An efficient method for computing the eigenfunctions of the dynamo equation

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ABSTRACT

Aims. We present an elegant method of determining the eigensolutions of the induction and the dynamo equation in a fluid embedded in a vacuum.

Methods. The magnetic field is expanded in a complete set of functions. The new method is based on the biorthogonality of the adjoint electric current and the vector potential with an inner product defined by a volume integral over the fluid domain. The advantage of this method is that the velocity and the dynamo coefficients do not have to be differentiated, and thus even numerically determined tabulated values of the coefficients produce reasonable results.

Results. We provide test calculations and compare with published results obtained by the classical treatment based on the biorthogonality of the magnetic field and its adjoint. We especially consider dynamos with mean-field coefficients determined from direct numerical simulations of the geodynamo and compare with initial value calculations and the full MHD simulations.

Key words. Magnetohydrodynamics (MHD) – magnetic fields – methods: numerical

1. Introduction

The generation and evolution of magnetic fields in cosmic bodies like the planets and stars is generally thought to be governed by induction processes due to motions in their electrically conducting fluid interior. The magnetic field $B$ is described by the induction equation

$$\frac{\partial B}{\partial t} = \nabla \times DB$$  (1)

where

$$DB = u \times B - \eta \nabla \times B.$$  (2)

Here $u$ represents the velocity and $\eta$ the magnetic diffusivity.

In the framework of mean-field theory (e.g. Moffatt 1978; Krause & Rädler 1980), $u$ and $B$ are considered as mean, e.g. ensemble averaged, quantities, whereas the action of the small-scale turbulent flow on the mean magnetic field is parametrised by the so-called dynamo coefficients, $\alpha$ and $\beta$. They are, in general, tensors of second and third rank, respectively. We use the following compact notation of the mean field coefficients, which include the so-called $\gamma$, $\delta$, and $\kappa$-effects, see e.g. Rädler (1980). Then, the operator $D$ reads

$$DB = u \times B + \alpha \cdot B - \beta : (\nabla B) - \eta \nabla \times B$$  (3)

instead of (2), and acts on the mean magnetic field. Except for the additional $\alpha$ and $\beta$ terms in the $D$ operator, the induction and the dynamo equation are formally equivalent. Thus, the new method presented here equally applies to both.

The dynamo region is located in a flow domain $V$ with exterior vacuum $E$. In this work we assume $V$ to be either a sphere or a spherical shell. The magnetic field $B$ is continuous through the boundary $\partial V$ and potential in $E$.

In kinematic dynamo theory, all coefficients ($u$, $\alpha$, $\beta$, and $\eta$) are assumed given and independent of the magnetic field. Thus the dynamo equation is linear in the magnetic field and can be solved by considering an eigenvalue problem

$$\lambda B = \nabla \times DB$$  (4)

with eigenvalues $\lambda$ describing the time evolution proportional to $\exp(\lambda t)$ of the magnetic field $B$.

Many studies have been made of the eigenvalues of the dynamo operator for various celestial bodies and with many forms of the dynamo coefficients (e.g. Bullard & Gellman 1954; Roberts 1960; Steenbeck & Krause 1969; Deinzer & Stix 1971; Roberts 1972; Roberts & Stix 1972; Gubbins 1973; Kumar & Roberts 1975; Schmitt & Schüssler 1989; Dudley & James 1989; Deinzer et al. 1993; Gubbins et al. 2000; Schubert & Zhang 2001; Livermore & Jackson 2004, 2005; Jiang & Wang 2006, 2007). Often the coefficients are approximated by simple analytical functions of position, and their tensorial character is disregarded. Recently, the test-field method, developed by Schrinner et al. (2005, 2007) (see also Ossendrijver et al. 2001, 2002), allows one to determine all tensorial components of $\alpha$ and $\beta$ directly from self-consistent numerical simulations (Brandenburg et al. 2008; Käpylä et al. 2009). These coefficients are sometimes strongly varying functions of position. This may introduce large errors because the dynamo operator $\nabla \times D$ involves differentiation of the dynamo coefficients, and these are only available as numerically determined tabulated values.

In this paper we present a new method that does not require differentiation, so it is also applicable to numerically determined dynamo coefficients. The method is based on the biorthogonality of the electric current and the vector potential with an inner prod-
The integration volume constitutes a biorthogonal set. For a given set of functions with the following inner product or the fluid domain presented in Sect. 3. In Sect. 4 we apply the new method to eigenmodes of the dynamo operator with coefficients obtained from geodynamo models. Our conclusion are drawn in Sect. 5.

2. Eigenvalue problem

We expand the field $B$ of the dynamo in a complete set of functions $b_i(r)$:

$$ B = \sum_i e_i b_i(r) \equiv e_i b_i .$$  \hspace{1cm} (5)

Here and in the following we make use of the summation convention for two identical indices. The expansion functions are often eigenfunctions of some differential operator. Since this operator is, in general, not self-adjoint, the functions are not orthogonal. This problem is handled by using the adjoint set $\hat{b}_i(r)$, with the following inner product

$$ (\hat{b}_k , b_i)_X \equiv \int_X \hat{b}_k \cdot b_i \, d^3r = \delta_{ki} .$$  \hspace{1cm} (6)

The integration volume $X$ can be either the whole space $V + E$ or the fluid domain $V$ alone. The base functions $b_k(r)$ and $b_i(r)$ constitute a biorthogonal set. For a given set of functions $b_i$, the adjoint set $\hat{b}_i$ depends on the choice of the integration domain $X$, so in principle we have two different sets $\hat{b}_i$, one for $X = V$ and one for $X = V + E$.

Later we adopt the free magnetic decay modes, for which the base functions $b_i$ and their adjoints $\hat{b}_i$ are known. But at this point there is no need to specify which set $b_i$ we actually use.

2.2. Classical eigenvalue method

Inserting the expansion (5) in the dynamo eigenvalue equation (4) yields

$$ \lambda e_i b_i = \nabla \times (D e_i b_i) .$$  \hspace{1cm} (8)

Subsequently, we take the inner product (6) based on $V$ with the adjoint magnetic field. This leads to

$$ \lambda e_i = M_{ki} \hat{e}_i \quad \text{with} \quad M_{ki} = (\hat{b}_k , \nabla \times D b_i)_V .$$  \hspace{1cm} (9)

A partial integration to shift the curl from the second to the first term, as done in (7) above and used in the new method below, is not possible because the surface term $\int_V (D \hat{b}_k \times \hat{b}_i) : \cdot d^2\sigma$ need not vanish here.

We mention as an aside that the magnetic field is often decomposed in its poloidal and toroidal components (see Appendix A) after which the dynamo equation is formulated in terms of the defining scalars $P$ and $T$. If the dynamo coefficients possess certain symmetry properties, the solutions can be split into two independent subsets, describing magnetic fields symmetric and antisymmetric with respect to the equator.

2.3. New eigenvalue method

We start again with (8), which we uncurl to obtain

$$ \lambda e_i a_i = D e_i b_i + \nabla \psi .$$  \hspace{1cm} (10)

Taking now the inner product (7) with the adjoint current results in

$$ \lambda e_i = N_{ki} a_i \quad \text{with} \quad N_{ki} = (j_k , D b_i)_V .$$  \hspace{1cm} (11)

The gradient term drops out as discussed in Sect. 2.1 above. The corresponding adjoint functions $\hat{b}_i$ here are different from those in Sect. 2.2 as they pertain to a different inner product.

The matrices $M_{ki}$ and $N_{ki}$ have the same eigenvalues $\lambda$. The advantage of the new method using $N_{ki}$ in (11) instead of $M_{ki}$ in (9) is that no differentiation of the operator $D$ is required, so even numerically computed or tabulated values of $u$, $\alpha$, and $\beta$ produce accurate results.

2.4. Choice of $b_i$ and numerical handling of (11)

For the set of base functions, we adopt the free magnetic decay modes whose magnetic fields $b_i$ are known analytically in $V + E$ in terms of the defining scalars $P$ and $T$ as described in Appendix A. The decay modes are continuous through $\partial V$ and potential in $E$, so they satisfy the boundary conditions. They are characterised by three numbers, the radial order $n$, the latitudinal degree $\ell$, and the azimuthal order $m$.

Another advantage of the decay modes is that they are self-adjoint on $V + E$ so that the adjoint functions are the complex conjugates $\hat{b}_i = b_i^\ast$ and likewise $\hat{j}_k = j_k^\ast$. Normalisation on $V + E$, i.e. $(b_i^\ast, b_i)_{V+E} = (j_k^\ast, a_i)_V = \delta_{ki}$, is thus readily achieved, see Appendix A.

The computation of the matrix elements $N_{ki}$ is now straightforward. Once we know the matrix elements, the eigenvalue problem (11) is solved numerically using LAPACK routines (http://www.netlib.org/lapack), and we obtain the eigenvalues $\lambda_k$ and eigenvectors $[e_{k,i}]$, such that

$$ B_k = e_{k,i} b_i .$$  \hspace{1cm} (12)
is eigenfunction of $\nabla \times \mathbf{D}$ with eigenvalue $\lambda_i$. Each mode $k$ contains, in general, a mixture of $n$, $l$, and $m$ values.

In the following we consider only velocities and dynamo coefficients that are independent of azimuth $\varphi$, but this is not a necessary constraint. Thus each value of $m$ can be treated separately. Although we present only results for $m = 0$ here, we have tested and applied other values of $m$ as well. We employ the robust Gauss-Legendre quadrature in $r$ and $\cos \theta$ to compute the matrix elements since the basis functions are heavily oscillatory in $r$ for high values of $n$ and in $\theta$ for high degree $l$. For the Gauss-Legendre integration we used 66 quadrature points here in the radial and 80 in the latitudinal direction, respectively. In general, this depends of course on the required resolution.

### 2.5. Adjoint eigenfunctions

We now show how one may construct the adjoint set of eigenfunctions $\mathbf{B}_p^*$ of a set of eigenfunctions $\mathbf{B}_i$ of the dynamo operator $\nabla \times \mathbf{D}$. Although these adjoints are not presented in the present paper, they appear in applications. For example, let $\mathbf{B}$ be the actual magnetic field of the dynamo, then it is often advantageous to expand $\mathbf{B}$ in eigenfunctions, i.e., $\mathbf{B}(r, t) = \sum_i c_i(t) \mathbf{B}_i(r)$. To find the coefficients $c_i(t)$ we use the adjoint set, to find $c_i = (\mathbf{B}_i, \mathbf{B})_{V+E} = (\mathbf{J}_i, \mathbf{A})_V$.

This illustrates that we need the adjoints $\mathbf{B}_p^*$, and these may be constructed as follows. Let $\mathbf{B}_k = e_k \mathbf{b}_k$ be the representation of $\mathbf{B}_k$ in terms of the self-adjoint magnetic decay modes as above. Then we write $\mathbf{B}_p^* = f_{pi}^* \mathbf{b}_i^*$ and $\mathbf{J}_p^* = f_{pj}^* \mathbf{j}_i^*$, and we require

$$\delta_{pk} = (\mathbf{J}_p, \mathbf{A}_k)_V = (f_{pj}^* \mathbf{j}_j^*, e_k \mathbf{a}_j)_V = f_{pi}^* e_k \delta_{ij} = f_{pi}^* e_{kj}. \quad (13)$$

Therefore $f_{pi}^* = e_{ji}^*$ in matrix notation, and we find a unique biorthogonal set. Here, $\dagger$ indicates the Hermitean adjoint, $f_i = (f_i^*)^\dagger$, where $T$ indicates the transposed and * complex conjugation.

Three important messages follow from this construction: (i) the adjoint of $\nabla \times \mathbf{B}$ is $\nabla \times \mathbf{B}^*$, that is, the adjoint operation commutes with $\nabla$; (ii) to obtain the adjoint eigenfunctions, it is not necessary to know the explicit form of the adjoint dynamo operator $\nabla \times \mathbf{D}$; and (iii) the eigenfunctions and their adjoints have the same boundary conditions because they are a linear combination of the decay modes and their complex conjugates, respectively.

### 3. Test results

#### 3.1. $\alpha^2$-sphere

We first consider the so-called $\alpha^2$-sphere of unit radius $r_0 = 1$, represented by $\mathbf{u} = \beta = 0$, $\alpha_{ij} = R_\alpha \delta_{ij}$, and $\eta = 1$, which also can be treated analytically (Krause & Rädler 1980, Chap. 14). The eigenvalues are independent of azimuth $m$, and the eigenfunctions decouple in latitudinal quantum number $l$. For $R_\alpha \neq 0$, the modes couple in radial number $n$, as they do between the poloidal and toroidal components.

For $R_\alpha = 4.493409458$, the first mode is a stationary dipole, while the overtones decay with the rates given by Hoyng & van Geffen (1993) (HvG93). We successfully reproduced the fundamental mode and the overtones. In Table 1 we consider the convergence in the eigenvalues as a function of the maximum radial number $n_{\text{max}}$ for some dipolar ($l = 1$) modes. Higher $l$ modes behave similarly. We also reproduced the eigenfunction plots as provided by Krause & Rädler (1980).

#### 3.2. Spherical flows

As a next test, we apply the spherical stationary $\ell_1 s_2$ (MDJ) flow of Livermore & Jackson (2004) which is given by

$$u = u_0 K^{-1} \nabla \times \left( r^2 (1 - r^2) P_{l0}^2 (\cos \theta) \hat{r} \right)$$
$$+ u_0 K^{-1} \epsilon \nabla \times \nabla \times \left( r^2 (1 - r^2) P_{l0}^2 (\cos \theta) \hat{r} \right)$$

with $K^{-1} = \sqrt{0.909/572}$ and $\epsilon = 0.5 \sqrt{143}/1008$ such that the rms poloidal to toroidal energy ratio is 0.5, and the flow has an rms value of $u_0$, $P_{l0}$ and $\hat{r}$ defined in Appendix A.

Like Livermore & Jackson (2005) we consider the axisymmetric ($m = 0$) and equatorially antisymmetric magnetic field solution for a unit sphere ($r_0 = 1$) embedded in a vacuum. Table 2 shows the convergence in the eigenvalue with the largest real part as a function of truncation $n_{\text{max}}$ and $l_{\text{max}}$ for two magnetic Reynolds numbers $R_m = u_0 r_0 / \eta = 10$ and $R_m = 100$, together with the converged values given by Livermore & Jackson (2005) (LJ05). There is a difference of about one permille between their value and ours for $R_m = 10$.

#### 3.3. $\alpha^2$ and $\alpha^2 \Omega$-dynamos

We reproduced the critical dynamo numbers $R_\alpha$ further for the dipolar ($l = 1$) mode of an isotropic $\alpha^2$-dynamo with $\alpha_{rr} = \alpha_{\theta \theta} = \alpha_{\phi \phi} = R_\alpha \sin N \pi (r - r_1)$ and $N = 1.2$ in a spherical shell of inner and outer radius $r_i$ and $r_0$ with $r_0 - r_i = 1$ and $r_0$.

The numerical value of $R_\alpha$ is equal to the first zero of the spherical Bessel function $j_1$. 

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Table 1. Eigenvalues of the fundamental dipolar mode and the fifth and tenth overtones of the $\alpha^2$-sphere.

| $n_{\text{max}}$ | $\lambda_0$ | $\lambda_5$ | $\lambda_{10}$ |
|------------------|-------------|-------------|---------------|
| 4                | -0.0241     | -115.936    | -349.489      |
| 8                | -0.0034     | -115.110    | -349.489      |
| 12               | -0.0010     | -115.058    | -349.097      |
| 16               | -0.0004     | -115.047    | -349.050      |
| 20               | -0.0002     | -115.043    | -349.036      |
| 24               | -0.0001     | -115.041    | -349.031      |
| 28               | -0.0001     | -115.040    | -349.028      |
| 32               | 0.0000      | -115.040    | -349.026      |
| 36               | 0.0000      | -115.039    | -349.026      |

HvG93 0.0000 -115.04 -349.02

Table 2. Eigenvalues of the fundamental mode for the $\ell_1 s_2$ flow for two magnetic Reynolds numbers $R_m$.

| $n_{\text{max}}$ | $\lambda_0$ ($R_m = 10$) | $\lambda_0$ ($R_m = 100$) |
|------------------|---------------------------|---------------------------|
| 4                | -8.02911                  | -5.88075                  |
| 8                | -8.02625                  | -6.84495                  |
| 12               | -8.02626                  | -6.92613                  |
| 16               | -8.02627                  | -6.92859                  |
| 20               | -8.02627                  | -6.92866                  |
| 24               | -8.02627                  | -6.92869                  |

LJ05 -8.01600 -6.92885

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Table 3. Eigenvalues of the first and fifth modes of the $a^2\Omega$-dynamo (see Sect. 3.3).

| $n_{\text{max}}$ | $l_{\text{max}}$ | New method | Classical method |
|------------------|------------------|-------------|------------------|
|                  |                  | $\lambda_0$ (S) | $\lambda_0$ (A) | $\lambda_2$ (S) | $\lambda_2$ (A) |
| 8                | 8                | (16.572, 0.0) | (16.577, 0.0) | (16.572, 0.0) | (16.577, 0.0) |
| 12               | 12               | (16.053, 0.0) | (16.053, 0.0) | (16.053, 0.0) | (16.053, 0.0) |
| 16               | 16               | (16.052, 0.0) | (16.052, 0.0) | (16.052, 0.0) | (16.052, 0.0) |
| 20               | 20               | (16.053, 0.0) | (16.053, 0.0) | (16.053, 0.0) | (16.053, 0.0) |

$r_i/r_0 = 0.35$ and 0.8 surrounded by a vacuum and either an insulating or a conducting inner core, as reported in Table 2 of Schubert & Zhang (2001). With this test we treated in particular two different aspect ratios of a spherical shell (a thick and a thin one) and two different molecular diffusivities (insulating or conducting) of the inner core.

Finally we applied our method to an $a^2\Omega$-dynamo of Jiang & Wang (2006) who employ the classical eigenvalue treatment for the poloidal and toroidal scalars $P$ and $T$ expanded in spherical harmonics in the angular coordinates and in Chebychev polynomials in $r$-direction. We set $u_\theta = u_q = 0$, $u_\varphi = R_g r^2 \sin^2 \varphi$, $\alpha_{ij} = \beta_{ij} = 0$, except $\alpha_{rr} = \alpha_{\varphi\varphi} = R_g \sin \varphi (r - r_i)/(r_0 - r_i) \cos \varphi$ with $r_i = 0.5$, $r_0 = 1$, embedded in a vacuum inside and outside, and $R_g = \alpha_0 r_0/\eta = 10$ and $R_g = u_q r_0/\eta = 1000$. Some results obtained by the new and the classical methods are compiled in Table 3. Numbers in parentheses (., ., .) are the real and imaginary parts of complex eigenvalues. The real part denotes the growth rate, the imaginary part the frequency of the mode in units of $\eta/r_G^2$. The modes are axisymmetric ($m = 0$), the fundamental mode is monotonously growing and symmetric (indicated by $S$) with respect to the equator, and the fourth overtone is damped, oscillatory and antisymmetric (indicated by $A$). Modes with higher $m$ are more strongly damped. $n_{\text{max}}$ refers to the maximum radial number of the decay modes (spherical Bessel functions) for the new method and to the maximum degree of the Chebychev polynomials for the code of Jiang & Wang (2006), respectively. Since the modes have smaller length scales in latitudinal than in radial direction, higher values of $l$ than of $n$ are required for convergence. We find remarkably similar convergence of the eigenvalues for both methods. This also applies to modes with higher $m$. Of course we have also verified that the eigenfunctions obtained with the two methods are identical.

4.1. Benchmark dynamo

We examine a quasi-steady geodynamo model which has been used before as a numerical benchmark dynamo (Christensen et al. 2001, case 1). The governing parameters are Ekman number $E = 10^{-3}$, Rayleigh number $Ra = 100$, Prandtl number $Pr = 1$, and magnetic Prandtl number $Pm = 5$. The convection pattern is columnar with a natural 4-fold azimuthal symmetry and is stationary except for an azimuthal drift. The intensity of the fluid motion is characterised by a magnetic Reynolds number of $R_m = 40$, defined with a characteristic flow velocity, the thickness of the convecting shell, and the molecular magnetic diffusivity. The magnetic energy density exceeds the kinetic one by a factor of 20.

In Schrinner et al. (2007), the mean-field coefficients are derived from the numerical simulation. We solved the dynamo equation with these mean-field coefficients by the new method and obtained the eigenvalues and eigenfunctions. Since the coefficients are spatially variable to a considerable degree, converged solutions require high truncation levels in $n$ and $l$. The eigenvalues of the first two modes are shown in Table 4. Beyond $n_{\text{max}} \approx 20$ and $l_{\text{max}} \approx 20$, the eigensolution of the first mode does not change significantly and is displayed in Fig. 1. The convergence of the second mode requires a larger $l_{\text{max}}$ of about 32. The results for high values of $n_{\text{max}}$ may be affected by the spatial variation of the mean-field coefficients and would require more than 66 radial quadrature points to compute the matrix elements.

A comparison of Fig. 1 with its counterpart Fig. 10 of Schrinner et al. (2007) shows that the field of the antisymmetric fundamental mode resembles the field of an initial-value mean-field dynamo calculation remarkably well as it does the axisymmetric component of the direct numerical simulation. The mode here grows slightly with a rate around $\lambda_0 \approx 4.2\eta/L^2$, the field of the initial value calculation decays slightly with a rate of approximately $-0.25\eta/L^2$, while the solution of the direct numerical simulation is stationary. Here $L = r_0 - r_i = 1$ is the thickness of the spherical shell. The difference in these rates between the eigenvalue and initial value calculation comes from the higher numerical diffusivity of the latter at the chosen resolution of 33 radial and 80 latitudinal grid points. The difference is actually small, much less than one effective decay rate, because the relevant turbulent diffusivity, described by the $\beta$ coefficient with values up to $33\eta$, is much higher than the molecular one.

Besides the true physical eigenmodes, we find growing unphysical spurious eigenmodes. Their eigenvalues depend strongly on the resolution, and their eigenfunctions are highly structured. We attribute their appearance to a locally confined inappropriate parametrisation of the mean electromotive force by

\footnote{In Schrinner et al. (2007) the mean flow entered with a sign error into the initial value calculation, leading to a stronger decay of $-3.5\eta/L^2$. We apologise and correct this value here.}

4. Geodynamo models

Having proven that the new method works correctly and efficiently, we now apply it to determine the eigensolutions of the dynamo operator with mean-field coefficients obtained from self-consistent numerical simulations of the geodynamo. For a recent review of numerical geodynamo simulations, see Christensen & Wicht (2007). Schrinner et al. (2007) developed an efficient method of calculating all tensorial mean-field coefficients $\alpha$ and $\beta$ and compared the results of mean-field and direct numerical simulations of the geodynamo. We plan to use the eigenmodes of the dynamo equation to decompose the magnetic field of the numerical simulations and to determine the statistical properties of the mode coefficients (Hoyng 2009) to analyse the working of the geodynamo.
the mean-field coefficients $\alpha$ and $\beta$ (Schrinner et al. 2007). The spurious modes are present neither in the initial value calculation nor in the following example of a time-dependent dynamo, because of a higher numerical and molecular diffusivity, respectively.

4.2. A time-dependent dynamo in the columnar regime

The next example has stronger forcing with parameters $E = 10^{-4}$, $Ra = 334$, $Pr = 1$, and $Pm = 2$. The numerical simulation by Olson et al. (1999, case 2) shows a highly time-dependent, but still dominantly columnar convection characterised by a magnetic Reynolds number of $R_m \approx 88$. The magnetic energy exceeds the kinetic energy by a factor of three. The magnetic field has a strong axial dipole contribution. Although chaotically time-dependent, the velocity field is symmetric and the magnetic field antisymmetric with respect to the equatorial plane.

The mean-field coefficients are obtained as before by the test-field method of Schrinner et al. (2007). The coefficients are now of course also highly time-dependent. A time average yields coefficients that roughly resemble those for the benchmark dynamo, although there are differences in some profiles and amplitudes.

For the time-averaged dynamo operator the eigenvalues of the first two antisymmetric eigenmodes for various values of $n_{\text{max}}$ and $l_{\text{max}}$ are shown in Table 5. It seems that a value of $n_{\text{max}} \approx 16$ is sufficient for convergence, while $l_{\text{max}} \approx 32$ is needed. Figure 2 shows the eigenfunctions of these modes. The eigenvalues for $\lambda_1 = (-6.298, 0.0)$ and $\lambda_2 = (-28.712, \pm 5.364)$, values for $n_{\text{max}} = 16$ and $l_{\text{max}} = 32$, are symmetric with respect to the equator.

An initial-value, mean-field dynamo calculation with the same mean velocity and dynamo coefficients shows a slightly decaying solution with a decay rate of approximately $-5.99\eta/L^2$ which is to be compared with the eigenvalue $\lambda_0 = -3.87\eta/L^2$ of the fundamental mode. Again, the turbulent diffusivity exceeds the molecular one by a factor of up to 23 in this case. The difference in the decay rates is therefore much less than one effective decay rate. As for the benchmark dynamo, the profile of the antisymmetric fundamental mode is again remarkably similar to the solution of the initial value calculation and to the axisymmetric component of the direct numerical simulation.

A decomposition of the actual magnetic field of the simulation by Olson et al. (1999, case 2) in eigenfunctions of the time-averaged dynamo operator, i.e., $B(r, t) = \sum c_i(t)B_i(r)$, shows that the antisymmetric fundamental mode contributes to about 75 percent and, together with the first antisymmetric overtone (see Table 5 and Fig. 2), to about 85 percent of the total magnetic energy. The variability in time of the magnetic field of the

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**Table 4. Eigenvalues in units of $\eta/L^2$ of the first two eigenmodes of the benchmark dynamo.**

| $n_{\text{max}}$ | $l_{\text{max}}$ | $a_0$ (Å) | $a_1$ (Å) |
|------------------|------------------|-----------|-----------|
| 12               | 12               | +4.960    | -8.605    |
| 16               | 16               | +4.235    | -8.110    |
| 20               | 20               | +4.180    | -8.362    |
| 24               | 24               | +4.195    | -6.620    |
| 28               | 28               | +4.255    | -7.275    |
| 32               | 32               | +4.275    | -6.015    |
| 32               | 32               | +4.382    | -6.777    |

**Notes.** The symmetry with respect to the equator is marked $A$ for antisymmetric and $S$ for symmetric. The eigensolution marked bold is displayed in Fig. 1.

**Table 5. Eigenvalues in units of $\eta/L^2$ of the first two antisymmetric eigenmodes of the temporally averaged dynamo operator obtained from the time-dependent dynamo (case 2, Sect. 4.2).**

| $n_{\text{max}}$ | $l_{\text{max}}$ | $a_0$ | $a_2$ |
|------------------|------------------|------|------|
| 12               | 12               | (-4.520, 0.0) | (-35.046, ±10.118) |
| 16               | 16               | (-4.278, 0.0) | (-35.090, ±10.256) |
| 20               | 20               | (-4.112, 0.0) | (-34.578, ±10.096) |
| 24               | 24               | (-3.930, 0.0) | (-34.770, ±9.874)  |
| 28               | 28               | (-3.880, 0.0) | (-34.766, ±10.236) |
| 32               | 32               | (-3.868, 0.0) | (-34.804, ±10.318) |
| 32               | 32               | (-3.874, 0.0) | (-34.830, ±10.310) |

**Notes.** The eigensolutions marked bold are displayed in Fig. 2.

**Fig. 1. Magnetic field structure of the fundamental antisymmetric eigenmode for the benchmark dynamo. Compare with Fig. 10 of Schrinner et al. (2007). For each plot the grey scale is separately adjusted to its maximum modulus with white as negative and black as positive. The contour lines correspond to ±0.1, ±0.3, ±0.5, ±0.7, and ±0.9 of the maximum modulus.**

**Fig. 2. Radial component of the first two antisymmetric eigenmodes for the case2 dynamo. Left: fundamental mode; middle: real part of the first overtone; right: imaginary part of the first overtone. Grey scales and contours as in Fig. 1.**
direct numerical simulation is reflected in the variability of the expansion coefficients. More details are presented in Schrinner et al. (2009).

5. Conclusions and outlook

We presented a new method for computing the eigenvalues and eigenfunctions of the induction and the dynamo equation. The method is based on the biorthogonality of the adjoint electric current and the vector potential with an inner product defined by a volume integral over the fluid domain. The advantage of the method is that the velocity and dynamo coefficients do not have to be differentiated. The method is therefore well-suited for spatially strongly variable dynamo coefficients.

We tested the new method against the classical treatment and proved that it works correctly and efficiently. We applied it to two cases with dynamo coefficients derived from direct numerical simulations of the geodynamo. The obtained dynamo eigenmodes are promising candidates for decomposing the magnetic field of the numerical simulations and for analysing the statistical properties of the mode coefficients as proposed by Hoyng (2009).

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Appendix A: Free magnetic decay modes in a sphere or spherical shell embedded in vacuum

We decompose the magnetic field in its poloidal and toroidal components

\[ B = \nabla \times \nabla \times \hat{P} \hat{r} + \nabla \times \hat{T} \hat{r} \]  

(A.1)

with defining scalars \( P(r, \theta, \varphi) \) and \( T(r, \theta, \varphi) \) and unit vector in radial direction \( \hat{r} = (1, 0, 0) \) in spherical coordinates \( (r, \theta, \varphi) \).

The equation of free magnetic decay

\[ \frac{\partial B}{\partial t} = -\eta \nabla \times \nabla \times B \]  

(A.2)

with constant magnetic diffusivity \( \eta \) then reads

\[ \frac{\partial}{\partial t} (P, T) = \eta \left( \Delta_H + \frac{\partial^2}{\partial r^2} \right) (P, T) \]  

(A.3)

where \( \Delta_H \) is the horizontal Laplacian

\[ \Delta_H = \frac{1}{r^2 \sin^2 \theta} \frac{\partial}{\partial \theta} (r^2 \sin \theta \frac{\partial}{\partial \theta}) + \frac{1}{r^2 \sin^2 \theta} \frac{\partial^2}{\partial \varphi^2}. \]  

(A.4)

The solutions are the free magnetic decay modes

\[ P_{mn} = \exp \left( i \lambda_{lm} r \right) x f_l(p_m x) Y_{lm}(\theta, \varphi), \]  

(A.5)

\[ T_{mn} = \exp \left( i \lambda_{lm} r \right) x g_l(t_m x) Y_{lm}(\theta, \varphi), \]  

(A.6)

where \( x = r/r_0 \) where \( r_0 \) is the radius of the sphere. The growth rates are given by

\[ \lambda_{lm}^0 = -\eta \frac{p_m^2}{p_0^2} r_0^2 \]  

and \( \lambda_{lm}^T = -\eta \frac{t_m^2}{t_0^2} r_0^2 \)

(A.7)

and are independent of the azimuthal degree \( m \). The constants \( p_m \) and \( t_m \) are

\[ p_m = j_{l-1,n} \]  

and \( t_m = j_{l,n} \)

(A.8)

where \( j_{l,n} \) is the \( n \)-th zero of \( j_l \). The \( Y_{lm}^n \) are the spherical harmonics and normalised to unity by taking

\[ Y_{lm}^n(\theta, \varphi) = \frac{4\pi}{2l+1} \frac{(l+m)!}{(l-m)!(l+1)!} P_l^m(\cos \theta) e^{im\varphi} \]  

(A.9)

using Ferrer’s definition of the Legendre functions of first kind \( P_l^m \) with degree \( l \) and order \( m \).

For a sphere embedded in vacuum the radial functions are given by

\[ x f_l(p_m x) = \begin{cases} a_{lm} j_l(p_m x) & 0 \leq x \leq 1 \\ a_{lm} j_l(p_m x)^{-1} & x \geq 1 \end{cases} \]  

(A.10)

\[ x g_l(t_m x) = \begin{cases} b_{lm} j_l(t_m x) & 0 \leq x \leq 1 \\ 0 & x \geq 1 \end{cases} \]  

(A.11)
with the spherical Bessel functions of first kind \( j_l \). This ensures regularity in the origin of the sphere, vanishing toroidal component at its outer boundary and smooth transition of the poloidal component to a potential field in the vacuum outside.

For a spherical shell with inner radius \( r_i \) (\( x_i = r_i / r_0 \)) and outer radius \( r_0 \) (\( x_0 = 1 \)) embedded in vacuum the radial functions inside the shell are given by

\[
f_l(p \ln x) = j_l(p \ln x) - y_l(p \ln x) j_{l+1}(p \ln x_i) / y_{l+1}(p \ln x_i) \quad (A.12)
\]

and

\[
g_l(t \ln x) = j_l(t \ln x) - y_l(t \ln x) j_{l+1}(t \ln x_i) / y_{l+1}(t \ln x_i), \quad (A.13)
\]

and the constants in the arguments are the roots of

\[
j_{l+1}(p \ln x_i) y_{l+1}(p \ln x) - j_l(p \ln x) y_{l+1}(p \ln x_i) = 0
\]

for \( p \ln \) and of

\[
j_l(t \ln x_i) y_l(t \ln x) - j_{l-1}(t \ln x) y_{l+1}(t \ln x_i) = 0
\]

for \( t \ln \). Here \( y_l \) are the spherical Bessel functions of second kind.

The magnetic field of the decay modes \( \mathbf{B}_i \) is obtained by inserting the spatial parts of the defining scalars \( P_{lmn} \) and \( T_{lmn} \), respectively, into (A.1). Here we have comprised the three indices into one. The decay modes are self-adjoint on \( V + E \), so that the adjoint functions are obtained simply by complex conjugation:

\[
\hat{B}_k = B_k^* \quad \text{and likewise} \quad \hat{J}_k = J_k^*.
\]

Normalisation on \( V + E \), i.e., \((B_k^*, B_i)_{V+E} = \delta_{ki}\), is thus straightforward. For a unit sphere the radial functions are normalised to unity by scaling the \( f_l \) with

\[
a_{ln} = \left( \frac{1}{2r_0} l(l+1) j_{l-1,n}^2 f_l(j_{l-1,n}) \right)^{-1/2} \quad (A.16)
\]

and the \( g_l \) with

\[
b_{ln} = \left( \frac{r_0}{2} l(l+1) j_{l+1,n}^2 (j_{l,n}) \right)^{-1/2} \quad . \quad (A.17)
\]

For a spherical shell the normalisation constants are more lengthy expressions, which we suppress here.

The free magnetic decay modes form a complete and orthogonal set of functions, and they obey the boundary conditions of the magnetic field between the dynamo volume \( V \) and the exterior vacuum \( E \).

We mention for completeness that the poloidal decay modes are not self-adjoint on \( V \), i.e., \((B_k^*, B_i)_V \neq \delta_{ki}\). If we like to work with an inner product defined on \( V \), the adjoint functions \( \hat{B}_k \) can be constructed by requiring \((B_k^*, B_i)_V = \delta_{ki}\), similar to the one described in Sect. 2.5.