Chaotic scattering on surfaces and collisional damping of collective modes

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The damping of hot giant dipole resonances is investigated. The contribution of surface scattering is compared with the contribution from interparticle collisions. A unified response function is presented which includes surface damping as well as collisional damping. The surface damping enters the response via the Lyapunov exponent and the collisional damping via the relaxation time. The former is calculated for different shape deformations of quadrupole and octupole type. The surface as well as the collisional contribution each reproduce almost the experimental value, therefore we propose a proper weighting between both contributions related to their relative occurrence due to collision frequencies between particles and of particles with the surface. We find that for low and high temperatures the collisional contribution dominates whereas the surface damping is dominant around the temperatures $\sqrt{3/2\pi}$ of the centroid energy.

I. INTRODUCTION

The damping mechanisms of collective motions in excited nuclei are a topic of continuing debate [1]. Mainly two lines of thought are pursued. In one line of thought it is assumed that collisions are the physical reason for damping only as developed via a Fermi liquid approach with bulk matter properties [2–15]. The other line of thought considers new features of the finite nucleus, such as surface oscillations and a level density with finite spacing. Partially the investigations are performed without inertia [16–24] or by including inertia [22,25–30]; note that inertia is absent in infinite matter.

Both classes of models predict a comparable degree of damping necessary to reproduce the experimental data. Consequently, it is an open question which is the correct physical reason for damping. Of course, the correct description has to assume a finite nucleus consisting of nucleons which are bound via the mean field, through which the nucleons undergo mutual collisions and where the surface is formed by the particles themselves. These features are usually included in Boltzmann-Uehling-Uhlenbeck-(BUU) simulations [31–33] or in its nonlocal extensions [34,35]. In full simulations, however, we will not gain a simple insight into the physical origins of the damping mechanism, in particular, how much is due to surface contributions and how much is due to collisional contributions.

The aim of this article is to compare both pictures in the frame of linear response theory. Within the collision-free Vlasov equation the linear response of finite systems is well known [36] and allows one to calculate the strength function of finite nuclei. The damping, however, does not reproduce the experimental damping of giant resonances since collisions are absent.

To include damping, we can take into account on the one hand the collisional damping in infinite matter and find a scaling for finite size effects in the sense of a Thomas-Fermi local density approximation. On the other hand we can consider the boundary of the finite nucleus as a fixed surface such that the nucleons simply bounce off this wall like in a billiard-type situation. Provided we accept this simple picture, it is possible to compare the damping caused by this chaotic scattering off the wall with the damping from collisional contributions in infinite matter. We will derive a response function which includes the additional chaotic process and find a total damping rate $\Gamma = \Gamma_{\text{coll}} + \Gamma_{\text{surf}}$ similar to the Matthiessen rule in metals [37] but with an additional weight between the different processes according to the ratio of the corresponding collision frequencies. Moreover the hypotheses by Swiatecki [38] is questioned that negative curved surfaces of octupole deformations would induce an additional chaotic mechanism of dissipation and consequently octupole modes should be overdamped. We will find that the octupole deformation does not lead to any special enhancement of the Lyapunov exponent in comparison with the quadrupole mode. Moreover, any deformation will cause a contribution to the damping of collective modes.

The outline of the paper is as follows. In the next chapter II the largest Lyapunov exponent for a nucleon in a deformed nucleus is calculated. Then in chapter III we derive a unified response function which combines both the collisional damping and the contribution from surface scattering. In chapter IV both damping contributions are compared with the experimental values of hot isovector giant dipole resonances (IVGDR). A proper weighting
factor corresponding to the relative collision frequencies yields a unified picture which describes the data rather well. We find that the collisional contributions dominate for low temperatures and for very high temperatures while around the highest experimental achievable temperatures the surface effects are important. Chapter V will summarize the results.

II. LARGEST LYAPUNOV EXPONENT OF DEFORMED NUCLEI

We will consider in the following only classical 3-D closed billiards and use the largest Lyapunov exponent as a relevant measure to characterize chaoticity. The surface of these billiards is chosen to resemble the surface deformation of a nucleus undergoing quadrupole or octupole deformations.

The Lyapunov exponent can be given as the deviations of the difference of the trajectories

\[ |\vec{r}_1(t) - \vec{r}_2(t)| = |\vec{r}_1(0) - \vec{r}_2(0)| e^{\lambda t} \]

as

\[ \lambda = \lim_{t \to \infty} \lim_{\epsilon \to 0} \frac{1}{t} \ln \left( \frac{|\vec{r}_1(t) - \vec{r}_2(t)|}{|\vec{r}_1(0) - \vec{r}_2(0)|} \right). \]

For \( \lambda > 0 \) the difference in phase space trajectories grows exponentially leading to chaotic behavior.

We solve the Hamilton equations for one particle in finite nuclear matter with an infinite potential

\[ V(r, \theta) = V_0 \Theta (r - R(\theta)) \]

modeling the deformation of a axial symmetric nuclear surface

\[ R_\lambda(\theta) = R_0 \left( 1 + \alpha_{00} + \alpha_\lambda P_\lambda(\cos(\theta)) \right) \]

with the nuclear radius \( R_0 = 1.13A^{1/3} \text{fm} \), and where \( \lambda = 2 \) corresponds to the quadrupole and \( \lambda = 3 \) to the octupole deformation [39]. The coefficient \( \alpha_{00} \) is adjusted to conserve the volume corresponding to incompressible nuclear matter.

\[ V(r, \theta) \approx V_0 \left( \arctan \left( \frac{r - R(\theta)}{\sigma} \right) + \frac{\pi}{2} \right) \]

resembling an infinite step function.

The Lyapunov exponent is calculated by considering the time evolution of small deviations from a reference trajectory due to infinitesimal initial changes. We have used the Brandstaetter method resetting the deviations of the reference trajectory repeatedly after a certain time to the initial infinitesimal difference. This corresponds to an averaging and the largest mean Lyapunov exponent is obtained.

\[ \lambda \text{ [c/fm]} \]

\[ \alpha_{00} = 0.3 \]

\[ \alpha_{00} = 0.15 \]

\[ \alpha_{00} = 0.6 \]

\[ \alpha_{00} = 0.45 \]

FIG. 1. Overview of the geometry of modes

FIG. 2. The largest Lyapunov exponent of a spherical deformed billiard versus deformation parameter \( \alpha \). The regions of \( \alpha \) where the surface starts to become negatively curved \( (k < 0) \) via (5) are indicated by dotted lines \( \alpha_{k=0} \) for octupole and quadrupole deformation, respectively.

Figure 2 shows the Lyapunov exponent for different
octupole and quadrupole deformations according to (4). The error bars are taken from averaging 50 runs of different initial conditions indicating 95% confidence level. We see that the quadrupole deformation leads to an immediate increase of the Lyapunov exponent while the Lyapunov exponent for the octupole deformation increases at larger deformation parameters and proceeds more slowly with increasing deformation. Let us note that the characteristic time scales from the Lyapunov exponent are similar to the damping rate of the IVGDR which will be presented in figure 4. This motivates us to consider a unified response from collisions and surface contributions.

Following [38] a surface with negative curvature should induce a new chaotic mechanism analogous to the Sinai billiard. We discuss therefore the curvature of a rotating body given by the curvature of the boundary curve \( R(\theta) \) of (4) via

\[
k = \frac{R^2 + 2R'^2 - RR''}{(R^2 + R'^2)^{3/2}}
\]

where \( R' \) and \( R'' \) denote the derivatives with respect to \( \theta \). In figure 2 we see that the octupole deformation shows parts with negative curvature at smaller deformation parameter than the quadrupole deformation shows. However, the Lyapunov exponent of octupole deformation remains always smaller. We therefore conclude that the octupole deformation compared with the quadrupole deformation leads to no significant enhancement of chaos/tization.

III. DAMPING OF COLLECTIVE OSCILLATIONS

We would like to focus on a unified description of the linear response including collisional contributions and chaotic scattering with the surfaces. Therefore let us briefly sketch the response function formalism starting from appropriate kinetic equations. For infinite matter this procedure will result in the known Mermin response function which has been used to describe the IVGDR in symmetric [11] and asymmetric nuclear matter [13]. Then we derive the response function for finite nuclei and show that in a local density approximation chaotic scattering from the surface can be incorporated. The result will be a Mermin-like response function in local density approximation where the largest Lyapunov exponent appears as an imaginary shift in the frequency.

### A. Infinite matter response

In order to consider the collisional damping we start from the kinetic equation for the quasiclassical distribution function. For neutrons this kinetic equation reads

\[
f_n(p, r, t) + \frac{p}{m} \partial_t f_n(p, r, t) - \partial_r (V_n + V_{\text{ext}}) \partial_p f(p, r, t) = I
\]

with the self-consistent mean-field potential given by a schematic Skyrme type [40,41]

\[
V_n(r, t) = t_0 \left\{ \left( 1 + \frac{x_0}{2} \right) (n_n(r, t) + n_p(r, t)) - \left( x_0 + \frac{1}{2} \right) n_n(r, t) \right\} + \frac{t_3}{4} \left\{ (n_n(r, t) + n_p(r, t))^2 - n_n^2(r, t) \right\}.
\]

The neutron and proton densities are \( n_n = \int \frac{dp}{(2\pi\hbar)^3} f_n \) and \( n_p = \int \frac{dp}{(2\pi\hbar)^3} f_p \), respectively. The parameters are \( x_0 = 0.48, \ t_0 = -983.4 \text{ MeV fm}^3 \) and \( t_3 = 13106 \text{ MeV fm}^4 \). The kinetic equation and mean field for protons are obtained from the linearization of the Vlasov equation \( \partial_t f + \left( \partial_r V_{\text{ext}} + \frac{p}{m} \right) \partial_p f = 0 \). The neutron and proton densities are

\[
\delta n = \Pi_0(\delta V_{\text{ext}} + V_0 \delta n)
\]

with [41]

\[
V_0 = 2(V_n - V_p) = -\frac{t_0}{2} \left( x_0 + \frac{1}{2} \right) - \frac{t_3}{8} n_0
\]

where \( n_0 = \int \frac{dp}{(2\pi\hbar)^3} f_0 = 0.16 \text{ fm}^{-3} \) is the nuclear saturation density and \( \Pi_0 \) is the Lindhard-function

\[
\Pi_0^{\text{inf}}(q, \omega) = 4 \int \frac{dp'}{(2\pi\hbar)^3} \frac{q \partial_{p'} f_0(p')}{p' - \omega - i0}.
\]

Here we first consider the homogeneous equilibrium \( f_0(p) \) or infinite matter.

The response function \( \Pi \) connects the induced density fluctuation \( \delta n \) with the external potential via

\[
\delta n = \Pi V_{\text{ext}}
\]

and the response function follows from (8) for infinite matter

\[
\Pi^{\text{inf}}(q, \omega) = \frac{\Pi_0^{\text{inf}}(q, \omega)}{1 - V_0 \Pi_0^{\text{inf}}(q, \omega)} = \frac{\Pi_0^{\text{inf}}(q, \omega)}{\epsilon(q, \omega)}.
\]
The zeros $\omega = \Omega + i\gamma$ of the dielectric function $\epsilon(q, \omega) = 0$ in (12) determine the collective modes with the energy $\Omega$ and the Landau damping $\gamma$ of the collective excitation. Within the Steinwedel-Jensen model for IVGDR [42], the wave vector scales like $q = \frac{2\pi}{R_0}$ where the nuclear radius is $R_0 = 1.13A^{1/3}$fm.

2. Collisional model

In order to take collisions into account as a further damping effect beyond Landau damping, we start from a kinetic equation analogous to (6) with an additional collisional term $I[p, r, t]$ on the right hand side. In [11] we have derived a collision integral in a non-Markovian relaxation time approximation

$$I(p, r, t) = \frac{1}{\tau(t-t)} \int_0^t \frac{\tilde{f}(p, r, \bar{t}) - f(p, r, \bar{t})}{\tau(t-t)} dt$$

(13)

with the dynamical non-Markovian relaxation time

$$\frac{1}{\tau_m(\omega)} = \frac{1}{\tau_B} \left[ 1 + \frac{3}{4} \left( \frac{\omega}{\pi T} \right)^2 \right].$$

(14)

The Markovian relaxation time is given by $\tau_B^{-1} = \frac{8\pi m c^2}{3R_T} T^2$, where $\sigma$ is the averaged spin-isospin proton-neutron cross section. This collision integral holds for low temperatures compared to the Fermi energy. The non-Markovian relaxation time arises from the coupling of collective modes regarding two-particle scattering and consequently describes the effect of zero sound damping. The local equilibrium distribution $\tilde{f}$ is determined through the conservation of the local current. The linearization of the kinetic equation leads then to the extended response function of Mermin [43]

$$\Pi_0^M(q, \omega) = \frac{\Pi_0^M(q, \omega + \frac{\omega}{2})}{1 - \frac{1}{\omega T} \left( 1 - \frac{\omega}{\omega_T} \right)},$$

(15)

where the selfconsistency leads to the replacement of $\Pi_0^M$ by $\Pi_0^M$ in (12). Generalizations to asymmetric nuclear matter can be found in [13].

The energy and damping rates are determined by the zeros of the (Mermin) response function [11]

$$e^M(q, \Omega + i\gamma) = 1 - V_0(q) \Pi_0^M(q, \Omega + i\gamma) = 0.$$  

(16)

Here the damping rate represents Landau and collisional damping.

B. Finite matter response

In the next step we will present the finite matter response function in terms of a memory integral over all trajectories. It allows us to introduce the local density approximation as a first order memory effect in the trajectories. For this purpose we rewrite the Vlasov equation (6) in a slightly different way. Introducing the Lagrange picture by following the trajectory $x(t), p(t)$ of a particle we linearize the Vlasov equation (6) according to $f(x, p, t) = f_0(x, p) + \delta f(x, p, t)$ as

$$\frac{d}{dt}\delta f(x(t), p(t), t) = \partial_p f_0 \partial x(t) \delta V$$

(17)

and get with one integration

$$\delta f(x, p, t) = -2m \int_0^\infty dt' \int d\epsilon \epsilon \partial\epsilon f_0(p, \epsilon) \delta\epsilon f_0(p', \epsilon') \delta\epsilon \epsilon.$$  

(18)

The density variation caused by the external potential is obtained as

$$\delta n(x, \omega) = -2ms \int dx' \int dp' \partial p' \partial x' f_0(p', x') \delta\epsilon f_0(p, \epsilon) \delta\epsilon \epsilon.$$  

(19)

where $s$ denotes the spin-isospin degeneracy. Comparing this expression with the definition of the polarization function $\Pi_0,$

$$\delta n(x, \omega) = \int dx' \Pi_0(x, x', \omega) \delta V(x', \omega),$$  

(20)

we are able to identify the polarization of finite systems as

$$\Pi_0(x, x', \omega) = -2ms \int dp' \partial p' \partial x' f_0(p', x') \delta\epsilon f_0(p, \epsilon) \delta\epsilon \epsilon.$$  

(21)

Further simplifications are possible if we focus on the ground state $f_0(p^2) = \Theta(p^2 - p^2)$. The modulus integration of momentum can be carried out and the Kirzhnitz-formula [44,45] for the polarization function appears

$$\Pi_0(x, x', \omega) = \frac{-msf(x)\epsilon}{4\pi^2h^4} \left[ -2m \int_0^\infty dt' \frac{-i\epsilon}{dt'} \int d\Omega_{p} \frac{1}{4\pi} \delta(x' - x(t')) \right]$$

$$= \frac{-msf(x)\epsilon}{4\pi^2h^4} \left[ \delta(x' - x(0)) \right.$$

$$+ \frac{1}{4\pi} \int_0^\infty dt' \epsilon^2 \int d\Omega_{p} \frac{1}{4\pi} \delta(x' - x(t')) \right].$$  

(22)

4
This formula represents the ideal free part and a contribution which arises by the trajectories \( x(t) \) averaged over the direction at the present time \( \vec{n}_p, p_f = m \vec{v}(0) \). In principle, the knowledge of the evolution of all trajectories is necessary to evaluate this formula. Avoiding the latter expense, we discuss two approximations which will give us an insight into the physical processes behind. First the most restrictive one shows how the local density approximation emerges. As an extension we consider then the influence of chaotic scattering on a surface.

1. Local density approximation

The local density approximation appears from (22) when two simplifications are performed. Introducing Wigner coordinates \( R = (x + x')/2 \), \( r = x - x' \) one has to assume

1. gradient expansion
   
   \[
   p_f(R + r/2) \approx p_f(R) + O(\partial R)
   \]  

2. expansion of the trajectories up to first order history
   
   \[
   x' - x(t') \approx -r - t' \dot{x} = -r - t' \frac{p_f}{m} \vec{n}_p + O(t'^2).
   \]  

With these two assumptions we obtain from (22) after trivial integrations

\[
\Pi^\text{LDA}_0(q, R, \omega) = -\frac{m_{sp} f(R)}{4 \pi^2 \hbar^3} \int_0^\infty dy \exp \left( \frac{i \zeta y}{y} \right)
\]

where \( \zeta = \frac{m_{sp}}{m_{sp} f(R)} \). This can be further integrated with the help of

\[
\int_0^\infty dy \exp \left( \frac{i \zeta y}{y} \right) = \frac{\pi}{2} \left( \frac{1 + \zeta}{1 - \zeta} \right)
\]

leading to the standard Lindhard result for (25).

C. Influence of chaotic damping with surface on damping

Now we want to assume an additional chaotic scattering which will be caused e.g. by the curved surface. In order to investigate this effect we add to the regular motion (24) a small irregular part \( \Delta x \)

\[
x' - x(t') \approx -r - t' \frac{p_f}{m} \vec{n}_p + \Delta x.
\]

This irregular part of the motion has the direction of the velocity, \( \vec{n}_p \), and lasts a time \( \Delta t \). During this time an exponential increase in phase-space occurs controlled by the largest Lyapunov exponent \( \lambda \). Therefore we assume \( [t' < 0] \)

\[
\Delta x \approx \frac{p_f \vec{n}_p}{m} \Delta t \exp[-\lambda (t' - \Delta t)] + \text{const.}
\]

Since we are looking for the upper bound of Lyapunov exponent we can take (29) at the maximum \( \Delta t = -1/\lambda \). Furthermore, in the case of vanishing Lyapunov exponent, the regular motion (24) should be recovered. This determines the constant. We obtain finally

\[
x' - x(t') \approx -r - \frac{p_f}{m} \vec{n}_p \left[ 1 - \exp\left(-\frac{\lambda t'}{\lambda}\right) \right].
\]

With this ansatz one derives from (22) the result

\[
\Pi(q, R, \omega) = \frac{m_{sp} f(R)}{4 \pi^2 \hbar^3} \int_0^\infty dy \frac{\sin y}{y} \left( 1 + \frac{\zeta y}{\omega} \right)^{\frac{i \omega}{\lambda} - 1}
\]

which for \( \lambda \to 0 \) resembles exactly (25). The further integration could be given in terms of hypergeometric functions but this is omitted here.

With formula (31) we have derived the main result of this section, i.e. a polarization function due to many particle effects including the influence of an additional chaotic process characterized by the Lyapunov exponent \( \lambda \).

For small values of \( \frac{m_{sp}}{m_{sp} f(R)} \zeta \lambda \) which corresponds to relative small Lyapunov exponents, we can use \( \lim_{x \to \infty} (1 + \frac{a}{x})^x = \exp(a) \) in the integral of (31) and the final integration can be performed with the result of (27) and a complex shift in the frequency

\[
\Pi^\text{surf}_0(q, R, \omega) = \Pi^\text{inf}_0(q, p_f(R), \omega + i \lambda).
\]

Solving the dispersion relation we obtain in this way the Matthiessen rule which states that different damping mechanisms are additive in the total damping. However, the collisional damping of infinite matter, \( 2 \gamma \), and the
the largest Lyapunov exponent, $\lambda$, are not yet the appropriate values of finite matter. Instead, these two damping mechanisms have to be added by a proper relative weight which will take into account the relative occurrence of the processes, the collisions with other particles and the collisions with the surface.

**IV. COMPARISON OF DAMPING MECHANISMS**

**A. Connection between surface curvature and temperature**

The damping of excited nuclei can be understood by two different mechanisms. Besides the collisional damping we also have to consider the shape fluctuations. In the absence of inertia which will be assumed in the following, the driving force for shape fluctuations is the temperature within a statistical model. We use as a measure for the mean deformation $\alpha$ to the temperature within a statistical model. We use as a measure for the mean deformation $\alpha$ to

$$\langle \alpha \rangle = \frac{\int da |\alpha| \exp\left(-E_B(\alpha)/T\right)}{\int da \exp\left(-E_B(\alpha)/T\right)}$$

(33)

where the surface dependent energy $E_B(\alpha)$ is given by the Bethe-Weizsäcker formula

$$E_B(\alpha) = -a_1 + \frac{a_2}{A^{1/3}} + \frac{a_3Z^2}{A^{4/3}} + a_4\delta + a_5A^{2/3}S(\alpha)/S(0)$$

(36)

with the volume energy $a_1 = 15.68$ MeV, Coulomb energy $a_3 = 0.717$ MeV, the symmetry energy $a_4 = 28.1$ MeV and the surface energy $a_5 = 18.56$ MeV.

The surface of rotational symmetric nuclei (4) is given by

$$S = 2\pi \int_{-1}^{1} dx R_\lambda(x) \sqrt{R_\lambda(x)^2 + (1 - x^2)R'_\lambda(x)^2}$$

(37)

This represents the lowest order expansion in $\alpha$, however the next term gives already corrections in fractions of percents for the highest deformations considered here. By this way, the statistical model (33) leads to a connection between temperature and deformation as

$$T = c\pi a_5A^{2/3}/\alpha^2$$

(39)

where the constant $c$ is given by the coefficient of $\alpha^2$ in Eq. (38) for the corresponding quadrupole or octupole deformation.

**B. Finite matter scaling for collisional damping**

According to the local density approximation (27) we want to investigate the collisional damping in finite matter. Therefore we have to replace all densities in the dispersion relation (15) by the local density which are parameterized by a Woods-Saxon potential

$$n(r) = \frac{n_0}{\exp[(r-R_0)/0.545 \text{ fm}]+1}.$$  

(40)

Solving the dispersion relation (16) within the local density approximation according to (27) we obtain a spatial dependent damping rate $\gamma(r)$.

In the following we derive a finite size scaling factor, $\xi_A$, of the damping

$$\xi_A = \frac{\langle \gamma(r) \rangle}{\gamma},$$

(41)

which approximates the local ($r$-dependent) dispersion relation by an averaged one

$$\langle \gamma \rangle = \frac{3}{R_0^3} \int_0^{R_0} r^2 \ldots \, dr$$

(42)

assuming radial symmetry. The factor (41) has to be applied to the infinite matter bulk collisional damping and is found $\xi_{\text{Sn}} = 0.9199$ and $\xi_{\text{Pb}} = 0.9250$ for $^{120}\text{Sn}$ and $^{208}\text{Pb}$ respectively. Therefore, the bulk matter value of damping is diminished by finite size effects.

**C. Comparison of collisional damping and surface damping**

Using Eq. (39) we can translate the Lyapunov exponent $\lambda$ calculated as a function of deformation in figure 2
into a function of the temperature. In figure 3 the contribution to the damping of IVGDR for $^{120}$Sn (circles) and $^{208}$Pb (squares) is presented for different shape deformations versus temperature. If we add quadrupole and octupole deformations we come up with a damping curve very similar to the paper of Ormand [22]. The damping starts at zero and increases rapidly with increasing temperature. We see that the main contribution comes from the quadrupole deformation while the octupole deformation is only sizeable at higher temperature. Let us note that the qualitative difference between Sn and Pb is reproduced by surface scattering as well as collisional damping.

The collisional contribution 2γ scaled to finite sizes via (41) is compared with the damping according to the chaotic scattering from the surface of quadrupole and octupole deformed shapes for $^{120}$Sn and $^{208}$Pb from figure 2.

![Graph showing damping vs. temperature for Sn and Pb](image)

FIG. 3. The collisional damping scaled with finite size according to (41) is compared with the damping according to the chaotic scattering from the surface of quadrupole and octupole deformed shapes for $^{120}$Sn and $^{208}$Pb from figure 2.

We use the ratio of these two frequencies to weight properly the two damping mechanisms, the surface collisional, λ, and interparticle collisional, $< \gamma(r) >$, contributions. Consequently the full - width - half - maximum reads (FWHM)

$$\Gamma_{\text{FWHM}} = 2(\zeta < \gamma(r) > + (1 - \zeta) \lambda)$$

$$\equiv \Gamma_{\text{coll}} + \Gamma_{\text{surf}}.$$  (44)

With the help of (14) and (43) the weighting factor ζ is given by

$$\zeta(T) = \frac{1}{\frac{1}{\tau_{\text{surf}}(T)} + \nu_{\text{surf}}(T)}.$$  (45)

One sees that for zero and high temperatures ζ = 1 and due to Eq. (44) only the collisional contributions matter. Since νsurf is linear in the temperature and 1/τ depends quadratically on the temperature, the weighting factor ζ has a minimum at temperatures around $T_c = \frac{\sqrt{2}}{\pi} \omega$ and the surface contributions become important. In the case of the IVGDR this corresponds to a temperature of $T \approx 3.7 \text{MeV}$, which is the upper limit of current experimental achievable temperatures. Therefore we can state that at low and high temperatures the collisional damping is dominant while for temperatures around $T_c$ the surface contribution becomes significant.

In figure 4 we compare the effective damping according to Eq. (44) with the experimental data. We find a reasonable quantitative agreement.

![Graph showing effective damping vs. temperature for Sn and Pb](image)

FIG. 4. The effective damping consisting of collisional and surface damping together with the experimental data (filled symbols) from Ref. [48] (Sn) and from Ref. [49] (Pb).

While the collisional damping by itself can reproduce the increase at small temperatures well enough it fails to reproduce the sharp rise at higher temperatures, see figure 3. The surface damping by itself fails at zero temperature and leads to an unphysical shape. Only the connection of both effects, collisional and surface damping seems to be able to reproduce the data correctly.
At this point we like to stress that we have adopted a simplified model of finite Fermi liquids and introduced a phenomenological weighting between surface and interparticle collision contributions proportional to their relative occurrence. The discussion of damping due to finite size effects within shell model considerations can be found in literature, e.g. the review [1] where this problem has been addressed from the point of doorway states.

V. SUMMARY

We have considered the influence of collisional damping as well as surface scattering for the damping of hot isovector giant dipole resonances (IVGDR). We find that the relative importance of both effects is dependent on the temperature scale. While for low and high temperatures the collisional contribution dominates, the surface scattering is the one of significant importance at temperatures around 28% of the centroid energy.

The surface contribution has been calculated from the Lyapunov exponent of surface scattering for different deviations of the nucleus shape. The deviations are related to the temperature by a statistical model.

The influence of both the surface and collisional damping is described by a generalized response function for finite nuclear matter which takes into account the chaotic processes of scattering with the surface. We derive by this way a response function similar to the Lindhard response in local density approximation which is now modified by the Lyapunov exponent of the surface scattering. In an approximative way we show that the total damping is described by the sum of the two contributions to the damping in analogy to the Matthiessen rule.

Comparing the collision frequencies for one particle with other particles and for one particle with the surface, we derive a proper relative weight for both processes, the collisional and the surface damping. The resulting effective damping reproduces the experimental data rather well in the whole accessible range.

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