

9 Computational aspects of astrophysical MHD and turbulence

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The advantages of high-order finite difference scheme for astrophysical MHD and turbulence simulations are highlighted. A number of one-dimensional test cases are presented ranging from various shock tests to Parker-type wind solutions. Applications to magnetized accretion discs and their associated outflows are discussed. Particular emphasis is placed on the possibility of dynamo action in three-dimensional turbulent convection and shear flows, which is relevant to stars and astrophysical discs. The generation of large scale fields is discussed in terms of an inverse magnetic cascade and the consequences imposed by magnetic helicity conservation are reviewed with particular emphasis on the issue of α -quenching.

9.1. Introduction

Over the past 20 years multidimensional astrophysical gas simulations have become a primary tool to understand the formation, evolution, and the final fate of stars, galaxies, and their surrounding medium. The assumption that those processes happen smoothly and in a non-turbulent manner can at best be regarded as a first approximation. This is evidenced by the ever improving quality of direct imaging techniques, e.g. using space telescopes. At the same time not only have computers become large enough to run three-dimensional simulations with relatively little effort, there have also been substantial improvements in the algorithms that are used. In fact, there is now a vast literature on numerical astrophysics. An excellent book was published recently by LeVeque *et al.* (1998) where both numerical methods and astrophysical applications were discussed in great detail. Most of the applications focused however on rather more “violent” processes such as supersonic jets, supernova explosions, core collapse, and on radiative transfer problems, while hydromagnetic phenomena and turbulence problems were only touched upon briefly. Meanwhile, hydromagnetic turbulence simulations have become crucial for understanding viscous dissipation in accretion discs (Hawley *et al.*, 1995), and for understanding magnetic field generation by dynamo action in discs (Brandenburg *et al.*, 1995, 1996a; Hawley *et al.*, 1996; Stone *et al.*, 1996), stars (Nordlund *et al.*, 1992; Brandenburg *et al.*, 1996b), and planets (Glatzmaier and Roberts, 1995, 1996).

Much of the present day astrophysical hydrodynamic work is based on the ZEUS code, which has been documented in great detail and described with a number of test cases in a series of papers by Stone and Norman (1992a,b). The main advantage is its flexibility in dealing with arbitrary orthogonal coordinates which makes the code applicable to a wide variety of astrophysical systems. The code, which is freely available on the net, uses artificial viscosity for stability and shock capturing, and is based on an operator split method with

2nd-order finite differences on a staggered mesh. Another approach used predominantly in turbulence research are spectral methods (e.g. Canuto *et al.*, 1988), which have the advantage of possessing high accuracy. Although these methods are most suitable for incompressible flows (imposing the solenoidality condition is then straightforward), they have also been applied to compressible flows (e.g. Passot and Pouquet, 1987). As a compromise one may resort to high-order finite difference methods, which have the advantage of being easy to implement and yet have high accuracy. Compact methods (e.g. Lele, 1992) are a special variety of high-order finite difference methods, but the truncation error is smaller than for an explicit scheme of the same order. For example, compact schemes have been used by Nordlund and Stein (1990) in simulations of solar convection (Stein and Nordlund, 1989, 1998) and convective dynamos (Nordlund *et al.*, 1992; Brandenburg *et al.*, 1996b).

The use of compact methods involves solving tridiagonal matrix equations, making this method essentially non-local in that all points are now coupled at once. This is problematic for massively parallel computations, which is why Nordlund and Galsgaard (1995, see also Nordlund *et al.*, 1994) began to use explicit high-order schemes for their work on coronal heating by reconnection (Galsgaard and Nordlund, 1996, 1997a,b). In their code the equations are solved in a semi-conservative fashion using a staggered mesh. This code was also used by Padoan *et al.*, (1997) and Padoan and Nordlund (1999) in models of isothermal interstellar turbulence in molecular clouds, and by Rögnvaldsson *et al.* (2001) in simulations of cooling flows and galaxy formation.

A somewhat different code was used by Brandenburg (1999) and Bigazzi (1999) in simulations of the inverse magnetic cascade, by Kerr and Brandenburg (1999) in a work on the possibility of a singularity of the non-resistive and inviscid MHD equations, and by Sanchez-Salcedo and Brandenburg (1999, 2001) in simulations of dynamical friction. A two-dimensional version of the code modelling outflows from magnetized accretion discs has been described by Brandenburg *et al.* (2000). This code uses 6th-order explicit finite differences in space and 3rd-order Runge–Kutta timestepping. It employs central finite differences, so the extra cost of recentering a large number of variables between staggered meshes each timestep is avoided.

Apart from high numerical accuracy, another important requirement for astrophysical gas simulations is the capability to deal with a large dynamical range in density and temperature. This requirement favors the use of non-conservative schemes, because then logarithmic variables can be used which vary much less than linear density and energy density per unit volume. Solving the non-conservative form of the equations can be more accurate than solving the conservative form. The conservation properties can then be used as an indicator for the overall accuracy.

In this chapter we concentrate on numerical astrophysical turbulence aspects starting with a discussion of different numerical methods and a description of the results of various numerical test problems. This is a good way of assessing the quality of a numerical scheme and of comparing with other methods; see Stone and Norman (1992a,b) for a series of tests using the ZEUS code. After that we discuss particular astrophysical applications including stellar convection, accretion disc turbulence and associated outflows, as well as the generation of magnetic fields (small and large scale) from turbulence in various astrophysical settings.

9.2. The Navier–Stokes equations

The discussion of magnetic fields will be postponed until later, because the inclusion of the Lorentz force in the momentum equation is straightforward. We begin by writing down

the Navier–Stokes equations in non-conservative form and rewrite them such that the main thermodynamical variables are entropy and either logarithmic density or potential enthalpy. These variables have the advantage of varying spatially much less than, e.g. linear pressure and density.

The primitive form of the continuity equation is

$$\frac{\partial \rho}{\partial t} = -\nabla \cdot (\rho \mathbf{u}), \quad (9.1)$$

which means that the local change of density is given by the divergence of the mass flux at that point. The Navier–Stokes equation can be written as

$$\rho \frac{D\mathbf{u}}{Dt} = -\nabla p - \rho \nabla \Phi + \mathbf{F} + \nabla \cdot \boldsymbol{\tau}, \quad (9.2)$$

where $D/Dt = \partial/\partial t + \mathbf{u} \cdot \nabla$ is the advective derivative, p is the pressure, Φ is the gravitational potential, \mathbf{F} is a body force (e.g. the Lorentz force), and $\boldsymbol{\tau}$ is the stress tensor.

The Navier–Stokes equation is here written in terms of forces per unit volume. As argued above, if the density contrast is large it is advantageous to write it in terms of forces per unit mass and to divide by ρ . Before we can replace p and ρ by entropy and logarithmic density or potential entropy we first have to define some thermodynamic quantities.

Internal energy, e , and specific enthalpy, h are related to each other by

$$h = e + pv, \quad (9.3)$$

where $v = 1/\rho$ is the specific volume and ρ the density. The specific entropy is defined by

$$T ds = de + p dv, \quad (9.4)$$

where T is temperature. The specific heats at constant pressure and constant volume are defined as $c_p = dh/dT|_p$ and $c_v = de/dT|_v$, their ratio is $\gamma = c_p/c_v$, and their difference is $\mathcal{R}/\mu = c_p - c_v$, where \mathcal{R} is the universal gas constant and μ the specific molecular weight.

In the following we assume c_p and c_v to be constant for all processes considered. Ionization and recombination processes are therefore ignored here, although this is not a major obstacle; see, e.g. simulations of Nordlund (1982, 1985), Steffen *et al.* (1989), Stein and Nordlund (1989, 1998), Rast *et al.* (1993), and Rast and Toomre (1993a,b) where realistic equations of state have been used.

We now assume that c_p and c_v are constant, so internal energy and specific enthalpy are given by

$$h = c_p T \quad \text{and} \quad e = c_v T. \quad (9.5)$$

This allows us then to write the specific entropy (up to an additive constant) as

$$s = c_v \ln p - c_p \ln \rho. \quad (9.6)$$

The pressure gradient term in the momentum equation can then be written as

$$\frac{1}{\rho} \nabla p = \frac{p}{\rho} \nabla \ln p = \frac{\gamma p}{\rho} (\nabla \ln \rho + \nabla s/c_p) = c_s^2 (\nabla \ln \rho + \nabla s/c_p), \quad (9.7)$$

where we have used

$$c_s^2 = \gamma p / \rho = \frac{\gamma p_0}{\rho_0} \exp[(\gamma - 1) \ln(\rho/\rho_0) + \gamma s/c_p] = c_{s0}^2 \left(\frac{\rho}{\rho_0} \right)^{\gamma-1} \exp(\gamma s/c_p), \quad (9.8)$$

where c_s is the adiabatic sound speed, and $c_{s0}^2 = \gamma p_0 / \rho_0$.

With these preparations the evolution of velocity \mathbf{u} , logarithmic density $\ln \rho$, and specific entropy s can be expressed as follows:

$$\frac{D\mathbf{u}}{Dt} = -c_s^2 (\nabla \ln \rho + \nabla s/c_p) - \nabla \Phi + \mathbf{f} + \frac{1}{\rho} \nabla \cdot (2\nu\rho\mathbf{S}), \quad (9.9)$$

$$\frac{D \ln \rho}{Dt} = -\nabla \cdot \mathbf{u}, \quad (9.10)$$

$$T \frac{Ds}{Dt} = 2\nu\mathbf{S}^2 + \Gamma - \rho\Lambda, \quad (9.11)$$

where $\mathbf{f} = \mathbf{F}/\rho$ is the body force per unit mass, Γ and Λ are heating and cooling functions, ν kinematic viscosity and \mathbf{S} is the (traceless) strain tensor with the components

$$S_{ij} = \frac{1}{2} (u_{i,j} + u_{j,i} - \frac{2}{3} \delta_{ij} u_{k,k}). \quad (9.12)$$

In the presence of an additional kinematic bulk viscosity, ζ , the term $2\nu S_{ij}$ under the divergence in (9.9) would need to be replaced by $2\nu S_{ij} + \zeta \delta_{ij} \nabla \cdot \mathbf{u}$, and the viscous heating term, $2\nu\mathbf{S}^2$, in (9.11) would need to be replaced by $2\nu\mathbf{S}^2 + \zeta(\nabla \cdot \mathbf{u})^2$.

Instead of using ρ as a dependent variable one can also use the specific enthalpy h , which allows us to write the pressure gradient as

$$-\frac{1}{\rho} \nabla p = -\nabla h + T \nabla s. \quad (9.13)$$

This formulation is particularly useful if the entropy is nearly constant (or if the gas is barotropic, i.e. $p = p(\rho)$) and if there is a gravitational potential Φ , so that the potential enthalpy $H \equiv h + \Phi$ can be used as dependent variable. In order to express (9.10) in terms of h we write down the total differential of the specific entropy,

$$ds/c_p = \frac{1}{\gamma} d \ln p - d \ln \rho = \frac{1}{\gamma} d \ln h - \left(1 - \frac{1}{\gamma} \right) d \ln \rho, \quad (9.14)$$

so

$$\frac{D \ln h}{Dt} = \gamma \frac{Ds/c_p}{Dt} + (\gamma - 1) \frac{D \ln \rho}{Dt}. \quad (9.15)$$

Furthermore, $T \nabla s = h \nabla s / c_p$, and so the final set of equations is

$$\frac{D\mathbf{u}}{Dt} = -\nabla H + h \nabla s / c_p + \mathbf{f} + \nabla \cdot (2\nu\rho\mathbf{S}), \quad (9.16)$$

$$\frac{Ds/c_p}{Dt} = \frac{1}{h} (2\nu\mathbf{S}^2 + \Gamma - \rho\Lambda), \quad (9.17)$$

$$\frac{DH}{Dt} = \mathbf{u} \cdot \nabla \Phi + \gamma h \frac{Ds/c_p}{Dt} - c_s^2 \nabla \cdot \mathbf{u}, \quad (9.18)$$

where we have absorbed Φ in the potential enthalpy $H = h + \Phi$. In this formulation the density can be recovered as

$$\rho = \rho_0 \left[\left(1 - \frac{1}{\gamma} \right) \left(\frac{h}{c_{s0}^2} \right) e^{-\gamma s / c_p} \right]^{1/(\gamma-1)} \quad (9.19)$$

(in dimensional form) or, for $\gamma = 5/3$ and in non-dimensional form (where $\rho_0 = p_0 = c_p = 1$),

$$\rho = (0.4h)^{1.5} e^{-2.5s}. \quad (9.20)$$

We shall use either of the two sets of the equations, (9.9)–(9.11) or (9.16)–(9.18), in some of the following sections, especially in connection with shock tests and stellar wind problems. In these cases the gravity potential Φ is important and it turns out that the potential enthalpy $H \equiv h + \Phi$ varies only very little near the central object even though Φ itself tends to become singular.

The heating and cooling terms (Γ and Λ) are important, e.g. in the case of interstellar turbulence which is driven primarily by supernova explosions which inject a certain amount of thermal energy ($\int \rho \Gamma dV$) with each supernova explosion. MHD turbulence simulations of this type were performed recently by Korpi *et al.* (1999). At the same time there is cooling through various processes (e.g. bremsstrahlung at high temperatures) which transports energy either non-locally via a cooling term $\Lambda(T)$, or locally via thermal conduction or radiative diffusion. In the radiative diffusion approximation we express Λ as $-\rho^2 \Lambda = \nabla \cdot K \nabla T$, where K is the radiative conductivity which is in general a function of temperature and density. The radiative diffusivity (which has the same dimensions as the kinematic viscosity ν) is given by $\chi = K / (\rho c_p)$, so

$$-\rho \Lambda / (c_p T) = \frac{1}{\rho c_p T} \nabla \cdot \rho \chi c_p \nabla T. \quad (9.21)$$

Since we shall use a non-conservative scheme with centered finite differences it is important to isolate second derivative terms, so

$$-\rho \Lambda / h = \chi (\nabla^2 \ln T + \nabla \ln p \cdot \nabla \ln T), \quad (9.22)$$

where we have assumed for simplicity that χ is constant. In terms of s/c_p and $\ln \rho$ we have

$$-\rho \Lambda / h = \chi \gamma [\nabla^2 s / c_p + \nabla_{ad} \nabla^2 \ln \rho + \gamma (\nabla s / c_p + \nabla \ln \rho) \cdot (\nabla s / c_p + \nabla_{ad} \nabla \ln \rho)], \quad (9.23)$$

where $\nabla_{\text{ad}} = 1 - 1/\gamma$ is a commonly used abbreviation in stellar astrophysics. For $\gamma = 5/3$ we have $\nabla_{\text{ad}} = 2/5 = 0.4$. We shall use (9.23) later in connection with shock and wind calculations. However, we begin by discussing first a suitable numerical scheme which will be used in most of the cases presented below.

9.3. The advantage of higher-order derivative schemes

Spectral methods are commonly used in almost all studies of ordinary (usually incompressible) turbulence. The use of this method is justified mainly by the high numerical accuracy of spectral schemes. Alternatively, one may use high-order finite differences that are faster to compute and that can possess almost spectral accuracy. Nordlund and Stein (1990) and Brandenburg *et al.* (1995) use high-order finite difference methods, e.g. 4th- and 6th-order compact schemes (Lele 1992).¹

In this section we demonstrate, using simple test problems, some of the advantages of high-order schemes. We begin by defining various schemes including their truncation errors and their high wavenumber characteristics. We consider centered finite differences of 2nd-, 4th-, 6th-, 8th-, and 10th-order, which are given respectively by the formulae

$$f'_i = (-f_{i-1} + f_{i+1})/(2\delta x), \quad (9.24)$$

$$f'_i = (f_{i-2} - 8f_{i-1} + 8f_{i+1} - f_{i+2})/(12\delta x), \quad (9.25)$$

$$f'_i = (-f_{i-3} + 9f_{i-2} - 45f_{i-1} + 45f_{i+1} - 9f_{i+2} + f_{i+3})/(60\delta x), \quad (9.26)$$

$$\begin{aligned} f'_i = & (3f_{i-4} - 32f_{i-3} + 168f_{i-2} - 672f_{i-1} + 672f_{i+1} \\ & - 168f_{i+2} + 32f_{i+3} - 3f_{i+4})/(840\delta x), \end{aligned} \quad (9.27)$$

$$\begin{aligned} f'_i = & (-2f_{i-5} + 25f_{i-4} - 150f_{i-3} + 600f_{i-2} - 2100f_{i-1} \\ & + 2100f_{i+1} - \dots)/(2520\delta x), \end{aligned} \quad (9.28)$$

for the first derivative, and

$$f''_i = (f_{i-1} - 2f_i + f_{i+1})/(\delta x^2), \quad (9.29)$$

$$f''_i = (-f_{i-2} + 16f_{i-1} - 30f_i + 16f_{i+1} - f_{i+2})/(12\delta x^2), \quad (9.30)$$

$$\begin{aligned} f''_i = & (2f_{i-3} - 27f_{i-2} + 270f_{i-1} - 490f_i \\ & + 270f_{i+1} - 27f_{i+2} + 2f_{i+3})/(180\delta x^2), \end{aligned} \quad (9.31)$$

$$\begin{aligned} f''_i = & (-9f_{i-4} + 128f_{i-3} - 1008f_{i-2} + 8064f_{i-1} - 14350f_i \\ & + 8064f_{i+1} - 1008f_{i+2} + \dots)/(5040\delta x^2), \end{aligned} \quad (9.32)$$

$$\begin{aligned} f''_i = & (8f_{i-5} - 125f_{i-4} + 1000f_{i-3} - 6000f_{i-2} + 42000f_{i-1} \\ & - 73766f_i + 42000f_{i+1} - \dots)/(25200\delta x^2), \end{aligned} \quad (9.33)$$

for the second derivative. The expressions for one-sided and semi-one-sided finite difference formulae are given in Appendix A.

9.3.1. High wavenumber characteristics

The chief advantage of high-order schemes is their high fidelity at high wavenumber. Suppose we differentiate the function $\sin kx$, we are supposed to get $k \cos kx$, but when k is close to

the Nyquist frequency, $k_{\text{Ny}} \equiv \pi/\delta x$, where δx is the mesh spacing, numerical schemes yield *effective* wavenumbers, k_{eff} , that can be significantly less than the actual wavenumber k . Here we calculate k_{eff} from

$$(\cos kx)'_{\text{num}} = -k_{\text{eff}} \sin kx. \quad (9.34)$$

When $k = k_{\text{Ny}}$, every centered difference scheme will give $k_{\text{eff}} = 0$, because then the function values of $\cos kx$ are just $-1, +1, -1, \dots$, so the function values on the left and the right are the same, and the difference that enters the scheme gives therefore zero.

It is useful to mention at this point that for a staggered mesh, where the first derivative is evaluated *between* mesh points, the value of the first derivative remains finite at the Nyquist frequency, provided one does not need to remesh back to the original mesh. Especially in the context with magnetic fields, however, remeshing needs to be done quite frequently, which therefore diminishes the advantage of a staggered mesh.

In Fig. 9.1 we plot effective wavenumbers for different schemes. Apart from the different *explicit* finite difference schemes given above, we also consider a *compact* scheme of sixth order, which can be written in the form

$$\frac{1}{3}f'_{i-1} + f'_i + \frac{1}{3}f'_{i+1} = (f_{i-2} - 28f_{i-1} + 28f_{i+1} - f_{i+2})/(36\delta x), \quad (9.35)$$

for the first derivative, and

$$\frac{2}{11}f''_{i-1} + f''_i + \frac{2}{11}f''_{i+1} = (3f_{i-2} + 48f_{i-1} - 102f_i + 48f_{i+1} + 3f_{i+2})/(44\delta x^2), \quad (9.36)$$

for the second derivative. As we have already mentioned in the introduction, this scheme involves obviously solving tridiagonal matrix equations and is therefore effectively non-local.

In the second panel of Fig. 9.1 we have plotted effective wavenumbers for second derivatives, which were calculated as

$$(\cos kx)''_{\text{num}} = -k_{\text{eff}}^2 \cos kx. \quad (9.37)$$

Of particular interest is the behavior of the second derivative at the Nyquist frequency, because that is relevant for damping zig-zag modes. For a 2nd-order finite difference scheme k_{eff}^2 is only 4, which is less than half the theoretical value of $\pi^2 = 9.87$. For 4th-, 6th-, and 10th-order schemes this value is respectively 5.33, 6.04, and 6.83. The last value is almost the same as for the 6th-order compact scheme, which is 6.86. Significantly stronger damping at the Nyquist frequency can be obtained by using hyperviscosity, which Nordlund and Galsgaard (1995) treat as a quenching factor that diminishes the value of the second derivative for wavenumbers that are small compared with the Nyquist frequency. Accurate high-order second derivatives (with no quenching factors) are important when calculating the current \mathbf{J} in the Lorentz force $\mathbf{J} \times \mathbf{B}$ from a vector potential \mathbf{A} using $-\mu_0 \mathbf{J} = \nabla^2 \mathbf{A} - \nabla \nabla \cdot \mathbf{A}$. This will be important in the MHD calculations presented below.

9.3.2. The truncation error

One can express f_{i-1} , f_{i+1} , etc., in terms of the derivatives of f at point i , so

$$f_{i-1} = f_i - \delta x f'_i + \frac{1}{2} \delta x^2 f''_i - \frac{1}{6} \delta x^3 f'''_i + \dots \quad (9.38)$$

$$f_{i+1} = f_i + \delta x f'_i + \frac{1}{2} \delta x^2 f''_i + \frac{1}{6} \delta x^3 f'''_i + \dots \quad (9.39)$$

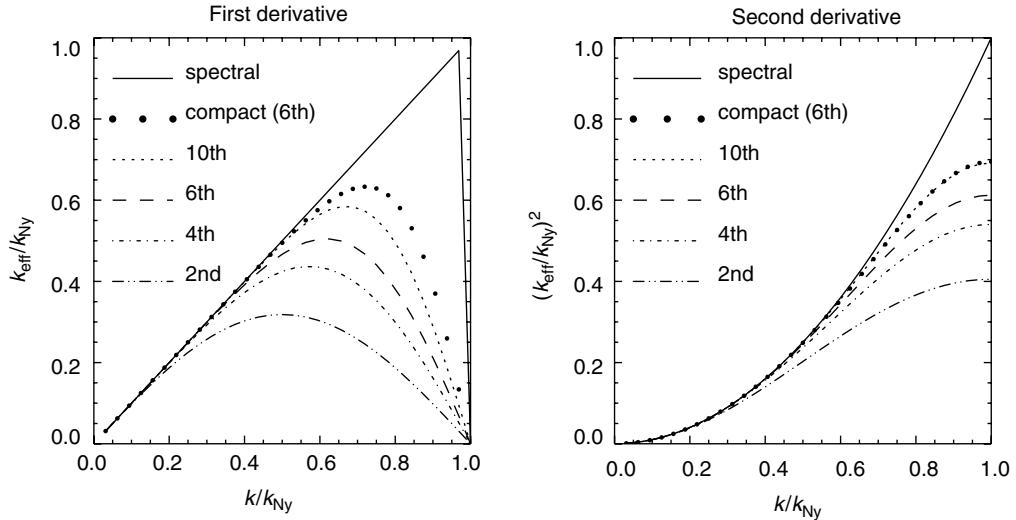


Figure 9.1 Effective wavenumbers for first and second derivatives using different schemes. Note that for the second derivatives the 6th-order compact scheme is almost equivalent to the 10th-order explicit scheme. For the first derivative the 6th-order compact scheme is still superior to the 10th-order explicit scheme.

Inserting this into the finite difference expressions yields for the 2nd-order formula

$$(f'_i)_{\text{2nd}} \equiv (f_{i+1} - f_{i-1})/(2\delta x) = f'_i + \frac{1}{6}\delta x^2 f'''_i. \quad (9.40)$$

The error scales quadratically with the mesh size, which is why the method is called second order. The truncation error is proportional to the third derivative of the function. Because this is an odd derivative it corresponds to a dispersive (as opposed to diffusive) error. Schemes that are only first order (or of any odd order) have diffusive errors, and it is this what is sometimes referred to as *numerical diffusivity*, which is not to be confused with artificial diffusivity that is sometimes used for stability and shock capturing. For the other schemes given in (9.25)–(9.28) the truncation errors are

$$(f'_i)_{\text{4th}} = f'_i + 3 \times 10^{-2} \delta x^4 f_i^{(\text{v})}, \quad (9.41)$$

$$(f'_i)_{\text{6th}} = f'_i + 7 \times 10^{-3} \delta x^6 f_i^{(\text{vii})}, \quad (9.42)$$

$$(f'_i)_{\text{10th}} = f'_i + 3 \times 10^{-4} \delta x^{10} f_i^{(\text{x})}. \quad (9.43)$$

For the 6th-order compact scheme the error scales like for the 6th-order explicit scheme, but the coefficient in front of the truncation error is about 10 times smaller, so

$$(f'_i)_{\text{6th}}^{\text{compact}} = f'_i + 5.1 \times 10^{-4} \delta x^6 f_i^{(\text{vii})}. \quad (9.44)$$

For the second derivatives we have

$$(f_i'')_{2\text{nd}} = f_i'' + 8 \times 10^{-2} \delta x^2 f_i^{(\text{iv})}, \quad (9.45)$$

$$(f_i'')_{4\text{th}} = f_i'' - 1 \times 10^{-2} \delta x^4 f_i^{(\text{vi})}, \quad (9.46)$$

$$(f_i'')_{6\text{th}} = f_i'' + 2 \times 10^{-3} \delta x^6 f_i^{(\text{viii})}, \quad (9.47)$$

$$(f_i'')_{10\text{th}} = f_i'' - 5 \times 10^{-5} \delta x^{10} f_i^{(\text{xii})}. \quad (9.48)$$

Again, for the 6th-order compact scheme the scaling is the same as for the 6th-order explicit scheme, but the coefficient in front of the truncation error is about five times less, so

$$(f_i'')_{6\text{th}}^{\text{compact}} = f_i'' + 3.2 \times 10^{-4} \delta x^6 f_i^{(\text{viii})}. \quad (9.49)$$

This information about the accuracy of schemes would obviously be of little use if the various schemes did not perform well when applied to real problems. For this reason we now begin by carrying out various tests, including advection and shock tests.

9.3.3. Advection tests

As a first test we compare the various schemes by performing inviscid advection tests and solve the equation $Df/Dt = 0$, i.e.

$$\dot{f} = -uf', \quad (9.50)$$

on a periodic mesh. It is advantageous to use a relatively small number of meshpoints (here we use $N_x = 8$ meshpoints), because that way we see deficiencies most clearly. This case is actually also relevant to real applications, because in practice one will always have small scale structures that are just barely resolved.

After some time an initially sinusoidal signal will suffer a change in amplitude and phase. We have calculated the amplitude and phase errors for schemes of different spatial order. For the time integration we use high-order Runge–Kutta methods of third- or fourth-order, RK3 and RK4, respectively. In most cases considered below we use the RK3 scheme that allows reasonable use of storage. It can be written in three steps (Rogallo, 1981)

$$\begin{aligned} \text{1st step: } & f = f + \gamma_1 \delta t \dot{f}, & g = f + \zeta_1 \delta t \dot{f}, \\ \text{2nd step: } & f = g + \gamma_2 \delta t \dot{f}, & g = f + \zeta_2 \delta t \dot{f}, \\ \text{3rd step: } & f = g + \gamma_3 \delta t \dot{f}, \end{aligned} \quad (9.51)$$

where

$$\gamma_1 = \frac{8}{15}, \quad \gamma_2 = \frac{5}{12}, \quad \gamma_3 = \frac{3}{4}, \quad \zeta_1 = -\frac{17}{60}, \quad \zeta_2 = -\frac{5}{12}, \quad (9.52)$$

where, f and g always refer to the current value (so the same space in memory can be used), but \dot{f} is evaluated only once at the beginning of each of the three steps at $t = t_0$, $t_{1/3} = t_0 + \gamma_1 \delta t \approx t_0 + 0.5333 \delta t$, and at $t_{2/3} = t_0 + (\gamma_1 + \zeta_1 + \gamma_2) \delta t = t_0 + (2/3) \delta t$. Even more memory-effective are the so-called $2N$ -schemes that require one set of variables less to be held in memory. Such schemes work for arbitrarily high order, although not all Runge–Kutta schemes can be written as $2N$ -schemes (Williamson, 1980; Stanesco and Habashi, 1998). These schemes work iteratively according to the formula

$$w_i = \alpha_i w_{i-1} + \delta t F(t_{i-1}, u_{i-1}), \quad u_i = u_{i-1} + \beta_i w_i. \quad (9.53)$$

Table 9.1 Possible coefficients for different $2N$ -RK3 schemes

Label	α_2	α_3	β_1	β_2	β_3	$(t_1 - t_0)/\delta t$	$(t_2 - t_0)/\delta t$
Symmetric (i)	-2/3	-1	1/3	1	1/2	1/3	2/3
Symmetric (ii)	-1/3	-1	1/3	1/2	1	1/3	2/3
Predictor–corrector	-1/4	-4/3	1/2	2/3	1/2	1/2	1
Inhomogeneous	-17/32	-32/27	1/4	8/9	3/4	1/4	2/3
Quadratic	-0.367	-1.028	0.308	0.540	1	0.308	0.650
Williamson (1980)	-5/9	-153/128	1/3	15/16	8/15	1/3	3/4

For a three-step scheme we have $i = 1, \dots, 3$. In order to advance the variable u from $u^{(n)}$ at time $t^{(n)}$ to $u^{(n+1)}$ at time $t^{(n+1)} = t^{(n)} + \delta h$ we set in (9.53)

$$u_0 = u^{(n)} \quad \text{and} \quad u^{(n+1)} = u_3, \quad (9.54)$$

with u_1 and u_2 being intermediate steps. In order to be able to calculate the first step, $i = 1$, for which no $w_{i-1} \equiv w_0$ exists, we have to require $\alpha_1 = 0$. Thus, we are left with five unknowns, $\alpha_2, \alpha_3, \beta_1, \beta_2$, and β_3 . Three conditions follow from the fact that the scheme be third order, so we have to have two more conditions. One possibility is to choose the fractional times at which the right-hand side is evaluated, e.g. $(0, 1/3, 2/3)$ or even $(0, 1/2, 1)$. In the latter case the right-hand side is evaluated twice at the same time. It is therefore some sort of “predictor–corrector” scheme. In the following, these two schemes are therefore referred to as “symmetric” and “predictor–corrector” schemes. Yet another possibility is to require that inhomogeneous equations of the form $\dot{u} = t^n$ with $n = 1$ and 2 are solved exactly. Such schemes are abbreviated as “inhomogeneous” schemes. The detailed method of calculating the coefficients for such 3rd-order Runge–Kutta schemes with $2N$ -storage is discussed in detail in Appendix B. Several possible sets of coefficients are listed in Table 9.1 and compared with the favorite scheme of Williamson (1980). Note that the 1st-order Euler scheme corresponds to $\beta_1 = 1$ and the classic second-order to $\alpha_2 = -1/2$, $\beta_1 = 1/2$, and $\beta_2 = 1$.

We estimate the accuracy of these schemes by solving the homogeneous differential equation

$$\dot{u} = nu^{1-1/n}, \quad u(1) = 1. \quad (9.55)$$

The exact solution is $u = t^n$. In Table 9.2 we list the rms error with respect to the exact solution, for the range $1 < t \leq 4$ and fixed timestep $\delta t = 0.1$ using $n = -1, 2$, or 3 .

The length of the timestep must always be a certain fraction of the Courant–Friedrich–Levy condition, i.e. $\delta t = k_{\text{CFL}} \delta x / U_{\max}$, where $k_{\text{CFL}} = \mathcal{O}(1)$ and U_{\max} is the maximum transport speed in the system (taking into account advection, sound waves, viscous transport, etc.). Too long a timestep can not only lead to instability, but it also increases the error.

In Table 9.3 we give amplitude and phase errors for the various schemes. The most important conclusion to be drawn from this is the fact that low-order spatial schemes result in large *phase errors*. In the case of a 2nd-order scheme the phase error is 36° after a single passage of a barely resolved wave through a periodic mesh. Higher-order schemes have easily a hundred times smaller phase errors. The amplitude error, on the other hand, is virtually not affected by the spatial order of the scheme. The amplitude error is mainly affected both by the temporal order of the scheme and by the length of the timestep; see also Table 9.4. Therefore,

Table 9.2 Errors (in units of 10^{-6}) for different $2N$ -RK3 schemes, obtained by solving (9.55) in the range $1 < t \leq 4$ with $\delta t = 0.1$ and different values of n

Label	$n = -1$	$n = 2$	$n = 3$
Symmetric (i)	69	103	193
Symmetric (ii)	226	119	411
Predictor–corrector	469	346	1068
Inhomogeneous	84	6	91
Quadratic	197	94	339
Williamson (1980)	68	10	123
For comparison: RK3	66	13	134

Table 9.3 Amplitude and phase errors for inviscid advection of the function $f = \cos kx$ with $k = 2\pi$ and $N = 8$ meshpoints after $t = 20$, corresponding to 20 revolutions in a periodic mesh. The amplitude error is counted positive when the amplitude decreases. A positive phase error means that the solution lags behind the theoretical one

δt scheme	2nd-order	4th-order	6th-order	10th-order	Spectral
RK4	1%	2%	2%	2%	2%
$c_{\delta t} = 0.6$	720°	87°	13°	2.9°	2.8°
RK3	10%	14%	14%	15%	15%
$c_{\delta t} = 0.4$	716°	83°	8°	-2.1°	-2.3°
RK3	4%	6.3%	6.6%	6.6%	6.6%
$c_{\delta t} = 0.3$	717°	84°	10°	-0.6°	-0.8°

Table 9.4 Dependence of the amplitude and phase errors on the length of the timestep and the scheme used for the timestep. In all cases spectral x -derivatives are used

δt scheme	RK3	RK3	RK4	RK4	RK4
$c_{\delta t}$	0.3	0.4	0.4	0.6	1.0
Amplitude error	6.6%	15%	0.3%	2%	21%
Phase error	-0.8°	-2.3°	0.6°	2.8°	18°

high-order schemes with low dissipation and dispersion are particularly important in computational acoustics (Stanescu and Habashi, 1998). However, in applications to turbulence a certain amount of viscosity is always necessary. This would decrease the amplitude of the wave further and would eventually be even more important. (This additional viscosity could be the real one, an explicit artificial, or an implicit numerical viscosity that would result from the discretisation error or the numerical scheme; see Section 9.3.2.)

A common criticism of high-order schemes is their tendency to produce Gibbs phenomena (ripples) near discontinuities. Consequently one needs a small amount of diffusion to damp out the modes near the Nyquist frequency. Thus, one needs to replace (9.50) by the equation

$$\dot{f} = -uf' + vf''. \quad (9.56)$$

The question is now how much diffusion is necessary, and how this depends on the spatial order of the scheme.

A perfect step function would produce large start-up errors; it is better to use a smoothed profile, e.g. one of the form

$$f(x) = \tanh\left(\frac{\cos x}{\delta x}\right), \quad (9.57)$$

where δx is the mesh width. For a periodic mesh of length L one would obviously use $f = f(kx)$, where $k = 2\pi/L$. In that case the step width would be $k\delta x$. In the following we consider a periodic domain of size $L_x = 1$ with $N_x = 60$ meshpoints, so we use $k = 2\pi/L_x = 2\pi$.

In Fig. 9.2 we plot the result of advecting the periodic step-like function, $f(kx)$, over five wavelengths, corresponding to a time $T = L/u$. The goal is to find the minimum diffusion coefficient ν necessary to avoid wiggles in the solution. In the first two panels one sees that for a 6th-order scheme the diffusion coefficient has to be approximately $\nu = 0.01 u\delta x$. For $\nu = 0.005 u\delta x$ there are still wiggles. For a 10th-order scheme one can still use $\nu = 0.005 u\delta x$ without producing wiggles, while for a spectral scheme of nearly infinite order one can go down to $\nu = 0.002 u\delta x$ without any problems.

We may thus conclude that all these schemes need some diffusion, but that the diffusion coefficient can be much reduced when the spatial order of the scheme is high. In that sense it is therefore not true that high-order schemes are particularly vulnerable to Gibbs phenomena, but rather the contrary!

In Fig. 9.3 we compare the corresponding results of advection tests for 2nd- and 4th-order schemes with the 6th-order scheme. It is evident that a 2nd-order scheme requires a relatively high diffusion coefficient, typically around $\nu = 0.05 u\delta x$, but this leads to rather unacceptable distortions of the original profile. (It may be noted that, if one uses at the same time a 1st-order temporal scheme, which has antidiffusive properties, and a timestep which is not too short, then the antidiffusive error of the timestep scheme would partially compensate the actual diffusion and one could reduce the value of ν , but this would be a matter of tuning and hence not generally useful for arbitrary profiles.)

9.3.4. Burgers equation

In the special case where the velocity itself is being advected, i.e. $f = u$, (9.56) turns into the Burgers equation,

$$\dot{u} = -uu' + vu''. \quad (9.58)$$

In one dimension there is an analytic solution for a kink,

$$u = U \left(1 - \tanh \frac{x - Ut}{2\delta} \right), \quad (9.59)$$

where $\delta = \nu/U$ is the shock thickness (e.g. Dodd *et al.*, 1982). Note that the amplitude of the kink is twice its propagation speed. Expressed in terms of the Reynolds number, $Re = UL/\nu$, we have $\delta = LRe^{-1}$. (We note in passing that the dissipative cutoff scale in ordinary turbulence is somewhat larger; $\delta = LRe^{-3/4}$.)

In order to have a stationary shock we use the initial condition

$$u = -U \tanh(x/2\delta). \quad (9.60)$$

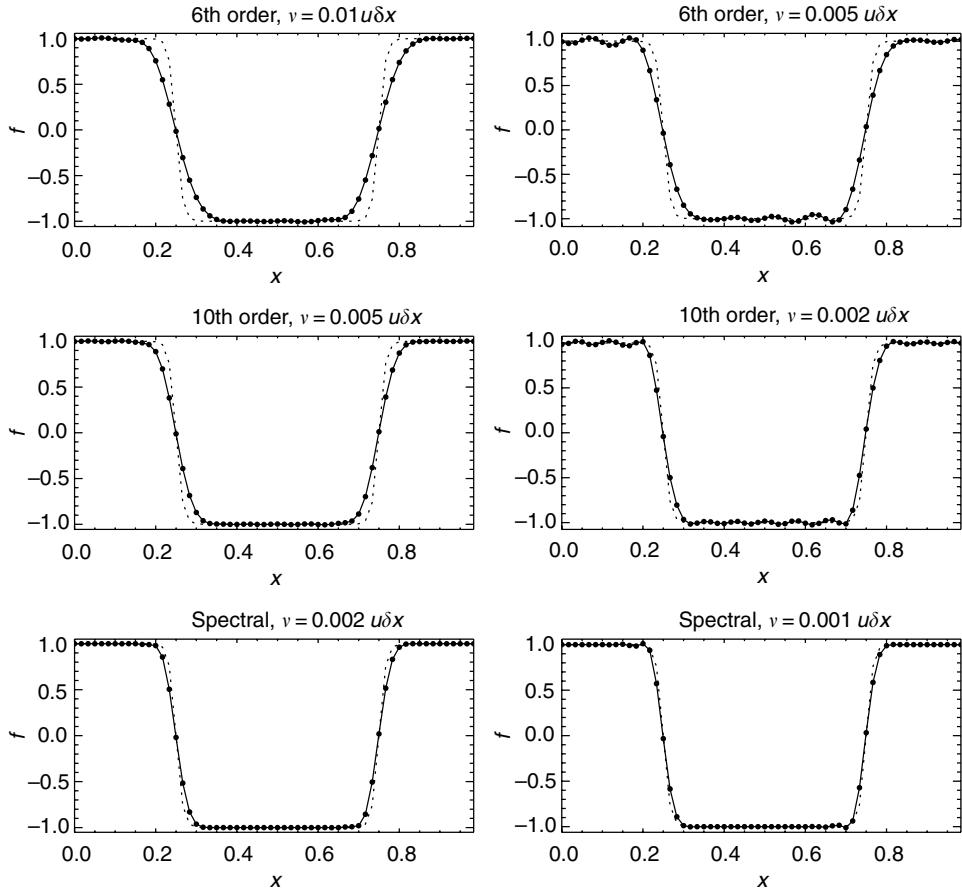


Figure 9.2 Resulting profile after advecting a step-like function five times through the periodic mesh. The dots on the solid line give the location of the function values at the computed mesh-points and the dotted line gives the original profile. For the panels on the right-hand side the diffusion coefficient is too small and the profile shows noticeable wiggles. $\delta x = 1/60$.

In Fig. 9.4 we present numerical solutions using the 6th-order explicit scheme with different values of the mesh Reynolds number, $\delta x U/v$, which was varied by changing the value of v . Here we used $N_x = 100$ meshpoints in the range $-1 \leq x \leq 1$. Note that the overall error, defined here as $\max(|f - f_{\text{exact}}|)$, decreases with decreasing mesh width like δx^5 .

The test cases considered so far were not directly related to the Navier–Stokes equation, e.g. which permits sound waves that can pile up to form shocks. This will be considered in the next section.

9.3.5. Shock tube tests

A popular test problem for compressible codes is the shock tube problem of Sod (1951). On the one hand, one can assess the sharpness of the various fronts. On the other hand, and

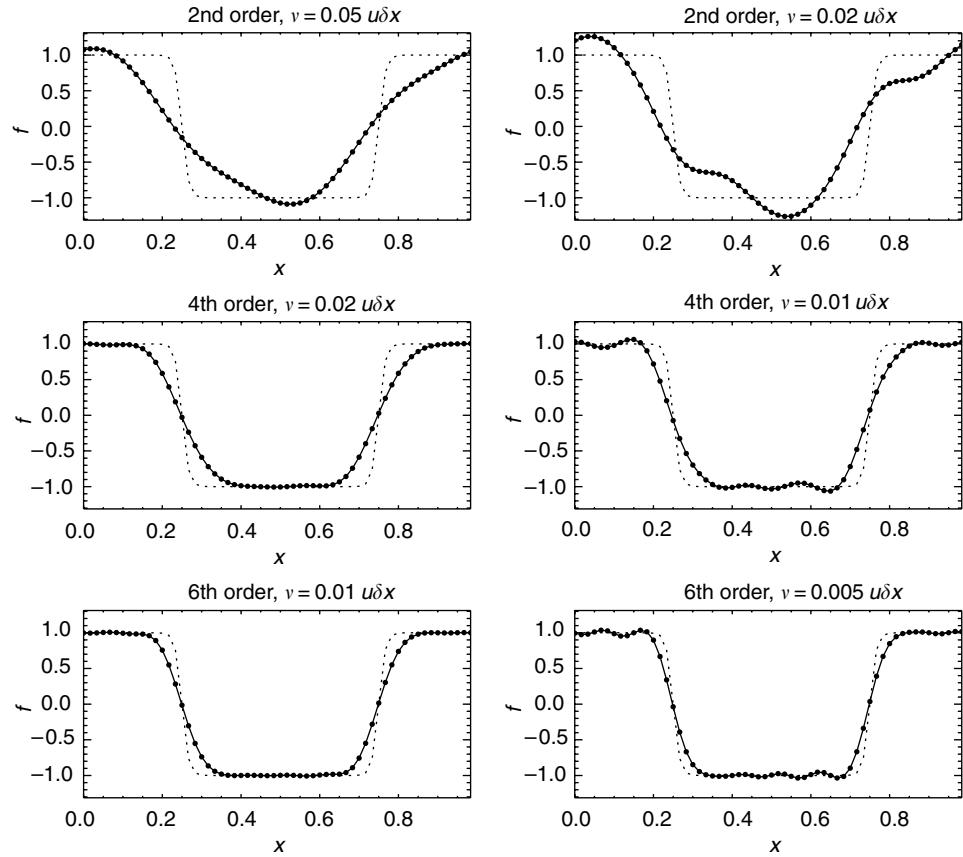


Figure 9.3 Like Fig. 9.2, but for low-order schemes. Again, for the panels on the right-hand side the diffusion coefficient is too small and the profile shows noticeable wiggles. For the 2nd-order scheme one needs a viscosity of $\nu = 0.05 u \delta x$ to prevent wiggles, but then the resulting distortion of the original profile becomes rather unacceptable. $\delta x = 1/60$.

perhaps most importantly, it allows one to test the conversion of kinetic energy to thermal energy via viscous heating.

In the following we use the formulation of the compressible Navier–Stokes equations in terms of entropy and enthalpy (9.16)–(9.18). We use units where $p_0 = \rho_0 = c_p = 1$ and adopt the abbreviations $\Lambda = \ln \rho$ (not to be confused with the cooling function used in Section 9.2). In one dimension (with $\nu = \text{const}$) these equations reduce to

$$\dot{u} = -uu' - c_s^2(\Lambda' + s') + \tilde{\nu}(u'' + \Lambda'u'), \quad (9.61)$$

$$\dot{s} = -us' + (\gamma - 1)\tilde{\nu}u'^2/c_s^2 + Q_s, \quad (9.62)$$

$$\dot{\Lambda} = -u\Lambda' - u', \quad (9.63)$$

where dots and primes refer respectively to time and space derivatives, Q_s describes the change of entropy due to radiative diffusion, and $\Lambda = \ln \rho$ is the logarithmic density. In

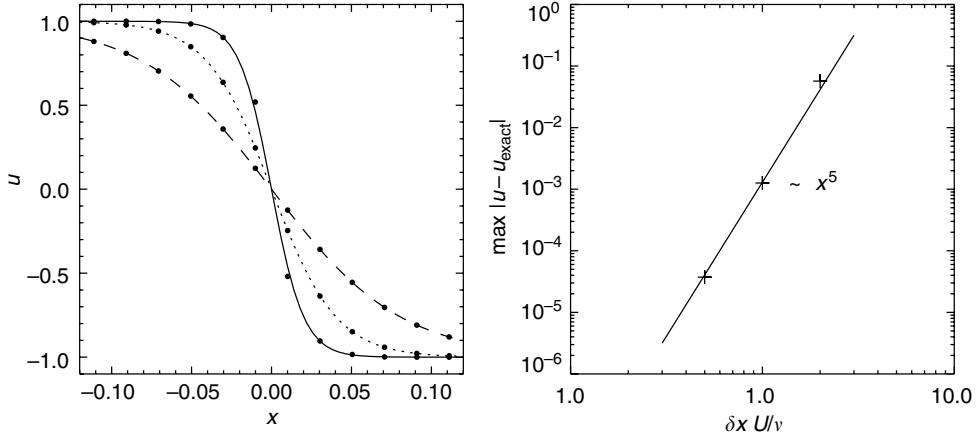


Figure 9.4 Numerical solution of the Burgers equation using the 6th-order explicit scheme with (9.60) as initial condition. In the left-hand panel the lines give the exact solution (9.60) and the dots give the numerical solution for the corresponding value of the mesh Reynolds number: $\delta x U/v = 0.5$ (solid line), 1.0 (dotted line), and 2.0 (dashed line). In the right-hand panel the scaling of the error with δx is shown.

(9.61)–(9.63) we have used the abbreviation

$$c_s^2 = \gamma p / \rho = \gamma \exp[(\gamma - 1)\Lambda + \gamma s] \quad (9.64)$$

for the adiabatic sound speed squared, and $\tilde{\nu} = \frac{4}{3}\nu$ is the effective viscosity for compressive motions. This 4/3 factor comes from the fact that in one-dimensional

$$\mathbf{S} = \text{diag}\left(\frac{2}{3}, -\frac{1}{3}, -\frac{1}{3}\right) u_{x,x}, \quad (9.65)$$

and therefore

$$\frac{1}{\rho} \nabla \cdot (2\nu\rho \mathbf{S}) = \frac{4}{3}\nu [u_{x,xx} + (\ln \rho)_{,x} u_{x,x}], \quad (9.66)$$

so $\mathbf{S}^2 = \frac{2}{3}u_{x,x}^2$, or $2\nu\mathbf{S}^2 = \tilde{\nu}u_{x,x}^2$. In the radiative diffusion approximation we have $Q_s = -\Lambda_{\text{cond}}/(c_p T)$, and so (9.23) gives in one dimension

$$Q_s = \chi\gamma[s'' + \nabla_{\text{ad}}\Lambda'' + \gamma(\Lambda' + s')(s' + \nabla_{\text{ad}}\Lambda')] \quad (9.67)$$

In Fig. 9.5 we show the solution for an initial density and pressure jump of 1 : 10 and the viscosity is now $\nu = 0.6\delta x c_s$. In this case a small amount of thermal diffusion (with Prandtl number $\chi/\nu = 0.05$) has been adopted to remove wiggles in the entropy.

For stronger shocks velocity and entropy excess increase; see Figures 9.6 and 9.7, where the initial pressure jumps are 1 : 100 and 1 : 1000, respectively, and the viscosities are chosen to be $\nu = 1.6\delta x c_s$ and $\nu = 2.4\delta x c_s$. (For the stationary shock problem considered below we also find that the viscosity must increase with the Mach number and, moreover, that the two should be proportional to each other.) In the cases shown in Figs. 9.6 and 9.7 we were able to put $\chi = 0$ without getting any wiggles in s . However, in the case of strong shocks

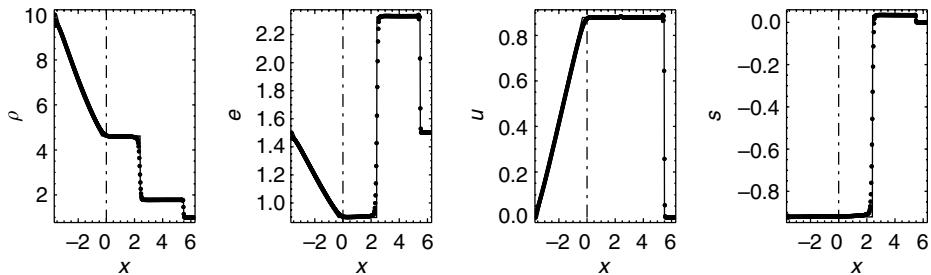


Figure 9.5 Standard shock tube test with an initial pressure jump of 1:10 and $\nu = 0.6\delta x c_s$ and $\chi/\nu = 0.05$. The solid line indicates the analytic solution (in the limit $\tilde{\nu} \rightarrow 0$) and the dots the numerical solution. Note the small entropy excess on the right of the initial entropy discontinuity.

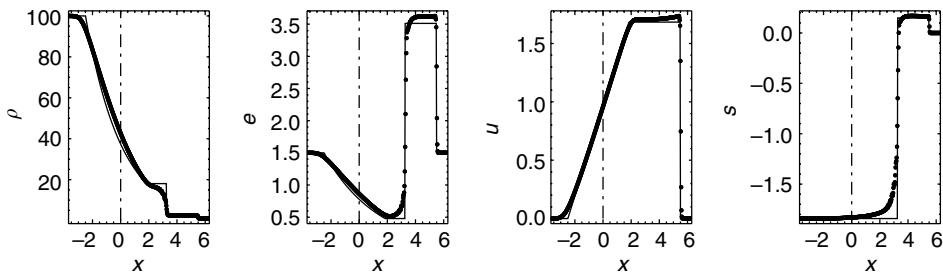


Figure 9.6 Same as Fig. 9.5, but for an initial pressure jump of 1:100 and $\nu = 1.6\delta x c_s$.

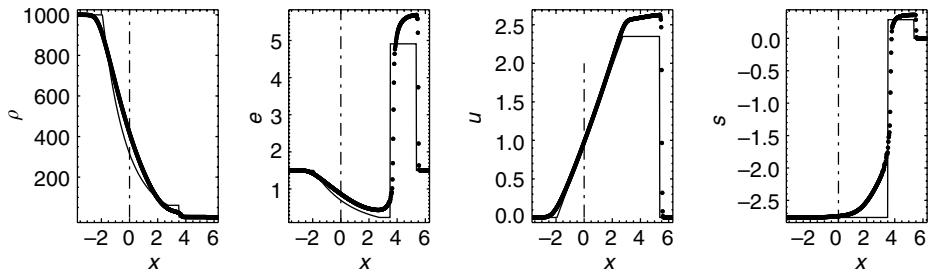


Figure 9.7 Same as Fig. 9.6, but for an initial pressure jump of 1:1000 and $\nu = 2.4\delta x c_s$.

(pressure ratio 1:1000) the discrepancy between numerical and analytical solutions becomes quite noticeable.

In many practical applications shocks occur only in a small portion of space. One can therefore reduce the viscosity outside shocks or, conversely, use a small viscosity everywhere except in the locations of shock, i.e. where the flow is convergent (negative divergent). This

leads to the concept of an artificial (Neumann–Richtmyer) shock viscosity,

$$v_{\text{shock}} = c_{\text{shock}} \delta x^2 \langle (-\nabla \cdot \mathbf{u})_+ \rangle_{\text{n.n.}}, \quad (9.68)$$

where $\langle \dots \rangle_{\text{n.n.}}$ indicates averaging over nearest neighbors and the subscript + means that only the positive part is taken.

The last panel in Figures 9.6 and 9.7 shows quite clearly how the entropy increases behind the shock. This entropy increase is just a consequence of the viscous heating term, $\tilde{\nu} u'^2/h$. Without this term the solution would obviously be wrong everywhere behind the shock, especially when the shock is strong.

A somewhat simpler situation is encountered with standing shocks. In Fig. 9.8 we give an example of a numerically determined solution at $\text{Ma} = 100$. The agreement in the jump for the numerically determined solution (dotted line and dots) and the theoretical solution (solid lines) is very good, although the position of the jump has moved away somewhat from the initial location ($x = 0$), but this is merely a consequence of having used more-or-less arbitrarily a tanh profile to smooth the initial jumps. After some initial adjustment phase the profiles do indeed remain stationary. Note also, however, that the entropy profile is slightly shifted relative to the profiles of u and Λ .

It is interesting to note that when solving the Rankine–Hugonot jump conditions for shocks one is allowed to use the inviscid equations provided they are written in conservative form. Sometimes one finds in the literature the inviscid Navier–Stokes equations written in non-conservative form. This is not strictly correct, because without viscosity there would be no viscous heating and hence no entropy increase behind the shock. Moreover, it is quite common to consider a polytropic equation of state, $p = K\rho^\Gamma$. Again, in this case the entropy is constant, and so energy conservation is violated. Nevertheless, given that polytropic equations of state are often considered in astrophysics we consider this case in more detail in the next subsection.

9.3.6. Polytropic and isothermal shocks

For polytropes with $p = K\rho^\Gamma$, but $\Gamma \neq \gamma$ in general, we can write

$$-\nabla h + h \nabla s = -\frac{1}{\rho} \nabla p = -\nabla \left(\frac{\Gamma K}{\Gamma - 1} \rho^{\Gamma-1} \right) \equiv \nabla \tilde{h}, \quad (9.69)$$

so we can introduce a pseudo enthalpy \tilde{h} as

$$\tilde{h} = \frac{\Gamma K}{\Gamma - 1} \rho^{\Gamma-1} = \left[\left(1 - \frac{1}{\gamma} \right) / \left(1 - \frac{1}{\Gamma} \right) \right] h. \quad (9.70)$$

This is consistent with a fixed entropy dependence, where s only depends on ρ like

$$s = s(\rho) = \frac{1}{\gamma} \ln(p/\rho^\gamma) = \frac{1}{\gamma} \ln(K \rho^{\Gamma-\gamma}) = \frac{1}{\gamma} \ln K + \left(\frac{\Gamma}{\gamma} - 1 \right) \ln \rho, \quad (9.71)$$

which implies that in the polytropic case (9.62) is discarded. In the adiabatic case, $\Gamma = \gamma$, entropy is constant. In the isothermal case, $p = c_s^2 \rho$, we have $\Gamma \rightarrow 1$, so entropy is not constant, but it varies only in direct relation to $-\ln \rho$ and not as a consequence of viscous heating behind the shock.

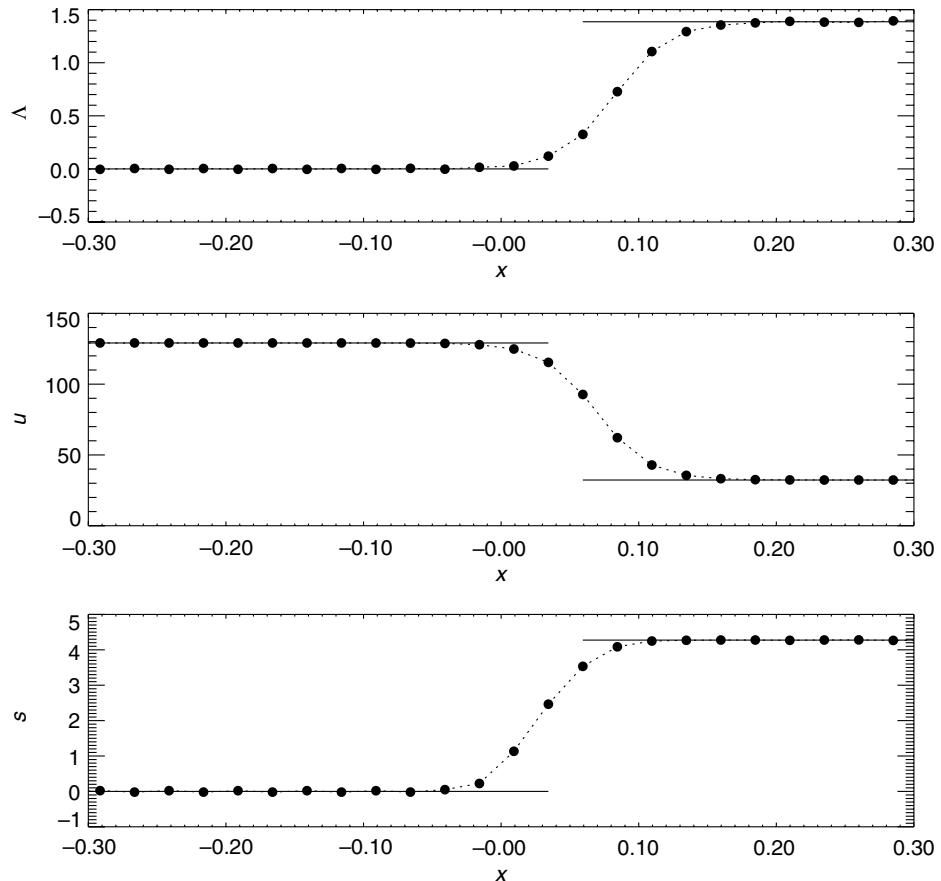


Figure 9.8 Example of a very strong standing shock with $\text{Ma} = 100$. Note the relative shift of the position where s increases relative to where $\Lambda = \ln \rho$ increases. The viscosity is chosen to be $\nu = \text{Ma} \times \delta x$.

In deriving the Rankine–Hugoniot jump conditions one uses the conservation of mass, momentum, and energy in a comoving frame, where the following three quantities are constants of motion:

$$J = \rho u, \quad I = \rho u^2 + p, \quad E = \frac{1}{2} f u^2 + \frac{\gamma}{\gamma - 1} \frac{p}{\rho}. \quad (9.72)$$

The values of these three constants can be calculated when all three variables, u , p , and ρ , are known on one side of the shock. For polytropic equations of state, with $p = K\rho^\gamma$, the energy equation is no longer used, so there are only the following two conserved quantities,

$$J = \rho u, \quad I = \rho u^2 + K\rho^\gamma. \quad (9.73)$$

The dependence of the velocity, density, pressure, and entropy jumps on the upstream Mach number is plotted in Fig. 9.9 for the case $\gamma = 5/3$ and compared with the polytropic case using $\Gamma = \gamma$ (Fig. 9.10).

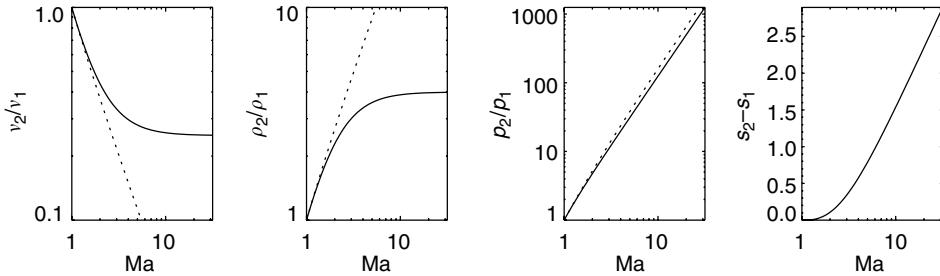


Figure 9.9 The dependence of the velocity, density, pressure, and entropy jumps on the upstream Mach number for the case $\gamma = 5/3$ (solid line) and comparison with the polytropic case using $\Gamma = \gamma$ (dotted line). Note that the velocity and density jumps saturate at $1/4$ and 4 , respectively, while there is no such saturation for the polytropic shock.

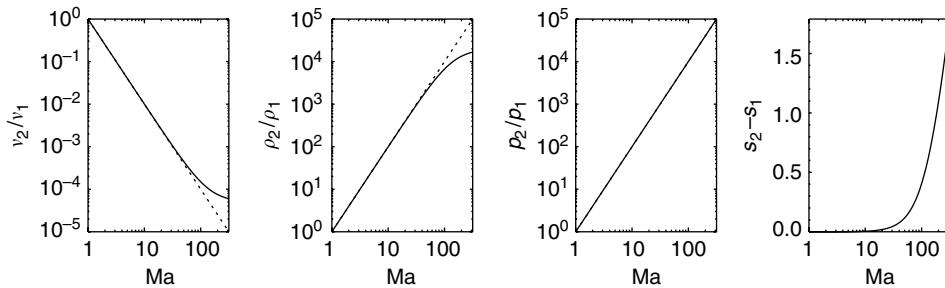


Figure 9.10 Like Fig. 9.9, but for $\gamma = \Gamma = 1.0001$. Note that the velocity and density jumps saturate at much more extreme values than for $\gamma = 5/3$. Thus, the results using polytropic and perfect gas equations agree up to much larger Mach numbers than for $\gamma = 5/3$.

Note that the pressure jump, p_2/p_1 , is almost independent of the value of γ and does also not significantly depend on the polytropic assumption.

9.4. Non-uniform and lagrangian meshes

In many cases it is useful to consider non-uniform meshes, either by adding more points in places where large gradients are expected, or by letting the points move with the flow (lagrangian mesh). The lagrangian mesh is particularly useful in one-dimensional cases, because then the mesh topology (i.e. the ordering of mesh points) remains unchanged. This approach is related to the method of Smooth Particle Hydrodynamics; see Maron and Howes (2001) for references and an improvement of this method. Another method that gains constantly in popularity is adaptive mesh refinement (e.g. Grauer *et al.*, 1998), which will not be discussed here.

9.4.1. Non-uniform topologically cartesian meshes

The implementation of non-uniform meshes can be relatively easy when each of the new coordinates depend on only one variable, e.g. when $x = x(\tilde{x})$, $y = y(\tilde{y})$, and $z = z(\tilde{z})$. Here, \tilde{x} , \tilde{y} ,

and \tilde{z} are cartesian coordinates on a uniform mesh. In the more general case, however, we have

$$x = x(\tilde{x}, \tilde{y}, \tilde{z}), \quad y = y(\tilde{x}, \tilde{y}, \tilde{z}), \quad z = z(\tilde{x}, \tilde{y}, \tilde{z}), \quad (9.74)$$

so that x , y , and z derivatives of a function f can be calculated using the chain rule,

$$\frac{\partial f}{\partial x} = \frac{\partial f}{\partial \tilde{x}} \frac{\partial \tilde{x}}{\partial x} + \frac{\partial f}{\partial \tilde{y}} \frac{\partial \tilde{y}}{\partial x} + \frac{\partial f}{\partial \tilde{z}} \frac{\partial \tilde{z}}{\partial x} \equiv \frac{\partial f}{\partial \tilde{x}_i} \frac{\partial \tilde{x}_i}{\partial x}. \quad (9.75)$$

Corresponding formulae apply obviously for the other two directions, so in general we can write

$$\frac{\partial f}{\partial x_j} = \frac{\partial f}{\partial \tilde{x}_i} J_{ij}, \quad \text{where} \quad J_{ij} = \frac{\partial \tilde{x}_i}{\partial x_j} \quad (9.76)$$

is the jacobian of this coordinate transformation. This method allows one to have high resolution, e.g. near a central object, without however having high resolution anywhere else far away from the central object. This is useful in connection with outflows from jets.

We discuss here one particular application that is relevant for simulating flows in a sphere. It is possible to transform a cartesian mesh to cover a sphere without a coordinate singularity. It will turn out, however, that there is a discontinuity in the jacobian. We discuss this here in two-dimension. We denote the coordinate mesh by a tilde, so (\tilde{x}, \tilde{y}) are the coordinates in a uniform cartesian mesh. We want to stretch the mesh such that points on the \tilde{x} and \tilde{y} axes are not affected, and that the distance of points on the diagonal is reduced by a factor $1/\sqrt{2}$ (or by $1/\sqrt{3}$ in three-dimension). This can be accomplished by introducing new coordinates (x, y) as

$$\begin{pmatrix} x \\ y \end{pmatrix} = \begin{pmatrix} \tilde{x} \\ \tilde{y} \end{pmatrix} \frac{(\tilde{x}^n + \tilde{y}^n)^{1/n}}{(\tilde{x}^2 + \tilde{y}^2)^{1/2}}, \quad (9.77)$$

where n is a large even number. In the limit $n \rightarrow \infty$ we may substitute

$$(\tilde{x}^n + \tilde{y}^n)^{1/n} \rightarrow \max(|\tilde{x}|, |\tilde{y}|). \quad (9.78)$$

Examples of the resulting meshes for two different values of n are given in Fig. 9.11.

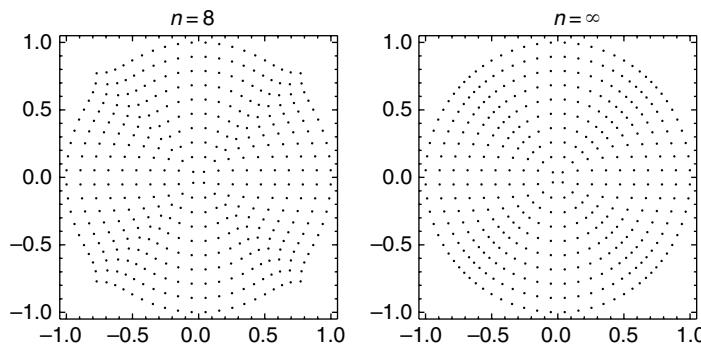


Figure 9.11 Examples of the resulting meshes for $n = 8$ and $n \rightarrow \infty$ in which case (9.78) is used.

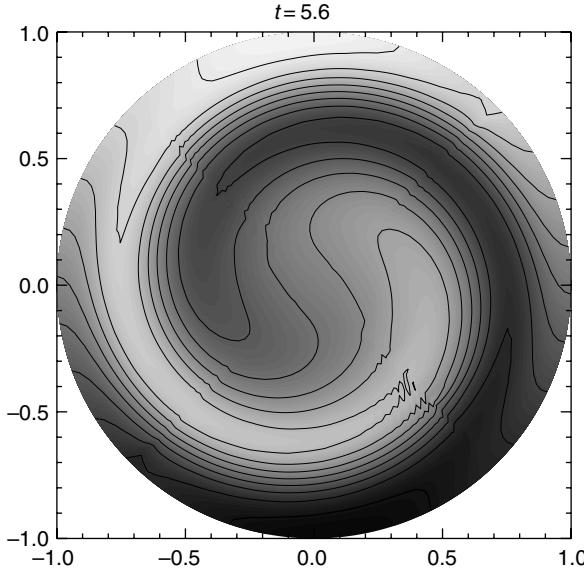


Figure 9.12 Example of an advection experiment on a $n = 8$ mesh.

In order to obtain the jacobian of this transformation, $\partial \tilde{x}_i / \partial x_j$, we have to consider separately the cases $\tilde{x} \geq \tilde{y}$ and $\tilde{x} \leq \tilde{y}$. The derivation is given in Appendix C, and is most concisely expressed in terms of the logarithmic derivative, so

$$\begin{pmatrix} \frac{\partial \ln \tilde{x}}{\partial \ln x} & \frac{\partial \ln \tilde{x}}{\partial \ln y} \\ \frac{\partial \ln \tilde{y}}{\partial \ln x} & \frac{\partial \ln \tilde{y}}{\partial \ln y} \end{pmatrix} = \begin{pmatrix} +1 - \left(\frac{\tilde{y}}{\tilde{r}}\right)^2 & +1 - \left(\frac{\tilde{x}}{\tilde{r}}\right)^2 \\ -1 + \left(\frac{\tilde{x}}{\tilde{r}}\right)^2 & +1 + \left(\frac{\tilde{y}}{\tilde{r}}\right)^2 \end{pmatrix} \quad \text{if } |\tilde{x}| \geq |\tilde{y}|, \quad (9.79)$$

$$\begin{pmatrix} \frac{\partial \ln \tilde{x}}{\partial \ln x} & \frac{\partial \ln \tilde{x}}{\partial \ln y} \\ \frac{\partial \ln \tilde{y}}{\partial \ln x} & \frac{\partial \ln \tilde{y}}{\partial \ln y} \end{pmatrix} = \begin{pmatrix} +1 + \left(\frac{\tilde{x}}{\tilde{r}}\right)^2 & -1 + \left(\frac{\tilde{y}}{\tilde{r}}\right)^2 \\ +1 - \left(\frac{\tilde{y}}{\tilde{r}}\right)^2 & +1 - \left(\frac{\tilde{x}}{\tilde{r}}\right)^2 \end{pmatrix} \quad \text{if } |\tilde{x}| \leq |\tilde{y}|, \quad (9.80)$$

where $\tilde{r}^2 = \tilde{x}^2 + \tilde{y}^2$. Note that the jacobian is discontinuous on the diagonals. This is a somewhat unfortunate feature of this transformation. It is not too surprising however that something like this happens, because the diagonals are the locations where a rotating flow must turn direction by 90° in the coordinate mesh. Nevertheless, it is possible to obtain reasonably well behaved solutions; see Fig. 9.12 for an advection experiment using a prescribed differentially rotating flow.

The fluid equations are still solved in rectangular cartesian coordinates, so, e.g. the equation $Ds/Dt = 0$ is solved in the form

$$\frac{\partial s}{\partial t} = -u_x \frac{\partial s}{\partial x} - u_y \frac{\partial s}{\partial y}, \quad (9.81)$$

where the spatial derivatives are evaluated according to (9.75). For the velocity field, stress-free boundary conditions, e.g. would be written in the form

$$\hat{r}_j u_j = 0, \quad \hat{\phi}_i S_{ij} \hat{r}_j = 0, \quad (9.82)$$

where S_{ij} is the rate of strain tensor, \hat{r}_j and $\hat{\phi}_i$ are the cartesian components ($i = x, y, z$) of the radial and azimuthal unit vectors, i.e.

$$\hat{\mathbf{r}} = \frac{1}{r} \begin{pmatrix} x \\ r \end{pmatrix} y \quad \text{and} \quad \hat{\boldsymbol{\phi}} = \frac{1}{r} \begin{pmatrix} -y \\ x \end{pmatrix} \quad (9.83)$$

are unit vectors in the r and ϕ directions and $r = \sqrt{x^2 + y^2}$ is the distance from the rotation axis. The stress-free boundary conditions are then

$$x u_x + y u_y = 0 \quad (9.84)$$

and

$$(x^2 - y^2)(u_{x,y} + u_{y,x}) - 2xy(u_{x,x} - u_{y,y}) = 0. \quad (9.85)$$

9.4.2. Lagrangian meshes

We now consider a simple one-dimensional lagrangian mesh problem. Assume that ℓ labels the particle, then the lagrangian derivative is

$$\frac{Ds}{Dt} \equiv \left(\frac{\partial s}{\partial t} \right)_{\ell=\text{const}} = \left(\frac{\partial s}{\partial t} \right)_{x=\text{const}} + \left(\frac{\partial x}{\partial t} \right)_{\ell=\text{const}} \frac{\partial s}{\partial x}. \quad (9.86)$$

Now, because

$$\left(\frac{\partial x}{\partial t} \right)_{\ell=\text{const}} = \frac{Dx}{Dt} = u, \quad (9.87)$$

we have the well-known equation

$$\frac{Ds}{Dt} = \frac{\partial s}{\partial t} + u \frac{\partial s}{\partial x}. \quad (9.88)$$

As an example we now consider the Burgers equation,

$$\frac{Du}{Dt} = \tilde{v} \frac{\partial^2 u}{\partial x^2}. \quad (9.89)$$

We now take $u(x, t) = u(\ell(x), t)$ to be a function of the coordinate variable ℓ which, in turn, is a function of x . The x -derivatives are obtained using the chain rule, i.e.

$$\frac{\partial u}{\partial x} = \frac{\partial \ell}{\partial x} \frac{\partial u}{\partial \ell} = \frac{u'}{x'}, \quad (9.90)$$

and likewise for the second derivative

$$\frac{\partial^2 u}{\partial x^2} = \frac{u''x' - u'x''}{x'^3}. \quad (9.91)$$

Thus, the Burgers equation can then be written as

$$\frac{\partial u}{\partial t} = \tilde{v} \frac{u''x' - u'x''}{x'^3}, \quad (9.92)$$

where the x variable is given by

$$\frac{\partial x}{\partial t} = u. \quad (9.93)$$

A solution of these two equations is given in Fig. 9.13.

In the test problem above the initial meshpoint distribution was uniform. Although this is not quite suitable for this problem, it shows that subsequently the mesh spacing became narrower still, which means that the timestep is now governed by viscosity, $\delta t \leq 0.06\delta x_{\min}^2/\tilde{v}$, where the numerical factor is empirical. However, the mesh spacing does not need to be governed by (9.93), so it is quite possible to come up with other prescriptions for the mesh spacing.

Consider as another example the isothermal eulerian equations

$$\frac{Du}{Dt} = -c_s^2 \frac{\partial \ln \rho}{\partial x} + \tilde{v} \left(\frac{\partial^2 u}{\partial x^2} + \frac{\partial \ln \rho}{\partial x} \frac{\partial u}{\partial x} \right), \quad (9.94)$$

$$\frac{D \ln \rho}{Dt} = -\frac{\partial u}{\partial x}. \quad (9.95)$$

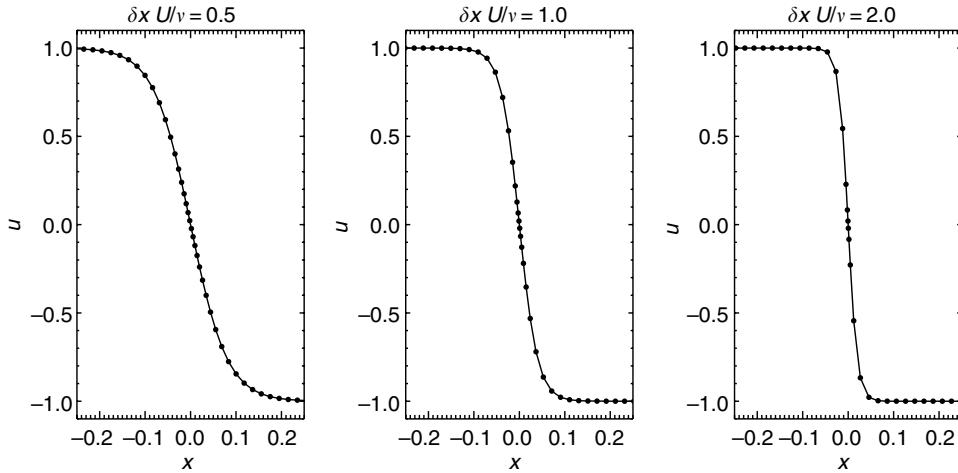


Figure 9.13 Solution of the Burgers equation using a lagrangian mesh combined with a 6th-order explicit scheme. The values of the mesh Reynolds number vary between $\delta x U/v = 0.5$ and 2.0, where δx refers to the initially uniform mesh spacing.

In lagrangian form they take the form

$$\frac{\partial u}{\partial t} = -c_s^2 \frac{(\ln \rho)'}{x'} + \tilde{v} \left(\frac{u''x' - u'x''}{x'^3} + \frac{(\ln \rho)'}{x'} \frac{u'}{x'} \right), \quad (9.96)$$

$$\frac{\partial \ln \rho}{\partial t} = -\frac{u'}{x'}, \quad (9.97)$$

$$\frac{\partial x}{\partial t} = u. \quad (9.98)$$

The example above demonstrates clearly the problem that lagrangian mesh points can continue to pile up near convergence points of the flow. This is a general problem with fully lagrangian schemes. One possible alternative is to use *lagrangian-eulerian* schemes (e.g. Benson, 1992; Peterkin *et al.*, 1998; Arber *et al.*, 2001), which combine the advantages of lagrangian and eulerian codes, but involve obviously some kind of interpolation. Another alternative is to use a semi-lagrangian code which advects the mesh points not with the actual gas velocity \mathbf{u} , but with a more independent mesh velocity \mathbf{U} . Clearly, we want to avoid too small distances between neighboring points, so one could artificially lower the effective mesh velocity by involving, e.g. the modulus of the jacobian, $|\mathbf{J}|$, which becomes large when the concentration of mesh points is high. Thus, one could choose, e.g. $\mathbf{U} = \mathbf{u}/(1 + |\mathbf{J}|)$. In the present case, $|\mathbf{J}| = |x'|^{-1}$. In the following we discuss the formalism that needs to be invoked in order to calculate first and second derivatives on an advected mesh.

9.4.3. Non-lagrangian mesh advection

The main advantage of a lagrangian mesh is that it allows higher resolution locally. Another advantage, which is however less crucial, is that the nonlinear advection term drops out. The main disadvantage is however that a lagrangian mesh may become too distorted and overconcentrated, as seen in the previous section. In this subsection we address the possibility of advecting the mesh with a velocity \mathbf{U} that can be different from the fluid velocity. This way one can remove the swirl of the mesh by taking a velocity that is the gradient of some other quantity, i.e.

$$\mathbf{U} = -\nabla \Phi_{\text{mesh}}, \quad (9.99)$$

where Φ_{mesh} should be large in those regions where many points are needed. One possible criterion would be to require that the number of scale heights per meshpoint, $|\delta \mathbf{x} \cdot \nabla \ln \rho|$, does not exceed an empirical value of $1/3$, say. Thus, $3|\delta \mathbf{x} \cdot \nabla \ln \rho| < 1$ would be a necessary condition. Another possibility would be to let Φ_{mesh} evolve itself according to some suitable advection-diffusion equation. However, no generally satisfactory method seems to be available as yet. In order to calculate the jacobian for the coordinate transformation one can make use of the fact that the mesh evolves only gradually from one timestep to the next. For a more extended discussion of mesh advection schemes we refer to the article by Dorfi in the book by LeVeque *et al.* (1998).

9.4.3.1. Calculating the jacobian

Initially, at $t = 0$, we have $\mathbf{x} = \tilde{\mathbf{x}}$. After the n th timestep, at $t = n\delta t$, we calculate the new \mathbf{x} -mesh, $\mathbf{x}^{(n+1)}$, from the previous one, $\mathbf{x}^{(n)}$, i.e.

$$\mathbf{x}^{(n+1)} = \mathbf{x}^{(n)} + \mathbf{U}(\mathbf{x}^{(n)}, t) \delta t. \quad (9.100)$$

Here, $\mathbf{x}^{(0)} = \tilde{\mathbf{x}}$ is just the original coordinate mesh. Differentiating the i th component with respect to the j th component, as we have done in Section 9.4.1, we obtain

$$\delta_{ij} = \frac{\partial x_i^{(n+1)}}{\partial x_j^{(n+1)}} = \frac{\partial x_i^{(n)}}{\partial x_j^{(n+1)}} + \frac{\partial U_i^{(n)}}{\partial x_j^{(n+1)}} \delta t, \quad (9.101)$$

where $U_i^{(n)} = U_i(\mathbf{x}^{(n)}, t)$. In the expression above we have U_i on the mesh $\mathbf{x}^{(n)}$, but we need to differentiate with respect to the new mesh $\mathbf{x}^{(n+1)}$. This can be fixed by another factor $\partial x_i^{(n)}/\partial x_j^{(n+1)}$. Thus, we have

$$\delta_{ij} = \frac{\partial x_i^{(n)}}{\partial x_j^{(n+1)}} + \frac{\partial x_k^{(n)}}{\partial x_j^{(n+1)}} \frac{\partial U_i(\mathbf{x}^{(n)}, t)}{\partial x_k^{(n)}} \delta t = \left[\delta_{ik} + \frac{\partial U_i^{(n)}}{\partial x_k^{(n)}} \delta t \right] \frac{\partial x_k^{(n)}}{\partial x_j^{(n+1)}}. \quad (9.102)$$

This can be written in matrix form,

$$\delta_{ij} = \mathbf{M}_{ik} \mathbf{J}_{kj}^{(n)}, \quad (9.103)$$

where

$$\mathbf{M}_{ik} = \delta_{ik} + \frac{\partial U_i(\mathbf{x}^{(n)}, t)}{\partial x_k^{(n)}} \delta t \quad (9.104)$$

is a transformation matrix and

$$\mathbf{J}_{ij}^{(n)} = \frac{\partial x_i^{(n)}}{\partial x_j^{(n+1)}} \quad (9.105)$$

is the incremental jacobian, so $\mathbf{J}^{(n)} = \mathbf{M}^{-1}$. To obtain the jacobian at $t = 2\delta t$, e.g. we calculate

$$\frac{\partial x_i^{(0)}}{\partial x_j^{(2)}} = \frac{\partial x_i^{(0)}}{\partial x_k^{(1)}} \frac{\partial x_k^{(1)}}{\partial x_j^{(2)}} = \mathbf{J}_{ik}^{(0)} \mathbf{J}_{kj}^{(1)} \equiv (\mathbf{J}^{(0)} \mathbf{J}^{(1)})_{ij}. \quad (9.106)$$

The jacobian at $t = n\delta t$ is then obtained by successive matrix multiplication from the right, so

$$\mathbf{J}_{ij}^{(0 \rightarrow n+1)} = \mathbf{J}_{ik}^{(0 \rightarrow n)} \mathbf{J}_{kj}^{(n)}, \quad (9.107)$$

where $\mathbf{J}_{ij}^{(0 \rightarrow n+1)}$ and $\mathbf{J}_{ij}^{(0 \rightarrow n)}$ are the full (as opposed to incremental) jacobians at the new and previous timesteps, respectively.

9.4.3.2. Calculating the 2nd-order jacobian

A corresponding calculation (see Appendix D) for the second derivatives of a function f shows that

$$\frac{\partial^2 f}{\partial x_i \partial x_j} = \frac{\partial^2 f}{\partial \tilde{x}_p \partial \tilde{x}_q} \mathbf{J}_{pi} \mathbf{J}_{qj} + \frac{\partial f}{\partial \tilde{x}_k} \mathbf{K}_{kij}, \quad (9.108)$$

where

$$\mathbf{K}_{kij} = \frac{\partial^2 \tilde{x}_k}{\partial x_i \partial x_j} \quad (9.109)$$

is the 2nd-order jacobian. Like for the first derivative the 2nd-order jacobian can be obtained by successive tensor multiplication,

$$\mathbf{K}_{kij}^{(0 \rightarrow n+1)} = \mathbf{K}_{kpq}^{(0 \rightarrow n)} \mathbf{J}_{pi}^{(n)} \mathbf{J}_{qj}^{(n)} + \mathbf{J}_{kl}^{(0 \rightarrow n)} \mathbf{K}_{lij}^{(n)}, \quad (9.110)$$

where $\mathbf{K}_{kij}^{(0 \rightarrow n+1)}$ and $\mathbf{K}_{kij}^{(0 \rightarrow n)}$ are the 2nd-order jacobians at the new and previous timesteps, respectively, and

$$\mathbf{K}_{kij}^{(n)} = \frac{\partial^2 x_k^{(n)}}{\partial x_i^{(n+1)} \partial x_j^{(n+1)}} \quad (9.111)$$

is the incremental 2nd-order jacobian, which is calculated at each timestep as

$$\mathbf{K}_{kij}^{(n)} = -\delta t \left(\mathbf{M}^{-1} \right)_{kl} U_{l,pq} \mathbf{J}_{pi}^{(n)} \mathbf{J}_{qj}^{(n)}, \quad (9.112)$$

where \mathbf{M} was defined in (9.104) and

$$U_{l,pq} = \frac{\partial^2 U_l(\mathbf{x}^{(n)}, t)}{\partial x_p^{(n)} \partial x_q^{(n)}} \quad (9.113)$$

is the 2nd-order velocity gradient matrix on the physical mesh. Since $\mathbf{M}^{-1} = \mathbf{J}^{(n)}$ is just the incremental jacobian, we can write (9.112) as

$$\mathbf{K}_{kij}^{(n)} = -\mathbf{J}_{kl}^{(n)} \delta t U_{l,pq} \mathbf{J}_{pi}^{(n)} \mathbf{J}_{qj}^{(n)}. \quad (9.114)$$

Since the expressions (9.110) and (9.114) involve both a multiplication with $\mathbf{J}_{pi}^{(n)} \mathbf{J}_{qj}^{(n)}$, we can simplify (9.110) to give

$$\mathbf{K}_{kij}^{(0 \rightarrow n+1)} = \left[\mathbf{K}_{kpq}^{(0 \rightarrow n)} - \mathbf{J}_{kl}^{(0 \rightarrow n)} \mathbf{J}_{lm}^{(n)} \delta t U_{m,pq} \right] \mathbf{J}_{pi}^{(n)} \mathbf{J}_{qj}^{(n)}. \quad (9.115)$$

Here the expression $\mathbf{J}_{kl}^{(0 \rightarrow n)} \mathbf{J}_{lm}^{(n)}$ is of course the new jacobian, $\mathbf{J}_{kl}^{(0 \rightarrow n+1)}$.

So in summary, the new 1st- and 2nd-order jacobians are obtained from the previous ones via the formulae

$$\mathbf{J}_{ij}^{(0 \rightarrow n+1)} = \mathbf{J}_{ik}^{(0 \rightarrow n)} \mathbf{J}_{kj}^{(n)}, \quad (9.116)$$

$$\mathbf{K}_{kij}^{(0 \rightarrow n+1)} = \left[\mathbf{K}_{kpq}^{(0 \rightarrow n)} - \mathbf{J}_{kl}^{(0 \rightarrow n+1)} \delta t U_{l,pq} \right] \mathbf{J}_{pi}^{(n)} \mathbf{J}_{qj}^{(n)}, \quad (9.117)$$

$$\frac{\partial f}{\partial x_i} = \frac{\partial f}{\partial \tilde{x}_p} \mathbf{J}_{pi}, \quad (9.118)$$

$$\frac{\partial^2 f}{\partial x_i \partial x_j} = \frac{\partial^2 f}{\partial \tilde{x}_p \partial \tilde{x}_q} \mathbf{J}_{pi} \mathbf{J}_{qj} + \frac{\partial f}{\partial \tilde{x}_k} \mathbf{K}_{kij}, \quad (9.119)$$

where $\mathbf{J} \equiv \mathbf{J}^{(0 \rightarrow n+1)}$ and $\mathbf{K} \equiv \mathbf{K}^{(0 \rightarrow n+1)}$ has been assumed.

Since now the mesh is moving in time with the local speed \mathbf{U} which is different from the gas velocity \mathbf{u} , the lagrangian derivative is

$$\frac{D}{Dt} = \frac{\partial}{\partial t} + (\mathbf{u} - \mathbf{U}) \cdot \nabla. \quad (9.120)$$

In all other respects the basic equations, written in cartesian form, are still unchanged, provided all x , y , and z derivatives (first and second) are evaluated, as in (9.75) and (9.108), with the components of the jacobian. As an example we show in Fig. 9.14 the result of a kinematic collapse calculation where $D\mathbf{u}/Dt = -\nabla\phi$ and $Ds/Dt = 0$ with a smoothed but localized gravitational potential ϕ . In Fig. 9.15 we compare the results of an eulerian and a lagrangian calculation using the same number of meshpoints. Already after some short time the eulerian calculation begins to become underresolved and develops wiggles while the lagrangian calculation proceeds without problems.

9.4.4. Unstructured meshes

We now discuss how we can calculate spatial derivatives of our variables from a non-uniformly spaced ensemble of points. Consider the function $f(x, y, z)$, which stand for one of the components of a vector (velocity or magnetic vector potential) or a scalar, such as $\ln \rho$. We approximate the function $f(x, y, z)$ in the neighborhood of the point $\mathbf{x}_i = (x_i, y_i, z_i)$ by a multidimensional polynomial of degree N ,

$$f(x, y, z) = f(x_i, y_i, z_i) + \sum_{l+m+n \leq N} c_{lmn} (x - x_i)^l (y - y_i)^m (z - z_i)^n, \quad (9.121)$$

where l , m , and n are non-negative integers and c_{lmn} are coefficients that are to be determined separately for each point by applying (9.121) to all neighboring points \mathbf{x}_j . Note that $c_{000} = 0$ and does not need to be considered. Thus, for each point j we have a system of equations

$$f_{ij} = \sum_{l+m+n \leq N} \frac{1}{l! m! n!} c_{lmn} x_{ij}^l y_{ij}^m z_{ij}^n, \quad (9.122)$$

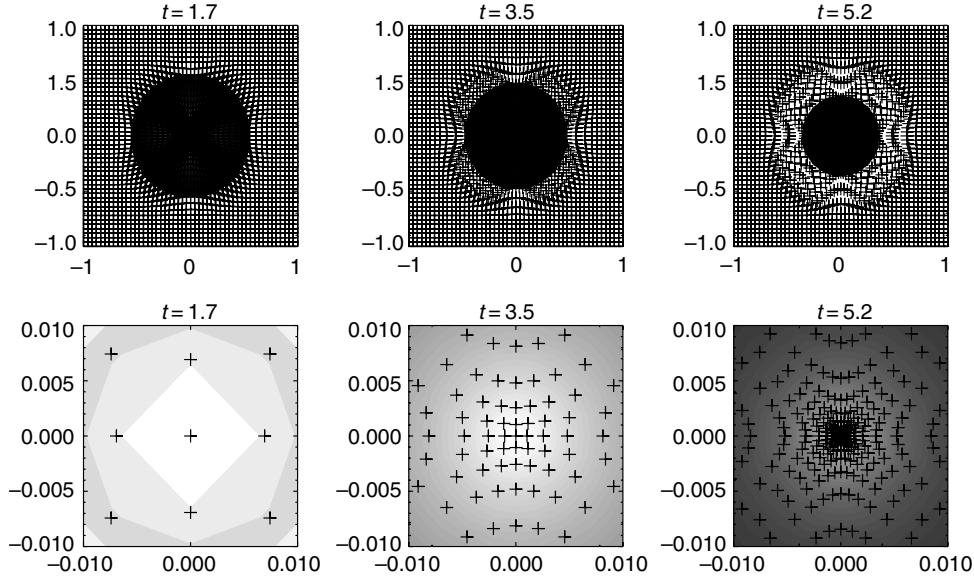


Figure 9.14 Example of a collapse calculation. The second row shows only the inner parts with $|x|, |y| \leq 0.01$ at the same times as in the upper row.

where $f_{ij} = f(x_i, y_i, z_i) - f(x_j, y_j, z_j)$ and $\mathbf{x}_{ij} = \mathbf{x}_i - \mathbf{x}_j$. This system of equations can be written in matrix form

$$\mathbf{f}_\alpha = \mathbf{M}_{\alpha\beta} \mathbf{C}_\beta, \quad (9.123)$$

where $1 \leq (\alpha, \beta) \leq M$ and M is the spatial dimension of the matrix, which is related to N and the dimension as follows:

$$M = \begin{cases} N & \text{in one-dimension,} \\ (N+1)(N+2)/2 - 1 & \text{in two-dimension,} \\ (N+1)(3N/2) & \text{in three-dimension.} \end{cases} \quad (9.124)$$

When $N = 2$ the matrix \mathbf{M} is given by

$$\mathbf{M} = \begin{pmatrix} x_{ij_1} & y_{ij_1} & z_{ij_1} & \frac{1}{2}x_{ij_1}^2 & x_{ij_1}y_{ij_1} & \frac{1}{2}y_{ij_1}^2 & y_{ij_1}z_{ij_1} & \frac{1}{2}z_{ij_1}^2 & z_{ij_1}x_{ij_1} \\ x_{ij_2} & y_{ij_2} & z_{ij_2} & \frac{1}{2}x_{ij_2}^2 & x_{ij_2}y_{ij_2} & \frac{1}{2}y_{ij_2}^2 & y_{ij_2}z_{ij_2} & \frac{1}{2}z_{ij_2}^2 & z_{ij_2}x_{ij_2} \\ \dots & \dots \\ x_{ij_M} & y_{ij_M} & z_{ij_M} & \frac{1}{2}x_{ij_M}^2 & x_{ij_M}y_{ij_M} & \frac{1}{2}y_{ij_M}^2 & y_{ij_M}z_{ij_M} & \frac{1}{2}z_{ij_M}^2 & z_{ij_M}x_{ij_M} \end{pmatrix} \quad (9.125)$$

and

$$\mathbf{C} = (c_{100}, c_{010}, c_{001}, c_{200}, c_{110}, c_{020}, c_{011}, c_{002}, c_{101})^T. \quad (9.126)$$

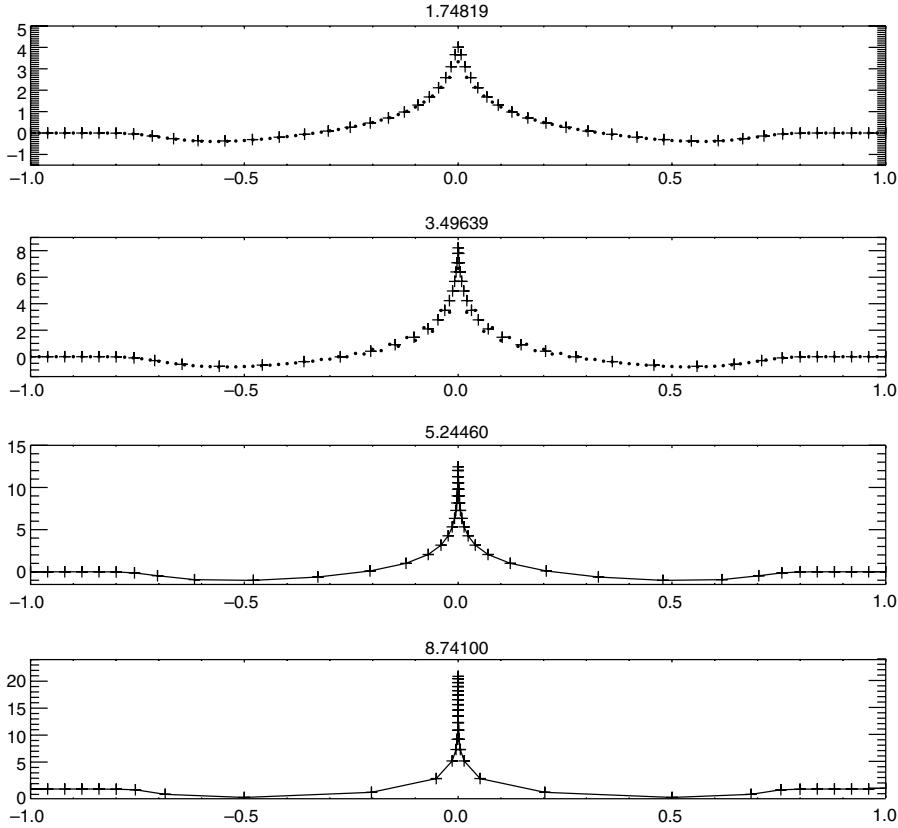


Figure 9.15 Comparison of lagrangian (+ signs) and eulerian (dots) calculations in the first two plots, and later development (last two plots) where the eulerian no longer works. Note that already in the second plot the eulerian calculation has developed noticeable wiggles which the lagrangian proceeds without problems.

Here, j_n ($n = 1, 2, \dots, M$) are the M nearest neighbors of the point i . In general the matrix can be written in the form

$$M_{\alpha\beta}^{(J)} = x_{\alpha J}^{l(\beta)} y_{\alpha J}^{m(\beta)} z_{\alpha J}^{n(\beta)}, \quad (9.127)$$

where J is the index of the point at which the derivative is to be calculated. The set of exponents $l(\beta)$, $m(\beta)$, and $n(\beta)$ is given here for the case $N = 4$ in three-dimensional

$$l(\beta) = (1, 0, 0|2, 1, 0, 0, 0, 1|3, 2, 1, 0, 0, 0, 0, 1, 2|4, 3, 2, 1, 0, 0, 0, 0, 0, 1, 2, 3), \quad (9.128)$$

$$m(\beta) = (0, 1, 0|0, 1, 2, 1, 0, 0|0, 1, 2, 3, 2, 1, 0, 0, 0|0, 1, 2, 3, 4, 3, 2, 1, 0, 0, 0, 0, 0), \quad (9.129)$$

$$n(\beta) = (0, 0, 1|0, 0, 0, 0, 1, 2, 1|0, 0, 0, 0, 1, 2, 3, 2, 1|0, 0, 0, 0, 0, 1, 2, 3, 4, 3, 2, 1), \quad (9.130)$$

where the vertical bars separate the sets of exponents that correspond to increasing orders. Once the C vector has been obtained, the first derivatives of f are simply given by

$$\frac{\partial f}{\partial x} = c_{100}, \quad \frac{\partial f}{\partial y} = c_{010}, \quad \frac{\partial f}{\partial z} = c_{001}. \quad (9.131)$$

Likewise, the second derivatives are given by

$$\frac{\partial^2 f}{\partial x^2} = c_{200}, \quad \frac{\partial^2 f}{\partial y^2} = c_{020}, \quad \frac{\partial^2 f}{\partial z^2} = c_{002}, \quad (9.132)$$

and the mixed second derivatives are given by

$$\frac{\partial^2 f}{\partial x \partial y} = c_{110}, \quad \frac{\partial^2 f}{\partial y \partial z} = c_{011}, \quad \frac{\partial^2 f}{\partial z \partial x} = c_{101}. \quad (9.133)$$

Although this method can be used for meshes that are static in time, it can also be used in connection with multidimensional lagrangian schemes. In that case there may arise the problem that neighboring points get very close together, and so small errors strongly affect the coefficients. A good way out of this is to use a few more points and to solve the linear matrix equation using singular value decomposition. An example of such a calculation is shown in Fig. 9.16, where a passive scalar, with the initial distribution $A(\mathbf{x}, 0) = x$, is advected by the velocity, $\mathbf{u} = \dot{\mathbf{r}}$, which in turn is obtained by solving Kepler's equation, $\ddot{\mathbf{r}} = -GM\mathbf{r}/r^3$, using the normalization $GM = 1$. This windup problem corresponds to the windup of initially horizontal magnetic field lines.

In diffusivity used in Fig. 9.16 was $\eta = 0.02$, but due to the coarse resolution and the implicit smoothing resulting from the singular value decomposition technique the effective diffusivity is somewhat larger.

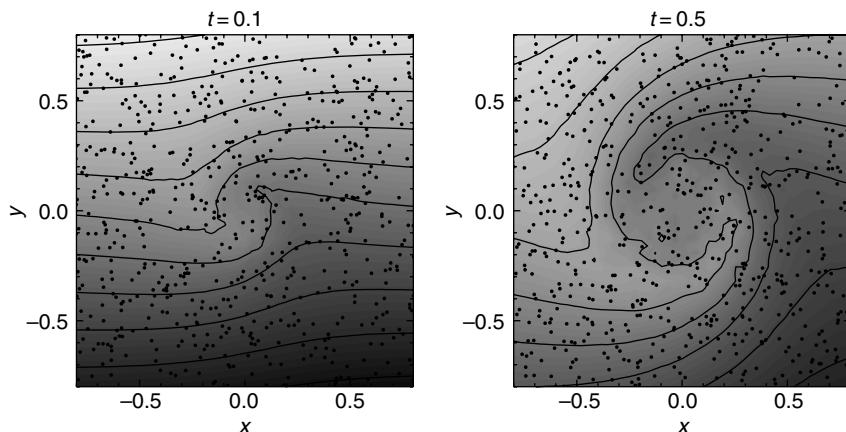


Figure 9.16 Two-dimensional advection problem on an unstructured lagrangian mesh. The dots indicate the 1000 lagrangian particles which constitute the unstructured mesh.

9.5. Implementing magnetic fields

As mentioned in Section 9.2, implementing magnetic fields is relatively straightforward. On the one hand, the magnetic field causes a Lorentz force, $\mathbf{J} \times \mathbf{B}$, where \mathbf{B} is the flux density, $\mathbf{J} = \nabla \times \mathbf{B}/\mu_0$ is the current density, and μ_0 is the vacuum permeability. Note, however, that $\mathbf{J} \times \mathbf{B}$ is the force per unit volume, so in (9.9) we need to add the term $\mathbf{J} \times \mathbf{B}/\rho$ on the right-hand side. On the other hand, \mathbf{B} itself evolves according to the Faraday equation,

$$\frac{\partial \mathbf{B}}{\partial t} = -\nabla \times \mathbf{E}, \quad (9.134)$$

where the electric field \mathbf{E} can be expressed in terms of \mathbf{J} using Ohm's law in the laboratory frame, $\mathbf{E} = -\mathbf{u} \times \mathbf{B} + \mathbf{J}/\sigma$, where $\sigma = (\eta\mu_0)^{-1}$ is the electric conductivity and η is the magnetic diffusivity.

In addition we have to satisfy the condition $\nabla \cdot \mathbf{B} = 0$. This is most easily done by solving not for \mathbf{B} , but instead for the magnetic vector potential \mathbf{A} , where $\mathbf{B} = \nabla \times \mathbf{A}$. The evolution of \mathbf{A} is governed by the uncurred form of (9.134),

$$\frac{\partial \mathbf{A}}{\partial t} = -\mathbf{E} - \nabla\phi, \quad (9.135)$$

where ϕ is the electrostatic potential, which takes the role of an integration constant which does not affect the evolution of \mathbf{B} . The choice $\phi = 0$ is most advantageous on numerical grounds. (By contrast, the Coulomb gauge $\nabla \cdot \mathbf{A} = 0$, which is very popular in analytic considerations, would obviously be of no advantages, since one still has the problem of solving the a solenoidality condition.)

Solving for \mathbf{A} instead of \mathbf{B} has significant advantages, even though this involves taking another derivative. However, the total number of derivatives taken in the code is essentially the same. In fact, when centered finite differences are employed, Alfvén waves are better resolved when \mathbf{A} is used, because then the system of equations for one-dimensional Alfvén waves in the presence of a uniform B_{x0} field in a medium of constant density ρ_0 reduces to

$$\dot{u}_z = \frac{1}{\mu_0\rho_0} B_{x0} A_y'', \quad \dot{A}_y = B_{x0} u_z, \quad (9.136)$$

where a second derivative is taken only once (primes denote x -derivatives). If, instead, one solves for the B_z field, one has

$$\dot{u}_z = \frac{1}{\mu_0\rho_0} B_{x0} B_z', \quad \dot{B}_z = B_{x0} u_z', \quad (9.137)$$

where a first derivative is applied twice, which is far less accurate at small scales if a centered finite difference scheme is used. At the Nyquist frequency, e.g. the first derivative is zero and applying an additional first derivative gives still zero. By contrast, taking a second derivative once gives of course not zero. The use of a staggered mesh circumvents this difficulty. However, such an approach introduces additional complications which hamper the ease with which the code can be adapted to other problems.

Another advantage of using \mathbf{A} is that it is straightforward to evaluate the magnetic helicity $\langle \mathbf{A} \cdot \mathbf{B} \rangle$, which is a particularly important quantity to monitor in connection with dynamo and reconnection problems.

The main advantage of solving for \mathbf{A} is of course that one does not need to worry about the solenoidality of the \mathbf{B} -field, even though one may want to employ irregular meshes or complicated boundary conditions.

As we have emphasized before, when centered meshes are used, it is usually a good idea to avoid taking first derivatives of the same variable twice, because it is more accurate to take instead a second derivative only once. For this reason we calculate the current not as $\mathbf{J} = \mu_0^{-1} \nabla \times (\nabla \times \mathbf{A})$, but as

$$\mathbf{J} = \mu_0^{-1} \left[-\nabla^2 \mathbf{A} + \nabla(\nabla \cdot \mathbf{A}) \right]. \quad (9.138)$$

Taking the gradient of $\nabla \cdot \mathbf{A}$ involves of course also taking first derivatives of the same variable twice, but these contributions are cancelled by corresponding components of the $\nabla^2 \mathbf{A}$ term. There are some advantages relying here on the numerical cancellation, which is of course not exact. The reason is that the full $\nabla^2 \mathbf{A}$ term is important when used in the magnetic diffusion term. If the diagonal terms, $\partial^2 A_x / \partial x^2$, $\partial^2 A_y / \partial y^2$, and $\partial^2 A_z / \partial z^2$, which would all drop out analytically, were taken out there would be no diffusion of \mathbf{A} in the direction of \mathbf{A} .

There is one more aspect that is often useful keeping in mind. There is a particular gauge that allows one to rewrite the uncircled induction equation in such a form that the evolution of \mathbf{A} is controlled by the advective derivative of \mathbf{A} . The calculation is easy. Write the induction term $\mathbf{u} \times \mathbf{B}$ in component form and express \mathbf{B} in terms of \mathbf{A} , so

$$(\mathbf{u} \times \nabla \times \mathbf{A})_i = u_j (\partial_i A_j - \partial_j A_i) = \partial_i(u_j A_j) - A_j \partial_i u_j - u_j \partial_j A_i. \quad (9.139)$$

Here the last term contributes to the advective derivative, the first term can be removed by a gauge transformation and the middle term is a modified stretching term, so the induction equation takes the form

$$\frac{D A_i}{Dt} = -A_j \partial_i u_j - \eta \mu_0 J_i. \quad (9.140)$$

This gauge was used by Brandenburg *et al.* (1995) in order to treat a linear velocity shear using pseudo-periodic (shearing box) boundary conditions. The formulation (9.140) can also be useful when solving the induction equation using lagrangian methods. Note, however, that the non-resistive evolution of \mathbf{A} differs from that of \mathbf{B} in that the indices of the matrix $U_{ij} \equiv \partial u_i / \partial x_j$ are interchanged and that the sign is different; positive for the \mathbf{B} -equation,

$$\frac{DB_i}{Dt} = +U_{ij} B_j + \text{other terms}, \quad (9.141)$$

and negative for the \mathbf{A} -equation,

$$\frac{DA_i}{Dt} = -A_j U_{ji} + \text{other terms}. \quad (9.142)$$

These two formulations are particularly advantageous when the velocity has a constant gradient, as in the case of linear shear. In local simulations of accretion discs, e.g. the shear component is $u_y(x) = -\frac{3}{2}\Omega x$, so $U_{yx} = -\frac{3}{2}\Omega$, and all other U_{ij} vanish. Hence

$$\frac{DA_x}{Dt} = +\frac{3}{2}\Omega A_y + \text{other terms} \quad (9.143)$$

for the **A**-formulation, or

$$\frac{DB_y}{Dt} = -\frac{3}{2}\Omega B_x + \text{other terms} \quad (9.144)$$

for the **B**-formulation. In these two formulations all dependent variables are clearly periodic (or rather pseudo-periodic), so there is no term that is explicitly non-periodic such as $u_y(x) = -\frac{3}{2}\Omega x$. In the following, whenever magnetic fields are present, we use the **A**-formulation, mainly because it guarantees the solenoidality of **B** everywhere (including the boundaries), and also because it is easy to use.

9.5.1. Cache-efficient coding

Unlike the CRAY computers that dominated supercomputing in the eighties and early nineties, all modern computers have a cache that constitutes a significant bottleneck for many codes. This is the case if large three-dimensional arrays are constantly used within each timestep. The advantage of this way of coding is clearly the conceptual simplicity of the code. A more cache-efficient way of coding is to calculate an entire timestep (or a corresponding substep in a three-stage $2N$ -Runge–Kutta scheme) only along a one-dimensional pencil of data within the box. On Linux and Irix architectures, e.g. this leads to a speed-up by 60%. An additional advantage is a drastic reduction in temporary storage that is needed for auxiliary variables within each timestep.

9.6. Application to astrophysical outflows

9.6.1. The isothermal Parker wind

Before discussing outflows from accretion discs it is illuminating to consider first the one-dimensional example of pressure-driven outflows in spherical geometry. A particularly simple case is the *isothermal* wind problem, which is governed by the equations

$$\frac{\partial u}{\partial t} + u \frac{\partial u}{\partial r} = -c_s^2 \frac{\partial \ln \rho}{\partial r} - \frac{\partial \Phi}{\partial r}, \quad (9.145)$$

$$\frac{\partial \ln \rho}{\partial t} + u \frac{\partial \ln \rho}{\partial r} = -\frac{1}{r^2} \frac{\partial}{\partial r}(r^2 u) + \frac{\dot{M} \xi(r)}{\rho}, \quad (9.146)$$

where c_s is the isothermal sound speed (assumed constant), \dot{M} is the mass loss rate, and $\xi(r)$ is a prescribed function of position, normalized such that $\int 4\pi r^2 \xi(r) dr = 1$, and non-vanishing only near $r = 0$. For a point mass the gravity potential Φ would be $-GM/r$, but this becomes singular at the origin. Therefore we use the expression $\Phi = -GM/(r^n + r_0^n)^{1/n}$ instead, where we choose $n = 5$ in all cases, and $1/r_0$ gives the depth of the potential well. In Fig. 9.17 we show radial velocity and density profiles for different values of \dot{M} . Note that the velocity profile is independent of the value of \dot{M} , but the density profile changes by a constant factor. In the steady case the equations can be combined to

$$(u^2 - c_s^2) \frac{d \ln |u|}{dr} = \frac{2c_s^2}{r} - \frac{GM}{r^2}, \quad (9.147)$$

so the sonic point, $|u| = c_s$, is at $r = r^* = GM/2c_s^2$. In Fig. 9.17 we have chosen $GM = 2$ and $c_s = 1$, so $r^* = 1$, which is consistent with the graph of u .

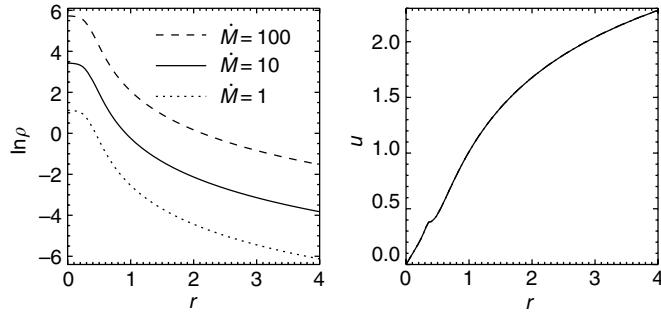


Figure 9.17 Isothermal Parker wind solutions for different values of \dot{M} . Note that the u profile is independent of the value of \dot{M} . $GM = 2$, $c_s = 1$, $r_0 = 0.4$.

9.6.2. The polytropic or adiabatic wind

In the following we make the assumption that the entropy is constant. In that case it is particularly useful to solve for the *potential enthalpy*, $H \equiv h + \Phi$, which varies much less than either h or Φ . Using H as dependent variable is particularly useful if one solves the equations all the way to the origin, $r = 0$, where Φ tends to become singular (or at least strongly negative if a smoothed potential is used). In terms of H the governing equations are

$$\frac{\partial u}{\partial t} + u \frac{\partial u}{\partial r} = -\frac{\partial H}{\partial r} - u \dot{M} \xi, \quad (9.148)$$

$$\frac{\partial H}{\partial t} + u \frac{\partial H}{\partial r} = u \frac{\partial \Phi}{\partial r} + c_s^2 \left[-\frac{1}{r^2} \frac{\partial}{\partial r} (r^2 u) + \frac{\dot{M} \xi(r)}{\rho} \right], \quad (9.149)$$

where $H = h + \Phi$ is the potential enthalpy, $h = p/\rho + e$ is the enthalpy, and $c_s^2 = (\gamma - 1)h = (\gamma - 1)(H - \Phi)$ for a perfect gas, where c_s is the adiabatic sound speed and $h = c_p T$ is the enthalpy. These equations are also valid in the nonisothermal case ($\gamma \neq 1$). The isothermal case may be recovered by putting $\gamma = 1$ and replacing h by $c_s^2 \ln \rho$. In Fig. 9.18 we show solutions for different values of \dot{M} and $\gamma = 5/3$. Again we put $GM = 2$ and $c_{s0} = 1$.

We note that, depending on the strength of the mass source, the polytropic wind problem allows a variety of different velocity and Mach number profiles, whereas for the isothermal wind problem there was only one solution possible, independent of the strength of the mass source. The velocity profile was always the same and also the density was the same up to some scaling factor that changes with \dot{M} . This is connected with the additional degree of freedom introduced through the polytropic constant $K = p/\rho^\gamma$. Since c_s is no longer constant, the position of the sonic point is no longer fixed and different solutions are possible.

In Fig. 9.19 we show solutions where \dot{M} is kept constant, but the depth of the potential well, GM/r_0 , is changed by varying the value of r_0 . Note that the deeper the potential well, the higher the wind speed. The density far away from the source is then correspondingly smaller, so as to maintain the same mass flux.

As we have seen in Section 9.3.6, a polytropic equation of state is unphysical. Therefore we now consider the case where the energy equation is included. To be somewhat more general we consider first the basic equations in conservative form with mass, momentum, and energy

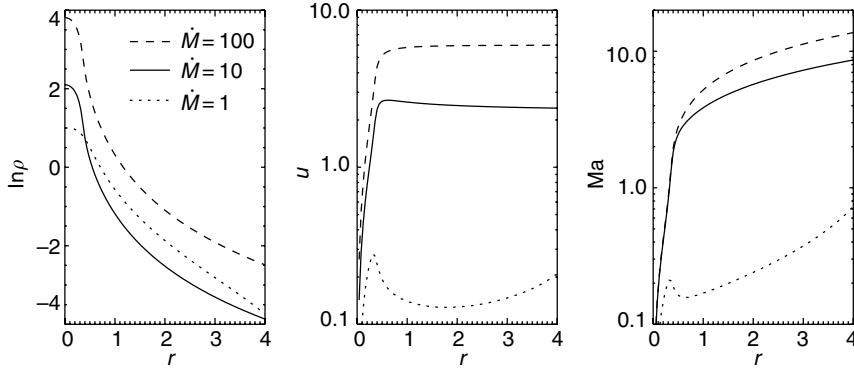


Figure 9.18 Polytropic Parker wind solutions for different values of \dot{M} . $GM = 2$, $c_{s0} = 1$, $r_0 = 0.4$.

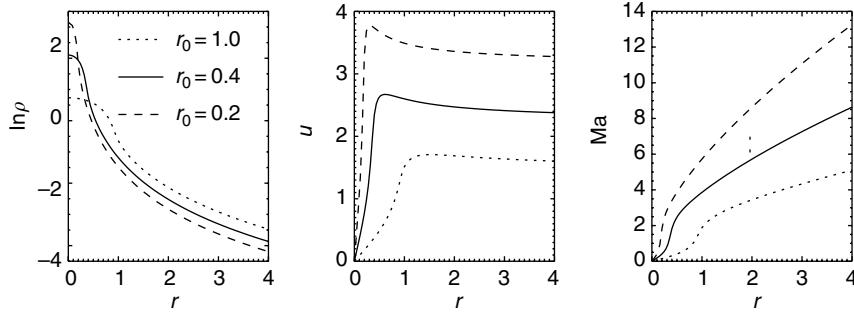


Figure 9.19 Density ρ , velocity u , and Mach number $Ma = u/c_s$ for the polytropic Parker wind solutions for different values of r_0 . $GM = 2$, $c_{s0} = 1$, $\dot{M} = 10$.

sources included, i.e.

$$\frac{\partial \rho}{\partial t} + \frac{\partial}{\partial x_j} (\rho u_j) = \dot{M} \xi, \quad (9.150)$$

$$\frac{\partial}{\partial t} (\rho u_i) + \frac{\partial}{\partial x_j} (\rho u_i u_j + p \delta_{ij} - \tau_{ij}) = \dot{I}_i \xi, \quad (9.151)$$

$$\frac{\partial}{\partial t} \left(\frac{1}{2} \rho \mathbf{u}^2 + \rho e \right) + \frac{\partial}{\partial x_j} \left[u_j \left(\frac{1}{2} \rho \mathbf{u}^2 + \rho h \right) - u_i \tau_{ij} \right] = \dot{E} \xi, \quad (9.152)$$

where \dot{M} , $\dot{\mathbf{I}}$, and \dot{E} are the rates of mass, momentum, and energy injection into the system, $\tau_{ij} = 2\nu\rho \mathbf{S}_{ij}$ is the viscous stress tensor, and \mathbf{S}_{ij} is the (traceless) rate of strain tensor; see (9.12). Rewriting the energy equation in non-conservative form we have

$$\frac{De}{Dt} + \frac{p}{\rho} \nabla \cdot \mathbf{u} \equiv T \frac{Ds}{Dt} = 2\nu \mathbf{S}^2 + \left[\dot{E} - \mathbf{u} \cdot (\dot{\mathbf{I}} - \mathbf{u} \dot{M}) - \left(\frac{1}{2} \mathbf{u}^2 + e \right) \dot{M} \right] \frac{\xi(r)}{\rho}, \quad (9.153)$$

which can also be rewritten in terms of entropy, so the final system of non-conservative equations with source terms is

$$\frac{D \ln \rho}{Dt} + \nabla \cdot \mathbf{u} = \dot{M} \frac{\xi(r)}{\rho}, \quad (9.154)$$

$$\frac{D \mathbf{u}}{Dt} + c_s^2 (\nabla \ln \rho + \nabla s) = \frac{1}{\rho} \nabla \cdot (2\nu\rho \mathbf{S}) + (\dot{\mathbf{I}} - \mathbf{u}\dot{M}) \frac{\xi(r)}{\rho}, \quad (9.155)$$

$$T \frac{Ds}{Dt} = 2\nu \mathbf{S}^2 + [\dot{E} - \mathbf{u} \cdot \dot{\mathbf{I}} + (\frac{1}{2}\mathbf{u}^2 - e)\dot{M}] \frac{\xi(r)}{\rho}, \quad (9.156)$$

where T can be replaced by $c_s^2/(\gamma - 1)$ (remember that $c_p = 1$), and c_s^2 is given by (9.8).

In Fig. 9.20 we present solutions of (9.154)–(9.156) for different values of \dot{E} and \dot{M} . The main effect of varying the value of \dot{E} is to change the value of the entropy in the wind. Outside the acceleration region, however, the value of the entropy is fairly constant, so the polytropic assumption appears to be reasonably good here.

While outflows of some very early-type stars are driven mostly by the $\dot{\mathbf{I}}$ term (resulting from the radiation pressure in lines), the winds of cool stars are driven mostly by the \dot{E} term (resulting from the hot coronae). Similar differences may also explain why some jets are massive (e.g. stellar jets), whilst others are not (jets from active galactic nuclei, e.g. or those anticipated in gamma-ray bursters).

9.6.3. Relevance to outflows and jets

The pressure-driven outflows discussed in the previous section may take the form of more collimated outflows once a magnetic field is involved. This applies to the case of magnetized accretion discs. These discs are generally magnetized both because of dynamo action within

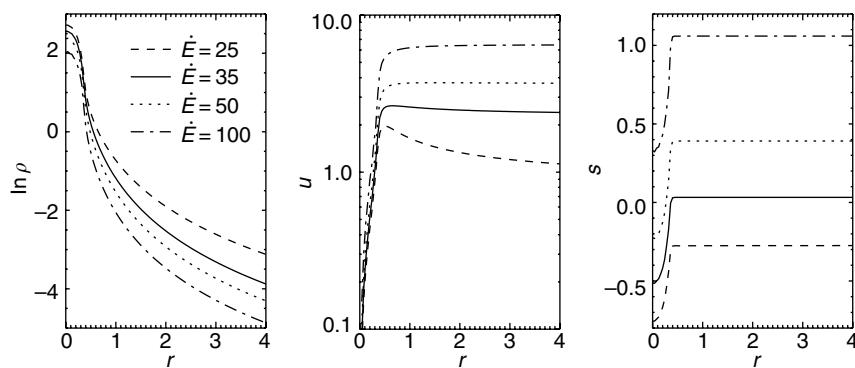


Figure 9.20 Wind solutions for different values of \dot{E} and $\dot{M} = 10$. Note that the solution with $\dot{E} = 35$ is quantitatively very similar to the polytropic solution with the same value of \dot{M} . $GM = 2$, $c_{s0} = 1$, $r_0 = 0.4$.

the disc and because of external fields that were dragged into the disc from outside due to the accretion flow.

At least in some types of jets the outflows may be driven by hot coronae. Other possibilities for driving outflows involve the magneto-centrifugal effect. It is well known that outflows can be driven from a magnetized disc if the angle between the field and the disc is less than 60° (Blandford and Payne, 1982). Recent work in this field was directed to the question whether this angle is the result of some self-regulating process (Ouyed *et al.*, 1997; Ouyed and Pudritz, 1997a,b, 1999) and whether it can be obtained automatically from a dynamo operating within the disc (Campbell, 1999, 2000; Dobler, *et al.*, 1999; Rekowski *et al.*, 2000). This latter question is particularly interesting in view of the fact that jets in star-forming regions are not really pointing in a similar direction (e.g. Hodapp and Ladd, 1995), as one might expect from jet models that start off with a prescribed large scale field.

In Fig. 9.21 we present a particular model of Dobler *et al.* (1999) and Brandenburg (2000); see Brandenburg *et al.* (2000) for a full account of this work. In these models the outflow is driven by mass sources whose strength is proportional to the local density deficit relative to that in the original equilibrium solution of the disc. Such a density deficit was initially caused by slow gas motions that resulted from an instability of the initial equilibrium solution, because a cool disc embedded in a hot corona is non-rotating outside the disc, and it is the resulting vertical shear profile that causes the instability (cf. Urpin and Brandenburg, 1998). At later times, of course, the outflow makes the corona corotating, but by that time the outflow is driven by a persistent density deficit in the disc relative to the initial references solution.

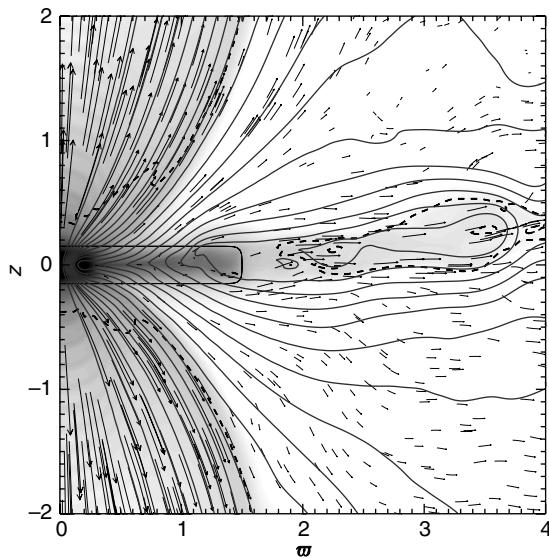


Figure 9.21 Poloidal velocity vectors and poloidal magnetic field lines superimposed on a gray-scale image of the logarithmic density. Dark means high density. The thick dashed line denotes the location where the poloidal flow speed equals the fast magnetosonic speed. The thin solid line gives the location of the disc surface. The slight asymmetry in the field is a relic from the mixed-parity initial condition. (Adapted from Brandenburg, 2000.)

In this model the magnetic field was generated by an $\alpha-\Omega$ dynamo operating within the disc. However α is negative in the upper disc plane (see Brandenburg *et al.*, 1995), and then the most preferred field geometry is dipolar (Campbell, 1999; Rekowski *et al.*, 2000). The field parity is sensitive to details in the disc physics assumed in the particular model (aspect ratio, disc thickness, the presence of outflows, and the conductivity in the disc and the exterior). Nevertheless, both dipolar and quadrupolar fields are equally well able to contribute to wind launching, at least in the outer parts of the disc where the angle between the field and the disc plane is less than 60° , the critical angle for magneto-centrifugal wind launching (Blandford and Payne, 1982). We note, however, that the more detailed analysis of Campbell (1999) suggests that the critical angle can be significantly smaller.

In our models the outflow is only weakly collimated (if at all). This is probably connected with the fact that here the fast magnetosonic surface is rather close to the disc surface, making it difficult for the field to become strong enough to channel the magnetic field. Instead, the field lines themselves are still being controlled too strongly by the outflow. However, outflows with rather large opening angles are actually seen in some star-forming regions; see Greenhill *et al.* (1998).

While most of the disc mass is ejected in a cone of half-opening angle around 25° , most of the disc angular momentum is ejected at rather low latitudes, almost in the direction of the disc plane away from the central object. The timescales for these various processes are comparable. In Fig. 9.22 we show the azimuthally integrated mass flux, angular momentum flux, and magnetic (Poynting) flux as a function of polar angle, and compare with a non-magnetic run. We find that in the magnetic run the outflow is more strongly concentrated towards the axis. Also, the amount of angular momentum loss (dash-dotted line) is larger when the disc is magnetized. We emphasize in particular that in the magnetic run significant amounts of magnetic field are ejected from the system. In the following section we discuss the significance of such magnetic flux ejection for magnetizing the interstellar medium into which the outflow is streaming. This discussion is similar to a corresponding discussion for the contamination of the intergalactic medium via outflows from active galactic nuclei (Brandenburg, 2000).

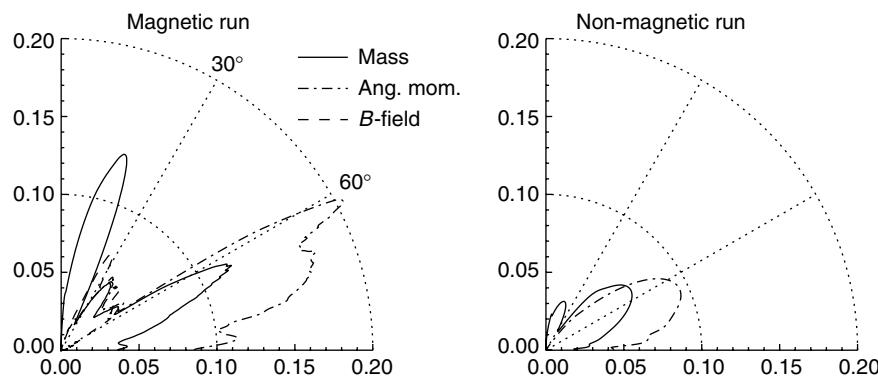


Figure 9.22 Comparison of the angular dependence of azimuthally integrated fluxes for magnetic and non-magnetic outflows. The solid line refers to mass flux, the dashed line to angular momentum flux, and the dash-dotted line (in the second panel) corresponds to the Poynting flux. The units of all quantities are thus $1/[t]$.

9.6.4. Magnetic contamination from outflows

It may at first appear somewhat unrealistic to expect significant magnetization of the interstellar medium from outflows. However, the following calculation shows that the effect may be quite significant. Assume that every star did undergo a phase of strong accretion with associated outflows, so $N = 10^{11}$ for the whole galaxy. The duration of intense outflow activity is 10^5 years, say, but it could even be 10^6 years. The magnetic luminosity is $L_{\text{mag}} = 0.05 \dot{M}_w c_s^2$ (Brandenburg *et al.*, 2000), where $c_s \approx 10$ km/s is the average sound speed of the interstellar medium, and $\dot{M}_w = 0.1 \dot{M}_d$ (see Pelletier and Pudritz, 1992), where $\dot{M}_d \approx 10^{-7} M_\odot/\text{year}$ is a conservative estimate for the disc accretion rate. Again this value may be larger. With the above numbers the magnetic luminosity from all $N = 10^{11}$ sources is then $N L_{\text{mag}} = 7 \times 10^{39}$ erg/s and the total energy output delivered from all stars at some early point in the life time is therefore $E_{\text{mag}} = \tau N L_{\text{mag}} = 2 \times 10^{52}$ erg. Diluting this over a volume of a galaxy of 300 kpc^3 (radius 10 kpc, height 1 kpc) gives 2×10^{-15} erg/cm³. Multiplying this by 8π and taking the square root gives $0.2 \mu\text{G}$. Expressed more concisely in a formula we have for the rms magnetic field strength

$$\langle \mathbf{B}^2 \rangle^{1/2} \approx \left(8\pi \frac{F_{\text{Poy}}}{F_{\text{kin}}} \frac{N \dot{M}_w c_s^2}{V} \delta t \right)^{1/2}, \quad (9.157)$$

where the efficiency factor $F_{\text{Poy}}/F_{\text{kin}}$ ($=0.05$ in our model) may be lower in systems where the disc dynamo is less strong.

The parameters for a corresponding estimate for outflows from young galactic discs (active galactic nuclei) are as follows. Assuming $N \sim 10^4$ galaxies per cluster, each with $\dot{M}_w \approx 0.1 M_\odot/\text{year} = 10^{25}$ g/s, and $c_s = 1000$ km/s for the sound speed in the intracluster gas, the rate of magnetic energy injection for all galaxies together is $L_{\text{mag}} = 10^{44}$ erg/s. Distributing this over the volume of the cluster of $V \sim 1 \text{ Mpc}^3$, and integrating over a duration of $\delta t = 1 \text{ Gyr}$, this corresponds to a mean magnetic energy density of $\langle \mathbf{B}^2 / 8\pi \rangle \approx 10^{-13}$ erg/cm³, so $\langle \mathbf{B}^2 \rangle^{1/2} \approx 10^{-6}$ G, which is indeed of the order of the field strength observed in galaxy clusters. We note that our estimate has been rather optimistic in places (e.g. \dot{M}_w could be lower, or the relevant δt could be shorter), but it does show that outflows are bound to produce significant magnetization of the intracluster gas and the interstellar medium (see also Völk and Atoyan, 1999). In the latter case it will provide a good seed field for the galactic dynamo. A dynamo is still necessary to shape the magnetic field and to prevent it from decaying in the galactic turbulence. Similarly, many galaxy clusters undergo merging and this too can enhance and reorganize the magnetic field. The necessity for a recent merger event would also be consistent with the fact that not all halos are observed to have strong magnetic fields. Recent simulations by Roettiger *et al.* (1999) suggest that after a merger the field strength may increase by a factor of at least 20 (and this value increases with improving observational resolution).

As an alternative consideration for causing the magnetization in clusters of galaxies, *primordial* magnetic fields are sometimes discussed. There are numerous mechanisms that could generate relatively strong fields at an early time, e.g. during inflation (age $\sim 10^{-36}$ s) or during the electroweak phase transition (age $\sim 10^{-10}$ s). Such fields would now still be at a very small scale if one considers only the cosmological expansion. However, depending on the degree of magnetic helicity in this primordial field, the magnetic energy can be transferred to larger scales that are now on the scale of galaxies. For a recent discussion of these results see Brandenburg (2001a).

9.7. Hydromagnetic turbulence and dynamos

As mentioned in the beginning, accurate high-order schemes are essential in all applications to turbulent flows. Nevertheless, we should mention that one often attempts solutions of the inviscid and nonresistive equations using low-order finite differences combined with monotonicity schemes that result in some kind of effective diffusion. The piece-wise parabolic method (PPM) of Colella and Woodward (1984) is an example. However, unlike the Smagorinsky scheme (see Chan and Sofia, 1986, 1989; Steffen *et al.*, 1989; Fox *et al.*, 1991 for applications to convection simulations), PPM and similar methods cannot be proven to converge to the original Navier–Stokes equation in the limit of infinite resolution. Nevertheless, they are rather popular in astrophysical gas simulations. These schemes are rather robust and have also been applied to high-resolution simulations of compressible turbulence (Porter *et al.*, 1992, 1994). While the results from those simulations are generally quite plausible, the power spectrum shows a k^{-1} subrange at large wavenumbers, which is still not fully understood. This was sometimes regarded as an artifact of PPM, and should therefore only occur at small scales. However, as the resolution was increased further (up to 1024^3), the k^{-1} subrange just became more extended.

A similar feature was found in cascade models of turbulence when the ordinary ∇^2 diffusion operator was replaced by a $-\nabla^4$ “hyperdiffusion” operator (Lohse and Müller-Groeling, 1995). Whatever the outcome of this puzzle is, it is clear that with schemes that cannot be proven to converge to the actual Navier–Stokes equations in the limit of infinite resolution, there would always remain some uncertainty and debate. On the other hand, especially in the incompressible case the use of hyperviscosity does generally allow the exploration of larger Reynolds numbers and broader inertial ranges.

MHD simulations with the highest resolution to date have been performed by Biskamp and Müller (1999), who considered decaying turbulence with and without magnetic helicity. They found that in the presence of magnetic helicity the magnetic energy decay is significantly slower. In particular, they found the magnetic energy decays like $t^{-1/2}$, as opposed to t^{-1} found earlier by Mac Low *et al.* (1998) for compressible turbulence.

Before we start discussing dynamo action in turbulence simulations representative of more astrophysical settings, such as accretion discs and the solar convection zone, let us first illustrate the mechanism of the inverse cascade that is believed to be an important ingredient of large scale magnetic field generation.

9.7.1. Isotropic MHD turbulence

Most developments in the theory of turbulence have been carried out under the assumptions of homogeneity and isotropy. This is certainly true of the work on the inverse cascade (or turbulent cascades in general), but it is also true of much of the work on the α -effect which – like the inverse cascade – describes the generation of large scale fields. However, unlike the inverse cascade process, the energy comes here directly from the velocity field at the scales of the energy-carrying eddies and not from the velocity and magnetic field at successively smaller scales, which are usually larger than the scale of the energy-carrying eddies.

It is not easy to see whether any of these effects is actually responsible for the large scale field generation in astrophysical bodies or even the simulations. In simulations of accretion disc turbulence there is certainly some evidence for the presence of an α -effect, but it is extremely noisy (Brandenburg *et al.*, 1995; Brandenburg and Donner, 1997; Ziegler and Rüdiger, 2000). Evidence for the inverse magnetic cascade comes mostly from the magnetic

energy spectra (Balsara and Pouquet, 1999; Brandenburg, 2001b), which show a marked peak at large scales, but this is convincing only in cases where the flow is driven at a wavenumber that is clearly larger than the smallest wavenumber in the box. In practice, e.g. in convectively driven turbulence, the flow is driven at all scales including the large scale making it difficult to see a marked peak at the smallest wavenumber (see a corresponding discussion in Meneguzzi and Pouquet, 1989).

From the seminal papers of Frisch *et al.* (1975) and Pouquet *et al.* (1976) it is clear that amplification of large scale fields can also be explained by an inverse cascade of magnetic helicity. In those papers the authors also showed that the inverse cascade is a consequence of the fact that the magnetic helicity $\langle \mathbf{A} \cdot \mathbf{B} \rangle$, is conserved by the nonresistive equations. (\mathbf{A} is the magnetic vector potential giving the magnetic field as $\mathbf{B} = \nabla \times \mathbf{A}$.) The inverse magnetic cascade effect too is rather difficult to isolate in simulations of astrophysical turbulence. However, under somewhat more idealized conditions, e.g. when magnetic energy is injected at high wavenumbers, one clearly sees how the magnetic energy increases at large scales; see Fig. 9.23. Further details of this model have been published in the proceedings of the helicity meeting in Boulder (Brandenburg, 1999).

In the model considered above the flow was forced magnetically. This may be motivated by the recent realization that strong magnetic field generation in accretion discs can be facilitated by *magnetic instabilities*, such as the Balbus–Hawley instability. Other examples of magnetic instabilities include the magnetic buoyancy instability, which can lead to an α -effect (e.g. Brandenburg and Schmitt, 1998; Thelen, 2000), and the reversed field pinch which also leads to a dynamo effect (e.g. Ji *et al.*, 1996). Before returning to the accretion disc dynamo in Section 9.7.9 we should emphasize that strong large scale field generation is also possible with purely hydrodynamic forcing. Simulations in this type were considered recently by Brandenburg (2001b). There are many similarities compared with the case of

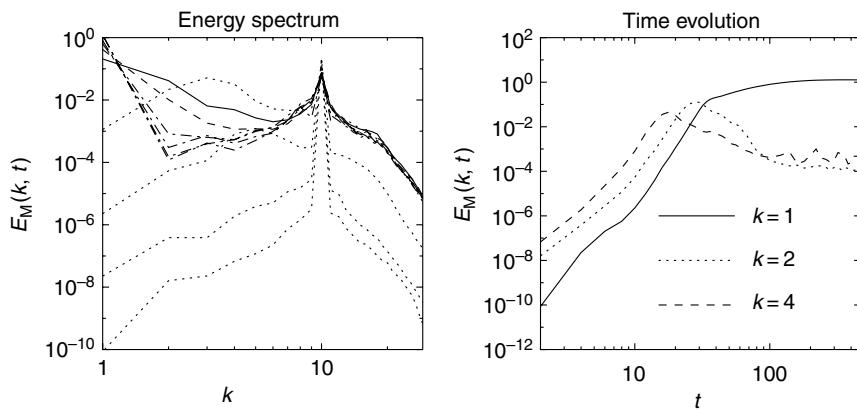


Figure 9.23 Spectral magnetic energy, $E_M(k, t)$, as a function of wavenumber k for different times: dotted lines are for early times ($t = 2, 4, 10, 20$), the solid and dashed lines are for $t = 40$ and 60 , respectively, and the dotted-dashed lines are for later times ($t = 80, 100, 200, 400$). Here magnetic energy is injected at wavenumber 10. Note the occurrence of a sharp secondary peak of spectral magnetic energy at $k = 10$. By the time the energy at $k = 1$ has reached equipartition the energies in $k = 2$ and $k = 4$ become suppressed.

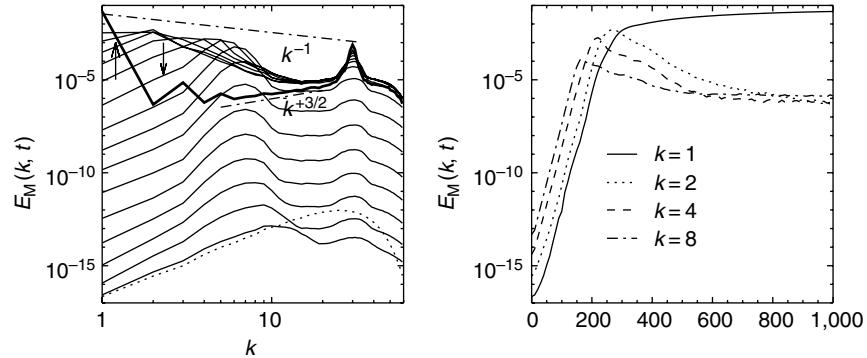


Figure 9.24 Left: Magnetic energy spectra for a run with forcing at $k = 30$. The times range from 0 (dotted line) to 10, 30, ..., 290 (solid lines). The thick solid line gives the final state at $t = 1000$. Note that at early times the spectra peaks at $k_{\max} \approx 7$. The k^{-1} and $k^{+3/2}$ slopes are given for orientation as dash-dotted lines. Right: Evolution of spectral magnetic energy for selected wavenumbers in a simulation with hydrodynamical forcing at $k = 30$.

magnetic forcing. The evolution of magnetic energy spectra in the presence of hydrodynamic forcing is shown in Fig. 9.24. Like in the case of magnetic forcing (Fig. 9.23) there are marked peaks both at the forcing scale and at the largest scale of the box. Furthermore, the evolution of spectral energy at the largest scales shows similar behavior: the magnetic energy with wavenumber $k = 8$ increases, reaches a maximum, and begins to decrease when the magnetic energy at $k = 4$ reaches a maximum. The same happens for the next larger scales (wavenumbers $k = 4$ and 2, until the scale of the box (with $k = 1$) is reached.

The *suppression* of magnetic energy at intermediate scales, $2 \leq k \leq 8$, is quite essential for the development of a well-defined large scale field. In a recent letter Brandenburg and Subramanian (2000) showed that this type of *self-cleaning* effect can also be simulated by using ambipolar diffusion as nonlinearity and ignoring the Lorentz force altogether. Without any nonlinearity, however, there would be no interaction between different scales and the magnetic energy would increase at all scales, especially at small scales, which would soon swamp the large scale field structure with small scale fields.

The model presented in Fig. 9.24 has large scale separation in the sense that there is a large gap between the forcing wavenumber ($k = k_f = 30$) and the wavenumber of the box ($k = k_1 = 1$). One sees that during the growth phase there is a clear secondary maximum at $k = 7$. This is indeed expected for an α^2 dynamo, whose maximum growth rate is at $k_{\max} = \frac{1}{2}\alpha/\eta_T$, where η_T is the total (turbulent plus microscopic) magnetic diffusion coefficient.

The disadvantage of a high forcing wavenumber is that for modest resolution (here we used 120^3 meshpoints) no inertial range can develop. This is different if once forces at $k_f = 5$, keeping otherwise the same resolution. In Fig. 9.25 we show spectra for different cases with $k_f = 5$ where we compare the results for different values of the magnetic Reynolds and magnetic Prandtl number. In Fig. 9.26 we show cross-sections of one field component at different times. In this model (Run 3 of Brandenburg, 2001b) the forcing is at $k_f = 5$, so there is now a clear tendency for the build-up of an inertial range in $8 \leq k \leq 25$.

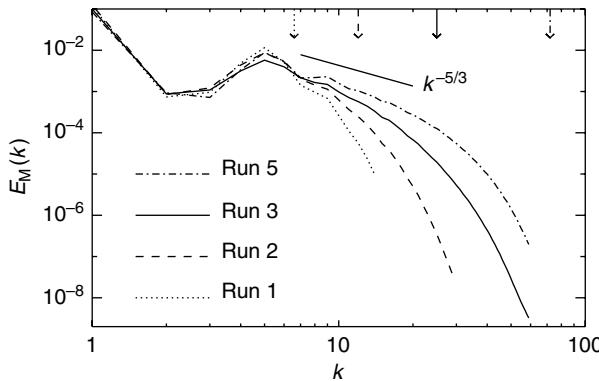


Figure 9.25 Comparison of time averaged magnetic energy spectra for Runs 1–3 ($t = 600 - 1000$) with a non-averaged spectrum for Run 5 (large magnetic Prandtl number) taken at $t = 1600$. To compensate for different field strengths and to make the spectra overlap at large scales, two of the three spectra have been multiplied by a scaling factor. There are clear signs of the gradual development of an inertial subrange for wavenumbers larger than the forcing scale. The $k^{-5/3}$ slope is shown for orientation. The dissipative magnetic cutoff wavenumbers, $\langle \mathbf{J}^2 / \eta^2 \rangle^{1/4}$, are indicated by arrows at the top.

9.7.2. The inverse cascade in decaying turbulence

We now turn to the case of *decaying* turbulence, which is driven only by an initial kick to the system. There are several circumstances in astrophysics where this could be relevant: early universe, neutron stars, and mergers of galaxy clusters. In all those cases one is interested in the development of large scale fields. In the context of the early universe the possibility of energy conversion from small to large scale fields was pointed out by Brandenburg *et al.* (1996) who found that fields generated at the horizon scale of 3 cm after the electroweak phase transition would now have a scale on the order of kiloparsecs, even though the cosmological expansion alone would only lead to scales on the order of 1 AU. These results were only based on either two-dimensional simulations or three-dimensional cascade model calculation (e.g. Biskamp, 1994). Therefore we now turn to fully three-dimensional simulations.

In the absence of any forcing and with no kinetic energy initially an initial magnetic field can only decay. However, if initially most of the magnetic energy is in the small scales, there is the possibility that magnetic helicity and thereby also magnetic energy is transferred to large scales. This is exactly what happens (Fig. 9.27), provided there is initially some net helicity. The inset of Fig. 9.27 shows that in the absence of initial net helicity the field at large scales remains unchanged, until diffusion kicks in and destroys the remaining field at very late times.

If the magnetic field has the possibility to tap energy also from the large scale velocity the situation is somewhat different again and there is the possibility that a large scale magnetic field can also be driven without net helicity. In that case the large scale field can increase due to dynamo action from the incoherent $\alpha-\Omega$ -effect (Vishniac and Brandenburg, 1997). In astrophysical settings there is usually large scale shear from which energy can be tapped.

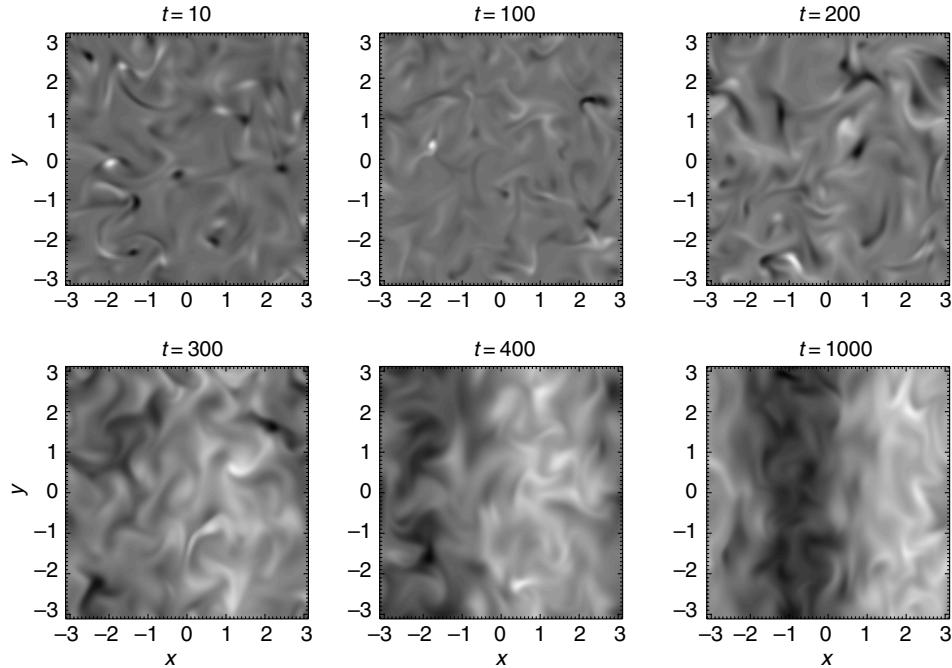


Figure 9.26 Gray-scale images of cross-sections of $B_x(x, y, 0)$ for Run 3 of Brandenburg (2001b) at different times showing the gradual build-up of the large scale magnetic field after $t = 300$. Dark (light) corresponds to negative (positive) values. Each image is scaled with respect to its min and max values.

Before we discuss simulations with imposed shear in more detail we first present a simple argument that makes the link between the inverse cascade and helicity conservation.

9.7.3. The connection with magnetic helicity conservation

In the following we give a simple argument due to Frisch *et al.* (1975) that helps to understand why the magnetic helicity conservation property leads to the occurrence of an inverse cascade. We define in the following magnetic energy and helicity spectra, $M(k)$ and $H(k)$, respectively. Now, because of Schwartz inequality, we have

$$|\hat{\mathbf{B}}(k)|^2 = |(i\mathbf{k} \times \hat{\mathbf{A}}) \cdot \hat{\mathbf{B}}| \geq |\mathbf{k}| |\hat{\mathbf{A}} \cdot \hat{\mathbf{B}}| \quad (9.158)$$

we have a lower bound on the spectral magnetic energy at each wavenumber $k = \mathbf{k}$. In terms of shell integrated magnetic energy and helicity spectra this corresponds to

$$M(k) \geq \frac{1}{2} k |H(k)|, \quad (9.159)$$

where the 1/2-factor comes simply from the 1/2-factor in the definition of the magnetic energy. Assuming that two wavenumbers q and p interact such that they produce power at a new wavenumber k , then

$$M(p) + M(q) = M(k), \quad H(p) + H(q) = H(k). \quad (9.160)$$

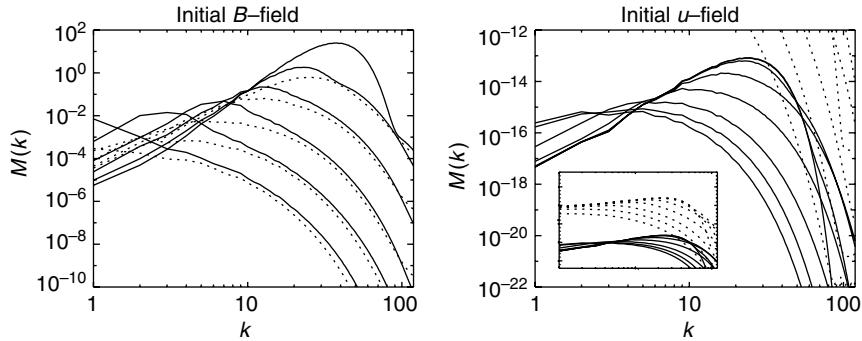


Figure 9.27 Power spectra of magnetic energy (solid lines) and kinetic energy (dotted lines) in a decay run with helicity. The left-hand panel is for a case where the flow is only driven by an initial helical magnetic field. In the right-hand panel the field is weak and governed by strong decaying fluid turbulence. The inset shows both velocity and magnetic spectra in the same plot. The Prandtl number ν/η is equal to one in both cases, but the mesh Reynolds number, which is kept constant at all times, is different in the two cases: 20 in the left-hand panel and 50 in the right-hand panel. The times are 0, 0.01, 0.1, etc., till $t = 10^2$ in the left-hand panel and $t = 10^3$ in the right-hand panel.

For simplicity we consider the case $p = q$, so

$$2M(p) = M(k), \quad 2H(p) = H(k). \quad (9.161)$$

Assume also that initially the constraint was sharp (maximum helicity), then

$$M(p) = \frac{1}{2}pH(p). \quad (9.162)$$

Now, from the constrain again we have

$$\frac{1}{2}kH(k) \leq M(k) = 2M(p) = pH(p) = \frac{1}{2}pH(k), \quad (9.163)$$

so

$$k \leq p, \quad (9.164)$$

that is the wavenumber of the target result must be larger or equal to the wavenumbers of the initial field.

The argument given above is of course quite rough, because it ignores, e.g. the detailed angular dependence of the wave vectors. This was taken into account properly already in the early paper by Pouquet *et al.* (1976), but this approach was based on closure assumptions for the higher moments, which is in principle open to criticism. Thus, numerical simulations, like those presented above, are necessary for an independent confirmation that the inverse cascade really works. In this connection one should mention that there are some parallels with the inverse cascade of enstrophy in two-dimensional hydrodynamic (non-magnetic) turbulence. In that case the enstrophy (i.e. the mean squared vorticity) is conserved because of the absence of vortex stretching in two dimensions. The inverse hydrodynamic cascade has some significance in meteorology and perhaps in low aspect ratio convection experiments,

where one finds a peculiar energy and entropy spectrum that is referred to as Bolgiano scaling; see Brandenburg (1992) and Suzuki and Toh (1995) for corresponding shell model calculations and Toh and Iima (2000) for direct simulations.

9.7.4. Inverse cascade or α -effect?

In Section 9.7.1 we made a distinction between inverse cascade and α -effect in the sense that, although both lead to large scale field generation, in the inverse cascade there is a gradual transfer of magnetic helicity and energy to ever larger scales, whereas the α -effect produces large scale magnetic fields directly from small scale fields. Thus, the distinction is really one between local and non-local inverse cascades.

In Fig. 9.28 we show the normalized spectral energy transfer function $T(k, p, t)$ for $k = 1$ and 2 as a function of p , and at different times t . The index k signifies the gain or losses of the field at wavenumber k , and the index p indicates the wavenumber of the velocity from which the energy comes from. This function shows that most of the energy of the large scale field at $k = 1$ comes from velocity and magnetic field fluctuations at the forcing scale, which is here $k = k_f = 5$. At early times this is also true of the energy of the magnetic field at $k = 2$, but at late times, $t = 1000$, the gain from the forcing scale, $k = 5$, has diminished, and instead there is now a net loss of energy into the next larger scale, $k = 3$, suggestive of a direct cascade operating at $k = 2$, and similarly at $k = 3$.

Based on these results we may conclude that in the saturated state the magnetic energy at $k = 1$ is sustained by a *non-local* inverse cascade from the forcing scale directly to the largest scale of the box. This is characteristic of the α -effect of mean-field electrodynamics, except that here nonlinearity plays an essential role in isolating the large scale from the small scale “magnetic trash”, as Parker used to say.

A closer look at Fig. 9.24, where $k = k_f = 30$, suggests that once the scale separation is large enough the energy is at first transferred not to the scale of the box, but instead to a somewhat smaller scale (here at wavenumber $k = 7$). Following the corresponding discussion in Brandenburg (2001b), this wavenumber is close to the wavenumber, $k_{\max} = \frac{1}{2}|\alpha|/\eta_T$, where the α^2 dynamo grows fastest.

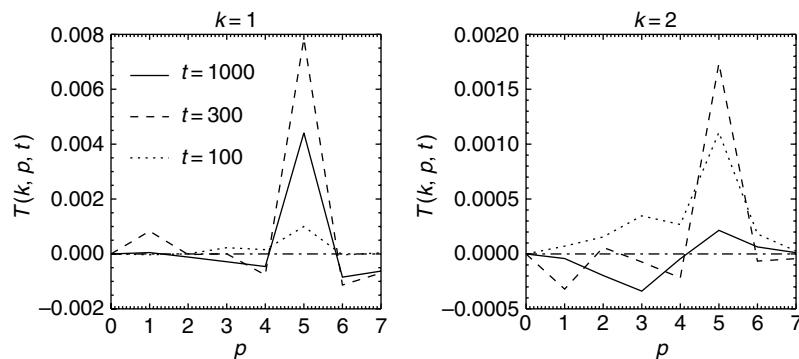


Figure 9.28 Spectral energy transfer function $T(k, p, t)$, normalized by $\langle \mathbf{B}^2 \rangle$ for three different times, for $k = 1$ and 2. Run 3 of Brandenburg (2001b).

In the following section we address the issue of magnetic helicity conservation which has important consequences for the timescale after which the large scale field begins to develop. This has also a bearing on the widely discussed controversy of the so-called “catastrophic α -quenching” of Vainshtein and Cattaneo (1992).

9.7.5. Approximate helicity conservation

The magnetic helicity, $H = \langle \mathbf{A} \cdot \mathbf{B} \rangle$, is conserved by the non-resistive MHD equations. For a closed or periodic box $\langle \mathbf{A} \cdot \mathbf{B} \rangle$ satisfies the equation

$$\frac{d}{dt} \langle \mathbf{A} \cdot \mathbf{B} \rangle = -2\eta\mu_0 \langle \mathbf{J} \cdot \mathbf{B} \rangle, \quad (9.165)$$

where $\langle \mathbf{J} \cdot \mathbf{B} \rangle$ is the current helicity, and angular brackets denote volume averages. Note that for a periodic box $\langle \mathbf{A} \cdot \mathbf{B} \rangle$ is gauge invariant, i.e. $\langle \mathbf{A} \cdot \mathbf{B} \rangle$ does not change after a gauge transformation, $\mathbf{A} \rightarrow \mathbf{A} + \nabla\varphi$. This is a direct consequence of the solenoidality of the magnetic field, because $\langle \nabla\varphi \cdot \mathbf{B} \rangle = -\langle \varphi \nabla \cdot \mathbf{B} \rangle = 0$ owing to $\nabla \cdot \mathbf{B} = 0$.

In order to judge whether $\langle \mathbf{A} \cdot \mathbf{B} \rangle$ is small or large we calculate the length scale

$$\ell_H = |\langle \mathbf{A} \cdot \mathbf{B} \rangle| / \langle \mathbf{B}^2 \rangle. \quad (9.166)$$

In Fig. 9.29 we see that the evolution of ℓ_H proceeds in three distinct phases: (i) a very short period ($t < 1$) where ℓ_H is very small and comparable to the numerical noise level, so magnetic helicity almost perfectly conserved, (ii) an intermediate interval ($2 < t < 200$) where ℓ_H is much larger, but still only roughly equal to the mesh size of the calculation, and then (iii) a regime where ℓ_H is of order unity. The latter is only possible because of the presence of helicity in the system, which leads to a large scale magnetic field configuration that is nearly force-free.

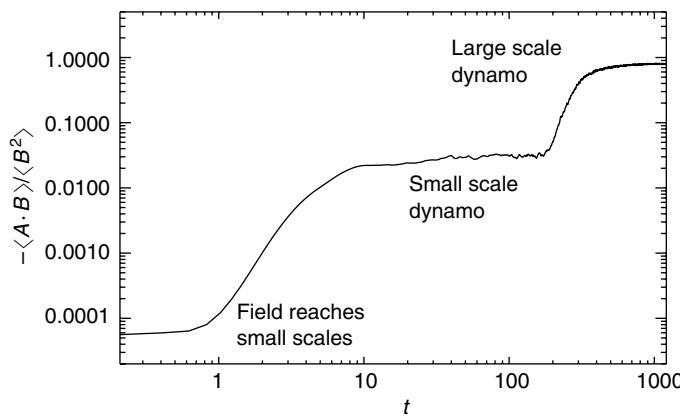


Figure 9.29 Evolution of the (negative) magnetic helicity length scale in a double-logarithmic plot. Note the presence of three distinct phases: very approximate helicity conservation near zero, followed by a phase of larger magnetic helicity scale (three orders of magnitude), and finally a phase where the magnetic helicity scale has reached the scale of the box.

9.7.6. Resistively limited growth of the large scale field

The approximate conservation of magnetic helicity has an important consequence for the generation of large scale fields: in order to build up a large scale field with magnetic helicity one has to *change* the value of $\langle \mathbf{A} \cdot \mathbf{B} \rangle$ from its initial value of zero to its certain final value. This final value of $\langle \mathbf{A} \cdot \mathbf{B} \rangle$ is such that the length scale ℓ_H is close to the maximum value possible for a certain geometry. It also implies that when interpreting the results in terms of mean-field electrodynamics (α -effect and turbulent diffusion), the α -effect must be quenched early on, well before the field has reached its final value. Nevertheless, this strong quenching, which was first anticipated by Vainshtein and Cattaneo (1992) and later confirmed numerically by Cattaneo and Hughes (1996), does not prevent the field from reaching its final super-equipartition field strength.

Before we come to the details we mention already now that the helicity constraint is probably too severe to be acceptable for astrophysical conditions, so one must look for possible escape routes. The most plausible way of relaxing the helicity constraint is in allowing for open boundary conditions (Blackman and Field, 2000; Kleeorin *et al.*, 2000), but the situation can still be regarded as inconclusive. Given that much of the work on large scale dynamos so far assumes periodic boundaries, we shall now consider this particular case in more detail. In the periodic case the final field geometry can be, e.g. of the form $\bar{\mathbf{B}} = B_0(\cos k_1 z, \sin k_1 z, 0)$, where $k_1 = 1$ is the smallest possible wavenumber in the box, B_0 is the field amplitude, and $\langle \bar{\mathbf{B}}^2 \rangle = B_0^2$. Alternatively, the field may vary in the x or y direction, and there may be an arbitrary phase shift; examples of these possibilities have been reported in Brandenburg (2001b). Anyway, for $\bar{\mathbf{B}} = B_0(\cos k_1 z, \sin k_1 z, 0)$ the corresponding vector potential is $\bar{\mathbf{A}} = -(B_0/k_1)(\cos k_1 z, \sin k_1 z, 0) + \nabla\phi$, where ϕ is an arbitrary gauge which does not affect the value of $\langle \bar{\mathbf{A}} \cdot \bar{\mathbf{B}} \rangle$. In this example we have

$$\langle \bar{\mathbf{A}} \cdot \bar{\mathbf{B}} \rangle = -\langle \bar{\mathbf{B}}^2 \rangle / k_1, \quad (9.167)$$

where we have included the k_1 factor, even though in the present case $k_1 = 1$. (The minus sign in (9.167) would turn into a plus if the forcing had negative helicity.) The mean current density is given by $\bar{\mathbf{J}} = -B_0 k_1 (\cos k_1 z, \sin k_1 z, 0)$, so the current helicity of the mean field is given by

$$\mu_0 \langle \bar{\mathbf{J}} \cdot \bar{\mathbf{B}} \rangle = -k_1 \langle \bar{\mathbf{B}}^2 \rangle. \quad (9.168)$$

Before we can use (9.167) and (9.168) in (9.165) we need to relate the magnetic and current helicities of the mean field to those of the actual field. We can generally split up the two helicities into contributions from large and small scales, i.e.

$$\langle \mathbf{A} \cdot \mathbf{B} \rangle = \langle \bar{\mathbf{A}} \cdot \bar{\mathbf{B}} \rangle + \langle \mathbf{a} \cdot \mathbf{b} \rangle, \quad (9.169)$$

$$\langle \mathbf{J} \cdot \mathbf{B} \rangle = \langle \bar{\mathbf{J}} \cdot \bar{\mathbf{B}} \rangle + \langle \mathbf{j} \cdot \mathbf{b} \rangle. \quad (9.170)$$

As the large scale magnetic field begins to saturate, the magnetic helicity has to become constant and so (9.165) dictates that $\langle \mathbf{J} \cdot \mathbf{B} \rangle$ must go to zero in the steady state. Consequently, the contribution from $\langle \mathbf{j} \cdot \mathbf{b} \rangle$ must be as large as that of $\langle \bar{\mathbf{J}} \cdot \bar{\mathbf{B}} \rangle$, and of opposite sign, so that the two cancel. This, together with (9.168), allows us immediately to write down an expression for the equilibrium strength of the mean field;

$$\langle \bar{\mathbf{B}}^2 \rangle = \mu_0 |\langle \mathbf{j} \cdot \mathbf{b} \rangle| / k_1, \quad (9.171)$$

which is now valid for both signs of the helicity of the forcing. The *residual* helicity (Pouquet *et al.*, 1976),

$$H_{\text{res}} = \langle \boldsymbol{\omega} \cdot \mathbf{u} \rangle - \langle \mathbf{j} \cdot \mathbf{b} \rangle / \rho_0, \quad (9.172)$$

is small in the nonlinear saturated state and nearly vanishing. [We mention that this is also the case in the models of Brandenburg and Subramanian (2000).] Furthermore, for forced turbulence with a well defined forcing wavenumber the kinetic helicity may be estimated as $\langle \boldsymbol{\omega} \cdot \mathbf{u} \rangle \approx k_f \langle \mathbf{u}^2 \rangle$. Together with (9.171) we have

$$\langle \bar{\mathbf{B}}^2 \rangle = \frac{k_f}{k_1} B_{\text{eq}}^2, \quad (9.173)$$

where $B_{\text{eq}}^2 = \mu_0 \langle \rho \mathbf{u}^2 \rangle$ and B_{eq} is the equipartition field strength. Thus, the mean field can exceed (!) the equipartition field by the factor $(k_f/k_1)^{1/2}$. This estimate agrees well with the results of the simulations; see Brandenburg (2001b).

Using (9.169) and (9.170) together with (9.167) and (9.168) we can rewrite (9.165) in the form

$$\frac{d}{dt} \langle \bar{\mathbf{B}}^2 \rangle = -2\eta k_1^2 \langle \bar{\mathbf{B}}^2 \rangle + 2\eta k_1 \mu_0 |\langle \mathbf{j} \cdot \mathbf{b} \rangle|, \quad (9.174)$$

where we have taken into account the contribution of the small scale current helicity which is of similar magnitude as the large scale current helicity. For the magnetic helicity, on the other hand, the small scale contribution is negligible, because

$$|\langle \mathbf{a} \cdot \mathbf{b} \rangle| \approx \mu_0 |\langle \mathbf{j} \cdot \mathbf{b} \rangle| / k_f^2 \approx \mu_0 |\langle \bar{\mathbf{J}} \cdot \bar{\mathbf{B}} \rangle| / k_f^2 \approx |\langle \bar{\mathbf{A}} \cdot \bar{\mathbf{B}} \rangle| \left(\frac{k_1}{k_f} \right)^2 \ll |\langle \bar{\mathbf{A}} \cdot \bar{\mathbf{B}} \rangle|. \quad (9.175)$$

After the saturation at small and intermediate scales the small scale current helicity is approximately constant and can be estimated as

$$|\langle \mathbf{j} \cdot \mathbf{b} \rangle| \approx \rho_0 |\langle \boldsymbol{\omega} \cdot \mathbf{u} \rangle| \approx k_f \langle \rho \mathbf{u}^2 \rangle = k_f B_{\text{eq}}^2 / \mu_0. \quad (9.176)$$

The solution of (9.174) is given by

$$\langle \bar{\mathbf{B}}^2 \rangle = \epsilon_0 B_{\text{eq}}^2 \left[1 - e^{-2\eta k_1^2 (t - t_{\text{sat}})} \right], \quad (9.177)$$

where ϵ_0 is a coefficient which, in the present model with a well-defined forcing wavenumber, can be approximated by $\epsilon_0 \approx k_f / k_1$.

This is indeed also the limiting behavior found for α^2 -dynamos with simultaneous α and η quenching of the form

$$\alpha = \frac{\alpha_0}{1 + \alpha_B \bar{\mathbf{B}}^2 / B_{\text{eq}}^2}, \quad \eta_t = \frac{\eta_{t0}}{1 + \eta_B \bar{\mathbf{B}}^2 / B_{\text{eq}}^2}, \quad (9.178)$$

where $\alpha_B = \eta_B$ is assumed. Assuming that the magnetic energy density of the mean field, $\bar{\mathbf{B}}^2$, is approximately uniform (which is well satisfied in the simulations) we can obtain the solution $\bar{\mathbf{B}} = \bar{\mathbf{B}}(t)$ of (9.301) in the form

$$\bar{\mathbf{B}}^2 / (1 - \bar{\mathbf{B}}^2 / B_{\text{fin}}^2)^{1+\lambda/\eta k_1^2} = B_{\text{ini}}^2 e^{2\lambda t}, \quad (9.179)$$

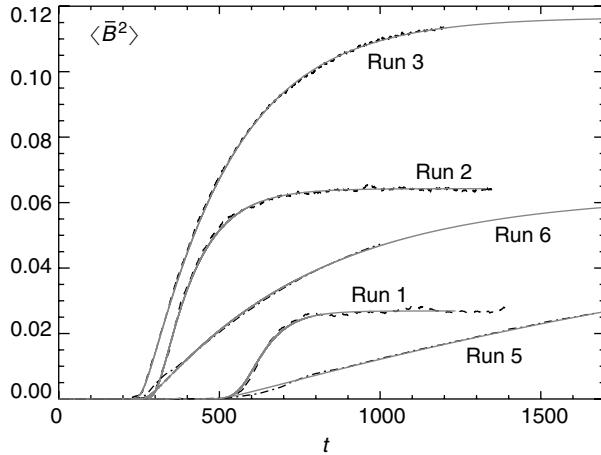


Figure 9.30 Evolution of $\langle \bar{B}^2 \rangle$ for Runs 1–3, 5, and 6, compared with the solution (9.179) of the dynamo equations using (9.179).

where

$$\lambda = |\alpha_0|k_1 - \eta_{T0}k_1^2 \quad (9.180)$$

is the kinematic growth rate of the dynamo, B_{ini} is the initial field strength, and B_{fin} is the final field strength of the large scale field, which is related to α_B and η_B via

$$\alpha_B = \eta_B = \frac{\lambda}{\eta k_1^2} \left(\frac{B_{\text{eq}}}{B_{\text{fin}}} \right)^2. \quad (9.181)$$

The full derivation is given in Appendix E. The significance of this result is that it provides an excellent fit to the numerical simulations; see Fig. 9.30 where we present the evolution of $\langle \bar{B}^2 \rangle$ for the different runs of Brandenburg (2001b). Equation (9.179) can therefore be used to extrapolate to astrophysical conditions. The time it takes to convert the small scale field generated by the small scale dynamo to a large scale field, τ_{eq} , increases linear with the magnetic Reynolds number, R_m . Apart from some coefficients of order unity the ratio of τ_{eq} to the turnover time is therefore just R_m . For the sun this ratio would be 10^8 – 10^{10} . However, before interpreting this result further one really has to know whether or not the presence of open boundary conditions could alleviate the issue of very long timescales for the mean magnetic field. Furthermore, it is not clear whether the long timescales discussed above have any bearing on the cycle period in the case of oscillatory solutions. The reason this is not so clear is because for a cyclic dynamo the magnetic helicity in each hemisphere stays always of the same sign and is only slightly modulated. It is likely that this modulation pattern is advected precisely with the meridional circulation, in which case the helicity could be nearly perfectly conserved in a lagrangian frame. This could provide an interesting clue for why the solar dynamo is migrating. The relation between meridional circulation and dynamo wave propagation has been advocated by Durney (1995) and Choudhuri *et al.* (1995), but helicity conservation would strongly lock the two aspects.

It is clear that virtually all astrophysical bodies are open, allowing for constant loss of magnetic helicity. In the case of the sun significant amounts of magnetic helicity are indeed

observed at the solar surface (Berger and Ruzmaikin, 2000). Significant losses of magnetic helicity are particularly obvious in the case of accretion discs which are almost always accompanied by strong outflows that can sometimes be collimated into jets. Thus, dynamo action from accretion disc turbulence would be a good candidate for clarifying the significance of open boundaries on the nature of the dynamo. Another reason why accretion disc turbulence is a fruitful topic for understanding dynamo action is because the shear is extremely strong. In the following we discuss some recent progress that has been made in this field.

9.7.7. Joule dissipation from mean and fluctuating fields

In an MHD flow the mean magnetic Joule dissipation per unit volume is given by

$$Q_{\text{Joule}} = \eta \mu_0 \langle \mathbf{J}^2 \rangle. \quad (9.182)$$

Whilst in many astrophysical flows η may be very small, $|\mathbf{J}|$ can be large so that Q_{Joule} remains finite even in the limit $\eta \rightarrow 0$. One example where this is very important is accretion discs, where Joule dissipation (together with viscous dissipation) are important in heating the disc. These viscous and resistive processes are indeed the only significant sources of energy supply in discs, and yet the luminosities of discs that result from the conversion of magnetic and kinetic energies into heat and radiation can be enormous. Much of the work on discs involves mean-field modelling, so it would be interesting to see how the Joule dissipation, $Q_{\text{Joule}}^{(\text{mf})}$, predicted from a mean-field model,

$$Q_{\text{Joule}}^{(\text{mf})} = \eta_t \mu_0 \langle \bar{\mathbf{J}}^2 \rangle, \quad (9.183)$$

relates to the actual Joule dissipation. In Fig. 9.31 we show the evolution of actual and mean-field Joule dissipation and compare with an estimate for the rate of total energy dissipation, B_{eq}^2/τ , where τ is the turnover time. Here we have taken into account that η_t is “catastrophically” quenched using the formulae of Brandenburg (2001a) with the parameters for Run 3.

There is no reason *a priori* that the magnetic energy dissipations from the mean-field model should agree with the actual one. It turns out that the mean-field dissipation is a fourth of the actual one, so it is definitely significant. It would therefore be interesting to see how those two dissipations compare with each other in other models.

9.7.8. Possible pitfalls in connection with hyperresistivity

In many astrophysical applications hyperresistivity and hyperviscosity are sometimes used in order to concentrate the effects of magnetic diffusion and viscosity to the smallest possible scale. The purpose of this section is to highlight possible spurious artifacts associated with this procedure. As we have seen above, large scale dynamos can depend on the microscopic magnetic diffusivity and must therefore be affected when it is replaced by hyperresistivity. The resulting modifications that are to be expected are easily understood: on the right hand side of (9.165) the term $\langle \mathbf{J} \cdot \mathbf{B} \rangle$ needs to be replaced by $\langle (\nabla^4 \mathbf{A}) \cdot \mathbf{B} \rangle$. This leads to a change of the relative importance of small and large scale contributions, which therefore changes (9.173) to

$$\langle \bar{\mathbf{B}}^2 \rangle = \left(\frac{k_f}{k_l} \right)^3 B_{\text{eq}}^2. \quad (9.184)$$

Thus, the final field strength will be even larger than before: instead of a factor of 5 super-equipartition (for $k_f = 5$) one now expects a factor of 125. Recent simulations by Brandenburg and Sarson (2001) have indeed confirmed this tendency. The main conclusion is that hyper-resistivity can therefore be used to address certain issues regarding large magnetic Reynolds numbers that are otherwise still inaccessible. On the other hand, the results are in some ways distorted and need therefore be interpreted carefully.

9.7.9. Remarks on accretion disc turbulence

We have already mentioned the possibility of dynamo action in accretion discs. Accretion discs have been postulated some 30 years ago in order to explain the incredibly high luminosities of quasars. Only in the past few years has direct imaging of accretion discs become possible, mostly due to the Hubble Space Telescope. Accretion discs form in virtually all collapse processes, such as galaxy and star formation. In the latter case the central mass is of the order of one solar mass, while in the former it is around 10^8 solar masses and is concentrated in such a small volume that it must be a black hole. If the surrounding matter was nonrotating, it would fall radially towards the center. But this is unrealistic and even the slightest rotation relative to the central object would become important eventually as matter falls closer to the center.

If there was no effective diffusive process in discs, the angular momentum of the matter would stay with the gas parcels, and since the gravity force is balanced by the corresponding centrifugal force, the gas would never accrete. However, the angular velocity of the gas follows a $r^{-3/2}$ Kepler law, so the gas is differentially rotating and one may expect shear instabilities to occur that would drive turbulence and hence turbulent dissipation. Unfortunately, however, the story is not so simple. Discs are both linearly stable (Stewart, 1975) and probably also nonlinearly stable (Hawley *et al.*, 1996). Nevertheless, in the presence of a magnetic field there is a powerful *linear* instability (Balbus and Hawley, 1991), and subsequent work has shown that this instability is indeed capable of driving the instability and hence turbulence.

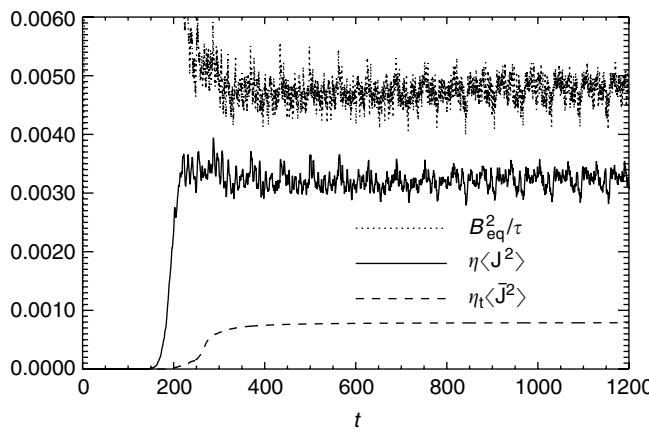


Figure 9.31 Joule dissipation for Run 3 (solid line), compared with the Joule dissipation estimated for a corresponding mean-field model (dashed line). An estimate for the rate of total energy dissipation, B_{eq}^2/τ , is also given.

One of the key outcomes of such simulations is the rate of turbulent dissipation, which determines the rate of angular momentum transport and correspondingly the rate at which orbital kinetic and potential energy is liberated in the form of heat. This is normally expressed in terms of a turbulent viscosity (e.g. Frank *et al.*, 1992), but it may equally well be expressed in terms of the horizontal components of the Reynolds and Maxwell stress tensors. The stress may then be normalized by $\langle \rho c_s^2 \rangle$ to give a non-dimensional measure (called α_{SS}) of the ability of the turbulence to transport angular momentum outward (if $\alpha_{SS} > 0$). This α_{SS} is indeed always positive, see Fig. 9.32, but it fluctuates significantly about a certain mean value. These fluctuations are in fact correlated with the energy in the mean magnetic field, $\langle \mathbf{B} \rangle^2$, as is shown in the right hand panel of Fig. 9.32. This mean magnetic field shows regular reversals combined with a migration away from the midplane, as can be seen in Fig. 9.33.

The evolution of the mean magnetic field found in the simulations is reminiscent of the behavior known from mean-field $\alpha-\Omega$ dynamos. Further details regarding this correspondence (relation between the value of α and cycle period, field parity for different boundary conditions, etc.) can be found in recent reviews of the subject (e.g. Brandenburg, 1998, 2000).

9.7.10. Connection with the solar dynamo problem

The disc simulations have shown that a global large scale field can be obtained even in cartesian geometry. The detailed behavior of this large scale field depends of course on the boundary conditions adopted (Brandenburg, 1998), and will therefore be different in different geometries. Nevertheless, the very fact that large scale dynamo action is possible already in simple cartesian geometry is interesting.

In Fig. 9.34 we show the evolution of magnetic and kinetic energies as well as the magnitudes of the large scale field for a simulation of a convectively driven dynamo in the presence of large scale shear (Brandenburg *et al.*, 2001). It turns out that the ratio of the magnetic energies in large scale fields relative to the total field, $\langle \mathbf{B} \rangle^2 / \langle \mathbf{B}^2 \rangle$, which is a measure of the filling factor of the magnetic field, is around 15% when the field has reached saturation, i.e. when the field growth has stopped. This is similar to the case of isotropic nonmirror-symmetric turbulence considered in Section 9.7.1. On the average, however, the magnetic field is then directed

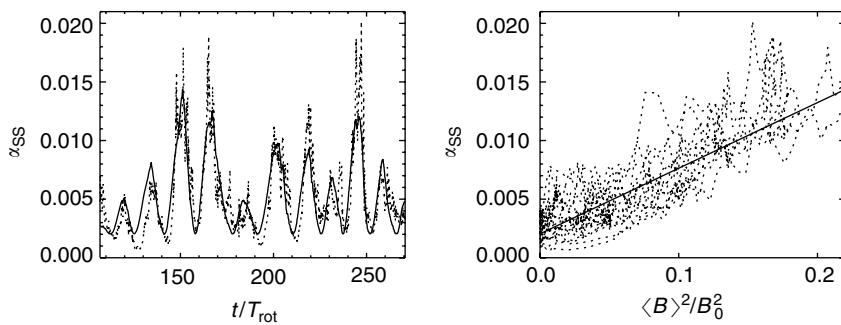


Figure 9.32 Dependence of α_{SS} on time and the mean magnetic field strength for a local accretion disc model (Run B of Brandenburg *et al.*, 1996a). Here $B_0 = (\mu_0 \rho c_s^2)^{1/2}$ is the thermal equipartition field strength and $T_{\text{rot}} = 2\pi/\Omega$ the local rotation period. In the left hand panel the dotted line represents the actual data and the solid line gives the fit obtained by correlating α_{SS} with the mean magnetic field (right-hand panel).

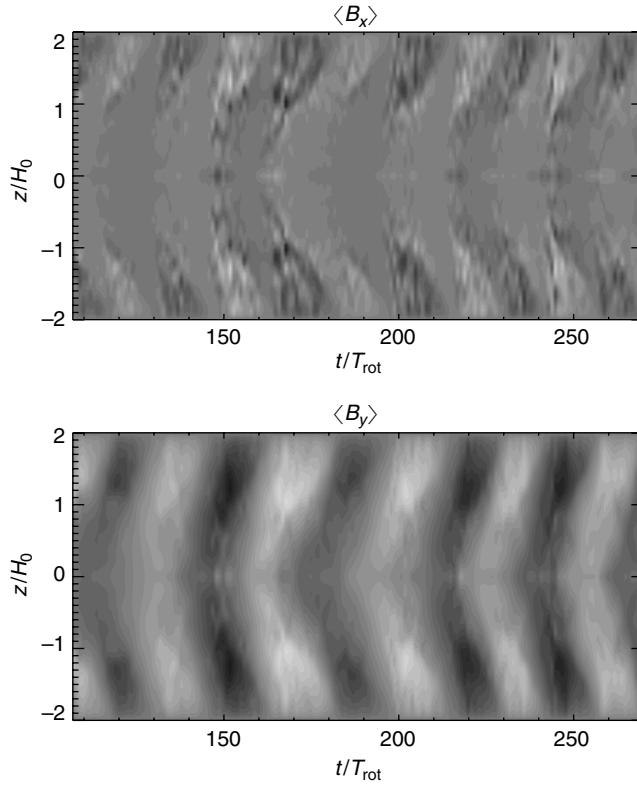


Figure 9.33 Butterfly (space-time) diagram of the poloidal and toroidal magnetic field components averaged over the two horizontal (x and y) directions for the local accretion disc model of (Brandenburg *et al.*, 1996a, Run B). Note that the poloidal field is much more noisy than the toroidal field, and that there is a clear outward migration of magnetic field.

into the negative y -direction (corresponding to the negative azimuthal direction in spherical geometry), but there is a weak and more noisy cross-stream field component directed in the positive x -direction (pointing north).

When the field has reached saturation, the mean field direction is approximately constant. Although the magnitude of this mean field fluctuates somewhat, the sign is always the same. Thus, this simulation shows no cycles, which are so characteristic of the solar dynamo. However, since those features, including the field geometry depend strongly on boundary condition and on the location of the boundary conditions, this disagreement is to be expected, and one would really need to resort to global simulations in spherical geometry.

9.7.11. Dynamics of the overshoot layer

Late-type stars with outer convection zones have an interface between the convection zone proper and the radiative interior. This leads to some additional dynamics that is important to include, especially in connection with the dynamo problem. This interface is the layer where magnetic flux can accumulate, i.e. *not* necessarily the layer where the dynamo operates;

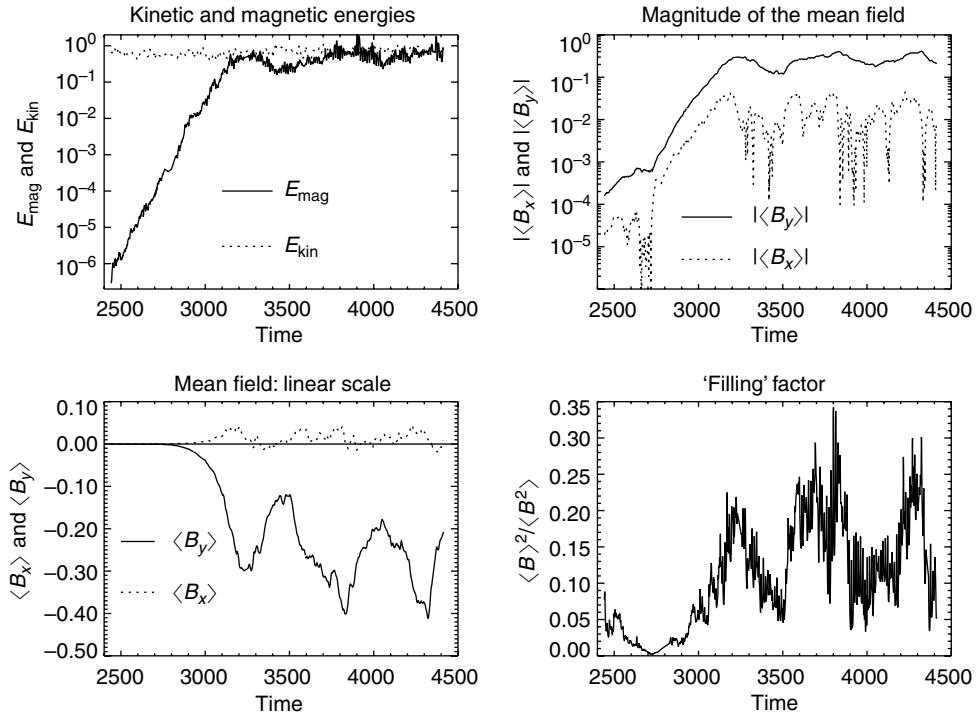


Figure 9.34 Evolution of several quantities for a convective dynamo model with shear: kinetic and magnetic energies (dotted and solid lines; first panel), mean latitudinal and toroidal fields (dotted and solid lines; second panel), mean magnetic field in a linear scale (third panel), and the filling factor (fourth panel). Energies and mean magnetic fields are given in units of the equipartition value, $B_{\text{eq}} = 4\pi \langle \mathbf{u}^2 \rangle$.

see the discussion in Brandenburg (1994). The accumulation is a consequence of turbulent pumping down the turbulence intensity gradient and the effect was seen clearly in video animations reported by Brandenburg and Tuominen (1991) and was analyzed in detail by Nordlund *et al.* (1992). Tobias *et al.* (1998) have studied the effect in isolation starting with an initial magnetic field distribution as opposed to a dynamo-generated field.

The flow dynamics changes drastically as one enters this overshoot layer. The stabilizing buoyancy effect provides a restoring force on a downward moving element, which can give rise to gravity waves that could be driven by individual plumes. This leads to a marked wavy pattern that can extend deep into the lower overshoot layer, as seen in Fig. 9.35 where we have plotted the vertical rms velocity as a function of depth and time. These waves extend a major fraction into the stably stratified layer beneath the convection zone, but are damped eventually. The typical period of such events is seen to be around 20 (in units of $\sqrt{d/g}$), where d is the depth of the unstable (convective) layer. This is comparable with the mean Brunt–Väisälä frequency,

$$N_{\text{BV}}^2 = -\mathbf{g} \cdot \nabla(s/c_p). \quad (9.185)$$

which is around 0.3 in the overshoot layer; see Fig. 9.36.

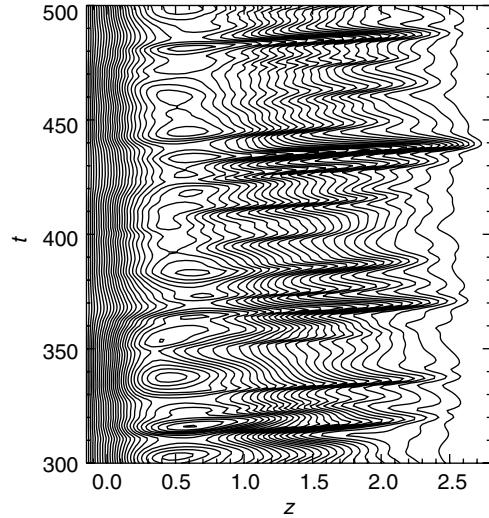


Figure 9.35 Space-time diagram of the vertical rms velocities for a nonmagnetic convection zone model of Brandenburg *et al.* (2001). Note the propagation of isolated plumes in more or less regular time intervals. Note also that the wavy pattern extends well into the convection zone proper ($0.5 \leq z \leq 1$), and that the plumes appear to propagate at an approximately constant speed towards the bottom. This speed is around 0.1, which is comparable to the rms velocity in the runs.

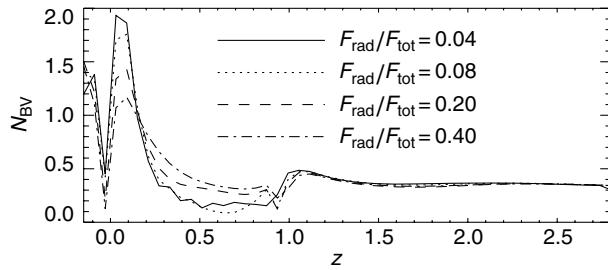


Figure 9.36 Modulus of the Brunt–Väisälä frequency for a run with polytropic index $m = -0.9$ and a nondimensional input flux $\mathcal{F}_{\text{conv}} = 0.01$. The various curves are for different values of the radiative flux, but fixed nominal convective flux. Small values of $F_{\text{rad}}/F_{\text{tot}}$ are typical for the upper layers of the solar convection zone.

Technically the presence of a lower overshoot layer provides a formidable challenge, because there the dynamics is governed by the very slow thermal timescale. This can lead to problems if the properties of the upper surface layers change which could affect the entropy in deeper parts of the convection zone. This in turn will also affect the stratification in the rest of the radiative interior. Since this can only happen on a thermal timescale (there is no turbulence in these layers to speed things up) it takes very long before one arrives at a new statistically stable state. This problem is of course also encountered when starting with initial conditions that are derived under too unrealistic assumptions. There are essentially two

different approaches to this: one either considers a toy model where dynamical and thermal timescales are artificially brought closer together or one adopts an implicit code which allows the use of somewhat longer timesteps. The former approach implies that one adopts input fluxes that exceed those in real stars, but the good news is that turbulent velocities and temperature fluctuations vary with changing flux in exactly the way that is expected from mixing length theory; see Brandenburg *et al.* (2001) for details.

9.8. Conclusions

Many phenomena in astrophysics show direct manifestations of turbulence. As in the case of accretion discs, without turbulence there would be no enhanced dissipation, no heating of the disc, and hence no emission. Magnetic fields are of major importance as is evidenced again by the example of accretion discs, where the turbulence is a direct consequence of the presence of magnetic fields (Balbus–Hawley instability). Although magnetic fields in discs could have primordial origin and could just have been compressed during their formation, it is also clear that discs are actually a favorable candidate for producing strong large scale magnetic fields, as shown by the local simulations discussed above.

Other bodies where strong dynamo action is possible are stars. Simulations of convection have shown that strong small scale magnetic fields are possible (Meneguzzi and Pouquet, 1989; Nordlund *et al.*, 1992; Brandenburg *et al.*, 1996b; Cattaneo, 1999), but if there is strong shear an intense large scale field can also emerge.

The use of non-conservative high-order schemes has proved useful in all those applications. They are easy to implement and to modify, but they are also reasonably accurate. In this chapter we have illustrated the behavior of such schemes using various test problems. Using potential enthalpy and entropy as the main thermodynamic variables has a number of advantages, especially in connection with strongly stratified flows near a central object with a deep potential well, which is relevant to studying outflow phenomena. Contrary to common belief, high-order schemes are not more vulnerable to Gibbs phenomena near discontinuities. Instead, in simple advection tests high-order methods are able to produce smoother solutions with less viscosity, which is important for accurate modeling of turbulence.

In the last part of this chapter we have briefly mentioned some astrophysical applications of simulations using high-order schemes where hydromagnetic turbulence played an important role. In the next few years we may expect a dramatic increase in the quality and predictive power of such simulations, as larger computers become available. Already now a number of very promising simulations are emerging. There is important work addressing the stability of astrophysical jets in three dimensions (Ouyed *et al.*, 2000). Also worthwhile mentioning are recent high resolution simulations by Hawley (2000) of three-dimensional accretion tori in global geometry. What remains to be done in this field is a proper connection between disc physics and the launching mechanism of jets. This would require incorporating proper thermodynamics allowing for radiative cooling and magnetic heating in particular. Global simulations would also be highly desirable to address the global stellar dynamo problem. For example, it would be interesting to see how the dynamo works in fully convective stars. This problem is in some ways simpler than the solar dynamo problem, because one does not need to worry about the lower overshoot layer where the relevant timescales are much longer than in the convection zone proper.

Appendix A. Centered, one-sided and semi-one-sided derivatives

In Section 9.3 we gave the centered finite difference formulae for schemes of different order. Here we first describe the method for determining the finite difference formulae for second order, but the generalization to higher order is straightforward. We also give the corresponding expressions for one-sided and semi-one-sided finite difference formulae.

We want to write the derivative $f'(x)$ at the point x_i as

$$f'_i = a_{-1} f_{i-1} + a_0 f_i + a_1 f_{i+1}, \quad (9.186)$$

where $f'_i = f'(x_i)$, $f_{i-1} = f(x_i - \delta x)$, and $f_{i+1} = f(x_i + \delta x)$. To determine the coefficients a_{-1} , a_0 , and a_1 we expand $f(x)$ up to second order

$$f(x) = c_0 + c_1 x + c_2 x^2. \quad (9.187)$$

The first derivative is then

$$f'(x) = c_1 + 2c_2 x. \quad (9.188)$$

In particular, the value at $x = 0$ is just $f'(0) = c_1$. Likewise, we have $f''(0) = 2c_2$. To determine all coefficients we make use of our knowledge at the neighboring points around x_i , i.e. we use the function values $f(x_i - \delta x) \equiv f_{i-1}$, $f(x_i) \equiv f_i$, and $f(x_i + \delta x) \equiv f_{i+1}$, so we have

$$f_{i-1} = c_0 + c_1(-\delta x) + c_2(-\delta x)^2, \quad (9.189)$$

$$f_i = c_0, \quad (9.190)$$

$$f_{i+1} = c_0 + c_1(+\delta x) + c_2(+\delta x)^2. \quad (9.191)$$

This can be written in matrix form

$$\begin{pmatrix} f_{i-1} \\ f_i \\ f_{i+1} \end{pmatrix} = \begin{pmatrix} (-1)^0 & (-1)^1 & (-1)^2 \\ 0^0 & 0^1 & 0^2 \\ 1^0 & 1^1 & 1^2 \end{pmatrix} \begin{pmatrix} c_0 \\ c_1 \delta x \\ c_2 \delta x^2 \end{pmatrix} \quad (9.192)$$

(where $(-1)^0 = 0^0 = 1^0 = 1$), or

$$\mathbf{f} = \mathbf{M}\mathbf{c}, \quad (9.193)$$

and so we obtain the coefficients as

$$\mathbf{c} = \mathbf{M}^{-1}\mathbf{f}. \quad (9.194)$$

To calculate f' we need the value of c_1 , see (9.188), and so the coefficients a_n needed to express the derivative are $a_{-1} = (\mathbf{M}^{-1})_{10}$, $a_0 = (\mathbf{M}^{-1})_{11}$, and $a_1 = (\mathbf{M}^{-1})_{12}$, i.e. all points of the inverted matrix in the second row. The resulting formula for f'_i is well known,

$$f'_i = (-f_{i-1} + f_{i+1})/(2\delta x). \quad (9.195)$$

The corresponding result for the second derivative is

$$f''_i = (f_{i-1} - 2f_i + f_{i+1})/(\delta x^2). \quad (9.196)$$

On the boundaries we have to calculate for derivative using only points inside the domain, which is explained in the next subsection for second order accuracy, but again the generalization to higher order is straightforward and only the results will be listed.

A.1. One-sided 2nd-order derivatives

Again, we want to write the derivative $f'(x)$ as

$$f'_i = a_0 f_0 + a_1 f_1 + a_2 f_2, \quad (9.197)$$

but now

$$\begin{pmatrix} f_i \\ f_{i+1} \\ f_{i+2} \end{pmatrix} = \begin{pmatrix} 0^0 & 0^1 & 0^2 \\ 1^0 & 1^1 & 1^2 \\ 2^0 & 2^1 & 2^2 \end{pmatrix} \begin{pmatrix} c_0 \\ c_1 \delta x \\ c_2 \delta x^2 \end{pmatrix}. \quad (9.198)$$

Thus, one arrives at

$$f'_i = (-3 f_i + 4 f_{i+1} - f_{i+2}) / (2 \delta x). \quad (9.199)$$

Correspondingly, for the second derivative we have

$$f''_i = (2 f_i - 5 f_{i+1} + 4 f_{i+2} - f_{i+3}) / \delta x^2. \quad (9.200)$$

A.2. 4th-order derivatives

First derivatives

$$f'_i = (f_{i-2} - 8 f_{i-1} + 8 f_{i+1} - f_{i+2}) / (12 \delta x), \quad (9.201)$$

$$f'_i = (-3 f_{i-1} - 10 f_i + 18 f_{i+1} - 6 f_{i+2} + f_{i+3}) / (12 \delta x), \quad (9.202)$$

$$f'_i = (-25 f_i + 48 f_{i+1} - 36 f_{i+2} + 16 f_{i+3} - 3 f_{i+4}) / (12 \delta x). \quad (9.203)$$

Second derivatives

$$f''_i = (-f_{i-2} + 16 f_{i-1} - 30 f_i + 16 f_{i+1} - f_{i+2}) / (12 \delta x^2), \quad (9.204)$$

$$f''_i = (11 f_{i-1} - 20 f_i + 6 f_{i+1} + 4 f_{i+2} - f_{i+3}) / (12 \delta x^2), \quad (9.205)$$

$$f''_i = (35 f_i - 104 f_{i+1} + 114 f_{i+2} - 56 f_{i+3} + 11 f_{i+4}) / (12 \delta x^2). \quad (9.206)$$

A.3. 6th-order derivatives

First derivatives

$$f'_i = (-f_{i-3} + 9 f_{i-2} - 45 f_{i-1} + 45 f_{i+1} - 9 f_{i+2} + f_{i+3}) / (60 \delta x), \quad (9.207)$$

$$f'_i = (2 f_{i-2} - 24 f_{i-1} - 35 f_i + 80 f_{i+1} - 30 f_{i+2} + 8 f_{i+3} - f_{i+4}) / (60 \delta x), \quad (9.208)$$

$$\begin{aligned} f'_i = & (-10 f_{i-1} - 77 f_i + 150 f_{i+1} - 100 f_{i+2} + 50 f_{i+3} \\ & - 15 f_{i+4} + 2 f_{i+5}) / (60 \delta x), \end{aligned} \quad (9.209)$$

$$\begin{aligned} f'_i = & (-147 f_i + 360 f_{i+1} - 450 f_{i+2} + 400 f_{i+3} - 225 f_{i+4} \\ & + 72 f_{i+5} - 10 f_{i+6}) / (60 \delta x). \end{aligned} \quad (9.210)$$

Second derivatives

$$f_i'' = (2f_{i-3} - 27f_{i-2} + 270f_{i-1} - 490f_i + 270f_{i+1} - 27f_{i+2} + 2f_{i+3})/(180\delta x^2), \quad (9.211)$$

$$f_i'' = (-13f_{i-2} + 228f_{i-1} - 420f_i + 200f_{i+1} + 15f_{i+2} - 12f_{i+3} + 2f_{i+4})/(180\delta x^2), \quad (9.212)$$

$$f_i'' = (137f_{i-1} - 147f_i - 255f_{i+1} + 470f_{i+2} - 285f_{i+3} + 93f_{i+4} - 13f_{i+5})/(180\delta x^2), \quad (9.213)$$

$$f_i'' = (812f_i - 3132f_{i+1} + 5265f_{i+2} - 5080f_{i+3} + 2970f_{i+4} - 972f_{i+5} + 137f_{i+6})/(180\delta x^2). \quad (9.214)$$

Appendix B. The $2N$ -RK3 scheme

If N is the number of variables to be updated from one timestep to the next, the $2N$ -schemes require only $2 \times N$ variables to be stored in memory at any time. This is better than for the standard Runge–Kutta schemes. The general iteration formula is

$$w_i = \alpha_i w_{i-1} + hF(t_{i-1}, u_{i-1}), \quad u_i = u_{i-1} + \beta_i w_i. \quad (9.215)$$

For 3-step scheme we have $i = 1, \dots, 3$. In order to advance the variable u from $u^{(n)}$ at time $t^{(n)}$ to $u^{(n+1)}$ at time $t^{(n+1)} = t^{(n)} + h$ we set in (9.215)

$$u_0 = u^{(n)} \quad \text{and} \quad u^{(n+1)} = u_3, \quad (9.216)$$

with u_1 and u_2 being intermediate steps. In order to be able to calculate the first step, $i = 1$, for which no $w_{i-1} \equiv w_0$ exists, we have to require $\alpha_1 = 0$. Thus, we are left with five unknowns, $\alpha_2, \alpha_3, \beta_1, \beta_2$, and β_3 . We write down (9.215) in explicit form for $i = 1, \dots, 3$,

$$w_1 = hF(t_0, u_0), \quad u_1 = u_0 + \beta_1 w_1, \quad (9.217)$$

$$w_2 = \alpha_2 w_1 + hF(t_1, u_1), \quad u_2 = u_1 + \beta_2 w_2, \quad (9.218)$$

$$w_3 = \alpha_3 w_2 + hF(t_2, u_2), \quad u_3 = u_2 + \beta_3 w_3. \quad (9.219)$$

Written in explicit form, we have for $i = 1$

$$u_1 = u_0 + \beta_1 hF(t_0, u_0). \quad (9.220)$$

The $i = 2$ step yields

$$w_2 = \alpha_2 hF(t_0, u_0) + hF(t_1, u_1), \quad (9.221)$$

$$u_2 = u_0 + h[(\beta_1 + \beta_2 \alpha_2)F(t_0, u_0) + \beta_2 F(t_1, u_1)]. \quad (9.222)$$

and the $i = 3$ step gives

$$w_3 = h[\alpha_2 \alpha_3 F(t_0, u_0) + \alpha_3 F(t_1, u_1) + F(t_2, u_2)], \quad (9.223)$$

$$u_3 = u_0 + [\beta_1 + \alpha_2(\beta_2 + \alpha_3\beta_3)]hF(t_0, u_0) \\ + (\beta_2 + \alpha_3\beta_3)hF(t_1, u_1) + \beta_3hF(t_2, u_2). \quad (9.224)$$

The corresponding times can be calculated by putting $F = 1$. This yields

$$t_1 = t_0 + \beta_1 h, \quad (9.225)$$

$$t_2 = t_0 + h[\beta_1 + \beta_2(1 + \alpha_2)], \quad (9.226)$$

$$t_3 = t_0 + h[\beta_1 + \beta_3 + (1 + \alpha_2)(\beta_2 + \alpha_3\beta_3)]. \quad (9.227)$$

The last expression can also be written in the form

$$t_3 = t_0 + h\{\beta_1 + \beta_2(1 + \alpha_2) + \beta_3[1 + (1 + \alpha_2)\alpha_3]\}. \quad (9.228)$$

Next we need to determine the conditions that the scheme is indeed of third order. This can be done by considering the differential equation

$$\frac{du}{dt} = u, \quad u(0) = u_0, \quad (9.229)$$

for $u = u(t)$, where u_0 is the initial value of u . The exact solution of (9.229) is $u_0 e^t$. Its Taylor expansion for $t = t_0 + h$ is

$$u(t_0 + h) = u_0 \left[1 + h + \frac{1}{2}h^2 + \frac{1}{6}h^3 + \dots \right]. \quad (9.230)$$

The solution based on (9.224) is

$$u_3 = u_0 + [\beta_1 + \alpha_2(\beta_2 + \alpha_3\beta_3)]hu_0, + (\beta_2 + \alpha_3\beta_3)hu_1 + \beta_3hu_2. \quad (9.231)$$

In order to compare with (9.230) we need the explicit expressions for u_1 and u_2 , which are

$$u_1 = (1 + h\beta_1)u_0, \quad (9.232)$$

$$u_2 = \left\{ 1 + h[\beta_1 + (1 + \alpha_2)\beta_2] + h^2\beta_1\beta_2 \right\} u_0. \quad (9.233)$$

and so we can write

$$u_3 = u_0 + \gamma_1 h + \gamma_2 h^2 + \gamma_3 h^3 \quad (9.234)$$

with

$$\gamma_1 = \beta_1 + \beta_3 + (1 + \alpha_2)(\beta_2 + \alpha_3\beta_3), \quad (9.235)$$

$$\gamma_2 = \beta_1\beta_2 + \beta_3[\beta_2(1 + \alpha_2) + \beta_1(1 + \alpha_3)], \quad (9.236)$$

$$\gamma_3 = \beta_1\beta_2\beta_3. \quad (9.237)$$

In order for the scheme to be third order we have to require $\gamma_1 = 1$, $\gamma_2 = 1/2$, and $\gamma_3 = 1/6$; see (9.230). Thus, we have now *three* equations for *five* unknowns. We now have to come up with two more equations to solve for the five unknowns.

If we assume that the intermediate timesteps are evaluated in equidistant time intervals, we have to require that the time increments in (9.225) and (9.226) are $1/3$ and $2/3$, respectively. This yields

$$\beta_1 = \beta_2(1 + \alpha_2) = 1/3, \quad (9.238)$$

with two particular solutions²

$$\begin{aligned} \alpha_2 &= -2/3, & \alpha_3 &= -1, & \beta_1 &= 1/3, & \beta_2 &= 1, & \beta_3 &= 1/2, \\ \alpha_2 &= -1/3, & \alpha_3 &= -1, & \beta_1 &= 1/3, & \beta_2 &= 1/2, & \beta_3 &= 1. \end{aligned} \quad (9.239)$$

These are in fact the simplest $2N$ -RK3 schemes that also lead to comparatively small residual errors.

Alternatively, one can move the times closer to the end time of the timestep and evaluate the right-hand side at times $t_1 - t_0 = \frac{1}{2}h$ and $t_2 - t_0 = h$. This gives the particular solution

$$\alpha_2 = -1/4, \quad \alpha_3 = -4/3, \quad \beta_1 = 1/2, \quad \beta_2 = 2/3, \quad \beta_3 = 1/2. \quad (9.240)$$

Again, there could be other solutions.

Another possibility is to require that the inhomogeneous equation

$$du/dt = t^n, \quad u(0) = 0, \quad (9.241)$$

is solved exactly up to some n . The exact solutions for $t = h$ are $u = \frac{1}{2}h^2$ for $n = 1$ and $u = \frac{1}{3}h^3$ for $n = 2$.

The case $n = 0$ was already considered in (9.225)–(9.227). For $n = 1$ we have $F(t_1, u_1) = \frac{1}{3}h$ and $F(t_2, u_2) = \frac{2}{3}h$, so (9.224) gives

$$u_3 = u_0 + (\beta_2 + \beta_3\alpha_3)\frac{1}{3}h^2 + \beta_3\frac{2}{3}h^2, \quad (9.242)$$

or

$$u_3 = u_0 + \frac{1}{3}[\beta_2 + \beta_3(\alpha_3 + 2)]h^2. \quad (9.243)$$

Comparing with the exact solution this yields the additional equation

$$\beta_2 + \beta_3(\alpha_3 + 2) = \frac{3}{2}. \quad (9.244)$$

For $n = 2$ we have $F(t_1, u_1) = \frac{1}{9}h^2$ and $F(t_2, u_2) = \frac{4}{9}h^2$, so (9.224) gives

$$u_3 = u_0 + (\beta_2 + \beta_3\alpha_3)\frac{1}{9}h^3 + \beta_3\frac{4}{9}h^3, \quad (9.245)$$

or

$$u_3 = u_0 + \frac{1}{9}[\beta_2 + \beta_3(\alpha_3 + 4)]h^3. \quad (9.246)$$

Again, comparing with the exact solution one obtains

$$\beta_2 + \beta_3(\alpha_3 + 4) = 3. \quad (9.247)$$

This gives the solution

$$\alpha_2 = -17/32, \quad \alpha_3 = -32/27, \quad \beta_1 = 1/4, \quad \beta_2 = 8/9, \quad \beta_3 = 3/4, \quad (9.248)$$

which implies that the right hand sides are evaluated at the times $t_1 - t_0 = \frac{1}{4}h$ and $t_2 - t_0 = \frac{2}{3}h$. In tables 9.1 and 9.2 this scheme is referred to as “inhomogeneous.”

Yet another idea (W. Dobler, private communication) is to obtain the additional two equations by requiring that the quadratic differential equation $du/dt = u^2$ with $u_0 = 1$ is solved exactly. The solution is $u = (1-t)^{-1}$, of which we only need the expansion up to h^2 , so we have $u_3 \approx 1 + h + h^2$. Again, we use (9.224), but now with $F = u^2$,

$$u_3 = 1 + [\beta_1 + \alpha_2(\beta_2 + \alpha_3\beta_3)]h + (\beta_2 + \alpha_3\beta_3)hu_1^2 + \beta_3hu_2^2. \quad (9.249)$$

We need u_1^2 and u_2^2 only up to the term linear in h . Using (9.232) and (9.233) we have

$$u_1^2 = 1 + 2\beta_1h + \mathcal{O}(h^2), \quad u_2^2 = 1 + 2[\beta_1 + (1 + \alpha_2)\beta_2]h + \mathcal{O}(h^2). \quad (9.250)$$

Inserting this in (9.249) yields

$$u_3 = 1 + \delta_1h + \delta_2h^2 + \mathcal{O}(h^3), \quad (9.251)$$

with

$$\delta_1 = \beta_1 + \alpha_2(\beta_2 + \alpha_3\beta_3) + (\beta_2 + \alpha_3\beta_3) + \beta_3 \quad (9.252)$$

and

$$\delta_2 = 2\beta_1(\beta_2 + \alpha_3\beta_3) + 2\beta_3[\beta_1 + (1 + \alpha_2)\beta_2]. \quad (9.253)$$

Thus, the two additional equations are

$$\beta_1 + (1 + \alpha_2)(\beta_2 + \alpha_3\beta_3) + \beta_3 = 1, \quad (9.254)$$

$$\beta_1[\beta_2 + (1 + \alpha_3)\beta_3] + (1 + \alpha_2)\beta_2 = \frac{1}{2}. \quad (9.255)$$

The numerical solution is

$$\begin{aligned} \alpha_2 &= -0.36726297, & \alpha_3 &= -1.0278248, \\ \beta_1 &= 0.30842796, & \beta_2 &= 0.54037434, & \beta_3 &= 1. \end{aligned} \quad (9.256)$$

which implies that the right-hand sides are evaluated at the times $t_1 - t_0 = 0.308h$ and $t_2 - t_0 = 0.650h$. In tables 9.1 and 9.2 this scheme is referred to as “quadratic.”

Appendix C. Derivation of the jacobian for transformation on a sphere

Here we give the explicit derivation of (9.79) and (9.80). We first use the transformation in the form

$$x = \tilde{x}^2/\tilde{r}, \quad y = \tilde{x}\tilde{y}/\tilde{r}, \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.257)$$

$$x = \tilde{x}\tilde{y}/\tilde{r}, \quad y = \tilde{y}^2/\tilde{r}, \quad \text{if } \tilde{x} \leq \tilde{y}. \quad (9.258)$$

To obtain the jacobian we differentiate with respect to x , so

$$1 = 2 \frac{\partial \ln \tilde{x}}{\partial \ln x} - \frac{\partial \ln \tilde{r}}{\partial \ln x}, \quad 0 = \frac{\partial \ln \tilde{x}}{\partial \ln x} + \frac{\partial \ln \tilde{y}}{\partial \ln x} - \frac{\partial \ln \tilde{r}}{\partial \ln x}, \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.259)$$

$$1 = \frac{\partial \ln \tilde{x}}{\partial \ln x} + \frac{\partial \ln \tilde{y}}{\partial \ln x} - \frac{\partial \ln \tilde{r}}{\partial \ln x}, \quad 0 = 2 \frac{\partial \ln \tilde{y}}{\partial \ln x} - \frac{\partial \ln \tilde{r}}{\partial \ln x}, \quad \text{if } \tilde{x} \leq \tilde{y}. \quad (9.260)$$

We now differentiate with respect to y ,

$$0 = 2 \frac{\partial \ln \tilde{x}}{\partial \ln y} - \frac{\partial \ln \tilde{r}}{\partial \ln y}, \quad 1 = \frac{\partial \ln \tilde{x}}{\partial \ln y} + \frac{\partial \ln \tilde{y}}{\partial \ln y} - \frac{\partial \ln \tilde{r}}{\partial \ln y}, \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.261)$$

$$0 = \frac{\partial \ln \tilde{x}}{\partial \ln y} + \frac{\partial \ln \tilde{y}}{\partial \ln y} - \frac{\partial \ln \tilde{r}}{\partial \ln y}, \quad 1 = 2 \frac{\partial \ln \tilde{y}}{\partial \ln y} - \frac{\partial \ln \tilde{r}}{\partial \ln y}, \quad \text{if } \tilde{x} \leq \tilde{y}. \quad (9.262)$$

The derivatives of \tilde{r} can be written as

$$\frac{\partial \ln \tilde{r}}{\partial \ln x} = \left(\frac{\tilde{x}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{x}}{\partial \ln x} + \left(\frac{\tilde{y}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{y}}{\partial \ln x}, \quad (9.263)$$

$$\frac{\partial \ln \tilde{r}}{\partial \ln y} = \left(\frac{\tilde{x}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{x}}{\partial \ln y} + \left(\frac{\tilde{y}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{y}}{\partial \ln y}. \quad (9.264)$$

In all cases we have

$$\frac{x}{y} = \frac{\tilde{x}}{\tilde{y}}, \quad (9.265)$$

so

$$+1 = \frac{\partial \ln \tilde{x}}{\partial \ln x} - \frac{\partial \ln \tilde{y}}{\partial \ln x}, \quad (9.266)$$

$$-1 = \frac{\partial \ln \tilde{x}}{\partial \ln y} - \frac{\partial \ln \tilde{y}}{\partial \ln y}, \quad (9.267)$$

and so

$$1 = 2 \frac{\partial \ln \tilde{x}}{\partial \ln x} - \left(\frac{\tilde{x}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{x}}{\partial \ln x} - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2 \left(\frac{\partial \ln \tilde{x}}{\partial \ln x} - 1 \right), \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.268)$$

$$0 = \frac{\partial \ln \tilde{x}}{\partial \ln x} + \frac{\partial \ln \tilde{y}}{\partial \ln x} - \left(\frac{\tilde{x}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{x}}{\partial \ln x} - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2 \left(\frac{\partial \ln \tilde{x}}{\partial \ln x} - 1 \right), \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.269)$$

so

$$1 = \frac{\partial \ln \tilde{x}}{\partial \ln x} + \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.270)$$

and so

$$\frac{\partial \ln \tilde{x}}{\partial \ln x} = 1 - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \frac{\partial \ln \tilde{y}}{\partial \ln x} = - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.271)$$

and correspondingly

$$1 = \frac{\partial \ln \tilde{x}}{\partial \ln x} + \frac{\partial \ln \tilde{y}}{\partial \ln x} - \left(\frac{\tilde{x}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{x}}{\partial \ln x} - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2 \left(\frac{\partial \ln \tilde{x}}{\partial \ln x} - 1 \right), \quad \text{if } \tilde{x} \leq \tilde{y}, \quad (9.272)$$

$$0 = 2 \frac{\partial \ln \tilde{x}}{\partial \ln x} - \left(\frac{\tilde{x}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{x}}{\partial \ln x} - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2 \left(\frac{\partial \ln \tilde{x}}{\partial \ln x} - 1 \right), \quad \text{if } \tilde{x} \leq \tilde{y}, \quad (9.273)$$

so

$$1 = \frac{\partial \ln \tilde{y}}{\partial \ln x} + \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \text{if } \tilde{x} \leq \tilde{y}, \quad (9.274)$$

and so

$$\frac{\partial \ln \tilde{x}}{\partial \ln x} = 2 - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \frac{\partial \ln \tilde{y}}{\partial \ln x} = 1 - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \text{if } \tilde{x} \leq \tilde{y}. \quad (9.275)$$

Hence note that there is a discontinuity of the jacobian along the diagonals. Now for the y -derivatives we have

$$0 = 2 \frac{\partial \ln \tilde{x}}{\partial \ln y} - \left(\frac{\tilde{x}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{x}}{\partial \ln y} - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2 \left(\frac{\partial \ln \tilde{x}}{\partial \ln y} + 1 \right), \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.276)$$

$$1 = \frac{\partial \ln \tilde{x}}{\partial \ln y} + \frac{\partial \ln \tilde{y}}{\partial \ln y} - \left(\frac{\tilde{x}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{x}}{\partial \ln y} - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2 \left(\frac{\partial \ln \tilde{x}}{\partial \ln y} + 1 \right), \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.277)$$

so

$$0 = \frac{\partial \ln \tilde{x}}{\partial \ln y} - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.278)$$

and so

$$\frac{\partial \ln \tilde{x}}{\partial \ln y} = + \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \frac{\partial \ln \tilde{y}}{\partial \ln y} = 1 + \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.279)$$

and correspondingly

$$0 = \frac{\partial \ln \tilde{x}}{\partial \ln y} + \frac{\partial \ln \tilde{y}}{\partial \ln y} - \left(\frac{\tilde{x}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{x}}{\partial \ln y} - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2 \left(\frac{\partial \ln \tilde{x}}{\partial \ln y} + 1 \right), \quad \text{if } \tilde{x} \leq \tilde{y}, \quad (9.280)$$

$$-1 = 2 \frac{\partial \ln \tilde{x}}{\partial \ln y} - \left(\frac{\tilde{x}}{\tilde{r}} \right)^2 \frac{\partial \ln \tilde{x}}{\partial \ln y} - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2 \left(\frac{\partial \ln \tilde{x}}{\partial \ln y} + 1 \right), \quad \text{if } \tilde{x} \leq \tilde{y}, \quad (9.281)$$

so

$$0 = \frac{\partial \ln \tilde{x}}{\partial \ln y} + 1 - \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \text{if } \tilde{x} \leq \tilde{y}, \quad (9.282)$$

and so

$$\frac{\partial \ln \tilde{x}}{\partial \ln y} = -1 + \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \frac{\partial \ln \tilde{y}}{\partial \ln y} = + \left(\frac{\tilde{y}}{\tilde{r}} \right)^2, \quad \text{if } \tilde{x} \leq \tilde{y}. \quad (9.283)$$

So, in summary

$$\begin{pmatrix} \frac{\partial \ln \tilde{x}}{\partial \ln x} & \frac{\partial \ln \tilde{x}}{\partial \ln y} \\ \frac{\partial \ln \tilde{y}}{\partial \ln x} & \frac{\partial \ln \tilde{y}}{\partial \ln y} \end{pmatrix} = \begin{pmatrix} +1 - \left(\frac{\tilde{y}}{\tilde{r}}\right)^2 & +1 - \left(\frac{\tilde{x}}{\tilde{r}}\right)^2 \\ -1 + \left(\frac{\tilde{x}}{\tilde{r}}\right)^2 & +1 + \left(\frac{\tilde{y}}{\tilde{r}}\right)^2 \end{pmatrix} \quad \text{if } \tilde{x} \geq \tilde{y}, \quad (9.284)$$

$$\begin{pmatrix} \frac{\partial \ln \tilde{x}}{\partial \ln x} & \frac{\partial \ln \tilde{x}}{\partial \ln y} \\ \frac{\partial \ln \tilde{y}}{\partial \ln x} & \frac{\partial \ln \tilde{y}}{\partial \ln y} \end{pmatrix} = \begin{pmatrix} +1 + \left(\frac{\tilde{x}}{\tilde{r}}\right)^2 & -1 + \left(\frac{\tilde{y}}{\tilde{r}}\right)^2 \\ +1 - \left(\frac{\tilde{y}}{\tilde{r}}\right)^2 & +1 - \left(\frac{\tilde{x}}{\tilde{r}}\right)^2 \end{pmatrix} \quad \text{if } \tilde{x} \leq \tilde{y}. \quad (9.285)$$

Appendix D. Derivation of the incremental jacobian for second derivatives

Here we present the explicit derivation of (9.110). To calculate the second derivative of a function f that is represented on a coordinate mesh $\tilde{\mathbf{x}}$, is given by

$$\frac{\partial^2 f}{\partial x_i \partial x_j} = \frac{\partial}{\partial x_i} \left(\frac{\partial f}{\partial x_j} \right) = \frac{\partial}{\partial x_i} \left(\frac{\partial f}{\partial \tilde{x}_k} \frac{\partial \tilde{x}_k}{\partial x_j} \right) \quad (9.286)$$

so

$$\frac{\partial^2 f}{\partial x_i \partial x_j} = \frac{\partial^2 f}{\partial \tilde{x}_l \partial \tilde{x}_k} \frac{\partial \tilde{x}_l}{\partial x_i} \frac{\partial \tilde{x}_k}{\partial x_j} + \frac{\partial f}{\partial \tilde{x}_k} \frac{\partial^2 \tilde{x}_k}{\partial x_i \partial x_j}, \quad (9.287)$$

or

$$\frac{\partial^2 f}{\partial x_i \partial x_j} = \frac{\partial^2 f}{\partial \tilde{x}_p \partial \tilde{x}_q} \mathbf{J}_{pi} \mathbf{J}_{qj} + \frac{\partial f}{\partial \tilde{x}_k} \mathbf{K}_{kij}, \quad (9.288)$$

which is just (9.108), using (9.109) for the definition of \mathbf{K}_{kij} of the 2nd-order jacobian.

To obtain the 2nd-order jacobian by successive tensor multiplication we differentiate twice the evolution equation for \mathbf{x} ,

$$x_k^{(n+1)} = x_k^{(n)} + u_k^{(n)} \delta t, \quad (9.289)$$

so

$$\frac{\partial^2 x_k^{(n+1)}}{\partial x_i^{(n+1)} \partial x_j^{(n+1)}} = \frac{\partial^2 x_k^{(n)}}{\partial x_i^{(n+1)} \partial x_j^{(n+1)}} + \frac{\partial^2 u_k^{(n)}}{\partial x_i^{(n+1)} \partial x_j^{(n+1)}} \delta t. \quad (9.290)$$

The expression on the left-hand side is just the derivative of a Kronecker delta, see (9.102), so it is zero. Thus we have

$$0 = \frac{\partial^2 x_k^{(n)}}{\partial x_i^{(n+1)} \partial x_j^{(n+1)}} + \frac{\partial}{\partial x_i^{(n+1)}} \left(\frac{\partial u_k^{(n)}}{\partial x_q^{(n)}} \frac{\partial x_q^{(n)}}{\partial x_j^{(n+1)}} \right) \delta t, \quad (9.291)$$

or

$$0 = \frac{\partial^2 x_q^{(n)}}{\partial x_i^{(n+1)} \partial x_j^{(n+1)}} \left(\delta_{kq} + \frac{\partial u_k^{(n)}}{\partial x_q^{(n)}} \delta t \right) + \frac{\partial^2 u_k^{(n)}}{\partial x_p^{(n)} \partial x_q^{(n)}} \mathbf{J}_{pi}^{(n)} \mathbf{J}_{qj}^{(n)} \delta t, \quad (9.292)$$

which can be written as

$$0 = \mathbf{M}_{kq} \mathbf{K}_{qij}^{(n)} + u_{k,pq} \mathbf{J}_{pi}^{(n)} \mathbf{J}_{qj}^{(n)} \delta t, \quad (9.293)$$

so

$$\mathbf{K}_{kij}^{(n)} = -\delta t \left(\mathbf{M}^{-1} \right)_{kl} u_{lpq} \mathbf{J}_{pi}^{(n)} \mathbf{J}_{qj}^{(n)}, \quad (9.294)$$

which is just (9.112).

We now need to derive the equation that relates the incremental 2nd-order jacobians to the 2nd-order jacobian of the previous timestep. To this end we begin with the 2nd-order jacobian at time $2\delta t$, so

$$\mathbf{K}_{kij}^{(0 \rightarrow 2)} \equiv \frac{\partial^2 x_k^{(0)}}{\partial x_i^{(2)} \partial x_j^{(2)}} = \frac{\partial}{\partial x_i^{(2)}} \left(\frac{\partial x_k^{(0)}}{\partial x_q^{(1)}} \frac{\partial x_q^{(1)}}{\partial x_j^{(2)}} \right) \quad (9.295)$$

or

$$\mathbf{K}_{kij}^{(0 \rightarrow 2)} \equiv \frac{\partial^2 x_k^{(0)}}{\partial x_p^{(1)} \partial x_q^{(1)}} \frac{\partial x_p^{(1)}}{\partial x_i^{(2)}} \frac{\partial x_q^{(1)}}{\partial x_j^{(2)}} + \frac{\partial^2 x_q^{(1)}}{\partial x_i^{(2)} \partial x_j^{(2)}} \frac{\partial x_k^{(0)}}{\partial x_q^{(1)}} \quad (9.296)$$

or

$$\mathbf{K}_{kij}^{(0 \rightarrow 2)} = \mathbf{K}_{kpq}^{(0 \rightarrow 1)} \mathbf{J}_{pi}^{(1)} \mathbf{J}_{qj}^{(1)} + \mathbf{J}_{kl}^{(0 \rightarrow 1)} \mathbf{K}_{lij}^{(1)}. \quad (9.297)$$

For the next step we have

$$\mathbf{K}_{kij}^{(0 \rightarrow 3)} \equiv \frac{\partial^2 x_k^{(0)}}{\partial x_i^{(3)} \partial x_j^{(3)}} = \frac{\partial}{\partial x_i^{(3)}} \left(\frac{\partial x_k^{(0)}}{\partial x_q^{(1)}} \frac{\partial x_q^{(1)}}{\partial x_p^{(2)}} \frac{\partial x_p^{(2)}}{\partial x_j^{(3)}} \right), \quad (9.298)$$

so

$$\mathbf{K}_{kij}^{(0 \rightarrow 3)} = \mathbf{K}_{kpq}^{(0 \rightarrow 1)} \mathbf{J}_{pr}^{(1)} \mathbf{J}_{qs}^{(1)} \mathbf{J}_{ri}^{(2)} \mathbf{J}_{sj}^{(2)} + \mathbf{J}_{kl}^{(0 \rightarrow 1)} \mathbf{K}_{lrs}^{(1)} \mathbf{J}_{ri}^{(2)} \mathbf{J}_{sj}^{(2)} + \mathbf{J}_{kr}^{(0)} \mathbf{J}_{rs}^{(1)} \mathbf{K}_{sij}^{(2)}. \quad (9.299)$$

This can be written as

$$\mathbf{K}_{kij}^{(0 \rightarrow 3)} = \mathbf{K}_{kpq}^{(0 \rightarrow 2)} \mathbf{J}_{pi}^{(2)} \mathbf{J}_{qj}^{(2)} + \mathbf{J}_{kl}^{(0 \rightarrow 2)} \mathbf{K}_{lij}^{(2)} \quad (9.300)$$

which, for the general step from 0 to n , becomes (9.110).

Appendix E. Solution for α and η_t quenched α^2 -dynamo

Here we present the explicit derivation of (9.181). According to mean-field theory for non-mirror symmetric isotropic homogeneous turbulence with no mean flow the mean magnetic field is governed by the equation

$$\frac{\partial}{\partial t} \bar{\mathbf{B}} = \nabla \times (\alpha \bar{\mathbf{B}} - \eta_T \eta_0 \bar{\mathbf{J}}), \quad (9.301)$$

where bars denote the mean fields and $\eta_T = \eta + \eta_t$ is the total (microscopic plus turbulent) magnetic diffusion. Both α -effect and turbulent diffusion are assumed to be quenched in the same way, so

$$\alpha = \frac{\alpha_0}{1 + \alpha_B \bar{B}^2 / B_{\text{eq}}^2}, \quad \eta_t = \frac{\eta_{t0}}{1 + \eta_B \bar{B}^2 / B_{\text{eq}}^2}. \quad (9.302)$$

In the following we assume $\alpha_B = \eta_B$ and denote

$$\alpha_B / B_{\text{eq}}^2 = 1 / B_0^2 = \eta_B / B_{\text{eq}}^2. \quad (9.303)$$

We emphasize that only the *turbulent* magnetic diffusivity is quenched, not of course the total one. It is only because of the presence of microscopic diffusion that saturation is possible.

In the simulations \bar{B}^2 is to a good approximation spatially uniform. Defining the magnetic energy as $M(t) = \frac{1}{2}\langle \bar{B}^2 \rangle$ we have

$$\bar{B}^2 = 2M, \quad (9.304)$$

which is only a function of time.

Consider the particular example where the large scale field varies only in the z direction (9.301) becomes

$$\dot{\bar{B}}_x = -\alpha \bar{B}'_y + \eta_T \bar{B}''_x, \quad (9.305)$$

$$\dot{\bar{B}}_y = +\alpha \bar{B}'_x + \eta_T \bar{B}''_y, \quad (9.306)$$

where dots and primes denote differentiation with respect to t and z , respectively. Since $\alpha < 0$, the solution can be written in the form

$$\bar{B}_x(y, t) = b_x(t) \cos(z + \phi), \quad (9.307)$$

$$\bar{B}_y(y, t) = b_y(t) \sin(z + \phi), \quad (9.308)$$

where $b_x(t)$ and $b_z(t)$ are positive functions of time that satisfy

$$\dot{b}_x = |\alpha| b_y - \eta_T b_x, \quad (9.309)$$

$$\dot{b}_y = |\alpha| b_x - \eta_T b_y. \quad (9.310)$$

We now choose the special initial condition, $b_x = b_y \equiv b$, so we have only one equation for the variable b . Note also that in the quenching factor $\bar{B}^2 = b_x^2 + b_y^2 = 2b^2$. Thus, we have

$$\frac{db}{dt} = \frac{\alpha_0 k_1 - \eta_{t0} k_1^2}{1 + 2b^2 / B_0^2} b - \eta k_1^2 b. \quad (9.311)$$

Multiplying with b yields

$$\frac{1}{2} \frac{db^2}{dt} = \frac{\alpha_0 k_1 - \eta_{t0} k_1^2}{1 + 2b^2 / B_0^2} b^2 - \eta k_1^2 b^2. \quad (9.312)$$

Using the definition $M = \frac{1}{2}b^2$ we have

$$\frac{1}{2} \frac{1}{M} \frac{dM}{dt} = \frac{\alpha_0 k_1 - \eta_{t0} k_1^2}{1 + 2M/M_0} - \eta k_1^2, \quad (9.313)$$

where $M_0 = \frac{1}{2}B_0^2$. Thus, we have

$$\left(1 + 2\frac{M}{M_0}\right) \frac{1}{2} \frac{1}{M} \frac{dM}{dt} = \alpha_0 k_1 - \eta_{T0} k_1^2 - 2\eta k_1^2 \frac{M}{M_0}, \quad (9.314)$$

where we have defined $\eta_{T0} = \eta_{t0} + \eta$. We define the abbreviations $\lambda = \alpha_0 k_1 - \eta_{T0} k_1^2$ for the kinematic growth rate of the dynamo and $\lambda_{t0} = \eta_{t0} k_1^2$ for the turbulent decay rate if there were no dynamo action, and arrive thus at the integral

$$\int_{M_{\text{ini}}}^M \frac{1 + 2M'/M_0}{1 - (2\eta k_1^2/\lambda) M'/M_0} \frac{dM'}{M'} = 2\lambda t. \quad (9.315)$$

We now also define the abbreviation

$$\mathcal{M} = 2M'/M_0 \quad (9.316)$$

and have

$$\int_{M_{\text{ini}}}^M \frac{1 + \mathcal{M}}{1 - (\eta k_1^2/\lambda) \mathcal{M}} \frac{d\mathcal{M}}{\mathcal{M}} = 2\lambda t. \quad (9.317)$$

which can be split into two integrals,

$$\int_{M_{\text{ini}}}^M \frac{d\mathcal{M}}{[1 - (\eta k_1^2/\lambda) \mathcal{M}] \mathcal{M}} + \int_{M_{\text{ini}}}^M \frac{d\mathcal{M}}{1 - (\eta k_1^2/\lambda) \mathcal{M}} = 2\lambda t. \quad (9.318)$$

To solve these integrals we note that

$$\int \frac{dx}{1 - x/x_0} = -x_0 \ln(x_0 - x), \quad (9.319)$$

$$\int \frac{dx}{(1 - x/x_0)x} = \int \frac{dx}{x_0 - x} + \int \frac{dx}{x} = \ln x - \ln(1 - x/x_0) = \ln \left(\frac{x}{x_0 - x} \right), \quad (9.320)$$

so (9.318) becomes

$$\ln \left(\frac{\mathcal{M}}{\lambda/\eta k_1^2 - \mathcal{M}} \right) + \frac{\lambda}{\eta k_1^2} \ln \left(\frac{1}{\lambda/\eta k_1^2 - \mathcal{M}} \right) = 2\lambda(t - t_0), \quad (9.321)$$

where t_0 is an integration constant. Exponentiation yields

$$\frac{\mathcal{M}}{(\lambda/\eta k_1^2 - \mathcal{M})^{1+\lambda/\eta k_1^2}} = e^{2\lambda(t-t_0)}, \quad (9.322)$$

which is, in terms of the original variables, $\bar{\mathbf{B}}^2/B_0^2 = 2b^2/B_0^2 = 2M/M_0 = \mathcal{M}$, we have

$$\frac{\bar{\mathbf{B}}^2}{B_0^2} \left/ \left(\frac{\lambda}{\eta k_1^2} - \frac{\bar{\mathbf{B}}^2}{B_0^2} \right)^{1+\lambda/\eta k_1^2} \right. = e^{2\lambda(t-t_0)}. \quad (9.323)$$

The final field strength,

$$B_{\text{fin}} = \lim_{t \rightarrow \infty} |\bar{\mathbf{B}}(t)|, \quad (9.324)$$

is given by requiring the denominator to vanish, which yields

$$\frac{B_{\text{fin}}^2}{B_0^2} = \frac{\lambda}{\eta k_1^2}. \quad (9.325)$$

Rewriting (9.323) in terms of B_{fin} we have

$$\frac{\bar{\mathbf{B}}^2}{B_{\text{fin}}^2} \left/ \left(1 - \frac{\bar{\mathbf{B}}^2}{B_{\text{fin}}^2} \right)^{1+\lambda/\eta k_1^2} \right. = \left(\frac{\lambda}{\eta k_1^2} \right)^{1+\lambda/\eta k_1^2} e^{2\lambda(t-t_0)}. \quad (9.326)$$

We can express t_0 in terms of the initial field strength, $B_{\text{ini}} = |\bar{\mathbf{B}}(0)|$, and if the initial field strength is much weaker than the final field strength, i.e. $B_{\text{ini}} \ll B_{\text{fin}}$, then we can rewrite (9.326), in the form

$$\bar{\mathbf{B}}^2 / (1 - \bar{\mathbf{B}}^2/B_{\text{fin}}^2)^{1+\lambda/\eta k_1^2} = B_{\text{ini}}^2 e^{2\lambda t}. \quad (9.327)$$

Thus, for early times we have the familiar relation

$$|\bar{\mathbf{B}}| \approx B_{\text{ini}} e^{\lambda t} \quad (|\bar{\mathbf{B}}| \ll B_{\text{fin}}), \quad (9.328)$$

whereas for late times near the final field strength we have

$$(1 - \bar{\mathbf{B}}^2/B_{\text{fin}}^2)^{-(1+\lambda/\eta k_1^2)} \approx (B_{\text{ini}}/B_{\text{fin}})^2 e^{2\lambda t} \quad (|\bar{\mathbf{B}}| \approx B_{\text{fin}}), \quad (9.329)$$

or

$$\bar{\mathbf{B}}^2 \approx B_{\text{fin}}^2 \left[1 - e^{-2\lambda(t-t_{\text{sat}})} \right]^{\eta k_1^2/(\lambda+\eta k_1^2)} \quad (|\bar{\mathbf{B}}| \approx B_{\text{fin}}), \quad (9.330)$$

where $t_{\text{sat}} = \lambda^{-1} \ln(B_{\text{fin}}/B_{\text{ini}})$ is the time it takes to reach saturation. If the Reynolds number is large we have $\lambda \gg \eta k_1^2$, so

$$\bar{\mathbf{B}}^2 \approx B_{\text{fin}}^2 \left[1 - e^{-2\eta k_1^2(t-t_{\text{sat}})} \right] \quad (|\bar{\mathbf{B}}| \approx B_{\text{fin}}), \quad (9.331)$$

which is identical to the result obtained from helicity conservation.

Note that the solution (9.327) is governed by four parameters: B_{ini} , B_{fin} , λ , and ηk_1^2 . The latter is known from the input data to the simulation, B_{ini} and λ can be determined from the

linear growth phase of the dynamo (characterized by properties of the small scale dynamo!) and so B_{fin} is the only parameter that is determined by the nonlinearity of the dynamo and can easily be determined from the simulations. Once B_{fin} is measured from numerical experiments we know immediately the quenching parameters

$$\alpha_B = \eta_B = \frac{\lambda}{\eta k_1^2} \left(\frac{B_{\text{eq}}}{B_{\text{fin}}} \right)^2, \quad (9.332)$$

and since $B_{\text{fin}}^2/B_{\text{eq}}^2 \approx k_f/k_1$ we have

$$\alpha_B = \eta_B \approx \frac{\lambda}{\eta k_1 k_f}, \quad (9.333)$$

which shows that α_B and η_B are proportional to the magnetic Reynolds number.

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Notes

- 1 The 4th-order compact scheme is really identical to calculating derivatives from a cubic spline, as was done in Nordlund and Stein (1990). In the book by Collatz (1966) the compact methods are also referred to as *Hermitian methods* or as *Mehrstellen-Verfahren*, because the derivative in one point is calculated using the derivatives in neighboring points.
- 2 I thank Petri Käpylä from Oulu for pointing out the second of these solutions.

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