A NEW MULTI-DIMENSIONAL GENERAL RELATIVISTIC NEUTRINO HYDRO_DYNAMIC CODE FOR CORE-COLLAPSE SUPERNOVAE. I. METHOD AND CODE TESTS IN SPHERICAL SYMMETRY

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ABSTRACT

We present a new general relativistic code for hydrodynamical supernova simulations with neutrino transport in spherical and azimuthal symmetry (one dimension and two dimensions, respectively). The code is a combination of the CoCoNuT hydro module, which is a Riemann-solver-based, high-resolution shock-capturing method, and the three-flavor, fully energy-dependent VERTEX scheme for the transport of massless neutrinos. VERTEX integrates the coupled neutrino energy and momentum equations with a variable Eddington factor closure computed from a model Boltzmann equation and uses the “ray-by-ray plus” approximation in two dimensions, assuming the neutrino distribution to be axially symmetric around the radial direction at every point in space, and thus the neutrino flux to be radial. Our spacetime treatment employs the Arnowitt–Deser–Misner 3+1 formalism with the conformal flatness condition for the spatial three metric. This approach is exact for the one-dimensional case and has previously been shown to yield very accurate results for spherical and rotational stellar core collapse. We introduce new formulations of the energy equation to improve total energy conservation in relativistic and Newtonian hydro simulations with grid-based Eulerian finite-volume codes. Moreover, a modified version of the VERTEX scheme is developed that simultaneously conserves energy and lepton number in the neutrino transport with better accuracy and higher numerical stability in the high-energy tail of the spectrum. To verify our code, we conduct a series of tests in spherical symmetry, including a detailed comparison with published results of the collapse, shock formation, shock breakout, and accretion phases. Long-time simulations of proto-neutron star cooling until several seconds after core bounce both demonstrate the robustness of the new CoCoNuT-VERTEX code and show the approximate treatment of relativistic effects by means of an effective relativistic gravitational potential as in PROMETHEUS-VERTEX to be remarkably accurate in spherical symmetry.

Key words: hydrodynamics – methods: numerical – neutrinos – radiative transfer – relativistic processes – supernovae: general

1. INTRODUCTION

The collapse and explosion of massive stars has been an active topic in astrophysical research for decades, and has provided a major impetus for advances in computational relativistic astrophysics, in particular ever since the first numerical simulations of core collapse were conducted in the 1960s (Colgate & White 1966). The general outline of this rich astrophysical phenomenon is fairly well established by now: after the collapse and bounce of the central iron core to a compact neutron star (\(M/R \approx 0.1–0.2\) in relativistic units) a shock wave is formed, which stalls a few milliseconds after bounce. Afterward, accretion onto the newly formed neutron star continues typically for a few hundreds of milliseconds until the shock is re-energized either by neutrino heating (the “delayed neutrino-driven mechanism,” Wilson 1985), probably aided by multi-dimensional hydrodynamical instabilities like convection (Epstein 1979; Herant et al. 1992, 1994; Burrows et al. 1995, 2006a; Janka & Müller 1995, 1996; Buras et al. 2006b, 2006a; Bruenn et al. 2006) and the standing accretion-shock instability (SASI; Blondin et al. 2003), or by some alternative mechanism, such as MHD effects (e.g., Burrows et al. 2007 and references therein), or energy deposition by acoustic waves (Burrows et al. 2006b). The ongoing quest for the explosion mechanism, with various research groups pursuing a number of different avenues, is a testimony to the complexity of the problem, which involves not only multi-dimensional (magneto-)hydrodynamics, but also kinetic theory (i.e., neutrino transport), nuclear and particle physics, and general relativistic (GR) gravity. Even for the few hundreds of milliseconds from the collapse to the onset of the explosion, there is no unified approach that fully accounts for all these intricacies. Instead, a number of quite disparate approaches have been used in the past to investigate different aspects of core-collapse supernovae, all of which are arguably deficient in the sense that they sacrifice some essential physical ingredients in favor of others.

Thus, there have been repeated attempts to solve the problem of the explosion mechanism by means of an accurate treatment of neutrino transport in spherical symmetry, combined with reasonably accurate microphysics (i.e., the nuclear equation of state (EOS), nuclear burning, and neutrino–matter interactions). These endeavors culminated in the development of solvers for the neutrino Boltzmann equation (Mezzacappa & Messer 1999; Burrows et al. 2000; Liebendörfer et al. 2001b, 2004; Rampp & Janka 2002; Yamada et al. 1999); but the exclusion of multi-dimensional hydrodynamical effects is obviously a serious shortcoming. On the other hand, the restriction to spherical symmetry allows for a full GR treatment with relatively little additional effort: relativistic supernova simulations with gray or multi-group flux-limited diffusion (MGFLD) have been conducted since the 1980s in one dimension (Baron et al. 1989; Bruenn et al. 2001), and relativistic Boltzmann transport became feasible after another few years (Yamada et al. 1999; Liebendörfer et al. 2001b, 2004). As expected, GR effects were shown to be anything but negligible: comparisons with the Newtonian case (see, e.g., Bruenn et al.
Another complementary approach has focused primarily on the interplay between neutrino heating/cooling and multi-dimensional (magneto-)hydrodynamic effects, which has turned out to be no less important for the understanding of the explosion mechanism than the proper treatment of neutrino transport and general relativity: convective instabilities were found to operate both inside the proto-neutron star (Epstein 1979) and the “hot-bubble region” behind the stalled shock (Bethe 1990). Hot-bubble convection in particular was shown to be helpful for increasing the neutrino heating efficiency already in the 1990s (Herant et al. 1992, 1994; Burrows et al. 1995; Janka & Müller 1995, 1996), and another recently identified hydrodynamic instability of the shock, known as the SASI (Blondin et al. 2003), was found to have a similar effect. Nowadays, the most advanced computational tools used in this approach combine multi-dimensional Newtonian hydrodynamics with an up-to-date treatment of the microphysics with multi-group neutrino transport methods of similar sophistication as in the one-dimensional case (Livne et al. 2004; Buras et al. 2006b; Bruenn et al. 2006; Liebendörfer et al. 2009). But some weak spots remain to be criticized. First, multi-dimensional neutrino transport has not been treated in its full complexity yet; the approaches now in use either assume the neutrino flux vector to point in the radial direction (the ray-by-ray approach; Buras et al. 2006b; Bruenn et al. 2006), neglect energy redistribution in phase space (Livne et al. 2004; Liebendörfer et al. 2009), or involve even more radical approximations. Second, all serious attempts to model the development of multi-dimensional instabilities during the post-bounce phase, while treating the neutrino transport with reasonable accuracy, have been carried out in the framework of Newtonian hydrodynamics, which is rather unsatisfactory considering the importance of GR effects established in one-dimensional studies. So far, attempts to include relativistic corrections have been limited to the use of modified gravitational potentials that can be heuristically derived from the Tolman–Oppenheimer–Volkov (TOV) equation of stellar structure (Rampp & Janka 2002; Marek et al. 2006). Although this method can be justified on the basis of comparisons with results from relativistic one-dimensional neutrino transport codes and multi-dimensional relativistic hydrodynamics codes without neutrino transport (Liebendörfer et al. 2005; Marek et al. 2006; Müller et al. 2008), it comes along with several drawbacks: as the equations of Newtonian hydrodynamics are left unaltered, it may lead to incorrect predictions for some dynamical properties of the neutron star such as the oscillation eigenfrequencies (Müller et al. 2008), which may also result in a distortion of the gravitational wave spectrum. Moreover, the “effective potential approach” ignores the effects of relativistic kinematics and still allows superluminal velocities—which have actually occurred in Newtonian simulations of the explosions of O–Ne–Mg cores. Finally, it should be noted that, different from the purely Newtonian case, a conservation law for the total (i.e., kinetic, internal, and potential) energy cannot be formulated when a modified gravitational potential is used. As a consequence, it is difficult to exclude the possibility of non-physical energy loss or generation, in particular as a cumulative effect over long evolution periods.

On the other hand, the last few years have seen the development of genuinely relativistic hydrodynamics codes for multi-dimensional simulations of core collapse, often based on a rather crude treatment of the microphysics, and without any up-to-date scheme for the neutrino transport. Typically, simple analytic EOSs, such as the hybrid EOS of Janka et al. (1993) were employed until recent years; and the polytropic models for the progenitor had only a limited resemblance to iron cores from stellar evolution calculations. These limitations have naturally led the numerical relativity community to focus on the one hand on the only phases of the supernova for which such severe approximations may be deemed sufficient, i.e., the collapse and bounce of the core, and on the other hand on those issues such as black hole formation (Sekiguchi & Shibata 2005), gravitational wave emission from rotational core collapse (Dimmelmeier et al. 2002a, 2002b; Shibata & Sekiguchi 2004), and the development of non-axisymmetric instabilities (Shibata & Sekiguchi 2005) that do not lend themselves readily to a Newtonian analysis, at least not with the same accuracy. Obviously, such an approach, which neglects physics that is unquestionably of paramount importance for the stellar core-collapse problem, is beset with large uncertainties. However, the first steps toward a more sophisticated treatment of the microphysics in core-collapse supernovae were taken by Dimmelmeier et al. (2007) and Ott et al. (2007a), who studied the collapse of rotating iron cores with a nuclear EOS and a simple parameterized scheme for the deleptonization during collapse (Liebendörfer 2005). The fact that their findings are rather different from earlier studies with more radical approximations—they exclusively find gravitational wave signals of the so-called type I (Zwerger & Müller 1997)—underlines the need for a reliable treatment of the microphysics also in the field of relativistic stellar core collapse.

Clearly, the most satisfactory approach to modeling supernovae would be to sacrifice as little of the essential ingredients (neutrino transport, EOS and other microphysics, general relativity, and multi-dimensional hydrodynamics) as possible in the numerical models. In this paper, we therefore describe a relativistic generalization of the Newtonian ray-by-ray-plus method of Buras et al. (2006b) for the neutrino transport, which we have combined with the so-called conformal flatness condition (CFC) or Isenberg–Wilson–Mathews approximation (Isenberg 2008; Wilson et al. 1996) for the gravitational field equations. The method presented here has been implemented by building on the existing radiation transport code VERTEX (Rampp & Janka 2002; Buras et al. 2006b) and the relativistic hydrodynamics code CoCoNuT (Dimmelmeier et al. 2002a, 2005), and has already been applied to two-dimensional simulations of the collapse and post-bounce phase, which will be the subject of future papers. At this stage, however, we shall confine ourselves to presenting thorough tests of our new VERTEX-CoCoNuT code in spherical symmetry.

This paper is organized as follows. In Section 2, we discuss the governing equations of hydrodynamics and neutrino transport (“neutrino hydrodynamics”) for the ray-by-ray-plus approximation, and comment on their implementation in CoCoNuT and VERTEX. Tests of the new code in spherical symmetry are presented in Section 3, most notably a detailed comparison of one-dimensional simulations of the 15 $M_\odot$ progenitor model of Woosley & Weaver (1995) with VERTEX-CoCoNuT, the relativistic Boltzmann code AGILE of Liebendörfer et al. (2004), and the Newtonian VERTEX code of Rampp & Janka (2002), in which GR effects are taken into account approximately by using a modified gravitational potential. In this context, the issue of
energy and lepton number conservation also receives particular attention. In order to demonstrate the ability of VERTEX-CoCoNuT to evolve supernova models stably well beyond 1 s after core bounce, we also present results for a one-dimensional explosion model of an O–Ne–Mg core (Nomoto 1984, 1987; Kitaura et al. 2006).

Further technical details of our numerical implementation, which may also be of relevance for similar radiation hydrodynamics codes, are given in the appendices. In Appendix A we address the problem of total energy conservation for a self-gravitating fluid in Newtonian and relativistic hydrodynamics, and in Appendix B we outline an efficient finite-difference scheme for the neutrino moment equations that guarantees the simultaneous conservation of energy and lepton number.

Note that geometrized units are used throughout Section 2, i.e., both the speed of light and the gravitational constant are set to unity: \( c = \epsilon = 1 \). Greek indices run from 0 to 3, Latin indices from 1 to 3.

2. GENERAL RELATIVISTIC NEUTRINO RADIATION HYDRODYNAMICS

2.1. Hydrodynamics

In general relativity, the hydrodynamic evolution of a perfect fluid is governed by two conservation equations for the baryonic rest-mass current \( J^\mu \) and the stress–energy tensor \( T^{\mu\nu} \):

\[
\nabla_\mu J^\mu = 0, \quad \nabla_\nu T^{\mu\nu} = 0, \quad (1)
\]

where \( \nabla_\nu \) denotes the covariant derivative. For a perfect fluid, \( J^\mu \) and \( T^{\mu\nu} \) can be expressed in terms of the baryon rest-mass density \( \rho \) in the local fluid frame, the four-velocity \( u^\mu \), the pressure \( P \), the specific enthalpy \( h = 1 + \epsilon + P/\rho \) (where \( \epsilon \) is the specific internal energy), and the four-metric \( g_{\mu\nu} \):

\[
J^\mu = \rho u^\mu, \quad T^{\mu\nu} = \rho hu^\mu u^\nu + P g^{\mu\nu}. \quad (2)
\]

For the metric, we adopt the Arnowitt–Deser–Misner (ADM) 3 + 1 formalism (Lichnerovicz 1944) of general relativity to foliate the spacetime into spacelike hypersurfaces. In this approach, the four-dimensional line element \( ds^2 = g_{\mu\nu} dx^\mu dx^\nu \) is written as follows:

\[
d s^2 = -\alpha^2 dt^2 + \gamma_{ij}(dx^i + \beta^j dt)(dx^j + \beta^j dt), \quad (3)
\]

where \( \alpha \) is the lapse function, \( \beta^i \) is the shift vector, and \( \gamma_{ij} \) is the induced three-metric on each hypersurface. Using this decomposition of the four-metric, the equations of hydrodynamics can be formulated in the frame of an Eulerian observer moving orthogonally to the spacelike hypersurfaces, i.e., with a four-velocity \( n^\mu = (\alpha^{-1}, \alpha^{-1}\beta^i) \). As in Banyuls et al. (1997), we introduce the following conserved variables:

\[
D = \rho W, \quad S^i = \rho h W^2 v_i, \quad \tau = \rho h W^2 - P - D. \quad (4)
\]

where \( v^i = u^i/(\alpha u^0) + \beta^i/\alpha \) is the three-velocity in the Eulerian frame, and \( W = 1/\sqrt{1 - v_i v^i} \) is the corresponding Lorentz factor. The equations of GR hydrodynamics in flux-conservative form then read

\[
\frac{\partial}{\partial t} \sqrt{\gamma} \rho W + \frac{\partial}{\partial x^i} \sqrt{\gamma} g \rho W v^i = 0, \quad (5)
\]

\[
\frac{\partial}{\partial t} \sqrt{\gamma} \rho h W^2 v_j \frac{\partial}{\partial x^j} \left( \frac{\partial}{\partial t} \sqrt{\gamma} S^i \right) + \frac{\partial}{\partial x^i} \left( \frac{\partial}{\partial t} \sqrt{\gamma} S^i \right) = \frac{1}{2} \left( \frac{\partial}{\partial x^i} T^{\mu\nu} \frac{\partial}{\partial x^j} + \left( \frac{\partial}{\partial t} \sqrt{\gamma} S^i \right) \right) \quad (6)
\]

\[
\frac{\partial}{\partial t} \sqrt{\gamma} \tau \frac{\partial}{\partial x^i} \left( \frac{\partial}{\partial t} \sqrt{\gamma} T^{\mu\nu} \frac{\partial}{\partial x^j} + \left( \frac{\partial}{\partial t} \sqrt{\gamma} S^i \right) \right) = \alpha \sqrt{-g} \left( T^{\mu\nu} \frac{\partial}{\partial x^\mu} - T^{0\mu} \frac{\partial}{\partial x^\nu} \right) \quad (7)
\]

Here, \( \hat{v}^i \) is defined as \( \hat{v}^i = v^i - \beta^i/\alpha \), and \( g \) and \( \gamma \) are the determinants of the four-metric \( g_{\mu\nu} \) and the three-metric \( \gamma_{ij} \), respectively. Different from the purely hydrodynamic case discussed by Banyuls et al. (1997), the right-hand sides of Equations (6) and (7) contain source terms for the exchange of momentum and energy with neutrinos (which can be related to angular moments of the collision integral and are therefore denoted by a subscript “C”) in addition to the gravitational and geometric source terms. These source terms will be specified in Section 2.3.5.

Equations (5)–(7) must be supplemented by an additional equation expressing the conservation of electron number in the absence of weak interactions (which will be discussed separately):

\[
\frac{\partial}{\partial t} \sqrt{\gamma} \rho Ye + \frac{\partial}{\partial x^i} \sqrt{\gamma} g \rho Ye v^i = \left( \frac{\partial}{\partial t} \sqrt{\gamma} Ye \right) \quad (8)
\]

where \( Ye \) is the electron fraction (number of electrons minus number of positrons per baryon). The change of the electron number due to neutrino emission or absorption is accounted for by the source term on the right-hand side, in a similar fashion as energy and momentum exchange with neutrinos in Equations (6) and (7).

Since we are dealing with a multi-component fluid whose composition is not uniquely determined by the thermodynamic state \( (\rho, T, Y_e) \) unless nuclear statistical equilibrium (NSE) applies, additional conservation equations for the mass fractions \( X_i \) of protons, neutrons, \( \alpha \)-particles, and heavier nuclei are also needed:

\[
\frac{\partial}{\partial t} \sqrt{\gamma} \rho X_k \frac{\partial}{\partial x^i} + \frac{\partial}{\partial t} \sqrt{\gamma} g \rho X_k \hat{v}^i = 0. \quad (9)
\]

Finally, the pressure and the composition (in the NSE regime) in Equations (6), (7), and (9) have to be provided by an EOS.

We rely on the CoCoNuT code (Dimmelmeier et al. 2002a, 2005) to solve the equations of relativistic hydrodynamics in spherical polar coordinates by means of a high-resolution shock-capturing (HRSC) scheme, employing piecewise parabolic (PPM) reconstruction (Colella & Woodward 1984), an approximate Riemann solver, and second-order Runge–Kutta time stepping. Our method of choice for the Riemann solver is a new hybrid HLLC/HLL approximation scheme in CoCoNuT along the lines of
Quirk (1994), combining the high resolution of the HLLC solver (Toro et al. 1994; Mignone & Bodo 2005) with the robustness of the HLLE solver (Harten et al. 1983; Einfeldt 1988) against odd-even decoupling near grid-aligned shocks (Quirk 1994; Liou 2000; Kifonidis et al. 2003; Sutherland et al. 2003).

In order to treat the advection equations for the nuclear mass fractions, we have added a simplified version of the consistent multi-fluid advection (CMA) method by Plewa & Müller (1999). Moreover, the computational performance of the code has been improved significantly by enhancing parallelism in the axisymmetric (two-dimensional) mode.

Most notably, however, our most recent simulations also use a reformulated energy equation instead of Equation (7) to improve total energy conservation. The new scheme, which is described in detail in Appendix A, is designed to alleviate a well-known problem with the gravitational source term in the energy equation: while there is a global conservation law for the ADM energy (or the total energy in the Newtonian case), the gravitational source term in the standard form of the energy equation (7) used in Dimmelmeier et al. (2002a, 2005, 2008) does not have the form of a flux divergence, and therefore a standard finite-volume discretization will lead to a numerical violation of global energy conservation. Nonetheless, as we demonstrate in Appendix A, the gravitational source term can partly be absorbed in a flux divergence term to considerably reduce this conservation error. In the case of a stationary spacetime, or in the Newtonian limit, it is even possible to achieve total energy conservation up to machine precision.

2.2. Metric Equations—The Conformal Flatness Condition

The CoCoNuT code solves Einstein’s field equation using the CFC, which was introduced by Isenberg (2008) and first used in a pseudodynamical context by Wilson et al. (1996). In the CFC approximation, the spatial three-metric is assumed to be conformally flat, i.e., it is obtained from the flat three-metric \( \hat{g}_{ij} \) by a conformal transformation \( g_{ij} = \phi^4 \hat{g}_{ij} \), where \( \phi \) is the conformal factor, thus reducing the number of independent metric quantities to five. As a consequence, the four constraint equations for the metric in the ADM formalism, combined with a slicing condition, determine the metric completely. In our case, we use maximal slicing, i.e., we require the trace of the extrinsic curvature \( K^{ij} \) to vanish: \( K = K^i_i = 0 \). We then obtain a set of five nonlinear elliptic equations for \( \alpha, \phi, \) and \( \beta^i \):

\[
\hat{\Delta} \phi = -2\pi \phi^5 \left( E + \frac{K_{ij} K^{ij}}{16\pi} \right),
\]

\[
\hat{\Delta} (\alpha \phi) = -2\pi \alpha \phi^5 \left( E + 2S + \frac{7K_{ij} K^{ij}}{16\pi} \right),
\]

\[
\hat{\Delta} \beta^i = 16\pi \alpha \phi^4 S^i + 2 \phi^{10} K^{ij} \hat{\nabla}_j \left( \frac{\alpha}{\phi^4} \right) - \frac{1}{3} \hat{\nabla}^i \hat{\nabla}_j \beta^j,
\]

where \( \hat{\Delta} \) and \( \hat{\nabla} \) are the Laplace and covariant derivative operators for a flat three-space, respectively. The total matter-energy density \( E = \tau + D \) (for \( \tau \) and \( D \) see Section 2.1.1), the three-momentum density \( S_j \), and the trace \( S = g_{ij} S^i \) of the spatial components \( S_j \) of the stress–energy tensor (as measured by an Eulerian observer) appear in the nonlinear source terms on the right-hand side, and comprise contributions both from matter and (neutrino) radiation. For the matter component we obtain

\[
E = \rho \hbar W^2 - P,
\]

\[
S = \rho \hbar W^2 v^2 + 3P,
\]

\[
S^i = \rho \hbar W^2 v^i,
\]

while for neutrinos,

\[
E = S = 4\pi \hbar W^2 (J + 2v_r H + v_r^2 K),
\]

\[
S_1 = 4\pi \phi^{-2} \hbar W^2 \left[ H (1 + v^2_r) + v_r (J + K) \right],
\]

\[
S_2 = S_3 = 0,
\]

in our approximation. Here, \( J, H, \) and \( K \) denote the zeroth, first, and second moments of the neutrino radiation intensity in the comoving orthonormal frame (see Section 2.3), and \( v_r = \phi^2 v^i \) is the radial velocity component in the orthonormal basis of an Eulerian observer.

We solve the CFC Equations (10)–(12) in the reformulated version of Cordero-Carrión et al. (2009) using a fixed point iteration scheme as described in Dimmelmeier et al. (2005); i.e., during each iteration step, the nonlinear source terms are updated, and the Laplace operators on the left-hand side of the equations are inverted by means of a multipole expansion (cf. Müller & Steinmetz 1995). While this method suffers from a somewhat low convergence rate compared with the spectral nonlinear Poisson solver of Dimmelmeier et al. (2005), the underlying algorithm is more easily amenable to parallelization, which is a critical issue for multi-dimensional neutrino transport simulations.

2.3. Neutrino Transport in CFC Spacetime

2.3.1. Overview

Up to now, all self-consistent numerical simulations of neutrino transport in core-collapse supernovae are based on the assumptions (1) that a semi-classical treatment of kinetic theory is applicable, and (2) that neutrino oscillations need not be taken into account (see Cardall 2008; Yamada 2000; Strack & Burrows 2005 for a discussion of the more general case). In our approach to the transport problem, we retain those assumptions, although they may be invalid under certain specific circumstances (cf. Section 2.4). Under this proviso, the evolution of the invariant neutrino distribution function \( f \) (the number of neutrinos per phase-space volume \( d^3x \times d^3p \) ) is governed by the relativistic Boltzmann equation for massless or ultra-relativistic particles, which reads (see e.g., Lindquist 1966; Ehlers 1971; Stewart 1971)

\[
p^\mu \frac{\partial f}{\partial x^\mu} + \frac{dp^\mu}{dk} \frac{\partial f}{\partial p^\mu} = \mathcal{C}
\]

in any coordinate basis. Here, \( f \) and the collision term \( \mathcal{C} \) depend on the spacetime coordinates \( x^\mu \) and the four-momentum vector \( p^\mu \) (obeying the mass-shell constraint \( p_\mu p^\mu = 0 \) ) as measured in the associated holonomic base \( \partial_\mu ; \lambda \) is the affine parameter.

In general, Equation (19) describes a six-dimensional time-evolution problem, which can be scaled down considerably by using the ray-by-ray-plus approximation (Buras et al. 2006b) in the following manner. First, the equations of neutrino transport

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5 We note in passing that this assumption can always be fulfilled by an appropriate gauge choice for a spherically symmetric spacetime, and no approximations to the ADM equations are actually made by using CFC in that case. See Section 2.4 for more details.

6 Thus, in spherical polar coordinates we have \( \hat{g}_{ij} = \text{diag}(1, r^2, r^2 \sin^2 \theta) \).
are formulated for the spherically symmetric case (Section 2.3.2.2 to 2.3.5), thus reducing the dimensionality of the individual transport problems to be solved from six to three. By working with a finite number (i.e., two) of angular moments of the Boltzmann equation, yet another dimension can be eliminated.

The additional space dimensions are then taken into account by solving these equations independently on different radial “rays” (corresponding to the angular zones of the polar grid) with direction unit vector \( n \), assuming rotational symmetry around \( n \) for the neutrino distribution function. In addition, certain terms from the full transport equations in two dimensions (or three dimensions in general) are taken into account to avoid unphysical behavior in the optically thick regime (see Buras et al. 2006b and Section 2.3.6). We emphasize that the ray-by-ray-plus approximation does not mean that the transport of neutrinos is considered to be only in the radial direction. Instead, it is assumed that the neutrino distribution function is axially symmetric around the radial direction, and that the neutrino pressure tensor is diagonal.

### 2.3.2. Boltzmann Equation in Spherical Symmetry

The precise form of Equation (19) in a spherically symmetric spacetime has been discussed in various places in the literature (Lindquist 1966; Castor 1972; Schinder 1988; Mezzacappa & Matzner 1989); it depends on the adopted gauge and slicing conditions, and on the tetrad basis used in momentum space. In our case, the spacetime metric is of the CFC type with \( \beta^1 = \psi \) and \( \beta^2 = \beta^3 = 0 \),

\[
\begin{align*}
0 = & -\alpha(t, r)^2 dt^2 + \phi(t, r)^4 \times [(dr + \beta(t, r) dt)^2 + r^2 d^2 + r^2 \sin^2 \theta d\alpha^2].
\end{align*}
\]

Furthermore, we carry out a transformation to the orthonormal frame comoving with the fluid, in which the collision integral \( C \) can be expressed most conveniently; the expressions for the collision integral provided in Appendix A of Rampp & Janka (2002) thus remain valid. With the velocity field chosen as

\[
(v^1, v^2, v^3) = (\phi^{-2} v_\psi (t, r), 0, 0),
\]

a possible choice for the basis four-vectors of such a frame in terms of the coordinate derivatives \( \partial_t, \partial_r, \partial_\theta, \) and \( \partial_\phi \) (see Hawking & Ellis 1973; Misner et al. 1973; Straumann 2004 for this notation) is given by

\[
\begin{align*}
\mathbf{e}_0 &= \alpha^{-1} W \partial_t + W(\phi^{-2} v_\psi - \alpha^{-1} \beta^\psi) \partial_r, \\
\mathbf{e}_1 &= \alpha^{-1} v_\psi W \partial_t + W(\phi^{-2} - \alpha^{-1} \beta^\psi v_\psi) \partial_r, \\
\mathbf{e}_2 &= \phi^{-2} r^{-1} \partial_\theta, \\
\mathbf{e}_3 &= \phi^{-2} r^{-1} \sin^{-1} \theta \partial_\phi.
\end{align*}
\]

In spherical symmetry, the distribution function \( f \) depends only on \( t, r \), the comoving frame energy \( \varepsilon = \mathbf{p} \cdot \mathbf{e}_0 \) (where \( \mathbf{p} \) is the neutrino four-momentum), and the angle cosine \( \mu = \mathbf{p} \cdot \mathbf{e}_1 / \mathbf{p} \cdot \mathbf{e}_0 \), and obeys the following PDE:

\[
\begin{align*}
W \left[ \frac{\xi}{\alpha} \frac{\partial f}{\partial t} - \beta^r \frac{\partial f}{\partial r} \right] + v \frac{\partial f^*}{\partial \alpha} - \varepsilon W^3 \frac{\partial f}{r \alpha \phi^3 \varepsilon} \\
\times \left[ \beta^r \phi^3 \left( -\psi - r \mu \frac{\partial v_\psi}{\partial r} \right) + \phi^2 \phi_r \left( 2r \frac{\partial \phi}{\partial r} - \phi \phi_r \right) \right] + r \left( -\mu \frac{\partial \alpha}{\partial r} + \mu^2 \phi^2 \frac{\partial \beta^r}{\partial r} - \phi \frac{\partial \phi}{\partial r} \right) \\
+ v^3 \left( -\mu \frac{\partial \alpha}{\partial r} + \phi \frac{\partial \phi^2}{\partial r} - \phi \frac{\partial \phi^2}{\partial \alpha} \right) - \psi \alpha \frac{\partial \phi^2}{\partial r} \\
+ \phi \left( \frac{\partial \phi^2}{\partial t} - r \beta^r \frac{\partial \phi^2}{\partial r} \right) + v \alpha \left( \phi \frac{\partial \psi}{\partial r} + \mu \frac{\partial \phi^2}{\partial \alpha} \right) \\
+ \phi \left( \frac{\partial \phi^2}{\partial t} - r \beta^r \frac{\partial \phi^2}{\partial r} \right) + v \alpha \left( \phi \frac{\partial \psi}{\partial r} + \mu \frac{\partial \phi^2}{\partial \alpha} \right) \\
+ 2r \psi \frac{\partial \phi^2}{\partial r} + \phi \left( \frac{\partial \phi^2}{\partial \alpha} - \phi \frac{\partial \phi^2}{\partial \alpha} \right) \\
+ W^3 \left( 1 - \mu^2 \right) \frac{\partial f}{\partial \alpha} + \phi \left( \frac{\partial \phi^2}{\partial t} - r \beta^r \frac{\partial \phi^2}{\partial r} \right) \\
+ \phi \left( \beta^r \phi^2 \left( r \frac{\partial v_\psi}{\partial r} - v \right) - r \phi^2 \left( \frac{\partial \phi^2}{\partial \alpha} - \phi \frac{\partial \phi^2}{\partial \alpha} \right) \right) \right] = \mathcal{C} f.
\end{align*}
\]

where \( \mu = v + \psi, \xi = 1 + \mu v_\psi \) (bearing in mind that \( c = 1 \)) and \( \psi = 1 - \mu^2 \). It should be noted that our result is analytically equivalent to the one obtained by Mezzacappa & Matzner (1989) for the same gauge and slicing conditions. The transfer equation for the radiation intensity \( J = h^{-3} c^{-2} \varepsilon^3 \) can be obtained by exploiting the relation \( \varepsilon^3 \frac{\partial f}{\partial \varepsilon} = \partial (f \varepsilon^3) / \partial \varepsilon - 3 \varepsilon^2 \), and is omitted here for the sake of brevity.

### 2.3.3. Exact Moment Equations for CFC Metric in Spherical Symmetry

Multiplying Equation (26) by \( h^{-3} c^{-2} \varepsilon^3 \) and taking the zeroth and first angular moments yields the moment equations in a spherically symmetric CFC spacetime, resulting in

\[
\begin{align*}
\frac{\partial W}{\partial t} + \frac{\partial}{\partial r} \left[ \left( Wv_\psi - \beta^r v_\psi \right) \hat{H} + \left( Wv_\alpha - \beta^\alpha \right) \hat{J} \right] \\
- \frac{\partial}{\partial r} \left[ \frac{W}{r \phi^3} \left( \frac{1}{r} \beta^r - \alpha v_\psi \phi^2 \right) + 2 \beta^r \phi^2 \phi_\alpha \frac{\partial \ln \phi}{\partial r} \right] - 2 \phi_\psi \phi_r \frac{\partial \phi^2}{\partial r} + \phi \left( \beta^r \phi^2 \left( r \frac{\partial v_\psi}{\partial r} - v \right) - r \phi^2 \left( \frac{\partial \phi^2}{\partial \alpha} - \phi \frac{\partial \phi^2}{\partial \alpha} \right) \right) \right] \\
- W \hat{J} \left( \frac{1}{r} \beta^r - \alpha v_\psi \phi^2 \right) + 2 \beta^r \phi^2 \frac{\partial \ln \phi}{\partial r} - 2 \frac{\partial \ln \phi}{\partial r} \right] \\
- \frac{\partial}{\partial r} \left[ Wv_r \left( \phi_\alpha - \alpha v_\psi \phi^2 \right) \frac{\partial \ln \phi}{\partial r} - \alpha \phi \frac{\partial \ln \alpha W}{\partial r} \right] - \phi_\psi \phi_r \frac{\partial \phi^2}{\partial r} + \phi \left( \beta^r \phi^2 \left( r \frac{\partial v_\psi}{\partial r} - v \right) - r \phi^2 \left( \frac{\partial \phi^2}{\partial \alpha} - \phi \frac{\partial \phi^2}{\partial \alpha} \right) \right) \right] \\
- W \hat{J} \left( \frac{1}{r} \beta^r - \alpha v_\psi \phi^2 \right) + 2 \beta^r \phi^2 \frac{\partial \ln \phi}{\partial r} - 2 \frac{\partial \ln \phi}{\partial r} \right] \\
- \frac{\partial}{\partial r} \left[ Wv_r \left( \phi_\alpha - \alpha v_\psi \phi^2 \right) \frac{\partial \ln \phi}{\partial r} - \alpha \phi \frac{\partial \ln \alpha W}{\partial r} \right] + \hat{K} \left( \beta^r \phi^2 \left( r \frac{\partial v_\psi}{\partial r} - v \right) - r \phi^2 \left( \frac{\partial \phi^2}{\partial \alpha} - \phi \frac{\partial \phi^2}{\partial \alpha} \right) \right) \right] = \alpha \hat{C}^{(0)},
\end{align*}
\]

for the energy equation and

\[
\frac{\partial W}{\partial t} + \frac{\partial}{\partial r} \left[ \left( Wv_\psi - \beta^r v_\psi \right) \hat{H} + \left( Wv_\alpha - \beta^\alpha \right) \hat{J} \right] \]
for the momentum equation. Here, the angular moments $J$, $H$, $K$, and $L$ of the specific intensity are given by

$$\{J, H, K, L\} = \frac{1}{2} \int_{-1}^{1} d\mu \mu^{(1, 2, 3)} \mathcal{I},$$

while $C^{(0)}$ and $C^{(1)}$ are the corresponding zeroth and first moment of the collision integral. Note that Equations (27) and (28) are written in terms of the densitized moments $\hat{J}$, $\hat{H}$, $\hat{K}$, and $\hat{L}$ (where $\hat{X} = \sqrt{\gamma} X$ for any quantity $X$) which contain the square root of the determinant of the three-metric $\sqrt{\gamma} = \phi^{\alpha\beta}\sin\theta$ as an additional factor, in order to better reflect the underlying law of energy conservation. The virtue of such a formulation is more apparent in the equations governing the evolution of the neutrino number density and flux, 

$$\frac{\partial}{\partial t} \left[ W \alpha \frac{\partial}{\partial r} \right] + \frac{\partial}{\partial r} \left[ \left( W \alpha \frac{\partial}{\partial r} - \beta' v_r \right) \hat{J} \right] + \frac{\partial}{\partial t} \left[ \left( W \alpha \frac{\partial}{\partial r} - \beta' v_r \right) \hat{H} \right] + \frac{\partial}{\partial t} \left[ \left( W \alpha \frac{\partial}{\partial r} - \beta' v_r \right) \hat{K} \right] + \frac{\partial}{\partial t} \left[ \left( W \alpha \frac{\partial}{\partial r} - \beta' v_r \right) \hat{L} \right] = 0,$$

where $\{J, H, K, L\} = \{\epsilon^{-1} J, \epsilon^{-1} H, \epsilon^{-1} K, \epsilon^{-1} L\}$, $\epsilon^{-1} C$. Apart from the collision integral, no source terms appear in the neutrino number equation (30). Since the evolved quantity $W(\mathcal{J} + v_r \mathcal{H})$ is proportional to the spectral neutrino number density $N_{\text{eul}}$ in the Eulerian (coordinate/lab) frame,

$$N_{\text{eul}} = \frac{4\pi}{c} W(\mathcal{J} + v_r \mathcal{H}),$$

a finite-volume discretization of Equation (30) exactly conserves the total neutrino number measured in the laboratory (coordinate) frame,

$$\int d^3x \int ds \sqrt{\gamma} N_{\text{eul}} = \frac{4\pi}{c} \int d^3x \int ds \sqrt{\gamma} W(\mathcal{J} + v_r \mathcal{H}),$$

in the absence of neutrino–matter interactions. As a consequence, numerical conservation of lepton number can be achieved with an appropriate conservative treatment of the flux terms in the neutrino moment equations and the source term for the electron fraction.

With a few minor exceptions, the numerical solution of the neutrino moment equations in our code proceeds exactly as described by Rampp & Janka (2002). Equations (27) and (28) are written in a conservative finite-difference form using second-order space discretization on a staggered mesh and backward-differencing in time; only the neutrino advection terms

$$\frac{\partial}{\partial r} \left[ \left( W \alpha \frac{\partial}{\partial r} - \beta' v_r \right) \hat{J} \right], \quad \frac{\partial}{\partial r} \left[ \left( W \alpha \frac{\partial}{\partial r} - \beta' v_r \right) \hat{H} \right],$$

are discretized with upwind differences in space and central differences in time. The higher moments $K$ and $L$ are computed from the variable Eddington factors $f_K$ and $f_L$ obtained from the solution of a model Boltzmann equation (cf. Section 2.3.4) as $K = f_K J$ and $L = f_L J$ where necessary.

Our numerical scheme for the moment equations differs from that of Rampp & Janka (2002) in one important respect. Rampp & Janka (2002) also solve Equations (30) and (31) for the neutrino number density and flux in addition to Equations (27) and (28) in order to guarantee both lepton number conservation and energy conservation. However, this procedure has
two important drawbacks: solving another pair of equations is computationally expensive, and, more importantly, the numerical solution can become inconsistent in the sense that \((J, H) = (\epsilon, f, \epsilon \tau)\) no longer holds. In the high-energy tail of the spectrum \(J_f, J_{\bar{f}}\) and \(H_f/H_{\bar{f}}\) in a given zone in energy space can even assume values outside the zone boundaries. In order to circumvent this problem, we have developed a special second-order finite-difference representation for the terms governing advection in energy space accounting for gravitational redshift and Doppler shift. This allows us to guarantee both energy and lepton number conservation without solving the equations for neutrino number density and flux. Since our improved treatment of the redshift and Doppler terms is rather involved, we give a complete description of the algorithm in Appendix B.

Finally, the moment equations are supplemented by two additional equations for the operator-split update of the specific internal energy and the electron fraction of the matter due to neutrino–matter interactions,

\[
\left( \frac{\partial \epsilon}{\partial t} \right)_v = -\frac{4\pi}{\rho W} \int_0^\infty d\epsilon \sum_{i \in \{e, \bar{e}, e_{\nu}, \bar{e}_{\nu}\}} C_i^{(0)},
\]

\[
\left( \frac{\partial Y_e}{\partial t} \right)_v = -\frac{4\pi}{\rho W} \int_0^\infty d\epsilon \left( C_i^{(0)} - C_i^{(0)} \right),
\]

where the factor \(\alpha W^{-1}\) takes care of the conversion from proper time to coordinate time. The resulting nonlinear system of equations for \(J, H, \epsilon,\) and \(Y_e\) is then solved by Newton–Raphson iteration. As in Rampp & Janka (2002), we treat \(e\) and \(\bar{e}\) neutrinos and their antiparticles as single species, and solve the corresponding moment equations separately from the (fully coupled) moment equations for the transport of electron neutrinos and antineutrinos.

### 2.3.4. Model Boltzmann Equation

In the variable Eddington factor technique, the closure relations required for Equations (27) and (28) are extracted from the formal solution of a model Boltzmann equation, with the moments of the neutrino distribution function occurring in the collision integral taken from the solution of the moment equations in an iterative procedure (Mihalas & Weibel Mihalas 1984; Burrows et al. 2000; Rampp & Janka 2002). Since the model Boltzmann equation is only used to compute normalized moments of the radiation intensity, it is sufficient to consider a strongly simplified version of Equation (26). We therefore neglect terms containing the shift vector or derivatives of the metric functions \(\alpha\) and \(\phi\) in Equation (26), as well as higher-order terms \(O(v^4)\) in the radial velocity. Furthermore, we set \(v = \mu + v_r \rightarrow \mu\) and \(\xi = 1 + \mu v_r \rightarrow 1\) in the terms for advection in energy space. However, the radial advection term is kept exactly as in the full equation. The model transfer equation for the radiation intensity thus obtained reads

\[
\begin{align*}
\frac{\partial T}{\partial t} + \left( \frac{\alpha v_r}{\phi^2} - \beta^* \right) \frac{\partial f}{\partial r} + \frac{\alpha \mu}{\phi^2} \frac{\partial f}{\partial \mu} + \int \left[ (3 - \mu^2) \frac{\alpha v_r}{r \phi^2} \\
+ (1 + \mu^2) \alpha \phi \frac{\partial v_r}{\partial r} + 2 \alpha \mu \frac{\partial v_r}{\partial \mu} \right] - \frac{\partial}{\partial \epsilon} \left[ \epsilon \left( 1 - \mu^2 \right) \frac{\alpha v_r}{\phi^2} \right] \\
+ \mu^2 \frac{\alpha}{\phi^2} \frac{\partial v_r}{\partial r} + \alpha \frac{\partial v_r}{\partial \mu} \right] + \alpha \left( 1 - \mu^2 \right) \frac{\partial T}{\partial \mu} + \frac{\partial}{\partial \mu} \left[ \mu \left( \frac{\alpha v_r}{r \phi^2} - \frac{\alpha}{\phi^2} \frac{\partial v_r}{\partial r} \right) - \mu \alpha \frac{\partial v_r}{\partial \mu} \right] \right] = C[I].
\end{align*}
\]

Here, \(C[I]\) denotes the collision term in the transfer equation, which is related to the collision term \(C\) in the Boltzmann Equation (26) by \(C[I] = \epsilon \pi h^2 c^{-2} [f, \bar{f}]\). Aside from the appearance of prefactors containing the metric functions \(\alpha\) and \(\phi\) and of the radial shift \(\beta^*\) in the advection term, Equation (37) is identical to the flat-space transfer equation in the \(O(v/c)\) approximation. Therefore, the same solution method as in Rampp & Janka (2002) can be employed, i.e., the radial advection term and the energy derivatives in Equation (37) are treated by a conservative interpolation procedure and a time-explicit upwind scheme respectively, while the tangent-ray method (Yorke 1980; Mihalas & Weibel Mihalas 1984) is used for the remaining terms.

#### 2.3.5. Source Terms

Once the solution of the moment equations is known, the source terms \(Q_Y, Q_E,\) and \(Q_M\) for electron number, energy, and momentum in the local fluid frame can be computed from the zeroth and first angular moments \(C^{(0)}\) and \(C^{(1)}\) of the collision integral,

\[
\begin{align*}
Q_Y &= \left( \frac{d\rho Y_e}{d\lambda} \right)_C = -4\pi m_B \int_0^\infty d\epsilon \left( C_{\nu e}^{(0)} - C_{\bar{\nu} e}^{(0)} \right), \\
Q_E &= \left( \frac{d\rho e}{d\lambda} \right)_C = -4\pi \int_0^\infty d\epsilon \sum_{i \in \{e, \bar{e}, e_{\nu}, \bar{e}_{\nu}\}} C_i^{(0)}, \\
Q_M &= -4\pi \int_0^\infty d\epsilon \sum_{i \in \{e, \bar{e}, e_{\nu}, \bar{e}_{\nu}\}} C_i^{(1)}.
\end{align*}
\]

To obtain the correct transformation to the Eulerian frame (in which the equations of hydrodynamics are solved), we consider the equations of electron number conservation and energy–momentum conservation in their covariant form. Source terms due to neutrino interactions must enter as scalar and four-vector terms on the right side in these equations:

\[
\nabla_{\mu}(\rho Y_e u^\mu) = s, \quad \nabla_{\mu} T^{\mu\nu} = q^\nu.
\]

Here, \(s\) and \(q^\nu\) are the scalar electron number density source term and the covariant energy–momentum vector source term. Since the transformation of scalars and four-vectors is trivial, we only need to express \(s\) and \(q^\nu\) in terms of \(Q_Y, Q_E,\) and \(Q_M\) in the comoving frame to obtain the correct source terms in any desired frame:

\[
\begin{align*}
Q_Y &\rightarrow Q_{Ye}, \quad q^\nu = Q_{Ye} e_0 + Q_{Ye} e_1.
\end{align*}
\]

After carrying out a Lorentz transformation back to the Eulerian frame and recasting the resulting equations into conservative form, we obtain the following source terms for the conserved Eulerian quantities that appear in Equations (6)–(8):

\[
\begin{align*}
\left( \frac{\partial \sqrt{\gamma} S_i}{\partial t} \right)_c &= \sqrt{\gamma} W(v_r, Q_E + Q_M), \\
\left( \frac{\partial \sqrt{\gamma} T}{\partial t} \right)_c &= \sqrt{\gamma} W(Q_{Ye} + v_r Q_M), \\
\left( \frac{\partial \sqrt{\gamma} \rho W Y_e}{\partial t} \right)_c &= \sqrt{\gamma} Q_{Ye}.
\end{align*}
\]

\(^7\) Since we consider a purely radial velocity field at this stage, the Lorentz transformation of a four-vector from the comoving frame to the Eulerian frame simply reads \((q^0, q^1, q^2, q^3) \rightarrow (W(q^0 + v_r q^1, q^1 + v_r q^0, q^2, q^3)).\)
Note that we have multiplied the source terms by the metric \(\sqrt{g}\), because it is the densitized variables \(\sqrt{g}S_i, \sqrt{g}T_\tau\), and \(\sqrt{g}D\gamma\), that are actually evolved in Equations (6)–(8). This is possible because the metric factor \(\sqrt{g}\) is obviously not changed by neutrino interactions.

### 2.3.6. Ray-by-ray-plus Approach for Axisymmetric Problems

As noted by Rampp & Janka (2002), a treatment of neutrino transport in spherical symmetry (as described in Sections 2.3.3 to 2.3.5) can be extended to the multi-dimensional case by assuming that all the involved physical quantities depend only parametrically on the coordinates \(\theta, \phi\) (using spherical polar coordinates), while neglecting any lateral derivatives in the full three-dimensional moment equations. However, Buras et al. (2006b) pointed out that the lateral advection of neutrinos in the optically thick regime and the non-radial components of the neutrino pressure gradient need to be taken into account to avoid unphysical convective activity in the proto-neutron star. In general relativity, the additional advection terms in the moment equations are

\[
\left(\frac{\partial \sqrt{g}W (J + v_\phi H)}{\partial t}\right)_{\text{adv}} + \alpha \frac{\partial \sqrt{g}W \dot{v}_\phi^2 (J + v_\phi H)}{\partial \phi} = 0,
\]

\[
\left(\frac{\partial \sqrt{g}W (H + v_\phi K)}{\partial t}\right)_{\text{adv}} + \alpha \frac{\partial \sqrt{g}W \dot{v}_\phi^2 (H + v_\phi K)}{\partial \phi} = 0,
\]

(46)

(47)

which are solved in an operator-split step before the radial transport sweep. We also include terms for the compressional heating of the neutrinos due to lateral motions, which are added to the moment Equations (27) and (28) when performing the radial transport sweep,

\[
\left(\frac{\partial \sqrt{g}(J + v_\phi H)}{\partial t}\right)_{\text{comp}} = -\frac{\partial \alpha \sqrt{g}W \dot{v}_\phi^2 J - K}{\partial \phi} \left(\frac{\partial}{\partial \phi}\right) - \frac{\partial}{\partial \phi}
\]

\[
\times \left(\frac{1}{2} \sqrt{g}W \dot{v}_\phi^2 J - K \right),
\]

(48)

\[
\left(\frac{\partial \sqrt{g}(H + v_\phi K)}{\partial t}\right)_{\text{comp}} = -\frac{\partial \alpha \sqrt{g}W \dot{v}_\phi^2 K - L}{\partial \phi} \left(\frac{\partial}{\partial \phi}\right) - \frac{\partial}{\partial \phi}
\]

\[
\times \left(\frac{1}{2} \sqrt{g}W \dot{v}_\phi^2 H - L \right).
\]

(49)

It should be noted that gravitational redshift associated with lateral motion is still neglected here; however, this is a tiny effect in typical supernova environments due to the relatively small asphericity of the gravitational field.

The lateral component of the neutrino pressure gradient appears as a source term for the component \(S_2\) of the fluid momentum density,

\[
\left(\frac{\partial \sqrt{g}S_2}{\partial t}\right)_{\text{lat}} = -\sqrt{-g} \frac{\partial P_\nu}{\partial \phi},
\]

(50)

where we assume that the neutrino pressure is given by \(P_\nu = 4/3\pi J\) in the optically thick regime. Since the acceleration four-vector due to the neutrino pressure gradient must be orthogonal to the four-velocity of the fluid, the source term in the energy equation (work exerted on the fluid) is given by

\[
\left(\frac{\partial \sqrt{g}T_\tau}{\partial t}\right)_{\text{lat}} = -\sqrt{-g} v_\phi \frac{\partial P_\nu}{\partial \phi}.
\]

(51)

Finally, contrary to the Newtonian case, the covariant source vector \(q^{\mu}\) has to be transformed to the Eulerian frame by a Lorentz boost which includes the non-radial velocity components. \(Q_E\) and \(Q_M\) now enter differently into the source term for \(S_1\), and additional source terms for \(S_2\) and \(S_3\) appear,

\[
\left(\frac{\partial \sqrt{g}S_1}{\partial t}\right)_{c} = \sqrt{g} \left[ W v_1 Q_E + Q_M \left(1 + v_1 v_\phi \frac{W - 1}{v_\phi v_\phi} \right) \right],
\]

(52)

\[
\left(\frac{\partial \sqrt{g}S_2}{\partial t}\right)_{c} = \sqrt{g} \left[ W v_2 Q_E + Q_M v_2 v_\phi \frac{W - 1}{v_\phi v_\phi} \right],
\]

(53)

\[
\left(\frac{\partial \sqrt{g}S_3}{\partial t}\right)_{c} = \sqrt{g} \left[ W v_3 Q_E + Q_M v_3 v_\phi \frac{W - 1}{v_\phi v_\phi} \right].
\]

(54)

While the new terms in \(Q_M\) are of order \(\mathcal{O}(v^2)\) and can usually be neglected, all the terms in \(Q_E\) are of order \(\mathcal{O}(v)\) and need to be included to ensure the correct evolution of the specific internal energy \(\epsilon\). This is particularly crucial for rapidly rotating configurations, where the rotational velocity \(v_\phi\) can reach a significant fraction of \(c\).

### 2.4. Quality of Our Approximation

Our relativistic ray-by-ray-plus method entails several approximations, which fall into three categories, viz., approximations in the treatment (1) of the microphysics (EOS, neutrino interactions, etc.), (2) of the neutrino transport, and (3) of the gravitational field equations. Concerning the microphysics, we rely on the same up-to-date description as the non-relativistic VERTEX-PROMETHEUS code (Rampp & Janka 2002; Buras et al. 2006b), which includes a number of additional neutrino reactions compared to the “standard” set of neutrino opacities used by Bruenn (1985; see Burrows & Sawyer 1998; Hennestad & Raffelt 1998; Buras et al. 2003; Itoh et al. 2004; Marek et al. 2005; Langanke et al. 2003, 2008 for the extensions to the standard opacities). Neutrino oscillations are presently neglected despite the fact that the classical Mikheyev–Smirnov–Wolfenstein (MSW) effect and collective neutrino–antineutrino flavor oscillations may possibly have an important bearing at least on the observable neutrino properties. However, considering that our understanding of the latter phenomenon is by no means complete, and that the classical MSW effect is only relevant far outside the supernova core, it is not certain whether neutrino oscillations are dynamically relevant for the supernova problem at all. Moreover, no self-consistent dynamical simulations including these effects have been carried out so far; recent studies have contented themselves with post-processing progenitor or simulation data (Duan et al. 2008; Lunardini et al. 2008). Thus, our treatment of neutrino physics, while certainly not complete, is state of the art for dynamical supernova simulations, and the same is true for the different high-density EOSs (Lattimer & Swesty 1991; Shen et al. 1998; Hillebrandt & Wolff 1985) available in our code.

Handling the neutrino transport with the “ray-by-ray-plus” scheme of Buras et al. (2006b) introduces approximations on two different levels: first, the variable Eddington factor technique is not strictly equivalent to full Boltzmann transport
even in spherical symmetry, since only a simplified "model" Boltzmann equation is solved to provide the closure for the moment equations. This shortcoming is mitigated by the fact that comparisons with analytic solutions (Rampp & Janka 2002) and with Boltzmann solvers based on the $S_N$-method (Liebendörfer et al. 2005) show that the approximate solution is entirely adequate in many cases. Second, the "ray-by-ray-plus+" approach ignores the lateral propagation (but not the lateral advection) of neutrinos. However, the analysis of Buras et al. (2006b) at least suggests that this approximation does not affect the heating in the gain layer dramatically, and one may therefore expect that it quite adequately captures the overall dynamics during the proto-neutron star accretion phase and also during the subsequent wind phase. Unfortunately, no detailed comparison of the "ray-by-ray-plus" method with true multi-angle transport or alternative approaches, like MGFLD, which would allow us to better assess the quality of our method, is available as yet. So far, the different approaches to multi-dimensional transport hitherto pursued by supernova modelers all have their individual strengths and weaknesses: the only true multi-angle simulations (Livne et al. 2004; Ott et al. 2008), for example, have been carried out at the expense of neglecting inelastic scattering reactions, and MGFLD has been shown to smear out lateral variations of the radiation field quite considerably (Ott et al. 2008), while the "ray-by-ray-plus" method probably tends to overestimate such variations. Still, simulations with ray-by-ray transport show effects that are qualitatively similar to those seen by Ott et al. (2008) with multi-angle transport. Global asymmetries of the neutrino flux are preserved better than with MGFLD, for example. At any rate, ray-by-ray transport can be classified as a state-of-the-art method in the field, notwithstanding the fact that it is certainly limited to scenarios where the anisotropies in the matter distribution (at least out to the neutrinosphere) do not become exceedingly large. Compared to other radiative transfer methods used in multi-dimensional relativistic simulations up to now—like the deleptonization scheme of Liebendörfer (2005), which was first used in GR by Ott et al. (2007b), or the gray two-moment formalism (using the Eddington approximation) by Farris et al. (2008)—it is a much more ambitious and accurate approach to neutrino transport in core-collapse supernovae.

Given the limitations of the "ray-by-ray-plus" method, the use of the CFC approximation to the gravitational field equations is quite natural: CFC deteriorates in accuracy for strongly deformed, rapidly rotating matter configurations, where the "ray-by-ray-plus" method is not applicable anyway—there is little to be gained by a more accurate treatment of GR in that regime. It is noteworthy, however, that the CFC approximation has been found to be accurate for a wide range of (sometimes extreme) astrophysical scenarios, such as rapidly rotating neutron stars (Cook et al. 1996), rigidly rotating disks (Kley & Schäfer 1999), and rotational core collapse (Shibata & Sekiguchi 2004; Cerdá-Durán et al. 2005; Ott et al. 2007a, 2007b; Dimmelmeier et al. 2007). In the core-collapse scenario, CFC still appears to be adequate (Dimmelmeier et al. 2006) for the extremely rapidly rotating models of Shibata & Sekiguchi (2005); even toroidally deformed neutron stars obviously pose little problem to CFC. These facts indicate that in typical supernova simulations, where the neutron star rotates moderately at most, CFC will not be the weakest link in the chain of approximations, and that, at least for our purpose, GR effects are treated correctly in our code. All in all, VERTEX-CoCoNuT can thus be regarded as a truly relativistic radiation hydrodynamics code for core-collapse supernovae, featuring a sophisticated neutrino transport scheme and an up-to-date treatment of the microphysics.

3. CODE TESTS IN SPHERICAL SYMMETRY

3.1. Simple Test Problems—Radiating Spheres

Stationary solutions for the propagation of radiation from an extended spherical central source into a vacuum furnish one of the simplest kinds of test problems in radiation hydrodynamics. Although not overly challenging per se, they provide a useful consistency check for the numerical implementation of the relativistic neutrino moment equations. The principle underlying these tests may be illustrated by considering the neutrino energy Equation (27) in flat space, assuming stationarity and specializing to the laboratory frame (where $v_r \equiv 0$).

$$\frac{1}{r^2} \frac{\partial (r^2 H_{\text{rad}})}{\partial r} = C^{(0)}. \quad (55)$$

Outside the central source (which may be chosen as a homogeneous isothermal sphere of radius $R$ with vanishing scattering opacity and frequency-independent absorption opacity as a very simple case), the collision integral vanishes, and hence the lab frame luminosity remains constant,

$$L_{\text{rad}} = 16 \pi^2 H_{\text{rad}} r^2 = \text{const.} \quad (56)$$

Likewise, since no energy-exchanging reactions occur, and since all the Doppler terms in the moment equations vanish, the mean energy of the radiation is independent of the radius $r$ in the vacuum region:

$$\langle \epsilon \rangle_{\text{rad}} = H_{\text{rad}}/H_{\text{tot}} = \text{const.} \quad (57)$$

The luminosity and mean energy as measured by moving observers can be constructed from the transformation properties of the radiation moments $J$, $H$, and $K$ (see, e.g., Mihalas & Weibel Mihalas 1984). Since $J = H = K$ for $r \gg R$, we have

$$L W^2 (1 + v_r^2 + 2 v_r) = L \frac{1 + v_r}{1 - v_r} = \text{const}, \quad (58)$$

$$\langle \epsilon \rangle W (1 + v_r) = \text{const},$$

for the luminosity $L = 16 \pi^2 H/c$ and the mean energy $\langle \epsilon \rangle$ in the frame of a moving observer.

**Test problem 1.** We have compared the numerical solution computed with VERTEX-CoCoNuT in the comoving frame to this analytical solution for the case of a radiating sphere with $R = 4$ km and an observer velocity field with a peak value of $v = -0.2c$ at $r = 150$ km.

$$v_r = \begin{cases} 0, & r < 135 \text{ km} \\ -0.2c & 135 \text{ km} \leq r < 150 \text{ km} \\ -0.2c \left( \frac{150 \text{ km}}{r} \right)^2, & 150 \text{ km} \leq r. \end{cases} \quad (59)$$

This setup mimics the velocity profile typically encountered in the accretion phase of core-collapse supernovae with a strong shock discontinuity at radii of order $\sim 100$ km. The left panel of Figure 1 shows that the numerical solution is in excellent agreement with Equation (58), with a maximum deviation of less than 2%. Considering that the finite-difference representation of the moment equations has not been specifically tuned to
curved spacetime, assuming that the metric is diagonal and the propagation of radiation from an isothermal sphere in a redshift terms can be tested in a similar fashion: we now consider factor corrections can be taken into account analytically by including another convenience, coarse ($\Delta\varepsilon/\varepsilon \approx 10^{-2}$) reproduces the analytic result, and that the energy grid is rather seems more than adequate, and indicates that the velocity-dependent terms in the moment equations have been correctly implemented.

**Test problem 2.** The implementation of the gravitational redshift terms can be tested in a similar fashion: we now consider the propagation of radiation from an isothermal sphere in a curved spacetime, assuming that the metric is diagonal and isotropic, i.e., $ds^2 = -\alpha^2 dt^2 + \phi^4 (dr^2 + r^2 d\theta^2 + \sin^2 \theta d\phi^2)$. The frequency-integrated neutrino energy and number equations (obtained from Equations (27) and (28) after setting $\beta^r \equiv v_r \equiv 0$ and $W \equiv 1$) can then be solved directly in the vacuum region,

$$\alpha \phi^4 r^2 = \text{const.}, \quad \alpha^2 \phi^4 H r^2 = \text{const.}, \quad \alpha \langle \varepsilon \rangle = \text{const.}$$

(60)

To test our numerical scheme, we have chosen a lapse function $\alpha(r)$ from a pseudo-relativistic core-collapse simulation with the VERTEX-PROMETHEUS code (Rampp & Janka 2002; with a minimum value of $\alpha(0) \approx 0.72$), thus mimicking a potential well of similar strength as in real applications. For convenience, the conformal factor $\phi$ is set to $\phi = \sqrt{\alpha^{-1}}$ as in the weak-field limit. Again, the luminosity and mean energy are in very good agreement with the analytic results for this test case, as shown in the right panel of Figure 1. The redshift-corrected luminosity $L_{rs}$ is given by $16\pi \alpha \phi^4 H r^2$, which should be constant outside the radiating sphere for our specific choice for $\alpha$ and $\phi$.

**Test problem 3.** Test problems 1 and 2 can be easily combined to serve as a more stringent check for the implementation of the moment equations (as more and more non-vanishing terms enter), if we consider non-Eulerian observers in a curved spacetime. The redshift-corrected luminosity $L_{rs}$ and mean energy as measured by an observer moving with a velocity $v$, then obey the following simple equations:

$$\frac{1 + v_r}{1 - v_r} L_{rs} = 16\pi \alpha^2 \phi^4 \frac{1 + v_r}{1 - v_r} H r^2 = \text{const.}$$

$$W\alpha(1 + v_r)\langle \varepsilon \rangle = \text{const.}$$

(61)

With the same choice for $v_r$, $\alpha$, and $\phi$ as in test problems 1 and 2, VERTEX-CoCoNuT once more reproduces the analytic solution to very good accuracy (see Figure 2).

### 3.2. Core Collapse in Spherical Symmetry

#### 3.2.1. Aims and Test Setup

While the accuracy of VERTEX-CoCoNuT for the simple test problems presented in the last section is reassuring, only a comparison with existing one-dimensional GR neutrino transport codes can serve as a true shakedown test for future applications in “uncharted territory” (i.e., multi-dimensional GR neutrino transport). Fortunately, such codes are available (Yamada et al. 1999; Liebendörfer et al. 2001b, 2001a, 2004). A detailed comparison of the relativistic AGILE-BOLTZTRAN and the Newtonian VERTEX-PROMETHEUS code (AGILE and PROMETHEUS for short) has already been carried out by Liebendörfer et al. (2005) and Marek et al. (2006). Since
Liebendörfer et al. (2005) provide a useful framework for assessing the quality of a newly developed code, we have repeated their run G15 with CoCoNuT, and have compared our results to those obtained with AGILE. As the neutrino transport module in VERTEX-CoCoNuT is largely identical to that in VERTEX-PROMETHEUS (apart from the underlying moment equations), there is an additional payoff from this testing strategy: in a number of cases, the origin of the differences between the three codes can be pinpointed more accurately with a third available code. In addition to the PROMETHEUS run discussed in Liebendörfer et al. (2005), which was carried out using a potential derived from the TOV equation (“case R”), we also considered a PROMETHEUS run with a different gravitational potential (“case A”) proposed by Marek et al. (2006). In total, our analysis therefore encompasses four different simulations of model G15 with AGILE, CoCoNuT, and PROMETHEUS (with two different choices for the gravitational potential).

Following Liebendörfer et al. (2005) as closely as possible, we launch our simulation from the 15 $M_\odot$ progenitor s15s7b2 of Woosley & Weaver (1995), and use the same set of neutrino interactions rates and the same high-density EOS of Lattimer & Swesty (1991). Our treatment of the EOS differs slightly from both AGILE and PROMETHEUS for this test run: while using the EOS of Lattimer & Swesty (1991) above a threshold density of $6 \times 10^7$ g cm$^{-3}$, nuclear burning is switched off below $3 \times 10^7$ g cm$^{-3}$ (as in AGILE), and the composition of the progenitor is advected with the fluid (different from AGILE, which considers only one nucleus, $^{28}$Si, in that regime). Between those transition densities, the composition is taken from a 17-species NSE table. An Eulerian grid is used both for the hydrodynamics and the neutrino transport (400 and 234 zones, respectively), and the energy resolution is the same as in PROMETHEUS (17 zones).

Although the microphysical input is almost identical in all three codes, it is crucial to bear in mind that they differ considerably from each other in their approach to neutrino radiation hydrodynamics before proceeding with a comparison of their results: AGILE solves the relativistic Boltzmann equations directly by means of a discrete-angle ($\Sigma N$) method, and relies on an artificial viscosity scheme with an adaptive (moving) grid for the hydrodynamics. VERTEX-PROMETHEUS employs a variable Eddington factor technique for the neutrino transport and the PPM with an exact Riemann solver for evolving the equations of hydrodynamics; it uses a number of approximations to incorporate GR effects, while remaining essentially a Newtonian code. The key element of its GR treatment is a modification of the Newtonian gravitational potential that leads to the same hydrostatic stellar structure equation (TOV equation) as in general relativity (VERTEX-PROMETHEUS “case R”) or to a slightly modified TOV equation that improves the agreement with the relativistic case (“case A”). VERTEX-CoCoNuT, while based on the same general approach to neutrino transport and hydrodynamics as VERTEX-PROMETHEUS, accounts for GR effects exactly in spherical symmetry (apart from the computation of the Eddington factors). However, the approximate Riemann solver used in CoCoNuT (HILLE for this run) is somewhat less accurate than the exact solver in PROMETHEUS. Moreover, the two relativistic codes CoCoNuT and AGILE differ in their choice of gauge and slicing conditions: ambiguities arising from this fact must be handled carefully.

In most instances, the different gauges employed by CoCoNuT and AGILE do not pose any problem because the conversion of the gauge-dependent quantities is usually straightforward. Thus, the areal (circumferential) radius $r_{\text{cf}}$, which is used by Liebendörfer et al. (2005), can easily be computed from the isotropic radial coordinate $r$, which actually appears as $r$ in the CFC metric:

$$r_{\text{cf}} = \phi^2 r.$$  (62)

We also follow Liebendörfer et al. (2005) in choosing the rate of change of the areal radius per proper time as radial velocity $v$. However, the identification of time slices for the comparison of radial profiles of density, luminosity, etc. is non-trivial. In principle, a complete reconstruction of the entire spacetime would be necessary to allow for a rigorous comparison of the code output. Since the data provided in Liebendörfer et al. (2005) are insufficient for such a reconstruction, we are forced to identify time slices by the elapsed coordinate time $\tau$ (i.e., proper time for a non-moving observer at infinity). This is actually less critical than it might appear if we synchronize the different codes at bounce and use the post-bounce time $t_{\text{fb}}$ as time coordinate, because the system soon settles down to a quasi-stationary state for which the slicing conditions in AGILE and CoCoNuT are identical to very good approximation.

Fortunately, we can even form an estimate of the asynchronicity that develops between the CoCoNuT and AGILE runs, by considering the difference of the elapsed proper time $t_{\text{proper}}$ at the center of the proto-neutron star after a given interval $\delta t$ of coordinate time. Since the velocity and shift vector vanish at the center, $t_{\text{proper}}$ is given in terms of the central lapse function $\alpha_c$ by

$$t_{\text{proper}} = \int \alpha_c \, dt.$$  (63)

The elapsed proper time between the CoCoNuT and AGILE runs therefore differs by

$$\Delta t_{\text{proper}} = \int (\alpha_c^{\text{CoCoNuT}} - \alpha_c^{\text{AGILE}}) \, dt.$$  (64)

Taking the difference of the central lapse function at bounce (obtained by extrapolating to $r_{\text{cf}} = 0$) as an upper limit, we find that the two runs become asynchronous by at most another 2 ms every 100 ms at the center of the proto-neutron star. Since the $|\alpha_c^{\text{CoCoNuT}} - \alpha_c^{\text{AGILE}}|$ again becomes smaller after bounce, one may conclude that the two runs are never more than a few milliseconds out of sync until the end of our simulation 250 ms after bounce. Backed by these findings, we can undertake a detailed comparison of the simulations in CoCoNuT and AGILE (and, of course, also in PROMETHEUS) during the collapse, bounce, and post-bounce phases.

### 3.2.2. Collapse, Bounce, and Neutrino Burst Phase

Relativistic effects are of minor importance until the late stages of collapse when the density reaches several $10^{13}$ g cm$^{-3}$; only then do the infall velocity $v$ and the compactness parameter $M/R$ of the iron core reach values of the order of 0.1 (in

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5 The data from the AGILE and PROMETHEUS (case R) runs have been taken from the online material provided in the electronic version of Liebendörfer et al. (2005).

10 The details of the implementation differ, however: in particular, both VERTEX codes simplify the physics in the $\mu/r$-neutrino sector by not treating neutrinos and antineutrinos separately.

11 This procedure has been adopted to ensure that material in grid cells passing from the high-density to the low-density regime has a physically reasonable composition.

12 For a heuristic motivation of the modifications in case A, see Marek et al. (2006).
relativistic units). Consequently, the three codes are in excellent agreement during most of the collapse phase. Slight differences in the central electron and lepton fraction after trapping (i.e., at a density of a few $10^{12} \text{ g cm}^{-3}$) can be discerned in Figure 3. Since the VERTEX-based codes consistently produce somewhat lower values, it is likely that this difference stems from small differences in the treatment of the neutrino interaction rates. It should be noted that the new VERTEX-CoCoNuT code conserves the central lepton fraction and entropy to very good accuracy, thus passing another important test.

The good agreement between the different codes persists until shock formation (defined as the instant when the specific entropy inside the sonic point first reaches $3k_B/\rho$, which occurs at an enclosed mass of $0.53 M_\odot$ in AGILE, CoCoNuT, and PROMETHEUS (case A), and at a slightly smaller mass of $0.49 M_\odot$ in PROMETHEUS (case R). Interestingly, important special relativistic effects can be seen shortly after bounce (see Figure 4): within 3 ms, the shock has propagated to the mass shell $m \approx 1.0 M_\odot$. The velocity profile for the best effective potential (case A) of Marek et al. (2006) still matches the profiles from AGILE and CoCoNuT perfectly, but the entropy profile left behind by the shock differs appreciably. As long as the shock travels through an optically thick medium (to neutrinos), the post-shock entropy is higher by up to $0.5k_B$ in the relativistic case (AGILE, CoCoNuT), indicating that the shock is initially stronger. The post-shock velocity and the velocity jump across the shock indeed remain larger in CoCoNuT than in PROMETHEUS out until $m \approx 0.9 M_\odot$—velocity profiles from AGILE during the relevant phase are not available from Liebendörfer et al. (2005), unfortunately. Once the shock reaches $m \approx 0.75 M_\odot$, the combination of the relativistic jump conditions and the transition to the semi-transparent regime results in the formation of a shallow trough in the entropy profile, which is centered around $m \approx 0.9 M_\odot$. On the other hand, the specific entropy increases almost monotonically with $m$ out to the shock radius in that region in the PROMETHEUS run, and the post-shock entropy is even a little higher than in the GR runs.

Although a detailed analysis of the cause of these differences is rather complicated, it is at least possible to account for the higher post-shock entropies inside $m \approx 0.75 M_\odot$ in the relativistic case. In this region, the material is still optically thick, and the different jump conditions thus remain the only distinguishing factor from the Newtonian case. One can therefore estimate the relativistic correction to the post-shock entropy by comparing the Newtonian and relativistic form of the jump conditions and applying a perturbative analysis: the Newtonian Rankine–Hugoniot conditions and their relativistic analog (Taub 1948; Thorne 1973) can be recast into equations for the difference of the pre-shock and post-shock pressure ($P_1$ and $P_2$) and specific enthalpy ($h_1$ and $h_2$). In the relativistic case, $P_2 - P_1$ and $h_2 - h_1$ are given by

$$P_2 - P_1 = \frac{M^2}{h_1/\rho_1 - h_2/\rho_2},$$

and

$$h_2 - h_1 = \frac{M^2}{h_1 + h_2} \left[ \frac{(h_1)^2}{\rho_1} - \frac{(h_2)^2}{\rho_2} \right].$$

Here, $M$ is the baryonic mass flux through the discontinuity in the rest frame of the shock, which is given in terms of the three-velocities $\tilde{v}_1$ and $\tilde{v}_2$ and the Lorentz factors $\tilde{W}_1$ and $\tilde{W}_2$ on either side by

$$M = \rho_1 \tilde{W}_1 \tilde{v}_1 = \rho_1 \tilde{W}_2 \tilde{v}_2.$$  (67)

The three-velocities and Lorentz factors are measured in the rest frame of the shock (which is denoted by the tilde). The Newtonian limit of the jump conditions can then be found by setting $h_1 \to 1$, $h_2 \to 1$ in Equations (65) and (66) and $W_1 \to 1$, $W_2 \to 1$ in Equation (67). By linearizing Equations (65)–(67) around $h = 1$ and $W = 1$ and around a Newtonian solution for $h$ and $P$, one obtains an estimate for the relativistic post-shock enthalpy and pressure (as well as the other thermodynamic variables). For the post- and pre-shock conditions in the PROMETHEUS run, this analysis predicts post-shock entropies that are higher by up to $\Delta s = 0.4k_B$ in the relativistic case, mostly during a very brief period where the infall velocity $\tilde{v}_1$ is as large as $|\tilde{v}_1| = 0.34$ in the rest frame of shock. This prediction is in rough agreement with the actual simulation data. As the shock moves further out beyond $m = 0.75 M_\odot$, $|\tilde{v}_1|$ decreases and the relativistic correction to the post-shock entropy soon becomes negligible (e.g., $\Delta s \lesssim 0.03k_B$ for $|\tilde{v}_1| \lesssim 0.2$), resulting in a negative entropy gradient in the range $m \approx 0.8$–$0.9 M_\odot$. For $m \gtrsim 0.9 M_\odot$ the relativistic post-shock entropy then matches the Newtonian one very closely. Although neutrino losses from the post-shock region also start to play a role at this stage, this reduction of $\Delta s$ with decreasing $|\tilde{v}_1|$ thus provides a qualitative explanation for the observed trough in the entropy profile.

The effects of relativistic shock propagation are also visible in the density stratification between the proto-neutron star and the shock (Figure 4), which is more shallow in the region outside $r_{ct} \approx 25$ km ($m \approx 0.7$), i.e., just outside the region of the first entropy peak in the relativistic models. Such small differences may seem unimportant in the case of spherically symmetric problems—indeed particular when they are washed out after a few tens of milliseconds—but they can be of some relevance in multi-dimensional simulations, where they can affect the growth of hydrodynamic instabilities. Our one-dimensional simulations already give a strong indication that this is indeed the case: convective instability is expected if the Schwarzschild–Ledoux criterion $C_L$ (cf. Schwarzschild 1958) is positive. $C_L$ is given in terms of the density $\rho$, the pressure $P$, the specific internal energy density $\epsilon$, and the speed of sound $c_b$ by

$$C_L = \frac{\partial \rho (1 + \epsilon)}{\partial r_{ct}} \frac{1}{c_b^2} \frac{\partial P}{\partial r_{ct}},$$

in the relativistic case (Thorne 1966), or

$$C_L = \frac{\partial \rho}{\partial r_{ct}} \frac{1}{c_b^2} \frac{\partial P}{\partial r_{ct}},$$

in Newtonian hydrodynamics. The left panel of Figure 5 reveals a convectively unstable layer at $r_{ct} \approx 50$ km with a thickness of more than $10$ km at a time of $20$ ms after bounce in the CoCoNuT run, which is not present in the PROMETHEUS run.

Note that the Schwarzschild–Ledoux criterion reduces to the simple Schwarzschild criterion $C_s = ds/dr$ for a chemically homogeneous fluid. The alternative form $C_L = ds/dr (\partial(\rho + \epsilon)/\partial s)_{\rho,Y_{lep}} + dY_{lep}/dr (\partial(\rho + \epsilon)/\partial Y_{lep})$, $r$ (Buras et al. 2006b) can not be applied in the relativistic case. However, the relativistic Ledoux criterion can be expressed in terms of the spatial derivatives of $s$ and $Y_{lep}$ in the following manner,

$$C_L = ds/dr (\partial(\rho + \epsilon)/\partial s)_{\rho,Y_{lep}} + dY_{lep}/dr (\partial(\rho + \epsilon)/\partial Y_{lep}), r.$$
The stronger deleptonization in the region around \( m = 0.9 \, M_\odot \) also leads to a visibly different evolution of the electron neutrino luminosities during the neutronization burst (first panel of Figure 6). Although the peak luminosity \( L_{\text{burst}} \approx 3.8 \times 10^{53} \text{ erg s}^{-1} \) in PROMETHEUS (case A and R) agrees well with the one in AGILE, the radiated energy between \( t_{\text{pb}} = 1 \) ms and \( t_{\text{pb}} = 8 \) ms is smaller by around 10%. Conversely, the total energy radiated in \( \nu_e \)-s in CoCoNuT during the burst is in perfect agreement with AGILE despite the higher peak value of \( L_{\text{burst}} \approx 4.3 \times 10^{53} \text{ erg s}^{-1} \). The different shape of the burst in AGILE and CoCoNuT is probably due to the choice of the radial grid, or the different accuracy of the discretization scheme (first order versus second order in space), which both influence the numerical diffusivity of the transport scheme and thus lead to a different degree of broadening during the propagation of the neutrino burst from the decoupling radius to the observer radius at \( r_{\text{cf}} = 500 \) km (cf. Liebendörfer et al. 2004 for an analysis of the broadening of the burst). The treatment of the model Boltzmann equation in CoCoNuT may also be an important factor, either because of the approximations in the tangent-ray scheme (which may affect the transition from the optically thick to the optically thin regime) or because of the superior angular resolution of the tangent-ray scheme at large radii compared with the \( S_\omega \)-method used in AGILE, but these factors probably play a minor role.

### 3.2.3. Accretion Phase

The post-bounce evolution of the luminosities of all neutrino flavors (see also Figure 6) is again very similar in CoCoNuT and AGILE. During the first 150 ms after bounce, the luminosities in CoCoNuT tend to be slightly higher than in AGILE. The agreement is about as good as it is for PROMETHEUS (case A) and significantly better than it is for case R, where the strength of
the gravitational potential and hence the accretion luminosity is considerably overestimated. While PROMETHEUS (case A) is always very close to both CoCoNuT and AGILE, it is interesting to note that only the two last-named codes show a very abrupt transition from the fast rise in the \( \nu_\mu/\nu_\tau \) luminosity until \( t_{pb} = 20 \) ms to the subsequent plateau phase. The small differences between CoCoNuT and AGILE are well within the range expected for different codes (cf. the Newtonian run N13 in Liebendörfer et al. 2005) or even for one code at different resolutions (cf. Marek 2007). As in PROMETHEUS, there is a noticeable drop in the luminosities of all flavors (\( \nu_e \) and \( \bar{\nu}_e \) in particular) between \( t_{pb} = 150 \) ms and \( t_{pb} = 200 \) ms—i.e., at a time, when the shock reaches the oxygen-rich silicon shell—which is far less pronounced in AGILE due to the different treatment of the nuclear composition in the low-density regime and the possible superiority of the Riemann solver methods in PROMETHEUS and CoCoNuT in following the composition discontinuity at the Si–SiO interface.

There is also good agreement between the three codes concerning the spectral properties of the emitted neutrinos (see Figure 7). The root mean square (rms) energies of electron neutrino and antineutrinos as measured by an observer at \( r = 500 \) km in the comoving frame. Electron neutrino luminosities are shown in the upper panels, electron antineutrino and \( \mu/\tau \) neutrino luminosities in the lower left and lower right panels.
that the excellent matching of $v_{\nu}/v_{r}$ rms energies between case A and AGILE is due to a cancellation of the error introduced by the approximate treatment of general relativity (leading to lower rms energies) and differences in the implementation of $v_{\mu}/v_{r}$ interaction rates or the transport schemes (leading to higher rms energies). Conceivably, the higher peak value in CoCoNuT may originate from the different neutrino transport algorithms, finite-difference schemes, and numerical grids used in VERTEX and AGILE, which we already suggested as reason for the slightly higher peak luminosity of the neutrino burst.

Similar luminosities and spectral properties already suggest similar neutrino loss rates in the cooling zone and similar heating rates in the gain region; therefore the dynamical evolution of the proto-neutron star and the accretion shock in CoCoNuT and AGILE should not be very different either. The right panel of Figure 5 shows that this is indeed the case: until $t_{pb} = 150$ ms the accretion shock in CoCoNuT is very close to the one in AGILE; the radial deviation corresponds to only one or two grid zones. Interestingly, there is a brief transient period of shock stagnation at $t_{pb} \approx 7$ ms (lasting a few milliseconds) both in AGILE and CoCoNuT, which is absent or at least much less pronounced in PROMETHEUS (cases A and R).

It is conceivable that this feature is connected to the different relativistic and Newtonian form of the jump conditions, and to the slightly higher energy loss in electron neutrinos during the burst in the GR simulations. In all simulations, the maximum shock radius of $r_{cf,max} \approx 150$ km is reached at $t_{pb} = 60–80$ ms. Afterward, the shock position in CoCoNuT lies somewhere in between the two PROMETHEUS runs until $t_{pb} \approx 150$ ms, and is a little closer to AGILE than either of them. The CoCoNuT and PROMETHEUS runs then show a transient phase of shock expansion when the infalling Si–SiO interface reaches the accretion front; this feature is absent in AGILE (as explained before) because of the different nuclear composition of low-density material. At later times ($t_{pb} > 220$ ms), the shock in CoCoNuT again recedes to the same radius as the shock in AGILE. The overall shock trajectory in PROMETHEUS (case A, but not for case R) is very similar to the one in CoCoNuT, which indicates that the effects of general relativity can still be adequately captured by an effective gravitational potential in one-dimensional simulations.

The very close overall agreement between CoCoNuT, AGILE, and PROMETHEUS (case A) and the even better agreement between CoCoNuT and AGILE is also evident in Figures 8 and 9, which show radial profiles of selected physical quantities at 100 ms and 150 ms after bounce. There are only minute differences in the velocity profiles produced by the three codes; the shock front seems to be resolved a little more sharply in CoCoNuT, but this is probably incidental, since the sharpness of the jump changes slightly as the shock slowly moves from zone to zone. The effects of genuinely relativistic shock propagation in the early post-bounce phase, which had resulted in a somewhat different density stratification in the post-shock region in the GR case, have now been largely washed out. This can also be seen in the entropy profiles, which are very similar during this phase, except for the fact that PROMETHEUS (case A) seems to give slightly lower entropies in the gain region. Interestingly, we now find the electron fraction behind the shock to be somewhat lower in PROMETHEUS (case A), contrary to the situation near shock breakout. The luminosity and neutrino rms energy profiles show some visible, but not worrisome differences between the three codes; e.g., the Doppler jump at the shock is generally a little larger in CoCoNuT than in AGILE, which is to be expected, however, since there are small differences in the pre- and post-shock velocities. Considering that the three codes use different radial grids (and in the case of AGILE also a different grid in energy space), such small discrepancies seem unavoidable and are probably not to be ascribed to coding errors in any of them.

3.2.4. Long-time Tests

Unfortunately a direct verification of our new code beyond 250 ms by means of a comparison with AGILE-BOLTZTRAN is not easily possible, because data for model G15 are not provided by Liebendörfer et al. (2005) beyond that point. However, the mere demonstration that VERTEX-CoCoNuT can stably evolve supernova models considerably further would be valuable in its own right, and might also add some insights concerning the reliability of the effective potential approach used in PROMETHEUS at very late times. We have therefore also repeated one of the longest-running one-dimensional simulations carried out with VERTEX-PROMETHEUS, viz., a model (“G8.8”) of the explosion of the 8.8 $M_\odot$ progenitor of Nomoto (1984, 1987) with an O–Ne–Mg core. To this end, we have included the effect of electron captures on $^{20}$Ne and $^{24}$Mg (Takahara et al. 1989) and the same approximate burning

Figure 7. Left: neutrino rms energies for model G15 obtained with AGILE-BOLTZTRAN (solid), VERTEX-CoCoNuT (dashed), and VERTEX-PROMETHEUS (dotted and dash-dotted for the effective potentials R and A), sampled at $r = 500$ km in the comoving frame. The three panels show the rms energies for electron neutrinos, electron antineutrinos, and $\mu/\tau$ neutrinos (from left to right).
Figure 8. Radial profiles of the velocity $v$ (dashed lines), the density $\rho$ (solid lines, upper left), the electron fraction $Y_e$ (solid lines), the entropy $s$ per nucleon (dashed lines, bottom left), the neutrino luminosities $L_{\nu}$ (upper right), and the neutrino rms energies $\langle \varepsilon \rangle_{\text{rms}}$ (bottom right) in AGILE-BOLTZTRAN (black), VERTEX-CoCoNuT (green), and VERTEX-PROMETHEUS with the effective potential $A$ (blue) at a post-bounce time of 100 ms. Neutrino luminosities and energies are given in the comoving frame; solid lines are used for electron neutrinos, dashed lines for electron antineutrinos, and dash-dotted lines for $\mu/\tau$ neutrinos.

Figure 9. Radial profiles of the velocity $v$ (dashed lines), the density $\rho$ (solid lines, upper left), the electron fraction $Y_e$ (solid lines), the entropy $s$ per nucleon (dashed lines, bottom left), the neutrino luminosities $L_{\nu}$ (upper right), and the neutrino rms energies $\langle \varepsilon \rangle_{\text{rms}}$ (bottom right) in AGILE-BOLTZTRAN (black), VERTEX-CoCoNuT (green), and VERTEX-PROMETHEUS with the effective potential $A$ (blue) at a post-bounce time of 150 ms. Neutrino luminosities and energies are given in the comoving frame; solid lines are used for electron neutrinos, dashed lines for electron antineutrinos, and dash-dotted lines for $\mu/\tau$ neutrinos.

treatment as in Kitaura et al. (2006) in the CoCoNuT code. Different from the G15 run and from published long-time simulations of the same progenitor model (Fischer et al. 2009; Hüdepohl et al. 2010), we now use a new nuclear EOS kindly provided by J. Lattimer (2007, private communication) in the high-density regime; a more extensive discussion of this new EOS in the context of supernova simulations will be given in L. Hüdepohl et al. (2010, in preparation).

As for the more massive progenitor used in the G15 run, CoCoNuT and PROMETHEUS give very similar results throughout collapse and during the short accretion phase of model G8.8, and we therefore refrain from repeating the same detailed analysis of these stages as for model G15. Instead, we avail ourselves of the opportunity to compare a full GR treatment and the effective potential approach during the post-explosion phase. Due to the structure of the progenitor model, the accretion rate drops rapidly as early as 70–80 ms after bounce when the shock reaches the edge of the core, and neutrino heating becomes effective in unbinding the post-shock material, resulting in a weak ($\approx 10^{50}$ erg) explosion (cf. Kitaura et al. 2006; Janka et al. 2008). Interestingly, an important difference between the CoCoNuT and PROMETHEUS runs emerges at this point: as the shock propagates through the extremely steep density gradient at the edge of the core, it accelerates considerably until it reaches the dilute hydrogen envelope of the progenitor whose density in the innermost $10^7$ km is of the order of only...
10^{-8} – 10^{-7} \text{ g cm}^{-3}. In the PROMETHEUS run, the maximum post-shock velocity $v_{\text{ps}}$ becomes superluminal with a maximum value of $v_{\text{ps}} = 5.3 \, c$, which clearly implies that the propagation of the shock through the outer layers of the core and the hydrogen envelope cannot be captured correctly with a Newtonian code. In the CoCoNuT run, all velocities correctly remain subluminal with a maximum value of $v_{\text{ps}} = 0.9848 \, c$ (corresponding to a Lorentz factor of $W = 5.76$). However, the shock becomes relativistic only at densities below $\approx 10 \, \text{ g cm}^{-3}$, and at that point the post-shock temperatures are already far too low to allow for nuclear burning. The results of Janka et al. (2008) about the nucleosynthesis conditions behind the shock in O–Ne–Mg supernovae therefore remain valid.

Apart from the different propagation of the shock in the CoCoNuT and PROMETHEUS runs, the subsequent evolution of the newly formed remnant in model G8.8 is remarkably similar in both cases. The contraction of the proto-neutron star due to the continuing neutrino losses is almost in perfect agreement as illustrated by the left panel of Figure 10, which shows the evolution of the central density $\rho_c$. Even more than 4 s after bounce, $\rho_c$ is only $\approx 3\%$ higher in the PROMETHEUS run, and the neutron star radii (not shown) are almost identical as well, with a maximum deviation of $2\%$. Slightly larger differences are observed in the neutrino luminosities (see Figure 11), which are typically a little higher in the pseudo-Newtonian PROMETHEUS run from 0.5 to 2 s after bounce, but the difference to the relativistic simulation stays low ($\approx 5\%$).

The mean energies of the emitted neutrinos also agree to within $3\%$ during the post-explosion phase, and it is only after more than 2 s that the deviation starts to become larger, reaching a maximum of 0.4 MeV for the $\mu$/r neutrinos. Considering that the finest energy grid zone has a width of 1.3 MeV, this close agreement seems quite remarkable. However, despite the fact that the neutrino luminosities and mean energies, as well as the structure of the proto-neutron star remain extremely similar, there is a somewhat bigger discrepancy between the two codes concerning the properties of the neutrino-driven wind (see the right panel of Figure 10). While the electron fraction $Y_e$ in the wind differs by less than 0.003 at late times, the entropy $s$ reached asymptotically in the wind is roughly 7\% lower in the PROMETHEUS run. Since $s$ scales roughly with a small negative power of the luminosity ($s \propto L^{-1/6}$, see Qian & Woosley 1996), it seems unlikely that this is a result of the slightly stronger emission of neutrinos in the CoCoNuT run. Furthermore, we find the wind entropy to be lower in PROMETHEUS already at early times when the luminosities are still almost in perfect agreement with CoCoNuT. Even if the small differences in the mean neutrino energies of the order of less than 3\% are also taken into account, the different wind entropies still cannot be explained on the basis of the neutrino emission alone: assuming that the dependence of the final entropy on the mean energy of electron antineutrinos $\langle s \rangle_{\nu_e}$ is roughly given by $s \propto \langle s \rangle_{\nu_e}^{-1/3}$ (see again Qian & Woosley 1996), the combined effect of different neutrino energies and luminosities could only account for a relative difference of $\lesssim 2\%$ in $s$ in the most optimistic case. This indicates that the Newtonian treatment of the equations of hydrodynamics in PROMETHEUS is probably the cause for the observed discrepancy: although the effective potential used in PROMETHEUS reproduces the structure of static neutron stars very accurately, it is obviously somewhat less adequate for stationary (but non-static) flows. However, with a deviation of the specific entropy of the wind of less than 10\%, the best effective potential (case A) of Marek et al. (2006) can still be considered quite reliable even for the wind phase.

These results are very reassuring in two different respects: the VERTEX-CoCoNuT code has proved a reliable and robust tool even for long-time supernova simulations well into the proto-neutron star cooling phase. In addition, the effective potential approach, though unable to reproduce all the details of the relativistic G8.8 run exactly, has been shown to be very accurate in a regime where it has hitherto remained untested.

### 3.2.5. Energy and Lepton Number Conservation

Our test runs also provide an opportunity to check the ability of VERTEX-CoCoNuT to conserve total energy and lepton number. The usefulness of our improved numerical treatment of the relativistic energy equation presented in Appendix A and the finite-difference scheme for redshift and Doppler terms in the neutrino moment equation described in Appendix B can thus be tested beyond the level of simplified test problems. It should be noted that the very fact that the new code allows us to quantify the numerical violation of energy conservation unambiguously is a major advantage compared to the pseudo-relativistic approach of VERTEX-PROMETHEUS: for the effective potentials of Marek et al. (2006) and Müller et al. (2008), a conservation for the total energy cannot even be formulated. On the other hand, in the framework of the $3 + 1$ formalism there is a conservation law for the ADM energy (Landau & Lifschitz 1997; Misner et al. 1973), which is given by (cf. Equation (10))

$$E_{\text{ADM}} = \int_0^\infty \int_0^{\pi} \int_0^{2\pi} \phi^5 \left( E + \frac{K_{ij}K^{ij}}{16\pi} \right) r^2 \sin \theta \, d\phi \, d\theta \, dr,$$

(70)

in the case of a CFC spacetime. The accuracy of total energy conservation can therefore be checked by monitoring $E_{\text{ADM}}$ (taking into account surface flux terms where necessary) at least in spherical symmetry, where the CFC does not involve any approximations to the Einstein equations. In the multi-dimensional case, the CFC system (10–12) is not equivalent to the full Einstein equations, and $E_{\text{ADM}}$ is therefore not conserved in general. Even though, $E_{\text{ADM}}$ can still be computed and remains a valuable diagnostic quantity for two reasons: first, $E_{\text{ADM}}$ should still be conserved to good accuracy as long as the deviations from the conformal flatness remain sufficiently small, which is probably the case for the generic supernova problem (cf. Section 2.4). Second, whether the non-conservation of $E_{\text{ADM}}$ is indicative of a numerical problem with energy conservation or of some failure of the CFC approximation is actually of minor importance: an overly large deviation $\Delta E_{\text{ADM}}$ from the initial value would be a sign of trouble either way.

However, in the one-dimensional case, $E_{\text{ADM}}$ ought to remain constant (neglecting surface fluxes), and $\Delta E_{\text{ADM}}$ can be directly used to measure the numerical energy conservation error. The time evolution of $\Delta E_{\text{ADM}}$ is shown in Figure 12: after a very rapid increase by $0.5 \times 10^{51}$ erg around bounce, $\Delta E_{\text{ADM}}$ gradually grows by another $1.5 \times 10^{51}$ erg during the next 500 ms after bounce.\footnote{Interestingly, the evolution of $\Delta E_{\text{ADM}}$ is somewhat different from the relativistic test case in Appendix A, where $\Delta E_{\text{ADM}}$ becomes negative at bounce, and afterward settles down at around $-1 \times 10^{50}$ erg. Imperfect energy conservation in the neutrino sector as described in this section may partly account for this. However, the different dynamical evolution of the GR model in Appendix A (where neutrino energy losses are neglected completely) may also be of relevance.} In two-dimensional simulations to $\approx 300$ ms.
after bounce, we also achieve about the same accuracy. The significance of these numbers is not immediately clear without further analysis. On the one hand, the fact that $\Delta E_{\text{ADM}}$ is of the order of the typical explosion energy may seem worrisome, on the other hand $\Delta E_{\text{ADM}}$ is small compared to the gravitational binding energy. Fortunately, the proper scale of reference for $\Delta E_{\text{ADM}}$ can be worked out by determining the cause of the violation of energy conservation more precisely. As shown in Appendix A for the purely hydrodynamic case, $|\Delta E_{\text{ADM}}|$ should not exceed a few $10^{50}$ erg during the bounce phase for our improved implementation of the gravitational source term in the energy Equation (7), and the additional error accumulated after the formation of the proto-neutron star is even smaller. This suggests that the secular increase of $\Delta E_{\text{ADM}}$ during the accretion phase occurs in the neutrino transport sector. To verify this, we specialize to the case of a static background spacetime, where the relativistic evolution equations for the component $T_{00}$ of the stress–energy tensor $T_{\mu\nu}$ of the matter (or the neutrinos) can be written in a fully conservative form by means of the following manipulations:

\[
\frac{1}{\sqrt{-g}} \frac{\partial}{\partial t} (\sqrt{-g} T^{00}) + \frac{1}{\sqrt{-g}} \frac{\partial}{\partial x^j} (\sqrt{-g} T^{0j}) + T_{\nu\lambda}^{0} T^{\nu\lambda} = Q, \tag{71}
\]

\[
\frac{\partial}{\partial t} (\sqrt{-g} T^{00}) + \frac{\partial}{\partial x^j} (\sqrt{-g} T^{0j}) + \sqrt{-g} g^{00} \frac{\partial}{\partial x^i} (g_{00}) T^{0i} = \sqrt{\gamma} Q, \tag{72}
\]

\[
\frac{\partial}{\partial t} (\alpha \sqrt{-g} T^{00}) + \frac{\partial}{\partial x^j} (\alpha \sqrt{-g} T^{0j}) + \sqrt{-g} \frac{\partial}{\partial x^i} (g_{00}) T^{0i} = \alpha \sqrt{\gamma} Q. \tag{73}
\]

Here, $T^{\mu\nu}$ may denote either the stress–energy tensor of the matter or of the neutrinos, and $Q$ is the source term due
to neutrino–matter interactions, which enters with different signs in both cases \( (\mathcal{Q}_{\text{matter}} = -\mathcal{Q}_{\text{neutrino}}) \). The time-like components \( T_{\alpha \nu}^{\text{tot}} \) of the total stress–energy tensor therefore obey the following conservation law:

\[
\frac{\partial}{\partial t} \left( \alpha \sqrt{-g} T_{\alpha \nu}^{(0)} \right) + \frac{\partial}{\partial x^\nu} \left( \alpha \sqrt{-g} T_{\alpha \nu}^{(0)} \right) = 0. \quad (75)
\]

Since our implementation of the gravitational source term in the energy Equation (7) is essentially equivalent to a conservative discretization of Equation (74), any spurious energy loss or generation in the case of a static background spacetime must originate from our finite-difference representation of the neutrino moment equations. This can be verified by checking whether the integral version of Equation (74) for the neutrinos is fulfilled numerically. More specifically, we consider the rate of change of the total neutrino energy inside a given radius \( R \), which ought to be given by

\[
\frac{\partial}{\partial t} \int_0^R \alpha \sqrt{-g} T^{(0)} \, dr = -\int_0^R \alpha \sqrt{-g} \gamma Q \, dr. \quad (76)
\]

In other words, the rate of change of the total neutrino energy \( \int_0^R \alpha \sqrt{-g} T^{(0)} \, dr \) inside \( R \) is given by a surface flux term \(-\int_0^R \alpha \sqrt{-g} T^{(0)} \, \partial_\nu \gamma \) and a volume-integrated source term \( \int_0^R \alpha \sqrt{-g} \gamma Q \, dr \). By plotting the left-hand side and right-hand side of Equation (76) as functions of the radius \( R \) as in Figure 12, one may directly determine where energy conservation is actually violated: ideally, the numerically evaluated left-hand side and right-hand side (solid and dashed lines in Figure 12) should be identical. Inside the shock at \( R \approx 100 \, \text{km} \), the agreement is indeed very good, the maximum deviation being \( 5 \times 10^{-5} \, \text{erg} \, \text{s}^{-1} \). Numerical energy conservation is violated more strongly at the shock and in the pre-shock region, and the accumulated error amounts to approximately \( 2 \times 10^{-5} \, \text{erg} \, \text{s}^{-1} \), which corresponds nicely to the rate of change of the ADM energy at \( t_{\text{sh}} = 230 \, \text{ms} \).

The underlying reason for the numerical violation of neutrino energy conservation can also be worked out easily: when deriving Equation (74) from the frequency-integrated moment equations in the comoving frame on a static spacetime background, all the Doppler and gravitational redshift terms are canceled by corresponding terms arising from the Lorentz transformation to the lab frame and Equation (74). For the finite-differenced moment equations, these cancellations are no longer exact. Further analysis shows that the most critical terms in this respect are the velocity-dependent terms of \( \mathcal{O}(v) \), which explains why energy conservation is violated at the shock and the pre-shock region, where the absolute velocities are highest. In this region, however, neutrino heating and cooling are negligible (except for the early post-bounce phase, where there may be some pre-heating), and the dynamically relevant error in the heating rate resulting from energy non-conservation in the neutrino sector is therefore much smaller. Thus, the dynamical evolution of the accreting proto-neutron star and its environment should not be significantly affected by the secular increase of the ADM energy. Furthermore, since Equation (76) is violated mainly in the optically thin region, it is quite natural to interpret the numerical increase of \( E_{\text{ADM}} \) of the order of a few \( 10^{51} \, \text{erg} \, \text{s}^{-1} \) as a luminosity error of a few percent and not as spurious energy input. Considering that the uncertainties in the microphysics and the discretization error probably limit the accuracy of the neutrino luminosities to a similar level anyway, the conservation error observed in simulation G15 does not seem very disturbing.

Moreover, the problem with the finite-difference representation of the Doppler terms in the moment equations becomes less acute in the post-explosion phase once the shock has left the grid: in model G8.8, the spurious change of the ADM energy between \( t_{\text{sh}} = 1 \, \text{s} \) and \( t_{\text{sh}} = 4 \, \text{s} \) is only \( 3.7 \times 10^{51} \, \text{erg} \). However, since the velocities in the neutrino-driven wind can still reach a few percent of the speed of light, the energy error still does not vanish completely.

Somewhat surprisingly, lepton number conservation is also not fulfilled exactly in models G15 and G8.8; there is a secular drift of the order of \( 1\%-2\% \) of the net lepton number luminosity. At first sight, this seems puzzling, since the treatment of Doppler and gravitational redshift terms described in Appendix B allows us to solve the neutrino energy Equation (7) in such a way that lepton number is conserved numerically. However, this apparent contradiction can be easily resolved: in order to simplify the structure of the Jacobian of the discretized moment equations, we compute the flux terms \( \partial/\partial \epsilon (\cdots) \) using the flux factor \( f_{\text{H}} \) from the solution of the model Boltzmann equation and the zeroth moment of the radiation intensity \( J \), while the first moment \( H \) is used directly in the corresponding terms outside the derivative \( \partial/\partial \epsilon \). If \( f_{\text{H}} J \neq H \), the cancellation of these terms described in Appendix B would still work, and lepton number conservation would be guaranteed; but we find \( f_{\text{H}} J \neq H \) for two reasons: first, our Boltzmann equation is not fully consistent with the moment Equations (27) and (28) in the sense that some terms of the underlying Boltzmann Equation (26) are omitted, and hence the flux factors computed from the solutions of the model Boltzmann and moment equations do not agree perfectly, particularly in the vicinity of the shock. Second, \( f_{\text{H}} J \neq H \) are defined on cell centers and cell interfaces, respectively, so that the finite interpolation accuracy will, in general, lead to \( f_{\text{H}} J \neq H \). It should be pointed out, however, that such problems are far less severe in the optically thick regime, because there \( H \approx 0 \). These arguments suggest that the observed imperfect conservation of lepton number conservation is again essentially equivalent to a small error in the neutrino (number) luminosity.

4. CONCLUSIONS

We have presented a new method for GR neutrino radiation hydrodynamics in core-collapse supernovae that combines multi-group neutrino transport, an elaborate treatment of the microphysics, multi-dimensional GR hydrodynamics, and the conformal flatness (CFC) approximation for the gravitational field equations. Our approach is a generalization of the variable Eddington factor technique as implemented by Rampp & Janka (2002), and relies on the ray-by-ray-plus approach (Buras et al. 2006b) for treating the multi-dimensional case. As in the Newtonian approximation (Rampp & Janka 2002; Buras et al. 2006b), the equations of neutrino transport in spherical symmetry are solved on different angular bins ("rays") independently, assuming a radial neutrino flux and a parametric dependence of the matter and spacetime background on altitude. However, terms for lateral neutrino advection and compressional heating of the neutrino gas are also included in the correct relativistic form, as is the acceleration of the matter due to lateral neutrino pressure gradients. Although we have considered only the axisymmetric case in this paper, the method can be generalized to three dimensions in a straightforward manner.

Our approach involves a number of approximations to the full six-dimensional relativistic neutrino transport problem on different levels. While the variable Eddington factor technique...
is well tested in spherical symmetry in the context of core-collapse supernovae (Rampp & Janka 2002; Liebendörfer et al. 2004), no direct comparisons of the ray-by-ray-plus method with multi-angle transport in two dimensions or three dimensions are available as yet. However, there is at least qualitative agreement with multi-angle transport (see Ott et al. 2008 for results obtained with such an approach and Marek & Janka 2009 for a discussion of similarities/dissimilarities to ray-by-ray transport), even in situations where alternative approaches like multi-dimensional flux-limited diffusion suffer from excessive lateral smearing of the radiation field. Although ray-by-ray transport may not be the method of choice for extremely aspherical configurations (such as accretion tori around black holes), it is probably adequate for the typical core-collapse scenario where the proto-neutron star rotates only at moderate speeds and is therefore not overly aspherical. As far as the GR field equations are concerned, the introduction of CFC is well justified, since it is an excellent approximation even for rapidly spinning compact objects for which the ray-by-ray approach would break down. We also note that the relativistic ray-by-ray method could easily be generalized to the case where the Einstein equations are solved in spherical polar coordinates without any approximations.

We have implemented our relativistic ray-by-ray-plus method using the framework of two existing codes, the time-explicit relativistic hydrodynamics code CoCoNuT (Dimmelmeier et al. 2002a) and the implicit neutrino transport code VERTEX (Rampp & Janka 2002; Buras et al. 2006b). Re-utilizing most of the routines of VERTEX allows us to include the same up-to-date and well-tested treatment of the neutrino microphysics as in the recent pseudo-Newtonian simulations of our group (Buras et al. 2006b, 2006a; Kitaura et al. 2006; Marek et al. 2009; Marek & Janka 2009; Hüdepohl et al. 2010), which greatly facilitates comparisons between the Newtonian and relativistic cases.

A few improvements were made to the CoCoNuT code, such as the implementation of the approximate Riemann solver HLLC (Mignone & Bodo 2005). More significantly, an improved scheme for the energy equation was developed, which greatly reduces the numerical violation of total energy conservation for self-gravitating systems. The new scheme has also been added to the most recent version of the PROMETHEUS code, and may be of use for other grid-based Eulerian finite-volume codes as well. We also devised a more efficient and consistent treatment for the advection of neutrinos in energy space in the VERTEX module. The new method eliminates the need to solve both the neutrino energy and number equations to maintain energy and lepton number conservation.

In this paper, we have limited ourselves to code tests in spherical symmetry before moving on to the first multi-dimensional applications of VERTEX-CoCoNuT. While the analytically tractable radiation transport problems presented in Section 3.1 merely serve as a consistency check for the implementation of the neutrino moment equations, we also conducted a full simulation (“G15 run”) of the collapse, bounce, and post-bounce phases for the progenitor model s15x7b2 of Woosley & Weaver (1995), and compared our results to the relativistic one-dimensional Boltzmann solver AGILE-BOLTZTRAN (Liebendörfer et al. 2004), building on the detailed comparative study of Liebendörfer et al. (2005). Since our code is largely an adapted version of the Newtonian VERTEX-PROMETHEUS code of Rampp & Janka (2002), this also allows us to better assess the reliability of the effective potential approach for GR effects (Rampp & Janka 2002; Marek et al. 2006; now also used, e.g., by Messer et al. 2007). For this reason, we have also included PROMETHEUS runs with two different gravitational potentials (case R and case A) in our comparison. Moreover, we performed long-time proto-neutron star cooling simulations of the post-explosion remnant of the 8.8 M⊙ progenitor model of Nomoto (1984, 1987) with VERTEX-CoCoNuT and PROMETHEUS (“G8.8 run”).

The results of our tests may be summarized as follows: on the one hand, VERTEX-CoCoNuT is in very good agreement with analytic solutions to simple test problems and also with AGILE in the context of a fully fledged one-dimensional simulation from core collapse well into the accretion phase. The agreement with AGILE is slightly better than for the PROMETHEUS code with the best gravitational potential currently available. On the other hand, our simulations of the 8.8 M⊙ model of Nomoto (1984, 1987) show that PROMETHEUS still produces results extremely similar to CoCoNuT well beyond the point up to which a direct comparison to the full GR case was hitherto available in the work of Liebendörfer et al. (2005).

The conclusions that can be drawn from these tests are twofold. First, the VERTEX-CoCoNuT emerges as a reliable neutrino transport code. Its accuracy has been demonstrated by a thorough comparison of results with a well-established relativistic one-dimensional neutrino transport code. The G8.8 run, covering more than 4 s of the proto-neutron star cooling phase, also shows it to be sufficiently robust for long-time simulations. For an explicit hydro code with an implicit transport solver, the
ability to stably evolve a quasi-stationary proto-neutron star over thousands of dynamical timescales and to maintain total energy conservation adequately over such long evolution times cannot be taken for granted, but our simulations with CoCoNuT prove that these requirements can be met with an up-to-date higher-order Eulerian scheme for GR hydrodynamics. Naturally, the accuracy and robustness of CoCoNuT in the one-dimensional case will also serve as a point of reference for future multi-dimensional numerical studies of core-collapse supernovae in general relativity.

The very good agreement of the results of CoCoNuT and PROMETHEUS, both for the G15 and G8.8 run, is also noteworthy because it provides further evidence that, at least in one dimension, the evolution of proto-neutron stars can be adequately captured by approximating GR effects in a pseudo-Newtonian framework by means of an effective gravitational potential. The usefulness and accuracy of the best effective potential of Marek et al. (2006) has also now been established for the proto-neutron star cooling phase, and for much longer evolution times and significantly more compact neutron stars than in Liebendörfer et al. (2005) and Marek et al. (2006). Despite the smaller differences to the GR case seen during the early post-bounce phase, this also gives further credence to two-dimensional simulations using effective potentials (Buras et al. 2006b, 2006a; Marek et al. 2009; Marek & Janka 2009; Bruenn et al. 2006; Messer et al. 2007).

However, with the advent of VERTEX-CoCoNuT, the choice between rather crude approximations for GR effects or a drastic simplification of the microphysics in multi-dimensional supernova studies is now obsolete. The role of GR effects in the multi-dimensional case can now be investigated directly with up-to-date microphysics and a sophisticated neutrino transport scheme. There are several prominent topics in supernova physics where these new capabilities of our code may prove highly valuable, and which we plan to address in follow-up papers. In particular, we intend to reinvestigate the dynamics of rotational core collapse and the development of multi-dimensional hydrodynamic instabilities during the post-bounce phase. The determination of accurate gravitational wave signals—which is connected to these issues—is obviously another important goal.

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**APPENDIX A**

**ENERGY CONSERVATION IN NEWTONIAN AND RELATIVISTIC HYDRODYNAMICS**

A.1. Importance of Energy Conservation in the Supernova Problem

Conservative discretization schemes for the Euler equations of gas dynamics are one of the cornerstones in modern computational astrophysics, both because they have appealing mathematical properties (LeVeque 1998), and because they conserve important integral quantities (total rest mass, total energy, etc.) to machine precision. Unfortunately, the standard formulation of the Euler equations for self-gravitating fluids is not strictly conservative despite the fact that there exists an integral conservation law for the total kinetic, internal, and potential energy for any localized matter distribution.

In the supernova problem, the non-conservative formulation of gravitational source terms in the Euler equations raises some critical questions: Can a secular drift of the total energy on the order of $10^{51}$ erg over a few 100 ms (which is a typical value for many simulations; see Liebendörfer et al. 2004; Burrows et al. 2006a; Murphy & Burrows 2008) have a qualitative influence on the propagation of the shock and the development of the explosion, as the explosion energy is typically of the same order? Is there an appreciable effect in the first few seconds of the neutron star cooling phase? The former question may probably be answered in the negative, because the accumulated error in the gravitational energy liberated during the collapse of the iron core and the contraction of the proto-neutron star makes itself felt primarily as a small excess or deficit (on the percent level or below) in the internal energy of the compact central object (which contains most of the mass on the grid), and not as a large relative error in the region around the shock. On the other hand, a secular drift of the total energy can qualitatively affect neutron star cooling in the post-explosion phase: once spurious generation of energy due to numerical inaccuracies outweighs the slowly decreasing neutrino losses, the proto-neutron star does not cool any longer (such an effect was actually observed in one of our pseudo-Newtonian simulations).

In order to alleviate the problems associated with the non-conservative standard formulation of gravitational source terms in the energy equation, we have developed an alternative scheme for treating the energy equation that significantly decreases the secular drift of the total energy. The new scheme can easily be integrated into any Eulerian finite-volume method, and is applicable (with only small variations) both in Newtonian and relativistic hydrodynamics. In the Newtonian case, the new scheme also has clear advantages when working with effective gravitational potentials that do not obey the Poisson equation, despite the fact that a conservation law for the total energy cannot be formulated in this case: using a modified gravitational potential, it turns out that the energy released during the formation of the neutron star is in good agreement with the binding energy computed from the TOV equation, at least for neutron stars of moderate compactness (cf. Hüdepohl et al. 2010), while systematic energy conservation errors originating from the discretization of the gravitational source term may qualitatively change the outcome of numerical simulations.

A.2. Energy Equation in Newtonian Hydrodynamics

A.2.1. Description of the Algorithm

Usually the energy equation for a self-gravitating fluid is written as

$$\frac{\partial e}{\partial t} + \frac{\partial (e + P) v^i}{\partial x^i} = -\rho v^i \frac{\partial \Phi}{\partial x^i}. \quad (A1)$$

Here, $e = \rho v^2/2 + \rho \epsilon$ is the total energy density, $P$ is the pressure, $v_i$ is the fluid velocity, $\rho$ is the (baryonic) mass density, $\epsilon$ is the...
specific internal energy, and \( \Phi \) is the Newtonian gravitational potential. Although Equation (A1) is not written in the form of a pure conservation equation, there is a conservation law for the volume integral of the total (i.e., kinetic, internal, and potential) energy (see, e.g., Shu 1992): in order to obtain this integral conservation law, we absorb the source term \( -\rho \dot{v}^i \partial \Phi / \partial x^i \) into the flux term on the left-hand side,

\[
\frac{\partial e}{\partial t} + \frac{\partial (e + P + \rho \Phi)}{\partial x^i} v^i = \Phi \frac{\partial \rho v^i}{\partial t}, \tag{A2}
\]

and use the continuity equation \( \partial \rho / \partial t + \nabla \cdot (\rho \mathbf{v}) = 0 \) to eliminate the divergence of the mass flux in the new source term,

\[
\frac{\partial e}{\partial t} + \frac{\partial (e + P + \rho \Phi)}{\partial x^i} v^i = -\Phi \frac{\partial \rho}{\partial t}. \tag{A3}
\]

Integrating Equation (A3) over the entire spatial domain then yields

\[
\int \frac{\partial e}{\partial t} dV = - \int \Phi \frac{\partial \rho}{\partial t} dV = \frac{1}{4\pi G} \int \Phi \Delta \Phi dV
\]

\[
= - \frac{1}{4\pi G} \int \nabla \Phi \nabla \Phi dV
\]

\[
= - \frac{1}{8\pi G} \frac{\partial}{\partial t} \int \nabla \Phi \cdot \nabla \Phi dV
\]

\[
= \frac{1}{8\pi G} \frac{\partial}{\partial t} \int \Phi \Delta \Phi dV = -\frac{1}{2} \frac{\partial}{\partial t} \int \rho \Phi dV. \tag{A4}
\]

Here, we have used Green’s first identity, the Poisson equation \( \Delta \Phi = 4\pi G \rho \), and the fact that the surface integral,

\[
\int \Phi \nabla \Phi dA,
\]

vanishes if the domain of integration is extended to infinity. Equation (A4) implies that the total energy is conserved,

\[
\int \left( e + \frac{1}{2} \rho \Phi \right) dV = \text{const.} \tag{A6}
\]

However, the \textit{discretized} representation of the total energy

\[
E_{\text{tot}} = \sum_{i,j,k} \left( e_{i,j,k} + \frac{1}{2} \rho_{i,j,k} \Phi_{i,j,k} \right) \Delta V_{i,j,k}, \tag{A7}
\]

will not be automatically conserved, if the numerical implementation of the gravitational source term is based on Equation (A1), because the discretized source term cannot be manipulated in the same fashion as in the steps from Equation (A1)–(A3). One way to eliminate any secular drift of the total energy resulting from this, and to avoid the unwanted consequences described in Section A.1, would be to use a strictly conservative form of the energy equation (Noonan 1984)

\[
\frac{\partial}{\partial t} \left[ e + \rho \Phi + \frac{1}{8\pi G} \left( \frac{\partial \Phi}{\partial x^i} \right)^2 \right] + \frac{\partial}{\partial x^i} \left[ (e + P + \rho \Phi) v^i + \frac{1}{4\pi G} \Phi \frac{\partial^2 \Phi}{\partial t \partial x^i} \right] = 0. \tag{A8}
\]

Unfortunately, this form of the energy equation has several drawbacks: the “gravitational self-energy” contribution \( 1/(8\pi G) (\partial \Phi / \partial x^i)^2 \) to the total energy density can exceed the kinetic, internal, and binding energy of the matter by a large factor, resulting in considerable round-off errors in these quantities. Furthermore, the flux term contains a mixed spatial and temporal derivative, which may introduce numerical inaccuracies. It is also impossible to formulate an analog of Equation (A8) for the effective relativistic potentials commonly used in VERTEX-PROMETHEUS (Marek et al. 2006; Müller et al. 2008) because these do not obey the Poisson equation.

However, there is a simple and efficient solution to the problem of numerical energy conservation that avoids these shortcomings and requires only a minimal modification of the “standard” scheme based on Equation (A1). The starting point is the reformulation of the energy equation in Equation (A2): using the continuity equation, the divergence of \( \rho \mathbf{v} \) can be replaced by the time derivative of the density, which is in turn partially absorbed in the time derivative of \( \Phi \).

\[
\frac{\partial}{\partial t} \left( e + \rho \Phi \right) + \frac{\partial (e + P + \rho \Phi)}{\partial x^i} v^i = \rho \frac{\partial \Phi}{\partial t}. \tag{A9}
\]

The remaining source term now contains the time derivative of the gravitational potential. Equation (A9) can be conveniently solved with an operator-split scheme in three steps.

\textit{Step 1.} First, the equations of hydrodynamics are solved with the additional potential energy terms \( \rho \Phi \), where \( \Phi \) itself is kept constant,

\[
\frac{\partial (e + P + \rho \Phi)}{\partial t} + \frac{\partial (e + P + \rho \Phi)}{\partial x^i} v^i = 0, \quad \Phi = \text{const.} \tag{A10}
\]

Since \( \rho \Phi \) may be considerably larger than the internal energy density \( e \), using \( e + \rho \Phi \) as conserved quantity instead of \( e \) may lead to large round-off errors in \( e \) and produce unwanted numerical noise (in particular spurious entropy loss or production). To circumvent this problem, the time derivative \( \partial \rho \Phi / \partial t \) can be re-expressed in terms of the mass flux (as \( \Phi = \text{const.} \), i.e., \( \partial \rho \Phi / \partial t = (\Phi \partial \rho / \partial t) = -\Phi \nabla \cdot (\rho \mathbf{v}) \)),

\[
\frac{\partial e}{\partial t} + \frac{\partial (e + P + \rho \Phi)}{\partial x^i} v^i = \frac{\partial \rho v^i}{\partial x^i} \Phi = 0, \quad \Phi = \text{const.} \tag{A11}
\]

The value \( e^{(n+1/3)} \) of the energy density after the first fractional step can be obtained by integrating Equation (A11) over the cell volume \( V \). It is instructive to compare the final formula for the update of \( e \) with “standard” implementations of the gravitational source term. Working with forward differences in time,\(^{16}\) we now have

\[
e^{(n+1)} = e^{(n)} - \frac{\Delta t}{V} \int \left[ \int_{\partial V} (e + P) \mathbf{v} \cdot dA + \sum_k \int_{A_k} (\Phi_c - \Phi_k) \rho \mathbf{v} \cdot dA_k \right] \tag{A12}
\]

(where \( A_k \) denotes the \( k \)th cell interface, while \( \Phi_c \) and \( \Phi_k \) are the cell-center and cell-interface values of \( \Phi \)) instead of

\[
e^{(n+1/3)} = e^{(n)} - \frac{\Delta t}{V} \left[ \int_{\partial V} (e + P) \mathbf{v} \cdot dA \int \nabla \Phi \cdot \rho \mathbf{v} dV \right] \tag{A13}
\]

\(^{16}\)First-order forward differences are only adopted for illustration at this point. The implementation in higher-order time integration schemes is straightforward.
for the standard formulation. Since \( \int (\Phi_k - \Phi_i) \, dA_k / V \) is the component of \( \nabla \Phi \) pointing in the direction of \( dA_k \), the update prescription (A12) can be viewed as an alternative version of Equation (A13), in which the cell-averaged mass flux \( \rho \nu \) has been replaced by an average of the mass flux at cell interfaces. Thus, gravitational energy is only released or stored when mass is actually exchanged between different cells. This is the key ingredient for eliminating a secular drift of the total energy.

When the gravitational field is stationary, no further fractional steps are required, and the value of \( e \) after one full time step is just \( e^{(n+1)} = e^{(n+1/3)} \). If the gravitational potential varies with time, two additional steps are necessary once \( \Phi \) has been updated from its old value \( \Phi^{(n)} \) to \( \Phi^{(n+1)} \) at the end of the time step.

Step 2. In order to ensure that Equation (A10) is fulfilled after the update of the gravitational potential \( \Phi \), the energy density \( e^{(n+2/3)} \) after the second fractional step must be calculated according to

\[
e^{(n+2/3)} = e^{(n+1/3)} + \frac{1}{2} (\rho^{(n)} + \rho^{(n+1)}) (\Phi^{(n+1)} - \Phi^{(n)}).
\]

(A14)

Step 3. Finally, the source term \( \rho \partial \Phi / \partial t \) has to be accounted for in a third step,

\[
\frac{\partial (e + \rho \Phi)}{\partial t} = \frac{\partial \Phi}{\partial t}.
\]

(A15)

If centered differences in time are used to obtain second-order accuracy, the energy density at the end of the time step reads

\[
e^{(n+1)} = e^{(n+2/3)} + \frac{1}{2} (\rho^{(n)} - \rho^{(n+1)}) (\Phi^{(n+1)} - \Phi^{(n)}).
\]

(A16)

Steps 2 and 3 can easily be combined into a compact formula for the update of \( e \) after each recalculation of the gravitational potential:

\[
e^{(n+1)} = e^{(n+1/3)} + \frac{1}{2} (\rho^{(n)} + \rho^{(n+1)}) (\Phi^{(n+1)} - \Phi^{(n)}).
\]

(A17)

### A.2.2. Condition for the Exact Conservation of Energy

Under certain conditions, the scheme described in Section A.2.1 conserves the total energy exactly. To demonstrate this for the special case of a spherical polar grid, we consider the total energy balance after each sub-step of the algorithm: Steps 1 and 2 conserve the (discretized) total volume integral of \( e + \rho \Phi \), provided that the flux \( (e + P + \rho \Phi) \nu \) vanishes at the boundaries of the computational domain,

\[
\sum_{i,j,k} \left( e_{i,j,k}^{(n+1)} + \rho_{i,j,k}^{(n)} \Phi_{i,j,k}^{(n)} \right) \Delta V_{i,j,k} = \sum_{i,j,k} \left( e_{i,j,k}^{(n)} + \rho_{i,j,k}^{(n)} \Phi_{i,j,k}^{(n)} \right) \Delta V_{i,j,k}.
\]

(A18)

If \( e \) is then updated according to Equation (A16), we obtain, by combination with Equation (A18),

\[
\sum_{i,j,k} \left( e_{i,j,k}^{(n+1)} + \rho_{i,j,k}^{(n+1)} \Phi_{i,j,k}^{(n+1)} \right) \Delta V_{i,j,k} = \sum_{i,j,k} \left[ e_{i,j,k}^{(n)} + \frac{1}{2} \rho_{i,j,k}^{(n)} \Phi_{i,j,k}^{(n)} - \rho_{i,j,k}^{(n+1)} \Phi_{i,j,k}^{(n)} + \rho_{i,j,k}^{(n+1)} \Phi_{i,j,k}^{(n+1)} \right] \Delta V_{i,j,k}.
\]

(A19)

Hence, the total energy after the \((n + 1)\)th time step is according to the definition in Equation (A7),

\[
E_{\text{tot}}^{(n+1)} = E_{\text{tot}}^{(n)} + \sum_{i,j,k} \rho_{i,j,k}^{(n)} \Phi_{i,j,k}^{(n+1)} \Delta V_{i,j,k} - \sum_{i,j,k} \rho_{i,j,k}^{(n+1)} \Phi_{i,j,k}^{(n)} \Delta V_{i,j,k}.
\]

(A20)

The sums on the right-hand side quantify the violation of total energy conservation during one time step. If they were replaced by integrals, we could manipulate them using Green's first identity and the Poisson equation \( \nabla \cdot \nabla \Phi = 4\pi G \rho \) to obtain an exact cancellation:

\[
\int \rho^{(n)} \Phi^{(n)} \, dV - \int \rho^{(n+1)} \Phi^{(n+1)} \, dV = \frac{1}{4\pi G} \int \nabla \Phi^{(n)} \cdot (\nabla \Phi^{(n+1)}) \, dV - \frac{1}{4\pi G} \int (\nabla \Phi^{(n+1)}) \cdot (\nabla \Phi^{(n)}) \, dV = 0.
\]

(A21)

Using a special finite-difference representation of the Poisson equation, we can manipulate the sums in Equation (A20) in very much the same way as by applying Green's first identity. Let us assume that the source density \( 4\pi G \rho \) is expressed as the divergence of the gravitational acceleration \( \vec{g} \) in the following natural manner,

\[
4\pi G \rho \Delta A_{i,j,k} = g_{i+1/2,j,k} \Delta A_{i,j,k} - g_{i-1/2,j,k} \Delta A_{i,j,k} + g_{i,j+1/2,k} \Delta A_{i,j,k} - g_{i,j-1/2,k} \Delta A_{i,j,k} + g_{i,j,k+1/2} \Delta A_{i,j,k} - g_{i,j,k-1/2} \Delta A_{i,j,k}.
\]

(A22)

Here, \( \Delta A_{i+1/2,j,k} \) denotes the area of the interface between the cells \((i, j, k)\) and \((i + 1, j, k)\), and \( g_{i+1/2,j,k} \) denotes the perpendicular component of \( \vec{g} \) thereon; \( \Delta A_{i,j+1/2,k} \), etc. are defined analogously. Let us further suppose that the values of \( \vec{g} \) on any cell interface are obtained by multiplying the difference of \( \Phi \) in the adjacent cells by a time-independent factor \( \zeta_{i,j,k} \), as in the following natural finite-difference representation:

\[
g_{i+1/2,j,k} = \zeta_{i+1/2,j,k} \Phi_{i+1,j,k} - \Phi_{i+1,j,k} = 1 - \frac{1}{r_{i+1} - r_i} \langle \Phi_{i,j,k} - \Phi_{i+1,j,k} \rangle,
\]

(A23)

\[
g_{i,j+1/2,k} = \zeta_{i,j+1/2,k} \Phi_{i,j+1,k} - \Phi_{i,j+1,k} = \frac{1}{r_i(r_{i+1} - r_i)} \langle \Phi_{i,j,k} - \Phi_{i,j+1,k} \rangle,
\]

(A24)

\[
g_{i,j,k+1/2} = \zeta_{i,j,k+1/2} \Phi_{i,j,k} - \Phi_{i,j,k} = \frac{1}{r_i \sin \theta_j(r_{i+1} - r_i)} \langle \Phi_{i,j,k} - \Phi_{i,j,k+1} \rangle.
\]

(A25)
At the outer boundary, we may set $g_{(M+1/2,j,k)} = \zeta_{M+1/2,j,k} \Phi_{(M,j,k)} =: \Phi_{(M,j,k) + 1/2}$ to obtain the correct monopole moment of the gravitational field.

The first sum term in the energy balance Equation (A20) can now be written as

$$\sum_{i,j,k} \rho_{(i,j,k)} \Phi_{(i,j,k)+1} V_{i,j,k} = \frac{1}{4\pi G} \sum_{i,j,k} \left( g_{(i+1/2,j,k)} \Delta V_{i,j,k} - g_{(i-1/2,j,k)} \Delta V_{i,j,k} \right) \Phi_{(i,j,k)}^{(n+1)}$$

(A26)

The sums on the right-hand side of Equation (A26) can be rearranged via summation by parts using the finite-difference representation for $g$ in Equations (A23)–(A25). For the sum containing the radial components $g_{(i+1/2,j,k)}$ of the gravitational acceleration $g$ on cell interfaces, we have (bearing in mind that $A_{(1/2,j,k)} = 0$, as $i = 1/2$ corresponds to the origin)

$$\sum_{i,j,k} \left( g_{(i+1/2,j,k)} \Delta A_{(i+1/2,j,k)} - g_{(i-1/2,j,k)} \Delta A_{(i-1/2,j,k)} \right) \Phi_{(i,j,k)}^{(n+1)}$$

(A28)

In effect, we have obtained a finite-difference analog of Green’s first identity with non-vanishing surface contributions. The second sum in Equation (A20) can be rearranged in precisely the same way and thus cancels the first sum exactly. Hence, the total internal, kinetic, and gravitational energy is conserved:

$$E^{(n+1)} = E^{(n)}.$$  (A29)

A.3. Energy Equation in General Relativity

In GR, the situation is somewhat different from the Newtonian case: at first sight, the fact that no sources appear in the local conservation law $\nabla_{\mu} T^{\mu \nu} = 0$ for the stress–energy tensor $T^{\mu \nu}$ might suggest that the energy and momentum equations can be formulated in a strictly flux-conservative form (i.e., without sources). However, an integral conservation law can only be formulated (via Gauss’ theorem) for divergence-free vector fields on a differentiable and orientable manifold (Wald 1984; Straumann 2004); such fields can in general be constructed from the energy–momentum tensor only if there exists a Killing vector field.18 On the other hand, a conservation law can be formulated for the sum of $T^{\mu \nu}$ and the Landau–Lifshitz pseudo-tensor $\tau^{\mu \nu}$ of the gravitational field (Landau & Lifschitz 1997; Straumann 2004)

$$\frac{\partial \left(-g\right) T^{\mu \nu} + \tau^{\mu \nu}}{\partial x^\nu} = 0.$$  (A30)

Unfortunately, the practical use of Equation (A30) is limited because $t$ is a complicated (and non-unique) function of the metric and its derivatives. Furthermore, it is impossible to interpret $t^{\mu \nu}$ as the local energy–momentum vector density of the gravitational field due to its non-uniqueness and non-tensorial character; only the integral $E_{ADM} = \int \left( T^{\mu \nu} + \tau^{\mu \nu} \right) dV$ (known as ADM energy) over the entire spatial domain in an asymptotically flat spacetime has a well-defined physical meaning. In general, Equation (A30) will only be useful for special gauge choices and symmetry assumptions, where the resulting source-free energy

17 Note that in spherical polar coordinates further “surface contributions” vanish since $A_{(1/2,j,k)} = A_{(N+1/2,j,k)} = 0$, $\Phi_{(0,j,0)} = \Phi_{(j,j,O)}$, and $\Phi_{(j,j,O+1)} = \Phi_{(j,j)}$.

18 Special cases comprise (1) the flat Minkowski spacetime with ten Killing fields (implying the conservation of energy, momentum, and angular momentum), (2) stationary spacetimes with a time-like Killing field (implying energy conservation), and (3) axisymmetric spacetimes with a space-like Killing field that creates an $SO(2)$ isometry group (implying the conservation of one angular momentum component).
equation is not overly complicated (see e.g., Romero et al. 1996; Liebendörfer et al. 2001a). The total volume integral $E_{\text{ADM}}$ (which can also be expressed in different ways) is, of course, an important diagnostic quantity for assessing the quality of energy conservation in a numerical code.

Because of these obstacles we construct an improved scheme for the energy equation in a similar way as in the Newtonian case and do not attempt to eliminate source terms from the energy equation altogether. We start from the formulation used by Dimmelmeier et al. (2008),

$$\frac{\partial \sqrt{\gamma} \tau}{\partial t} + \frac{\partial}{\partial x^i} \left( \sqrt{-g} (\tau \dot{v}^i + P v^i) \right) = -\sqrt{-g} [\tau D \dot{v}^i + P v^i] + \frac{\partial}{\partial x^i} \left( 2 K_{ij} \beta^j - \frac{\partial \alpha}{\partial x^i} \right) + T^{ij} K_{ij}, \quad \text{(A31)}$$

For convenience, we add the continuity equation, and subsume the first source term on the right-hand side as in the Newtonian potential. This source term can be eliminated by multiplying Equation (A32) by $\alpha$,

$$\alpha \frac{\partial \sqrt{\gamma} (\tau + D)}{\partial t} + \frac{\partial}{\partial x^i} \left( \sqrt{-g} (\tau \dot{v}^i + D \dot{v}^i + P v^i) \right) = -\alpha \sqrt{-g} [(\tau + D) \dot{v}^i + P v^i] \frac{\partial \alpha}{\partial x^i} + \alpha Q, \quad \text{(A33)}$$

and by pushing the lapse function into the space and time derivatives,

$$\frac{\partial \sqrt{\gamma} \alpha (\tau + D)}{\partial t} + \frac{\partial}{\partial x^i} \left( \sqrt{-g} \alpha (\tau \dot{v}^i + D \dot{v}^i + P v^i) \right) = \sqrt{\gamma} (\tau + D) \frac{\partial \alpha}{\partial t} + \alpha Q, \quad \text{(A34)}$$

Again, the new source term $(\tau + D) \partial \alpha / \partial t$ corresponds closely to $\rho \partial \Phi / \partial t$ in the Newtonian case. Since it is numerically advantageous to separate the baryonic mass contributions to the total energy (due to the reduction of round-off errors), we subtract the continuity equation, and finally arrive at the following alternative energy equation:

$$\frac{\partial}{\partial t} [\sqrt{\gamma} \alpha (\tau + D) - \sqrt{\gamma} D] + \frac{\partial}{\partial x^i} \left[ \sqrt{-g} (\alpha \tau \dot{v}^i + D \dot{v}^i + \alpha P v^i) \right] = \sqrt{\gamma} (\tau + D) \frac{\partial \alpha}{\partial t} + \alpha Q. \quad \text{(A35)}$$

The numerical solution within a finite-volume scheme can again be split into three sub-steps, the first of which consists in solving the energy equation (including only $\alpha Q$ as source term) while keeping the lapse function $\alpha$ fixed,

$$\frac{\partial}{\partial t} \left[ \sqrt{\gamma} \alpha \tau + \sqrt{\gamma} (\alpha - 1) D \right] + \frac{\partial}{\partial x^i} \left[ \sqrt{-g} (\alpha \tau \dot{v}^i + D \dot{v}^i + \alpha P v^i) \right] = \alpha Q. \quad \text{(A36)}$$

In order to avoid working with $\sqrt{\gamma} \alpha \tau + \sqrt{\gamma} (\alpha - 1) D$ as conserved quantity, we note that the change in $\sqrt{\gamma} (\alpha - 1) D$ inside a cell volume $V$ during the time step $\Delta t$ can be expressed in terms of the integral of the flux $D \dot{v}^i$ over the cell boundary $A = \partial V$. The finite-volume version of the energy equation for this cell can therefore be written as

$$\int_V \alpha \sqrt{\gamma} \tau^{(n+1/3)} dV = \int_V \alpha \sqrt{\gamma} \tau^{(n)} dV + \Delta t \int_V Q dV - \int_V \sqrt{\gamma} \alpha (\alpha - 1) D \dot{v}^i dA + \int_V \left[ \sum_k \sqrt{\gamma} \alpha_k (\alpha_c - \alpha_k) dA \right]. \quad \text{(A37)}$$

The surface integrals in the second line of Equation (A37) can be combined into a single term containing the difference between the lapse function $\alpha_c$ at the cell center and the lapse function $\alpha_k$ at the $k$th cell interface:

$$\int_V \sqrt{\gamma} \tau^{(n+1/3)} dV = \int_V \sqrt{\gamma} \tau^{(n)} dV + \Delta t \int_V Q dV - \frac{\Delta t}{\alpha} \left[ \int_V \sqrt{\gamma} \alpha^2 (\tau \dot{v}^i + P v^i) dA + \int_V \alpha Q dV \right] + \sum_k \int_{A_k} \sqrt{\gamma} \alpha_k (\alpha_c - \alpha_k) dA. \quad \text{(A38)}$$

This formulation requires only a minimal modification of the scheme for the unmodified energy Equation (A31).

In a second step, the lapse function is updated, while the conserved quantity $\sqrt{\gamma} [\alpha \tau + (\alpha - 1) D]$ is kept constant,

$$\frac{\partial}{\partial t} \left[ \alpha \dot{\tau} + (\alpha - 1) \dot{D} \right] = 0. \quad \text{(A39)}$$

Note that it is convenient to work with the densitized quantities $\hat{\tau} = \sqrt{\gamma} \tau$ and $\hat{D} = \sqrt{\gamma} D$ here because $\sqrt{\gamma}$ may be updated at the same time as $\alpha$. The discretized version of Equation (A39),

$$\left( \alpha^{(n+1)} \right)^{\hat{\tau}} + (\alpha^{(n+1)} - 1) \hat{D}^{(n+1)} = \alpha^{(n)} \hat{\tau}^{(n+1/3)} + (\alpha^{(n)} - 1) \hat{D}^{(n+1/3)}$$

has the solution

$$\sqrt{\gamma} \hat{\tau}^{(n+2/3)} = \sqrt{\gamma} \hat{\tau}^{(n+1/3)} \left( \alpha^{(n)} \hat{\tau}^{(n+1/3)} + (\alpha^{(n)} - 1) \hat{D}^{(n+1/3)} \right). \quad \text{(A41)}$$

In a third operator-split step, the source term containing $\sqrt{\gamma} (\tau + D) \partial \alpha / \partial t$ in Equation (A35) is taken care of

$$\frac{\partial}{\partial t} \left[ \sqrt{\gamma} \alpha \tau + \sqrt{\gamma} (\alpha - 1) D \right] = \sqrt{\gamma} (\tau + D) \frac{\partial \alpha}{\partial t}. \quad \text{(A42)}$$
As in the Newtonian case, the second and third steps can be merged, and if centered differences in time are used, \( \tau \) has to be updated according to

\[
\bar{\tau}^{(n+1)} = \frac{3}{2} \left( \bar{\tau}^{(n+1/3)} + \bar{\tau}^{(n)} + \frac{D^{(n+1)} + D^{(n)}}{2} \frac{\sigma^{(n)} - \sigma^{(n-1)}}{\alpha^{(n)}} \right).
\]

### A.4. Numerical Tests

To illustrate the virtues of the new scheme for the energy equation, we have conducted several numerical tests in spherical symmetry with different codes. Newtonian core-collapse simulations have been carried out with the Newtonian version of CoCoNuT and with a self-written second-order accurate hydrodynamics code employing total variation diminishing (TVD) reconstruction and the Kurganov–Tadmor central scheme (Kurganov & Tadmor 2000).

For the CoCoNuT run, we employ the deleptonization scheme of Liebendörfer (2005), and take neutrino energy losses into account in the total energy budget, i.e., we subtract the (negative) time- and volume-integrated neutrino energy source term from the total energy measured at any given time. The left upper panel of Figure 13 shows that the new implementation still produces a residual violation of energy conservation of about 2 \times 10^{39} \text{ erg} around bounce. However, the total energy changes by less than 1.7 \times 10^{40} \text{ erg} after \( t_{pb} = 10 \text{ ms} \), which is a considerable improvement compared to the old implementation, where the accumulated error exceeds the absolute value of the initial binding energy (\( \approx 5.8 \times 10^{50} \text{ erg} \)) within a few tens of milliseconds after bounce. Nonetheless, the reduction of the numerical error is of little relevance dynamically: the protoneutron settles down at a slightly higher central density, which is consistent with the smaller total energy obtained with the new scheme. The total kinetic energy in the post-shock region (right panels in Figure 13) is essentially identical in both cases as well.

We also note that by evolving \( e \) instead of \( e + \rho \phi \), round-off errors in the internal energy and the specific entropy can indeed be kept small; we did not find any unphysical entropy production during collapse.

Although a change of the total energy by around 0.3% as in the CoCoNuT simulation is probably more than satisfactory for core-collapse simulations, this level of accuracy does not exhaust the full potential of our new scheme, which can in principle conserve the total energy exactly, as explained in Section A.2.2. In order to verify this, we have simulated the collapse of a 4/3-polytrope with our self-written TVD test code using the hybrid EOS of Janka et al. (1993). Different from CoCoNuT, the gravitational potential and acceleration are calculated according to Equations (A22)–(A25), hence the condition for exact energy conservation is satisfied. Figure 14 (left panel) shows that the total energy changes by less than 10^{39} \text{ erg}, which corresponds to a relative error of less than 10^{-12} in terms of the initial energy (\( E_{tot,init} = -1.08 \times 10^{53} \text{ erg} \)) and less than 10^{-14} in terms of the internal energy at the end of the simulation. Thus our improved scheme can indeed preserve the total energy almost to machine precision provided that the gravitational potential is determined appropriately.

In addition to these tests in Newtonian gravity, we have also simulated the collapse of model s20.0 in GR using CoCoNuT. We employ the Liebendörfer scheme for deleptonization with one minor modification: while reducing \( Y_e \) according to a \( Y_e(\rho) \)-parameterization, we neglect energy losses due to neutrino emission, because this simplifies the total energy budget considerably. Ideally, we would expect the ADM energy \( E_{ADM} \), which

19 Setting the neutrino energy source term to zero while retaining a source term for \( Y_e \) is of course unphysical, but, as far as our test problem is concerned, the prescription adopted here is only employed as a means of initiating the collapse. A non-zero neutrino energy source term would complicate the energy budget considerably: any update of the internal energy of the fluid (or, equivalently, the entropy) would not only change the terms \( h \) and \( P \) in the integrand in Equation (A44) locally, but would also affect the conformal factor \( \phi \) and the extrinsic curvature \( K_{ij} \) non-locally. On the other hand, this problem does not arise with full neutrino transport, because the neutrino source terms do not change the total (matter and neutrino) energy--momentum tensor \( T^{\text{tot}}_{\mu\nu} \) in that case, and therefore do not affect the gravitational field.
can be written as
\[
E_{\text{ADM}} = -2\pi \int \left( \phi^5 \rho W^2 - P + \frac{K_{ij} K^{ij}}{16\pi} \right) dV
\] (A44)
in a CFC spacetime, to be conserved in such a situation. Figure 14 (right panel) shows that neither the “standard” scheme nor our new scheme manage to conserve \(E_{\text{ADM}}\) perfectly. In both cases, conservation is violated by \(|\Delta E_{\text{ADM}}| \approx 6 \times 10^{50}\) erg around bounce, which is not surprising considering the complicated nonlinear nature of the field equation and the appearance of gravitational self-energy terms in the sources and in the ADM energy. However, \(|\Delta E_{\text{ADM}}|\) settles down to a value of less than \(10^{50}\) erg within 1 ms after bounce in the simulation using our new scheme, and the secular drift is only \(10^{50}\) erg for the next 100 ms afterward. The old scheme, on the other hand, leads to a violation of around \(9 \times 10^{50}\) erg and a small but clearly visible drift of \(E_{\text{ADM}}\) during the post-bounce phase.

The relativistic test thus confirms our findings in Newtonian gravity: our reformulation considerably improves numerical energy conservation in simulations of self-gravitating systems. The secular drift during the quasi-stationary proto-neutron star evolution is not completely eliminated in all cases, but can be easily reduced by an order of magnitude, which demonstrates the usefulness of our new scheme particularly for long-time simulations covering many dynamical timescales.

**APPENDIX B**

**SIMULTANEOUS CONSERVATION OF ENERGY AND LEPTON NUMBER**

While the monochromatic moment Equations (27), (28) and (30), (31) for neutrino energy and number transport are analytically equivalent, a flux-conservative discretization of either pair of equations does not necessarily guarantee neutrino number (or, in the presence of neutrino–matter interactions, lepton number) and energy conservation simultaneously. The reason for this problem, which is inherent to spectral neutrino transport in the comoving frame, can be understood by comparing the analytic and discretized forms of the Doppler and gravitational redshift terms (denoted by a subscript “D”) in the moment equations. In the neutrino energy equation these terms can be written as
\[
\left( \frac{\partial J}{\partial t} \right)_D = \frac{\partial w J}{\partial \varepsilon} - w J,
\] (B1)
where the advection “velocity” in energy space \(w\) is a function of the velocities, metric functions, and the Eddington factors \(f_H = H/J, f_K = K/J, f_L = L/J\). To obtain the corresponding term in the equation for the monochromatic neutrino number density \(J = J/\varepsilon\), we divide by \(\varepsilon\) and apply the product rule,
\[
\left( \frac{\partial J}{\partial t} \right)_D = \frac{\partial w J}{\partial \varepsilon} + \frac{w J}{\varepsilon} \frac{\partial w J}{\partial \varepsilon} = \frac{\partial w J}{\partial \varepsilon} + \frac{\partial w J}{\partial \varepsilon}.
\]
The Doppler term in the neutrino number equation is thus a pure flux derivative term and cancels out if the equation is integrated over \(\varepsilon\) to obtain the evolution equation for the total neutrino number density. On the other hand, if the energy equation is discretized with backward differences in time and a flux-conservative form for the energy advection term,
\[
\left( J^{n+1}_i \varepsilon_i - J^n_i \varepsilon_i \right)_D = \frac{\varepsilon_i^{n+1/2} w_i^{n+1/2} J^{n+1}_i}{\varepsilon_i} - \frac{\varepsilon_i^{n-1/2} w_i^{n-1/2} J^{n-1}_i}{\varepsilon_i} - \frac{\varepsilon_i^{n+1/2} w_i^{n+1/2} J^{n+1}_i}{\varepsilon_i} + \frac{\varepsilon_i^{n-1/2} w_i^{n-1/2} J^{n-1}_i}{\varepsilon_i},
\] (B2)
the discretized Doppler term for \(J\) reads
\[
\left( \frac{\partial J}{\partial t} \right)_D = \frac{\partial w J}{\partial \varepsilon} + \frac{w J}{\varepsilon} \frac{\partial w J}{\partial \varepsilon}
\]
\[
\left( J^{n+1}_i \varepsilon_i - J^n_i \varepsilon_i \right)_D = \frac{\varepsilon_i^{n+1/2} w_i^{n+1/2} J^{n+1}_i}{\varepsilon_i} - \frac{\varepsilon_i^{n-1/2} w_i^{n-1/2} J^{n-1}_i}{\varepsilon_i} - \frac{\varepsilon_i^{n+1/2} w_i^{n+1/2} J^{n+1}_i}{\varepsilon_i} + \frac{\varepsilon_i^{n-1/2} w_i^{n-1/2} J^{n-1}_i}{\varepsilon_i},
\] (B3)
Here, \(\varepsilon_i\) and \(\varepsilon_i^{n+1/2}\) denote the values of the energy at cell centers and interfaces, respectively, and \(\Delta \varepsilon_i = \varepsilon_i^{n+1/2} - \varepsilon_i^{n-1/2}\) is the width of the \(n\)th zone in energy space. The source term in the neutrino number equation thus does not vanish in the discretized form of the equations, unless the last three terms in Equation (B3) cancel. Different methods have been proposed to recover the property of neutrino number conservation (Bruenn 1985; Liebendörfer et al. 2004). In the original version of VERTEX (Rampp & Janka 2002) the problem is overcome by solving both the monochromatic moment equations for the neutrino energy density and flux \((J, H)\), and the number
density and flux \((J, \mathcal{H})\) for electron neutrinos and antineutrinos *simultaneously*, but there are significant drawbacks to this approach: first, the mean energy \(J / \mathcal{J}\) within a given energy zone \([\varepsilon_{i-1/2}, \varepsilon_{i+1/2}]\) is not constrained to the cell-center value \(\varepsilon_i\) and may even move beyond the cell boundaries if \(\mathcal{J}\) is treated as an independent quantity. Moreover, the solution of the equations by the Newton–Raphson method involves the inversion of square matrix blocks of size \(4 \times N_\nu + 2\) instead of \(2 \times N_\nu + 2\) for neutrino energy transport alone, which makes the Newton–Raphson iteration almost eight times more expensive.

Fortunately, the Doppler terms in the moment equations for neutrino energy transport can be discretized in such a way as to guarantee neutrino energy and number conservation *simultaneously*. For example, in a rather simple approach the way as to guarantee neutrino energy and number conservation for neutrino energy transport can be discretized in such a Newton–Raphson iteration almost eight times more expensive.

\[
\varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right) = \frac{\varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right)}{\varepsilon_i \varepsilon} = \frac{\varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right)}{\varepsilon_i \varepsilon} - \frac{\varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right)}{\varepsilon_i \varepsilon}
\]

so that the last three terms in Equation (B3) would indeed cancel. However, since this scheme allows large violations of the monotonicity condition \((J_i < J_{i+1/2} < J_{i+1} \text{ or } J_i > J_{i+1/2} > J_{i+1})\), the resulting algorithm would be rather unstable.

To construct a stable scheme, we need to impose a weaker condition on the interface fluxes \(F = \varepsilon w \mathcal{J}\), i.e., we only require the source term in the frequency-integrated neutrino number equation to vanish:

\[
\frac{\varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right)}{\varepsilon_i \varepsilon} = \sum_i \left( \frac{\varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right)}{\varepsilon_i \varepsilon} \right) = 0.
\]

In order to find interface fluxes \(F_{i+1/2} = \varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right)\) that fulfill this condition, we first rearrange the sum in Equation (B5) in the following manner:

\[
\frac{\varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right)}{\varepsilon_i \varepsilon} = \frac{\varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right)}{\varepsilon_i \varepsilon} - \frac{\varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right)}{\varepsilon_i \varepsilon} = \sum_i \left( \frac{\varepsilon_i \varepsilon \left( J_{i+1/2} - J_{i+1/2} \right)}{\varepsilon_i \varepsilon} \right) = 0.
\]

Next we split the fluxes into “left” and “right” components which are assigned to the adjacent energy zones,

\[
F_{i+1/2} = F_{L}^i + F_{R}^i.
\]
\[ F_{i}^R = w_i \epsilon_i - \frac{1}{2} J_{i}^{n+1} (1 - \xi_i). \]  
(B15)

In regions where \( J_i \) varies mildly with \( i \), \( \xi \) is close to 1/2, and hence the total interface flux is approximately given by the arithmetic mean of the flux in the adjacent cells,

\[ F_{i+1/2} = \frac{F_{i}^L + F_{i}^R}{2}. \]  
(B16)

On the other hand, in the high-energy tail of the spectrum \( J_i \) drops rapidly with increasing \( i \), i.e., \( J_{i+1} \ll J_i \). Assuming that the spectrum can be approximated by a Fermi–Dirac distribution (with a small chemical potential in the sense that \( \mu \ll \epsilon_i \)), \( J_i \) is given in terms of the inverse temperature \( \beta \) and a normalization factor \( k \) as

\[ J_i \approx ke^{-\beta \epsilon_i} \epsilon_i^3, \]  
(B17)

and the weighting factor is approximately

\[ \xi \approx e^{-\beta \Delta \epsilon} \ll 1. \]  
(B18)

In this case, the flux components read

\[ F_{i}^L = w_i \epsilon_i + \frac{1}{2} k e^{-\beta \epsilon_i} \epsilon_i^3, \]  
(B19)

\[ F_{i}^R = w_i \epsilon_i - \frac{1}{2} k e^{-\beta \epsilon_i} \epsilon_i^3. \]  
(B20)

Bearing in mind that \( \epsilon^3 \) varies far more slowly than \( e^{-\beta \epsilon} \) in the steep tail of the spectrum, we obtain

\[ F_{i}^L \approx w_i \epsilon_i + \frac{1}{2} k J_{i+1/2}, \quad |F_{i}^R| \approx |w_i \epsilon_i - \frac{1}{2} k J_i| \ll |F_{i-1}|, \quad F_{i+1/2} \approx F_{i}^L. \]  
(B21)

for \( \sigma = 1/2 \), while for \( \sigma = 1 \) we have

\[ F_{i}^L \approx w_i \epsilon_i + \frac{1}{2} k J_{i+1}, \quad F_{i}^R \approx w_i \epsilon_i - \frac{1}{2} k J_i, \quad F_{i+1/2} \approx (w_i + w_{i+1}) \epsilon_i + \frac{1}{2} k J_{i+1}. \]  
(B22)

Thus, in the case of \( \sigma = 1/2 \), the effective interface value of \( J_i \) appearing in the fluxes is given by the weighted geometric mean of \( J_i \) in the adjacent zones; for \( \sigma = 1 \) it is twice the minimum value in either zone, which corresponds more closely (but not exactly) to the weighted harmonic mean. Since \( \sigma = 1 \) generally reduces the absolute values of the fluxes, the resulting scheme is somewhat more robust than for \( \sigma = 1/2 \), i.e., the Newton–Raphson iteration fails to convergence less frequently.

“Harmonic” interpolation with \( \sigma = 1 \) has another welcome side effect in addition to the improved robustness of the scheme: the method of geometric interpolation (also in the original scheme of Rampp & Janka 2002) suffers from a mild “kink” instability, that may produce an unphysical dip in the neutrino spectra. In a long-time simulation of the neutron star cooling phase with VERTEX-PROMETHEUS, in which our new scheme was switched on about 1 s after bounce, we found that “harmonic” interpolation eliminates such a dip, as shown in Figure 15. Because of the favorable stability properties and the smoother neutrino spectra, we preferably work with \( \sigma = 1 \), but \( \sigma = 1/2 \) is also a viable choice.

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