

Control of Physical Parameters in Binary Black Hole Initial Data

Iago Mendes^{1,2} 

Mentors: Nils Vu¹ , Mark Scheel¹ , Saul Teukolsky^{1,3} 

¹Theoretical Astrophysics 350-17, California Institute of Technology, Pasadena, CA 91125, USA

²Department of Physics and Astronomy, Oberlin College, Oberlin, Ohio 44074, USA

³Cornell Center for Astrophysics and Planetary Science, Cornell University, Ithaca, New York 14853, USA

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Abstract

When solving Einstein’s equations for Binary Black Holes (BBH) numerically, we split the four-dimensional spacetime into three-dimensional spatial slices and evolve them over time. The first slice is described by solving an Initial Data problem, which is cast as five elliptic partial differential equations in the extended conformal thin-sandwich (XCTS) decomposition. Prior to solving the XCTS equations, we must choose “free data” to impose boundary conditions and specify background quantities. This is the approach taken in **SpECTRE**, a parallel code that aims to simulate BBH for the new generations of gravitational wave detectors. After solving the XCTS system, **SpECTRE** runs a horizon finder that measures the masses and spins of the black holes. Additionally, our work extends **SpECTRE**’s capabilities to calculate total energy, momentum, and center of mass as infinite integrals using the Arnowitt-Deser-Misner (ADM) formalism. Typically, we want to choose these physical quantities before running a BBH simulation, but they can only be measured after numerically solving the XCTS equations. To address this, we implemented an iterative scheme in **SpECTRE** that drives the physical parameters to their desired values by adjusting the free data in a quasi-Newton-Raphson method that efficiently computes the Jacobian with Broyden’s method.

1. Introduction

In the early twentieth century, Einstein revolutionized the study of gravity by connecting spacetime geometry with physical dynamics. As John Wheeler says, “Spacetime tells matter how to move; matter tells spacetime how to curve” [1]. Being a highly complex theory, many problems of interest only have analytic solutions in special cases with symmetry. In this context, Numerical Relativity emerged as an essential field to solve these problems numerically, allowing us to explore general cases that can be found in

the universe. Specifically, simulations of Binary Black Holes (BBH) became very important as gravitational wave detectors were developed, needing to use numerical results to identify and characterize signals in their data [2].

Famously, the Einstein equations relate the curvature of spacetime to the stress-energy of matter, forming a system of ten nonlinear partial differential equations (PDEs). With the 3+1 formalism, we can rearrange these equations so that spacetime is described by spacelike three-dimensional slices of constant time [3]. In doing so, we find that four out of the ten equations do not involve time derivatives, implying that they are constraints that must be satisfied at all times. The remaining six equations describe an evolution of the constraint-satisfying fields. Using this formalism, the Spectral Einstein Code (SpEC) [4] runs BBH simulations by first finding initial data and then running an evolution on them. Over time, as SpEC faced more challenging BBH with high mass ratios and spins, several improvements had to be made to the initial data techniques, which are summarized in [5].

Despite its success in BBH simulations, SpEC shows its limitations in more challenging problems, such as binary neutron star mergers and BBH with extreme configurations. In this context, SpECTRE [6] was created as a codebase that follows a better parallelism model and aims to be more scalable [7]. Previous work has already shown that SpECTRE can be faster and more accurate than SpEC when performing similar tasks due to its use of parallelism [8]. This will be especially needed for the upcoming gravitational wave detectors with higher sensitivity, such as the Cosmic Explorer, the Einstein Telescope and LISA.

As part of an effort to allow researchers to fully simulate BBH in SpECTRE, an initial data procedure similar to the one in SpEC needs to be completed. This is greatly benefitted by a scalable elliptic solver that was recently developed [8], which can now be used to solve the initial data equations. Be that as it may, before the start of the SURF program, SpECTRE did not have a way to enforce specific masses and spins for the black holes or to avoid drifts in their orbital trajectories. This report describes how we addressed these issues.

2. Methods

2.1. Free data

To find initial data, SpECTRE uses the extended conformal thin-sandwich (XCTS) decomposition (see, e.g., [9] for a review), which transforms the constraint PDEs into a system of five elliptic PDEs. Before solving the XCTS equations, we need to specify the conformal spatial metric $\bar{\gamma}_{ij}$, the extrinsic curvature trace K , and their respective time derivatives $\partial_t \bar{\gamma}_{ij}$ and $\partial_t K$ [9]. There are many methods for specifying these quantities, but the one that has shown to be the most promising for BBH simulations is the superposed Kerr-Schild (SKS) approach [10], which enforces quasiequilibrium conditions by setting $\partial_t \bar{\gamma}_{ij} = 0$ and $\partial_t K = 0$ and specifies $\bar{\gamma}_{ij}$ and K by superposing two analytic solutions of Kerr-Schild black holes with masses M_A^{Kerr} and M_B^{Kerr} and with dimensionless spins $\vec{\chi}_A^{\text{Kerr}}$ and $\vec{\chi}_B^{\text{Kerr}}$.

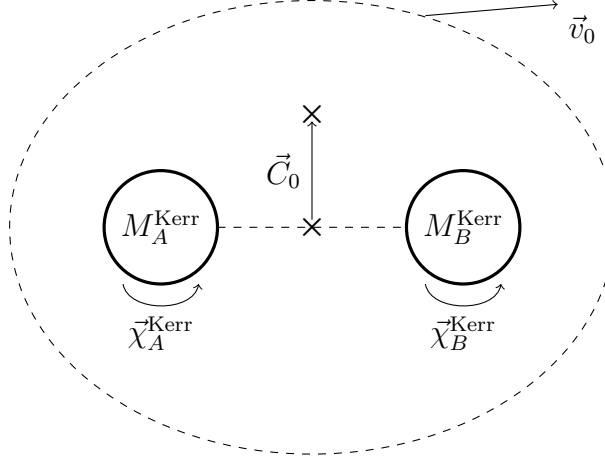


Figure 1: Schematic representation of the BBH free data.

We also need to choose the centers of the black holes \vec{c}_A and \vec{c}_B , which we set so that the Newtonian center of mass,

$$\vec{C}_{\text{Newtonian}} = \frac{M_A^{\text{Kerr}} \vec{c}_A + M_B^{\text{Kerr}} \vec{c}_B}{M_A^{\text{Kerr}} + M_B^{\text{Kerr}}}, \quad (1)$$

is zero and their relative distance $\vec{c}_A - \vec{c}_B = \vec{D}$, where \vec{D} is chosen when setting up the simulation. To give us more control of the placement of the black holes, we add a Newtonian center of mass offset \vec{C}_0 .

Having $\bar{\gamma}_{ij}$, K , $\partial_t \bar{\gamma}_{ij}$ and $\partial_t K$ specified, we can use the elliptic solver on the XCTS equations, three of which will be solved for the shift β^i [9]. Similar to any elliptic PDE problem, we have to set boundary conditions before solving these equations. At the domain outer boundary (with radius R), we can add a constant velocity \vec{v}_0 to the boundary conditions of the shift:

$$\beta^i|_R = (\vec{\Omega}_0 \times \vec{r})^i + \dot{a}_0 r^i + v_0^i, \quad (2)$$

giving us control over the initial kinematics of the binary system.

Figure 1 gives a schematic representation of these free data.

2.2. Calculation of asymptotic quantities

In Newtonian mechanics, it is typically straightforward to compute the center of mass and the linear momentum of a system. However, in general relativity, these are asymptotic quantities – defined by infinite surface integrals. Since the computational domains are coarser at their outer boundaries, these integrals tend to be very sensitivity to small errors, leading to less accurate results. To overcome this, we can cast these calculations as volume integrals using Gauss’ divergence theorem. Below, we explain the approaches we used to compute these asymptotic quantities in **SpECTRE**.

We can measure the total energy of a system by computing the ADM mass M_{ADM} . In terms of conformal quantities, it takes the form of

$$M_{\text{ADM}} = \frac{1}{16\pi} \oint_{S_\infty} \left(\bar{\gamma}^{jk} \bar{\Gamma}_{jk}^i - \bar{\gamma}^{ij} \bar{\Gamma}_j - 8\bar{D}^i \psi \right) d\bar{S}_i \quad (3)$$

[9, Eq. 3.139], where $\bar{\Gamma}_{jk}^i$ are the Christoffel symbols associated with $\bar{\gamma}_{ij}$ (see Eq. 2.44 in [9] for a definition) and $\bar{\Gamma}_j = \bar{\Gamma}_{ij}^i$. Applying Gauss' theorem to Eq. (3), we have

$$M_{\text{ADM}} = \frac{1}{16\pi} \int_{V_\infty} \bar{D}_i \left(\bar{\gamma}^{jk} \bar{\Gamma}_{jk}^i - \bar{\gamma}^{ij} \bar{\Gamma}_j - 8\bar{D}^i \psi \right) d\bar{V} \quad (4)$$

$$= \frac{1}{16\pi} \int_{V_\infty} \left(\partial_i \bar{\gamma}^{jk} \bar{\Gamma}_{jk}^i + \bar{\gamma}^{jk} \partial_i \bar{\Gamma}_{jk}^i + \bar{\Gamma}_l \bar{\gamma}^{jk} \bar{\Gamma}_{jk}^l - \partial_i \bar{\gamma}^{ij} \bar{\Gamma}_j - \bar{\gamma}^{ij} \partial_i \bar{\Gamma}_j - \bar{\Gamma}_l \bar{\gamma}^{lj} \bar{\Gamma}_j - 8\bar{D}^2 \psi \right) d\bar{V}, \quad (5)$$

where we can use the Hamiltonian constraint to replace $8\bar{D}^2 \psi$ with

$$8\bar{D}^2 \psi = \psi \bar{R} + \frac{2}{3} \psi^5 K^2 - \frac{1}{4} \psi^5 \frac{1}{\alpha^2} \left[(\bar{L}\beta)_{ij} - \bar{u}_{ij} \right] \left[(\bar{L}\beta)^{ij} - \bar{u}^{ij} \right] - 16\pi \psi^5 \rho \quad (6)$$

[9, Eq. 3.37].

Using the formalism developed in [11], we can find the center of mass of a system by evaluating the integral

$$C_{\text{CoM}}^i = \frac{3}{8\pi M_{\text{ADM}}} \oint_{S_\infty} \psi^4 n^i dA \quad (7)$$

[5, Eq. 26], where $n^i = x^i/r$ and $r = \sqrt{x^2 + y^2 + z^2}$. Applying Gauss' theorem to Eq. (7), we have

$$C_{\text{CoM}}^i = \frac{3}{8\pi M_{\text{ADM}}} \int_{V_\infty} \partial_j \left(\psi^4 n^i n^j \right) dV \quad (8)$$

$$= \frac{3}{4\pi M_{\text{ADM}}} \int_{V_\infty} \frac{1}{r^2} \left(2\psi^3 \partial_j \psi x^i x^j + \psi^4 x^i \right) dV. \quad (9)$$

We can also measure the total linear momentum of a system with the ADM linear momentum P_{ADM}^i using the expressions derived in [5, Eqs. 18–22]. As an infinite surface integral, it is expressed as

$$P_{\text{ADM}}^i = \frac{1}{8\pi} \oint_{S_\infty} \psi^{10} \left(K^{ij} - K \gamma^{ij} \right) dS_j. \quad (10)$$

Applying Gauss' theorem and using the momentum constraint, it becomes

$$P_{\text{ADM}}^i = \frac{1}{8\pi} \oint_{S_\infty} \partial_j \left[\psi^{10} \left(K^{ij} - K \gamma^{ij} \right) \right] dV \quad (11)$$

$$= -\frac{1}{8\pi} \int_{V_\infty} \left[\bar{\Gamma}_{jk}^i P^{jk} + \bar{\Gamma}_{jk}^j P^{jk} - 2\bar{\gamma}_{jk} P^{jk} \bar{\gamma}^{il} \partial_l (\ln \psi) \right] dV, \quad (12)$$

where $P^{ij} = \psi^{10}(K^{ij} - K\gamma^{ij})$ is the integrand of Eq. (10). Note that the matter sources are ignored when using the momentum constraint in [9], but it is straightforward to add the source term to Eq. (12) is needed in the future.

An attentive reader might have noticed that we have used Gauss' theorem in two different ways. That is, we used the conformal covariant derivative \bar{D}_i and the conformal volume element $d\bar{V}$ in Eq. (4), but the partial derivative ∂_i and the Euclidean volume element dV in Eq. (8) and Eq. (11). Both approaches are equivalent [12, 13], as long as we are consistent.

In practice, **SpECTRE** cannot evaluate volume integrals over all space because the interior of the black holes are excised from the domain. Hence, if we wish to use the infinite volume integrals, we must still perform finite surface integrals enclosing the excisions. For practicality, we choose to do such integrals over the excisions themselves. If we let S_0 be the combination of the excision surfaces, the resulting expressions are

$$\begin{aligned} M_{\text{ADM}} = & \frac{1}{16\pi} \oint_{S_0} \left(\bar{\gamma}^{jk} \bar{\Gamma}_{jk}^i - \bar{\gamma}^{ij} \bar{\Gamma}_j - 8\bar{D}^i \psi \right) d\bar{S}_i \\ & + \frac{1}{16\pi} \int_{V_\infty} \left(\partial_i \bar{\gamma}^{jk} \bar{\Gamma}_{jk}^i + \bar{\gamma}^{jk} \partial_i \bar{\Gamma}_{jk}^i + \bar{\Gamma}_l \bar{\gamma}^{jk} \bar{\Gamma}_{jk}^l \right. \\ & \quad \left. - \partial_i \bar{\gamma}^{ij} \bar{\Gamma}_j - \bar{\gamma}^{ij} \partial_i \bar{\Gamma}_j - \bar{\Gamma}_l \bar{\gamma}^{lj} \bar{\Gamma}_j - 8\bar{D}^2 \psi \right) d\bar{V}, \end{aligned} \quad (13)$$

$$\begin{aligned} C_{\text{CoM}}^i = & \frac{3}{8\pi M_{\text{ADM}}} \oint_{S_0} \psi^4 n^i dA \\ & + \frac{3}{4\pi M_{\text{ADM}}} \int_{V_\infty} \frac{1}{r^2} \left(2\psi^3 \partial_j \psi x^i x^j + \psi^4 x^i \right) dV, \end{aligned} \quad (14)$$

$$\begin{aligned} P_{\text{ADM}}^i = & \frac{1}{8\pi} \oint_{S_0} \psi^{10} \left(K^{ij} - K\gamma^{ij} \right) dS_j \\ & - \frac{1}{8\pi} \int_{V_\infty} \left[\bar{\Gamma}_{jk}^i P^{jk} + \bar{\Gamma}_{jk}^j P^{jk} - 2\bar{\gamma}_{jk} P^{jk} \bar{\gamma}^{il} \partial_l (\ln \psi) \right] dV. \end{aligned} \quad (15)$$

Because of **SpECTRE**'s task-based parallelism, we must first evaluate the integrals locally in each domain element. Then, we perform a reduction to combine the results from all elements into a single result for the entire domain.

2.3. Control scheme

Once the XCTS equations are solved, we have all the information that we need about the zero-time slice of spacetime. With this, we can use an apparent horizon finder to get measurements of the black holes in the constructed initial data. Let M_A , M_B , $\vec{\chi}_A$ and $\vec{\chi}_B$ be the masses and spins of the actual black holes that we have been able to create. In general, these values will differ from the “target” quantities M_A^* , M_B^* , $\vec{\chi}_A^*$ and $\vec{\chi}_B^*$. For example, suppose that we wish to simulate an equal-mass non-spinning case with $M_A^* = M_B^* = 0.5$ and $\vec{\chi}_A^* = \vec{\chi}_B^* = (0, 0, 0)$. In order to construct the SKS analytical expressions, it is natural to specify $M_{A,B}^{\text{Kerr}} = M_{A,B}^*$ and $\vec{\chi}_{A,B}^{\text{Kerr}} = \vec{\chi}_{A,B}^*$. With this, we can solve the XCTS equations and find horizons in the resulting initial data. Once we do it

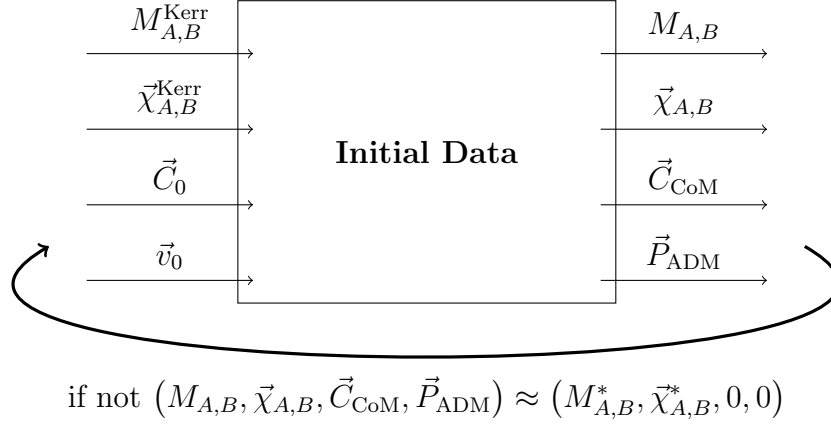


Figure 2: Schematic representation of the control loop.

in **SpECTRE**, we find that the resulting black hole parameters are actually $M_{A,B} \approx 0.6$ and $\vec{\chi}_{A,B} \approx (0, 0, -0.002)$.

If we measure the asymptotic quantities described in the previous subsection, we might find that they are also not desirable. Ideally, \vec{C}_{CoM} and \vec{P}_{ADM} should be equal to zero in order to minimize any drifts of the binary orbit, especially for long simulations. However, similar to the example above, we can only know these quantities after the initial data has been constructed.

Note that $M_{A,B}$, $\vec{\chi}_{A,B}$, \vec{C}_{CoM} and \vec{P}_{ADM} are the parameters that we wish to control, but we cannot directly set them to the desired values. We handle this issue by iterating over different choices of the free data $(M_{A,B}^{\text{Kerr}}, \vec{\chi}_{A,B}^{\text{Kerr}}, \vec{C}_0$ and $\vec{v}_0)$, trying to find the ones that result in the ideal physical parameters. Figure 2 shows a schematic representation of this “control loop”.

Let the choices of the free data be represented as

$$\mathbf{u} = \left(M_A^{\text{Kerr}}, M_B^{\text{Kerr}}, \chi_A^{\text{Kerr},x}, \chi_A^{\text{Kerr},y}, \chi_A^{\text{Kerr},z}, \right. \\ \left. \chi_B^{\text{Kerr},x}, \chi_B^{\text{Kerr},y}, \chi_B^{\text{Kerr},z}, C_0^x, C_0^y, C_0^z, v_0^x, v_0^y, v_0^z \right). \quad (16)$$

Also, let the difference between the current and target physical parameters be represented the “residual” function

$$\mathbf{F}(\mathbf{u}) = \left(M_A - M_A^*, M_B - M_B^*, \chi_A^x - \chi_A^{*,x}, \chi_A^y - \chi_A^{*,y}, \chi_A^z - \chi_A^{*,z}, \right. \\ \left. \chi_B^x - \chi_B^{*,x}, \chi_B^y - \chi_B^{*,y}, \chi_B^z - \chi_B^{*,z}, C_{\text{CoM}}^x, C_{\text{CoM}}^y, C_{\text{CoM}}^z, P_{\text{ADM}}^x, P_{\text{ADM}}^y, P_{\text{ADM}}^z \right). \quad (17)$$

As discussed in the example above, it is natural to use the target values as an initial guess for the free data:

$$\mathbf{u}_0 = \left(M_A^*, M_B^*, \chi_A^{*,x}, \chi_A^{*,y}, \chi_A^{*,z}, \right. \\ \left. \chi_B^{*,x}, \chi_B^{*,y}, \chi_B^{*,z}, 0, 0, 0, 0, 0, 0 \right). \quad (18)$$

With this, we can update our free data at every control iteration $k \geq 1$ using a Newton-Raphson scheme [14]:

$$\mathbf{u}_k = \mathbf{u}_{k-1} - \mathbb{J}_{k-1}^{-1} \cdot \mathbf{F}_{k-1}. \quad (19)$$

However, note that each evaluation of $\mathbf{F}(\mathbf{u})$ is computationally expensive as it requires an entire initial data solve. Therefore, computing the Jacobian in Eq. (19) analytically is unfeasible. That said, we can find it numerically using Broyden's method [14]:

$$\mathbb{J}_k = \mathbb{J}_{k-1} + \frac{\mathbf{F}_k \otimes \Delta \mathbf{u}_k}{\|\Delta \mathbf{u}_k\|^2}, \quad (20)$$

where $\Delta \mathbf{u}_k = \mathbf{u}_k - \mathbf{u}_{k-1}$.

Since we're iteratively computing the Jacobian with Eq. (20), we must choose an adequate initial guess \mathbb{J}_0 . Assuming that $M_{A,B}|_0 \approx M_{A,B}^{\text{Kerr}}|_0$ and $\vec{\chi}_{A,B}|_0 \approx \vec{\chi}_{A,B}^{\text{Kerr}}|_0$,

$$\left. \frac{\partial(M_{A,B} - M_{A,B}^*)}{\partial M_{A,B}^{\text{Kerr}}} \right|_0 \approx \left. \frac{\partial(\chi_{A,B}^i - \chi_{A,B}^{*,i})}{\partial \chi_{A,B}^{\text{Kerr},i}} \right|_0 \approx 1. \quad (21)$$

Using a Newtonian center of mass approximation,

$$\vec{C}_{\text{CoM}} \approx \vec{C}_0, \quad (22)$$

implying that

$$\left. \frac{\partial(C_{\text{CoM}}^i)}{\partial C_0^i} \right|_0 \approx 1. \quad (23)$$

We can also use a Newtonian linear momentum approximation with Eq. (2) as the center-of-mass velocity:

$$\vec{P}_{\text{ADM}} \approx (M_A + M_B) \left[(\vec{\Omega}_0 \times \vec{C}_{\text{CoM}}) + \dot{a}_0 \vec{C}_{\text{CoM}} + \vec{v}_0 \right], \quad (24)$$

where $\vec{\Omega}_0$ is the initial angular velocity and \dot{a}_0 is the initial radial expansion velocity. Typically, these initial orbit parameters are small and $M_A + M_B \approx 1$ in code units. Then,

$$\left. \frac{\partial(P_{\text{ADM}}^i)}{\partial v_0^i} \right|_0 \approx 1. \quad (25)$$

All the other partial derivatives that go into \mathbb{J}_0 are approximately zero under the assumptions imposed above. Therefore, a good initial guess for the Jacobian is

$$\mathbb{J}_0 \approx \mathbb{I}, \quad (26)$$

a 14-by-14 identity matrix.

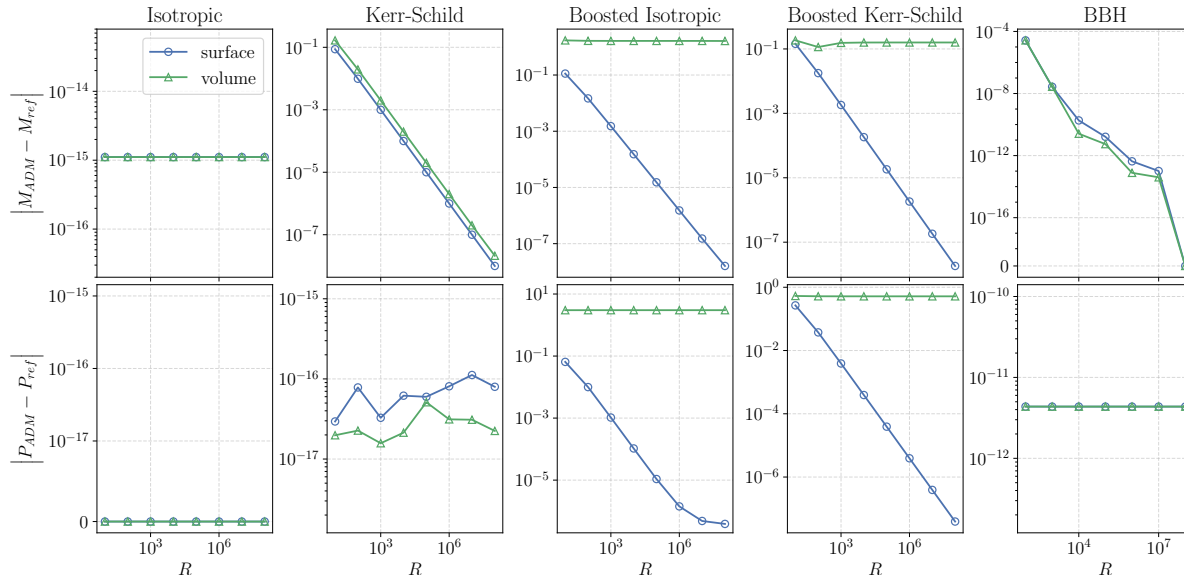


Figure 3: Convergence with distance for M_{ADM} and P_{ADM} . “Surface” indicates results from evaluating the infinite surface integrals defined in Eq. (3) and Eq. (10). “Volume” indicates results from evaluating the finite surface integrals and the infinite volume integrals defined in Eq. (13) and Eq. (15).

3. Results

3.1. Convergence of asymptotic quantities

We have tested our calculation of asymptotic quantities with analytic solutions of a single black hole, as well as with a numeric solution to the XCTS equations for BBH. The analytic solutions we used describe a Schwarzschild black hole of mass M in isotropic coordinates and in Kerr-Schild coordinates, both with and without a boost speed v . The expected values of M_{ADM} and P_{ADM} for the unboosted solutions are $M_{\text{ref}} = M$ and $P_{\text{ref}} = 0$, while the ones for the boosted solutions are $M_{\text{ref}} = \gamma M$ and $P_{\text{ref}} = \gamma M v$, where $\gamma = 1/\sqrt{1-v^2}$ is the Lorentz factor. We also use a BBH solution for an equal-mass non-spinning case, implying that the expected value of P_{ADM} is $P_{\text{ref}} = 0$. Since we do not know the expected value of M_{ADM} for the BBH solution, we use the result at the largest outer radius R as a reference: $M_{\text{ref}} = M_{\text{ADM}}|_{\text{max } R}$.

Figure 3 compares the difference between the measured M_{ADM} and P_{ADM} with the reference values M_{ref} and P_{ref} , as described in the previous paragraph. For each subplot, we see how this residual evolves as we increase the outer radius R of the domain, which serves as “infinity” in these calculations. Since these quantities are defined asymptotically, we should approach the reference values as R increases. Excluding trivial cases (in which the residuals already start very low), we see that the infinite surface integrals defined in Eq. (3) and Eq. (10) converge just as expected. However, this is not the case for the finite surface integrals and infinite volume integrals defined in Eq. (13) and Eq. (15). In fact, we see that the results do not converge at all for the boosted cases. Other than

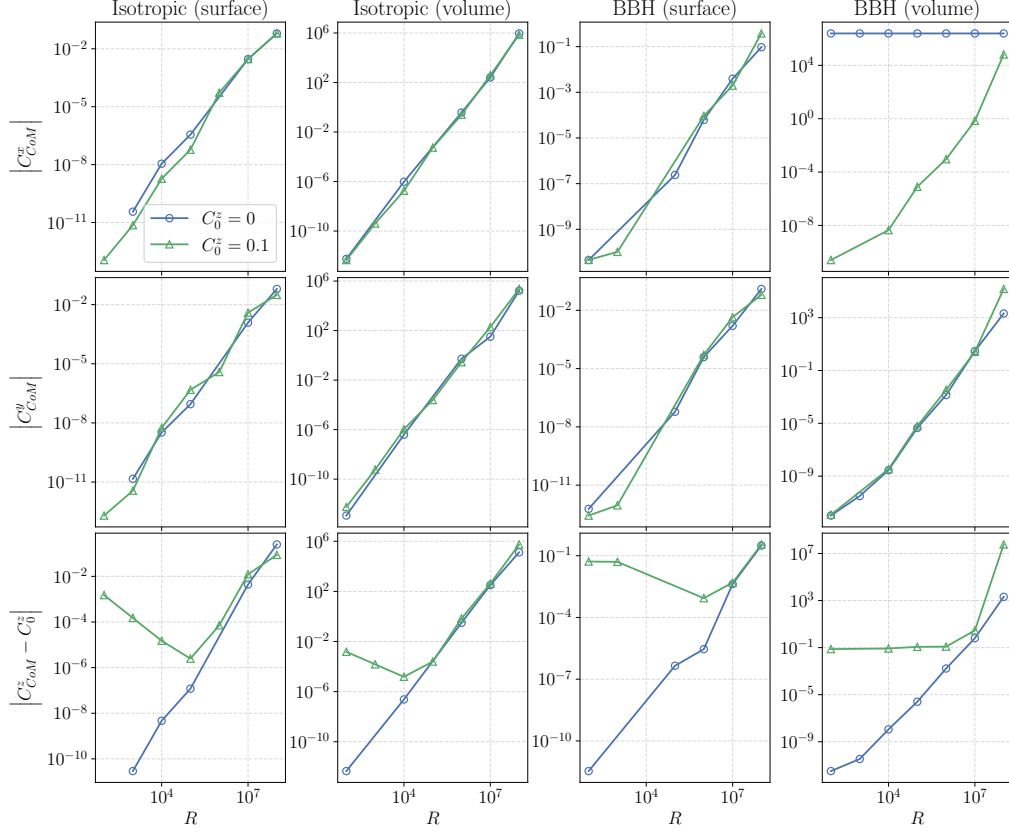


Figure 4: Convergence with distance for \vec{C}_{CoM} . “Surface” indicates results from evaluating the infinite surface integral defined in Eq. (7). “Volume” indicates results from evaluating the finite surface integral and the infinite volume integral defined in Eq. (14). Residuals for each component of \vec{C}_{CoM} are shown for an unshifted solution ($C_0^z = 0$) and for a shifted solution ($C_0^z = 0.1$).

that, this approach behaves as expected.

To study the center of mass calculations, we need a different analysis. First, the Kerr-Schild solutions do not fall off to flatness fast enough to satisfy the assumptions for Eq. (7) to hold. Second, the expected \vec{C}_{CoM} value for the isotropic Schwarzschild and BBH solutions are zero, so we need to shift the solution to get a non-vanishing result. Figure 4 shows some useful results taking these points into consideration.

An important feature in Figure 4 is that the results seem to diverge from the expected value as we increase R . We do not fully understand why this occurs, but our hypothesis is that there is a round-off error in the nature of this calculation that grows substantially as we increase R . One way to interpret Eq. (7) is that we are summing over the unit vectors n^i , rescaled by ψ^4 , in all directions. If $\psi(\vec{r})$ is constant, no rescaling happens and all the unit vectors cancel out. If $\psi(\vec{r})$ is not constant, then \vec{C}_{CoM} will emerge from the difference of large numbers. With larger and larger numbers being involved in this cancellation (i.e. with increasing R), we lose numerical accuracy. In other words, we

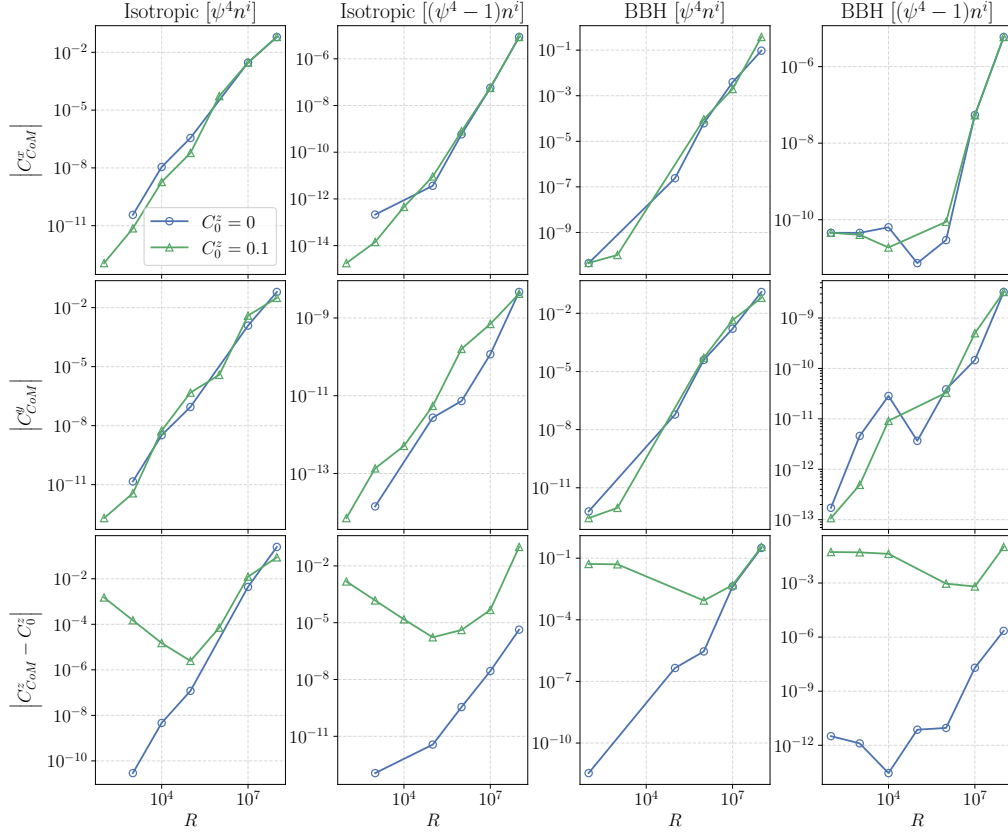


Figure 5: Comparison between different versions of the center-of-mass surface integrals. $[\psi^4 n^i]$ refers to Eq. (7), while $[(\psi^4 - 1)n^i]$ refers to Eq. (27).

are seeking the subdominant terms. This hypothesis makes sense when we look at the residuals of the shifted solutions. We see that the results converge as we increase R until when the residuals reach the level of round-off errors (indicated by the residuals of the unshifted solutions).

Comparing the “volume” residuals with the “surface” residuals in Figure 4, we see that Eq. (14) leads to higher round-off errors than Eq. (7). This also makes sense because the volume integral involves many more domain elements. However, we notice a strange behavior of the C_{CoM}^x “volume” residuals for the unshifted BBH solution. For that specific case, we see that the x component of Eq. (14) diverges since the beginning, which does not occur anymore if a z-shift is added to the solution. This might be an indication of a bug in the implementation of the method.

We can modify Eq. (7) in an attempt to address this issue with round-off errors. One way to do this is to transform the integrand $\psi^4 n^i$ to $(\psi^4 - 1)n^i$. Analytically, this should

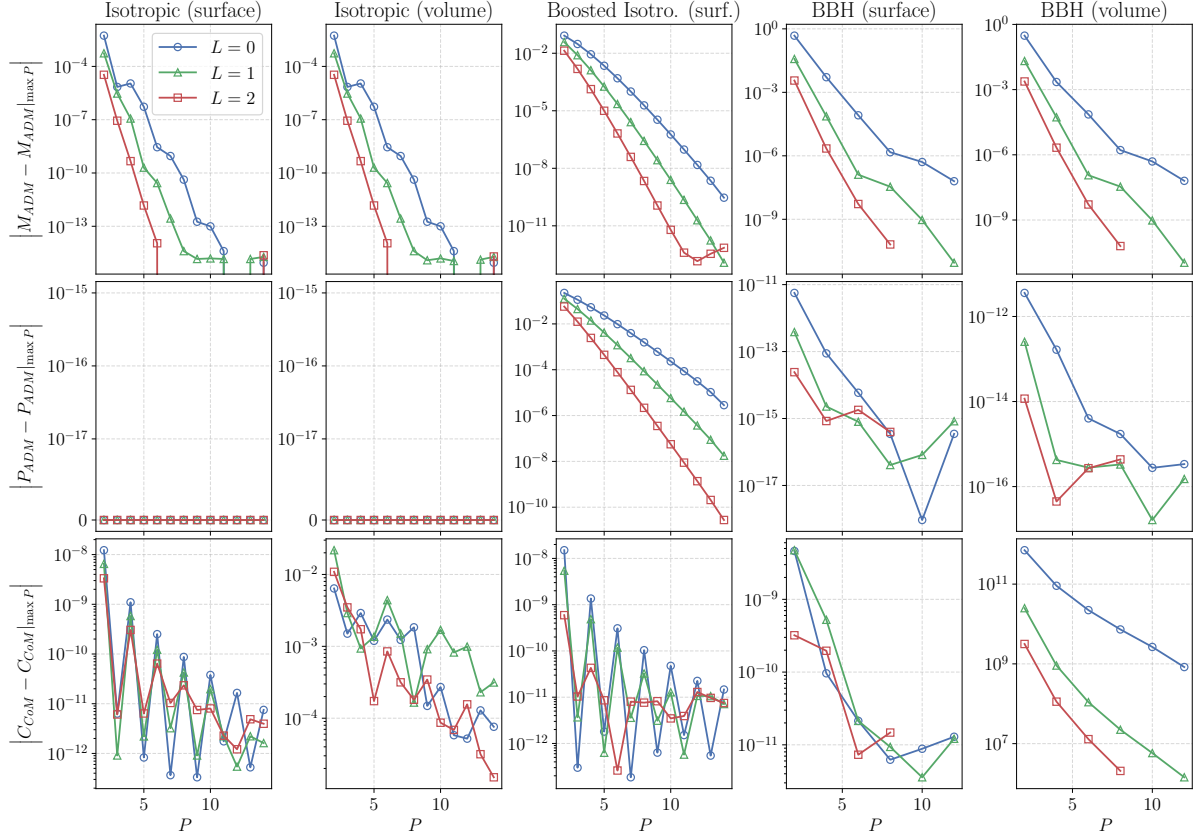


Figure 6: Convergence with resolution for M_{ADM} , P_{ADM} , and C_{CoM} . Each subplot shows the difference at each polynomial order P relative to the highest P result for multiple refinement levels L .

not change the result because

$$C_{\text{CoM}}^i = \frac{3}{8\pi M_{\text{ADM}}} \oint_{S_\infty} (\psi^4 - 1) n^i dA \quad (27)$$

$$= \frac{3}{8\pi M_{\text{ADM}}} \left(\oint_{S_\infty} \psi^4 n^i dA - \oint_{S_\infty} n^i dA \right) \quad (28)$$

$$= \frac{3}{8\pi M_{\text{ADM}}} \oint_{S_\infty} \psi^4 n^i dA. \quad (29)$$

Numerically, this makes the terms being cancelled in the calculation smaller. Figure 5 compares the residuals for each integrand. From this, it is clear that Eq. (27) significantly reduces the round-off errors. Therefore, we use this version for the remaining results of the center-of-mass surface integral.

Apart from knowing that we approach the expected values for increasing R , we must confirm that the results get more accurate with resolution. Figure 6 shows that the residuals in general converge exponentially with polynomial order P and converge faster with higher refinement level L , as expected for the Discontinuous Galerkin scheme used in SpECTRE [15]. The center-of-mass convergence plots do not follow this trend, which

Name	q	$\vec{\chi}_A^*$	$\vec{\chi}_B^*$
q1	1	(0,0,0)	(0,0,0)
q3	3	(0,0,0)	(0,0,0)
q3-spin	3	(0,0,0.9)	(0,0,0)

Table 1: Target parameters of the BBH runs. The masses $M_{A,B}^*$ are determined by the mass ratio $q = M_A^*/M_B^*$ and $M_A^* + M_B^* = 1$.

could be related to the round-off errors, but they still converge as we increase resolution. It is also important to note that the BBH center-of-mass “volume” integral has large residuals, which is expected given the C_{CoM}^x divergence in Figure 4.

The motivation to apply Gauss’ divergence theorem in the calculation of the asymptotic quantities was to achieve higher accuracy. Instead of relying on the coarse area elements at the outer boundary (i.e., Eqs. (3), (7), (10)), we can use the more-precise area elements at the inner boundaries along with the volume elements of the entire domain (i.e., Eqs. (13)–(15)). However, Figure 6 shows that we did not get much benefit out of this transformation. The ADM masses appear to be almost identical between the surface and volume results. The ADM linear momenta have slightly lower residuals for the BBH volume results compared to the surface ones, but nothing too significant. The centers of mass as volume integrals have higher round-off errors than their surface counterparts, and we see occasional blow-ups in the volume version.

Therefore, the volume integrals do not seem favorable for our purposes at the current stage. In **SpEC**, this transformation was used in the computation of the ADM linear momentum \vec{P}_{ADM} and the ADM angular momentum \vec{J}_{ADM} [5]. There, accuracy was improved by one order of magnitude for \vec{P}_{ADM} and by several orders of magnitude for \vec{J}_{ADM} [5]. That said, it is important to note that **SpEC**’s BBH outer boundary is $R \sim 10^{10}$ [5], while **SpECTRE**’s is $R \sim 10^5$ due to its better boundary conditions [16]. The smaller outer boundary might explain why we don’t see a significant improvement when using the \vec{P}_{ADM} volume integrals. We plan to compute \vec{J}_{ADM} in the future, so it is worth re-evaluating this at that time.

3.2. Control of BBH parameters

To test our control loop, we ran three BBH cases. Table 1 shows the target parameters for each of these runs. Figure 7 shows how many control iterations were needed for each run as well as how the control residual $\mathbf{F}(\mathbf{u})$ behaved. We finish the control loop when the largest component of $\mathbf{F}(\mathbf{u})$ reaches the residual tolerance of 10^{-6} . From these results, it is clear that our control scheme finished in very few iterations and was able to handle scenarios from a simple equal-mass non-spinning run to a non-trivial mass ratio with high spin.

We attempted to run the BBH cases shown in Table 1 of [5], but we did not succeed. For the **Spin0.9999** and **q10** runs, **SpECTRE**’s horizon finder failed after the first XCTS

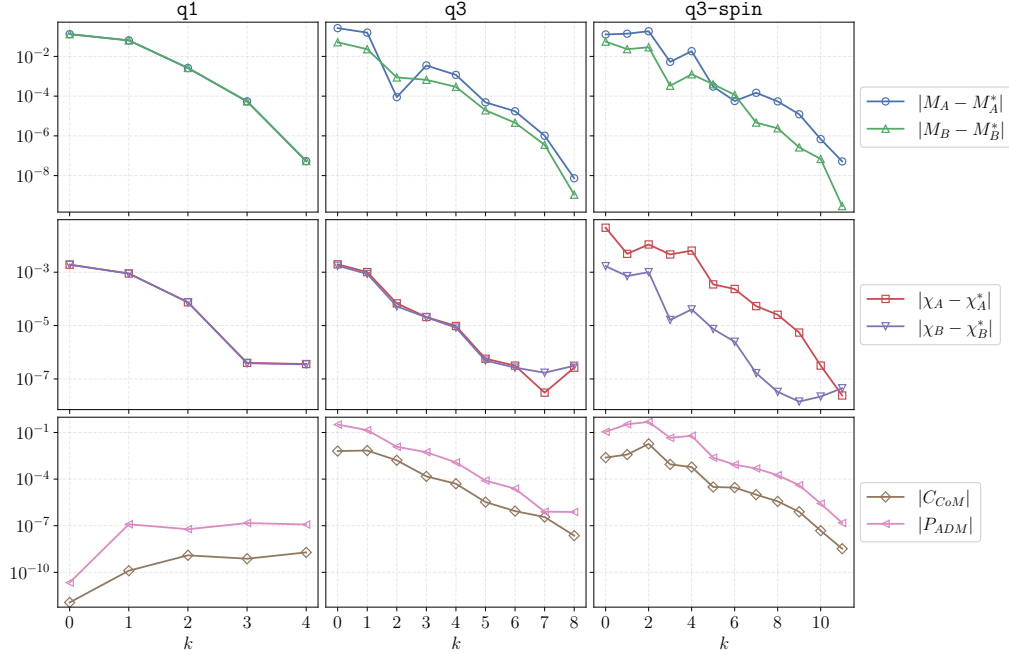


Figure 7: Control results for the BBH runs shown in Table 1. The control residuals $\mathbf{F}(\mathbf{u})$ are shown for each control iteration k . Here, \vec{C}_{CoM} and \vec{P}_{ADM} were computed using Eq. (27) and Eq. (10).

solve. There was a recent update to the **SpECTRE** BBH domain, which might have fixed such problems. For the **q3** and **q10** runs, the free data updated by Eq. (19) contained invalid values like negative masses and $\chi > 1$. This could be fixed by adding constraints to the iterative loop, making sure that the free data stays within an acceptable range.

4. Conclusion

SpECTRE is now capable of controlling physical parameters in BBH initial data. Our control scheme not only enforces specific masses and spins for the desired simulation, but also avoids drifts in the orbital trajectories by minimizing the center of mass and the total linear momentum. For this, we had to compute asymptotic quantities defined as infinite integrals, which can be cast as surface integrals or volume integrals. The integral results for the total energy and linear momentum convergence nicely. Although the integral results for the center of mass has some issues that must be investigated, the surface integral achieves acceptable residuals for the current applications.

There are many possible improvements to the work described in this report. For instance, we plan to find ways to get around the round-off errors observed in the center-of-mass integrals. This could include changes to the integral itself, like in Eq. (27), or alternative integration methods, such as the spectral integration done in **SpEC** [5]. We should also make the control scheme more rigorous with constraints for the free data and more efficient with optimizations like line search, allowing **SpECTRE** to reach more of the parameter space.

There is also a lot of interest in follow-up work. For example, it is important to understand whether we have made significant improvements in our control scheme relative to the one used by SpEC, either in terms of performance or parameter space. To this purpose, we need to do a careful comparison between the two codes. Additionally, by implementing the computation of total angular momentum \vec{J}_{ADM} , we would be able to control M_{ADM} and \vec{J}_{ADM} to generate initial data for BBH hyperbolic encounters.

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