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# Singularities Arising in the Linearized Magnetohydrodynamics Equations

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**ABSTRACT:** The magnetohydrodynamics equations in a cylinder linearized around an equilibrium state with vertical magnetic field reduce to a single singular second order partial differential equation where the time frequency enters as a parameter. Although there could exist whole families of solutions depending on arbitrary values at the singularity [1], only one of them is physically correct. We analyze theoretically and numerically the problem and the physical consequences of our results.

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## 1 THE SETTING OF THE PROBLEM.

Let a stationary plasma fill a hollow cylinder  $r_0 < r < r_1$ , and assume that its density  $\rho$  and magnetic field  $\mathbf{B} = (0, 0, B)$  depend only on the radius r, and that it satisfies the polytropic state equation  $p = A\rho^{\gamma}$ . For a fixed Fourier mode l in z, let u be the Fourier-Laplace transform  $(t \to \omega)$  of a small perturbation  $p_*$  of the total (kinetic plus magnetic) pressure. If the values of the perturbed velocity and field are zero at t = 0, u satisfies

div 
$$\left(\frac{1}{\omega^2 \rho - l^2 B^2} \nabla u\right) + \frac{1}{\gamma p} \frac{\omega^2 \rho - l^2 \gamma p}{\omega^2 \rho \left(1 + \frac{B^2}{\gamma p}\right) - l^2 B^2} u = 0.$$
 (1)

The points where the local Alfvén frequency  $\omega_A = \pm l B/\sqrt{\rho}$  matches the time frequency  $\omega$  are singular and very important in the study of magnetic oscillations, [2, 3, 4]. To simplify the study of this singularity, we will consider only the principal part of the differential operator and set the second term of (1) to zero; let  $F = \omega^2 \rho - l^2 B^2$ . After a Fourier transform  $\theta \to n$  in the angle variable, (1) becomes

$$u'' + \left(\frac{1}{r} - \frac{F'(r)}{F(r)}\right) u' - \frac{n^2}{r^2} u = 0.$$
 (2)

At any simple zero  $r_0$  of F, F'/F behaves as  $1/(r - r_0)$ . For further notational simplification, and without affecting the properties of the equation we will take  $l^2B^2 = 1$  and  $\rho(r) = r$ ; for  $\omega^2 = 1$ , (2) becomes

$$u'' + \left(\frac{1}{r} - \frac{1}{r-1}\right) \, u' - \frac{n^2}{r^2} \, u = 0 \, ,$$

which we recognize as an hypergeometric equation. Its classical theory [5, 6] allows us, after a cumbersome but essentially straightforward calculation, to write two independent solutions in B(1, 1) as

$$v_{1}(r) = r^{n}(1-r)^{2} \sum_{k=0}^{\infty} \frac{(\alpha(n))_{k}(\beta(n))_{k}}{(3)_{k}} \frac{(1-r)^{k}}{k!}$$
$$v_{2}(r) = -\frac{2r^{n}}{(\alpha(n)-2)_{2}(\beta(n)-2)_{2}} + r^{n}(1-r) \sum_{k=0}^{\infty} \frac{(\alpha(n))_{k}(\beta(n))_{k}}{(3)_{k}} \left(\Psi(\alpha(n)+k) + \Psi(\beta(n)+k) - -\Psi(3+k) - \Psi(1+k)\right) \frac{(1-r)^{k+2}}{k!} + \log(1-r)v_{1}(r),$$

where  $(x)_k = (x+k)/, (x), \alpha(n) = n+3/2 + \sqrt{n^2 + 1/4}, \beta(n) = n+3/2 - \sqrt{n^2 + 1/4},$ and  $\Psi$  is the logarithmic derivative of the , function.

## 2 MAIN RESULTS.

We see that  $v_2(r) = w_2(r) + \log(1-r)v_1(r)$ , with  $v_1$  and  $w_2$  real functions. Also notice that the wronskian determinant  $w'_2v_1 - w_2v'_1$  behaves as  $a(r-1) + O(r-1)^2$ near 1, with  $a = 4/(\alpha(n) - 2)_2 (\beta(n) - 2)_2 > 0$ . These function are  $C^1$  and, except for the term  $\log(1-r)$ , real valued in  $[r_0, r_1]$ . In the general case we would have  $\log(r(\omega) - r)$ , where  $r(\omega)$  is the (complex) zero of F for  $\omega$  near  $\pm 1$ . Hence, for  $\delta > 0$ ,  $\log(1 - (1 + \delta)) = \log(1 - (1 - \delta)) + i \sigma \pi$ , where  $\sigma = \pm 1$ , according to which branch of the logarithm is chosen. Let us study the correct determination of this imaginary part.

**Lemma 1** Let  $\omega_0 = \pm 1$ . Then  $\sigma$  is the sign of  $\omega_0/F'(1)$  (F'(1) = 1 in our case).

**Proof.** What we denote by u is in fact the Fourier-Laplace transform  $\int_0^\infty e^{i\omega t} p_*(t, \mathbf{x}) dt$ ; hence the solution of (1) for real  $\omega$  should be the limit of (damped) solutions for  $\omega + i \varepsilon, \varepsilon \downarrow 0$ . Let  $\omega(r)$  be the (complex) frequency which makes  $F(\omega(r), r) = 0$  near  $(\omega_0 = \pm 1, r = 1)$ . Then

$$\frac{\partial F}{\partial \omega} \,\omega' + F' = 0\,,$$

that is,  $2\omega(r)\rho(r)\omega'(r) + F'(r) = 0$ . This implies

$$\omega(r) - \omega(1) = -\frac{F'(\omega(1), 1)}{2\omega(1)\rho(1)} (r - 1) + O(r - 1)^2;$$

thus, if  $r(\omega)$  is the (complex) zero of  $F(\omega, r(\omega))$  near ( $\omega = \omega_0, r = 1$ ),

$$r(\omega) - r(\omega_0) = r(\omega) - 1 = -\frac{2\omega_0\rho(1)}{F'(\omega_0, 1)}(\omega - \omega_0) + O(\omega - \omega_0)^2.$$

Hence the sign of the imaginary part of  $r(\omega_0 + i\varepsilon)$  is the sign of  $-\omega_0/F'(1)$ . When positive,  $r(\omega)$  lies in the lower half plane and when r goes from  $1 - \delta$  to  $1 + \delta$ ,  $r(\omega) - r$ goes from  $r(\omega) - 1 + \delta$  to  $r(\omega) - 1 - \delta$  in the lower half plane, so that its argument changes from 0 to  $-\pi$ ; otherwise, from 0 to  $\pi$ . Always  $\sigma = \operatorname{sgn}(\omega_0/F'(1))$ .  $\Box$ 

How this fact relates to energy absorption? The increase of energy in X is given by

$$\frac{dE}{dt} = -\int_{\partial X} p_* \, \mathbf{v} \cdot \mathbf{n} \, d\sigma$$

(where **n** denotes the outer normal). It may be shown that the Fourier transforms P and V of  $p_*$  and  $\mathbf{v} \cdot \mathbf{n}$  satisfy  $V = (i \,\omega/F) P'$ . Since we are dealing with a single resonant sheet, assume that  $p_*$  has only the Fourier modes  $\pm l$  in z (so that  $F = \omega^2 \rho - l^2 B^2$  is fixed in the problem). We will study the integral above for a cylindrical surface r = a. As we know, under fairly general conditions, and with a somewhat relaxed notation,

$$\begin{split} &\int_{\mathbf{T}\times\mathbf{IR}} p_*(t,a,\theta,z) \,\mathbf{v}\cdot\mathbf{n}(t,a,\theta,z) \,d\theta \,dz = \\ &= \int_{\mathbf{Z}\times\mathbf{IR}} \hat{p}_*(t,a,n,l) \underbrace{\mathbf{v}\cdot\mathbf{n}}_{\mathbf{v}\cdot\mathbf{n}}(t,a,n,l) \,dn \,dl \\ &= \sum_{n\in\mathbf{Z}} \left( \left( P(t,a,n,l) \,V(t,a,-n,-l) + P(t,a,n,-l) \,V(t,a,-n,l) \right) \right). \end{split}$$

Hence the output of energy through r = a is

$$-\sum_{n \in \mathbb{Z}} \left( \left( P(t, a, n, l) \, V(t, a, -n, -l) + P(t, a, n, -l) \, V(t, a, -n, l) \right) \,. \tag{3}$$

Assume also that the boundary conditions determine a one-dimensional space of solutions depending only on  $\omega^2$  (such as decay at infinite,  $u(r_2 = \infty) = 0$ , or a solid wall  $u'(r_1) = 0$ , etc.; most real problems are in this class). We may construct a generator of the space of solutions as follows:

Let  $\alpha = \lambda_1 v_1 + \lambda_2 v_2$ ,  $\lambda_1, \lambda_2$  real;  $\beta = \delta \pi \lambda_2 v_1$ , with  $\delta = 0$  for r < 1 or  $\delta = 1$  for  $r \geq 1$ . Then

$$P = \lambda(1, n, l) (\alpha + i\beta) e^{it} + \lambda(-1, n, l) (\alpha - i\beta) e^{-it},$$
  

$$V = \frac{i}{F} \lambda(1, n, l) (\alpha' + i\beta') e^{it} - \frac{i}{F} \lambda(-1, n, l) (\alpha' - i\beta') e^{-it}.$$

Since  $p_*$  is real,  $P(-\omega, -n, -l) = \overline{P(\omega, n, l)}$ , so the same happens for the coefficient  $\lambda$ . Also dE/dt is real, so it is enough to take the real part of the terms in (3). The real part of the non-oscillatory terms (that is, the sums in (3) not affected by  $e^{2it}$  or  $e^{-2it}$ ; the time mean of these is zero and hence they are not relevant in the net energy increase) is therefore

$$\frac{1}{F} \left( |\lambda(-1,n,l)|^2 + |\lambda(1,n,l)|^2 \right) \left( \alpha'\beta - \alpha\beta' \right)$$

for a fixed mode *n*. The wronskian determinant  $\alpha'\beta - \alpha\beta'$  vanishes for r < 1, and satisfies W' = (F'/F - 1/r)W; hence W = AF/r for  $r \ge 1$ , A a constant. One easily gets

$$A = \frac{r}{F} \lambda_2^2 \pi (v_2' v_1 - v_2 v_1') = \frac{r}{F} \lambda_2^2 \pi (w_2' v_1 - w_2 v_1' + \frac{1}{r-1} v_1^2) \,.$$

which as shown before, is positive at r = 1. (Recall that F = r - 1,  $w'_2 v_1 - w_2 v'_1 = a(r-1) + O(r-1)^2$ , a > 0,  $v_1^2 = O(r-1)^4$ .)

Hence for a set X bounded for r = a and r = b, dE/dt is a continuous function of (a, b) whenever both are less than or greater than 1; however, if a < 1, b > 1, dE/dt jumps by a fixed amount. We have proved

**Theorem 1** Energy absorption due to alfvenic resonance takes place at the singular set.

Notice that the fact that  $\beta \neq 0$  (that is, that an imaginary part occurs at the solution after the singularity r = 1) is essential for the existence of this phenomenon. Any solution not taking this term in consideration is physically erroneous. This would happen if one applies directly any of a number of numerical methods of integration which tolerate the singular point, but do not get the solution analytic in  $\omega$ , as we will show next.

#### **3 NUMERICAL RESULTS.**

The integration of the initial value problem in a given interval (in our case  $r \in [0.5, 1.5]$ ) is delicate because of the singularity ar r = 1. If however, one takes  $r = x + i\varepsilon$ ,  $x \in [0.5, 1.5]$ ,  $\varepsilon > 0$ , not only the singularity disapears, but what we obtain is the correct solution in the limit  $\varepsilon = 0$ . This has been done in equation (2) for initial values u(0.5) = 1, u'(0.5) = 1, n = 1, 2, and  $\varepsilon = 10^{-1}, 10^{-2}, 10^{-3}$  (see figure 1). Behaviour for n = 2 is similar, as shown in figure 2. Observe that the



**Figure 1**: Left: real part of the solutions, for n = 1; right: imaginary part of the solutions, for n = 1.

behaviour of the imaginary part is as predicted: zero up to the singularity r = 1, and then a nontrivial function. By contrast, direct integration of (1) with  $\varepsilon = 0$ , even if the algorithm somewhat surpasses the critical point at r = 1, will never generate an



Figure 2: Left: real part of the solutions, for n = 2; right: imaginary part of the solutions, for n = 2.

imaginary part for a real equation with real initial value. We have checked that for  $\varepsilon = 0$ ; the real part is correctly obtained but the imaginary part of u is always zero, as one would expect.



Figure 3: Boundary value problems for n = 2; real and imaginary parts of the solution.

As for the boundary problem, the same method may be used. The regular boundary value problems obtained by considering  $r = x + i\varepsilon$ ,  $x \in [0.5, 1.5]$ , with the conditions u(0.5) = 0 = u(1.5) and  $\varepsilon = 10^{-1}, 10^{-2}, 10^{-3}$ , show the convergence of the solutions to the correct physical solution. For n = 2, we have the results of the figure 3.

These graphics have been obtained by using the shooting method with Mathematica, which allows a high degree of interactivity and is specially good at dealing with singularities. If we execute this method setting directly  $\varepsilon = 0$ , we still get an answer, but a completely wrong one (fig. 4).

In general, the behaviour of this integrator for  $\varepsilon = 0$  is erratic. For n = 1, one gets the correct answer for some try values of u'(0.5) and a wrong one for others. For n = 2 it is always false.

Finally, to emphasize the fact that one cannot choose the value of u at r = 1 arbitrarily and then solve the boundary problem at both intervals [0.5, 1], [1, 1.5], we have set u(1) = 10 and found the solutions for n = 2 for a slightly attenuated frequency  $\omega = 1 + 0.1i$  (fig. 5). The solutions are obviously not differentiable. This



**Figure 4**: The wrong answer for  $\varepsilon = 0$ .

shows that the problem should not be treated as an equation with a singular boundary; one cannot impose arbitrary Dirichlet conditions at r = 1. Those are determined by the original problem in a complex variable setting.



**Figure 5**: The solutions for u(1) = 10; real and imaginary part.

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