



# Article Aspects of Quantum Statistical Mechanics: Fractional and Tsallis Approaches

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**Abstract:** We investigated two different approaches, which can be used to extend the standard quantum statistical mechanics. One is based on fractional calculus, and the other considers the extension of the concept of entropy, i.e., the Tsallis statistics. We reviewed and discussed some of the main properties of these approaches and used the thermal Green function formalism to perform the developments, simultaneously allowing us to analyze each case's dynamics and thermodynamics aspects. In particular, the results allow us to understand how the extensions change the behavior of some quantities, particularly fluctuations related to the system.

Keywords: fractional dynamics; Tsallis entropy; power-law distributions; energy fluctuations

MSC: 82B03; 82B05; 82B10

# 1. Introduction

Statistical mechanics is one of the most intriguing fields of physics, where dynamics and probability are used for analyzing the behavior of systems in connection with thermodynamics. The foundations of this field resulted from intense discussions between adherents of different points of view, with the ideas first formulated by Boltzmann as a crucial starting point [1–4]. It has been successfully applied in many systems [5,6], allowing us to obtain a suitable description. On the other hand, the existence of many different phenomena, such as self-gravitating systems [7], turbulent systems [8], quantum systems with long-range interactions [9], non-Markovian systems [10,11], and anomalous diffusion [12], among others, which are characterized, for example, by fractal and multifractal aspects [13], has motivated the extension of the approach used in these contexts to obtain a suitable description. One of the possibilities for facing systems with unusual properties, such as long-range interactions, fractal and or multifractal aspects, was proposed in Ref. [14]. It considers the following extension for the Boltzmann–Gibbs entropy, namely Tsallis entropy,

$$S_q = k \frac{1 - \mathrm{Tr} \rho^q}{q - 1},\tag{1}$$

where the parameter q can be connected to different aspects of the system [15]. Equation (1) recovers the standard Boltzmann–Gibbs entropy for  $q \rightarrow 1$ , i.e.,  $S_q \rightarrow S_{BG} = -k_B \text{Tr} \rho \ln \rho$ , where  $k_B$  is the Boltzmann constant. Other entropies have also been considered to extend the standard approach, such as the Renyi entropy [16,17] and the Kaniadakis entropy [18,19]. The approach based on Equation (1) has been applied in different scenarios such as relativistic gas [20], Bose–Einstein condensation [21,22], seismology [23], and high energy

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**Copyright:** © 2023 by the authors. Licensee MDPI, Basel, Switzerland. This article is an open access article distributed under the terms and conditions of the Creative Commons Attribution (CC BY) license (https:// creativecommons.org/licenses/by/ 4.0/). multiparticle production [24]. Simultaneously with these developments, fractional calculus has been advanced in different fields as a powerful approach to incorporating different aspects of the system with extensions of the differential operators to a noninteger order. It has been applied in several scenarios, such as anomalous diffusion [25–27], anomalous charge transport in semiconductor [28], chaos [29], magnetic resonance imaging [30–34], and electrical impedance [35,36]. The fractional approach has also been used in thermodynamics [37–39] and/or in statistical mechanics [40–42]. In the first case, it extended the differential operator in the thermodynamic relations to a noninteger order fractional operator. In this context, ideal gas and gas with a weak interaction were analyzed and compared with experimental data [39]. Fractional calculus has also been used to extend statistical mechanics through fundamental equations such as the Liouville equation [43], with the possibility of applications in different scenarios, including long-range interactions and fractal and multifractal aspects, among others.

Here, we analyzed the quantum statistical mechanics in these approaches by focusing on fractional calculus and Tsallis formalism by extending and revisiting some results in the literature [44–46]. These results allow us to understand how the dynamic aspects and statistical weight changes the behavior of the thermodynamic quantities, particularly fluctuations related to the system.

## 2. Generalized Statistical Mechanics

Let us start our analysis of two possibilities for extending statistical mechanics, firstly by considering the case related to the fractional operators. Next, we considered the Tsallis formalism, which has been developed from the extension of the concept of entropy. After these developments, we analyzed the mixing between the fractional approach and a general entropy, which has the Tsallis entropy as a particular case. We focused on the quantum statistics for these cases and used the formalism employed in Ref. [47] regarding the thermal Green function method, which allowed us to evidence the differences among these cases.

#### 2.1. Fractional Dynamics and Statistical Mechanics

We shall consider the following effective Hamiltonian:

$$H_{\alpha} = \int d\mathbf{r} \int_{0}^{1} d\alpha' p(\alpha') \frac{1}{2m_{\alpha'}} \psi^{\dagger}(\mathbf{r}, t) \left(-\hbar^{2} \nabla^{2}\right)^{\alpha'} \psi(\mathbf{r}, t) + \int d\mathbf{r} \psi^{\dagger}(\mathbf{r}', t) Y(|\mathbf{r} - \mathbf{r}'|) \psi(\mathbf{r}', t) + \frac{1}{2} \int d\mathbf{r} \int d\mathbf{r}' \psi^{\dagger}(\mathbf{r}, t) \psi^{\dagger}(\mathbf{r}', t) V(|\mathbf{r} - \mathbf{r}'|) \psi(\mathbf{r}, t) \psi(\mathbf{r}', t), \qquad (2)$$

where  $\psi^{\dagger}(\mathbf{r}, \mathbf{t})$  and  $\psi(\mathbf{r}, \mathbf{t})$  are second quantized operators,  $m_{\alpha}$  is an effective constant,  $p(\alpha)$  is a distribution,

$$\left(-\hbar^2 \nabla^2\right)^{\alpha'} \psi(\mathbf{r}, t) \equiv \int \frac{d\mathbf{p}}{(2\pi\hbar)^3} e^{i\mathbf{p}/\hbar \cdot \mathbf{r}} |\mathbf{p}|^{2\alpha'} \psi(\mathbf{p}, t)$$
(3)

is the quantum Riesz operator [26,27],  $Y(|\mathbf{r}|, t)$  is a nonlocal term, and the last term is the interaction between the components of the system. The extension of the kinetic energy to fractional operators also incorporates a nonlocal behavior and recovers the usual one for  $p(\alpha') = \delta(\alpha' - 1)$ . It is also important to mention that the nonlocal term is such that  $Y(\mathbf{r}, t) = Y^{\dagger}(\mathbf{r}, t)$ , to preserve the interpretations concerning  $\psi(\mathbf{r}, r)$  as in the standard case in terms of the field operator. In addition, this Hamiltonian extends the one discussed in Refs. [44,45] with the presence of the fractional derivative in the space of distributed order and the nonlocal term. It should be stressed that Equation (2) can also be considered a phenomenological approach, where the parameters are adjusted to capture the behavior exhibited by the experimental data. In this sense, different systems may interact differently

and require different parameter values. The Heisenberg equation yields for  $\psi(\mathbf{r}, \mathbf{t})$  the following equation of motion:

$$i\hbar\frac{\partial}{\partial t}\psi(\mathbf{r},t) = \int_{0}^{1} d\alpha' p(\alpha') \frac{1}{2m_{\alpha'}} \left(-\hbar^{2}\nabla^{2}\right)^{\alpha'} \psi(\mathbf{r},t) + \int d\mathbf{r}\Upsilon(|\mathbf{r}-\mathbf{r}'|)\psi(\mathbf{r}',t) + \frac{1}{2}\int d\mathbf{r}' V(|\mathbf{r}-\mathbf{r}'|)\psi^{\dagger}(\mathbf{r}',t)\psi(\mathbf{r}',t)\psi(\mathbf{r},t).$$
(4)

Equation (4) is a Schrödinger-like equation with a nonlocal term and a fractional derivative applied to the spatial variable instead of the usual differential operators applied to the spatial variable. This equation in a first quantized perspective, without the interaction term, was analyzed in several scenarios [45,48,49]. In particular, the solution for the free case (i.e., absence of external fields and without interaction) reminds us of the Lévy distribution. By using the Fourier transform, we can simplify Equation (4) to

$$i\hbar \frac{\partial}{\partial t} \widetilde{\psi}(\mathbf{p}, \mathbf{t}) = \int_{0}^{1} d\alpha' p(\alpha') \frac{1}{2m_{\alpha'}} |\mathbf{p}|^{2\alpha'} \widetilde{\psi}(\mathbf{p}, t) + \widetilde{Y}(|\mathbf{p}|) \widetilde{\psi}(\mathbf{p}, t) + \frac{1}{2} \int d\mathbf{p}' \widetilde{V}(|\mathbf{p}' - \mathbf{p}|) \widetilde{\psi}^{\dagger}(\mathbf{p}', t) \widetilde{\psi}(\mathbf{p}', t) \widetilde{\psi}(\mathbf{p}, t).$$
(5)

From this equation, we can observe that the presence of the fractional derivative changes the exponent of the momentum, and the solution in the absence of interaction potentials is given by

$$\widetilde{\psi}(\mathbf{p},t) = e^{-it|\mathbf{p}|^{2\alpha}/(2m_{\alpha}\hbar)}e^{-itY(|\mathbf{p}|)/\hbar}\widetilde{\psi}(\mathbf{p},0),\tag{6}$$

for the case  $p(\alpha') = \delta(\alpha' - \alpha)$ . Equation (6) for  $\dot{Y}(|\mathbf{p}|) = 0$  formally reminds the Fourier transform of the Lévy distributions, which are characterized by a power law in the asymptotic behavior in the **r** representation. This feature can be related to extending the Brownian trajectories in the path integral formalism to the Lévy ones, which implies fractional differential operators [40].

Now, let us analyze the quantum statistics from the above scenario. To perform this analysis, we used the thermal Green function technique, which may be directly related to the dynamical aspects of the  $\psi(\mathbf{r}, t)$  and allowed us to obtain several physical quantities, such as the density of particles, specific heat, and other thermodynamical quantities. Following Ref. [47], we defined the one-particle Green function as follows:

$$\mathcal{G}(1,1') = \frac{1}{i\hbar} \left\langle \mathbf{T} \Big[ \psi(1)\psi^{\dagger}(1') \Big] \right\rangle, \tag{7}$$

where the thermodynamic averages,  $\langle \cdots \rangle$ , were evaluated by taking the grand canonical ensemble into account, i.e.,

$$\langle \cdots \rangle = \frac{\operatorname{Tr}(\cdots)e^{\beta(H-\mu N)}}{\operatorname{Tr}e^{\beta(H-\mu N)}};$$
(8)

T is the Dyson time-ordering operator and 1, and 1' correspond to the variables  $r_1$ ,  $t_1$  and  $r_{1'}$ ,  $t_{1'}$ , respectively. From this equation, we can define the correlation functions:

$$\mathcal{G}_{>}(1,1') = \frac{1}{i\hbar} \left\langle \psi(1)\psi^{\dagger}(1') \right\rangle \quad \text{and} \quad \mathcal{G}_{<}(1,1') = \pm \frac{1}{i\hbar} \left\langle \psi^{\dagger}(1')\psi(1) \right\rangle, \tag{9}$$

where > and < represent the Green function to  $t_1 > t_{1'}$ ,  $\mathcal{G} = \mathcal{G}_>$  and  $t_1 < t_{1'}$ ,  $\mathcal{G} = \mathcal{G}_<$ . The upper (lower) sign corresponds to the bosonic (fermionic) case. From Equations (7) and (9), it is possible to show that  $\mathcal{G}_<(1,1')|_{t_1=0} = \pm e^{\beta\mu}\mathcal{G}_>(1,1')|_{t_1=-i\beta\hbar}$  by using the cyclic invariance of the trace  $(\operatorname{Tr}(\hat{A}\hat{B}) = \operatorname{Tr}(\hat{B}\hat{A}))$ . This result shows that the above Green function satisfies the same periodic boundary condition of the non-fractional case [47], in contrast to the

one [46,50,51] formulated within the Tsallis formalism [52]. Similarly to what is performed in the usual case, we may introduce the spectral function,  $A(\mathbf{p}, \omega)$ , defined as follows:

$$\widehat{\widetilde{\mathcal{A}}}(\mathbf{p},\omega) = \widehat{\widetilde{\mathcal{G}}}_{>}(\mathbf{p},\omega) \mp \widehat{\widetilde{\mathcal{G}}}_{<}(\mathbf{p},\omega),$$
(10)

with

$$\widehat{\widetilde{\mathcal{G}}}_{>}(\mathbf{p},\omega) = \int d\mathbf{r}_1 \int_{-\infty}^{\infty} dt e^{-i\mathbf{p}\cdot\mathbf{r}} e^{i\omega t} \mathcal{G}_{>}(\mathbf{r},t)$$
(11)

and

$$\widehat{\widetilde{\mathcal{G}}}_{<}(\mathbf{p},\omega) = \pm i \int d\mathbf{r}_1 \int_{-\infty}^{\infty} dt e^{-i\mathbf{p}\cdot\mathbf{r}} e^{i\omega t} \mathcal{G}_{<}(\mathbf{r},t).$$
(12)

The form of the previous equations for the Green thermal functions results from the translational and rotational invariance of the Hamiltonian considered here for the absence of external potential. By using Equation (9) and the boundary condition obtained above, we can express  $\mathcal{G}^{<}$  and  $\mathcal{G}^{>}$  as follows:

$$\widehat{\widetilde{\mathcal{G}}}_{>}(\mathbf{p},\omega) = (1 \pm f(\omega,\beta,\mu))\widehat{\widetilde{\mathcal{A}}}(\mathbf{p},\omega) \quad \text{and} \quad \widehat{\widetilde{\mathcal{G}}}_{<}(\mathbf{p},\omega) = f(\omega,\beta,\mu)\widehat{\widetilde{\mathcal{A}}}(\mathbf{p},\omega), \tag{13}$$

i.e., in terms of the spectral function, with  $f(\omega, \beta, \mu) = 1/(e^{\beta(\omega-\mu)} \pm 1)$ . It is worth mentioning that  $f(\omega, \beta, \mu)$  is directly connected with the Fermi–Dirac or Bose–Einstein distribution. In this manner, the dynamic aspects are connected with the spectral function, which will give the dispersion of the energy related to the interaction term. The equation of motion satisfied by the Green function defined in Equation (7) is given by

$$i\hbar \frac{\partial}{\partial t} \mathcal{G}(1,1') = \delta(1-1') + \int_{0}^{1} d\alpha' p(\alpha') \frac{1}{2m_{\alpha'}} \left(-\hbar^{2} \nabla_{\mathbf{r}_{1}}^{2}\right)^{\alpha'} \mathcal{G}(1,1') \\ + \int d\mathbf{r}_{\bar{1}} Y(|\mathbf{r}_{1}-\mathbf{r}_{\bar{1}}|) \mathcal{G}(\bar{1},1') + \int d\mathbf{r}_{2} V(|\mathbf{r}_{1}-\mathbf{r}_{2}|) \mathcal{G}(1,2;1',2')|_{t_{2}=t_{1}}.$$
 (14)

It can be obtained by using the Heisenberg equation, where  $\mathcal{G}(1, 2; 1', 2')$  is given by

$$\mathcal{G}(12,1'2') = \frac{1}{\left(i\hbar\right)^2} \left\langle \mathbf{T}\left(\psi(1)\psi(2)\psi^{\dagger}(2')\psi^{\dagger}(1')\right) \right\rangle.$$
(15)

To obtain a solution for Equation (14), we may consistently approximate Equation (15) in accordance with the propagator interpretation of  $\mathcal{G}(1, 1')$ . The first approximation to be considered for Equation (15) to obtain information from Equation (14) is the Hartree approximation,  $\mathcal{G}(12, 1'2') = \mathcal{G}(1, 1')\mathcal{G}(2, 2')$ . The second is the Hartree–Fock approximation given by  $\mathcal{G}(12, 1'2') = \mathcal{G}(1, 1')\mathcal{G}(2, 2') \pm \mathcal{G}(1, 2')\mathcal{G}(2, 1')$ . Other approximations consider the collision process connected to the interaction potential. These approximations have a direct influence on the behavior of the spectral function. The two first cases are formally given in terms of the delta Dirac function, i.e., the energy is localized. For the other cases characterized by collision terms, the behavior is not localized and can be related to a power-law distribution.

By using the previous results, we may relate  $\tilde{\mathcal{G}}_{<}(\mathbf{p}, \omega)$  with the average of the particle density with momentum  $\mathbf{p}$  and energy  $\omega$ , i.e.,

$$\widehat{\widetilde{\mathcal{G}}}_{<}(\mathbf{p},\omega) = \frac{\widetilde{\widetilde{\mathcal{A}}}(\mathbf{p},\omega)}{e^{\beta(\omega-\mu)} \mp 1} = \langle \widehat{\widetilde{n}}(\mathbf{p},\omega) \rangle.$$
(16)

Equation (16) allows us to obtain the pressure for this system by using the equation  $P(\beta, \mu) = \int_{-\infty}^{\mu} d\mu' \langle \hat{\tilde{n}}(\mathbf{p}, \omega) \rangle$ . Another useful result that may be obtained in this framework

concerns the grand canonical potential,  $\Xi$ . By incorporating a coupling constant,  $\lambda$ , in front of the interaction energy, we have

$$\Xi(\beta, V, \mu; \lambda) = \Xi(\beta, V, \mu; 0) - \beta \Omega \int_0^1 \frac{d\lambda}{\lambda} \int \frac{d\mathbf{p}}{(2\pi\hbar)^3} \int \frac{d\omega}{2\pi} \frac{\omega - \widetilde{\Phi}(\mathbf{p})}{2} \widehat{\widetilde{\mathcal{A}}}_{\lambda}(\mathbf{p}, \omega) f(\omega), \quad (17)$$

in the absence of external potential, where

$$\widetilde{\Phi}(\mathbf{p}) = \int_0^1 d\alpha' p(\alpha') \frac{1}{2m_{\alpha'}} |\mathbf{p}|^{2\alpha'} - \widetilde{Y}(|\mathbf{p}|)$$
(18)

and  $\Omega$  is the volume where the system is defined. This result differs from the one presented in Refs. [44,45]. In Equation (18), we have the presence of a term connected with the nonlocal term in the Hamiltonian and the fact that the kinetic energy has a distribution of the moments.

Now, we consider the free particle case in this context in the Hartree approximation with the presence of the nonlocal term. Equation (14) can be written as follows:

$$i\hbar \frac{\partial}{\partial t} \mathcal{G}(1,1') = \delta(1-1') + \int_{0}^{1} d\alpha' p(\alpha') \frac{1}{2m_{\alpha'}} \left(-\hbar^{2} \nabla_{\mathbf{r}_{1}}^{2}\right)^{\alpha} \mathcal{G}(1,1') \\ + \int d\mathbf{r}_{\bar{1}} Y(|\mathbf{r}_{1}-\mathbf{r}_{\bar{1}}|) \mathcal{G}(\bar{1},1') + \int d\mathbf{r}_{2} V(|\mathbf{r}_{1}-\mathbf{r}_{2}|) \mathcal{G}(2,2) \mathcal{G}(1,1').$$
(19)

We can also simplify Equation (19) by approximating the interaction term with a constant, i.e.,

$$\int d\mathbf{r}_2 V(|\mathbf{r}_1 - \mathbf{r}_2|) \mathcal{G}(2, 2) = \int d\mathbf{r}_2 V(|\mathbf{r}_1 - \mathbf{r}_2|) \langle n(\mathbf{r}_2) \rangle = v \langle n \rangle,$$
(20)

for a translationally invariant system with  $\int d\mathbf{r} V(|\mathbf{r}|) = v$  and  $\langle n \rangle = \langle N \rangle / \Omega$ . By using the previous result and the equation

$$\mathcal{G}(1,1') = \frac{1}{-i\beta\hbar} \int \mathbf{r}_1 e^{-i\mathbf{p}(\mathbf{r}_1 - \mathbf{r}_{1'})} \sum_{\nu} e^{-iz_{\nu}(t_1 - t_{1'})} \widetilde{\overline{\mathcal{G}}}(\mathbf{p}, z_{\nu}), \tag{21}$$

where  $z_{\nu} = i[\pi\nu/(-i\beta\hbar) + \mu]$ , it is possible to obtain from Equation (19) the following result:

$$\overline{\overline{\mathcal{G}}}(\mathbf{p}, z_{\nu}) = 1/[z_{\nu} - \tilde{\Phi}(|\mathbf{p}|) - v\langle n \rangle],$$
(22)

with the spectral function in this approximation being given by

$$\widehat{\widetilde{\mathcal{A}}}(\mathbf{p},\omega) = 2\pi\delta\big(\omega - \tilde{\Phi}(|\mathbf{p}|) - v\langle n \rangle\big).$$
(23)

Equation (23) allows us to obtain the number of particles in terms of the chemical potential  $\mu$ , or vice versa, and the internal energy as follows:

$$\langle N \rangle = \Omega \int \frac{d\mathbf{p}}{(2\pi\hbar)^3} \frac{d\omega}{2\pi} \widehat{\widetilde{\mathcal{A}}}(\mathbf{p},\omega) f(\omega)$$

$$= \Omega \int \frac{d\mathbf{p}}{(2\pi\hbar)^3} \frac{1}{e^{\beta \left[\widetilde{\Phi}(|\mathbf{p}|) - v\langle n \rangle - \mu\right]} \mp 1}$$
(24)

and

These equations allow us to show that, in the limit of low density ( $\beta \mu \rightarrow -\infty$ ), the pressure can be found. It is given by  $P - (1/2)\langle n \rangle^2 = \langle n \rangle k_B T$ , which is essentially the van der Waals equation without the volume exclusion effect. This feature suggests that the presence of a fractional derivative in Equation (2) does not change the thermodynamic relations, as expected. The changes are expected in the quantities directly connected with the effect of the dynamics, such as in the low-temperature limit, in which the dynamic effects are pronounced. At this point, we can mention the developments performed in Refs. [37–39], where the thermodynamics relations were extended by incorporating fractional differential operators.

Figure 1 shows the behavior of the internal energy and the specific heat in the low-temperature limit for the Bose–Einstein distribution in the case  $\mu = 0$ . It shows that the fractional term present in the Hamiltonian has a pronounced effect on the low-temperature limit. This point is evidence that energy fluctuations are governed by the fractional terms in this limit when present in the system. It is worth mentioning that several systems present the behavior  $C \sim T^{\alpha}$  for the specific heat in this limit, such as two-dimensional Kagome antiferromagnet,  $Cu_3Zn(OH)_6FBr$  [53], and amorphous and crystalline particles at low temperatures [45]. In the limit of high temperatures, we recovered the standard behavior for the specific heat as shown in Figure 1.



**Figure 1.** Behavior of the internal energy (see (a)) and the specific heat (see (b)) for the Bose–Einstein distribution by taking  $\Phi(|\mathbf{p}|) = \mathbf{p}^2/(2m) + |\mathbf{p}|^{2\alpha}/(2m_{\alpha})$  into account, where  $C_{\alpha} = \xi C'_{\alpha}$  and  $T'' = \xi T$ . This choice combines the usual and fractional kinetic energies and shows that the fractional one has a pronounced effect on the asymptotic limit of low temperature. We considered, for simplicity,  $m_{\alpha} = (2m)^{\alpha/2}/2$  and  $\xi = \Omega(2m\pi k_B/h^2)^{3/2}$ .

Another case that has an unusual behavior in low temperatures is liquid Helium [54] as shown in Figure 2. The behavior in this limit for the specific heat is  $C \sim T^6$  as discussed in Ref. [54], which is the one obtained with the formalism presented in this section for the Bose–Einstein distribution. From a phenomenological point of view, the fractional approach captures the complex behavior of the interactions present in this system, which deserves a careful analysis.



**Figure 2.** This figure shows the behavior of the model for an ideal Bose–Einstein distribution with  $\Phi(|\mathbf{p}|) = |\mathbf{p}|^{2\alpha}/(2m_{\alpha})$  and the experimental data obtained from Ref. [55] for the specific heat of liquid Helium. The following relations are valid: C' = C/N and  $T' = T/T_c$ , similar to what was performed in Ref. [55].

# 2.2. Tsallis Approach and Thermal Green Functions

Let us start our discussion about quantum statistics in the context of Tsallis formalism. It will review and follow the developments performed in Ref. [50] in terms of the thermal Green function formalism.

The main difference to previous development concerns the statistical approach used, which is based on the Tsallis entropy [15], i.e.,

$$S_q = k_B \frac{1 - \mathrm{Tr}\rho^q}{q - 1},\tag{26}$$

instead of the usual entropy. Applying the maximum principle of entropy, we can obtain the density matrix by taking into account the normalized constraints [14],

$$\langle H \rangle_q = U_q = \operatorname{Tr} H \rho^q / \operatorname{Tr} \rho^q \quad , \quad \langle N \rangle_q = N_q = \operatorname{Tr} N \rho^q / \operatorname{Tr} \rho^q ,$$

$$\tag{27}$$

and the usual normalization of the density matrix  $\text{Tr } \rho = 1$ . It is given by

$$\rho = \exp_q(-X(\tilde{\beta}, \mu))/Z_q \quad \text{and} \quad Z_q = \operatorname{Tr}\left[\exp_q(-X(\tilde{\beta}, \mu))\right], \tag{28}$$

where

$$\exp_q(-X(\tilde{\beta},\mu)) \equiv (1-(1-q)X(\tilde{\beta},\mu))^{1/(1-q)}$$
 (29)

and  $X(\tilde{\beta}, \mu) = \tilde{\beta}((H - U_q) - \mu(N - N_q))$  with  $\tilde{\beta} = \beta / \operatorname{Tr} \tilde{\rho}^q$ . Here,  $\tilde{\beta}$  is the normalized temperature and  $\beta$  and  $\mu$  are the Lagrange multipliers associated with the normalized *q*-expectation values. Two self-consistency conditions follow from Equation (29): Tr  $\tilde{\rho}^q = \tilde{Z}_q^{1-q}$  and  $Z_q = \text{Tr}[e_q(-X(\tilde{\beta}, \mu))]^q$ . By using the previous results, the expression for a general normalized *q*-expectation value is then

$$\langle A \rangle_{q} = \operatorname{Tr} \{ A \rho^{q} \} / \operatorname{Tr} \rho^{q}$$
  
= 
$$\operatorname{Tr} \{ A \left[ \exp_{q} (-X(\tilde{\beta}, \mu)) \right]^{q} \} / \operatorname{Tr} \left[ \exp_{q} (-X(\tilde{\beta}, \mu)) \right]^{q}.$$
 (30)

From Equation (30),  $\langle H \rangle_q$ ,  $\langle N \rangle_q$  can be deduced. Equation (30) can also be written as follows:

$$\langle A \rangle_q = \operatorname{Tr} \left\{ A \left[ \exp_q(-X'(\beta',\mu)) \right]^q \right\} / \operatorname{Tr} \left[ \exp_q(-X'(\beta',\mu)) \right]^q, \tag{31}$$

where

$$X'(\beta',\mu) = \beta'(H-\mu N)$$
 and (32)

$$\beta' = \frac{\beta}{1 + (1 - q)\tilde{\beta}(U_q - \mu N_q)}.$$
(33)

In the previous equations of this section, we considered the Hamiltonian given by Equation (2) with  $p(\alpha') = \delta(\alpha' - 1)$  and  $Y(|\mathbf{r}|, t) = 0$ , which corresponds to the standard form of the Hamiltonian with an interaction term. In this approach to statistical mechanics, we revisited some of the developments performed in the previous section. The one-particle Green function is then with the normalized *q*-average as follows [50]:

$$\mathcal{G}^{(q)}(1,1') = \frac{1}{i\hbar} \left\langle \mathrm{T}\left(\psi(1)\psi^{\dagger}(1')\right) \right\rangle_{q}.$$
(34)

The previous equation recovers the standard case for q = 1, i.e.,  $\mathcal{G}^{(q)}(1, 1') \rightarrow \mathcal{G}^{(1)}(1, 1') = \mathcal{G}(1, 1')$ . By using the integral contour representation employed in Refs. [46,51],

$$\frac{i}{2\pi} \int_{C} du \exp(-ub)(-u)^{-z} = \frac{b^{1-z}}{\Gamma(z)},$$
(35)

with b > 0 and Re z > 0, where the contour *C* starts from  $+\infty$  on the real axis, encircles the origin once counterclockwise and returns to  $+\infty$ , Equation (34) is expressed in terms of the Green function with q = 1:

$$\mathcal{G}^{(q)}(1,1') = \int_{C} du K_{q}(u) Z_{1}(-u(1-q)\beta',\mu) \mathcal{G}^{(1)}(1,1';-u(1-q)\beta',\mu),$$
(36)

where

$$K_q(u) = \frac{i}{2\pi Z'_q} \Gamma\left(\frac{1}{1-q}\right) \exp(-u)(-u)^{-1/(1-q)},$$
(37)

with  $Z'_q = \text{Tr}\left[\exp_q(-X'(\beta', \mu))\right]^q$ . One particle's standard result for the Green function is recovered for q = 1. As before, we can introduce correlation functions [46,50]

$$\mathcal{G}_{>}^{(q)}(1,1') = \frac{1}{i\hbar} \left\langle \psi(1)\psi^{\dagger}(1') \right\rangle_{q} \quad \text{and} \quad \mathcal{G}_{<}^{(q)}(1,1') = \pm \frac{1}{i\hbar} \left\langle \psi^{\dagger}(1')\psi(1) \right\rangle_{q}.$$
(38)

The spectral weight function in frequency space by taking the Fourier transform concerning time differences,  $\widehat{\mathcal{A}}(\mathbf{r}_1, \mathbf{r}_{1'}; \omega)$ , can be introduced as before. We assumed a translational invariance in time at this stage. We can thus express  $\mathcal{G}_{>}^{(q)}$  and  $\mathcal{G}_{<}^{(q)}$  in terms of these in the following way [46,50]:

$$i\mathcal{G}_{>}^{(q)}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega) = \int_{C} du\tilde{K}_{q}(u)Z_{1}(-u(1-q)\beta',\mu)i\mathcal{G}_{>}^{(1)}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega;-u(1-q)\beta')$$
  
$$= \int_{C} du\tilde{K}_{q}(u)(1\pm f(\omega,-u(1-q)\beta',\mu))\hat{\mathcal{A}}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega)Z_{1}(-u(1-q)\beta',\mu)$$
(39)

and

$$i\mathcal{G}_{<}^{(q)}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega) = \int_{C} du \tilde{K}_{q}(u) \tilde{Z}_{1}(-u(1-q)\beta',\mu) i\mathcal{G}_{<}^{(1)}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega;-u(1-q)\beta',\mu)$$
  
$$= \pm \int_{C} du \tilde{K}_{q}(u) f(\omega,-u(1-q)\beta',\mu) \widehat{\mathcal{A}}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega) \tilde{Z}_{1}(-u(1-q)\beta',\mu).$$
(40)

The spectral function weight is a given system property not dependent on the ensemble as in the previous case. The results appear to be similar to those in the previous section with the modifications noted here. Using the Fourier representations of the step functions involved in the time-ordered Green function in Equation (36), an important result can be deduced [50]:

$$\mathcal{G}^{(q)}(1,1') = \int_{C} du K_{q}(u) Z_{1}(-u(1-q)\beta',\mu) \int_{-\infty}^{\infty} \frac{d\omega}{2\pi} e^{-i\omega(t_{1}-t_{1'})} \int_{-\infty}^{\infty} \frac{d\omega'}{2\pi} \widehat{\mathcal{A}}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega') \\ \times \left\{ P\left(\frac{1}{\omega-\omega'}\right) - i\pi\delta(\omega-\omega') \left[ \tanh\left(\frac{-u}{2}\beta'(1-q)(\omega-\mu)\right) \right]^{\mp 1} \right\}.$$

$$\tag{41}$$

For a uniform system, we can take Fourier transforms concerning  $r_1 - r_{1'}$  as in the previous section and express the one-particle momentum distribution function  $\langle \hat{\tilde{n}}(\mathbf{p}) \rangle_q$  in terms of the spectral weight function of the *N*-particle system as follows:

$$\langle \widehat{\widetilde{n}}(\mathbf{p}) \rangle_{q} = \int_{C} du K_{q}(u) \int_{-\infty}^{\infty} \frac{d\omega}{2\pi} \frac{Z_{1}(-u(1-q)\beta',\mu)}{e^{-u(1-q)\beta'(\omega-\mu)} \mp 1} \widehat{\widetilde{\mathcal{A}}}(\mathbf{p};\omega).$$
(42)

The chemical potential is determined by the following expression:

$$\langle N \rangle_{q} = \Omega \int_{C} du K_{q}(u) \int_{-\infty}^{\infty} \frac{d\omega}{2\pi} \int \frac{d\mathbf{p}}{(2\pi\hbar)^{3}} \frac{Z_{1}(-u(1-q)\beta',\mu)}{e^{-u(1-q)\beta'(\omega-\mu)} \mp 1} \widehat{\widetilde{\mathcal{A}}}(\mathbf{p};\omega).$$
(43)

It is worth mentioning that the spectral function in this context is determined by the thermal Green function, which emerges from the following equation:

$$i\hbar \frac{\partial}{\partial t} \mathcal{G}^{(q)}(1,1') + \frac{\hbar^2}{2m} \nabla^2_{\mathbf{r}_1} \mathcal{G}^{(q)}(1,1') = \delta(1-1') \\ + \int d\mathbf{r}_2 V(|\mathbf{r}_1 - \mathbf{r}_2|) \mathcal{G}^{(q)}(1,2;1',2') \Big|_{t_2 = t_1'}$$
(44)

which is similar to the one presented before. However, the simple form of the boundary condition  $\mathcal{G}^{<}(1,1')|_{t_1=0} = \pm e^{\beta\mu}\mathcal{G}^{>}(1,1')|_{t_1=-i\beta\hbar}$  cannot be applied due to the structure of the statistical distribution, which is a power law instead of an exponential. For the free particle case, it is possible to show that  $\widehat{\mathcal{A}}(\mathbf{p};\omega) = 2\pi\delta(\omega - \mathbf{p}^2/(2m))$  as in the previous case. This result for the free particle case implies that the internal energy is given by

$$\langle H \rangle_{q} = \Omega \int_{C} du K_{q}(u) \int \frac{d\mathbf{p}}{(2\pi\hbar)^{3}} \frac{|\mathbf{p}|^{2}}{2m} \frac{Z_{1}(-u(1-q)\beta',\mu)}{e^{-u(1-q)\beta'(\frac{|\mathbf{p}|^{2}}{2m}-\mu)} \mp 1}.$$
(45)

It is also possible to consider the Hartree approximation for Equation (44), with similar results for the spectral function. In fact, for the Hartree approximation, i.e.,  $\mathcal{G}^{(q)}(1,2;1',2') = \mathcal{G}^{(q)}(1,1')\mathcal{G}^{(q)}(2,2')$ , it is possible to show that the dynamics equation for the Green function is given by:

$$i\hbar \frac{\partial}{\partial t} \mathcal{G}^{(q)}(1,1') + \frac{\hbar^2}{2m} \nabla^2_{\mathbf{r}_1} \mathcal{G}^{(q)}(1,1') = \delta(1-1') \\ + \int d\mathbf{r}_2 V(|\mathbf{r}_1 - \mathbf{r}_2|) \mathcal{G}^{(q)}(2,2) \mathcal{G}^{(q)}(1,1').$$
(46)

Equation (46) may be simplified by using similar assumptions to those used for Equation (19) and allow us to obtain  $\widehat{\mathcal{A}}(\mathbf{p};\omega) = 2\pi\delta(\omega - |\mathbf{p}|^2/(2m) - \langle n \rangle_q v)$ , where  $\langle n \rangle_q = \langle N \rangle_q / \Omega$ . Note that considering collision terms implies using boundary conditions that will lead us to cumbersome calculations [56,57]. Thus, obtaining the corrections due to the collision terms is harder than the previous one in this context. Figure 3 illustrates the behavior of the specific heat for the formalism presented in this section. We observe that the behavior in the low and high temperatures is similar to that in the standard case. The difference is for intermediate temperatures, i.e., temperatures nearly T'' = 1, for the values of *q* considered in this figure.



**Figure 3.** Behavior of the internal energy (see (**a**)) and the specific heat (see (**b**)) for the Bose–Einstein distribution by taking the free particle case into account, where  $C_q = \xi' C'_q$  and  $T'' = \xi'^{2/3} T_q$ . We consider, for simplicity,  $\xi' = \left[\Omega(2m)^{3/2}/(2\pi\hbar^3)\right]^{2/3} k_B$ .

# 2.3. Fractional-Generalized Entropy and Thermal Green Functions

Now, we consider the fractional approach in the context of thermostatistics obtained from a generalized entropy, which can be related to the Tsallis entropy. For this, we considered the entropy [58],  $S_q = S_q(\text{Tr}\rho^q)$ , which can be related to the Tsallis and Renyi

entropies. Applying the maximum entropy principle, we can obtain the density matrix by considering the normalized constraints defined in the previous section, respectively. It is given by:

$$\rho = \exp_q(-X_{\mathcal{F}_q}(\beta \mathcal{F}_q(\operatorname{Tr}\rho^q), \mu))/Z_q \quad \text{and} \quad Z_q = \operatorname{Tr}\left[\exp_q(-X_{\mathcal{F}_q}(\beta \mathcal{F}_q(\operatorname{Tr}\rho^q), \mu))\right], \quad (47)$$

where  $X_{\mathcal{F}_q}(\beta \mathcal{F}_q(\mathrm{Tr}\rho^q),\mu) = \beta \mathcal{F}_q(\mathrm{Tr}\rho^q)((H-U_q)-\mu(N-N_q))$ , with

$$\mathcal{F}_{q}(\mathrm{Tr}\rho^{q}) = \frac{1}{(1-q)\mathrm{Tr}\rho^{q}} \frac{\partial \mathrm{Tr}\rho^{q}}{\partial S_{q}}.$$
(48)

In this context, the normalized temperature is represented by  $\beta_{\mathcal{F}_q} = \beta \mathcal{F}_q(\text{Tr}\rho^q)$ , instead of  $\tilde{\beta}$ , presented in the previous section. The parameters  $\beta$  and  $\mu$  are the Lagrange multipliers associated with the normalized *q*-expectation values as before. Similar to the previous section, we may consider the following form for the *q*-expectation values:

$$\langle A \rangle_q = \operatorname{Tr} \left\{ A \left[ \exp_q(-X'_{\mathcal{F}_q}(\beta \mathcal{F}_q(\operatorname{Tr}\rho^q), \mu) \right]^q \right\} / \operatorname{Tr} \left[ \exp_q(-X'_{\mathcal{F}_q}(\beta \mathcal{F}_q(\operatorname{Tr}\rho^q), \mu) \right]^q, \quad (49)$$

where

$$X'_{\mathcal{F}_q}(\beta \mathcal{F}_q(\mathrm{Tr}\rho^q),\mu) = \beta'_{\mathcal{F}_q}(H-\mu N)$$
(50)

and

$$\beta_{\mathcal{F}_q}' = \frac{\beta \mathcal{F}_q(\operatorname{Tr} \rho^q)}{1 + (1 - q)\beta \mathcal{F}_q(\operatorname{Tr} \rho^q) (U_q - \mu N_q)}.$$
(51)

In this approach to statistical mechanics, the one-particle Green function is given by:

$$\mathcal{G}^{(q)}(1,1') = \frac{1}{i\hbar} \left\langle \mathrm{T}\left(\psi(1)\psi^{\dagger}(1')\right) \right\rangle_{q'},\tag{52}$$

which is formally equal to the one presented in the previous section. In fact, Equation (52) has the same formal aspect of Equation (34). However, the *q*-average is given by Equation (49), obtained from the distribution given by Equation (47), which has a different dependence on the Lagrange parameter  $\beta$ . We may also use the integral contour representation employed in the previous section to obtain

$$\mathcal{G}^{(q)}(1,1') = \int_{C} du K_{q}(u) Z_{1}(-u(1-q)\beta'_{\mathcal{F}_{q}},\mu) \mathcal{G}^{(1)}(1,1';-u(1-q)\beta'_{\mathcal{F}_{q}},\mu),$$
(53)

where, in this case,  $Z'_q = \text{Tr}\left[\exp_q(-X'(\beta'_{\mathcal{F}_q},\mu))\right]^q$ . In terms of the definitions presented in this section for the q- expectation values and the thermal Green function, we can express  $\mathcal{G}^{(q)}_{>}$  and  $\mathcal{G}^{(q)}_{<}$  as follows:

$$i\mathcal{G}_{>}^{(q)}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega) = \int_{C} du \tilde{K}_{q}(u) Z_{1}(-u(1-q)\beta'_{\mathcal{F}_{q}},\mu) i\mathcal{G}_{>}^{(1)}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega;-u(1-q)\beta'_{\mathcal{F}_{q}})$$
(54)

and

$$i\mathcal{G}_{<}^{(q)}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega) = \int_{C} du \tilde{K}_{q}(u) \tilde{Z}_{1}(-u(1-q)\beta'_{\mathcal{F}_{q}},\mu) i\mathcal{G}_{<}^{(1)}(\mathbf{r}_{1},\mathbf{r}_{1'};\omega;-u(1-q)\beta'_{\mathcal{F}_{q}},\mu).$$
(55)

These results are formally equal to the previous ones for the Tsallis case and can also be expressed in terms of the spectral function  $\hat{A}(\mathbf{r}_1, \mathbf{r}_{1'}; \omega)$ . The spectral function in this context is determined by the thermal Green functions, which emerge from the following equation:

$$i\hbar\frac{\partial}{\partial t}\mathcal{G}^{(q)}(1,1') = \delta(1-1') + \int_{0}^{1} d\alpha' p(\alpha') \frac{1}{2m_{\alpha'}} \left(-\hbar^{2}\nabla_{\mathbf{r}_{1}}^{2}\right)^{\alpha'} \mathcal{G}^{(q)}(1,1') \\ + \int d\mathbf{r}_{\bar{1}} Y(|\mathbf{r}_{1}-\mathbf{r}_{\bar{1}}|) \mathcal{G}^{(q)}(\bar{1},1') + \int d\mathbf{r}_{2} V(|\mathbf{r}_{1}-\mathbf{r}_{2}|) \mathcal{G}^{(q)}(1,2;1',2')\Big|_{t_{2}=t_{1}},$$
(56)

which has a fractional spatial derivative of distributed order in the kinetic term and the presence of a nonlocal term. For the free particle case, it is possible to show that

$$\widehat{\widetilde{\mathcal{A}}}(\mathbf{p};\omega) = 2\pi\delta(\omega - \widetilde{\Phi}(|\mathbf{p}|)).$$
(57)

This result allows us to show that the internal energy for this case is given by

$$\langle H \rangle_{q} = \Omega \int_{C} du K_{q}(u) \int \frac{d\mathbf{p}}{(2\pi\hbar)^{3}} \widetilde{\Phi}(|\mathbf{p}|) \frac{Z_{1}(-u(1-q)\beta'_{\mathcal{F}_{q}},\mu)}{e^{-u(1-q)\beta'_{\mathcal{F}_{q}}(\widetilde{\Phi}(|\mathbf{p}|)-\mu)} \mp 1},$$
(58)

and the q-mean value for the number of particles is

$$\langle N \rangle_q = \Omega \int_C \, \mathrm{d} u K_q(u) \int \frac{\mathrm{d} \mathbf{p}}{(2\pi\hbar)^3} \frac{Z_1(-u(1-q)\beta'_{\mathcal{F}_q},\mu)}{\mathrm{e}^{-u(1-q)\beta'_{\mathcal{F}_q}(\tilde{\Phi}(|\mathbf{p}|)-\mu)} \mp 1}.$$
(59)

Figure 4 shows the trend of the specific heat for the previous cases and the mixing of the two cases obtained with the approach presented in this section. It is possible to verify that the fractional kinetic term, i.e.,  $\alpha \neq 0$ , has a pronounced effect on the limit of small and high temperatures. We observe the effect of  $q \neq 1$  for intermediate temperature as in Figure 3b.



**Figure 4.** Trend of the specific heat for the Bose–Einstein distribution by taking the free particle case into account with  $\mu = 0$  and  $\Phi(\mathbf{p}) = |\mathbf{p}|^{2\alpha}/(2m_{\alpha})$ , where  $C_{\alpha,q} = \xi'_{\alpha}C'_{\alpha,q}$  and  $T'' = \xi'_{\alpha}T_q$ . We consider, for simplicity,  $\xi'_{\alpha} = \left[\Omega(2m_{\alpha})^{3/(2\alpha)}/(2\pi\hbar^3)\right]^{2\alpha/3}k_B$ .

## 3. Discussions and Conclusions

We investigated two different extensions of the quantum statistical approach. The first one considers an effective Hamiltonian in terms of the fractional derivative in the space of distributed order with a nonlocal term. The second one considers the extension of the concept of entropy to a non-additive case, allowing us to obtain power-law distributions for the statistical weight instead of the Boltzmann–Gibbs distributions characterized by exponential behavior. These approaches were investigated using the thermal Green functions, which allowed us to work in a unified way with the Bose-Einstein and Fermi-Dirac distributions and establish some differences between the formalisms. In this case, we obtained the behavior with the temperature for internal energy and the specific heat for both extensions of the quantum statistics. The behavior for each case was illustrated in Figures 1 and 3. We observed for the first case that, depending on the choice of the fractional parameter in the distributed fractional space derivative, the behavior in the low-temperature limit is different from usual when  $\Phi(\mathbf{p}) = \mathbf{p}^2/(2m) + |\mathbf{p}|^{2\alpha}/(2m_{\alpha})$ . In the limit of high temperatures, the usual behavior is recovered. This feature reflects the mixing of the usual and the fractional differential operators connected to the different forms of the kinetic term. In particular, we considered this feature to fit the specific heat of He<sup>4</sup> from the phenomenological point of view (see Figure 2). For the case worked out in Section 2.2, we observed that the behavior for intermediate temperatures is different from usual for the free particle case. In the other limits, we see the same behavior as the temperature changes. These results have shown that different approaches present different behaviors. Choosing these or other extensions, such as the one presented in Section 2.3, which extends the results presented in Sections 2.1 and 2.2 (see, Figure 4), requires a careful system analysis to choose the suitable approach. We hope that the results presented here can be helpful in the discussion of systems with nonstandard behavior and statistical approaches.

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