

Numerical relativity in spherical coordinates: A new dynamical spacetime and general relativistic MHD evolution framework for the Einstein Toolkit

Vassilios Mewes^{1,2,3,*} Yosef Zlochower³ Manuela Campanelli³ Thomas W. Baumgarte⁴
Zachariah B. Etienne^{5,6} Federico G. Lopez Armengol^{3,7} and Federico Ciolletta^{8,9,3}

¹*National Center for Computational Sciences, Oak Ridge National Laboratory,*

P.O. Box 2008, Oak Ridge, Tennessee 37831-6164, USA

²*Physics Division, Oak Ridge National Laboratory, P.O. Box 2008,*

Oak Ridge, Tennessee 37831-6354, USA

³*Center for Computational Relativity and Gravitation, and School of Mathematical Sciences, Rochester Institute of Technology, 85 Lomb Memorial Drive, Rochester, New York 14623, USA*

⁴*Department of Physics and Astronomy, Bowdoin College, Brunswick, Maine 04011, USA*

⁵*Department of Physics and Astronomy, West Virginia University,*

Morgantown, West Virginia 26506, USA

⁶*Center for Gravitational Waves and Cosmology, West Virginia University,*

Chestnut Ridge Research Building, Morgantown, West Virginia 26505, USA

⁷*Instituto Argentino de Radioastronomía (IAR), C.C. No. 5, 1894 Buenos Aires, Argentina*

⁸*INFN-TIFPA, Trento Institute for Fundamental Physics and Applications,*

Via Sommarive 14, I-38123 Trento, Italy

⁹*Dipartimento di Fisica, Università di Trento, Via Sommarive 14, I-38123 Trento, Italy*



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We present `SphericalNR`, a new framework for the publicly available `Einstein Toolkit` that numerically solves the Einstein field equations coupled to the equations of general relativistic magneto-hydrodynamic (GRMHD) in a $3+1$ split of spacetime in spherical coordinates without symmetry assumptions. The spacetime evolution is performed using reference-metric versions of either the Baumgarte-Shapiro-Shibata-Nakamura equations or the fully covariant and conformal Z4 system with constraint damping. We have developed a reference-metric version of the *Valencia formulation* of GRMHD with a vector potential method, guaranteeing the absence of magnetic monopoles during the evolution. In our framework, every dynamical field (both spacetime and matter) is evolved using its components in an orthonormal basis with respect to the spherical reference metric. Furthermore, all geometric information about the spherical coordinate system is encoded in source terms appearing in the evolution equations. This allows for the straightforward extension of Cartesian high-resolution shock-capturing finite volume codes to use spherical coordinates with our framework. To this end, we have adapted `GRHydro`, a Cartesian finite volume GRMHD code already available in the `Einstein Toolkit`, to use spherical coordinates. We present the full evolution equations of the framework, as well as details of its implementation in the `Einstein Toolkit`. We validate `SphericalNR` by demonstrating it passes a variety of challenging code tests in static and dynamical spacetimes.

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I. INTRODUCTION

The detection of gravitational waves (GW) from binary black hole (BBH) mergers via the ground-based LIGO and VIRGO detectors [1–6] and the simultaneous detection of GW and electromagnetic (EM) radiation from binary neutron star (BNS) mergers [7–9] has opened a new window into the Universe. Accurate numerical simulations of compact binary mergers are crucial for estimating the physical parameters of detected systems [10] and for

informing physical models about the evolution of matter at nuclear densities in the postmerger remnant of BNSs and BHNSs [11,12]. However, the full self-consistent numerical simulation of a compact object merger through coalescence and subsequent postmerger evolution at high resolution is an extremely challenging computational problem involving vast differences in length and time-scales, as well as different approximate symmetries during the stages of the evolution. In particular, during the inspiral and merger, the absence of approximate axisymmetry lends itself to the use of Cartesian coordinates, while the postmerger remnant has approximate

*mewesv@ornl.gov

symmetries that are better captured by using spherical coordinates.

In the field of numerical relativity, it is now possible to do self-consistent simulations of compact object binary mergers: the first general relativistic hydrodynamics (GRHD) BNS merger simulation [13], the breakthrough simulations of BBH mergers [14–16], the first GRHD black hole neutron star (BHNS) merger simulation [17], the first general relativistic magnetohydrodynamics (GRMHD) BNS merger simulations [18,19], and the first GRMHD simulation of BHNS mergers [20]. Self-consistent simulations of these multimessenger sources require the accurate modeling of the dynamical spacetime evolution and magnetohydrodynamic (MHD) flows within and near compact objects. To this end, many codes have been written that solve the equations of GRMHD. Generally, such codes fall into two categories: GRMHD codes coupled to a $3 + 1$ dynamical spacetime solver (see, e.g., [21–34]) and GRMHD codes that adopt analytical expressions for the spacetime geometry—which could be either exact, if the spacetime is stationary, or approximate for some dynamical spacetimes (see, e.g., [35–45]). In general, the codes coupled to a $3 + 1$ spacetime solver perform the fluid evolution on Cartesian or multipatch grids, while many of the fixed background spacetime codes employ curvilinear coordinates. A notable exception are the codes of [26,27], which solve the fluid equations in curvilinear coordinates, coupled to a dynamical spacetime solver in an approximate, constrained evolution formulation of the Einstein field equations [46–48], which is a generalization of the conformal flatness condition [49,50].

Often, numerical error in conservation of momentum is smallest in the direction of coordinate lines. Accordingly, codes written in Cartesian coordinates conserve linear momentum well, while codes using spherical coordinates conserve angular momentum well. In GRMHD, momenta are only conserved when spacetime (rather than just coordinate) symmetries are present, due to the appearance of source terms in the evolution equations. Many astrophysical systems of interest to multimessenger astrophysics possess a natural axisymmetry at first approximation, so that one expects a better conservation of angular momentum in spherical coordinates, which allow the azimuthal coordinate to be aligned with the direction of this symmetry. Examples include (see references within the cited review articles): core-collapse supernovae [51,52], compact binary merger remnants [53–57], pulsars [58], magnetars [59–61], and self-gravitating accretion disks [62]. While there are techniques to mitigate the nonconservation of angular momentum in Cartesian coordinates (see, e.g., [63–65]), it would generally be more desirable to use numerical grids with spherical sampling, representing all tensors and vectors in the spherical basis. Evolving Einstein’s field equations in spherical coordinates introduces both conceptual and computational problems

associated with coordinate singularities, but many of these issues have been resolved in recent years.

Among the formalisms of Einstein’s field equations most commonly used in numerical simulations is the Baumgarte-Shapiro-Shibata-Nakamura (BSSN) formulation [66,67]. Choices made in the original version of this formulation are suitable in Cartesian coordinates only, but a generalization involving a reference-metric formalism allows for applications in any coordinate system (see, e.g., [68–72]). In the absence of spherical symmetry, coordinate singularities can be handled by properly rescaling components of tensors [73–75], which is equivalent to expressing all tensor components in a frame that is orthogonal with respect to the reference metric. In [76], the formalism was extended to the Z4 formalism (see, e.g., [77–79]). The SENR/NRPy+ code [80] provides a flexible computational framework for the implementation of formalism in a broad class of coordinate systems. Using this framework, we previously implemented this approach in the Einstein Toolkit [81].

In this work, we extend the framework presented in [81] to evolve the GRMHD equations in a reference-metric formalism and add a constraint-damping formulation for the spacetime evolution to the existing framework. Our new code applies this strategy by coupling the GRMHD equations with a fully dynamical spacetime evolution via the BSSN or fully covariant and conformal Z4 (fCCZ4) system [76] on three-dimensional spherical coordinate grids (i.e., no symmetry assumptions are made). The GRMHD evolution equations are evolved using a vector potential method. We also use the moving-puncture gauge conditions [15,16]. To our knowledge, SphericalNR is the first framework solving the coupled BSSN/fCCZ4 and GRMHD equations in three-dimensional spherical coordinate grids without symmetry assumptions.

The paper is organized as follows: In Sec. II, we describe the evolution equations for both spacetime and GRMHD in spherical coordinates. In Sec. III, we describe the implementation of the GRMHD reference-metric evolution formalism in the Einstein Toolkit [82]. Code tests are presented in Sec. IV. Finally, Sec. V contains conclusions and discussions. Throughout this paper and in the code implementation we use geometrized and rationalized (geometrized Heaviside-Lorentz) units in which $c = G = M_{\odot} = \epsilon_0 = \mu_0 = 1$, where c , G , M_{\odot} , ϵ_0 , and μ_0 are the speed of light, gravitational constant, solar mass, vacuum permittivity, and vacuum permeability, respectively. Latin indices denote spatial indices, running from 1 to 3; Greek indices denote spacetime indices, running from 0 to 3 (0 is the time coordinate); and the Einstein summation convention is used.

II. BASIC EQUATIONS: DYNAMICAL SPACETIME AND GRMHD EQUATIONS IN SPHERICAL COORDINATES

A. Spacetime evolution in spherical coordinates

We dynamically evolve the gravitational fields using a numerical relativity framework in spherical coordinates

implemented using the `EinsteinToolkit` infrastructure (see [81]). Our framework builds upon a reference-metric formulation [70–72] of the Baumgarte-Shapiro-Shibata-Nakamura (BSSN) formalism [66,67,83]. We appropriately scale out singular factors from components of tensors so that, for nonsingular spacetimes, all numerically evolved variables remain regular even at the origin and on the polar axis (see [73,75]). In this paper, we extend the spacetime evolution code described in [81] to include the fCCZ4 formalism (see [76]), which applies the same reference-metric formalism and rescaling approach to the CCZ4 evolution equations [79,84]. This represents a conformal reformulation of the constraint-damped Z4 system (see [77,85,86]; see also [78] for an alternative conformal reformulation of Z4). We have now implemented the fCCZ4 formalism in the `SphericalNR` framework, and therefore provide, as a reference, key equations below.

The constraint-damped Z4 system [77,85,86] replaces Einstein’s equations by

$$R_{\mu\nu} + \nabla_{\mu} \mathcal{Z}_{\nu} + \nabla_{\nu} \mathcal{Z}_{\mu} - \frac{\kappa_1}{\alpha} [n_{\mu} \mathcal{Z}_{\nu} + n_{\nu} \mathcal{Z}_{\mu} - (1 + \kappa_2) g_{\mu\nu} n_{\lambda} \mathcal{Z}^{\lambda}] = 8\pi \left(T_{\mu\nu} - \frac{1}{2} g_{\mu\nu} T \right), \quad (1)$$

where $R_{\mu\nu}$ is the (spacetime) Ricci tensor, $g_{\mu\nu}$ the spacetime metric, ∇_{μ} its associated covariant derivative, $T_{\mu\nu}$ the stress-energy tensor, $T \equiv g^{\mu\nu} T_{\mu\nu}$ its trace, \mathcal{Z}_{ν} a four-vector of constraints, and α the lapse function. We will shortly associate the timelike vector n^{μ} with the normal on spatial slices. Finally, κ_1 (units of inverse length) and κ_2 (dimensionless) are two damping coefficients, and all nonconstant constraint related modes are damped when $\kappa_1 > 0$ and $\kappa_2 > -1$ [86].

Following the discussion in [84] regarding the stability of the evolution system in the presence of black holes (BH), we have redefined $\kappa_1 \rightarrow \kappa_1/\alpha$. The Z4 system reduces to the Einstein equations when the constraint vector Z_{μ} vanishes.

We start with a 3 + 1 split of spacetime (see [87]) and foliate the four-dimensional spacetime with a set of nonintersecting spacelike hypersurfaces Σ . We denote the future-pointing, timelike normal on Σ as n_{μ} , and refer to the projection of the spacetime metric $g_{\mu\nu}$ onto Σ as the spatial metric

$$\gamma_{\mu\nu} = g_{\mu\nu} + n_{\mu} n_{\nu}. \quad (2)$$

Expressing the normal vector in terms of a lapse function α and a shift vector β^i ,

$$n^{\mu} = \left(\frac{1}{\alpha}, -\frac{\beta^i}{\alpha} \right) \quad (3)$$

or

$$n_{\mu} = (-\alpha, 0), \quad (4)$$

we may write the four-dimensional line element as

$$ds^2 = g_{\mu\nu} dx^{\mu} dx^{\nu} = -\alpha^2 dt^2 + \gamma_{ij} (dx^i + \beta^i dt)(dx^j + \beta^j dt) \quad (5)$$

and the spacetime metric $g_{\mu\nu}$ as

$$g_{\mu\nu} = \begin{pmatrix} -\alpha^2 + \beta_i \beta^i & \beta_j \\ \beta_i & \gamma_{ij} \end{pmatrix}. \quad (6)$$

As in the BSSN formalism, we conformally rescale the spatial metric according to

$$\bar{\gamma}_{ij} = e^{-4\phi} \gamma_{ij}, \quad (7)$$

where $\bar{\gamma}_{ij}$ is the conformally related metric and e^{ϕ} the conformal factor. The latter can be written as

$$e^{4\phi} = (\gamma/\bar{\gamma})^{1/3}, \quad (8)$$

where γ and $\bar{\gamma}$ are the determinants of the physical and conformally related metric, respectively. We will assume that

$$\partial_t \bar{\gamma} = 0, \quad (9)$$

meaning that $\bar{\gamma}$ remains equal to its initial value throughout the evolution. This choice, referred to as the ‘‘Lagrangian’’ choice in [71], simplifies some expressions in particular in the context of the GRMHD evolution, as explained below. We also rescale the trace-free part of the extrinsic curvature according to

$$\bar{A}_{ij} = e^{-4\phi} \left(K_{ij} - \frac{1}{3} \gamma_{ij} K \right), \quad (10)$$

where K_{ij} is the physical extrinsic curvature and $K \equiv \gamma^{ij} K_{ij}$ its trace.

The central idea of the reference-metric formalism¹ is to express the conformally related metric as the sum of a background metric $\hat{\gamma}_{ij}$ and deviations h_{ij} (which need not to be small),²

¹Splitting the metric into background and departures from the background (which need not to be small) is also done in bimetric formalisms [88–94] in general relativity, in which reference metrics are employed to give physical meaning to pseudotensors in curvilinear coordinates. We emphasize that we do not consider extensions of general relativity here; rather, we use the reference metric only as a convenient approach to express Einstein’s equations (see also [68,69,95,96]).

²Strictly speaking, it is sufficient to introduce a reference connection only (e.g., [71]), but it is convenient to assume that this connection is associated with a reference metric.

$$\bar{\gamma}_{ij} = \hat{\gamma}_{ij} + h_{ij}. \quad (11)$$

For our purposes, it is particularly convenient to choose as the reference metric the flat metric in spherical coordinates,

$$\hat{\gamma}_{ij} = \begin{pmatrix} 1 & 0 & 0 \\ 0 & r^2 & 0 \\ 0 & 0 & r^2 \sin^2 \theta \end{pmatrix}. \quad (12)$$

Another key ingredient is evolving vector and tensor components in the orthonormal basis with respect to $\hat{\gamma}_{ij}$ instead of components in the spherical coordinate basis. To this end, we introduce a set of basis vectors $\hat{\mathbf{e}}_i^{\{k\}}$ that are orthonormal with respect to the background metric $\hat{\gamma}_{ij}$,

$$\hat{\gamma}_{ij} = \delta_{\{k\}\{l\}} \hat{\mathbf{e}}_i^{\{k\}} \hat{\mathbf{e}}_j^{\{l\}}. \quad (13)$$

Since $\hat{\gamma}_{ij}$ is diagonal, the orthonormal vector basis tetrad and its inverse are given by

$$\hat{\mathbf{e}}_i^{\{k\}} = \text{diag}(1, r, r \sin \theta), \quad (14)$$

$$\hat{\mathbf{e}}_{\{i\}}^k = \text{diag}(1, 1/r, 1/(r \sin \theta)), \quad (15)$$

where we have adopted a notation involving plain Latin and Latin indices surrounded with curly braces: the components of a tensor \mathbf{T} in the standard coordinate basis will be denoted using the former, while the tensor components in the background orthonormal basis will be denoted by the latter, respectively. In this notation, we may write the deviation tensor h_{ij} as

$$h_{ij} = \hat{\mathbf{e}}_i^{\{k\}} \hat{\mathbf{e}}_j^{\{l\}} h_{\{k\}\{l\}} \quad (16)$$

and similarly write \bar{A}_{ij} as³

$$\bar{A}_{ij} = \hat{\mathbf{e}}_i^{\{k\}} \hat{\mathbf{e}}_j^{\{l\}} \bar{A}_{\{k\}\{l\}}. \quad (17)$$

While we will write most equations in terms of coordinate components (i.e., indices without curly braces), the code uses components in the orthogonal basis (i.e., with curly indices) as dynamical variables.

As in the original BSSN formalism, we introduce conformal connection functions $\bar{\Lambda}^i$ as independent variables. In the context of the reference-metric formalism, $\bar{\Lambda}^i$ satisfy the initial constraint

$$\bar{\Lambda}^i - \Delta\Gamma^i = 0, \quad (18)$$

³This is a novel notation since, in standard references, different symbols are used for rescaled quantities. For instance, in [73], the pairs $\{h_{ij}, h_{\{i\}\{j\}}\}$, $\{\bar{A}_{ij}, \bar{A}_{\{k\}\{l\}}\}$ are denoted as $\{\epsilon_{ij}, h_{ij}\}$, $\{\bar{A}_{ij}, a_{ij}\}$, respectively.

where

$$\Delta\Gamma^i \equiv \bar{\gamma}^{jk} \Delta\Gamma_{jk}^i \quad (19)$$

and

$$\Delta\Gamma_{jk}^i \equiv \bar{\Gamma}_{jk}^i - \hat{\Gamma}_{jk}^i. \quad (20)$$

Contrary to the Christoffel symbols themselves, differences between Christoffel symbols transform as rank-3 tensors. We compute the $\Delta\Gamma_{jk}^i$ from

$$\begin{aligned} \Delta\Gamma_{jk}^i &= \frac{1}{2} \bar{\gamma}^{il} (\hat{D}_j \bar{\gamma}_{kl} + \hat{D}_k \bar{\gamma}_{jl} - \hat{D}_l \bar{\gamma}_{jk}) \\ &= \frac{1}{2} \bar{\gamma}^{il} (\hat{D}_j h_{kl} + \hat{D}_k h_{jl} - \hat{D}_l h_{jk}), \end{aligned} \quad (21)$$

where \hat{D}_i is the covariant derivative associated with the reference metric $\hat{\gamma}_{ij}$ and where we have used $\hat{D}_i \hat{\gamma}_{jk} = 0$ in the second equality. Derivatives of coordinate components of tensors are evaluated by using the chain rule to analytically take derivatives of the basis vectors $\hat{\mathbf{e}}_i^{\{k\}}$, while the orthonormal components are finite differenced numerically in the code, e.g.,

$$\partial_k h_{ij} = \hat{\mathbf{e}}_i^{\{l\}} \hat{\mathbf{e}}_j^{\{m\}} \partial_k h_{\{l\}\{m\}} + h_{\{l\}\{m\}} \partial_k (\hat{\mathbf{e}}_i^{\{l\}} \hat{\mathbf{e}}_j^{\{m\}}). \quad (22)$$

Similar to our treatment of the metric and extrinsic curvature, we write

$$\bar{\Lambda}^i = \hat{\mathbf{e}}_{\{i\}}^i \bar{\Lambda}^{\{j\}} = \begin{pmatrix} \bar{\Lambda}^{\{r\}} \\ \bar{\Lambda}^{\{\theta\}}/r \\ \bar{\Lambda}^{\{\varphi\}}/(r \sin \theta) \end{pmatrix} \quad (23)$$

and evolve the orthonormal components $\bar{\Lambda}^{\{i\}}$ in our code.

One of the attractive features of the reference-metric formalism is that all quantities, including the conformal connection functions $\bar{\Lambda}^i$, transform as tensor densities of weight zero⁴ (see [71]).

We now extend the above formalism to the Z4 formulation, following [76]. We start with a 3 + 1 decomposition of the constraint vector \mathcal{Z}_μ ,

$$\mathcal{Z}_\mu = g_\mu^\nu \mathcal{Z}_\nu = \gamma_\mu^\nu \mathcal{Z}_\nu - n_\mu n^\nu \mathcal{Z}_\nu, \quad (24)$$

⁴A tensor density of weight w acquires a power J^w in a coordinate transformation, where $J \equiv \det |J^i_j|$ is the determinant of the Jacobian matrix of the coordinate transformation $J^i_j \equiv \frac{\partial x^i}{\partial x^j}$. For example, a rank (2, 0) tensor density of weight w transforms as $\mathcal{T}^{ij} = J^w J^i_k J^j_l \mathcal{T}^{kl}$. Tensor densities of weight zero therefore transform as ordinary (or absolute) tensors with familiar coordinate transformations.

and define

$$\Theta \equiv -n_\lambda Z^\lambda = \alpha Z^0, \quad (25)$$

$$Z_i \equiv \gamma_i^\lambda Z_\lambda. \quad (26)$$

In Eq. (29) below, we will absorb the spatial parts Z_i into the connection functions $\tilde{\Lambda}^i$, but we will evolve Θ as a new independent variable.

In order to write the evolution equations of the fCCZA system, we first define a new tensor

$$\bar{R}_{ij}^{Z^4} \equiv \bar{R}_{ij} + \mathcal{D}_i Z_j + \mathcal{D}_j Z_i, \quad (27)$$

where \mathcal{D}_i is the covariant derivative associated with the spatial metric γ_{ij} and \bar{R}_{ij} is the Ricci tensor associated with the conformal metric $\tilde{\gamma}_{ij}$,

$$\begin{aligned} \bar{R}_{ij} = & -\frac{1}{2} \tilde{\gamma}^{kl} \hat{\mathcal{D}}_k \hat{\mathcal{D}}_l \tilde{\gamma}_{ij} + \tilde{\gamma}_{k(i} \hat{\mathcal{D}}_{j)} \Delta \Gamma^k + \Delta \Gamma^k \Delta \Gamma_{(ij)k} \\ & + \tilde{\gamma}^{kl} (2\Delta \Gamma_{k(i}^m \Delta \Gamma_{j)ml} + \Delta \Gamma_{ik}^m \Delta \Gamma_{mj}), \end{aligned} \quad (28)$$

where parentheses around indices indicate the symmetric part of a tensor $T_{(ij)} \equiv \frac{1}{2}(T_{ij} + T_{ji})$. We next define new conformal connection functions according to

$$\tilde{\Lambda}^i \equiv \Delta \Gamma^i + 2\tilde{\gamma}^{ij} Z_j, \quad (29)$$

where we have used a tilde in order to distinguish these objects from $\tilde{\Lambda}^i$. We then have

$$Z^i = \frac{1}{2} e^{-4\phi} (\tilde{\Lambda}^i - \Delta \Gamma^i). \quad (30)$$

With these definitions, we may now write $R_{ij}^{Z^4}$ as

$$\begin{aligned} \bar{R}_{ij}^{Z^4} = & -\frac{1}{2} \tilde{\gamma}^{kl} \hat{\mathcal{D}}_k \hat{\mathcal{D}}_l \tilde{\gamma}_{ij} + \tilde{\gamma}_{k(i} \hat{\mathcal{D}}_{j)} (\tilde{\Lambda}^k - 2e^{4\phi} Z^k) \\ & + \mathcal{D}_i Z_j + \mathcal{D}_j Z_i + \Delta \Gamma^k \Delta \Gamma_{(ij)k} \\ & + \tilde{\gamma}^{kl} (2\Delta \Gamma_{k(i}^m \Delta \Gamma_{j)ml} + \Delta \Gamma_{ik}^m \Delta \Gamma_{mj}). \end{aligned} \quad (31)$$

Combining the terms $\tilde{\gamma}_{k(i} \hat{\mathcal{D}}_{j)} (-2e^{4\phi} Z^k)$ and $\mathcal{D}_i Z_j + \mathcal{D}_j Z_i$, it can be seen that all partial derivatives $\partial_i Z_j$ in $\bar{R}_{ij}^{Z^4}$ cancel out exactly, meaning that $R_{ij}^{Z^4}$ reduces to

$$\bar{R}_{ij}^{Z^4} = \bar{R}_{ij} - 8Z_{(i} \partial_{j)} \phi + 2\gamma_{k(i} (\Gamma_{j)l}^k - \hat{\Gamma}_{j)l}^k) Z^l. \quad (32)$$

With the above, and defining $\partial_0 \equiv \partial_t - \mathcal{L}_\beta$, where \mathcal{L}_β is the Lie derivative along the shift β^i , we arrive at the following set of coordinate basis evolution equations for the fCCZA system:

$$\begin{aligned} \partial_0 \tilde{\gamma}_{ij} = & -\frac{2}{3} \tilde{\gamma}_{ij} \bar{\mathcal{D}}_k \beta^k - 2\alpha \bar{A}_{ij}, \\ \partial_0 \phi = & \frac{1}{6} \bar{\mathcal{D}}_i \beta^i - \frac{1}{6} \alpha K, \\ \partial_0 \bar{A}_{ij} = & -\frac{2}{3} \bar{A}_{ij} \bar{\mathcal{D}}_k \beta^k - 2\alpha \bar{A}_{ik} \bar{A}_j^k + \alpha \bar{A}_{ij} (K - 2\Theta) \\ & + e^{-4\phi} [-2\alpha \bar{\mathcal{D}}_i \bar{\mathcal{D}}_j \phi + 4\alpha \bar{\mathcal{D}}_i \phi \bar{\mathcal{D}}_j \phi \\ & + 4\bar{\mathcal{D}}_{(i} \alpha \bar{\mathcal{D}}_{j)} \phi - \bar{\mathcal{D}}_i \bar{\mathcal{D}}_j \alpha \\ & + \alpha (\bar{R}_{ij}^{Z^4} - 8\pi S_{ij})]^{TF}, \\ \partial_0 K = & e^{-4\phi} [\alpha (\bar{R}^{Z^4} - 8\bar{\mathcal{D}}^i \bar{\mathcal{D}}_i \phi - 8\bar{\mathcal{D}}^2 \phi) \\ & - (2\bar{\mathcal{D}}^i \alpha \bar{\mathcal{D}}_i \phi + \bar{\mathcal{D}}^2 \alpha)] + \alpha (K^2 - 2\Theta K) \\ & - 3\kappa_1 (1 + \kappa_2) \Theta + 4\pi \alpha (S - 3E), \\ \partial_0 \Theta = & \frac{1}{2} \alpha [e^{-4\phi} (\bar{R}^{Z^4} - 8\bar{\mathcal{D}}^i \phi \bar{\mathcal{D}}_i \phi - 8\bar{\mathcal{D}}^2 \phi) \\ & - \bar{A}^{ij} \bar{A}_{ij} + \frac{2}{3} K^2 - 2\Theta K] \\ & - Z^i \partial_i \alpha - \kappa_1 (2 + \kappa_2) \Theta - 8\pi \alpha E, \\ \partial_0 \tilde{\Lambda}^i = & \tilde{\gamma}^{jk} \hat{\mathcal{D}}_j \hat{\mathcal{D}}_k \beta^i + \frac{2}{3} \Delta \Gamma^i \bar{\mathcal{D}}_j \beta^j + \frac{1}{3} \bar{\mathcal{D}}^i \bar{\mathcal{D}}_j \beta^j \\ & - 2\bar{A}^{jk} (\delta_j^i \partial_k \alpha - 6\alpha \delta_j^i \partial_k \phi - \alpha \Delta \Gamma_{jk}^i) \\ & - \frac{4}{3} \alpha \tilde{\gamma}^{ij} \partial_j K + 2\tilde{\gamma}^{ij} (\alpha \partial_j \Theta - \Theta \partial_j \alpha) - \frac{4}{3} \alpha K e^{4\phi} Z^i \\ & - 2\kappa_1 e^{4\phi} Z^i + 2\kappa_3 e^{4\phi} \left(\frac{2}{3} Z^i \hat{\mathcal{D}}_k \beta^k - Z^k \hat{\mathcal{D}}_k \beta^i \right) \\ & - 16\pi \alpha \tilde{\gamma}^{ij} S_j. \end{aligned}$$

Here \square^{TF} denotes the trace-free part of a tensor $T_{ij}^{TF} \equiv T_{ij} - \frac{1}{3} \gamma_{ij} T^k_k$, and κ_3 is a constant that determines the covariance of the equations, in particular $\kappa_3 = 1$ (the choice adopted for all simulations presented in this work), corresponds to full covariance [79]. Unlike in [76], we have absorbed all covariant derivatives of Z^i in $\bar{R}_{ij}^{Z^4}$, meaning that no derivatives of the constraint vector appear in the above equations.

During the time evolution, we continuously enforce $\partial_i \tilde{\gamma} = 0$, as well as the constraint $\tilde{A}_i^i = 0$. We have also implemented the $\chi = e^{4\phi}$ [15] and $W = e^{2\phi}$ [97] variants of the conformal factor evolution, resulting in the following evolution equations:

$$\partial_0 \chi = -\frac{2}{3} \chi \bar{\mathcal{D}}_i \beta^i + \frac{2}{3} \chi \alpha K, \quad (33)$$

$$\partial_0 W = -\frac{1}{3} W \bar{\mathcal{D}}_i \beta^i + \frac{1}{3} W \alpha K. \quad (34)$$

The choice between these variants can be made at runtime.

In order to close the evolution system, we need to choose a gauge. For all dynamical spacetime evolution simulations presented in this paper, we adopt the ‘‘standard gauge,’’ or ‘‘moving-puncture gauge,’’ meaning 1 + log slicing,

$$\partial_t \alpha = -2(K - 2\Theta) + \beta^i \partial_i \alpha \quad (35)$$

for the lapse (see [98]) and a Γ -driver,⁵

$$\partial_t \beta^i = B^i + \beta^j \hat{D}_j \beta^i, \quad (36)$$

$$\partial_t B^i = \frac{3}{4} (\partial_t \tilde{\Lambda}^i - \beta^j \hat{D}_j \tilde{\Lambda}^i) + \beta^j \hat{D}_j B^i - \eta B^i \quad (37)$$

for the shift (see [99]). Here η is a damping parameter with dimensions of inverse length, and we have adopted the covariant form of [71]. In the code, the inclusion of shift advection terms can be turned off by the user at runtime.

As noted above, the actual evolved tensors in the code are not in the coordinate basis $\{\tilde{\gamma}_{ij}, \tilde{A}_{ij}, \tilde{\Lambda}^i, \beta^i, B^i = \dot{\beta}^i\}$, but rather the orthonormal basis $\{h_{\{i\}\{j\}}, \tilde{A}_{\{i\}\{j\}}, \tilde{\Lambda}^{\{i\}}, \beta^{\{i\}}, B^{\{i\}} = \dot{\beta}^{\{i\}}\}$, respectively. The remaining evolved quantities $\{\alpha, K, \Theta, \phi\}$ are scalars and thus do not depend on choice of basis.

Finally, the matter sources in the spacetime evolution equations are given by projections of the stress-energy tensor, namely,

$$E \equiv n^\mu n^\nu T_{\mu\nu} = \frac{1}{\alpha^2} (T_{tt} - 2\beta^i T_{ti} + \beta^i \beta^j T_{ij}), \quad (38)$$

$$S_i \equiv -\gamma_{i\mu} n_\nu T^{\mu\nu} = -\frac{1}{\alpha} T_{ti} + \frac{1}{\alpha} \beta^j T_{ij}, \quad (39)$$

$$S_{ij} \equiv \gamma_{i\mu} \gamma_{j\nu} T^{\mu\nu}, \quad (40)$$

$$S \equiv \gamma^{ij} S_{ij}. \quad (41)$$

While we have only presented the fCCZ4 equations above, the user can select to evolve the BSSN system at runtime as well.

B. GRMHD in the reference-metric formalism

In this section, we review the reference-metric formalism for GRHD presented in [74,75] and extend it to GRMHD, using a vector potential evolution scheme to guarantee the absence of magnetic monopoles.

1. Conservation laws in four-dimensional form

The evolution of a magnetized fluid is governed by the conservation of baryon number

$$\nabla_\mu (\rho u^\mu) = 0, \quad (42)$$

which results in the continuity equation, and the conservation of energy-momentum

⁵The Γ -driver owes its name to the appearance of the conformal connection functions $\tilde{\Gamma}^i$ introduced in [67]. While we use $\tilde{\Lambda}^i$ here, rather than $\tilde{\Gamma}^i$, we still use the name Γ -driver for this gauge condition.

$$\nabla_\mu T^{\mu\nu} = \nabla_\mu (T_{\text{matter}}^{\mu\nu} + T_{\text{EM}}^{\mu\nu}) = 0, \quad (43)$$

which results in the (relativistic) Euler equation and the conservation of total energy. Assuming a perfect fluid, the fluid stress-energy tensor $T_{\text{matter}}^{\mu\nu}$ is given by

$$T_{\text{matter}}^{\mu\nu} = \rho h u^\mu u^\nu + P g^{\mu\nu}, \quad (44)$$

where ρ is the rest-mass density, P the fluid pressure, $h = 1 + \epsilon + P/\rho$ the specific enthalpy, ϵ the internal energy density, and u^μ the fluid four-velocity, respectively.

In terms of the Faraday tensor $F^{\mu\nu}$, the EM stress-energy tensor is⁶

$$T_{\text{EM}}^{\mu\nu} = F^{\mu\lambda} F^\nu{}_\lambda - \frac{1}{4} g^{\mu\nu} F^{\lambda\kappa} F_{\lambda\kappa}. \quad (45)$$

We next decompose the Faraday tensor $F^{\mu\nu}$ as

$$F^{\mu\nu} = U^\mu E^\nu{}_{(U)} - U^\nu E^\mu{}_{(U)} + \epsilon^{\mu\nu\lambda\kappa} U_\lambda B_{\kappa(U)}, \quad (46)$$

where

$$\epsilon^{\mu\nu\lambda\kappa} \equiv \frac{-1}{\sqrt{-g}} [\mu\nu\lambda\kappa], \quad (47)$$

and where $[\mu\nu\lambda\kappa]$ is the totally antisymmetric Levi-Civita symbol ($= + (-)1$ for even (odd) permutations of [0123], and 0 if any two indices are repeated). Here $E^\mu{}_{(U)}$ and $B^\mu{}_{(U)}$ are the electric and magnetic fields measured by an observer with *generic* four-velocity U^μ ,

$$E^\mu{}_{(U)} = F^{\mu\nu} U_\nu, \quad B^\mu{}_{(U)} = \frac{1}{2} \epsilon^{\mu\nu\kappa\rho} U_\nu F_{\rho\kappa}. \quad (48)$$

Both $E^\mu{}_{(U)}$ and $B^\mu{}_{(U)}$ are orthogonal to U^μ , i.e., $E^\mu{}_{(U)} U_\mu = B^\mu{}_{(U)} U_\mu = 0$.

In the following, we focus on two observers of particular interest, namely observers comoving with the fluid (i.e., with four-velocity $U^\mu = u^\mu$) and normal observers (with four-velocity $U^\mu = n^\mu$). Following convention, we denote the fields observed by the former with $E^\mu{}_{(u)}$ and $B^\mu{}_{(u)}$, but the latter simply with $E^\mu = E^\mu{}_{(n)}$ and $B^\mu = B^\mu{}_{(n)}$.

In the ideal MHD limit, we assume that the fluid acts as a perfect conductor, meaning that the electric field observed by an observer comoving with the fluid vanishes,

⁶The Heaviside-Lorentz (HL) units we adopt in this work are rationalized, as no explicit factors of 4π appear in the Maxwell equations in these units. Electric and magnetic fields in HL and Gauss units are therefore related by a factor of $\sqrt{4\pi}$: $E^\mu_{[\text{Gauss}]} = \sqrt{4\pi} E^\mu_{[\text{HL}]}$, $B^\mu_{[\text{Gauss}]} = \sqrt{4\pi} B^\mu_{[\text{HL}]}$. Consequently, in Gauss units, the EM stress-energy tensor is defined as $T_{\text{EM}}^{\mu\nu} = \frac{1}{4\pi} (F^{\mu\lambda} F^\nu{}_\lambda - \frac{1}{4} g^{\mu\nu} F^{\lambda\kappa} F_{\lambda\kappa})_{[\text{Gauss}]}$.

$$E_{(u)}^\mu = F^{\mu\nu} u_\nu = 0. \quad (49)$$

Thus, in this approximation $F^{\mu\nu}$, which generally depends on both electric and magnetic fields, can be expressed in terms of magnetic fields alone,

$$F^{\mu\nu} = \epsilon^{\mu\nu\lambda\kappa} u_\lambda B_\kappa^{(u)} \quad (50)$$

and $*F^{\mu\nu}$, the dual of the Faraday tensor, as

$$*F^{\mu\nu} = \frac{1}{2} \epsilon^{\mu\nu\lambda\kappa} F_{\lambda\kappa} = u^\mu B_{(u)}^\nu - u^\nu B_{(u)}^\mu. \quad (51)$$

In the ideal MHD limit, we can also write the magnetic field $B_{(u)}^\mu$ as a projection of B^μ along the fluid four-velocity u^ν ,

$$B_{(u)}^\mu = \frac{1}{W} P_\nu^\mu B^\nu, \quad (52)$$

where $P_\nu^\mu = g_\nu^\mu + u^\mu u_\nu$ and where we have introduced the Lorentz factor W between the fluid and normal observers, $W \equiv -n_\mu u^\mu = \alpha u^t$. Inserting (52) into (51) yields

$$*F^{\mu\nu} = \frac{1}{W} (u^\mu B^\nu - u^\nu B^\mu). \quad (53)$$

Introducing the abbreviation⁷

$$b^\mu \equiv B_{(u)}^\mu, \quad (54)$$

we can write the electromagnetic stress-energy tensor as

$$T_{\text{EM}}^{\mu\nu} = \left(u^\mu u^\nu + \frac{1}{2} g^{\mu\nu} \right) b^2 - b^\mu b^\nu, \quad (55)$$

where $b^2 \equiv b^\mu b_\mu$, so that the total stress-energy tensor $T^{\mu\nu}$ becomes

$$T^{\mu\nu} = \rho h^* u^\mu u^\nu + P^* g^{\mu\nu} - b^\mu b^\nu. \quad (56)$$

Here we have defined $h^* = 1 + \epsilon + (P + b^2)/\rho$ as the magnetically modified specific enthalpy and $P^* = P + b^2/2$ as the magnetically modified isotropic pressure. Finally, the evolution of the magnetic field is governed by the homogeneous Maxwell equations

$$\nabla_\nu (*F^{\mu\nu}) = 0. \quad (57)$$

⁷Similarly, to differences in the EM stress-energy tensor, treatments that adopt Gauss units ([22,29,36,38,41]) rather than Lorentz-Heaviside units ([21,23–28,30–35,37,39,40,42–45]) often use the definition $b^\mu \equiv B_{(u)}^\mu/\sqrt{4\pi}$ instead of (54). The resulting fields b^μ , however, are again identical in both treatments, so that expressions for the stress-energy tensor, for example, take the same form when written in terms of b^μ ; see (55).

2. The 3+1 GRMHD equations in the reference-metric formalism

We now recast the above conservation laws using both a 3 + 1 split and a reference-metric approach. The result will be a set of equations that is suitable for numerical integration in spherical coordinates and that meshes well with the form of the field equations as presented in Sec. II A. The fluid equations have been previously derived in [74,75], and we will extend the formalism to the GRMHD equations.

The key idea is to repeatedly use identities for divergences (see, e.g., Problem 7.7 in [100]). For the continuity equation (42), for example, we use

$$\nabla_\nu V^\nu = \frac{1}{\sqrt{|g|}} \partial_\nu (\sqrt{|g|} V^\nu) \quad (58)$$

to arrive at

$$\begin{aligned} 0 &= \nabla_\nu (\rho u^\nu) = \frac{1}{\sqrt{-g}} \partial_\nu (\sqrt{-g} \rho u^\nu) \\ &= \frac{1}{\sqrt{-g}} (\partial_t (\sqrt{-g} \rho u^t) + \partial_i (\sqrt{-g} \rho u^i)) \\ &= \partial_t (e^{6\phi} \sqrt{\tilde{\gamma}} \rho W) + \partial_i (\alpha e^{6\phi} \sqrt{\tilde{\gamma}} \rho W \bar{v}^i). \end{aligned} \quad (59)$$

Here we have defined

$$v^i \equiv \frac{1}{W} \gamma_\nu^i u^\nu = \frac{u^i}{W} + \frac{\beta^i}{\alpha} \quad (60)$$

as the fluid three-velocity and

$$\bar{v}^i \equiv v^i - \frac{\beta^i}{\alpha} = \frac{u^i}{W} \quad (61)$$

as the ‘‘advection velocity.’’ We have also written the square root of the determinant of the spacetime metric $\sqrt{-g}$ as

$$\sqrt{-g} = \alpha \sqrt{\tilde{\gamma}} = \alpha e^{6\phi} \sqrt{\tilde{\gamma}}. \quad (62)$$

Equation (59) is the continuity equation in a form that is often referred to as the ‘‘Valencia’’ form of the equations (see [101]). This version of the equations is well suited for simulations in Cartesian coordinates, but in curvilinear coordinates the vanishing of the determinant $\tilde{\gamma}$ may cause numerical problems. Following [74], we apply the identity (58) again, and convert the partial derivatives ∂_i to covariant derivatives \hat{D}_i associated with the background (reference) metric $\hat{\gamma}_{ij}$,⁸

$$0 = \partial_t (e^{6\phi} \sqrt{\tilde{\gamma}} / \hat{\gamma} \rho W) + \hat{D}_i (\alpha e^{6\phi} \sqrt{\tilde{\gamma}} / \hat{\gamma} \rho W \bar{v}^i). \quad (63)$$

⁸We will assume throughout that the reference metric $\hat{\gamma}_{ij}$ is independent of time.

Note that the combination $\bar{\gamma}/\hat{\gamma}$ remains finite for regular spacetimes. We now define the conserved density D ,

$$D \equiv e^{6\phi} \sqrt{\bar{\gamma}/\hat{\gamma}} \rho W \quad (64)$$

and the conserved density flux f_D ,

$$(f_D)^i \equiv \alpha D \bar{v}^i \quad (65)$$

to write the continuity equation in the form

$$\partial_t D + \hat{D}_i (f_D)^i = 0. \quad (66)$$

We note that a similar strategy is followed in the `CoCoNuT` code, without explicitly mentioning the reference metric (see, e.g., [102]). For reasons that will become apparent in Sec. II D below, we will implement the equation numerically as

$$\partial_t D + \partial_i (f_D)^i = -(f_D)^i \hat{\Gamma}_{ij}^j. \quad (67)$$

We proceed similarly for the conservation of stress energy (43), except that we now use the identity

$$\nabla_\lambda T_\mu^\lambda = \frac{1}{\sqrt{|g|}} \partial_\lambda (\sqrt{|g|} T_\mu^\lambda) - T_\sigma^\lambda \Gamma_{\lambda\mu}^\sigma \quad (68)$$

twice. The spatial projection of (43) then yields the relativistic Euler equation

$$\partial_t S_j + \hat{D}_i (f_S)^i_j = (s_S)_j \quad (69)$$

or

$$\partial_t S_j + \partial_i (f_S)^i_j = (s_S)_j - (f_S)^i_j \hat{\Gamma}_{ik}^k + (f_S)^i_k \hat{\Gamma}_{ij}^k, \quad (70)$$

where we have defined the conserved momentum

$$S_j \equiv e^{6\phi} \sqrt{\bar{\gamma}/\hat{\gamma}} T_j^i = e^{6\phi} \sqrt{\bar{\gamma}/\hat{\gamma}} (\rho h^* W^2 v_j - \alpha b^0 b_j), \quad (71)$$

the conserved momentum density fluxes

$$(f_S)^i_j \equiv \alpha (S_j \bar{v}^i + e^{6\phi} \sqrt{\bar{\gamma}/\hat{\gamma}} P^* \delta_j^i - e^{6\phi} \sqrt{\bar{\gamma}/\hat{\gamma}} b_j B^i / W) \quad (72)$$

and where the source term is given by

$$(s_S)_j = \alpha e^{6\phi} \sqrt{\bar{\gamma}/\hat{\gamma}} (-T^{00} \alpha \partial_j \alpha + T_i^0 \hat{D}_i \beta^i + \frac{1}{2} (T^{00} \beta^i \beta^k + 2T^{0i} \beta^k + T^{ik}) \hat{D}_j \gamma_{ik}) \quad (73)$$

(see [74] for a detailed derivation). The terms $\hat{D}_j \gamma_{ik}$ can be evaluated from

$$\hat{D}_j \gamma_{ik} = e^{4\phi} (4\bar{\gamma}_{ik} \partial_j \phi + \hat{D}_j \bar{\gamma}_{ik}), \quad (74)$$

where $\hat{D}_j \bar{\gamma}_{ik} = \hat{D}_j h_{ik}$ are computed already in (21).

Projecting the conservation of stress energy (43) along n_ν and subtracting the continuity equation (42) yields

$$\nabla_\mu (n_\nu T^{\nu\mu} - \rho u^\mu) = T^{\mu\nu} \nabla_\nu n_\mu. \quad (75)$$

We again apply the identity (58) twice to arrive at the energy equation

$$\partial_t \tau + \hat{D}_i (f_\tau)^i = s_\tau \quad (76)$$

or

$$\partial_t \tau + \partial_i (f_\tau)^i = s_\tau - (f_\tau)^i \hat{\Gamma}_{ij}^j, \quad (77)$$

where we have defined τ as the total conserved energy density subtracting the conserved density D ,⁹

$$\tau \equiv e^{6\phi} \sqrt{\bar{\gamma}/\hat{\gamma}} (\rho h^* W^2 - P^* - (\alpha b^0)^2) - D, \quad (78)$$

the conserved energy flux

$$(f_\tau)^i \equiv \alpha (\tau \bar{v}^i + e^{6\phi} \sqrt{\bar{\gamma}/\hat{\gamma}} P^* v^i - \alpha e^{6\phi} \sqrt{\bar{\gamma}/\hat{\gamma}} b^0 B^i / W) \quad (79)$$

and where the source term is given by [74]

$$s_\tau = \alpha e^{6\phi} \sqrt{\bar{\gamma}/\hat{\gamma}} (T^{00} (\beta^i \beta^j K_{ij} - \beta^i \partial_i \alpha) + T^{0i} (2\beta^j K_{ij} - \partial_i \alpha) + T^{ij} K_{ij}). \quad (80)$$

In the above equations, both B^i and $b^\mu = B_{(u)}^\mu$ make an appearance. The two fields are related by (52), so that we can always compute $B_{(u)}^\mu$ from B^μ . Specifically, we contract (52) with n_μ to obtain

$$b^0 = \frac{W B^i v_i}{\alpha}, \quad (81)$$

while a spatial projection of (52) yields

$$b^i = \frac{B^i}{W} + W (B^j v_j) \bar{v}^i. \quad (82)$$

We also have

⁹The motivation to subtract the continuity equation from the projection of the stress energy conservation along n_μ (75) was to arrive at an evolution for τ : it correctly recovers the Newtonian limit and is numerically more accurate than evolving the total conserved energy density (see, e.g., [23,103,104]).

$$b^2 = \frac{B^i B_i}{W^2} + (B^i v_i)^2. \quad (83)$$

We now adopt the same approach to rewrite Maxwell's equations (57). Since the Faraday tensor (as well as its dual) is antisymmetric, we now use the identity

$$\nabla_\nu A^{\mu\nu} = \frac{1}{\sqrt{|g|}} \partial_\nu (\sqrt{|g|} A^{\mu\nu}) \quad (84)$$

for antisymmetric tensors $A^{\mu\nu}$. Inserting (53) into (57) and using (84), we obtain

$$\partial_t (\sqrt{\gamma} B^\mu) = \partial_i \left(\frac{\alpha \sqrt{\gamma}}{W} (u^\mu B^i - u^i B^\mu) \right). \quad (85)$$

The temporal component of this equation results in the solenoidal constraint, stating the absence of magnetic monopoles,

$$\partial_i (\sqrt{\gamma} B^i) = \sqrt{\hat{\gamma}} \hat{D}_i (e^{6\phi} \sqrt{\hat{\gamma}} / \hat{\gamma} B^i) = 0. \quad (86)$$

We now define

$$\mathcal{B}^i \equiv e^{6\phi} \sqrt{\hat{\gamma}} / \hat{\gamma} B^i, \quad (87)$$

so that (86) reduces to

$$\hat{D}_i \mathcal{B}^i = 0. \quad (88)$$

For the spatial part of (85), we use (84) again to obtain the induction equation in 3 + 1 form

$$\partial_t \mathcal{B}^j = \hat{D}_i (\alpha (\bar{v}^j \mathcal{B}^i - \bar{v}^i \mathcal{B}^j)). \quad (89)$$

This form of the solenoidal constraint and continuity equation is very similar to the one presented in [26,105], but solved there using a constraint transport approach, while we evolve the vector potential of the magnetic field in our framework instead.

3. Vector potential evolution equations

Numerically evolving the induction equation (89) directly is generally problematic, since accumulating numerical error will typically result in the magnetic field having nonvanishing divergence. The resulting growth of spurious magnetic monopoles has severe consequences of the evolution, since it will result in nonphysical fluid acceleration in the direction of the magnetic field (see, e.g., [106]).

Various approaches have been implemented to avoid this growth of magnetic monopoles in (GR)MHD simulations

(see [107] for a comprehensive overview). The three most commonly adopted approaches in GRMHD codes are (i) hyperbolic divergence cleaning via a generalized Lagrange multiplier [108]; (ii) constrained transport (CT) [109] schemes in which the magnetic field is updated in such a way that the divergence (measured in a finite-difference stencil that is compatible with the base CT scheme) remains unchanged to round off during the evolution; and (iii) evolving not the magnetic field directly but rather its vector potential and taking the curl of the vector potential in order to compute the magnetic field [22,25,34]. As the divergence of the curl of a vector field is identically zero, the latter approach guarantees a solenoidal magnetic field to round-off error during the evolution.

In developing the GRMHD evolution framework in spherical coordinates, we opted to implement the latter, namely evolving the vector potential in a cell-centered fashion. We choose a vector-potential formulation for four reasons: (1) the resulting equations can be easily incorporated into our reference-metric formalism; (2) contrary to the hyperbolic divergence cleaning, the solenoidal constraint is automatically fulfilled to machine precision; (3) there is no need to extend our internal parity boundary conditions to deal with the staggered magnetic fields used in CT schemes (though see the CoCoNuT [26,105] and Aenus [110] codes for implementations of staggered CT schemes in spherical coordinates); and (4) it has recently been shown to be strongly hyperbolic [111].

Imposing the ideal MHD limit again, and taking a projection of Eq. (49) with the spatial metric (2) shows that the electric and magnetic field as observed by the normal observer are related by

$$E^\mu = \frac{-1}{W} e^{\mu\nu\lambda\kappa} u_\nu n_\lambda B_\kappa. \quad (90)$$

Defining the three-dimensional antisymmetric tensor as

$$\epsilon_{\mu\nu\lambda} = n^\kappa \epsilon_{\kappa\mu\nu\lambda}, \quad \text{or} \quad \epsilon^{\mu\nu\lambda} = n_\kappa \epsilon^{\kappa\mu\nu\lambda}, \quad (91)$$

so that $\epsilon_{ijk} = \sqrt{\gamma} [ijk]$, and we may rewrite (90) as

$$E_i = -\epsilon_{ijk} \bar{v}^j B^k, \quad (92)$$

where we have used $\epsilon_{tij} = -\beta^k \epsilon_{ikj}$ (see Eq. (32) in [112]).

We now introduce a four-vector potential

$$\mathcal{A}_\mu = \Phi n_\mu + A_\mu, \quad (93)$$

where Φ is the electromagnetic scalar potential and A_μ is purely spatial, $A_\mu n^\mu = 0$, so that $A_t = \beta^i A_i$ and $\mathcal{A}_t = -\alpha\Phi + \beta^i A_i$. Writing the Faraday tensor in terms of \mathcal{A}_μ yields

$$F_{\mu\nu} = \partial_\mu \mathcal{A}_\nu - \partial_\nu \mathcal{A}_\mu = n_\mu E_\nu + n_\nu E_\mu + \epsilon_{\mu\nu\lambda} B^\lambda. \quad (94)$$

Contracting this with $\epsilon^{\mu\nu\lambda}$ yields

$$\epsilon^{\mu\nu\lambda} (\partial_\mu \mathcal{A}_\nu - \partial_\nu \mathcal{A}_\mu) = \epsilon^{\mu\nu\lambda} \epsilon_{\mu\nu\kappa} B^\kappa = 2B^\lambda, \quad (95)$$

or

$$B^i = \epsilon^{ijk} \partial_j \mathcal{A}_k. \quad (96)$$

Inserting the definition (87) then results in

$$B^i = \hat{\epsilon}^{ijk} \partial_j A_k = \hat{\epsilon}^{ijk} \hat{\mathcal{D}}_j A_k, \quad (97)$$

where $\hat{\epsilon}_{ijk} = \sqrt{\hat{\gamma}} [ijk]$ and $\hat{\epsilon}^{ijk} = \hat{\gamma}^{-1/2} [ijk]$.

Finally, we may evaluate a mixed time-space component of (94) to find

$$\partial_t A_i = -\alpha E_i + \epsilon_{ij} B^j - \hat{\mathcal{D}}_i (\alpha \Phi - \beta^j A_j). \quad (98)$$

Inserting (92) and expressing the result in terms of \mathcal{B}^i result in

$$\partial_t A_i = \alpha \hat{\epsilon}_{ijk} \bar{v}^j \mathcal{B}^k - \hat{\mathcal{D}}_i (\alpha \Phi - \beta^j A_j). \quad (99)$$

In our code, we evolve the spatial components of the vector potential A_i , and compute, at each time step, the conserved magnetic field from (97).

We evolve the vector potential in a ‘‘generalized Lorenz gauge’’ [113],

$$\nabla_\mu A^\mu = -\zeta \Phi, \quad (100)$$

where ζ is a damping parameter with dimensions of inverse length. As before, we use the identity (58) twice to rewrite this as

$$\begin{aligned} \partial_t (e^{\delta\phi} \sqrt{\hat{\gamma}/\hat{\gamma}} \Phi) + \hat{\mathcal{D}}_i (\alpha e^{\delta\phi} \sqrt{\hat{\gamma}/\hat{\gamma}} A^i - e^{\delta\phi} \sqrt{\hat{\gamma}/\hat{\gamma}} \beta^i \Phi) \\ = -\zeta \alpha e^{\delta\phi} \sqrt{\hat{\gamma}/\hat{\gamma}} \Phi. \end{aligned} \quad (101)$$

We now define

$$\hat{\Phi} \equiv e^{\delta\phi} \sqrt{\hat{\gamma}/\hat{\gamma}} \Phi \quad (102)$$

and

$$(f_\Phi)^i \equiv \alpha e^{\delta\phi} \sqrt{\hat{\gamma}/\hat{\gamma}} A^i - \beta^i \hat{\Phi} \quad (103)$$

and evolve (101) as

$$\partial_t \hat{\Phi} + \partial_i (f_\Phi)^i = -\zeta \alpha \hat{\Phi} - (f_\Phi)^i \hat{\Gamma}_{ij}^j. \quad (104)$$

In all applications shown in this paper, we followed [29] and chose $\zeta = 1.5/\Delta t$,¹⁰ where Δt is the global time step of our numerical evolution.

We note that our algorithm is not staggered, i.e., the vector potential A_i lives at the cell centers as do all other variables. In order to update the magnetic field, we calculate the curl of A_i (97), where we apply the product rule and take derivatives of the scale factors analytically, as with all other fields. Initial data for magnetic fields are generated in the same way, namely by taking the curl of a prescribed initial vector potential A_i .

4. Summary

In summary, the GRMHD evolution system in the reference-metric formalism is composed of the following conserved quantities:

$$D \equiv e^{\delta\phi} \sqrt{\hat{\gamma}/\hat{\gamma}} \rho W, \quad (105)$$

$$S_j \equiv e^{\delta\phi} \sqrt{\hat{\gamma}/\hat{\gamma}} (\rho h^* W^2 v_j - \alpha b^0 b_j), \quad (106)$$

$$\tau \equiv e^{\delta\phi} \sqrt{\hat{\gamma}/\hat{\gamma}} (\rho h^* W^2 - P^* - (\alpha b^0)^2) - D, \quad (107)$$

$$\mathcal{B}^j \equiv e^{\delta\phi} \sqrt{\hat{\gamma}/\hat{\gamma}} B^j, \quad (108)$$

where we note that, unlike the corresponding conserved variables of the Valencia formulation, these are true scalars and vectors (tensor densities of weight zero). The GRMHD evolution equations in coordinate basis are:

$$\begin{aligned} \partial_t D + \partial_i (f_D)^i &= -(f_D)^i \hat{\Gamma}_{ij}^j, \\ \partial_t S_j + \partial_i (f_S)^i_j &= (s_S)_j - (f_S)^i_j \hat{\Gamma}_{ik}^k + (f_S)^i_k \hat{\Gamma}_{ij}^k, \\ \partial_t \tau + \partial_i (f_\tau)^i &= s_\tau - (f_\tau)^i \hat{\Gamma}_{ij}^j, \\ \partial_t A_i &= \alpha \hat{\epsilon}_{ijk} \bar{v}^j \mathcal{B}^k - \hat{\mathcal{D}}_i (\alpha \Phi - \beta^j A_j), \\ \partial_t \hat{\Phi} + \partial_i (f_\Phi)^i &= -\zeta \alpha \hat{\Phi} - (f_\Phi)^i \hat{\Gamma}_{ij}^j, \\ \mathcal{B}^i &= \hat{\epsilon}^{ijk} \partial_j A_k. \end{aligned}$$

Before proceeding with the description of the GRMHD evolution equations expressed in the orthonormal basis with respect to the spherical background metric, we note that the geometric source terms introduced by rewriting the equations in the reference-metric formalism break the

¹⁰Like the Γ -driving shift parameter η , ζ has units of $1/t$ (or equivalently $1/M$ in $G = c = 1$ units). Thus, the value of ζ is constrained by the Courant-Friedrichs-Lewy (CFL) condition (in precisely the same way as described by [114] for η). However, unlike η , we prefer the damping provided by $\zeta > 0$ to be as strong as possible everywhere. Our choice $\zeta = 1.5/\Delta t$ is quite strong, but should be stable for CFL factors of $\approx \frac{2}{3}$ or smaller, consistent with the required CFL factors of $\frac{1}{2}$ or smaller when solving the BSSN/CCZ4 equations. As a corollary, when convergence testing $\zeta = 1.5/\Delta t$ must correspond to the lowest-resolution simulations Δt to ensure the CFL condition is not violated.

round-off level conservation of baryon number when evolving D in a finite volume scheme. This is due to the fact that the resulting finite volume scheme is not well-balanced [115] unless the geometric source terms are evaluated in such a way as to numerically exactly cancel the fluxes through the cell surface (see [33] for developments toward well-balanced schemes in GRMHD; an extension that is beyond the scope of this work). This nonconservation is a drawback of the scheme. For the conserved momenta S_i and conserved energy τ , this problem is less severe, as those quantities are only strictly conserved in the presence of spacetime symmetries [116], due to the appearance of (spacetime)-geometric source terms. In a sense, both problems are similar: both non-conservations arise from rewriting covariant derivatives in terms of partial derivatives which are suitable for the numerical integration of the resulting evolution equations.

When numerically evolving the Euler equation in spherical coordinates, there is nonconservation of momentum in the θ coordinate, which is due to the presence of the “naked pressure term” $e^{6\phi} \sqrt{\bar{\gamma}} P^* \delta_j^a$ which causes the breaking of zero-force equilibria. This is due to the θ dependence of the spherical background metric introduces a pressure gradient even in the absence of forces, as the finite volume scheme is not well balanced. One key advantage of our reference-metric approach is its automatic conservation of θ momentum due to the absence of the reference-metric determinant. We note that there are various other strategies to deal with this problem in spherical coordinates; see, e.g., [63,117,118]. The φ coordinate is not affected by this, and angular momentum is therefore identically conserved in spherical coordinates.

C. Equations in orthonormal basis of spherical background metric

Before we continue to describe the implementation of the evolution system in a finite volume method in Sec. II D below, we first make the following choice for the initial determinant of the conformally related metric:

$$\bar{\gamma} = \hat{\gamma} \quad \text{at } t = 0. \quad (109)$$

Moreover, we adopt the “Lagrangian choice” (9), $\partial_t \bar{\gamma} = 0$, meaning that we have $\bar{\gamma} = \hat{\gamma}$ at all times (as noted in Sec. II A, this is continuously enforced in the spacetime evolution). Accordingly, the ratio $\bar{\gamma}/\hat{\gamma}$ is unity and our definitions of the GRMHD variables therefore reduce to

$$D = e^{6\phi} \rho W, \quad (110)$$

$$S_j = e^{6\phi} (\rho h^* W^2 v_j - ab^0 b_j), \quad (111)$$

$$\tau = e^{6\phi} (\rho h^* W^2 - P^* - (ab^0)^2) - D, \quad (112)$$

$$\mathcal{B}^j = e^{6\phi} B^j. \quad (113)$$

Using the orthonormal basis with respect to the background metric defined above (13), we write the continuity equation as

$$\partial_t D + \partial_i (\hat{\mathbf{e}}_{\{j\}}^i (f_D)^{\{j\}}) = -(f_D)^i \hat{\Gamma}_{ij}^j. \quad (114)$$

Note that the vector (f_D) is expressed in both bases in Eq. (114). We evaluate the flux divergence by using the product rule and analytically differentiating the scale factors,

$$\partial_i (\hat{\mathbf{e}}_{\{j\}}^i (f_D)^{\{j\}}) = (\partial_i \hat{\mathbf{e}}_{\{j\}}^i) (f_D)^{\{j\}} + \hat{\mathbf{e}}_{\{j\}}^i \partial_i (f_D)^{\{j\}}, \quad (115)$$

so that the continuity equation is given by

$$\partial_t D + \hat{\mathbf{e}}_{\{j\}}^i \partial_i (f_D)^{\{j\}} = -(f_D)^i \hat{\Gamma}_{ij}^j - (\partial_i \hat{\mathbf{e}}_{\{j\}}^i) (f_D)^{\{j\}}. \quad (116)$$

The momentum equation similarly becomes

$$\begin{aligned} \partial_t (\hat{\mathbf{e}}_j^{\{i\}} S_{\{i\}}) + \hat{\mathbf{e}}_j^{\{k\}} \hat{\mathbf{e}}_{\{i\}}^i \partial_i ((f_S)_{\{k\}}^{\{i\}}) \\ = -\partial_i (\hat{\mathbf{e}}_j^{\{k\}} \hat{\mathbf{e}}_{\{i\}}^i) (f_S)_{\{k\}}^{\{i\}} - (f_S)_j^i \hat{\Gamma}_{ik}^k + (f_S)_k^i \hat{\Gamma}_{ij}^k + (s_S)_j. \end{aligned} \quad (117)$$

Before we write out the rest of the GRMHD equations, we note that the diagonality of the spherical background metric allows us to simplify the notation in expressions containing the basis elements. For instance, in expressions such as $\partial_t (\hat{\mathbf{e}}_j^{\{i\}} \hat{S}_{\{i\}})$, $\hat{\mathbf{e}}_j^{\{i\}} = 0$ if $i \neq j$, which leads to element-wise multiplication of vector and tensor components in the orthonormal basis and the basis elements. We could therefore write $\hat{\mathbf{e}}_j^{\{j\}} \hat{S}_{\{j\}}$ where the summation convention does not apply on repeated coordinate (j) and orthonormal ($\{i\}$) indices, but this can lead to significant confusion when interpreting equations. Instead, we introduce a new notation defining the following vectors and matrices of rescale factors $\hat{\mathcal{R}}_{\{i\}}$ (corresponding to the scale factors h_i of the reference metric):

$$\hat{\mathcal{R}}_{\{i\}} \equiv \begin{pmatrix} 1 \\ r \\ r \sin \theta \end{pmatrix}, \quad (118)$$

$$\hat{\mathcal{R}}^{\{i\}} = 1/\hat{\mathcal{R}}_{\{i\}}, \quad (119)$$

$$\hat{\mathcal{R}}_{\{i\}\{j\}} = \hat{\mathcal{R}}_{\{i\}} \hat{\mathcal{R}}_{\{j\}}, \quad (120)$$

$$\hat{\mathcal{R}}_{\{j\}}^{\{i\}} = \hat{\mathcal{R}}^{\{i\}} \hat{\mathcal{R}}_{\{j\}}. \quad (121)$$

As explained above, the components of the orthonormal basis form a diagonal matrix, and the rescaling amounts to element-wise multiplication of vector and tensor components with the corresponding vectors and matrices of rescaling factors (essentially computing the Hadamard product). In

order to write this in a notation that can be used with the summation convention that we use throughout, we define the following symbols that will be used to express coordinate vectors and tensors in the orthonormal basis with respect to $\hat{\gamma}_{ij}$:

$$\sigma_{\{j\}\{k\}}^i = \sigma_i^{\{j\}\{k\}} \equiv \begin{cases} 1 & i = j = k \\ 0 & \text{otherwise.} \end{cases} \quad (122)$$

Note that $\sigma_{\{j\}\{k\}}^i$ and $\sigma_i^{\{j\}\{k\}}$ are *not* tensors and that the indices in both are not raised and lowered by the metric. In addition, indices of the two σ matrices can be contracted with both coordinate and orthonormal indices (i.e., indices surrounded by curly braces). As an example of using this notation for vectors and tensors, the coordinate three-velocity is written as

$$v^i = \hat{\mathbf{e}}_{\{j\}}^i v^{\{j\}} = \sigma_{\{j\}\{k\}}^i \hat{\mathcal{R}}^{\{j\}} v^{\{k\}} \quad (123)$$

and the coordinate conformally related metric as

$$\begin{aligned} \bar{\gamma}_{ij} &= \hat{\mathbf{e}}_i^{\{k\}} \hat{\mathbf{e}}_j^{\{l\}} (\delta_{\{k\}\{l\}} + h_{\{k\}\{l\}}) \\ &= \sigma_i^{\{k\}\{m\}} \sigma_j^{\{l\}\{n\}} \hat{\mathcal{R}}_{\{m\}} \hat{\mathcal{R}}_{\{n\}} (\delta_{\{k\}\{l\}} + h_{\{k\}\{l\}}). \end{aligned} \quad (124)$$

In this notation, the continuity, momentum, and energy equations are written as

$$\begin{aligned} \partial_t D + \sigma_{\{j\}\{k\}}^i \hat{\mathcal{R}}^{\{j\}} \partial_i (f_D)^{\{k\}} \\ = -(f_D)^i \hat{\Gamma}_{ij}^j - \sigma_{\{j\}\{k\}}^i (\partial_i \hat{\mathcal{R}}^{\{j\}}) (f_D)^{\{k\}}, \end{aligned} \quad (125)$$

$$\begin{aligned} \partial_t S_{\{j\}} + \sigma_{\{k\}\{l\}}^i \delta_j^n \hat{\mathcal{R}}^{\{k\}} \partial_i ((f_S)^{\{l\}}_{\{n\}}) \\ = [-\sigma_{\{k\}\{l\}}^i \sigma_q^{\{m\}\{n\}} (f_S)^{\{k\}}_{\{m\}} (\partial_i \hat{\mathcal{R}}^{\{l\}}_{\{n\}}) \\ - (f_S)_q^i \hat{\Gamma}_{ik}^k + (f_S)_k^i \hat{\Gamma}_{iq}^k + (s_S)_q] \sigma_{\{j\}\{p\}}^q \hat{\mathcal{R}}^{\{p\}}, \end{aligned} \quad (126)$$

$$\begin{aligned} \partial_t \tau + \sigma_{\{j\}\{k\}}^i \hat{\mathcal{R}}^{\{j\}} \partial_i (f_\tau)^{\{k\}} \\ = -(f_\tau)^i \hat{\Gamma}_{ij}^j - \sigma_{\{j\}\{k\}}^i (\partial_i \hat{\mathcal{R}}^{\{j\}}) (f_\tau)^{\{k\}} + s_\tau. \end{aligned} \quad (127)$$

Noting that $\sqrt{\bar{\gamma}} \hat{\mathcal{R}}^{\{r\}} \hat{\mathcal{R}}^{\{\theta\}} \hat{\mathcal{R}}^{\{\varphi\}} = 1$, the evolution equation for the vector potential takes a particularly simple form in our notation

$$\partial_t A_{\{i\}} = -E_{\{i\}} - \sigma_{\{i\}\{k\}}^l \hat{\mathcal{R}}^{\{k\}} \partial_l (\alpha \Phi - \beta^{\{j\}} A_{\{j\}}). \quad (128)$$

The evolution equation of the EM scalar potential is given by

$$\begin{aligned} \partial_t \hat{\Phi} + \sigma_{\{j\}\{k\}}^i \hat{\mathcal{R}}^{\{j\}} \partial_i (f_\Phi)^{\{k\}} \\ = -\zeta \alpha \hat{\Phi} - (f_\Phi)^i \hat{\Gamma}_{ij}^j - \sigma_{\{j\}\{k\}}^i (\partial_i \hat{\mathcal{R}}^{\{j\}}) (f_\Phi)^{\{k\}}, \end{aligned} \quad (129)$$

and finally, the conserved rescaled magnetic field is calculated from

$$\begin{aligned} \mathcal{B}^{\{i\}} &= \sigma_l^{\{i\}\{m\}} \hat{\mathcal{R}}_{\{m\}} \hat{\mathbf{e}}^{ljk} \partial_j A_k \\ &= \sigma_n^{\{i\}\{q\}} \hat{\mathcal{R}}_{\{q\}} \hat{\mathbf{e}}^{nj k} \sigma_k^{\{l\}\{m\}} (A_{\{m\}} \partial_j \hat{\mathcal{R}}_{\{l\}} \\ &\quad + \hat{\mathcal{R}}_{\{l\}} \partial_j A_{\{m\}}). \end{aligned} \quad (130)$$

Introducing the following generalized sources:

$$\Omega_D \equiv -(f_D)^i \hat{\Gamma}_{ij}^j - \sigma_{\{j\}\{k\}}^i (\partial_i \hat{\mathcal{R}}^{\{j\}}) (f_D)^{\{k\}}, \quad (131)$$

$$\begin{aligned} (\Omega_S)_{\{j\}} &\equiv [-\sigma_{\{k\}\{l\}}^i \sigma_q^{\{m\}\{n\}} (f_S)^{\{k\}}_{\{m\}} (\partial_i \hat{\mathcal{R}}_{\{n\}}^{\{l\}}) \\ &\quad - (f_S)_q^i \hat{\Gamma}_{ik}^k + (f_S)_k^i \hat{\Gamma}_{iq}^k + (s_S)_q] \sigma_{\{j\}\{p\}}^q \hat{\mathcal{R}}^{\{p\}}, \end{aligned} \quad (132)$$

$$\Omega_\tau \equiv -(f_\tau)^i \hat{\Gamma}_{ij}^j - \sigma_{\{j\}\{k\}}^i (\partial_i \hat{\mathcal{R}}^{\{j\}}) (f_\tau)^{\{k\}} + s_\tau, \quad (133)$$

$$(\Omega_A)_{\{i\}} \equiv -E_{\{i\}} - \sigma_{\{i\}\{k\}}^l \hat{\mathcal{R}}^{\{k\}} \partial_l (\alpha \Phi - \beta^j A_j), \quad (134)$$

$$\Omega_\Phi \equiv -\zeta \alpha \hat{\Phi} - (f_\Phi)^i \hat{\Gamma}_{ij}^j - \sigma_{\{j\}\{k\}}^i (\partial_i \hat{\mathcal{R}}^{\{j\}}) (f_\Phi)^{\{k\}}, \quad (135)$$

we can write the reference-metric GRMHD evolution system in the following compact form suitable for the integration in a finite volume scheme:

$$\begin{aligned} \partial_t D + \sigma_{\{j\}\{k\}}^i \hat{\mathcal{R}}^{\{j\}} \partial_i (f_D)^{\{k\}} &= \Omega_D, \\ \partial_t S_{\{j\}} + \sigma_{\{k\}\{l\}}^i \delta_j^n \hat{\mathcal{R}}^{\{k\}} \partial_i ((f_S)^{\{l\}}_{\{n\}}) &= (\Omega_S)_j, \\ \partial_t S_{\{j\}} + \sigma_{\{k\}\{l\}}^i \hat{\mathcal{R}}^{\{k\}} \partial_i ((f_S)^{\{l\}}_{\{j\}}) &= (\Omega_S)_{\{j\}}, \\ \partial_t \tau + \sigma_{\{j\}\{k\}}^i \hat{\mathcal{R}}^{\{j\}} \partial_i (f_\tau)^{\{k\}} &= \Omega_\tau, \\ \partial_t A_{\{i\}} &= (\Omega_A)_{\{i\}}, \\ \partial_t \hat{\Phi} + \sigma_{\{j\}\{k\}}^i \hat{\mathcal{R}}^{\{j\}} \partial_i (f_\Phi)^{\{k\}} &= \Omega_\Phi, \\ \mathcal{B}^{\{i\}} &= \sigma_l^{\{i\}\{m\}} \hat{\mathcal{R}}_{\{m\}} \hat{\mathbf{e}}^{ljk} \partial_j A_k. \end{aligned}$$

We note that in Cartesian coordinates these equations reduce to the equations of the Valencia formulation in a vector potential evolution. In the following section, we will turn to the finite volume implementation of the evolution system above, and in particular how to deal with the rescaling factors multiplying divergences of fluxes in the orthonormal basis with respect to the spherical background metric.

D. Evolution equations in integral form

Systems of nonlinear hyperbolic partial differential equations (PDEs) such as the GRMHD evolution system presented above are characterized by the fact that smooth initial data can develop discontinuities in the variables in finite time. The reason the evolution system is written in conservative form is that, in such a form, a numerical

scheme that converges guarantees the correct Rankine-Hugoniot conditions across discontinuities, which is called the *shock-capturing* property. This property is at the heart of high-resolution shock-capturing (HRSC) methods that guarantee that the physics of the flow will be correctly modeled by the numerical scheme in the presence of discontinuities in the fluid variables.

Moreover, finite-difference schemes written in conservation form guarantee that the convergence of the solution (if it exists) will be to one of the *weak solutions* of the system of PDEs [119]. Weak solutions are characterized by being solutions to the integral form of the conservation system. The set of all weak solutions is too large to be of practical use, as many (numerically) admissible weak solutions will not represent physically relevant solutions. Thus, there is need for an additional (thermodynamic) condition, the so-called *entropy condition* (namely that the entropy of a fluid element must increase when crossing a discontinuity) to guarantee that the numerical scheme will converge to the physical solution. The convergence of the numerical scheme is closely related to its stability, and one useful measure is the *total-variation* stability (see, e.g., [120] for a detailed discussion).

Additionally, numerical schemes written in conservation form guarantee that the conserved quantities of the system are numerically conserved in the absence of sources or sinks. This means that the change of the state vector U_A in time in a domain \mathcal{V} that does not contain sources or sinks will be given by the fluxes F_A^i through the boundaries of the domain $\partial\mathcal{V}$, a three-dimensional surface which is defined as the standard-oriented hyper-parallelepiped consisting of two spacelike surfaces $\{\Sigma_{x^0}, \Sigma_{x^0+\Delta x^0}\}$ and the timelike surfaces $\{\Sigma_{x^i}, \Sigma_{x^i+\Delta x^i}\}$ joining the two temporal slices together.

In a finite volume formulation, the evolution equations are integrated over the cell volumes. For D , and similarly the other fluid variables, this amounts to the following integrals that give the update of a conserved quantity in a given cell (see, e.g., [101,121]):

$$\partial_t(\langle D \rangle \Delta V) + \int_{\Delta V} \sigma_{\{j\}\{k\}}^i \hat{\mathcal{R}}^{\{j\}} \partial_i (f_D)^{\{k\}} d^3x = \langle \Omega_D \rangle \Delta V. \quad (136)$$

Here $d^3x = drd\theta d\phi$, and we have defined

$$\langle D \rangle = \frac{1}{\Delta V} \int_{\Delta V} D d^3x, \quad (137)$$

$$\langle \Omega_D \rangle = \frac{1}{\Delta V} \int_{\Delta V} \Omega_D d^3x, \quad (138)$$

$$\Delta V = \Delta r \Delta \theta \Delta \phi. \quad (139)$$

Notice the absence of the spherical background metric determinant $\sqrt{\hat{\gamma}}$ in the above expressions, as all knowledge

about the spherical coordinates has been moved to the background metric Christoffel symbols in the cell-centered source term Ω_D , together with our choice of $\bar{\gamma} = \hat{\gamma}$.¹¹ Up until now, the integration of the evolution equations over the cell volumes is exact. Approximating the value of D and the source Ω_D inside the cell volumes as piecewise constant, and being equal to their cell-centered value (which is a second-order accurate approximation), we obtain

$$\partial_t \langle D \rangle + \frac{1}{\Delta V} \int_V \sigma_{\{j\}\{k\}}^i \hat{\mathcal{R}}^{\{j\}} \partial_i (f_D)^{\{k\}} d^3x = \langle \Omega_D \rangle. \quad (140)$$

The integral $\int_V \sigma_{\{j\}\{k\}}^i \hat{\mathcal{R}}^{\{j\}} \partial_i (f_D)^{\{k\}} d^3x$ is not a true divergence due to the appearance of the rescaling vector $\hat{\mathcal{R}}^{\{i\}}$, and therefore we cannot use the divergence theorem to convert the volume integral into a surface integral over the cell surface to arrive at finite volume scheme. We therefore make a third-order approximation, setting $\hat{\mathcal{R}}^{\{i\}}$ to be piecewise constant and equal to its value at the cell center, denoted as $\langle \mathcal{R}^{\{i\}} \rangle_{ijk}$ (where the subscript ijk denotes a cell), so that the volume integral can then be converted to a surface integral of the fluxes through the cell faces,

$$\begin{aligned} \sigma_{\{m\}\{n\}}^l \frac{\langle \mathcal{R}^{\{m\}} \rangle_{ijk}}{\Delta V} \int_V \partial_l (f_D)^{\{n\}} d^3x \\ = \sigma_{\{m\}\{n\}}^l \frac{\langle \mathcal{R}^{\{m\}} \rangle_{ijk}}{\Delta V} \int_S f_D^{\{n\}} s_l dA, \end{aligned}$$

where s_i is the outward pointing unit normal to the cell surface and dA the surface element. Therefore, all volume integrals in a finite volume scheme in the reference-metric formalism will be ‘‘Cartesian’’ in the sense that we integrate over Cartesian volumes and surfaces in the spherical grid using fluxes in the orthonormal basis with respect to $\hat{\gamma}_{ij}$. In our second-order accurate approximation, the surface integrals are given by

$$\frac{\langle \mathcal{R}^{\{r\}} \rangle_{ijk}}{\Delta V} \int_S f_D^{\{r\}} s_r dA = \frac{\langle \mathcal{R}^{\{r\}} \rangle_{ijk}}{\Delta r} \Xi_r, \quad (141)$$

$$\frac{\langle \mathcal{R}^{\{\theta\}} \rangle_{ijk}}{\Delta V} \int_S f_D^{\{\theta\}} s_\theta dA = \frac{\langle \mathcal{R}^{\{\theta\}} \rangle_{ijk}}{\Delta \theta} \Xi_\theta, \quad (142)$$

$$\frac{\langle \mathcal{R}^{\{\phi\}} \rangle_{ijk}}{\Delta V} \int_S f_D^{\{\phi\}} s_\phi dA = \frac{\langle \mathcal{R}^{\{\phi\}} \rangle_{ijk}}{\Delta \phi} \Xi_\phi, \quad (143)$$

¹¹A different strategy could have been followed here, namely, the integration of Eq. (66) in spherical coordinates directly using the generalized Stokes theorem. This approach is followed in the CoCoNuT code; see [102,105] for details.

where the symbols Ξ_i are defined as

$$\Xi_r \equiv ((f_D)^{\{r\}}_{i+\frac{1}{2},j,k} - (f_D)^{\{r\}}_{i-\frac{1}{2},j,k}), \quad (144)$$

$$\Xi_\theta \equiv ((f_D)^{\{\theta\}}_{i,j+\frac{1}{2},k} - (f_D)^{\{\theta\}}_{i,j-\frac{1}{2},k}), \quad (145)$$

$$\Xi_\varphi \equiv ((f_D)^{\{\varphi\}}_{i,j,k+\frac{1}{2}} - (f_D)^{\{\varphi\}}_{i,j,k-\frac{1}{2}}). \quad (146)$$

We can then write the second-order accurate finite volume evolution equation for D as

$$\partial_t \langle D \rangle + \frac{\langle \mathcal{R}^{\{r\}} \rangle_{ijk}}{\Delta r} \Xi_r + \frac{\langle \mathcal{R}^{\{\theta\}} \rangle_{ijk}}{\Delta \theta} \Xi_\theta + \frac{\langle \mathcal{R}^{\{\varphi\}} \rangle_{ijk}}{\Delta \varphi} \Xi_\varphi = \langle \Omega_D \rangle \quad (147)$$

and similar for the Euler and energy equation. The evolution equations for the vector potential (128) and the EM scalar potential (129) are treated differently: the cell-centered electric field is calculated averaging the reconstructed velocity and magnetic field at the surrounding cell faces [see Eq. (148) below], and the divergence in the EM scalar potential evolution equation is evaluated using finite differences. This means that the time integration of the conserved variables is given by the interface fluxes of matter and energy momentum of the fluid, as well as the (cell-centered) sources. One can therefore approximate those *numerical* fluxes (which depend on the solution at the cell interfaces) as the time-averaged fluxes across cell interfaces during a time step. In general, the approximation to the real solution on a grid with finite resolution will be a piecewise continuous function, which means that the fluxes can be obtained by solving local Riemann problems at cell interfaces, an idea first described by Godunov [122].

Riemann problems are initial value problems with discontinuities in the solution. During the evolution, a discontinuity in the fluid variables decays into shock waves, rarefaction waves, and contact discontinuities. Shock waves move from the higher to lower density regions, while rarefaction waves move in the opposite direction. Contact discontinuities are characterized by a discontinuity in the density, while both pressure and velocity are constant across them. In order to solve the Riemann problem, we need to obtain the spectrum (eigenvalues and eigenvectors) of the first-order system. The fluid data at the cell interfaces needed to obtain the numerical fluxes via the solution of local Riemann problems need to be obtained from the cell averages. A wide variety of higher-order cell-reconstruction methods are available in the literature (see, e.g., [123]). Regardless of their spatial order for smooth solutions, these reconstruction techniques always reduce to first order in the presence of physical shocks and some reconstruction schemes even reduce to first order at local extrema of the fluid variables (such as the central density of an NS, for instance).

While the choice of variables is crucial to obtain the GRMHD evolution equations in conservative form, it is usually the primitive variables that are reconstructed at the cell interfaces. To do this, one needs a conservative-to-primitive scheme, involving numerical root finding. Once we have obtained the numerical fluxes via the solution of local Riemann problems, we update the solution of the conserved variables by one time step with the numerical fluxes and the sources. This is usually done employing high-order Runge-Kutta schemes [124].

From the structure of the equations in integral form, as noted above, we see that they are Cartesian by virtue of having written the conservation laws in the reference-metric formalism. The second-order accuracy is achieved by encoding all the geometric information about the underlying coordinate system in the cell-centered geometric source terms. Specifically, no care has to be taken to distinguish the coordinate center and centroid of volume of the computational cells. In general curvilinear coordinates, this is not the case, as replacing the average value of a cell quantity with a point value is only second-order accurate if the point is chosen to be the centroid of volume, not the coordinate center [125,126]. To this end, the second-order accurate prescription outlined could be applied to any existing Cartesian finite volume code by calculating the appropriate reference-metric source terms and incorporating them in the time integration of the evolution equations.

In the following section, we describe the necessary changes we performed to enable the use of spherical coordinates via the GRMHD reference-metric formalism presented above in GRHydro [28,127–129], a publicly available GRMHD code that comes with the Einstein Toolkit.

III. IMPLEMENTATION IN THE EINSTEIN TOOLKIT

The Einstein Toolkit [82] is an open source code suite for relativistic astrophysics simulations. It uses the modular Cactus framework [130] (consisting of general modules called “thorns”) and provides adaptive mesh refinement via the Carpet driver [131–133]. In our vacuum implementation of the BSSN equations [81], we have detailed how we enabled the use of spherical coordinates in the Einstein Toolkit, having supplied our own spacetime evolution thorn.

Enabling spherical coordinates in GRHydro to arrive at a GRMHD code in spherical coordinates in the Einstein Toolkit amounted to supplying a different metric determinant and the appropriate reference-metric source terms, so the changes to the existing code are minimally invasive and do not touch core algorithms of GRHydro. Perhaps the most substantial change involved using the NRPy+ code [134,135] to replace the Cartesian GRMHD source terms in GRHydro with the generalized source terms (131)–(135). NRPy+ converts these expressions—written in human-readable,

Einstein notation—into optimized C-code kernels, automatically constructing finite-difference derivatives at arbitrary order when needed. `NRPY+` was also used to fully construct the C-code kernels for Einstein's equations written in both BSSN (as described in [81]) and fCCZ4 formalisms.

At the interface between the spacetime and GRMHD evolution, the (physical) spacetime variables $\alpha, \beta^i, \gamma_{ij}$, and K_{ij} of the `ADMBase` thorn are passed to `GRHydro`. As outlined above, rewriting the equations to evolve the noncoordinate components of vectors and dividing every $\sqrt{\gamma}$ by $\sqrt{\hat{\gamma}}$ would amount to a great deal of rescaling and unrescaling (both vector components and determinants) in `GRHydro`. Instead, we follow a different route and pass the noncoordinate basis metric $\gamma_{\{i\}\{j\}} = e^{4\phi}(\delta_{\{i\}\{j\}} + h_{\{i\}\{j\}})$ and shift $\beta^{\{i\}}$ as the `ADMBase` variables in substeps of the method of lines integration. This means that $\sqrt{\det \gamma_{\{i\}\{j\}}} = e^{6\phi}$ and raising and lowering indices of noncoordinate vectors are achieved with the noncoordinate basis metric $v_{\{i\}} = \gamma_{\{i\}\{j\}} v^{\{j\}}$.

As outlined above, given that the evolution equations in integral form are Cartesian, the different reconstruction methods that are available in `GRHydro` may be used *without modification*. These include total variation diminishing (TVD) with minmod, superbee [136] and monotized central [137] limiters, the piecewise parabolic method (PPM) [138] and its enhanced version that retains higher order at smooth extrema [139,140], monotonicity-preserving fifth order (MP5) reconstruction [141], essentially nonoscillatory reconstruction [142], as well as weighted essentially nonoscillatory reconstruction (WENO [143] and its variant WENO-Z [144]). Using these reconstruction algorithms without modification would have been impossible had the code been written in spherical coordinates without the reference-metric formalism, as greater care must be taken when using these in spherical coordinates, especially for higher-order reconstruction methods (see, e.g., [126]). We use the HLLC (Harten-Lax-van Leer-Einfeldt) approximate Riemann solver [145,146] present in `GRHydro`, again without any changes to its implementation to calculate the numerical fluxes through cell faces.

In order to achieve magnetic flux conservation, the cell-centered electric field $E_{\{i\}}$ used in the update of the vector potential is calculated as the average of the nonzero magnetic fluxes given by the HLLC solver [107],

$$\begin{aligned} (E_{\{1\}})_{i,j,k} &= -\frac{1}{4}(\alpha v^{\{2\}} \mathcal{B}^{\{3\}} - \alpha v^{\{3\}} \mathcal{B}^{\{2\}})_{i,j-\frac{1}{2},k} \\ &\quad -\frac{1}{4}(\alpha v^{\{2\}} \mathcal{B}^{\{3\}} - \alpha v^{\{3\}} \mathcal{B}^{\{2\}})_{i,j+\frac{1}{2},k} \\ &\quad +\frac{1}{4}(\alpha v^{\{3\}} \mathcal{B}^{\{2\}} - \alpha v^{\{2\}} \mathcal{B}^{\{3\}})_{i,j,k-\frac{1}{2}} \\ &\quad +\frac{1}{4}(\alpha v^{\{3\}} \mathcal{B}^{\{2\}} - \alpha v^{\{2\}} \mathcal{B}^{\{3\}})_{i,j,k+\frac{1}{2}} \end{aligned}$$

and similarly for $E_{\{2\}}$ and $E_{\{3\}}$. As shown in [147,148], the cell-centered vector potential method we employ is identical to evolving the induction equation directly with the so-called flux-CD scheme [107], as the magnetic field is cell centered and the curl of the gradient $\hat{D}_i(\alpha\Phi - \beta^j A_j)$ in the RHS of A_i is zero.

To mitigate high-frequency oscillations in the cell-centered vector potential evolution, we add Kreiss-Oliger dissipation [149] to the RHSs of both A_i and $\hat{\Phi}$ [147,150].

One of the most delicate parts of GRMHD codes is the recovery of the primitive variables, which usually requires nonlinear inversion. `GRHydro` uses the conservative to primitive routines scheme of [40]. Some of the most problematic regimes for the inversion are in regions of very high Lorentz factors and in magnetically dominated plasmas, i.e., where $b^2/(2P) \gg 1$. In those regions, the inversion errors may become comparable to the truncation error and result in larger errors causing the evolution to fail eventually. The biggest problem is the violation of physical constraints such as the positivity of ρ and P during the recovery, as in this case the hyperbolicity of the evolution equations breaks down [151]. As is customary in GRMHD codes, we use a tenuous atmosphere, given that the GRMHD evolution equations break down in true vacuum. The atmosphere region is particularly difficult to handle, as even very small magnetic fields can result in very large ratios of magnetic to fluid pressure. We use the following checks prior to primitive recovery:

- (1) In cells where $D < e^{6\phi} \rho_{\text{atm}}$, reset the cell to atmosphere assuming a zero magnetic field (ignoring the contribution of the magnetic field to τ) and skip the primitive recovery. The magnetic field is fully evolved in the atmosphere and always calculated from the curl of the vector potential.
- (2) Following [26], when a BH is present, in regions where $b^2/(2P)$ is greater than a user-specified threshold, we raise the above criterion to reset to atmosphere, which avoids primitive inversion in cells that are just above the atmosphere threshold. Effectively, this results in a higher-density atmosphere, but in regions limited to high magnetic to fluid pressure ratios, while allowing the use of a low-density atmosphere in regions of small magnetic fields. This is important as a denser atmosphere can begin to affect the evolved physical system of interest [152].
- (3) Following [29], we check, after primitive recovery, whether ρ , P , and W exceed user-specified limits, and, if so, reset the primitives and then recalculate the conservatives.
- (4) Once an apparent horizon (AH) is found, we reset a small region deep inside the AH to atmosphere. In all the above steps, the magnetic field is never altered in any computational cell and always computed from the vector potential.

In future versions of the code, problems related to the primitive recovery could be handled by more modern algorithms, such as evolving the entropy S and using it to recover the pressure [42], or using different primitive recovery schemes; see [153] for an overview. Another attractive approach could be the use of physical-constraint-preserving methods [151,154,155].

A well-known problem of evolving hyperbolic PDEs in spherical coordinates is the severe CFL limitation due to the nonconstant cell volumes in space, which become smaller (therefore leading to smaller time steps) as the origin and axis are approached. There are several approaches to mitigate this problem (for an introduction, see, e.g., [156]), from various *multipatch* approaches which remove the polar axis [140,157–167], mesh coarsening in the azimuthal direction at high latitudes [168], radially dependent mesh coarsening in both polar and azimuthal angles [169], mesh coarsening as a conservative filter operation [170–173], focusing resolution of the polar angle at the equator [174,175], or the use of filters [176–178].

In order to avoid excessively small time steps in full three-dimensional (3D) simulations, we employ a radial and latitude dependent azimuthal fast Fourier transform (FFT) filter (using the FFTW3 library [179]) that will be described in detail in a companion publication [180]. In short, we expand all evolved fields in the azimuthal direction in a Fourier series and retain m modes such that the time step at the pole is limited by $n_\varphi = 6$ points. Higher-order m modes in the expansion are exponentially damped, which is sufficient to prevent instability by violating the CFL condition.

We use the boundary condition thorn described in [81]. In summary, internal boundary ghost zones for the r boundary at the origin, and the θ and φ boundaries are copied in from points in the physical domain, accounting for appropriate parity factors, which we list for rescaled and unrescaled vector and tensor components in Table I for completeness. The ghost zone to physical point mappings are as follows:

- (i) r boundary at the origin,

$$r \rightarrow -r, \quad \theta \rightarrow \pi - \theta, \quad \varphi \rightarrow \varphi + \pi,$$

- (ii) θ boundary at $\theta_{\min} = 0$,

$$r \rightarrow r, \quad \theta \rightarrow -\theta, \quad \varphi \rightarrow \varphi + \pi,$$

- (iii) θ boundary at $\theta_{\max} = \pi$,¹²

$$r \rightarrow r, \quad \theta \rightarrow 2\pi - \theta, \quad \varphi \rightarrow \varphi + \pi.$$

¹²We note there is a typo in the θ mapping of the $\theta_{\max} = \pi$ boundary in [81], the correct mapping is the one shown here.

TABLE I. Parity factors for rescaled and coordinate vector and tensor components at the origin and polar axis. The parity factors for contravariant components are the same as for the covariant components shown.

	Origin	Axis
$V_{\{r\}}$	–	+
$V_{\{\theta\}}$	+	–
$V_{\{\varphi\}}$	–	–
V_r	–	+
V_θ	–	–
V_φ	+	+
$T_{\{r\}\{r\}}$	+	+
$T_{\{r\}\{\theta\}}$	–	–
$T_{\{r\}\{\varphi\}}$	+	–
$T_{\{\theta\}\{\theta\}}$	+	+
$T_{\{\theta\}\{\varphi\}}$	–	+
$T_{\{\varphi\}\{\varphi\}}$	+	+
T_{rr}	+	+
$T_{r\theta}$	+	–
$T_{r\varphi}$	–	+
$T_{\theta\theta}$	+	+
$T_{\theta\varphi}$	–	–
$T_{\varphi\varphi}$	+	+

Finally, ghost zones for φ are set by imposing periodicity. We note that our boundary condition requires an even number of grid points in the φ direction in order to ensure that ghost zones lie at the exact locations of points in the physical domain.

IV. CODE TESTS

In the following, we show results for a series of code tests, ranging from special relativistic test problems in a fixed background Minkowski spacetime to fully dynamical spacetime evolutions of magnetized stable uniformly rotating neutron stars and the collapse of a magnetized uniformly rotating neutron star to a Kerr BH [181].

A. Tests in Minkowski spacetime

The first set of tests is performed in a fixed background Minkowski spacetime ($\{h_{\{i\}\{j\}} = \bar{A}_{\{i\}\{j\}} = 0, K = \Theta = 0, \beta^{\{i\}} = B^{\{i\}} = \tilde{\Lambda}^{\{i\}} = 0, \alpha = e^{4\phi} = 1\}$), allowing us to compare the performance of the MHD evolution with standard Newtonian tests. This enables us to validate our implementation of the GRMHD evolution equations with all metric terms set to flat space. To demonstrate the code is working correctly in this setting, we show two tests below. First, we solve a strong shock reflection problem without magnetic fields (evolving pure HD problems by simply setting the vector potential to zero everywhere initially). The second, more demanding test is an explosion test problem. As shown below, these tests are do not

exploit symmetries of the spherical coordinate system, demonstrating that the framework can, e.g., handle the passage of strong shocks through the origin and polar axis.

1. Relativistic spherical shock reflection test

Our first test is the relativistic spherical shock reflection problem [182–186]. The test consists of an initially cold ($\epsilon \approx 0$) fluid of unit density ($\rho = 1$) flowing in uniformly with a velocity of $v^r = v_{\text{in}} = -0.9$ toward the origin, where the fluid is compressed and heated up resulting in a shock that travels upstream through the inflow region. For numerical reasons, the problem is initialized with a small pressure of $p = 2.29 \times 10^{-5}(\Gamma - 1)$, where we use $\Gamma = 4/3$. The analytic solution to this problem is given by [182]

$$\rho(r) = \begin{cases} (1 + |v_{\text{in}}|t/r)^2 & r > v_s t \\ (1 + |v_{\text{in}}|/v_s)^2 \sigma & r < v_s t, \end{cases} \quad (148)$$

where the compression factor σ and the shock velocity v_s are given by

$$\sigma = \frac{\Gamma + 1}{\Gamma - 1} + \frac{\Gamma}{\Gamma - 1} (W_{\text{in}} - 1) \quad (149)$$

and

$$v_s = \frac{\Gamma - 1}{W_{\text{in}} + 1} W_{\text{in}} |v_{\text{in}}| \quad (150)$$

and where W_{in} is the Lorentz factor of the inflowing fluid at the outer boundary (≈ 2.29 for $v_{\text{in}} = -0.9$).

Behind the shock wave ($r < v_s t$), the fluid is at rest ($v^r = 0$), and internal energy is given by $\epsilon = W_{\text{in}} - 1$.

For this test, we used 800 radial points in the interval [0:1] and two points in the θ and φ directions, using the HLLE Riemann solver, TVD reconstruction with a Minmod limiter, and CFL factor of 0.4. At the outer radial boundary, ρ is set to the analytic solution (148), and v^r and p are kept fixed at their initial values. Figure 1 shows the radial profiles of ρ , P , and v^r of the numerical and analytic solution at $t = 4$.

The global relative error at $t = 4$ is 2.2%, 2.1% and 1.2%, for ρ , P , and v^r , respectively. In the density profile, a significant drop near the origin is present. This numerical effect is known as wall heating [187] and seems to be exacerbated in spherical coordinates due to the converging grid geometry [188].

We also note that we observed significant postshock oscillations behind the slowly moving shock when using higher-order reconstruction methods. This appears to be a known problem for HRSC schemes (see, e.g., [189] and references therein).

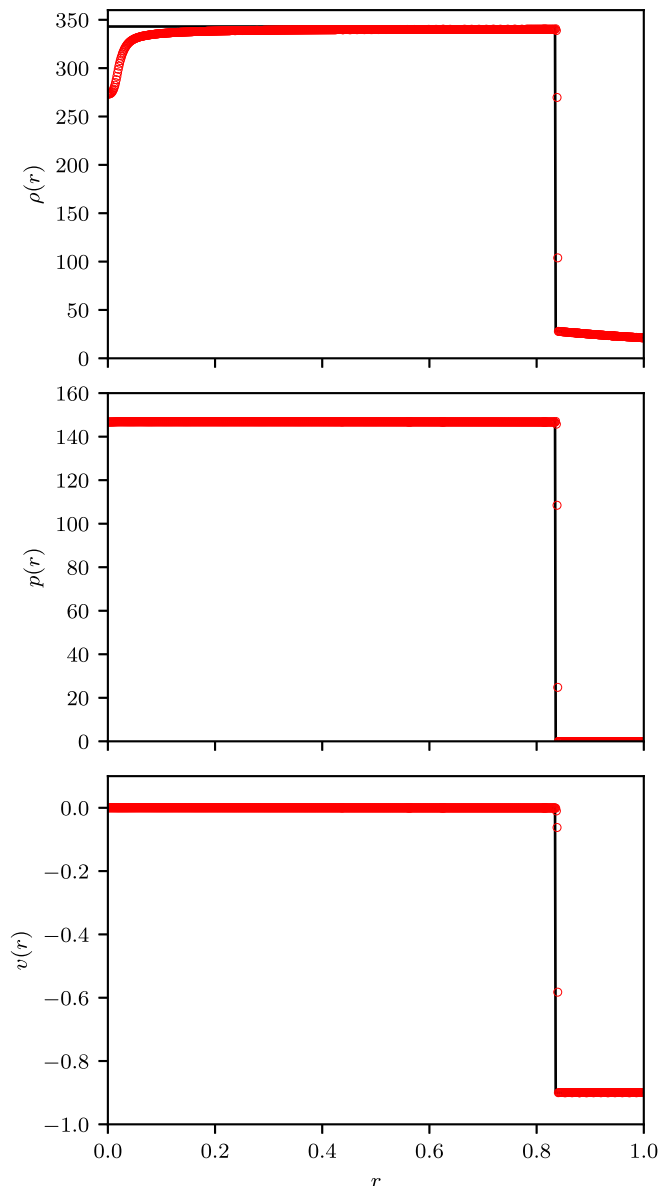


FIG. 1. Radial density, pressure, and velocity profiles (from top to bottom) of the relativistic spherical shock reflection problem at time $t = 4$. The red circles correspond to the numerical solution, while the analytic solution is shown as solid black lines.

2. Spherical explosion

Next, we test the relativistic MHD evolution with a magnetized spherical explosion problem [26]. This test is the natural extension of the cylindrical explosion test proposed in [190] to spherical coordinates, setting up spherically symmetric initial data. Using the same jump conditions as [190], the initial data consist of an overdense ($\rho = 1 \times 10^{-2}$, $p = 1.0$) ball of radius 1.0. From a radius of 0.8 onward, the solution is matched in an exponential decay to the surrounding medium ($\rho = 1 \times 10^{-4}$, $p = 3 \times 10^{-5}$). The initial pressure profile in the $y = 0$ plane is shown in the top left panel of Fig. 2. We are using a Γ – law equation of state (EOS) using $\Gamma = 4/3$. In the magnetized case, the

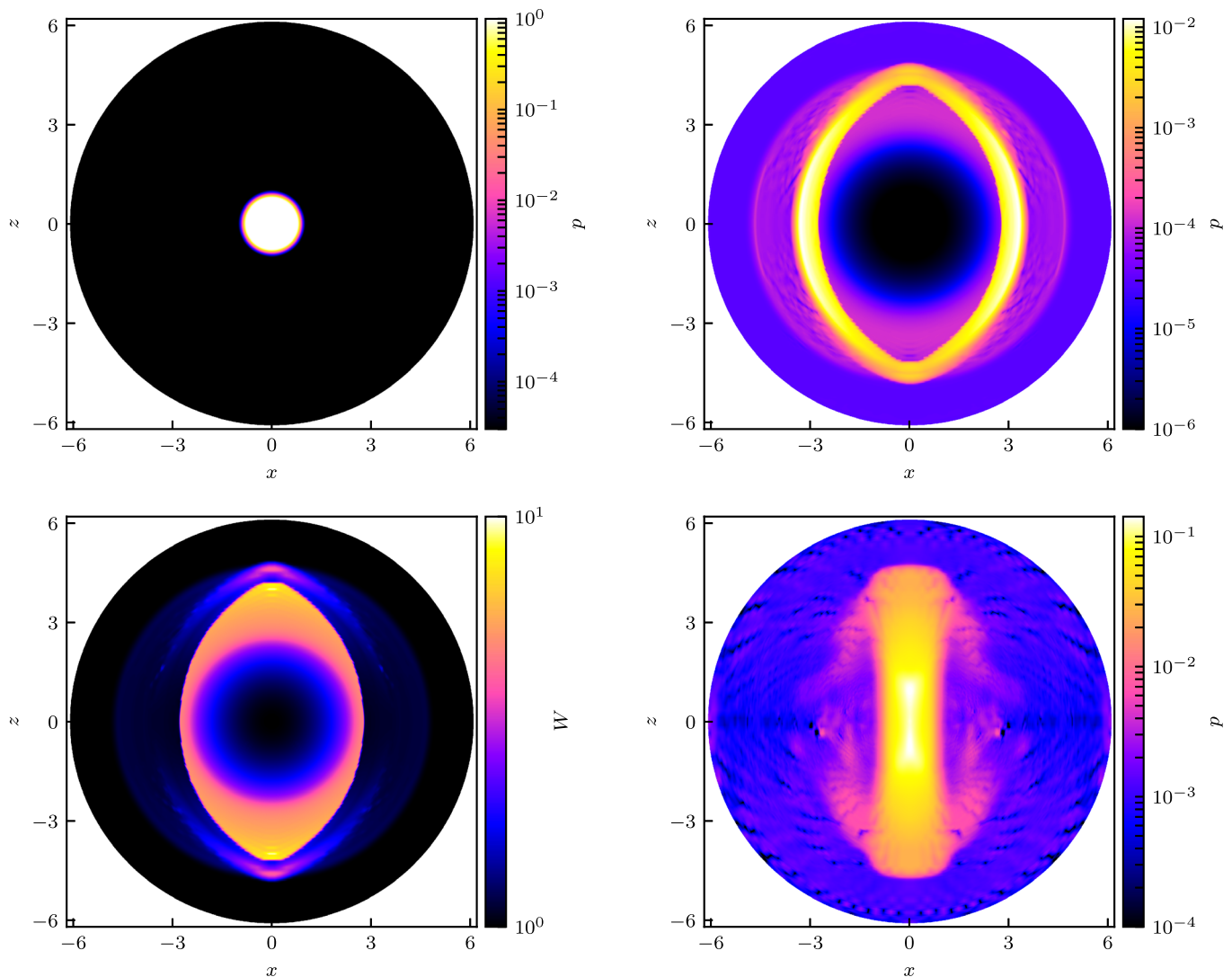


FIG. 2. Snapshots of magnetized spherical explosions in axisymmetry. Top left: initial pressure profile. Top right: pressure at $t = 4$, HLLC Riemann solver, and MP5 reconstruction. Bottom left: Lorentz factor at $t = 4$, HLLC Riemann solver, and MP5 reconstruction. Bottom right: pressure at $t = 4$, global Lax-Friedrichs fluxes, and TVD reconstruction, initial $B^z = 1.0$.

entire domain is initially threaded by a constant magnitude magnetic field parallel to the z axis ($B^z = 0.1$), and the fluid velocity is set to zero everywhere in the domain initially. As in the relativistic shock reflection problem, we use fixed background Minkowski spacetime for this test problem.

We first model a magnetized spherical explosion in axisymmetry, using $(n_r = 160, n_\theta = 80, n_\phi = 4)$ points, with the outer boundary $r_{\max} = 6.0$. We use the HLLC Riemann solver and different reconstruction methods for this test. The final distributions at $t = 4$ for the pressure P and Lorentz factor W in the $y = 0$ plane are shown in the top right and bottom panels of Fig. 2.

In the initially overdense explosion region, the fluid is only weakly magnetized, while being strongly magnetized in the ambient medium. This results in a rich flow morphology in which the fast magnetosonic wave travels out ahead in spherical symmetry at almost the speed of

light, while the Alfvén wave shows a $\cos \theta$ dependence in propagation speed, traveling close to the speed of light parallel to the initial magnetic field, while being significantly slowed down in the direction perpendicular to the magnetic field, shown in the top right panel of Fig. 2 (see the discussion in [26]).

During the explosion, the magnetic field is expelled from the initial explosion region, leaving a spherically symmetric low-density region behind in which the fluid is at rest, as evidenced by the plot of the Lorentz factor in the bottom left panel of Fig. 2. The results seem to be in very good qualitative agreement with the results presented in [26] (spherical coordinates and axisymmetry) and [34] (Cartesian coordinates).

As a final axisymmetric test, we perform the same explosion, but with an initial magnetic field of much larger strength $B^z = 1.0$. This is a very demanding test for which

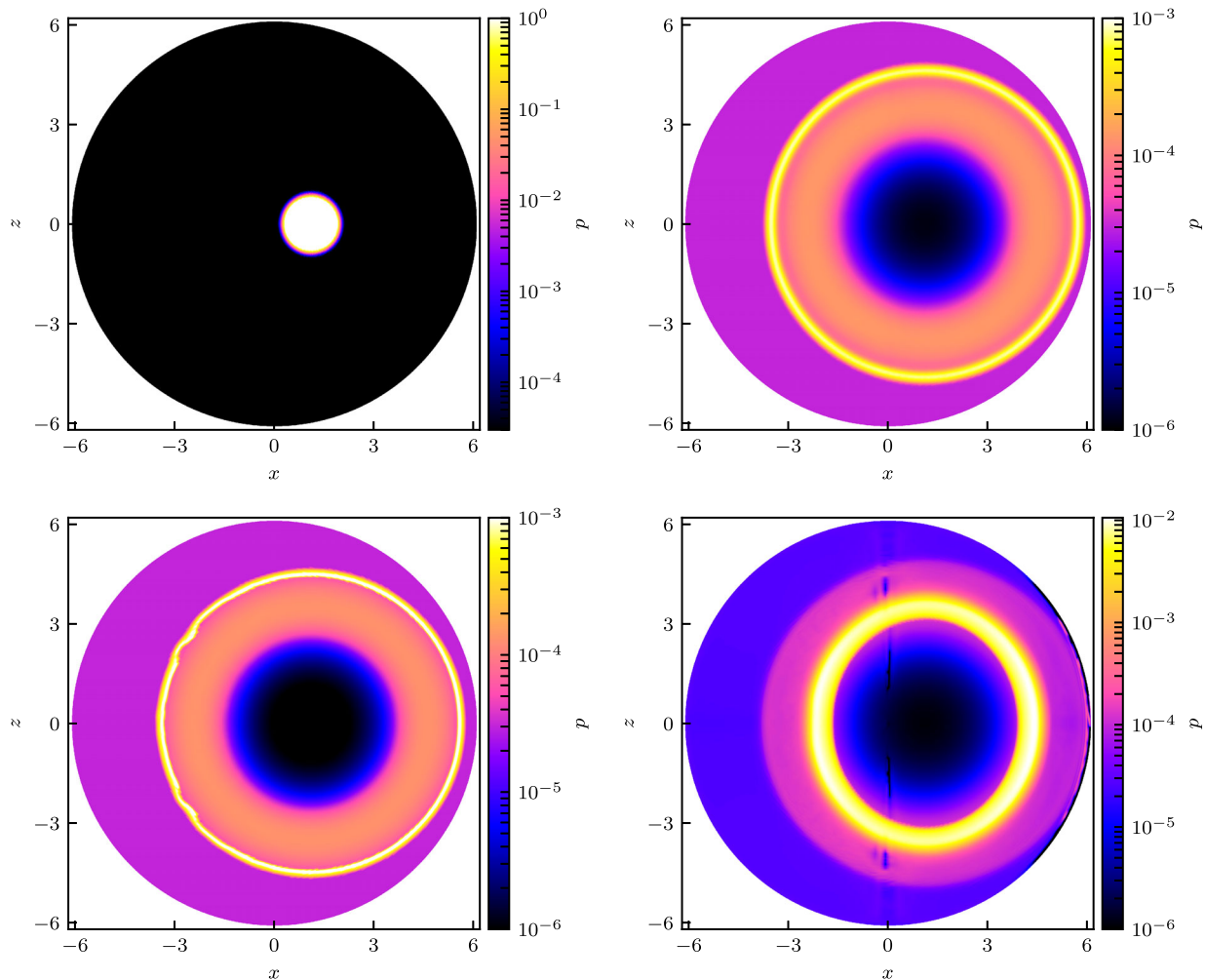


FIG. 3. Snapshots of off-centered spherical explosions. Top left: initial pressure profile. Top right: pressure at $t = 4$, HLLC Riemann solver, and TVD reconstruction, $B^i = 0$. Bottom left: pressure at $t = 4$, HLLC Riemann solver, and MP5 reconstruction, $B^i = 0$. Bottom right: pressure at $t = 4$, global Lax-Friedrichs fluxes, and TVD reconstruction with initial magnetic field $B^z = 0.1$ rotated by 45° about the x axis.

we have used global Lax-Friedrichs fluxes and TVD reconstruction with the Minmod limiter as they are more diffusive. In the bottom right panel of Fig. 2, we plot the pressure distribution at $t = 4$ in the $y = 0$ plane for this test. The morphology of the explosion changes completely and becomes bar shaped, as seen in Cartesian simulations of magnetized cylindrical and spherical explosions (see, e.g., [34,191]). Compared to the more weakly magnetized explosion, more noise can be seen in the ambient region. As discussed in [191], this test is most strenuous on the conservative to primitive solver, so we believe that the noise is due to inversion failures.

Next, we test the code by modeling an *off-center* spherical explosion, both in relativistic HD and relativistic MHD. As this test does not exploit the symmetries of our spherical coordinate system, we perform it in full 3D. The initial data are identical to the axisymmetric test described above, except the center of the explosion region has been

moved to $(x = 1.1, y = 0, z = 0)$. The resulting initial pressure profile in the $y = 0$ plane is shown in the top left panel of Fig. 3. In addition, compared to the axisymmetric explosion, the initial magnetic field $B^z = 0.1$ has been tilted by 45° about the x axis. This results in initial data that do not reflect the symmetries of the spherical coordinate system at all. For this full 3D test, we use $(n_r = 160, n_\theta = 80, n_\phi = 160)$ points, and use the azimuthal FFT filter to increase the time step to what it would have been, had the simulation been performed with $n_\phi = 6$ points.

We first perform two tests setting the magnetic field to zero initially, using TVD and MP5 reconstructions. The results for the two different reconstruction schemes are shown in the top right and bottom left panels of Fig. 3, respectively. When using TVD, there are no visible artifacts arising from the (in hydro only) spherically symmetric shock passing through the origin and axis. The rarefaction region is seen to be spherically symmetric as well, showing

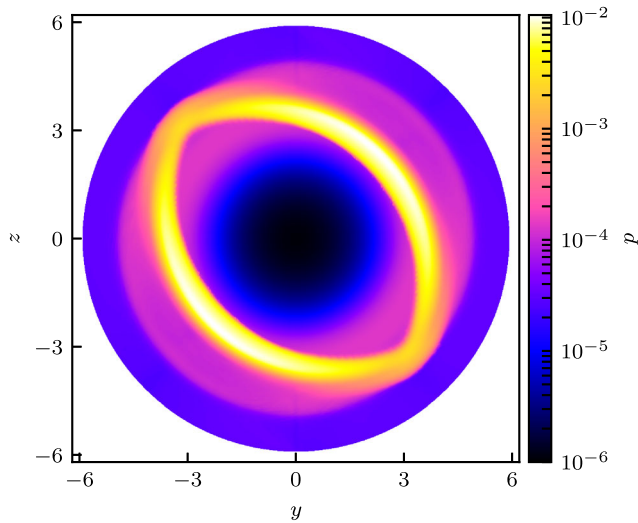


FIG. 4. Pressure distribution in spherical off-centered explosion with a magnetic field initially tilted by 45° about the x axis. The plane shown is the y - z plane centered at the initial center of the explosion region ($x = 1.1, y = 0, z = 0$). Simulation performed with global Lax-Friedrichs fluxes and TVD reconstruction with Minmod limiter.

no artifacts. When using MP5, the shock width is clearly reduced compared to TVD reconstruction, demonstrating a superior capture of the shock with the higher-order reconstruction, but small artifacts in those parts of the shock that have passed the origin and polar axis can be seen. A similar test in hydro is presented in [169]. The bottom right panel of Fig. 3 shows the final pressure distribution for the off-centered, tilted magnetic field spherical explosion, which was performed using global Lax-Friedrichs fluxes and TVD reconstruction with a Minmod limiter. This test displays more pronounced effects of the magnetized shock passing through origin and polar axis, showing primitive recovery failures at the polar axis. The global morphology is captured well nevertheless.

It is worth noting that the numerical artifacts that can be seen in the bottom panels of Fig. 3 are not concerning at all. The test setup was deliberately chosen to push the code to its limits by not exhibiting any (approximate) symmetries the framework was designed for and is a difficult test to pass even for Cartesian codes [34].

Finally, in Fig. 4, we plot the final pressure distribution in the $x = 1.1$ plane, i.e., the plane vertically cutting through the initial center of the explosion region. The flow morphology observed in the axisymmetric case (top right panel of Fig. 2) is clearly seen to be present and tilted by 45° , which is precisely the symmetry axis picked out by the tilted magnetic field initially threading the computational domain.

As explained in Sec. II B 4 above, the conserved rest-mass density D is not conserved to round off in our framework due to the appearance of geometric source

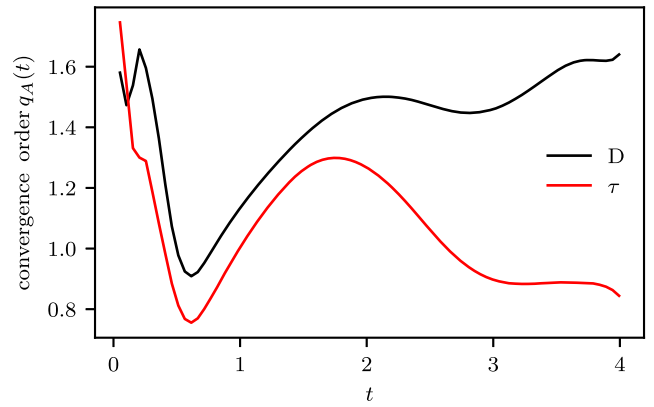


FIG. 5. Convergence order in conservation of D and τ for the spherical off-centered explosion with a magnetic field initially tilted by 45° about the x axis. See main text for details.

terms in the evolution equation for D (67). Furthermore, the FFT filter we employ to circumvent the severe CFL limit in full 3D simulations is inherently nonconservative as well. We therefore check for the convergence of the total rest mass and total energy conservation (s_τ , the source term in the evolution equation for τ (76) vanishes in Minkowski spacetime so that τ should be exactly conserved as well). To do so, we calculate $\epsilon_A(t)$, the volume integrated error arising from nonconservation of a quantity $A = (D, \tau)$ at time t as

$$\epsilon_A(t) = \int_{\Sigma} (A(t) - A(0)) \sqrt{\hat{\gamma}} dr d\theta d\phi \quad (151)$$

and calculate the convergence order $q_A(t)$ as [192]

$$q_A(t) = \frac{1}{\ln(f)} \ln \left(\frac{\|\epsilon_A(t)\|_{\text{low}}}{\|\epsilon_A(t)\|_{\text{high}}} \right), \quad (152)$$

where f is the ratio between the different resolutions used in the convergence test. We calculate $q_A(t)$ in the most demanding variant of the spherical explosion, the off-centered explosion with a tilted initial magnetic field, using two different resolutions of $(n_r, n_\theta, n_\phi) = (112, 56, 112)$ and $(160, 80, 160)$, corresponding to $f \approx \sqrt{2}$ and show the time evolution of $q_A(t)$ in Fig. 5. In our numerical scheme, we would expect the convergence order to be between 1 and 2, as our method is second-order accurate while reducing to first order in the presence of shocks. The convergence order of the conservation of total rest mass lies within that region, while for the total energy it drops below first order at the time the numerical artifacts at the polar axis seen in the bottom right panel of Fig. 3 start appearing. The maximum relative errors in the conservation of total rest mass and total energy in the high-resolution test are 0.0016 and 0.004, respectively.

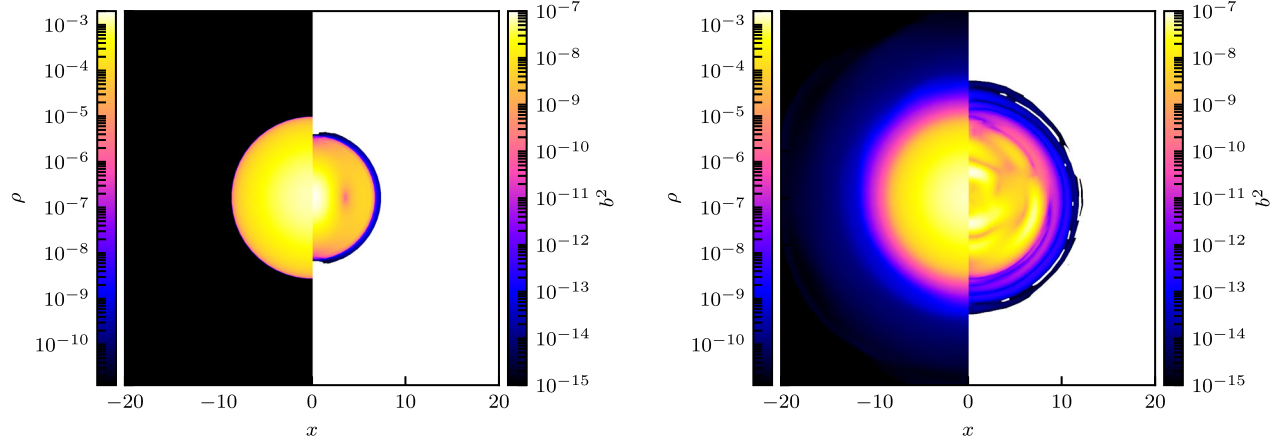


FIG. 6. ρ and b^2 evolution of tilted model BU2, shown at $t = 0$ (left) and $t = 25$ ms (right).

B. Dynamical spacetime tests

Next, we turn to dynamical spacetime evolutions of uniformly rotating neutron stars, testing the coupled spacetime and GRMHD evolution of the SphericalNR framework. To test the framework in this regime, we evolve two uniformly rotating polytropes, models BU2 [193] and D1 [127], adding a weak poloidal magnetic field initially. We perform tests of two important scenarios: The long-term evolution of a stable equilibrium model, as well as the gravitational collapse of a uniformly rotating polytrope to a BH. In the long-term evolution of model BU2, we initially tilt the star's rotation axis by 90° in order to test the evolution in full 3D without symmetry assumptions. In this test, the fluid rotates through the polar axis during the entire

TABLE II. Main properties of the relativistic polytrope models BU2 [193] and D1 [127]: central rest-mass density ρ_c , rest- and gravitational masses M_0 and M , the dimensionless angular momentum J/M^2 , the circumferential stellar radius R , the ratio of polar and equatorial radii of the star r_p/r_e , the ratio of kinetic energy and gravitational binding energy $T/|W|$, the adiabatic index Γ , the polytropic constant K , and the constants prescribing the initial magnetic field A_b , n_s , and P_{cut} (see main text for details).

	BU2	D1
ρ_c	1.28×10^{-3}	3.28×10^{-3}
M_0	1.58	1.83
M	1.47	1.67
J/M^2	3.19×10^{-1}	2.07×10^{-1}
R	10.11	7.74
r_p/r_e	0.9	0.95
$T/ W $	2.44×10^{-2}	1.17×10^{-2}
Γ	2	2
K	100	100
A_b	2	1
n_s	0	0
P_{cut}	6.55×10^{-6}	4.25×10^{-6}

simulation, dragging the magnetic field with it through the polar axis constantly. In the second test, we perform a simulation of the gravitational collapse of model D1, testing all aspects of our framework: the correct coupled evolution of the fCCZ4 and GRMHD equations leading to dynamical BH formation and ringdown to a Kerr BH.

1. Tilted, magnetized, uniformly rotating neutron star

The initial data are generated with the RNS code [194], which has been incorporated as a thorn named `Hydro_RNSID` in the Einstein Toolkit. As with all original Cartesian thorns present in the Einstein Toolkit, in order to interface with SphericalNR we need to coordinate transform the Cartesian initial data generated by the `Hydro_RNSID` code and then rescale the evolved fields. In order to test the FFT filter applied to both spacetime and GRMHD fields and nontrivial dynamics in full 3D, we initially tilt the rotation axis of the neutron star by 90° about the x axis so that the star's rotation axis is initially aligned with the y axis. After generating the tilted fluid and spacetime data, we add a small initial magnetic field, following the vector-potential-based prescription of [19]

$$A_r = 0, \quad (153)$$

$$A_\theta = 0, \quad (154)$$

$$A_\varphi = A_b (r \sin \theta)^2 \left(1 - \frac{\rho}{\rho_c}\right)^{n_s} \max(P_{\text{cut}} - P, 0), \quad (155)$$

$$\Phi = 0, \quad (156)$$

where values of A_b , ρ_c , n_s , and P_{cut} are provided in Table II. With this setup, the tilted, uniformly rotating star will constantly drag the magnetic field through the polar axis during the evolution. While the initial data are polytropic,

we evolve the star with a Γ -law EOS. We use a third-order strong stability-preserving Runge-Kutta (SSPRK3) method [124,195],¹³ implemented in the `MOL` thorn [129], the HLLC Riemann solver, and WENO-Z reconstruction for the simulations presented here. We have also tried ePPM and MP5 reconstruction, but found a large symmetry breaking at late times when using those. To check for convergence of our code, we evolve three different resolutions for 29 ms, which is more than 200 dynamical timescales of the star.

The $t = 0$ and $t = 25$ ms distributions of ρ and b^2 in the $y = 0$ plane are shown in Fig. 6. The star remains very stable and very contained, and there are no large outflows from the stellar surface into the atmosphere, demonstrating the code's capability to deal with the stellar surface. This is a difficult test, as the numerical dissipation at the stellar surface is minimal in spherical coordinates, due to the fact that the surface and computational cell surfaces are mostly aligned (in Cartesian coordinates, this effect is seen along the coordinate axes; see, e.g., Fig. 3 in [197]). During the evolution, the quantity b^2 develops a richer morphology than it has in the beginning, which we believe results from the fact that the initial poloidal field evolves into having poloidal and toroidal components (while the initial data are uniformly rotating, the misalignment between the star's rotation axis and the initial magnetic field dipole axis results in the generation of a toroidal magnetic field.). In order to quantify the error arising from the FFT filter and to check for the resolution dependence in the radial and angular coordinates, we plot the following diagnostics in Fig. 7: the relative error in central density and total rest mass in the top two panels, as well as the evolution of the L^2 norm of the Hamiltonian constraint in the bottom panel.

Additionally, in Fig. 8, we show a convergence study with three different resolutions, now increasing the resolution twice by a factor of $f = 1.5$ in all three coordinate directions. The top panel shows the evolution of the error in the central density evolution, where the medium and higher resolution errors have been multiplied by 1.8 and 3.375, respectively, assuming a convergence order of 1.5. Our code is formally second order, while reducing to first order in the presence of shocks. The surface of the neutron star is a discontinuity, so we would expect the order of convergence to be between 1 and 2. The bottom panel shows the convergence factor for the total rest mass M_0 (158), the ADM mass evaluated as a volume integral, and the L1 norm of the Hamiltonian. As in Sec. IV A 2, we calculate the convergence order for M_0 and M_{ADM} using (152), while the convergence order of the L1 norm of the Hamiltonian is calculated as [192]

¹³Strong stability-preserving time discretization methods have been called TVD methods historically [196].

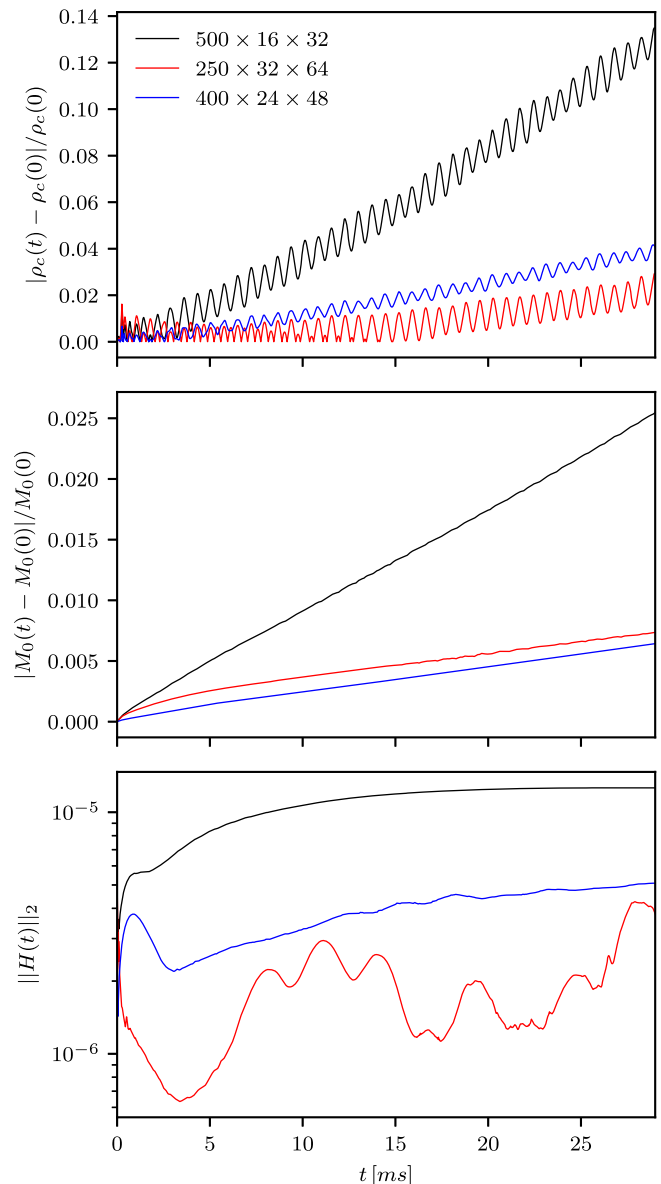


FIG. 7. Evolution of central density ρ_c (top panel), total rest mass M_0 (mid panel), and L^2 -norm of the Hamiltonian constraint (bottom panel) for magnetized model BU2. Three different resolutions are shown in each plot.

$$q_A(t) = \frac{1}{\ln(f)} \ln \left(\frac{\|H(t)\|_{\text{low}} - \|H(t)\|_{\text{med}}}{\|H(t)\|_{\text{med}} - \|H(t)\|_{\text{high}}} \right). \quad (157)$$

As the initial magnetic field is small, and the tilted rotation axis of the star should not affect its dynamics during evolution, we calculate the frequencies of oscillations in $\rho_{\text{max}}(t)$ as a power spectral density for the two higher resolution runs in Fig. 9. We overlay the expected frequencies of the fundamental quasiradial (F) and quadrupolar 2f modes and their first overtones (H_1 and 2p_1) (taken from Table 5 of [198]). In both resolutions, the fundamental modes F and 2f are in very good agreement

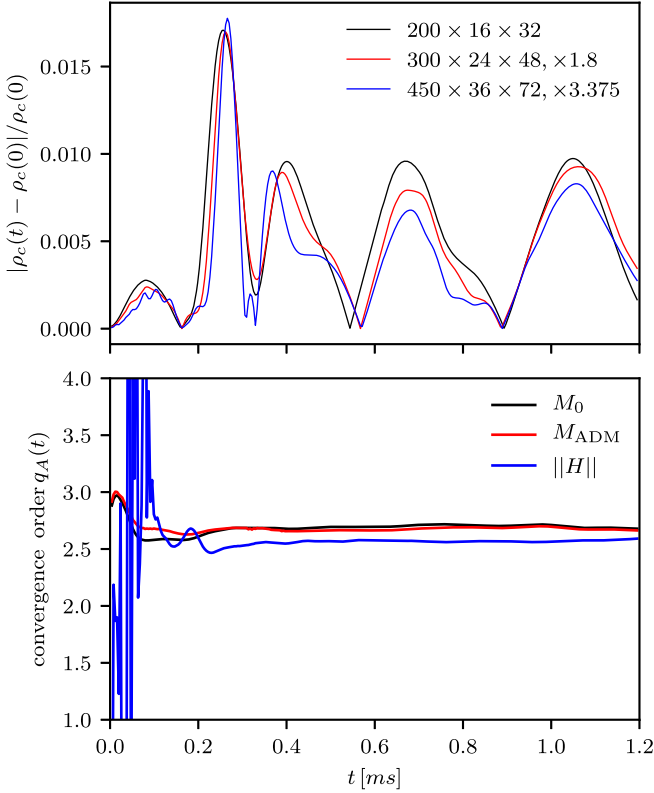


FIG. 8. Convergence study for magnetized model BU2. Top panel: relative error in $\rho_c(t)$ at multiple resolutions; medium- and high-resolution results are multiplied by the appropriate factor assuming a convergence order of 1.5. Bottom panel: convergence order for the total rest mass M_0 , ADM mass M_{ADM} , and L1 norm of the Hamiltonian constraint.

with [198]; however, the first overtones (H_1 and 2p_1) of both fundamental modes are not visible in the simulation with lower radial but higher angular resolution, while being slightly shifted in the higher radial resolution simulation.

These tests show that our code is capable of evolving equilibrium neutron stars with magnetic fields for many timescales, in a setup (constant fluid motion and magnetic field dragging through the polar axis) that was chosen to be particularly challenging for our framework.

2. Collapse of a magnetized uniformly rotating neutron star

As our last test, we present a very important test problem for numerical relativity simulations with matter: the collapse of a neutron star to a black hole (see, e.g., [127,140,199–202]). Using the `Hydro_RNSID` code thorn in the `Einstein Toolkit` again, we setup the uniformly rotating polytrope model D1 [127] with a weak poloidal magnetic field added initially and evolve its collapse to a Kerr BH. The initial data specifications of this model are listed in Table II. The simulation is performed in axisymmetry, and the collapse is induced by lowering the

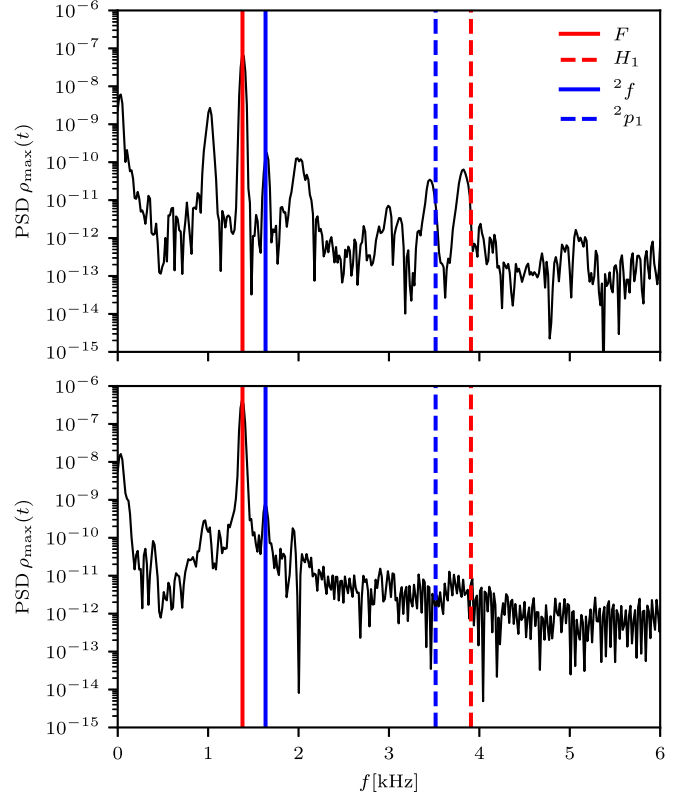


FIG. 9. Power spectral density of $\rho_{\text{max}}(t)$ evolution for resolutions of $(400 \times 24 \times 48)$ (top) and $(250 \times 32 \times 64)$ (bottom) panel. The vertical lines indicate the fundamental (F) and first overtone (H_1) of the fundamental quasiradial ($l=0$) mode and the fundamental (2f) and first overtone (2p_1) of the fundamental quadrupolar ($l=2$) mode (see Table 5 of [198]).

polytropic constant K everywhere in the star by 2% initially. The simulation is performed using $(n_r = 10000, n_\theta = 32, n_\phi = 2)$ points, with the outer boundary placed at $r_{\text{out}} = 500$. We evolve the conformal factor $W = e^{-2\phi}$ [203], use the `SSPRK3` method for time integration, and the `fCCZ4` damping parameters are set to $\kappa_1 = 0.06$, $\kappa_2 = 0$, $\kappa_3 = 1$. We use `WENO-Z` reconstruction, the `HLLC` Riemann solver, and a Γ -law EOS with $\Gamma = 2$.

The atmosphere value for ρ is set to be 10^{-8} times the initial density maximum ($\rho_{\text{min}} = 3.28 \times 10^{-11}$). This simulation requires the use of the higher atmosphere threshold in highly magnetized regions (described in Sec. III above), as the collapsing fluid leaves a highly magnetized atmosphere region behind. In these regions, we reset a cell to atmosphere if $e^{6\phi}\rho W < 100e^{6\phi}\rho_{\text{min}}$.

We use the `AHfinderDirect` thorn [204,205] to find the AH [206] once it has formed during collapse and the `QuasiLocalMeasures` thorn [207,208] to calculate the angular momentum of the AH during the evolution. The `SphericalNR` interface to these Cartesian diagnostic thorns in the `Einstein Toolkit` is described in [81]. The BH spin is measured using a surface integral on the AH [207] or the flat space rotational Killing vector method [209,210].

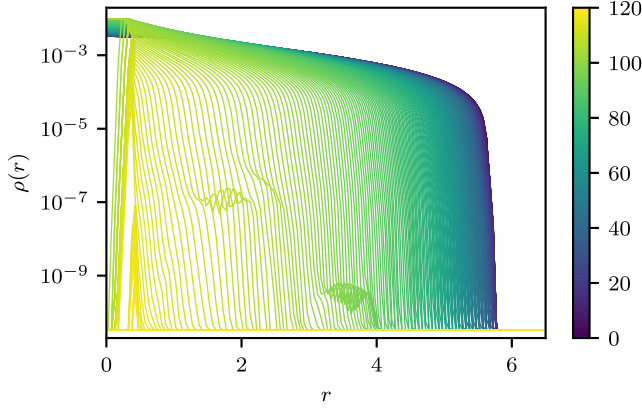


FIG. 10. Temporal evolution of radial profiles of ρ during the collapse of model D1, where the color bar indicates simulation time.

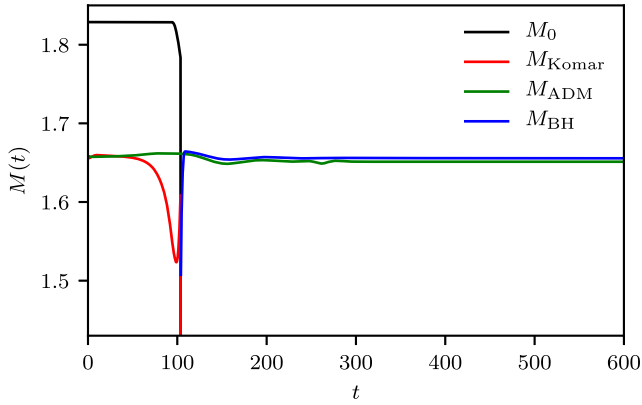


FIG. 11. Time evolution of various mass measurements during the collapse of model D1.

Deep inside the horizon, for points with coordinate radii $r < 0.2 \min(r_{\text{AH}})$, we set fluid variables to atmosphere values (the magnetic fields are evolved everywhere). Figure 10 shows the time evolution of radial profiles (along $\theta = d\theta/2$) for the density ρ . The evolution of ρ shows our modifications to the conservative to primitive solve in GRHydro: as the star collapses, ρ is first capped to a maximum value and then reset to atmosphere deep inside the AH once it has been found (AHFinderDirect reports the first finding of an AH at $t = 103.8$).

In Fig. 11, we plot different mass measurements for the duration of the simulation. Specifically, we monitor the total rest mass,

$$M_0 = \int_{\Sigma_t} \rho W e^{6\phi} \sqrt{\hat{\gamma}} d^3x, \quad (158)$$

the Komar mass [211] evaluated as a volume integral (see, e.g., [103]),

$$M_{\text{Komar}} = \int_{\Sigma_t} (\alpha(E + S) - 2\beta^i S_i) e^{6\phi} \sqrt{\hat{\gamma}} d^3x \quad (159)$$

and the ADM mass [212] of the spacetime, evaluated as the sum from contributions inside a finite radius r_{in} , evaluated as a surface integral, and those outside r_{in} , evaluated as a volume integral [213],

$$\begin{aligned} M_{\text{ADM}} = & \frac{1}{16\pi} \oint_{r_{\text{in}}} (\Delta\Gamma^r - 8e^\phi \bar{D}^r \phi) \sqrt{\hat{\gamma}} d\theta d\varphi \\ & + \frac{1}{16\pi} \int_{r>r_{\text{in}}} [e^{5\phi} (16\pi E + \bar{A}_{ij} \bar{A}^{ij} - \frac{2}{3} K^2) \\ & - \Delta\Gamma^{ijk} \Delta\Gamma_{jik} + (1 - e^\phi) \bar{R}] \sqrt{\hat{\gamma}} dr d\theta d\varphi. \end{aligned} \quad (160)$$

Comparing with the expressions found in [213] terms containing $\Delta\Gamma_{jk}^k$ are missing in the above expression, this is due to the fact that

$$\Delta\Gamma_{jk}^k = \frac{1}{\sqrt{\hat{\gamma}}} \hat{D}_j \sqrt{\hat{\gamma}}, \quad (161)$$

which, together with our choice of $\bar{\gamma} = \hat{\gamma}$, results in $\Delta\Gamma_{jk}^k = 0$. Finally, we compute the BH mass calculated as the Christodoulou mass [214],

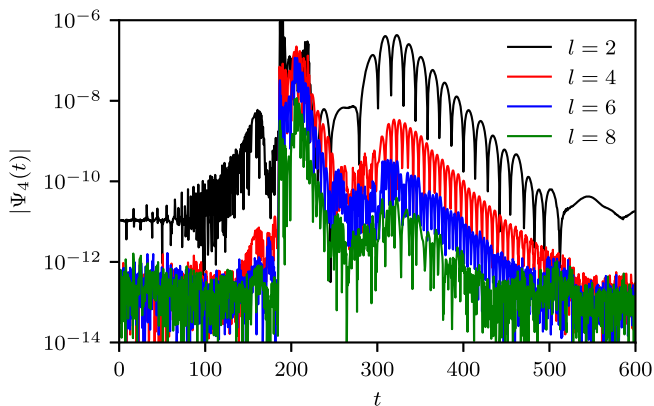
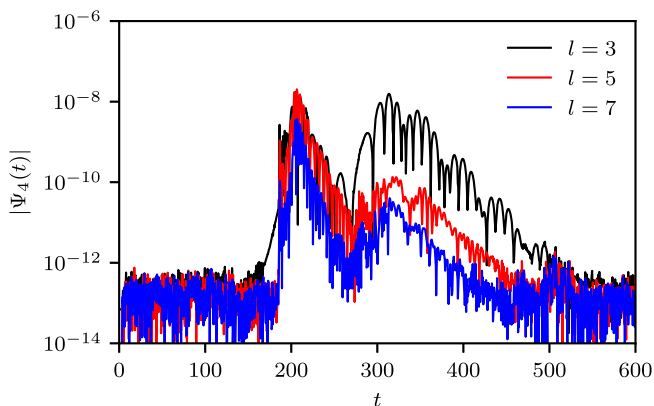
$$M_{\text{BH}} = \left(M_{\text{irr}}^2 + \frac{4\pi J^2}{A} \right)^{\frac{1}{2}}, \quad (162)$$

where M_{irr} is the BH irreducible mass, J the BH angular momentum, and A the AH area.

The mass measurements agree well with their initial value of the equilibrium neutron star. As the collapse proceeds, the total rest mass M_0 is seen to drop when the density deep inside the star is capped (see Fig. 10) and then quickly drops to zero once an AH has been found, as we exclude points within the horizon from volume integrals. The same drop is observed in the calculation of M_{Komar} , which also exhibits a stronger deviation from its initial value earlier, due to the fact that it is only defined for stationary spacetimes, and the collapse is an inherently dynamical process. Toward the end of the simulations, the calculation of the ADM mass via a surface integral shows oscillations related to the gravitational radiation leaving the domain and being partially reflected at the outer boundary.

After collapse, the newly formed BH is expected to quickly settle down to a Kerr BH via the ringdown of the BH's quasinormal modes (for a review, see [215]). To see if our simulation reproduces this expected behavior, in Figs. 12 and 13, we plot the $l = 2$ to 8 , $m = 0$ modes of the Weyl scalar Ψ_4 , split into even and odd l modes, respectively. The ringdown of all modes is clearly seen, as well beatings in the higher-order modes, whose origin (the equal m -mode mixing of spherical and spheroidal harmonics) we have explained in [81].

The simulation shows that our spherical GRMHD code is capable of capturing the relevant dynamics of the collapse of the magnetized uniformly rotating neutron star


 FIG. 12. Even $l = 2$ to 8 , $m = 0$ modes of the Weyl scalar Ψ_4 .

 FIG. 13. Odd $l = 3$ to 7 , $m = 0$ modes of the Weyl scalar Ψ_4 .

to a Kerr BH, capturing the postcollapse ringdown to Kerr with very high accuracy, with all modes dropping down to their initial background amplitudes.

V. CONCLUSIONS AND OUTLOOK

We have extended our vacuum numerical relativity code in spherical coordinates within the `Einstein Toolkit` [81] to a framework that numerically solves the coupled fCCZ4/BSSN and GRMHD equations in spherical coordinates without symmetry assumptions using a reference-metric formalism.

Extending the existing spacetime evolution thorn `SphericalBSSN` to evolve the fCCZ4 system with constraint damping as well enables future users of the framework with two distinct evolution systems for numerical relativity in spherical coordinates. The spacetime evolution thorn was written from scratch using `NRP+`, while the implementation of the reference-metric formalism GRMHD equations derived in this work was built as an extension of the `GRHydro` thorn, again using `NRP+`.

In our approach, the GRMHD equations in spherical coordinates acquire a Cartesian form, as all information about the underlying spherical coordinate system is

encoded in source terms of the equations. This has allowed us to use many of the core Cartesian building blocks of the HRSC finite volume implementation already present in `GRHydro` without modifications (and will enable the straightforward inclusion of Cartesian finite volume building blocks such as more sophisticated Riemann solvers in the future). Without the reference-metric approach, these building blocks would need to be adapted to spherical coordinates. Further, instead of evolving the magnetic field directly, the framework evolves the cell-centered vector potential in the generalized Lorenz gauge, guaranteeing the absence of magnetic monopoles to round-off error during the evolution by calculating the magnetic field as the curl of the vector potential.

We have tested our framework performing a set of demanding tests in flat background as well as fully dynamical spacetimes. We have chosen setups where the symmetries of the fluid are not aligned with the symmetries of the coordinate system (counter to our original motivation for developing the code). These tests include off-centered magnetized spherical explosions testing the passage of shocks and rarefaction waves through the coordinate origin and polar axis, as well as dynamical-spacetime simulations of a uniformly rotating neutron star with its rotation axis misaligned with the polar axis of the computational grid. Finally, we have shown that the code is able to perform simulations of the collapse of a magnetized uniformly rotating neutron star to a Kerr BH.

The `SphericalNR` framework will be made public and proposed to be included in a future official release of the `Einstein Toolkit`.

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