi_q \left[ \frac{(a_1, \rho_1), \ldots, (a_p, \rho_p)}{(b_1, \sigma_1), \ldots, (b_q, \sigma_q)} ; z \right] = \sum_{k=0}^{\infty} \frac{\prod_{r=1}^{p} \Gamma(a_r + \rho_r k) z^k}{\prod_{s=1}^{q} \Gamma(b_s + \sigma_s k) k!}, \]

where \( p \) and \( q \) are integers, \( \rho_r, a_r, \sigma_r, b_s \) are real or complex parameters such that \( \rho_r k + a_r \neq 0, -1, -2, \ldots \). Then, one can easily verify that \( E_{\alpha,\beta}^{\gamma}(z) \) is proportional to \( \Gamma(\gamma) p\Psi_1(\gamma,1) \).

\[ E_{\alpha,\beta}^{\gamma}(z) = \frac{1}{\Gamma(\gamma)} \Gamma(\gamma) \frac{1}{\Gamma(\gamma)} p\Psi_1 \left[ \frac{(\gamma,1)}{(\beta,\alpha)} ; z \right]. \]

Moreover, if we recall the relation between the Fox–Wright functions and the Fox \( H \)-function [102], i.e.

\[ p\Psi_q \left[ \frac{(a_1, \rho_1), \ldots, (a_p, \rho_p)}{(b_1, \sigma_1), \ldots, (b_q, \sigma_q)} ; z \right] = H_{p,q+1}^{1, p} \left[ \frac{(1-a_1, \rho_1), (1-a_2, \rho_2), \ldots, (1-a_p, \rho_p)}{(0,1), (1-b_1, \sigma_1), \ldots, (1-b_q, \sigma_q)} ; z \right], \]

then (4.8) can be rewritten in terms of the Fox \( H \)-function as (see [143])

\[ E_{\alpha,\beta}^{\gamma}(z) = \frac{1}{\Gamma(\gamma)} H_{1,2}^{1,1} \left[ \frac{(1-\gamma,1)}{(0,1), (1-\beta,\alpha)} ; z \right]. \]
Finally, it is worth mentioning that series in Prabhakar functions have been studied by J. Paneva–Konovska in [114–116].

4.2. Derivatives and integrals. A term-by-term derivation of the series expansion in (4.1) gives

\[
\frac{d^m}{dz^m} E_{\alpha,\beta}^\gamma (z) = \gamma (\gamma + 1) \cdots (\gamma + m - 1) E_{\alpha,\beta}^{\gamma + m} (z) \\
= \frac{\Gamma (\gamma + m)}{\Gamma (\gamma)} E_{\alpha,\beta}^{\gamma + m} (z),
\]

with \( m \in \mathbb{N} \), that reduces to the known result for the repeated derivative of the two-parameter ML function for \( \gamma = 1 \), \( i.e. \)

\[
(4.11) \quad \frac{d^m}{dz^m} E_{\alpha,\beta} (z) = m! E_{\alpha,\beta}^{m+1} (z).
\]

This last expression turns out to be particularly useful when combined with (4.4) or (4.5), since it allows to express derivatives of the two-parameter ML function as combination of different instances of the same function, a result widely exploited in [40] for numerical purposes.

Similarly, Dzhrbashyan’s formula [28, 47] for the derivative of the two-parameter ML function can be generalized to the Prabhakar case as (see [32])

\[
(4.12) \quad \frac{d}{dz} E_{\alpha,\beta}^\gamma (z) = \frac{E_{\alpha,\beta-1}^\gamma (z) + (1 - \beta) E_{\alpha,\beta}^\gamma (z)}{\alpha z}, \quad z \neq 0.
\]

Other important results concerning integrals and derivatives involving the Prabhakar function appear to be strictly related to the product \( t^{\beta-1} E_{\alpha,\beta}^\gamma (t^\alpha z) \). In details, it is easy to see that the \( m \)-th derivative and the integral of this particular combination read

\[
\frac{d^m}{dt^m} t^{\beta-1} E_{\alpha,\beta}^\gamma (t^\alpha z) = t^{\beta-m-1} E_{\alpha,\beta-m}^\gamma (t^\alpha z), \quad m \in \mathbb{N}
\]

and

\[
\int_0^t \tau^{\beta-1} E_{\alpha,\beta}^\gamma (\tau^\alpha z) d\tau = t^\beta E_{\alpha,\beta+1}^\gamma (t^\alpha z),
\]

respectively. Further, the \( m \)-th repeated integration of \( t^{\beta-1} E_{\alpha,\beta}^\gamma (t^\alpha z) \) gives

\[
J_0^m \left[t^{\beta-1} E_{\alpha,\beta}^\gamma (t^\alpha z)\right] = \int_0^t \int_0^{\tau_1} \cdots \int_0^{\tau_{m-1}} \int_0^{\tau_m} \tau_m^{\beta-1} E_{\alpha,\beta}^\gamma (\tau_m^\alpha z) d\tau_m \cdots d\tau_2 d\tau_1 = \frac{1}{(m-1)!} \int_0^t (t - \tau)^{m-1} \tau^{\beta-1} E_{\alpha,\beta}^\gamma (\tau^\alpha z) d\tau = 1^{\beta+m-1} E_{\alpha,\beta+m}^\gamma (t^\alpha z).
\]

We conclude this section by collecting some formulas for fractional integrals and derivatives of the Prabhakar function. To this aim, we first need to recall that for a function \( f \in L^1 [t_0, T] \) the Riemann-Liouville (RL) fractional integral of order \( \alpha > 0 \) is defined as [27, 81, 96, 106, 130]

\[
(4.14) \quad J_0^\alpha f(t) = \frac{1}{\Gamma (\alpha)} \int_{t_0}^t (t - u)^{\alpha-1} f(u) du.
\]

One of its left-inverse operators, known as the Riemann-Liouville fractional derivative, then reads

\[
(4.15) \quad \text{RL} D_{t_0}^\alpha f(t) := D^m J_{t_0}^{m-\alpha} f(t) = \frac{1}{\Gamma (m-\alpha)} \frac{d^m}{dt^m} \int_{t_0}^t (t - u)^{m-\alpha-1} f(u) du
\]
with $m = \lceil \alpha \rceil$ (the smallest integer greater than $\alpha$). A regularized version of the RL derivative, which still acts as left-inverse of the RL integral, is given by the Caputo fractional derivative. The latter is obtained by exchanging the operations of integration and differentiation in (4.15) and applies to functions $f \in AC^m[t_0, T]$ (the set of functions with absolutely continuous derivatives of order $m - 1$), i.e.

$$
(4.16) \quad CD^\alpha_{t_0} f(t) := J^m_0 f(t) = \frac{1}{\Gamma(m - \alpha)} \int_{t_0}^{t} (t - u)^{m-\alpha-1} f(u) \, du.
$$

All that being said, it is straightforward to verify that (see e.g. [47, 80])

$$
J^\rho_o \left[ t^{\beta-1} E_{\alpha,\beta}^\gamma (t^{\alpha} \lambda) \right] = t^{\beta+\rho-1} E_{\alpha,\beta+\rho}^\gamma (t^{\alpha} \lambda),
$$

$$
RI D^\rho_0 \left[ t^{\beta-1} E_{\alpha,\beta}^\gamma (t^{\alpha} \lambda) \right] = t^{\beta-\rho-1} E_{\alpha,\beta-\rho}^\gamma (t^{\alpha} \lambda),
$$

$$
CD^\rho_0 \left[ t^{\beta-1} E_{\alpha,\beta}^\gamma (t^{\alpha} \lambda) \right] = t^{\beta-\rho-1} E_{\alpha,\beta-\rho}^\gamma (t^{\alpha} \lambda), \quad \beta > \lceil \rho \rceil,
$$

with $\rho > 0$.

### 4.3. Integral transforms.

Let us consider, again, the product $t^{\beta-1} E_{\alpha,\beta}^\gamma (t^{\alpha} \lambda)$. Then it is not hard to show that (see e.g. [79])

$$
(4.17) \quad \mathcal{L} \left[ t^{\beta-1} E_{\alpha,\beta}^\gamma (t^{\alpha} z) ; s \right] = \frac{s^{\alpha-\beta}}{(s^\alpha - z)^\gamma}, \quad \text{Re}(s) > 0 \text{ and } |s| > |z|^{1/\gamma}.
$$

This result is particularly useful since it plays an important role in both physical applications and numerical computations of the Prabhakar function. Alternatively, proceeding in a more general fashion one finds (see [138, 144])

$$
(4.18) \quad \mathcal{L} \left[ t^{\beta-1} E_{\alpha,\beta}^\gamma (t^{\alpha} z) ; s \right] = \frac{s^{-\rho}}{\Gamma(\gamma)} 2 \Psi_1 \left[ (\rho, \sigma), (\gamma, 1) ; \frac{z}{s^\sigma} \right].
$$

Setting $t = 1$ in (4.17), one finds an integral representation of the Prabhakar function, i.e.

$$
E_{\alpha,\beta}^\gamma (z) = \frac{1}{2\pi i} \int_{C} e^{s} \frac{z^{\alpha-\beta}}{(s^\alpha - z)^\gamma} \, ds,
$$

where $C$ is understood as the Bromwich contour, for which all singularities lay on the left-hand side of a certain $\text{Re}(s) = \xi$ known as the abscissa of convergence. Similarly, a particular spectral representation of $t^{\beta-1} E_{\alpha,\beta}^\gamma (t^{\alpha} \lambda)$ has been computed in [12, 97, 149], and it reads

$$
(4.19) \quad t^{\beta-1} E_{\alpha,\beta}^\gamma (-t^{\alpha}) = \int_{0}^{\infty} e^{-rt} K_{\alpha,\beta}^\gamma (r) \, dr
$$

with

$$
K_{\alpha,\beta}^\gamma (r) = \frac{r^{\alpha-\beta} \sin(\gamma \theta_{\alpha}(r) + (\beta - \alpha \gamma) \pi)}{\pi (r^{2\alpha} + 2r^{\alpha} \cos(\alpha \pi) + 1)^{\gamma/2}} ,
$$

$$
\theta_{\alpha}(r) = \arctan \left( \frac{r^{\alpha} \sin(\pi \alpha)}{r^{\alpha} \cos(\pi \alpha) + 1} \right) \in [0, \pi],
$$

where a very specific branch of arctan($z$), treated as a multi-valued function, has been chosen in order to have $\theta_{\alpha}(r) \in [0, \pi]$ (see [97] for details). Further, this representation plays a key role in the determination of the range of parameters that yields a completely monotonic behavior of the Prabhakar function. Besides, this last representation is also particularly useful when studying estimates and bounds on the fractional Prabhakar integral, see e.g. [124].
In the light of the strict relation between Fox–Wright functions and the Prabhakar function, as shown in Eq. (4.8), one can easily derive the \textit{Mellin–Barnes integral representation} of the three-parameter ML function, \textit{i.e.}

\begin{equation}
E_{\alpha,\beta}^{\gamma}(z) = \frac{1}{\Gamma(\gamma)} \frac{1}{2\pi i} \int_{C} \frac{\Gamma(s)\Gamma(\gamma - s)}{\Gamma(\beta - \alpha s)} (-z)^{-s} ds, \quad |\arg z| < \pi,
\end{equation}

where $C$ is defined as above. Along this line, denoting by

\begin{equation}
M[f(t); s] = \int_{0}^{\infty} f(t) t^{s-1} dt,
\end{equation}

the \textit{Mellin transform} of $f(t)$, then from (4.20) one can immediately infer that

\begin{equation}
M[E_{\alpha,\beta}^{\gamma}(-t); s] = \frac{\Gamma(s)\Gamma(\gamma - s)}{\Gamma(\gamma)\Gamma(\beta - \alpha s)}.
\end{equation}

4.4. \textbf{Complete monotonicity.} A crucial role for the applications of special functions in relaxation models and probability theory is played by the notion of \textit{complete monotonicity}. We recall that a function $f : (0, +\infty) \to \mathbb{R}$ is completely monotonic (CM) if $f$ has derivatives of all orders on $(0, +\infty)$ and $(-1)^k f^{(k)}(t) \geq 0$, for any $k \in \mathbb{N} \cup \{0\}$ and $t > 0$.

The complete monotonicity of the Prabhakar function has been studied in several works \cite{12, 97, 149} and the current most general result states that the product $t^{\beta - 1} E_{\alpha,\beta}^{\gamma}(-t^\alpha)$ is CM if the three parameters satisfy the conditions

\begin{equation}
0 < \alpha \leq 1, \quad 0 < \alpha\gamma \leq \beta \leq 1
\end{equation}

which are obtained by combining the representation in (4.19) and the \textit{Bernstein theorem}. The latter, in particular, states that a function is CM if its spectral function is non-negative. Thus, the restrictions in (4.23) are obtained by inspecting the conditions according to which $K_{\alpha,\beta}^{\gamma}(r)$ is $\geq 0$ for all $r \geq 0$. This result is particularly relevant since $t^{\beta - 1} E_{\alpha,\beta}^{\gamma}(-t^\alpha)$ comes up very often in the study of anomalous dielectrics, as we shall discuss in more details in Section 6.

Finally, in \cite{51} it was shown that $E_{\alpha,\beta}^{\gamma}(-t)$ is CM if $0 < \alpha < 1$ and $0 < \alpha\gamma \leq \beta$, namely without the restriction $\beta \leq 1$. For similar results on another Mittag-Leffler type function with 3 parameters, the so-called \textquotedblleft Le Roy type\textquotedblright function, see \cite{52}.

4.5. \textbf{Asymptotic behavior.} The asymptotic behavior of the Prabhakar function, in the whole complex plane, is a non-trivial topic due to the dependence of the coefficients in the asymptotic expansions on the three parameters $\alpha$, $\beta$ and $\gamma$ (see \textit{e.g.} \cite{32} and references therein). On the other hand, rewriting the Prabhakar function as a Fox–Wright function, as in (4.8), allows one to take profit of the well-established results on the asymptotic behavior of this more general class of functions, see \textit{e.g.} \cite{7, 117, 118, 157–159}.

It turns out that the behavior of $E_{\alpha,\beta}^{\gamma}(z)$ for large values of $|z|$ varies from exponential to algebraical depending on the sector of the complex plane where $z$ lies. Furthermore, it is found that the parameter $\alpha$ is the one that (most prominently) controls the asymptotic properties of the Prabhakar function. Specifically, the lines $\arg z = \pm \alpha \pi/2$ are \textit{anti-Stokes lines}, where the function changes its behavior from increasing to decreasing. Whereas, $\arg z = \pm \alpha \pi$ are the \textit{Stokes lines}, where the exponential term quickly decays leaving just a predominant algebraic term \cite{118}. This behavior is depicted in Figure 1 for $0 < \alpha < 1$ (the acronyms \textquotedblleft E.S.	extquotedblright, \textquotedblleft E.L.	extquotedblright and \textquotedblleft Alg.	extquotedblright stand for \textit{exponentially small}, \textit{exponentially large}, and \textit{algebraic}, respectively). When $1 < \alpha < 2$ the Stokes lines $\arg z = \pm \alpha \pi$
collapse onto the negative real axis and there are no more regions in which the Prabhakar function presents a significant algebraic expansion.

\[
\arg z = \frac{\alpha}{2} \pi
\]

\[
\arg z = -\frac{\alpha}{2} \pi
\]

\[
\arg z = \alpha \pi
\]

\[
\arg z = -\alpha \pi
\]

**Figure 1.** Asymptotic behavior of \( E_{\alpha,\beta}^\gamma(z) \) in different regions of the complex plane for \( 0 < \alpha < 1 \). E.S.: exponentially small, E.L.: exponentially large and Alg.: algebraic

The explicit form of the exponential and algebraic terms, together with an algorithm for the evaluation of their coefficients, were described in [117,118]. In particular the algebraic expansion is

\[
A_{\alpha,\beta}^\gamma(z) = \frac{z^{-\gamma}}{\Gamma(\gamma)} \sum_{k=0}^{\infty} \frac{(-1)^k \Gamma(k + \gamma)}{k! \Gamma(\beta - \alpha(k + \gamma))} \frac{1}{k! \Gamma(\beta - \alpha(k + \gamma))} z^{-k}
\]

while the exponential expansion reads

\[
E_{\alpha,\beta}^\gamma(z) = \frac{1}{\Gamma(\gamma)} e^{\pm 1/\alpha} z^{-\alpha \gamma} \sum_{k=0}^{\infty} c_k z^{-\frac{k}{\alpha}}
\]

where \( c_k \) are obtained from the inverse factorial expansion of

\[
F_{\alpha,\beta}^\gamma(s) := \frac{\Gamma(\gamma + s) \Gamma(\alpha s + \psi)}{\Gamma(s + 1) \Gamma(\alpha s + \beta)} = \alpha^{1-\gamma} \left( 1 + \sum_{j=1}^{\infty} \frac{c_j}{(\alpha s + \psi)_j} \right)
\]

for \( |s| \to \infty \) in \( |\arg(s)| \leq \pi - \epsilon \) and any arbitrarily small \( \epsilon > 0 \). As usual, \((x)_j\) denotes the Pochhammer symbol and \( \psi = 1 - \gamma + \beta \).

One can now summarize the asymptotic behavior of the Prabhakar function, for \( |z| \to \infty \), as follows (see [118])

\[
0 < \alpha \leq 1 : \\
E_{\alpha,\beta}^\gamma(z) \sim \begin{cases} 
E_{\alpha,\beta}^\gamma(z) + A_{\alpha,\beta}^\gamma(ze^{\mp \pi i}) & |\arg z| < \frac{1}{2} \alpha \pi \\
A_{\alpha,\beta}^\gamma(ze^{\mp \pi i}) & \frac{1}{2} \alpha \pi < |\arg z| < \alpha \pi \\
A_{\alpha,\beta}^\gamma(ze^{\mp \pi i}) & \alpha \pi < |\arg z| \leq \pi 
\end{cases}
\]

\[
1 < \alpha < 2 : \\
E_{\alpha,\beta}^\gamma(z) \sim \begin{cases} 
E_{\alpha,\beta}^\gamma(z) + A_{\alpha,\beta}^\gamma(ze^{\mp \pi i}) & |\arg z| < \frac{1}{2} \alpha \pi \\
A_{\alpha,\beta}^\gamma(ze^{\mp \pi i}) + E_{\alpha,\beta}^\gamma(z) + E_{\alpha,\beta}^\gamma(ze^{\mp 2\pi i}) & \frac{1}{2} \alpha \pi < |\arg z| < \pi 
\end{cases}
\]
where the sign in $e^{∓πi}$ and $e^{∓2πi}$ is negative when $z$ lies in the upper complex half-plane and positive otherwise. Note that some of the cases discussed above differ just by the order in which $E^{γ}_{α,β}(z)$ and $A^{γ}_{α,β}(ze^{∓πi})$ appear. The reason for this choice comes from the fact that, for the sake of clarity, we prefer to give priority in the expression to the dominant term. The term $E^{γ}_{α,β}(ze^{∓2πi})$ can be usually neglected, except when $\arg z$ is close to $π$ since in this case it becomes comparable to $E^{γ}_{α,β}(z)$.

To complete the analysis we consider the case $α ≥ 2$. In detail, one finds

$$α ≥ 2 : E^{γ}_{α,β}(z) \sim \sum_{r=-P}^{P} E^{γ}_{α,β}(ze^{2πir}) + A^{γ}_{α,β}(ze^{∓πi}), \quad |\arg z| ≤ π,$$

where $P$ is an integer number such that $2P + 1$ is the smallest odd integer satisfying $2P + 1 > 1/α$. For the latter it is found that the anti-Stokes lines collapse onto the negative real axis, thus implying that the asymptotic behavior is always exponentially large. Indeed, we included the algebraic term just for the sake of completeness, though it turns out to be always negligible.

Providing an explicit formulation of the coefficients $c_k$ in the exponential expansion $E^{γ}_{α,β}(z)$ is not an easy task and a sophisticated algorithm is discussed in [117]. The first few values of the $c_k$’s have however been explicitly computed in [32] and [118]:

- $c_0 = 1$
- $c_1 = \frac{(γ - 1)}{2} (αγ + γ - 2β)$
- $c_2 = \frac{(γ - 1)(γ - 2)}{24} [3(α + 1)^2γ^2 - (α + 1)(α + 12β + 5)γ + 12β(1 + β)]$
- $c_3 = \frac{(γ - 1)(γ - 2)(γ - 3)}{48} [γ^3(1 + α)^3 - γ^2(1 + α)^2(5 + α + 6β) + 2γ(1 + α)(3 + α(1 + β) + 11β + 6β^2) - 8β(1 + β)(2 + β)]$.

Showing a longer list of these coefficients is not particularly instructive, hence we just refer again to the algorithm provided in [117].

Specific results for the asymptotic expansion along the negative real semi-axis, which is of most interest for the study of relaxation phenomena, have been provided, for instance, in [97]. In this case, focusing just on the leading term of the expansion, one can infer that

$$E^{γ}_{α,β}(-t^α) \sim \begin{cases} 
\frac{1}{Γ(β - αγ)}t^{-αγ} & β ≠ αγ \\
\frac{γ}{Γ(-α)}t^{-αγ-α} & β = αγ
\end{cases}$$

as $t \to +∞$.

### 5. Prabhakar fractional calculus

After an extensive review of the main properties of the Prabhakar function, we are now ready to introduce and discuss a generalization of fractional calculus based on this special function, that we shall call Prabhakar fractional calculus.

One of the main contributions of T. R. Prabhakar [126] consisted in the formulation of a new class of weakly-singular linear Volterra operators based on the function $e^{γ}_{α,β}(t; λ) = t^{β-1}E^{γ}_{α,β}(λt^α)$, $α > 0$, known in the literature as the Prabhakar kernel, with $λ$ being a real (or complex) number. Most notably, one finds that for $γ = 0$ or $λ = 0$ this kernel reduces to
shown that \( J^\gamma(5.8) \) i.e. Most importantly, in [80] a family of left-inverse operators of the Prabhakar fractional theory is given by

\[
(5.5) \quad J^\gamma_{\alpha,\beta,\lambda} f(t) := \left( e^{\gamma}_{\alpha,\beta}(\cdot, \lambda) * f \right) (t) = \int_0^t (t - u)^{\beta - 1} E^\gamma_{\alpha,\beta}(\lambda(t - u)^\alpha) f(u) du, \quad \alpha, \beta > 0,
\]

where the condition \( \alpha > 0 \) is necessary for the convergence of the series expansion (4.1), whereas the restriction \( \beta > 0 \) is imposed in order to guarantee the convergence of the integral. In view of (5.2), we have that the Prabhakar kernel in (5.3) reduces to the Gel’fand–Shilov distribution [42], which implies that

\[
(5.4) \quad J^0_{\alpha,\beta,\lambda} f(t) = J^\gamma_{\alpha,\beta,\lambda} f(t) = J^\gamma_{\alpha,\beta,\lambda} f(t) \quad \text{or,} \quad \text{i.e. (5.3) reduces to the RL integral for } \gamma = 0 \text{ or } \lambda = 0. \quad \text{Besides, another interesting relation between the RL and Prabhakar integrals is related to the series representation of the latter. Indeed, let } f \in L^1[t_0, T] \text{ and } \alpha, \beta > 0, \text{ then it is easy to see that (see [41])}
\]

\[
(5.5) \quad J^\gamma_{\alpha,\beta,\lambda} f(t) = \sum_{k=0}^{\infty} \frac{(\gamma)^k}{k!} J^{\gamma + k + \beta}_{\alpha,\beta} f(t).
\]

Additionally, a property that finds several applications when dealing with Prabhakar’s theory is given by

\[
(5.6) \quad J^\gamma_{\alpha,\beta,\lambda} (t - t_0)^{\mu - 1} E^{\mu}_{\alpha,\mu}(\lambda(t - t_0)^\alpha) = (t - t_0)^{\beta + \mu - 1} E^{\gamma + \mu}_{\alpha,\beta + \mu}(\lambda(t - t_0)^\alpha)
\]
or, in other terms,

\[
(5.7) \quad J^\gamma_{\alpha,\beta,\lambda} e^{\sigma}_{\alpha,\mu}(t - t_0 ; \lambda) = e^{\gamma + \sigma}_{\alpha,\beta + \mu}(t - t_0 ; \lambda).
\]

The Prabhakar fractional integral has been extensively studied in [80], where it was shown that \( J^\gamma_{\alpha,\beta,\lambda} : L_1[t_0, T] \rightarrow L_1[t_0, T] \) is a bounded operator (see [80, Theorem 4]). Most importantly, in [80] a family of left-inverse operators of the Prabhakar fractional integral was identified, i.e.

\[
(5.8) \quad D^\gamma_{\alpha,\beta,\lambda} f(t) = RL D^\mu_{t_0} J^{-\gamma}_{\alpha,\beta,\lambda} f(t) \quad \forall \mu \in \mathbb{R} \text{ with } \mu > \beta
\]

with \( f \in L_1[t_0, T] \). The presence of the fractional-order RL derivative in (5.8) makes this family of operators too convoluted for most practical applications. Hence, in most of the subsequent works (e.g., see [33, 37, 124]) the special case \( \mu = m = [\beta] \) is usually preferred. This then leads to what is now commonly known as the Prabhakar derivative of RL type, i.e.

\[
(5.9) \quad RL D^\gamma_{\alpha,\beta,\lambda} f(t) := D^m J^{-\gamma}_{\alpha,\beta,\lambda} f(t) = \frac{d^m}{dt^m} \int_0^t (t - u)^{m - \beta - 1} E^{-\gamma}_{\alpha,\beta,\lambda} f(u) du.
\]

In analogy with the classical theory of fractional differential operators, one can introduce a regularized version of the Prabhakar derivative (5.9). This is achieved by exchanging the order according to which the operations of integration and differentiation appear in (5.9). This procedure is known as the Caputo-like regularization of the fractional Prabhakar
derivative and it was first performed by M. D’Ovidio and F. Polito in [29]. Then, let 
\( f \in AC^m[t_0, T] \), the regularized Prabhakar derivative is defined as

\[
C^D_{\alpha,\beta,\lambda} f(t) := J^{-\gamma}_{\alpha,m-\beta,\lambda} D^m f(t) \\
= \int_{t_0}^{t} (t-u)^{m-\beta-1} E^{-\gamma}_{\alpha,m-\beta} (\lambda(t-u)^\alpha) f^{(m)}(u) \, du.
\]

(5.10)

This operator turns out to act as another left-inverse of the Prabhakar fractional integral, as shown in [42].

One reason why (5.10) is understood as a regularization of (5.9) is due to the fact that this procedure makes the solutions to the eigenvalue problem associated to (5.10) non-singular at \( t = t_0 \). Property which is not shared by the solutions to the eigenvalue problem associated to the (RL-type) Prabhakar derivative. Furthermore, again differently from the RL-type case, one can set up a well-posed Cauchy problem for the regularized Prabhakar derivative with initial conditions involving solely ordinary derivatives of the unknown function at \( t = t_0 \). Besides, one also has that

\[
C^D_{\alpha,\beta,\lambda} f(t) = 0, \quad k = 0, 1, \ldots, m-1, \quad m = \lceil \beta \rceil.
\]

(5.11)

It is now easy to see that for \( \gamma = 0 \) or \( \lambda = 0 \) (5.9) and (5.10) reduce to the standard RL and Caputo derivatives, namely (4.15) and (4.16). Moreover, it turns out that (5.9) and (5.10) satisfy

\[
C^D_{\alpha,\beta,\lambda} f(t) = \text{RL}^D_{\alpha,\beta,\lambda} \left[ f(t) - \sum_{k=0}^{m-1} \frac{(t-t_0)^k}{k!} f^{(k)}(t_0^+) \right],
\]

(5.12)

further strengthening their similarity to canonical fractional derivatives. Furthermore, using this last property A. Giusti [42] was then able to frame (to some extent) Prabhakar’s theory within the scheme of Kochubei’s general fractional calculus [87]. Note that, to this regard, some preliminary analysis had been carried out in [53] for a very specific problem.

Before concluding this general summary of the main definitions and properties of the Prabhakar fractional integral and derivatives, it is worth stressing that a similar result to the series expansion in (5.5) can be trivially computed also for the operators (5.9) and (5.10).

5.1. Laplace domain. It is now of paramount importance to recall a few results concerning the effects of integral transforms on Prabhakar operators.

First, let us consider the Laplace transform of the Prabhakar fractional integral, with \( t_0 = 0 \). Since (5.3) is a convolution-type integral, in the Laplace sense, it is easy to see that

\[
\mathcal{L} \left[ J_{\alpha,\beta,\lambda}^\gamma f(t) ; s \right] = \mathcal{L} \left[ e_{\alpha,\beta}(t ; \lambda) ; s \right] \tilde{f}(s) = \frac{s^{\alpha\gamma - \beta}}{(s^\alpha - \lambda)^\gamma} \tilde{f}(s),
\]

(5.12)

where in the last step we used the result in (4.17), with \( \text{Re}(s) > 0 \) and \( |s| > |\lambda|^\frac{1}{\gamma} \). Similarly, for the Prabhakar derivative (5.9) one finds

\[
\mathcal{L} \left[ \text{RL}^D_{\alpha,\beta,\lambda}^\gamma f(t) ; s \right] = s^{\beta - \alpha\gamma} (s^\alpha - \lambda)^\gamma \tilde{f}(s)
\]

(5.13)

\[- \sum_{k=0}^{m-1} s^{m-k-1} \left( J_{\alpha,m-\beta-k,\lambda}^\gamma f \right)(0^+),
\]

with \( m = \lceil \beta \rceil \), whose proof is simply based on the LT of the integer-order derivative, the result in (4.17), and that (see [80])
5.2. Eigenvalue problem. Let us consider the Cauchy problem

\[ \begin{align*}
& C^\gamma D_{\alpha,\beta,\lambda,0}^\gamma y(t) = Ay(t), \\
& y(0^+) = \xi_0, \quad y'(0^+) = \xi_1, \ldots, \quad y^{(m-1)}(0^+) = \xi_{m-1},
\end{align*} \]

with \( \alpha, \beta > 0, m = [\beta], \) and \( A, \xi_0, \ldots, \xi_{m-1} \in \mathbb{R}. \)

The Cauchy problem in (5.14) is well-posed and one can compute its solution by means of the Laplace transform method. Specifically, one finds

\[ y(t) = \sum_{j=0}^{m-1} \sum_{k=0}^{\infty} A^k t^j k^k \Gamma(\beta k + j + 1) E_{\alpha,\beta k+j+1}(\lambda t^\beta) \xi_j, \]

which thus gives the eigenfunction of \( C^\gamma D_{\alpha,\beta,\lambda,0}^\gamma \) corresponding to the eigenvalue \( A. \) Note that, if \( \gamma = 0 \) or \( \lambda = 0, \) then

\[ y(t) = \sum_{j=0}^{m-1} \sum_{k=0}^{\infty} A^k t^j k^k \Gamma(\beta k + j + 1) \xi_j = \sum_{j=0}^{m-1} j^j E_{\beta,j+1}(At^\beta) \xi_j, \]

which is indeed the eigenfunction of the Caputo derivative \( C^\gamma D_{0}^\gamma \) corresponding to the eigenvalue \( A \) (see [47]), as expected.

Moving on to the eigenvalue problem for the (RL-type) Prabhakar derivative, let us consider the Cauchy problem

\[ \begin{align*}
& RL^\gamma D_{\alpha,\beta,\lambda,0}^\gamma y(t) = Ay(t) \\
& (J_{\alpha,m—\beta—j,0}^\gamma y)(0^+) = \xi_j, \quad j = 0, 1, \ldots, m - 1.
\end{align*} \]

Following a procedure akin to the one presented above we find that the solution to this problem reads

\[ y(t) = \sum_{j=0}^{m-1} \sum_{k=0}^{\infty} A^k t^j k^k \Gamma(\beta k + j + 1) E_{\alpha,\beta k+j+1}(\lambda t^\beta) \xi_j \]

and hence the functions

\[ \sum_{k=0}^{\infty} A^k t^j k^k \Gamma(\beta k + j + 1) E_{\alpha,\beta k+j+1}(\lambda t^\beta), \quad j = 0, 1, \ldots, [\beta] - 1 \]
are eigenfunctions of $^{RLD}_{\alpha,\beta,\lambda,0} D_{0}^{\gamma}$. Besides, for $\gamma = 0$ or $\lambda = 0$, in light of (5.2) we have that Eq. (5.18) becomes

$$y(t) = \sum_{j=0}^{m-1} \sum_{k=0}^{\infty} A^{k+j} \xi_j$$

thus yielding the known eigenfunctions $t^{\beta-j-1} E_{\beta,\gamma}(A t^\beta)$, $j = 0, 1, \ldots, \lceil \beta \rceil - 1$ of the RL fractional derivative $^{RLD}_{0} D_{0}^{\beta}$ [47]. In particular, note that for $j = m - 1$ one has that $t^{\beta-j-1} = t^{\beta-\lfloor \beta \rfloor}$ which carries a weakly singular behavior if $\beta \notin \mathbb{N}$ since in this case $\beta - \lfloor \beta \rfloor < 0$.

5.3. Operators of Grünwald-Letnikov type. It is now worth mentioning that a Grünwald–Letnikov (GL) type formulation of the Prabhakar operators was proposed by R. Garrappa in [36]. Originally, this interpretation of the theory was conceived for the particular case $\beta = \alpha \gamma$ that turns out to be involved in the time-domain representation of the Havriliak–Negami relaxation. Its generalization to $\beta > \lfloor \beta \rfloor$ in [36]. Originally, this interpretation of the theory was conceived for the particular case $\beta = \alpha \gamma$ that turns out to be involved in the time-domain representation of the Havriliak–Negami relaxation. Its generalization to $\beta > \lfloor \beta \rfloor$ is however immediate and is achieved by following the same steps illustrated in [36] and by taking profit of Lubich’s theory [92–94] of generalized convolution quadratures. Indeed, one can find that, under suitable conditions, the Prabhakar integral (5.3) is equivalent to the GL integral

$$\mathcal{J}_{\alpha,\beta,\lambda,0}^{\gamma} f(t) = \lim_{h \to 0} \frac{h^{\beta}}{(1 - h^\alpha \lambda)^{\gamma}} \sum_{j=0}^{\lfloor t/h \rfloor} W_{n-j}(-) f(t - jh),$$

where the coefficients $W_{n}(-)$ depend on $\alpha$, $\beta$, $\gamma$, $\lambda$ and $h$. For $\gamma \neq 0$, $W_{n}(-)$ can be evaluated recursively as

$$W_{0}(-) = 1, \quad W_{k}(-) = \sum_{j=1}^{k} \left( \frac{1 - \gamma}{k} j - 1 \right) \bar{\omega}_j W_{k-j}(-),$$

with

$$\bar{\omega}_j = \frac{\omega_{j}^{(\beta/\gamma)} - h^\alpha \lambda \omega_{j}^{(\beta/\gamma - \alpha)}}{1 - h^\alpha \lambda},$$

while the $\omega_{j}^{(\alpha)}$ are related to the classical binomial coefficients according to

$$\omega_{j}^{(\alpha)} = (-1)^{j} \binom{\alpha}{j} \left( \frac{\Gamma(\alpha + 1)}{\Gamma(j + 1) \Gamma(\alpha - j + 1)} \right).$$

Clearly, the GL–Prabhakar integral in (5.19) generalizes the classical GL operator corresponding to the RL integral (4.14). The RL integral is instead recovered from (5.19) by setting $\lambda = 0$ since, in this case, one has $W_{j}(-) = \omega_{j}^{(-\beta)}$.

In a similar fashion, a GL operator corresponding to Prabhakar derivative $^{RLD}_{\alpha,\beta,\lambda,0} D_{0}^{\gamma}$ can also be obtained as

$$^{RLD}_{\alpha,\beta,\lambda,0} D_{0}^{\gamma} f(t) = \lim_{h \to 0} \frac{(1 - h^\alpha \lambda)^{\gamma}}{h^{\beta}} \sum_{j=0}^{\lfloor t/h \rfloor} W_{n-j}^{(+)} f(t - jh),$$

with

$$W_{0}^{(+)} = 1, \quad W_{k}^{(+)} = \sum_{j=1}^{k} \left( \frac{1 + \gamma}{k} j - 1 \right) \bar{\omega}_j W_{k-j}^{(+)}.$$
where the coefficients $\bar{\omega}_j$ are the same as the one introduced above. It is worth noting, again, that the classical GL derivative, equivalent to the RL derivative (4.15), is recovered for $\lambda = 0$.

Finally, in light of (5.11) a GL operator related to the regularized Prabhakar derivative $\mathcal{D}_{\alpha,\beta,\lambda}^{\gamma,\delta}$ can be easily obtained. The importance of these operators lies in their peculiar proneness toward their use as numerical approximation of the Prabhakar operators, once the step-size $h > 0$ has been fixed.

6. Physical applications

The aim of this section is to briefly summarize the main results and physical implications of Prabhakar calculus. Of course, the literature on this subject is rather vast and, at times, dispersive, hence our plan is to offer a general overview of the key physical features of these models that fall in the realm of anomalous phenomena. Specifically, we will start off by describing how this new framework can describe dielectric relaxation processes such as the one in the Davidson–Cole and the Havriliak–Negami models. Then, we will show how the electro-mechanical correspondence theorized by B. Gross and R. M. Fuoss paves the way to a brand new class of viscoelastic models. Finally, we summarize the main results concerning the role of the Prabhakar function (and the corresponding operators) in the theory of anomalous diffusion on complex media.

6.1. Anomalous relaxation in dielectrics. It is fairly well known in the literature [18, 19, 24, 71, 151–153] that materials that display glass-liquid transitions or amorphous polymers tend to deviate substantially from the standard Debye relaxation model. Setting aside for the moment the Cole–Cole relaxation, that can be described in terms of standard fractional calculus [12,37,49], let us focus on the Havriliak–Negami model. First, recalling that we are dealing with a causal linear system, we can recast Eq. (3.15) in the Laplace domain, thus the response function will read [12,37,50]

$$\tilde{\phi}_{HN}(s) = \tilde{\chi}_{HN}(s) = \frac{1}{1 + (s\tau^*)^\alpha}, \quad 0 < \alpha \leq 1, \quad \gamma > 0,$$

then, taking profit of Eq. (4.17), one gets [37]

$$\phi_{HN}(t) = \frac{1}{\tau^*} \left( \frac{t}{\tau^*} \right)^{\alpha\gamma-1} E_{\alpha,\alpha\gamma}^{\gamma} \left[ - \left( \frac{t}{\tau^*} \right)^\alpha \right],$$

that, in view of Eq. (3.8), leads to the relaxation function [37]

$$\Psi_{HN}(t) = 1 - \left( \frac{t}{\tau^*} \right)^{\alpha\gamma} E_{\alpha,\alpha\gamma+1}^{\gamma} \left[ - \left( \frac{t}{\tau^*} \right)^\alpha \right].$$

A few comments are now in order. First of all, it is worth mentioning the asymptotic behaviors of $\Psi_{HN}(t)$, namely

$$\Psi_{HN}(t) \sim \begin{cases} 
1 - \frac{(t/\tau^*)^{\alpha\gamma}}{\Gamma(1 + \alpha\gamma)}, & \text{for } t \ll \tau^*, \\
\gamma(t/\tau^*)^{-\alpha} \frac{\Gamma(1 - \alpha)}{\Gamma(1 - \alpha)}, & \text{for } t \gg \tau^*.
\end{cases}$$

Besides, it is also straightforward to prove that the relaxation equation for the Havriliak-Negami model is given by

$$\mathcal{D}_{\alpha,\alpha\gamma,\tau^*}^{\gamma,\delta} \Psi_{HN}(t) = -\frac{1}{\tau^*}, \quad \Psi_{HN}(0) = 1,$$
highlighting the strict connection between this model and Prabhakar calculus.

Finally, one can easily infer that the Davidson–Cole and the Cole–Cole models are particular realizations of the Havriliak–Negami relaxation. Indeed, the Davidson–Cole dielectric response is obtained from the Havriliak–Negami one by setting $\alpha = 1$ in Eq. (3.15), leading to a relaxation function that reads

$$\Psi_{DC}(t) = 1 - \left(\frac{t}{\tau^*}\right)^\gamma E_{1,\gamma+1}^\gamma \left(-\frac{t}{\tau^*}\right) = \frac{\Gamma(\gamma, t/\tau^*)}{\Gamma(\gamma)} ,$$

with $\Gamma(a, z) = \int_z^\infty u^{a-1} e^{-u} du$ denoting the incomplete gamma function, and a relaxation equation

$$C_D^{\gamma_{1,\gamma,\gamma+1}} \Psi_{DC}(t) = -\frac{1}{\tau^*} \Psi_{DC}(0) = 1 .$$

Analogously, it is easy to see that the Cole–Cole model can be derived from Eq. (3.15) by setting $\gamma = 1$. This, in turn, leads the relaxation function in Eq. (3.13) that corresponds to an evolution equation governed by Caputo derivatives [37,49], i.e.

$$C_D^{\alpha_{0,\gamma,\gamma+1}} \Psi_{CC}(t) = -\frac{1}{\tau^*} \Psi_{CC}(t) , \quad \Psi_{CC}(0) = 1 ,$$

which is indeed the canonical fractionalization of the standard relaxation problem, namely $\Psi'(t) = -(1/\tau) \Psi(t)$ with $\Psi(0) = 1$.

Before concluding this Section we wish to stress that the models discussed here clearly do not exhaust the full realm of theoretical set-ups that have been proposed, over the years, to describe the features of anomalous dielectrics. In this work we simply report the most known modifications of Debye’s theory that display a strict connection with the Prabhakar function and calculus. For a more detailed analysis of the literature on anomalous dielectric relaxation and its relation to the notion of complete monotonicity, we refer the interested reader to [12,37,97].

6.2. Linear Viscoelasticity. It is rather well known in the literature (see e.g. [13,14,20, 21, 44, 128] and references therein) that fractional calculus and the theory of completely monotone functions play a pivotal role and lie at heart of the modern mathematical formulation of linear viscoelasticity. To this regard, we refer the readers to [22,60,96,101,119] for a detailed description of both ordinary and fractional viscoelasticity.

Beside the clear effectiveness of Prabhakar calculus in serving as a fundamental tool for modeling anomalous dielectrics, this novel mathematical structure plays an interesting role also in fractional viscoelasticity. As discussed in [43], if one recalls the formal duality between viscoelastic models and electrical systems, originally introduced by Gross and Fuoss in [56–58] and then revisited in [45], it is easy to see that the anomalous relaxation processes highlighted in Section 6.1 have an analogous representation in linear viscoelasticity. It is therefore natural to wonder about what might happen if one writes the constitutive equation for a certain material in terms of Prabhakar derivatives.

Let us begin with a brief recap of the basics of the linear theory of viscoelasticity. Denoting by $\mathcal{H}^N$ the Heaviside class, namely the set of causal functions $f(t)$ such that $f \in C^N(\mathbb{R}^*_+)$ with $N \in \mathbb{N}$, then we denote by $\sigma \in \mathcal{H}^N$ and $\varepsilon \in \mathcal{H}^N$ the uniaxial stress and strain functions for a given material, respectively. The key physical information about the system are then stored in the stress-strain relation, which is known as the constitutive equation for the material, and it can be represented equivalently in differential form or as an integral equation. In particular, the distinctive feature of the integral form is that it relates the stress and the strain through either the creep compliance $J(t)$ or the relaxation
modulus $G(t)$. In detail, if one considers a quiescent viscoelastic body for $t < 0$ and assumes some sufficiently well behaved causal histories, then the general integral form of the constitutive equation reads

$$
\varepsilon(t) = J(0^+) \sigma(t) + \int_0^t J(t-\tau) \sigma'(\tau) \, d\tau,
$$

or equivalently,

$$
\sigma(t) = G(0^+) \varepsilon(t) + \int_0^t G(t-\tau) \varepsilon'(\tau) \, d\tau.
$$

Since the latter are convolution-type equations, they can be easily manipulated via the standard Laplace transform method. Specifically, these equations can be recast, in the Laplace domain, as

$$
\tilde{\varepsilon}(s) = s \tilde{J}(s) \tilde{\sigma}(s), \quad \tilde{\sigma}(s) = s \tilde{G}(s) \tilde{\varepsilon}(s),
$$

from which one can infer the so called reciprocity relation

$$
s \tilde{J}(s) = \frac{1}{s \tilde{G}(s)}.
$$

Furthermore, it is useful to recall that the creep compliance is a non-negative and non-decreasing function of time on $t \geq 0$, whereas the relaxation modulus is a non-negative and non-increasing function of time on the same domain.

All that being said, in [43] it was proposed a model representing a straightforward generalization of the renowned Maxwell model of linear viscoelasticity, where the ordinary derivatives were replaced with Prabhakar’s one. In detail, this new proposal is customarily referred to as Maxwell–Prabhakar model (or, alternatively, as the Giusti–Colombaro model) of linear viscoelasticity and it is described by the constitutive equation

$$
\sigma(t) + a C_D^{\gamma,\alpha,\beta,\lambda} \sigma(t) = b C_D^{\gamma,\alpha,\beta,\lambda} \varepsilon(t)
$$

where $\alpha > 0$, $0 < \beta < 1$ and $\gamma, \lambda \in \mathbb{R}$, in general. Besides, since $[\sigma] = \text{Force} \cdot \text{Area}^{-1}$ and $[\varepsilon] = 1$, one has that $[a] = \text{time}^\beta$, $[b] = \text{Force} \cdot \text{Area}^{-1} \cdot \text{time}^\beta$, and $[\lambda] = \text{time}^{-\alpha}$. Bringing Eq. (6.13) to the Laplace domain one gets

$$
s \tilde{J}(s) = \frac{a b}{b s^\beta (1 - \lambda s^{-\alpha})^\gamma} = \frac{1}{s \tilde{G}(s)}
$$

which, in turn, yields

$$
J(t) = \frac{a}{b} + \frac{t^\beta}{b} E_{\alpha,\beta+1}^\gamma(\lambda t^\alpha),
$$

and

$$
G(t) = \frac{b}{a} \sum_{n=0}^{\infty} \left( -\frac{t^\beta}{a} \right)^n E_{\alpha,\beta n+1}^{\gamma n}(\lambda t^\alpha),
$$

where the last series is absolutely convergent on $t > 0$ for the whole domain of parameters considered here.

It is also worth mentioning that this model is able to reproduce, as particular cases, the Maxwell, Voigt, and Zener models in both their fractional and ordinary realizations. For details and constraints on the parameter space see [43]. Furthermore, a thorough study of the processes of storage and dissipation of energy in materials described by (6.13) has been carried out by I. Colombaro, A. Giusti, and S. Vitali in [23].
6.3. Stochastic Processes and Diffusion. It is now worth discussing the implications of Prabhakar operators for stochastic processes, and their relation with anomalous diffusion phenomena. For a detailed review on this topic we refer the interested reader to [137,140,145].

It is well-known that the probability density function (PDF) for the location $x$ at time $t$ of a Brownian particle satisfies the classical heat equation. In this case, the mean squared displacement (MSD) is linear in time. Different alternative non-Brownian models have been studied in the literature in order to describe anomalous diffusions in complex media. A widely used and general scheme is given by the Continuous-Time Random Walk (CTRW) approach, first introduced in [109]. The structure of this stochastic jump process is described by specifying the distributions of independent random jump lengths and waiting times (see e.g. [86]).

In [133], a general class of integro-differential diffusion-type equations given by

$$
\frac{\partial}{\partial t} W(t,x) = \frac{\partial}{\partial t} \int_0^t \eta(t-t') \frac{\partial^2}{\partial x^2} W(t',x) \, dt',
$$

is considered. Here, $W(t,x)$ is the PDF relative to the location of a CTRW admitting jumps of finite variance and waiting times with a specific distribution. In particular, the memory kernel $\eta(t)$ in (6.17) is connected to the PDF $\psi(t)$ of the random waiting times through their Laplace transforms according to

$$
\tilde{\psi}(s) = \frac{\tilde{\eta}(s)}{1 + \tilde{\eta}(s)},
$$

(6.18)

Summarizing, Eq. (6.17) is the governing equation of a family of CTRWs, given some suitable constraints on the memory kernel $\eta(t)$.

One of the interesting cases considered in [137] is based on the application of the Prabhakar integral. In this respect, let us consider, for instance, the case of the tempered time-fractional diffusion equation with $\eta(t)$ such that

$$
\gamma(t) = e^{-bt} \frac{t^{-\alpha}}{\Gamma(1 - \alpha)},
$$

(6.19)

with $0 < \alpha < 1$ and $b > 0$ denoting the truncation parameter. The corresponding memory kernel then reads

$$
\eta(t) = t^{\alpha-1} E_{1,\alpha}^{-1} (-bt),
$$

(6.20)

leading to

$$
\frac{\partial}{\partial t} W(t,x) = R_t \mathcal{D}^{1-\alpha}_{\alpha,-1} \left[ \frac{\partial^2}{\partial x^2} W(t,x) \right].
$$

(6.21)

The MSD turns out to be $2t^\alpha E_{1,\alpha+1}^{\alpha-1} (-bt)$, from which one can immediately infer the emergence of a crossover from subdiffusion to normal diffusion.

If instead we reverse the problem by considering the case with

$$
\eta(t) = e^{-bt} \frac{t^{-\alpha}}{\Gamma(1 - \alpha)},
$$

(6.22)

then one finds

$$
C_{\alpha,-1} \mathcal{D}^{\alpha}_{\alpha+1,-1} W(t,x) = \frac{\partial^2}{\partial x^2} W(t,x).
$$

(6.23)

Correspondingly, the related MSD becomes $2t^\alpha E_{1,\alpha+1}^{\alpha-1} (-bt)$, implying a crossover from subdiffusion to a constant behavior.
Working in the framework of the theory of CTRWs, it was also shown [135] that one can build models aimed at characterizing the transition among anomalous diffusion scenarios based on the Prabhakar function and its derivative operators. Finally, it is also worth mentioning that a study of the generalized Langevin Equation with both the Prabhakar and tempered Prabhakar kernels as memory kernels has been carried out by T. Sandev in [131].

7. Applications in renewal processes

Generalizations of renewal processes based on the applications of time-fractional differential operators, or special functions related to fractional calculus, have recently attracted much attention. Different (non equivalent) approaches to build fractional counting processes have been developed in the literature. The motivation for the increasing interest on this topic is obviously related to the pervasive applications of classical counting processes, like the Poisson process, in real-world models.

The Mittag–Leffler function appeared as residual waiting time between events in renewal processes already in the 1960s, namely processes with properly scaled thinning out the sequence of events in a power law renewal process (see [46] and [100]). A renewal process with Mittag–Leffler–distributed waiting times is in essence a fractional Poisson process. Hilfer and Anton [70] were the first to introduce the Mittag–Leffler waiting-time density

$$f_{\mu}(t) = -\frac{d}{dt}E_{\mu}(\mu t) = \mu^{\mu-1}E_{\mu,\mu}(\mu t)$$

in the context of continuous time random walk. Specifically, they were able to show that such a waiting time density is required in order for the evolution of the sojourn density to be governed by a fractional Kolmogorov–Feller equation. In other words, they time-changed a random walk with an independent fractional Poisson process. Systematic studies of both analytic and probabilistic aspects of this topic were then triggered, to the best of our knowledge, in 2000 by the work of O. N. Repin and A. I. Saichev [127]. Many (mostly independent) contributions investigated the feature of this new renewal process following the path paved by the previously mentioned pivotal works, see e.g. [3, 4, 90, 103, 122] and references therein. To this regard, it is worth stressing that the Prabhakar function naturally arises in the context of fractional Poisson processes. Indeed, the state probabilities of a time-fractional Poisson process can be expressed in terms of Prabhakar functions, see e.g. [4,90] for details.

Recently, several generalizations of the Poisson process based on the Prabhakar function have appeared in the literature. These extensions clearly offer more flexibility when attempting to capture the main features of real-world renewal processes, also including as special cases the time-fractional and the classical Poisson processes. Essentially, the generalization of the standard Poisson process can be achieved in two different ways: i) deriving the state probabilities of the generalized counting process by solving an infinite system of time-fractional difference-differential equations involving regularized Prabhakar derivatives; ii) considering a renewal process with inter-event time density function involving a Prabhakar function. The first approach was originally developed by R. Garra, et al. in [33], whereas the second one was first proposed by D. O. Cahoy and F. Polito in [9] (see also [104,105]). Further, an alternative proposal by T. K. Pogány and Ž. Tomovski was provided in [121], where the generalization was obtained in the sense of weighted Poisson distributions.
Let us now analyze the first approach, namely the one based on the replacement of the first-order derivative with the regularized Prabhakar derivative. In other words, this procedure has to be implemented onto the infinite system of difference–differential equations governing the state probability of the Poisson process, \( i.e. \)

\[
\begin{cases}
\mathcal{C}D_{\rho,\mu,-\phi}^{\gamma} p_k(t) = -\lambda p_k(t) + \lambda p_{k-1}(t), & k \geq 0, \ t > 0, \ \lambda > 0,

p_k(0) = \begin{cases}
1, & k = 0,
0, & k \geq 1,
\end{cases}
\end{cases}
\]

(7.2)

where \( \phi > 0, \ \gamma \geq 0, \ 0 < \rho \leq 1, \ \text{and} \ 0 < \mu \leq 1. \) If \( \gamma \neq 0 \) one also has that \( 0 < \mu[\gamma]/\gamma - r \rho < 1, \ \forall r = 0, \ldots, \lceil \gamma \rceil. \) These constraints on the parameters are required in order to ensure non-negativity of the solution. Multiplying both terms of (7.2) by \( v_k \), with \( v \) denoting an auxiliary variable such that \( |v| \leq 1, \) and summing over all \( k, \) we obtain the fractional Cauchy problem for the probability generating function \( G(v, t) = \sum_{k=0}^{\infty} v^k p_k(t) \) of the counting number \( N(t), \ t \geq 0, \)

\[
\begin{cases}
\mathcal{C}D_{\rho,\mu,-\phi}^{\gamma} G(v, t) = -\lambda (1 - v) G(v, t), & |v| \leq 1,

G(v, 0) = 1,
\end{cases}
\]

(7.3)

whose solution reads (see [33] for details)

\[
G(v, t) = \sum_{k=0}^{\infty} (-\lambda t^\mu)^k (1 - v)^k E_{\rho,\mu k+1}^{\gamma k}(-\phi t^\rho), \quad |v| \leq 1.
\]

(7.4)

It is easy to see that, for \( \gamma = 0, \) the latter reduces to

\[
G(v, t) = \sum_{k=0}^{\infty} (-\lambda t^\mu)^k (1 - v)^k \frac{(\mu k)!}{(\mu k + 1)!} E_{\mu,1}^{1}(-\lambda(1 - v)t^\mu),
\]

(7.5)

that coincides with the probability generating function of the (standard) fractional Poisson process, see e.g. [90].

From the probability generating function (7.4) it is possible to infer the probability distribution at fixed time \( t \geq 0 \) of \( N(t), \) governed by the evolution equation (7.2). Indeed, a simple binomial expansion leads to

\[
G(v, t) = \sum_{k=0}^{\infty} v^k \sum_{r=k}^{\infty} (-1)^r-k \binom{r}{k} (\lambda t^\mu)^r E_{\rho,\mu r+1}^{\gamma r}(-\phi t^\rho),
\]

(7.6)

that implies

\[
p_k(t) = \sum_{r=k}^{\infty} (-1)^r-k \binom{r}{k} (\lambda t^\mu)^r E_{\rho,\mu(r+1)}^{\gamma r}(-\phi t^\rho), \quad k \geq 0, \ t \geq 0.
\]

(7.7)

From (7.2), one can also infer the mean value of \( N(t). \) To this aim, it suffices to differentiate Eq. (7.3) with respect to \( v \) and set \( v = 1, \) this leads to

\[
\begin{cases}
\mathcal{C}D_{\rho,\mu,-\phi}^{\gamma} E N(t) = \lambda, & t > 0,

E N(t) \bigg|_{t=0} = 0,
\end{cases}
\]

(7.8)

whose solution is simply given by

\[
E N(t) = \lambda t^\mu E_{\mu,1+\mu}^{\gamma}(-\phi t^\rho), \quad t \geq 0.
\]

(7.9)
It is also possible to prove that the generalized fractional Poisson process \( N(t) \) described so far can be constructed as a renewal process with specific waiting times. Indeed, consider \( k \) independent and identically distributed (i.i.d.) random variables \( T_j, j = 1, \ldots, k \), representing the inter-event waiting times, with probability density functions

\[
 f_{T_j}(t_j) = \lambda t_j^{\mu - 1} \sum_{r=0}^{\infty} (-\lambda t_j)^r E_{\rho,\mu r+\mu}^\gamma (-\phi t_j^\mu), \quad t_j \geq 0, \mu \in (0,1).
\]

The probability distribution \( Pr(N(t) = k), k \geq 0 \), can then be obtained by making the renewal structure explicit and comparing the Laplace transform of the state probabilities obtained from the two approaches (see [33] for details).

Lastly, it is worth remarking that the generalized fractional Poisson process \( N(t) \), similarly to the classical fractional Poisson process, can be represented as a time-changed Poisson process in which the random change of time is given by a suitable independent stochastic process with increasing paths. See Section 5.3.1 of [33] for details.

The second approach, based on the construction of a renewal process with inter-event time density function involving a Prabhakar function, was originally proposed in [9]. In this case the generalized Poisson process \( N_{\nu,\delta}(t) \) denotes a renewal process with i.i.d. random waiting times distributed as

\[
 f_{\nu,\delta}(t) = \lambda \nu^\delta \nu^{-1} E_{\nu,\delta
u}^\delta (-\lambda t^\nu), \quad \nu > 0, \delta \in \mathbb{R}, \nu \in (0,1].
\]

Clearly, the latter is a generalization of the (canonical) time-fractional Poisson process, recovered for \( \delta = 1 \), and of the classical Poisson process, which is instead reproduced when setting \( \delta = \nu = 1 \). Then, by means of the Laplace transform method, it is possible to prove that the state probabilities \( p_{\nu,\delta}^k(t) = Pr(N_{\nu,\delta}(t) = k) \) read

\[
 p_{\nu,\delta}^k(t) = \lambda^\delta \nu^\delta \nu^{-1} E_{\nu,\delta
u}^\delta (-\lambda t^\nu) - \lambda^{\delta(k+1)} \nu^\delta(k+1) E_{\nu,\delta
u}^\delta(k+1) (-\lambda t^\nu)
\]

and satisfy the Volterra equation

\[
 p_{\nu,\delta}^k(t) = \lambda^\delta \left( \mathcal{J}_{\nu,\delta
u}^{\nu,\delta} \left( p_{\nu,\delta}^{k-1}(t) \right) \right)
\]

involving the Prabhakar fractional integral.

Finally, we briefly recall that a further proposal for the construction of a generalized Prabhakar-Poisson distribution, based on the Poisson distribution approach, was developed in [121]. The probability mass function of a weighted Poisson process is of the form (we adopt the notation in [1] and references therein)

\[
 P\{N^w(t) = n\} = \frac{w(n)p(n,x)}{E[w(N)]}, \quad n \geq 0,
\]

where \( N \) is a random variable with a Poisson distribution \( p(n,x) = (x^n/n!) e^{-x} \), \( w(\cdot) \) is a non-negative weight function with non-zero finite expected value, i.e.

\[
 0 < E[w(N)] = \sum_{n=0}^{\infty} w(n)p(n,x) < \infty.
\]

Therefore, it is possible to build a Prabhakar-based generalization of the Poisson distribution by choosing the weights as \( w(n, \gamma, \alpha, \beta) = \Gamma(n+\gamma)/\Gamma(\gamma)\Gamma(\alpha n + \beta) \). This procedure provides a generalization of the Poisson distribution, but it does not represent a renewal process unlike the other approaches. On the other hand, this last generalization can be useful in the context of sub- or super-Poissonian distribution applications.
8. Numerical aspects

The numerical evaluation of the Prabhakar function is a quite delicate and involved topic that appears to have been loosely discussed in the current literature, except in a couple of works in [35, 146].

The simplest way to numerically evaluate the Prabhakar function consists in relying on its definition (4.1). Namely, for such an approach it would be sufficient to fix a sufficiently large, though finite, number $K \in \mathbb{N}$ and approximate the function by the truncated series

$$E^{\gamma}_{\alpha,\beta}(z) \approx \frac{1}{\Gamma(\gamma)} \sum_{k=0}^{K} \frac{\Gamma(\gamma + k) z^k}{k! \Gamma(\alpha k + \beta)}.$$  

(8.1)

However, some critical issues emerge when this simple scheme is implemented. Indeed, the Gamma function grows very fast as its argument gets larger. This means, from a numerical perspective, that a restriction on $K$ has to be imposed to avoid overflow, i.e. the generation of numbers exceeding the range that can be represented with a given number of digits. In the standard IEEE-754 double precision arithmetic of a commonly used computer $1.8 \times 10^{308}$ is approximately the largest floating-point number that can be represented. Thus, since $\Gamma(171.624) \approx 1.8 \times 10^{308}$, the maximum number of terms in the truncated series (8.1) is bounded by $K < (171.624 - \beta)/\alpha$. This bound heavily restricts the applicability of this approach since the series (4.1) converges very slowly for arguments with moderate or large modulus. Furthermore, additional problems arise when trying to deal with arguments outside the positive real axis, in particular when $|z| > 1$. In this case, as the value of $k$ increases the sum (8.1) has consecutive terms with large modulus but opposite sign whose sum is extremely ill-conditioned in the finite-precision arithmetic of computers. As a consequence, one often incurs in catastrophic numerical cancellation. Summing up, the approximation (8.1) can be applied only for arguments with small, or just very moderately large, modulus. Otherwise, this approach turns out to be unreliable. This is, after all, the same conclusion one reaches by carrying out a similar analysis for the Mittag–Leffer function and its derivatives (see e.g. [40, 48]).

A viable strategy for the numerical evaluation of the Prabhakar function for arguments with large modulus is to relay on some of the asymptotic expansions presented in Section 4.5. Whereas, the remaining cases can be addressed by working backward from the Laplace domain.

Regarding the method based on the Laplace transform, the very simple expression given in (4.17) suggests to numerically compute the Prabhakar function by means of the inversion of its Laplace transform. J. A. C. Weideman and L. N. Trefethen paved the way for the current developments of this approach, analyzing the case $\beta = \gamma = 1$ [154]. Later on, the analysis was extended to any $\beta$ and $\gamma = 1$ [39] and, successively, to any $\gamma > 0$ for $|\arg z| > \alpha \pi$ and $0 < \alpha < 1$ [35]. Then, these results were further extended to more general arguments $z$, with $\gamma \in \mathbb{N}$ in [40].

To give just a general idea of this last technique, the starting point is represented by the expression

$$E^{\gamma}_{\alpha,\beta}(z) = \frac{1}{2\pi i} \int_{\hat{C}} e^{s} H(s; z) \, ds,$$

(8.2)

obtained from the formula for the inversion of the Laplace transform, where the considered contour $\hat{C}$ is a parabola that begins and ends in the left half-plane, such that $\text{Re}(s) \to -\infty$ at each end. In this way, the integrand in (8.2) rapidly decays as $\text{Re}(s) \to -\infty$, being forced by the exponential factor. Moreover, the contour $\hat{C}$ must encompass all the singularities of
\( H(s; z) \) (that would all be lying on the left-hand side of the vertex of \( \bar{C} \)) and the branch-cut imposed on the negative real semi axis to make \( H(s; z) \) single-valued. All that being said, the trapezoidal quadrature rule can now be profitably applied for the numerical evaluation of (8.2). However, this procedure requires to accurately fix some parameters. Indeed, in order to define the contour \( \bar{C} \) one parameter needs to be set, while the trapezoidal rule requires fixing the number of nodes and the spacing between them. In practice, if
\[
w(u) = \mu(iu + 1)^2, \quad -\infty < u < \infty,
\]
describes the parabolic contour, \( h > 0 \) is a given step-size and the quadrature nodes are defined as \( u_k = kh \) for \( k = -N, \ldots, N \), then the resulting approximation to (8.2) reads
\[
(8.3) \tilde{E}^{[N,h]}(t) = \frac{h}{2\pi i} \sum_{k=-N}^{N} e^{w(u_k)} H(w(u_k); t)w'(u_k).
\]
To choose the optimal parameters \( N \), \( h \) and \( \mu \) a deep error analysis must be carried out by requiring that the global error is proportional to the required accuracy (usually something close to the machine precision to force the method to provide the highest possible accuracy). Specifically, the influence of the parameters \( N \), \( h \) and \( \mu \) on the round-off, truncation, and discretization errors have to be considered. Basically, this means that one needs to find the most suitable contour \( \bar{C} \). Incidentally, such a contour, roughly speaking, turns out to be the one whose branches are the farthest from the singularities of the integrand. Now, once the required accuracy is fixed, the algorithm finds the optimal parameters and then computes (8.3). In particular, the value \( N \) is kept as small as possible and, since it determines the computational cost of the procedure, the resulting algorithm turns out to be fast and very accurate (see [35] for details).

If the singularities of \( H(s; t) \) turn out to be too scattered in the complex plane and too far away from both the origin and the branch-cut imposed on the negative real semi axis, it is essential to allow the contour to go over some of the singularities and apply to (8.2) the residue subtraction formula
\[
E^\gamma_{\alpha,\beta}(z) = \sum_{s^* \in S^*_C} \text{Res}(e^s H(s; z), s^*) + \frac{1}{2\pi i} \int_{\bar{C}} e^s H(s; z) \, ds,
\]
with \( S^*_C \) the set of the singularities of \( H(s; z) \) laying on the right-hand side of the contour \( \bar{C} \). This correction can, however, be implemented just for \( \gamma \in \mathbb{N} \), otherwise the more involved branch-cut due to the two real powers in \( H(s, z) \) does not allow the contour to pass over any of the singularities. This is the reason for which the method described in [35], and corresponding Matlab code [34], are restricted to \( 0 < \alpha < 1 \) and \( |\arg z| > \alpha \pi \), whilst in [40] it is discussed the case \( \gamma \in \mathbb{N} \) and a formula for the residues \( \text{Res}(e^s H(s; z), s^*) \) is given.

9. Discussion and outlook

And here is where this Hitchhiker’s Guide to the wonders of Prabhakar fractional calculus comes to an end. In Section 2 we have offered a brief recap of the main events that have led to the discovery of the Prabhakar function and the associated fractional operators. In Section 3 we have shown how this generalization was not just a mathematical curiosity, but rather something that physics was begging for in order to explain some peculiar phenomena. After that, in Section 4 we have provided an extensive description of the mathematical features of the Prabhakar function. Then, in Section 5 we combined the results in Section 4 with the general wisdom of the standard (Riemann–Liouville and
approaches to fractional calculus in order to reconstruct Prabhakar’s theory. In Section 6 we have then framed the physical problems raised in Section 3 within this new scheme, thus summarizing the main physical application of this general set-up. In Section 7 we summarized how Prabhakar’s theory can be applied in probability theory, with particular regard for the theory of renewal processes. Finally, in Section 8 we have discussed the state of the art of the research on the numerical evaluation of the Prabhakar function.

From a physical perspective, setting aside the many possible scenarios in which Prabhakar’s theory might potentially emerge as a powerful modeling tool, it is worth stressing that much of the fundamental physics connected with the phenomena discussed in Section 6 is still uncharted territory. First, while this mathematical framework seems to be the key for the mathematical description of anomalous dielectrics of the Havriliak–Negami type, a first-principle derivation of these physical behaviors is still missing. Second, even though the study of viscoelastic applications of Prabhakar calculus, initiated by A. Giusti and I. Colombaro with [43], has attracted some attention in the literature, most physical aspects of this extension of linear viscoelasticity are yet to be unveiled. These two research lines represent the main open questions concerning the applications of Prabhakar fractional calculus to “real-world problems”.

From a probabilistic point of view, we can recall that in [104,105] the authors analyzed generalized fractional diffusion models on undirected graphs and infinite lattices. This application, making use of Prabhakar random variables and therefore of Prabhakar functions, seems to be a promising topic of research in the context of non-Markovian models on networks. Moreover, further investigations on applications of Prabhakar integrals and derivatives to more general or different point processes, such as birth-death processes or point processes related to macroevolutionary models (see e.g. [123]), are surely worth pursuing. Finally, according to [88], the applications of fractional renewal processes based on Prabhakar functions to models for neural spiking events could represent compelling research subjects.

From a numerical viewpoint, a further boost should be given toward the development of methods for the computation of the Prabhakar function, in order to cover a larger range of parameters. Moreover, another stimulating research topic, that supposedly will attract much attention in the future, concerns the possibility of developing highly efficient methods for solving differential equations with Prabhakar derivatives, particularly for applications in computational electromagnetism (see e.g. [6,15,38]).

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Authors’ Contributions

Andrea Giusti has conceived the idea, organized, and coordinated the work. He has collected the contributions of the authors and he actively worked on Sections 3, 4, 5, and 6. He has also prepared the final version of the text. Ivano Colombaro has contributed to the content of Sections 3 and 6. Roberto Garra has coordinated and contributed to the work on Sections 4 and 7. He also offered valuable insights on the content of Section 6.3. Roberto Garrappa has coordinated and worked on Sections 8. He has also contributed to the work on Sections 2, 4, and 5. Federico Polito has contributed to the work on Sections 4, 7, and he has offered valuable insights on the content of Section 6.3. Marina Popolizio has worked on Section 8. Francesco Mainardi has contributed to Section 2 and he has offered his guidance and insight to all the parts appearing in this work. Finally, all authors have approved the content of the manuscript.

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