ON THE SCALING OF THE DAMPING TIME FOR RESONANTLY DAMPED OSCILLATIONS IN CORONAL LOOPS

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ABSTRACT

There is not as yet full agreement on the mechanism that causes the rapid damping of the oscillations observed by TRACE in coronal loops. It has been suggested that the variation of the observed values of the damping time as function of the corresponding observed values of the period contains information on the possible damping mechanism. The aim of this Letter is to show that, for resonant absorption, this is definitely not the case unless detailed a priori information on the individual loops is available.

Subject headings: MHD — Sun: corona — Sun: magnetic fields — waves

1. INTRODUCTION

Transverse oscillations in coronal loops have been detected in observations made with the EUV telescope on board the Transition Region and Coronal Explorer (TRACE) in 1999 by Aschwanden et al. (1999) and Nakariakov et al. (1999). Since then, the detection of these oscillations has been confirmed, and in addition damped oscillations have been observed in hot coronal loops by the SUMER instrument on board SOHO (Wang et al. 2002; Kliem et al. 2002). The TRACE observations have periods of the order of ≈2–10 minutes and comparatively short damping times of the order of ≈3–20 minutes. There is general consensus that these oscillations are fast standing kink mode oscillations. However, there is still debate about the mechanism that causes the observed fast damping. Different mechanisms have been suggested: phase mixing (Ofman & Aschwanden 2002), resonant absorption (Hollweg & Yang 1988; Goossens et al. 2002; Ruderman & Roberts 2002), lateral and footpoint wave leakage (Smith et al. 1997; De Pontieu et al. 2001), and drag due to the ambient plasma (Chen & Schuck 2007). In order to discriminate between different damping mechanisms, Ofman & Aschwanden (2002) suggested the study of how the observed damping times vary as a function of the corresponding observed periods. In particular, Ofman & Aschwanden (2002) claimed that the observed values are compatible with phase mixing if the damping time increases with period as \( T^{4/3} \) and with resonant absorption if it increases as \( T \). The aim of the present Letter is to show that random observations of oscillation events in coronal loops are very unlikely to produce any particular relation between damping times and periods whatever the actual mechanism is that causes the damping. In particular, we focus on resonant absorption and use synthetic data for periods and damping times to show that various samples of pairs of periods and damping times can lead to various and widely different scaling laws.

2. ANALYTICAL THEORY

The suggestion that the damping time increases linearly with period, \( \tau_d \sim T \), was triggered by analytical asymptotic expressions for the damping of the quasi-mode given by, e.g., Goossens et al. (1992), Ruderman & Roberts (2002), and Goossens et al. (2002) (see reviews Goossens et al. 2006 and Goossens et al. 2008). These asymptotic expressions are derived in the approximation that the nonuniform layer is thin. This is the so-called thin-boundary (TB) approximation. Jump conditions are used to connect the solution over the ideal singularity and to avoid solving the nonideal MHD wave equations (see, e.g., Sakurai et al. 1991; Goossens et al. 1995). The asymptotic expression for the damping time in the TB approximation is

\[
\tau_d = \frac{F}{l} \frac{R \omega_i^2 + 1}{\xi - 1},
\]

where \( \omega_i = k/v_A \) is the local Alfvén frequency and \( v_A = B/\mu \rho \) the local Alfvén velocity (the subscripts \( i \) and \( e \) refer to internal and external, respectively). Now we note that for the...
fundamental mode \( k_z = \pi/L \), with \( L \) the length of the loop, and we convert frequencies to periods and rewrite equation (2) as

\[
\frac{T}{\tau_{a,i}} = \sqrt{2} A(\xi),
\]

where \( \tau_{a,i} = L/v_{A,i} \) is the internal Alfvén travel time and the function \( A(\xi) \) is defined as

\[
A(\xi) = \left( \frac{\xi + 1}{\xi} \right)^{\nu/2}.
\]

It is convenient to adopt a reference loop with magnetic field strength \( B_0 \), length \( L_0 \), and internal density \( \rho_0 \). This reference loop allows us to consider loops of different dimensions and to introduce the normalized period \( T^* \) defined as

\[
T^* = \frac{T}{(\sqrt{2})\tau_{a,i0}},
\]

\[ T^* \] can be interpreted as the period measured in the unit \( \tau_{a,i0}/\sqrt{2} \). With this normalized period equation (3) takes the simple form

\[
T^* = a A(\xi),
\]

with \( a = \tau_{a,i}/\tau_{a,i0} \). From equation (6) we learn that, for a given value of \( a \), \( T^* \) varies continuously from \( \sqrt{2}a \) to \( \infty \) when \( \xi \) is allowed to vary from \( 1 \) (no loop) to \( \infty \) (outside vacuum).

Let us now turn back to equation (1) and rewrite it as

\[
\frac{T}{\tau_{a,i}} = \frac{2}{\pi} B(\xi) \frac{1}{l/R},
\]

with the function \( B(\xi) \) defined as

\[
B(\xi) = \left( \frac{\xi + 1}{\xi} \right)^{\nu/2} = \left( \frac{\xi + 1}{\xi - 1} \right).
\]

In what follows we shall refer to the combination of the TB approximation to compute the damping time and the TT approximation for the period as the TTTB approximation.

In the same way as for the period it is convenient to introduce the normalized decay time \( \tau_d^* \) as

\[
\tau_d^* = \frac{\tau_d}{(2\sqrt{2}/\pi)\tau_{a,i0}}.
\]

\( \tau_d^* \) can be interpreted as the decay time measured in the unit \( \tau_{a,i0}/(2\sqrt{2}/\pi) \). With the use of this normalized decay time we can write equation (7) as

\[
\tau_d^* = a B(\xi) \frac{1}{l/R},
\]

which tells us that, for a given value of \( a \), \( \tau_d^* \) depends on \( \xi \) and \( l/R \). The dependency on \( l/R \) is straightforward. The dependency on \( \xi \) is slightly more complicated. Basically, \( \tau_d^* \) is a decreasing function of \( \xi \) varying from \( +\infty \) for \( l/R \to 0 \) to \( h_0(l/R) \) for \( l/R \to +\infty \). In summary, \( \tau_d^* \) is a decreasing function of both \( l/R \) and \( \xi \).

We now combine equations (6) and (10). In Figure 1 we have plotted \( \tau_d^* \) versus \( T^* \) for values of \( \xi \) from 1.5 up to 10 and of \( l/R \) from 0.01 to 2. For every value of \( a \) the pairs \( (T^*, \tau_d^*) \) define a vertical strip in the \( (T^*, \tau_d^*) \)-plane with a maximum horizontal extent of \( (\sqrt{2} - 1)a \). As a matter of fact, \( a \) can take on any positive real value, but for clarity we have limited the values to \( a = 1, 2, 3, 4, 5 \). The inclusion of intermediate values of \( a \), such as \( a = 4/3, 5/3, 7/3 \), produces vertical strips overlapping those shown in Figure 1. The isolines corresponding to a constant value of \( \xi \) are vertical lines since the period is independent of the inhomogeneity length scale, in the thin-tube approximation. The isolines corresponding to a constant value of \( l/R \) are not straight and are defined by the equation

\[
\tau_d^* = \frac{(T^*)^3}{2a^2 - (T^*)^2} \frac{1}{l/R} \quad \text{for} \quad T^* \in [a, a\sqrt{2}].
\]

It is clear from Figure 1 that the model of resonant absorption in its most simple mathematical formulation using the TTTB approximation produces a wide variety of combinations of \( (T^*, \tau_d^*) \). Now let us see what happens when a collection of pairs \( (T^*, \tau_d^*) \) is drawn from this reservoir.

3. **Scaling Laws: How Many Do We Want?**

The aim of this section is to show that different collections of data can be produced by the simple TTTB mathematical model of resonant absorption with each of them leading to scaling laws, \( \tau_d^*/(T^*)^{n} = C \) (with \( C \) a function of the equilibrium parameters), with different indices, \( n \). We use equations (6) and (10), together with the particular relation between \( \tau_d^* \) and \( T^* \) for each index \( n \), and derive those collections of data. As a first example we have plotted in Figure 2 a collection of \( (T^*, \tau_d^*) \) combinations drawn from the big reservoir depicted in Figure 1. These combinations define a scaling law with index \( n = -1 \). It is clear that we have engineered the data presented in Figure 2 so as to fit the scaling law with such an index. Actually, the engineering is relatively straightforward. We require \( T^* \tau_d^* = C \). The left-hand side is the value of the function \( C \) for a starting configuration with \( a = a_s, \xi = \xi_s \) and \( l/R = (l/R)_s \). We fix a loop to start with by prescribing the values \( a_s = 1, \xi_s = 1.5, \) and \( (l/R)_s = 0.1 \) and compute values of \( a, \xi \), and \( l/R \) that satisfy the equation

\[
\frac{l}{R} = \left( \frac{1}{R} \right) \left( \frac{a}{a_s} \right)^2 \left( \frac{f(\xi)}{f(\xi_s)} \right) \quad \text{for} \quad f(\xi) = A(\xi)B(\xi).
\]
Likewise in Figure 2 we have also plotted a second collection of data that now define a scaling law with index \( n = 2 \). To obtain these data, we start from a loop with prescribed values \( a_s = 1 \), \( \xi_s = 5 \), and \((ll/R)_s = 1\) and compute values of \( a \), \( \xi \), and \( ll/R \) that satisfy the equation

\[
\frac{l}{R} = \left(\frac{l}{R}\right)_s \frac{a_g(\xi)}{a_g(\xi_s)} ; \quad g(\xi) = \frac{B(\xi)}{A(\xi)}^2 ;
\]

(13)

A third collection of data that now define a scaling law with index \( n = 4/3 \) is also shown in Figure 2. According to Ofman & Aschwanden (2002), this value of the index singles out phase mixing as damping mechanism. Here it is obtained for a collection of data that are produced by the theoretical prediction for resonant absorption. We start from a loop with prescribed values \( a_s = 1 \), \( \xi_s = 3 \), and \((ll/R)_s = 0.5\) and compute values of \( a \), \( \xi \), and \( ll/R \) that satisfy the equation

\[
\frac{l}{R} = \left(\frac{l}{R}\right)_s \frac{a_h(\xi)}{a_h(\xi_s)} ; \quad h(\xi) = \frac{B(\xi)}{A(\xi)}^{4/3}.
\]

(14)

Finally, in Figure 2 we have also plotted a collection of data that define a scaling law with the canonical value \( n = 1 \) for the index. We start from a loop with prescribed values \( a_s = 1 \), \( \xi_s = 2.5 \), and \((ll/R)_s = 0.2\) and compute values of \( a \), \( \xi \), and \( ll/R \) that satisfy the equation

\[
\frac{l}{R} = \left(\frac{l}{R}\right)_s \frac{c(\xi)}{c(\xi_s)} ; \quad c(\xi) = \frac{B(\xi)}{A(\xi)} = \frac{\xi + 1}{\xi - 1}.
\]

(15)

Note that \( a \) is absent from equation (15), meaning that a solution of equation (15) can combined with any value of \( a > 0 \). An obvious solution to equation (15) is

\[
\xi = \xi_s, \quad \frac{l}{R} = \left(\frac{l}{R}\right)_s, \quad a > 0.
\]

(16)

Equation (16) defines a straight line in the \((T^*, \tau^*_{\xi})\)-plane. Note that equation (16) is not the only solution in the \((T^*, \tau^*_{\xi})\)-plane. Any combination \((T^*, \tau^*_{\xi})\) that satisfies equation (15) combined with \( a \in [0, \infty[ \) produces the same straight line in the \((T^*, \tau^*_{\xi})\)-plane as solution (16). Since the function \( c(\xi) \) is a decreasing function of its argument, a value \( \xi > \xi_s \) is combined with a value \( ll/R < (ll/R)_s \), and conversely a value \( \xi < \xi_s \) is combined with a value \( ll/R > (ll/R)_s \).

It is clear from Figure 2 that almost any scaling law can be obtained from data produced by the simple TTTTB mathematical model of resonant absorption. All the periods and damping times plotted in Figure 2 correspond to combinations of \( \xi \) and \( ll/R \), obtained from equations (12), (13), (14), and (15), that are reasonable values of these quantities in the ranges \( \xi \in [1.5, 10] \) and \( ll/R \in [0.016, 1.59] \). The discrete sets of values in Figure 2 arise due to the considered discrete values of \( a \), but loops with different internal travel times, with respect to the reference loop, would give the full set of values for \( \tau^*_\xi \) and \( T^* \) along the solid lines depicted in Figure 2.

4. BEYOND THE TTTTB APPROXIMATION

The TTTTB approximation as a mathematical model for resonant absorption has its clear limitations. First, the values of the period defined by equation (6) are independent of the radius \( R \) and the inhomogeneity length scale \( ll/R \). In reality, the period is a function of the density contrast \( \xi \), \( ll/R \), and \( R \). As a consequence, the isolines corresponding to a constant value of \( \xi \) in the \((T^*, \tau^*_{\xi})\)-plane are no longer straight vertical lines. For a given value of \( a \), this dependency of the period on \( \xi \) produces additional spread in the original vertical strips. Second, the TTTTB approximation is an accurate approximation as long as the damping time is sufficiently longer than the period, since this is an inherent assumption for carrying out the asymptotic analysis leading to equation (1). We can expect this equation to be inaccurate for large values of the density contrast \( \xi \) and large values of the inhomogeneity length scale \( ll/R \). Figure 3 is the twin version of Figure 1 with the values of \( T^* \) and \( \tau^*_\xi \) now computed for fully nonuniform 1-D loops (Van Doorsselaere et al. 2004). Figure 3 is very similar in appearance to Figure 1. The differences occur where they are expected to occur. The vertical strips of Figure 1 are now replaced with strips slightly oblique to the vertical axis, and the largest differences appear at the short values of the damping times corresponding to the large values of the density contrast \( \xi \) and large values of the inhomogeneity length scale \( ll/R \). The basic message from both figures is the same. The model of resonant absorption produces a wide variety of combinations in the \((T^*, \tau^*_{\xi})\)-plane. If we draw a curve defined by a given relation between \( T^* \) and \( \tau^*_\xi \), for instance \( \tau^*_\xi(T^*)^2 = C \), in this plane we can graphically determine as many points on this curve as we want. If we select the data determined in this way the result is a scaling law with index \( n = 2 \) as the one in Figure 2. The only difference is that now the procedure is numerical all the way.
5. CONCLUSION

In this Letter we have explained that scaling laws for a group of observations of oscillating loops to discriminate between different damping mechanisms are of not much use if there is no a priori information on the properties of the loops. The analytical expressions obtained by the TTTB approximation enable us to show in a straightforward and easy-to-follow manner that in the framework of resonant absorption collections of synthetic data can be produced that follow almost any scaling law. Then, we have backed up our findings by numerical eigenvalue computations that do not suffer from the TTTB limitations but are on their own rather less instructive.

It might be argued that nature is not as cunning as the authors of this Letter and does not attempt to engineer data according to prescribed rules as those defined in these equations. On the other hand, there is no reason why nature would want to produce coronal loops that all have the same values of $\xi$ and $l/R$. So, unless there is accurate a priori information on the coronal loops available, the use of scaling laws to discriminate between different damping mechanisms is questionable, to say the least.

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