We utilize the dynamical renormalization group formalism to calculate the real space trajectory of a compact binary inspiral for long times via a systematic resummation of secularly growing terms. This method generates closed form solutions without orbit averaging, and the accuracy can be systematically improved. The expansion parameter is $v^5 t \Omega(t - t_0)$ where $t_0$ is the initial time, $t$ is the time elapsed, and $\Omega$ and $v$ are the angular orbital frequency and initial speed, respectively. $\nu$ is the binary’s symmetric mass ratio. We demonstrate how to apply the renormalization group method to resum solutions beyond leading order in two ways. First, we calculate the second-order corrections of the leading radiation reaction force, which involves highly nontrivial checks of the formalism (i.e., its renormalizability). Second, we show how to systematically include post-Newtonian corrections to the radiation reaction force. By avoiding orbit averaging, we gain predictive power and eliminate ambiguities in the initial conditions. Finally, we discuss how this methodology can be used to find analytic solutions to the spin equations of motion that are valid over long times.

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I. INTRODUCTION

The recent detections of gravitational waves from a binary black hole coalescence [1,2] provide the first measurements of the dynamics of compact binaries and strong gravitational fields. To date, the measured events involved relatively large black hole masses. For the first detection, several gravitational wave cycles were observed in LIGO’s frequency band corresponding to a handful of orbits before merger. If compact binaries with lower total masses are observed, then these sources will evolve for much longer times within the detector’s bandwidth and will be amenable to analytic calculations via the post-Newtonian (PN) expansion.

In order to extract the most information from such long waveforms requires using templates that have been computed with the highest possible accuracy. The lengthy inspiral regime of the binary’s evolution is found by solving the post-Newtonian equations of motion. A challenge to this program is the calculation of the radiation reaction forces, the leading piece of which starts at 2.5PN [3,4]. The PN corrections at 1PN beyond leading order were calculated in Refs. [5–8]. At 1.5PN, there is a contribution from the “tail effect,” which was calculated in Refs. [9,10]. Higher-order corrections have yet to be calculated; however, the leading contributions from spin-orbit and spin-spin effects are known [11,12] and first appear at 1.5PN and 2PN orders, respectively, beyond the leading radiation reaction force. As such, solving the equations of motion, even numerically, would lead to errors which are of order the first unknown radiation reaction contribution. This would bound the accuracy of the prediction given that the conservative pieces of the equations of motion are known to higher order. To avoid this issue, one can utilize the power loss to account for radiation reaction by using the adiabatic approximation (more on this below), which requires one to average over the orbit, thereby generating information loss.

In addition to this aforementioned limitation, solving for the binary’s orbital motion is often achieved numerically because there are nine (once we include spin) nonlinearly coupled, ordinary differential equations that need to be solved accurately in order to follow the orbit’s inspiral. The orbit needs to be sampled at a sufficiently high rate so that the corresponding waveform is sampled appropriately. These accuracy requirements and/or high sample rate are often a computational bottleneck for gravitational wave data analysis applications that require many waveforms to be generated and hence many numerical solutions of the PN equations of motion. Such applications include template generation for gravitational wave searches and parameter estimation studies using Markov-chain Monte Carlo algorithms.

Recent developments in implementing a “precession-averaging” procedure [13,14] to the equations of motion, which utilize the separation of time scales present in precessing binaries, help alleviate some of the computational pressures mentioned above, depending on the specific application. However, accurate, globally valid, analytical
approximate solutions for the binary’s evolution would certainly remove any such computational bottleneck while simultaneously providing useful analytical expressions for studying the complicated physics of precession dynamics (such as recent evidence for precessional instability [15]).

Currently, there are two standard methods used to provide analytic solutions for the restricted case of compact binaries with component masses that are spinning but nonprecessing. The first is the “adiabatic approximation” and is based on equating the time-averaged flux of gravitational waves to the mechanical power lost by the binary (see, e.g., Refs. [16,17]). This method is often used to calculate the approximate gravitational wave phase of the (l, m) = (2, 2) spherical harmonic mode in the “restricted approximation” where the waveform is constructed from the orbital phase up to a definite order in the PN expansion, but the amplitude is taken to be the leading-order quantity with no PN corrections. The second is the “improved variation of constants method” [18]. Here, the idea is to assume that the integration constants of the conservative part of the binary’s dynamics exhibit a long-time evolution relative to the orbital period. Equations for these integration “constants” are then found using the method of variation of constants, which are then solved. This approach is also known as the “method of osculating orbits” and is very closely related to “multiple scale analysis.” Recent work in Ref. [19] builds off of Ref. [13] and uses multiple scale analysis to find accurate analytic approximations for the orbit and the frequency-domain waveform of a precessing binary inspiral.

These approaches are based on averaging the PN equations of motion over the orbital period to help simplify the differential equations being solved. However, there are potential shortcomings with averaging that have been raised in Refs. [20,21]. In particular, the initial conditions used to solve the PN equations of motion and the orbit-averaged version are not the same. The initial conditions set in the adiabatic approximation are fewer than those needed for solving the PN equations of motion because the orbit averaging tends to eliminate explicit dependence on the orbital phase while also constraining the relationship between orbital radius and frequency. Orbit averaging results in a simplified description of compact binary inspirals so that comparing the adiabatic approximation solutions to those of the full PN equations can be ambiguous. In addition, there is an ambiguity in the period to use for the averaging procedure for eccentric orbits because there are different ways to characterize the time scale of the orbit (such as the orbital period, eccentric anomaly, true anomaly, and mean anomaly). Using a different measure for the averaging can lead to different predictions over sufficiently long times.

Our work introduces a formalism that allows for the systematic solution of the PN equations of motion for a binary inspiral including radiation reaction forces and spin effects to an arbitrary order in the PN expansion. The method is based on applying ideas from renormalization group theory. We do not implement any averaging procedures so that our solutions describe the binary’s real-time orbital configuration at every instant of time. The approach starts with a background (e.g., circular or eccentric) orbit and treats the radiation reaction force as a perturbation. These perturbations grow secularly with time but can be resummed using the Dynamical Renormalization Group (DRG) method [22]. The DRG method subsumes several approaches for the global analysis of differential equations [23], including multiple scale analysis, boundary layer theory, and the Wentzel–Kramers–Brillouin method. See Ref. [22] and the Appendix of Ref. [24] for pedagogical examples using the DRG method. The DRG resummation is not to be confused with Padé “resummation,” which attempts to improve the radius of convergence of a perturbative expression. Padé resummation is not a systematic expansion as it does not resum any leading-order pieces in a systematic way. Different Padé approximants often give different predictions at the same scale. Conversely, DRG literally resums higher-order terms in the perturbation theory with a consistent power counting.

Our focus here is to present the DRG method and demonstrate the internal consistency of the approach for nonspinning compact binary inspirals via higher-order perturbative calculations. We will show how to find closed form solutions for the inspiral without recourse to the adiabatic approximation or orbit averages. In a future paper, we will incorporate spin effects to obtain accurate and globally valid, real-time approximate solutions for the generic case of precessing compact binaries.

II. DEFINING THE SYSTEMATICS

We will perturbatively solve the PN equations of motion for compact binary inspirals, which are derived in an expansion where the binary’s relative speed v is small compared to the speed of light. However, there is another power counting parameter for inspirals, namely,

$$\epsilon \equiv v / c \Omega (t - t_0).$$  \hspace{1cm} (2.1)

Here, t is the time elapsed since the initial time to, \Omega is the initial angular orbital frequency, v is the initial orbital speed, and \nu is the binary’s symmetric mass ratio. The parameter \epsilon arises as a consequence of the secular growth due to the radiation reaction force which is treated as a perturbation of the circular solution (or, more generally, energy-conserving motions) to the equations of motion. By performing a resummation, the accuracy of the perturbative solutions will be extended to much later times even when \epsilon is of order 1. Without such a resummation, the perturbative solutions would be of minimal utility. By resumming powers of \epsilon, we are able to make precision predictions with well-defined systematics such that the result for the
orbit is valid until the PN expansion breaks down as the plunge is approached. Our formalism allows us to go to arbitrary order in $\varepsilon$ and allows for the systematic inclusion of PN corrections. In this paper, we will demonstrate how to work to second order in $\varepsilon$. If there were no higher-order PN corrections, then the resulting resummed solution would be valid up to times when

$$v^{10}v^2\Omega(t - t_0) \sim 1. \quad (2.2)$$

However, in reality, we must consider PN corrections that would contribute at lower orders. We will also demonstrate how to include contributions from the radiation reaction force that are at higher PN orders by calculating the 1PN correction to the orbital motion. It is important to realize that none of the results in this paper include all of the effects at a given order because our purpose is to present the method here. If we wished to perform the calculation including all 2PN effects, for instance, we would need to include the conservative potential up to 2PN, which is of course known, but we would also need to include the 2PN correction to the radiation reaction force which is presently unknown.

### III. REVIEW OF RENORMALIZATION GROUP METHODOLOGY

For completeness, we present a lightning review of the logic behind the renormalization group (RG). The DRG applies the logic of the RG to differential equations, but the basic idea is the same. Canonical RG applications are formulated within the context of a Lagrangian which will not be the case for the DRG, though it is a simple exercise to embed the DRG into a Lagrangian formalism (necessarily for generic nonconservative systems [25,26]). However, doing so does not lead to any new insights (that we can see, at least). Thus, we will eschew such a treatment.

The basic algorithm is given as follows:

1. Write down a background solution around which to perturb. This solution is written in terms of “bare” parameters [i.e., $A_R(t_0)$]. These parameters implicitly depend upon the initial time $t_0$, away from which we flow.

2. Use this background to calculate perturbatively the first correction to the equations of motion. The perturbation will in general have secular “divergences,” that is, terms that grow as $(t - t_0)$.

3. Take this solution, and write the bare parameters as renormalized parameters [i.e., $A_R(\tau)$] plus “counterterms.” These counterterms will be proportional to $(\tau - t_0)$ and are chosen to eliminate the $t_0$ dependence of the aforementioned solution. $\tau$ is known as the subtraction point. This step yields the “renormalized” solution.

(4) The renormalized solution must be independent of the choice of subtraction point $\tau$. The explicit dependence on $\tau$ in the solution is cancelled by the implicit dependence of the renormalized parameters on $\tau$. One then uses this fact to derive a first-order differential equation (called the RG equation) for the renormalized parameter. The right-hand side of the RG equation is called the “beta function.”

(5) Solve the RG equation for the parameter, and choose $\tau = t$, the observation time. In so doing, all of the secularly growing terms are resummed at this order.

The ability to absorb divergences into the initial data, in the context of DRG, is called “renormalizability.” The renormalizability of the theory can be put on firmer mathematical ground using envelope theory as discussed in Ref. [27]. The basic notion is that a perturbative solution defines a family of curves parametrized by $t_0$. Each of these solutions is only valid locally for times near $t_0$. A global solution is then found by determining the envelope of this set of curves, which is defined as the curve of which the intersection with each curve in the family is tangent to the given curve.

The connection between the RG and global analysis is also manifest in holography. Solving the equations of motion for a scalar field in anti-de Sitter spacetime via the DRG leads to a first-order equation for the boundary data that exactly corresponds to the beta function for the coupling in the dual quantum field theory [28].

### IV. LEADING ORDER INSPIRAL

The equations of motion in the harmonic gauge in the center-of-mass frame through leading order in the potential (i.e., Newtonian) and radiation reaction forces are [3,4,17]

$$a = -\frac{M}{r^3}r + \frac{M^2}{15r^4}r^4 \left(\frac{136M}{r} + 72v^2\right)$$

$$-\frac{8M^2}{5r^3} \left(\frac{3M}{r} + v^2\right)v. \quad (4.1)$$

In terms of polar coordinates, Eq. (4.1) is expressed as

$$\ddot{r} - r\ddot{\omega} = -\frac{M}{r^2} + \frac{64M^3}{15r^4}r + \frac{16M^2}{5r}r^3 + \frac{16M^2}{5} - \frac{r}{r^5} \omega$$

$$r\dot{\omega} + 2\dot{r}\omega = -\frac{24M^2}{5r^3} - \frac{8M^3}{5r^2}r^2 - \frac{8M^2}{3} - \frac{r}{5} - \omega^3, \quad (4.2)$$

where $\omega(t) = \dot{\phi}(t)$ is the binary’s orbital angular frequency. The orbital plane does not precess, and the motion is described fully by the binary’s separation $r(t)$ and the orbital phase $\phi(t)$.

We will solve these equations perturbatively in the PN expansion. Of course, we are ignoring the 1PN and 2PN conservative forces that should be included for a consistent description through 2.5PN order. Nevertheless, it is
sufficient to use (4.2) for our purpose of demonstrating the DRG method.

A. Perturbations of a background circular orbit

We begin by considering the radiation reaction force to be negligible so that the background orbital motion is nearly circular. We have chosen these conditions because it is widely expected that many compact binary sources will have circularized by the time their radiated gravitational waves enter the frequency band of ground-based detectors. However, it is straightforward to incorporate eccentircity waves enter the frequency band of ground-based detectors.

The leading-order background circular orbit is described by a constant radius $R_B$ and constant angular frequency $\Omega_B$ with

$$\Omega_B^2 = \frac{M}{R_B^3}. \quad (4.3)$$

We next calculate the deviations of this background orbit due to the leading-order radiation reaction force (4.2) by writing $r(t) = R_B + \delta r(t)$ and $\omega(t) = \Omega_B + \delta \omega(t)$ where the perturbations scale with the relative velocity at the initial time $t_0$ so that the background orbital motion is be negligible so that the background orbital motion is

$$\delta r(t) = -3\Omega_B^2 \delta r(t) - 2R_B \delta \Omega_B \delta \omega(t) = 0$$

$$R_B \delta \omega(t) + 2\Omega_B \delta r(t) = -\frac{32\nu}{5} R_B^5 \Omega_B^2 \delta \omega(t). \quad (4.4)$$

Solving for $\delta \omega$ and substituting back into the $\delta r$ equation in (4.4) gives

$$\delta r(t) + \Omega_B^2 \delta \omega(t) = -\frac{64\nu}{5} \Omega_B^4 R_B^6 (t-t_0). \quad (4.5)$$

This equation is simple to solve using the retarded Green’s function

$$G_{ret}(t-t') = \frac{\sin \Omega_B(t-t')}{\Omega_B} \quad (4.6)$$

and results in the following general solution:

$$r(t) = R_B - \frac{64\nu}{5} \Omega_B^6 R_B^6 (t-t_0) + \frac{64\nu}{5} \Omega_B^6 R_B^6 \sin \Omega_B(t-t_0)$$

$$+ A \sin(\Omega_B(t-t_0) + \Phi)$$

$$\omega(t) = \Omega_B + \frac{96\nu}{5} R_B^4 \Omega_B^4 (t-t_0) - \frac{128\nu}{5} R_B^4 \Omega_B^4 \sin \Omega_B(t-t_0)$$

$$- \frac{2\Omega_B^4}{R_B} \sin(\Omega_B(t-t_0) + \Phi). \quad (4.7)$$

The last two terms are solutions to the homogeneous equation of (4.5) and come with two initial condition parameters, $A$ and $\Phi$. As such, we will redefine our background solution to be

$$r(t) = R_B + A_B \sin(\Omega_B(t-t_0) + \Phi_B)$$

$$\omega(t) = \Omega_B - \frac{2\Omega_B A_B}{R_B} \sin(\Omega_B(t-t_0) + \Phi_B), \quad (4.8)$$

where $A_B$ is related to a small orbital eccentricity, $e_B \sim v^5$, through

$$A_B = e_B R_B. \quad (4.9)$$

The perturbations consist of two types of pieces. The first are secularly growing in time. Since, at a time

$$t-t_0 \sim \frac{1}{\nu \Omega_B^6 R_B^5} \sim \frac{1}{\nu v^5 \Omega_B^6} \quad (4.10)$$

the perturbation becomes $O(1)$, these terms will need to be resummed in order to determine the long-time behavior of the system. The remaining terms will be perturbatively small for all times.

B. Renormalization

The first step in the resummation procedure is renormalization. This involves absorbing all of the $t_0$ dependence into the bare parameters, i.e., those constants labelled by a subscript $B$. We write our bare solution as

$$r(t) = R_B - \frac{64\nu}{5} R_B^6 \Omega_B^6 (t-t_0) + \frac{64\nu}{5} R_B^6 \Omega_B^6 \sin \Omega_B(t-t_0)$$

$$+ A_B \sin(\Omega_B(t-t_0) + \Phi_B)$$

$$\omega(t) = \Omega_B + \frac{96\nu}{5} R_B^4 \Omega_B^4 (t-t_0) - \frac{128\nu}{5} R_B^4 \Omega_B^4 \sin \Omega_B(t-t_0)$$

$$- \frac{2\Omega_B^4}{R_B} \sin(\Omega_B(t-t_0) + \Phi_B), \quad (4.11)$$

where we have promoted the integration constants $A$ and $\Phi$ to the status of bare parameters. Notice that $A_B \sim v^5 R_B$, which implies that $\dot{r}(t_0) \sim v^5 R_B \Omega_B$.

Furthermore, we may drop the nonsecularly growing sinusoidal terms (which are solutions to the homogeneous first-order equations of motion) in the solution for $r(t)$ and $\omega(t)$. This amounts to a shift in the initial conditions, which can be accomplished by the following replacement,

$$A_B \rightarrow A_B - \frac{64\nu}{5} R_B^6 \Omega_B^6 \cos(\Phi_B)$$

$$\Phi_B \rightarrow \Phi_B + \frac{64\nu}{5} R_B^6 \Omega_B^6 A_B \sin(\Phi_B). \quad (4.12)$$
The bare solution becomes

\[ r(t) = R_B - \frac{64\nu}{R_B} R_b^6 \omega_B^6 (t - t_0) + A_B \sin(\omega_B(t - t_0) + \phi_B) \]

\[ \omega(t) = \omega_B + \frac{96\nu}{5} R_B^5 \omega_B^5 (t - t_0) - \frac{2\Omega_B^A B}{R_B} \sin(\omega_B(t - t_0) + \phi_B), \quad (4.13) \]

which satisfies the equations of motion. For completeness, the orbital phase \( \phi(t) \) is computed from \( \omega(t) \) via a simple integration,

\[ \phi(t) = \Phi_B + \Omega_B(t - t_0) + \frac{48\nu}{5} R_B^5 \omega_B^5 (t - t_0)^2 \]

\[ + \frac{2A_B}{R_B} \cos(\omega_B(t - t_0) + \phi_B). \quad (4.14) \]

Notice that \( \Phi_B \) is not the initial phase, \( \phi(t_0) \). Overall constants can be dropped since they can be removed by a coordinate change without affecting the equations of motion.

The four quantities \( R_B, \Omega_B, A_B \), and \( \Phi_B \) are parameters fixed by the initial data of the problem. However, the initial time \( t_0 \) is completely arbitrary, and we could have performed the perturbative expansion at a slightly later time, \( t'_0 = t_0 + \delta t \), for instance. The formal expression of the perturbative solution would have the same form as in (4.8) except with a new set of initial conditions \( R'_B, \Omega'_B, A'_B \), and \( \Phi'_B \) and with \( t_0 \) replaced by \( t'_0 \). If \( \delta t \) is small, then it is straightforward to see that the initial conditions at \( t_0 \) are related to those at \( t'_0 \). This time shift can be compensated for by redefining the bare parameters as

\[ R'_B = R_B - \frac{64\nu}{5} R_B^6 \omega_B^6 \delta t + \mathcal{O}(\delta t^2) \]

\[ \Omega'_B = \Omega_B + \frac{96\nu}{5} R_B^5 \omega_B^5 \delta t + \mathcal{O}(\delta t^2) \]

\[ \Phi'_B = \Phi_B + \Omega_B \delta t + \mathcal{O}(\delta t^2). \quad (4.15) \]

Therefore, the perturbative solution at \( t'_0 \) is related to that at \( t_0 \) by redefining the initial conditions in such a way as to preserve the form of the perturbative solution in (4.8). In this way, the perturbative solutions may be pieced together from one time to any other, thereby generating the long-time inspiral dynamics, in the limit that \( \delta t \to 0 \), up to the PN accuracy of the original perturbative solution [27]. This process of redefining, or renormalizing, the initial conditions to ensure the form invariance of the perturbative solution at different times is at the heart of the DRG method [22] and, more generally, renormalization group theory.

We regard \( R_B, \Omega_B, A_B \), and \( \Phi_B \) as bare parameters that depend on the initial time \( t_0 \), as suggested in (4.15). We may think of \( t_0 \) as the cutoff in the usual Wilsonian sense.

All physical renormalized quantities are independent of \( t_0 \). We relate the bare parameters to their renormalized values \( R_R, \Omega_R, A_R \), and \( \Phi_R \) through the relations

\[ R_B(t_0) = R_R(\tau) + \delta_R(\tau, t_0) \quad (4.16) \]

\[ \Phi_B(t_0) = \Phi_R(\tau) + \delta_\phi(\tau, t_0) \quad (4.17) \]

\[ \Omega_B(t_0) = \Omega_R(\tau) + \delta_\Omega(\tau, t_0) \quad (4.18) \]

\[ A_B(t_0) = A_R(\tau) + \delta_A(\tau, t_0), \quad (4.19) \]

where \( \delta_R, \delta_\phi, \delta_\Omega, \) and \( \delta_A \) are quantities called counterterms that are to be determined order by order in the process of renormalizing the perturbative solutions in (4.8) and (4.14). The new time parameter \( \tau \) is the renormalization scale and is arbitrary. The initial time \( t_0 \) is like a cutoff scale when regularizing the divergences of a field theory. However, the perturbative solutions are independent of \( \tau \) at any given order in perturbation theory.

In terms of the renormalized initial parameters, the one-loop result becomes

\[ r(t) = R_R + \delta_R - \frac{64\nu}{5} R_R^6 \omega_R^6 (t - t_0) \]

\[ + (A_R + \delta_A) \sin((t - t_0) \omega_R + \Phi_R + \delta_\phi) \quad (4.20) \]

\[ \omega(t) = \omega_R + \delta_\Omega + \frac{96\nu}{5} R_R^5 \omega_R^5 (t - t_0) \]

\[ - \frac{2\Omega_R(A_R + \delta_A)}{R_R} \sin((t - t_0) \omega_R + \Phi_R + \delta_\phi) \quad (4.21) \]

\[ \phi(t) = \Phi_R + \delta_\phi + (t - t_0)(\omega_R + \delta_\Omega) + \frac{48\nu}{5} R_R^5 \omega_R^5 (t - t_0)^2 \]

\[ + \frac{2(A_R + \delta_A)}{R_R} \cos((t - t_0) \omega_R + \Phi_R + \delta_\phi). \quad (4.22) \]

where we have dropped terms of order \( \nu^{10} \).

We introduce the renormalization scale into the above solutions through \( t - t_0 = (t - \tau) + (\tau - t_0) \) so that (4.20)–(4.22) become

\[ r(t) = R_R + \delta_R - \frac{64\nu}{5} R_R^6 \omega_R^6 (t - \tau) \]

\[ - \frac{64\nu}{5} R_R^6 \omega_R^6 (t - t_0) + A_R \sin((t - \tau) \omega_R) \]

\[ + (\tau - t_0) \omega_R + \Phi_R + \delta_\phi \quad (4.23) \]

\[ \omega(t) = \omega_R + \delta_\Omega + \frac{96\nu}{5} R_R^5 \omega_R^5 (t - \tau) + \frac{96\nu}{5} R_R^5 \omega_R^5 (t - t_0) \]

\[ - \frac{2\Omega_R A_R}{R_R} \sin((t - \tau) \omega_R + (\tau - t_0) \omega_R + \Phi_R + \delta_\phi) \quad (4.24) \]
\[ \phi(t) = \Phi_R + \delta \phi + (t - \tau)\Omega_R + (t - \tau_0)\Omega_R + (t - \tau)\delta \Omega + (t - \tau_0)\delta \Omega + 48\frac{\nu}{5} R^5_k \Omega^4_R(t - \tau)^2 + \frac{96\nu}{5} R^5_k \Omega^4_R(t - \tau_0)^2 + 2A_R \cos((t - \tau)\Omega_R + (t - \tau_0)\Omega_R + \Phi_R + \delta \phi). \]

Renormalization proceeds by fixing the counterterms at this order in \( \epsilon \) to cancel the pieces that are proportional to powers of \((\tau - t_0)\). For instance, inspection of (4.23) shows that the counterterm \( \delta R \) is fixed at one-loop order to be \(^2\)

\[ \delta R^2(\tau, t_0) = \frac{64\nu}{5} R^5_R \Omega^6_R(\tau - t_0), \]  

where we have written \( \delta R = \delta R^2 + \mathcal{O}(\epsilon^2) \). The \( \mathcal{O}(\epsilon^2) \) term is a two-loop contribution that will be calculated in the next section. Likewise, the counterterm \( \delta \Omega \) is found from (4.24) to cancel the term proportional to \( \tau - t_0 \) so that

\[ \delta \Omega^3(\tau, t_0) = -\frac{96\nu}{5} R^5_k \Omega^4_R(\tau - t_0), \]  

where we have again written \( \delta \Omega = \delta \Omega^3 + \mathcal{O}(\epsilon^2) \).

Then, substituting these counterterms into (4.25), we find that the perturbative solution for \( \phi(t) \) becomes

\[ \phi(t) = \Phi_R + \delta \phi + (t - \tau)\Omega_R + (t - \tau_0)\Omega_R + \frac{48\nu}{5} R^5_k \Omega^4_R(\tau - \tau_0)^2 + \frac{48\nu}{5} R^5_k \Omega^4_R(\tau - \tau_0)^2 + 2A_R \cos((t - \tau)\Omega_R + (t - \tau_0)\Omega_R + \Phi_R + \delta \phi). \]

Notice that the term proportional to \((t - \tau)(\tau - t_0)\) automatically cancels out of the equation, which turns out to be an important check of self-consistency as we shall see in the next section. We see that we will need an additional counterterm for \( \phi_B \).

Choosing \( \delta \phi(\tau, t_0) = -\Omega_R(\tau - t_0) + \frac{48\nu}{5} R^5_k \Omega^4_R(\tau - \tau_0)^2 + \mathcal{O}(\epsilon^2) \)

removes the last remaining secular terms, even those appearing inside the oscillating terms in (4.23) and (4.24) to this order in \( \epsilon \). We are then left with the renormalized perturbative solutions

\[ r(t) = R_R - \frac{64\nu}{5} R^5_k \Omega^6_R(\tau - \tau) + A_R \sin((t - \tau)\Omega_R + \Phi_R) \]

\[ \omega(t) = \Omega_R + \frac{96\nu}{5} R^5_k \Omega^4_R(\tau - \tau) - \frac{2A_R \cos((t - \tau)\Omega_R + \Phi_R)}{R_R}. \]

Since \( \tau \) is arbitrary, we will choose it to equal \( t \) when we consider the physical solution so as to minimize all of the secular terms giving

\[ r(t) = R_R(t) + A_R(t) \sin \Phi_R(t) \]

\[ \omega(t) = \Omega_R(t) - \frac{2\Omega_R(t)A_R(t)}{R_R(t)} \sin \Phi_R(t) \]

\[ \phi(t) = \Phi_R(t) + \frac{2A_R(t)}{R_R(t)} \cos \Phi_R(t). \]

This step is akin to scale setting in the context of canonical RG flows.

**C. Renormalization group solution**

The time dependence of the renormalized initial data is found by noting that the bare parameters are independent of the arbitrary scale \( \tau \) so that \( dR_B(t_0)/d\tau = 0 \) and likewise for the other three initial parameters. For example, recall that the bare parameter \( R_B \) is given in (4.16) by

\[ R_B(t_0) = R_R(\tau) + \delta R(\tau, t_0) \]

\[ = R_R + \frac{64\nu}{5} R^5_k \Omega^6_R(\tau - t_0) + \mathcal{O}(\epsilon^2) \]

so that

\[ 0 = \frac{dR_B(t_0)}{d\tau} = \frac{dR_R(\tau)}{d\tau} + \frac{64\nu}{5} R^5_k \Omega^6_R \]

\[ + \frac{384\nu}{5} R^5_k \Omega^4_R(\tau - t_0) \frac{dR_R(\tau)}{d\tau} + \frac{384\nu}{5} R^5_k \Omega^4_R(\tau - t_0) \frac{d\Omega_R(\tau)}{d\tau} + \mathcal{O}(\epsilon^2). \]

It is easy to see that solving perturbatively for \( dR_R/d\tau \) leaves us with

\[ \frac{dR_R(\tau)}{d\tau} = \frac{64\nu}{5} R^5_k \Omega^6_R(\tau - \tau) + \frac{96\nu}{5} R^5_k \Omega^4_R(\tau - \tau) - \frac{2A_R \cos((t - \tau)\Omega_R + \Phi_R)}{R_R}. \]
\[ \frac{d}{d\tau} R_R(\tau) = -\frac{64\nu}{5} R_R^6(\tau) \Omega_R^5(\tau) + \cdots. \]  

(4.38)

since the last two terms in (4.37) are higher-order corrections. Repeating these steps for the remaining initial data yields a total of four renormalization group equations describing the RG flow, or trajectory, of the initial conditions

\[ \frac{d}{d\tau} \Omega_R(\tau) = \frac{96\nu}{5} R_R^5(\tau) \Omega_R^4(\tau), \]

(4.39)

\[ \frac{d}{d\tau} \Phi_R(\tau) = \Omega_R(\tau), \]

(4.40)

\[ \frac{d}{d\tau} A_R(\tau) = 0. \]

(4.41)

The right sides of these equations are called beta (\(\beta\)) functions in field theory. The solutions to the RG equations (4.38)–(4.41) are easily found by integrating from \(\tau = t_i\) to \(\tau = t\),

\[ R_R(t) = \left( R_R^4(t_i) - \frac{256\nu}{5} M^3(t - t_i) \right)^{1/4}, \]

(4.42)

\[ \Omega_R(t) = \Omega_R(t_i) \left( \frac{R_R(t_i)}{R_R(t)} \right)^{3/2}, \]

(4.43)

\[ \Phi_R(t) = \Phi_R(t_i) + \frac{R_R^{3/2}(t_i) - R_R^{3/2}(t)}{32\nu M^{5/2}}. \]

(4.44)

\[ A_R(t) = A_R(t_i). \]

(4.45)

These are nothing but the textbook orbit-averaged solutions (see, for instance, Sec. IV.1 in Ref. [16]). Thus, the difference between the DRG solutions and the orbit-averaged solutions are the sinusoidal terms in (4.33)–(4.35). Note that these terms do not have constant periods and thus orbit averaging will not set them strictly to zero. The lack of a definite period is another weakness of the averaging procedure [20,21]

As can be seen from (4.43), the quantity \(R_R^4(t) \Omega_R^2(t)\) is an invariant along the RG trajectory. This constant is just equal to \(M\). Other RG invariants can be found from these relations that are not so trivial, including

\[ R_R^4(t) + \frac{256\nu}{5} M^3 t = \text{constant} \]

(4.46)

\[ \Phi_R(t) + \frac{R_R^{3/2}(t)}{32\nu M^{5/2}} = \text{constant}. \]

(4.47)

The expressions in (4.42)–(4.45), combined with the renormalized solutions in (4.33)–(4.35), give the resummed solution to the 0PN inspiral dynamics valid up to times \(t - t_i\) of order \(1/(\nu v^5(t) \Omega_R(t))\). Note that the initial radial velocity depends on \(A_R(t_i)\) and \(\Phi_R(t_i)\) at this order via the relation

\[ \dot{r}(t_i) = A_R(t_i) \cos \Phi_R(t_i) - \frac{64\nu}{5} R_R^6(t_i) \Omega_R(t_i)^6. \]

(4.48)

For the purposes of comparison, we next find the numerical solution of an equal-mass compact binary inspiral where the total mass is \(M = 1\). Specifically, we choose the following initial data at \(t_i = 0\) for demonstration purposes:

\[ \phi(0) = 0 \]

\[ \omega(0) = 10^{-2}/M \]

\[ r(0) = (M/\omega(0)^2)^{1/3} = 10^{4/3} M \]

\[ \dot{r}(0) = 0. \]

(4.49)

Notice that the typical speed scale is \(v \sim r(0)\omega(0) \approx 0.2\) and \(v^5 \sim 5 \times 10^{-4}\), which are manageable numbers for numerical studies, which is why we have chosen them. To relate these initial conditions to the parameters \(R_R(t_i), \Omega_R(t_i), \Phi_R(t_i),\) and \(A_R(t_i)\), we set (4.30)–(4.32) and the time derivative of (4.30) at \(t_i = 0\) equal to the above initial data. This yields four equations in the four parameters, which we solve numerically. Recall that \(A_R(0) = e_R(0) R_R(0)\) is proportional to the initial eccentricity \(e_R(0)\), which we took to be \(O(v^5)\).

In Fig. 1, we compare the numerical solution (black) to our RG resummed solution (orange). The top left (right) panel shows these solutions for the orbital radius (phase). The bottom panels show the fractional errors for the orbital radius and phase solutions, respectively. The orange (blue) curves show the fractional errors between the RG resummed (adiabatic, orbit-averaged) and numerical solutions.

The adiabatic solutions come from solving the flux-balance equations, which are averaged over the orbital period [17]. It should be noted that the adiabatic solutions contain an ambiguity in specifying the initial data because of the orbit-averaging procedure, as discussed in Refs. [20,21], which can be seen at early times in the bottom right panel of Fig. 1. In addition, for orbits with larger eccentricities, it is not clear which oscillations the adiabatic approximation should remove (e.g., those parameterized by coordinate time, eccentric anomaly, true anomaly, or mean anomaly), which becomes important for periastron advance when PN corrections are included. As such, comparisons to the adiabatic approximation should be regarded as more qualitative rather than quantitative, perhaps. With these comments in mind, we remark that the DRG method provides a systematic procedure for deriving unambiguous predictions for the compact binary’s real-time evolution.

In Sec. V, we will improve the accuracy of the resummed perturbative solution to (4.2), especially at late times, by including second-order corrections in \(e\), which will induce...
an RG flow for the renormalized oscillation amplitude $A_R$ (i.e., the orbital eccentricity).

**D. Estimating errors of the resummed solutions**

The bare perturbative solutions in (4.13)–(4.14) are accurate up to $O(v^{10}_B)$ corrections when ignoring higher-order PN corrections that we did not originally include in the equations of motion in (4.2). When renormalizing the integration constants, the error being made in the perturbation theory is $O(v^{10}_R)$ because all bare parameters are written in terms of their renormalized values plus higher-order counterterms.

Next, we recall that $r(t)$, $\omega(t)$, and $\phi(t)$ are independent of $\tau$ so that differentiating the radial solution, for example, implies that

$$
0 = \frac{dR}{d\tau} + \frac{64}{5} v R^6_R \Omega^6_R + A_R \left( \frac{d\Phi}{d\tau} - \Omega_R \right) \cos(\Omega_R(t-\tau) + \Phi_R) + O(v^{10}_R R^6_R \Omega_R),
$$

(4.50)

where we have included the error term. The extra factor of $R_R$ in the error term is to ensure the correct dimensions and scaling for the radial solution, while the factor of $\Omega_R$ is the reciprocal of the orbital time scale from the $\tau$ derivative. Of course, the RG equations in (4.38) and (4.40) tell us that this is satisfied identically, but the error term implies that the RG equations should be written more completely as

$$
\frac{dR}{d\tau} = -\frac{64}{5} v R^6_R \Omega^6_R + O(v^{10}_R R^6_R \Omega_R)
$$

(4.51)

and similarly for the other ones. Therefore, the RG solutions are determined up to $O(v^{10}_R R^6_R (t-t_i))$ corrections, and the resummed perturbative solutions are valid until times

$$
t - t_i \sim \frac{1}{v^{10}_R \Omega_R} \sim \frac{1}{R^6_R \Omega_R^{11}}.
$$

(4.52)

Notice that this elapsed time of validity is measured with respect to the renormalized integration constants at the initial time $t_i$.

**V. GOING TO TWO LOOPS: TWO INSERTIONS OF RADIATION REACTION**

We now show how to include two insertions of the leading-order radiation reaction force. In doing so, we will show how to renormalize to $O(v^2)$ in the DRG formalism. Following the field theory terminology, we call this a “two-loop” calculation, despite the fact that all our calculations are more akin to “tree” level Feynman diagrams. Indeed, the DRG calculations can be couched in terms of these diagrams by thinking of the background (circular) orbit as a source insertion and treating $r(t)$ and $\omega(t)$ as two distinct one-dimensional fields. However, it is not clear that Feynman diagrams are of much utility for us, though they may help keep track of the systematics as one goes to higher orders.

We work in what is known as “bare perturbation theory.” In this way of organizing the calculation, we work with only bare parameters at arbitrary order and then fix the counterterms a posteriori, as we did in the previous section. We will see that at second order there will be a nontrivial set of consistency checks of the calculation.

\footnote{At leading order, the distinction between bare and renormalized perturbation theory is nominal.}
We begin by introducing the second-order notation

\[
\begin{align*}
  r & = R_B + \delta r(t) + \delta \kappa(t) \\
  \omega & = \Omega_B + \delta \omega(t) + \delta \rho(t),
\end{align*}
\]  

(5.1)

where \( \delta \kappa \sim v_B^{10} R_B \) and \( \delta \rho \sim v_B^{10} \Omega_B \sim v_B^{11} / R_B \) and the first-order solutions were calculated in the previous section. The equations of motion for \( \delta \kappa \) and \( \delta \rho \) are

\[
\begin{align*}
  \delta \kappa(t) - 3 \Omega_B^2 \delta \kappa(t) & = \frac{112}{15} \nu R_B^5 \Omega_B^6 \delta \dot{r}(t) - \frac{3 \Omega_B^2}{R_B} \delta r^2(t) + R_B \delta \omega^2(t) + 2 R_B \Omega_B \delta \rho(t) + 2 \Omega_B^3 \delta \rho(t) \delta \omega(t) \\
  \delta \rho(t) + \frac{2 \Omega_B}{R_B} \delta \kappa(t) & = -\frac{48}{5} \nu R_B^4 \Omega_B^5 \delta \omega(t) + 8 \nu R_B^4 \Omega_B^6 \delta r(t) - \frac{2}{R_B} \delta \omega(t) \delta \dot{r}(t) - \frac{2}{R_B} \delta r(t) \delta \dot{\omega}(t) - \frac{2 \Omega_B}{R_B} \delta r(t) \delta \dot{r}(t). 
\end{align*}
\]  

(5.2)

(5.3)

The solution for the second-order radial perturbation \( \delta \kappa \) is given by

\[
\begin{align*}
  \delta \kappa(t) & = -\frac{3 A_B^2}{2 R_B} + \frac{29696}{75} \nu^2 R_B^{11} \Omega_B^{10} - \frac{6144}{25} \nu^2 R_B^{11} \Omega_B^{12} (t-t_0)^2 + \frac{272}{5} \nu A_B R_B^5 \Omega_B^5 \cos \Phi_B + \frac{3 A_B^2}{R_B} \cos 2 \Phi_B \\
  & + \frac{3 A_B^2}{2 R_B} \cos \Omega_B (t-t_0) - \frac{29696}{75} \nu^2 R_B^{11} \Omega_B^{10} \cos \Omega_B (t-t_0) - \frac{32}{3} \nu A_B R_B^5 \Omega_B^5 \cos (\Phi_B - \Omega_B (t-t_0)) \\
  & - \frac{5 A_B^2}{4 R_B} \cos (2 \Phi_B - \Omega_B (t-t_0)) - \frac{656}{15} \nu A_B R_B^5 \Omega_B^5 \cos (\Phi_B + \Omega_B (t-t_0)) \\
  & + \frac{48}{5} \nu A_B R_B^5 \Omega_B^7 (t-t_0)^2 \cos (\Phi_B + \Omega_B (t-t_0)) + \frac{1 A_B^2}{2 R_B} \cos (2 \Phi_B + 2 \Omega_B (t-t_0)) \\
  & - \frac{9 A_B^2}{4 R_B} \cos (2 \Phi_B + \Omega_B (t-t_0)) - \frac{496}{15} \nu A_B R_B^5 \Omega_B^5 (t-t_0) \sin (\Phi_B + \Omega_B (t-t_0)),
\end{align*}
\]  

(5.4)

while that for \( \delta \rho \) is

\[
\begin{align*}
  \delta \rho(t) & = \frac{3 A_B^2 \Omega_B}{R_B} - \frac{59392}{75} \nu^2 R_B^{10} \Omega_B^{11} + \frac{16896}{25} \nu^2 R_B^{10} \Omega_B^{13} (t-t_0)^2 - \frac{408}{5} \nu A_B R_B^4 \Omega_B^6 \cos \Phi_B - \frac{9 A_B^2 \Omega_B}{2 R_B^2} \cos 2 \Phi_B \\
  & - \frac{3 A_B^2 \Omega_B}{R_B} \cos \Omega_B (t-t_0) + \frac{59392}{75} \nu^2 R_B^{10} \Omega_B^{11} \cos \Omega_B (t-t_0) + \frac{64}{3} \nu A_B R_B^4 \Omega_B^6 \cos (\Phi_B - \Omega_B (t-t_0)) \\
  & + \frac{5 A_B^2 \Omega_B}{2 R_B^2} \cos (2 \Phi_B - \Omega_B (t-t_0)) + \frac{904}{15} \nu A_B R_B^4 \Omega_B^6 \cos (\Phi_B + \Omega_B (t-t_0)) \\
  & - \frac{96}{5} \nu A_B R_B^4 \Omega_B^6 (t-t_0)^2 \cos (\Phi_B + \Omega_B (t-t_0)) - \frac{5 A_B^2 \Omega_B}{2 R_B^2} \cos (2 \Phi_B + 2 \Omega_B (t-t_0)) \\
  & + \frac{9 A_B^2 \Omega_B}{2 R_B^2} \cos (2 \Phi_B + 2 \Omega_B (t-t_0)) + \frac{32}{15} \nu A_B R_B^4 \Omega_B^7 (t-t_0) \sin (\Phi_B + \Omega_B (t-t_0)).
\end{align*}
\]  

(5.5)

As in the previous section for the one-loop calculation, we can shift the initial parameters so as to remove the redundant pieces that are finite (i.e., nonsecular) homogenous solutions. It is straightforward to show with some algebra and trigonometric identities that the shift...
removes the redundant, finite terms. We are then left with the following expressions for the perturbative solutions at $\mathcal{O}(\nu^{10})$,

\begin{align}
\delta \kappa(t) &= \frac{1}{2} \frac{A_B^2}{R_B^2} - \frac{29696}{75} \nu^2 R_{B}^{11} \Omega_{B}^{10} - \frac{6144}{25} \nu^2 R_{B}^{11} \Omega_{B}^{12} (t - t_0)^2 - \frac{656}{15} \nu A_B R_{B}^{5} \Omega_{B}^{5} \cos(\Phi_B + \Omega_B(t - t_0)) \\
&\quad + \frac{48}{5} \nu A_B R_{B}^{5} \Omega_{B}^{5} (t - t_0)^2 \cos(\Phi_B + \Omega_B(t - t_0)) + \frac{1}{2} \frac{A_B^2}{R_B^2} \cos(2\Phi_B + 2\Omega_B(t - t_0)) \\
&\quad - \frac{496}{15} \nu A_B R_{B}^{5} \Omega_{B}^{5} \sin(\Phi_B + \Omega_B(t - t_0)),
\end{align}

\begin{align}
\delta \rho(t) &= \frac{16896}{25} \nu^2 R_{B}^{10} \Omega_{B}^{13} (t - t_0)^2 + \frac{904}{15} \nu A_B R_{B}^{4} \Omega_{B}^{6} \cos(\Phi_B + \Omega_B(t - t_0)) \\
&\quad - \frac{96}{5} \nu A_B R_{B}^{4} \Omega_{B}^{6} (t - t_0)^2 \cos(\Phi_B + \Omega_B(t - t_0)) - \frac{5}{2} \frac{A_B^2}{R_B^2} \cos(2\Phi_B + 2\Omega_B(t - t_0)) \\
&\quad + \frac{32}{15} \nu A_B R_{B}^{4} \Omega_{B}^{6} \sin(\Phi_B + \Omega_B(t - t_0)),
\end{align}

which are easily shown to satisfy the equations of motion to the order in which we are working.

### A. Renormalization

Starting from the bare perturbative solutions for $\delta \kappa(t)$ and $\delta \rho(t)$, we next renormalize the initial parameters of the system to absorb the secular divergences as we did above at one loop.

Additional contributions enter at $\mathcal{O}(\nu^{10})$ that come from expanding out the bare parameters of the one-loop contribution to $r(t)$ as well as the $\delta r^{10}$ counterterm that comes from the background piece, $R_B$. The totality of those pieces together with the expression for $\delta \kappa(t)$ gives the full renormalized $\nu^{10}$ contribution to $r(t)$, which we call $r^{10}(t)$. Using the expressions for the one-loop counterterms given in (4.26), (4.27), and (4.29) and introducing the renormalization scale $\tau$ through $t - t_0 = (t - \tau) + (\tau - t_0)$, we find that the full $\mathcal{O}(\nu^{10})$ contribution to the perturbative radial solution is

\begin{align}
r^{10}(t) &= \frac{1}{2} \frac{A_B^2}{R_B^2} - \frac{29696}{75} \nu^2 R_{B}^{11} \Omega_{B}^{10} - \frac{6144}{25} \nu^2 R_{B}^{11} \Omega_{B}^{12} [(t - \tau)^2 - (t - t_0)^2] - \frac{656}{15} \nu A_B R_{B}^{5} \Omega_{B}^{5} \cos(\Phi_R + \Omega_R(t - \tau)) \\
&\quad + \frac{48}{5} \nu A_B R_{B}^{5} \Omega_{B}^{5} [(t - \tau)^2 \cos(\Phi_R + \Omega_R(t - \tau)) + \frac{1}{2} \frac{A_B^2}{R_B^2} \cos(2\Phi_R + 2\Omega_R(t - \tau)) \\
&\quad - \frac{496}{15} \nu A_B R_{B}^{5} \Omega_{B}^{5} \sin(\Phi_R + \Omega_R(t - \tau)) + \delta_R^{10} + \delta_A^{10} \sin(\Phi_R + (t - \tau)\Omega_R),
\end{align}

where, as usual, we are ignoring terms that are beyond $\nu^{10}$.

\footnote{These shifts have some freedom parametrized by a constant $\mu$ that should be fixed. For the shifts in (5.6), we have chosen a scheme so as to keep the resulting two-loop RG equations as simple as possible, which is equivalent to choosing $\mu$ so as to remove all of the finite, $t$-independent pieces in $\delta \rho(t)$. Of course, one is free to choose other values for $\mu$, which changes the ensuing RG equations and perturbative expressions but in a way that does not change the predictions for the physical quantities, $r(t)$ and $\phi(t)$.}
The $O(t^{10})$ counterterms for $R$ and $A$ are given by

$$
\delta\tau_R^{10} = -\frac{6144}{25} \nu^2 R^4_R \Omega_R^3 (t-t_0)^2,
$$

$$
\delta\tau_A^{10} = \frac{496}{15} A_R \nu R^5_R \Omega_R^6 (t-t_0).
$$

In calculating (5.9), we encounter terms proportional to $(t-t)(\tau-t_0)$ between the linear and quadratic terms. The fact that such cross-terms cancel when using the expressions for the one-loop counterterms constitutes a consistency check because otherwise there would be residual $t_0$ contributions surviving that would be akin to having a “nonrenormalizable” field theory. The renormalized, finite contribution to the second-order radial perturbation is then given by

$$
\omega_{\Omega}^{(2)}(t) = \frac{16896}{25} \nu^2 R^4_R \Omega_R^3 [(t-t)^2 - (\tau-t_0)^2] + \frac{904}{15} A_R R^4_R \nu \Omega_R^6 \cos(\Phi_R + \Omega_R (t-t))
$$

$$
- \frac{96}{5} \nu A_R R^4_R \Omega_R^6 (t-t)^2 \cos(\Phi_R + \Omega_R (t-t)) - \frac{5 A_R^2 \Omega_R}{2 R^2_R} \cos(2\Phi_R + 2 \Omega_R (t-t))
$$

$$
+ \frac{32}{15} \nu A_R R^4_R \Omega_R^6 (t-t) \sin(\Phi_R + \Omega_R (t-t)) + \delta\tau^{10}_\Omega.
$$

Notice, again, that in adding these contributions to (5.14) we encounter nontrivial cancellations. In particular, the terms proportional to sinusoids cancel exactly, as they must since there is no counterterm of this form. The only remaining secular divergence appears in the third term of the first line and is quadratic in $(\tau-t_0)^2$. We identify this term with $\delta\tau^{10}_\Omega$, which gives the full renormalized $t^{10}$ contribution to the perturbative angular frequency solution is

$$
\delta\tau^{10}_\Omega = \frac{16896}{25} \nu^2 R^4_R \Omega_R^3 (t-t_0)^2.
$$

so that we are left with, after scale setting $\tau = t$,

$$
\omega_{\Omega}^{(2)}(t) = \frac{904}{15} \nu A_R R^4_R \nu \Omega_R^6 \cos(\Phi_R + \Omega_R (t-t)) - \frac{5 A_R^2 \Omega_R}{2 R^2_R} \cos(2\Phi_R).
$$

As with the one-loop calculation, we can calculate the second-order contribution to the orbital phase, $\delta\sigma(t)$, by integrating $\delta\phi$ over time,

$$
\delta\sigma(t) = -\frac{504}{5} \nu A_R R^4_R \nu \Omega_R^6 \sin(\Phi_B) + \frac{5 A^2_R}{4 R_B^2} \sin(\Phi_B) + \frac{504}{5} \nu A_R R^4_R \Omega_R^6 \sin(\Phi_B + \Omega_B (t-t_0))
$$

$$
+ \frac{5632}{25} \nu^2 R^4_R \Omega_B^3 (t-t_0)^3 - \frac{608}{15} \nu A_R R^4_R \Omega_R^6 (t-t_0) \cos(\Phi_B + \Omega_B (t-t_0))
$$

$$
- \frac{96}{5} \nu A_R R^4_R \Omega_R^6 (t-t_0)^2 \sin(\Phi_B + \Omega_B (t-t_0)) - \frac{5 A^2_R}{4 R_B^2} \sin(2\Phi_B + 2 \Omega_B(t-t_0)).
$$

Proceeding as before, we find the $O(\nu^{10})$ contribution to the phase to be
\[ \phi_v(t) = \delta_\phi^{(0)} + \frac{504}{5} \nu A_R R_R^4 \Omega_R^5 \sin \Phi_B(t_0) + \frac{5 A_R^2}{4 R_R^2} \sin 2 \Phi_B(t_0) + \frac{5632}{25} \nu^2 R_R^{10} \Omega_R^{13} (\tau - t_0)^3 + \frac{5632}{25} \nu^2 R_R^{10} \Omega_R^{13} (t - \tau)^3 - \frac{608}{15} \nu A_R R_R^4 \Omega_R^5 (t - \tau) \cos (\Phi_R + \Omega_R (t - \tau)) - \frac{96}{5} \nu A_R R_R^4 \Omega_R^5 (t - \tau)^2 \sin (\Phi_R + \Omega_R (t - \tau)) + \frac{504}{5} \nu A_R R_R^4 \Omega_R^5 \sin (\Phi_R + \Omega_R (t - \tau)) - \frac{5 A_R^2}{4 R_R^2} \sin (2 \Phi_R + 2 \Omega_R (t - \tau)). \] (5.16)

We choose the \( O(v^{10}) \) phase counterterm \( \delta_\phi^{(0)} \) to cancel the last three terms in the first line of the equation above,

\[ \delta_\phi^{(0)} (\tau, t_0) = \frac{504}{5} \nu A_R R_R^4 \Omega_R^5 \sin \Phi_B(t_0) - \frac{5 A_R^2}{4 R_R^2} \sin 2 \Phi_B(t_0) + \frac{5632}{25} \nu^2 R_R^{10} \Omega_R^{13} (\tau - t_0)^3. \] (5.17)

The resulting expression for the \( O(v^{10}) \) phase at \( \tau = t \) is then given by

\[ \phi_v(t) = \frac{504}{5} \nu A_R R_R^4 \Omega_R^5 \sin \Phi_R - \frac{5 A_R^2}{4 R_R^2} \sin 2 \Phi_R. \] (5.18)

**B. Renormalization group solution**

Putting together the order \( \epsilon \) and \( \epsilon^2 \) counterterms, we have

\[ \delta_R = \frac{64 \nu}{5} R_R^6 \Omega_R^6 (\tau - t_0) - \frac{644}{25} \nu^2 R_R^{11} \Omega_R^{12} (\tau - t_0)^2 \]

\[ \delta_\Omega = - \frac{96 \nu}{5} R_R^6 \Omega_R^7 (\tau - t_0) + \frac{16896}{25} \nu^2 R_R^{10} \Omega_R^{13} (\tau - t_0)^2 \]

\[ \delta_A = \frac{496}{15} A_R \nu R_R^4 \Omega_R^6 (\tau - t_0) \]

\[ \delta_\phi = - \Omega_R (\tau - t_0) + \frac{48 \nu}{5} R_R^5 \Omega_R^7 (\tau - t_0)^2 - \frac{5632}{25} \nu^2 R_R^{10} \Omega_R^{13} (\tau - t_0)^3 + \frac{504}{5} \nu A_R R_R^4 \Omega_R^5 \sin \Phi_B(t_0) - \frac{5 A_R^2}{4 R_R^2} \sin 2 \Phi_B(t_0). \] (5.19)

From the expressions relating the bare parameters to the renormalized quantities and counterterms, we derive the RG equations through two loops. The RG equation for \( R_R \) through two loops is given by

\[ 0 = \frac{d}{d \tau} R_R(t_0) \]

\[ = \frac{dR_R(\tau)}{d \tau} + \frac{64 \nu}{5} R_R^6 \Omega_R^6 + \frac{384}{5} \nu R_R^5 \Omega_R^7 (\tau - t_0) \times \left( \Omega_R \frac{dR_R(\tau)}{d \tau} + R_R \frac{d\Omega_R(\tau)}{d \tau} \right) - \frac{12288}{25} \nu^2 R_R^{11} \Omega_R^{12} (\tau - t_0) + O(v^{15}). \] (5.20)

\[ \frac{d}{d \tau} A_R(t_0) = - \frac{496}{15} A_R \nu R_R^4 \Omega_R^6. \] (5.24)

Through \( O(v^{10}) \), the beta function for the amplitude of oscillation receives a two-loop correction and induces a nontrivial RG flow described by

\[ \frac{d}{d \tau} A_R(t) = - \frac{496}{15} A_R \nu R_R^4 \Omega_R^6, \]

which has the solution

\[ A_R(t) = A_R(t_0) \left( \frac{R_R(t)}{R_R(t_0)} \right)^{31/12} \Rightarrow e_R(t) = \frac{A_R(t)}{R_R(t)} \]

\[ = e_R(t_0) \left( \frac{R_R(t)}{R_R(t_0)} \right)^{19/12}. \] (5.25)

\[ ^5 \text{In the sense that the result depends upon the cutoff, } t_0. \]
where \( e_R(t) \) is the \( O(v_R^5) \) time-dependent eccentricity of the binary’s orbit from (4.9). The power-law relation between \( e_R \) and \( R_R \), namely, \( R_R \sim e_R^{12/19} \) agrees with the well-known result from Peters [29] when \( e_R \ll 1 \).

Finally, the RG equation through \( O(v^{10}) \) for the phase parameter satisfies

\[
0 = \frac{d}{d\tau} \Phi_R(t_0) = \frac{d\Phi_R}{d\tau} - \Omega_R - (\tau - t_0) \frac{d\Omega_R}{d\tau} + \frac{96}{5} \nu R_R^4 \Omega_R^5 (\tau - t_0)
\]

\[
+ 48 \nu R_R^4 \Omega_R^5 (\tau - t_0)^2 \left( 5 \Omega_R \frac{dR_R}{d\tau} + 7 R_R \frac{d\Omega_R}{d\tau} \right)
\]

\[
- \frac{16896}{25} \nu^2 R_R^{10} \Omega_R^{13} (\tau - t_0)^2 + O(v^{15}).
\] (5.26)

Note that the last two terms in (5.19) for \( \delta_\phi \) do not contribute at this order. Again, solving this iteratively, we find that the beta function is independent of the regulator \( t_0 \), leaving us with

\[
\frac{d\Phi_R}{d\tau} = \Omega_R.
\] (5.27)

Therefore, through two loops, we see that the RG equations for \( R_R \), \( \Omega_R \), and \( \Phi_R \) are the same as at one loop. However, the beta function at two loops for the eccentricity receives a nontrivial contribution that induces an RG flow for \( e_R \) in time.

Given our solutions to the RG equations, we may now write down the result for the resummed orbital coordinates through \( O(v^{10}) \),

\[
r(t) = R_R(t) \left( 1 + e_R(t) \sin \Phi_R(t) + \frac{1}{2} e_R^2(t) - \frac{29696}{75} \nu^2 R_R^{10} (t) \Omega_R^{10}(t) \right)
\]

\[
- \frac{656}{15} \nu e_R(t) R_R^5(t) \Omega_R^5(t) \cos \Phi_R(t) + \frac{1}{2} e_R^2(t) \cos 2\Phi_R(t) + O(v_R^{15}) \right) \] (5.28)

\[
o(t) = \Omega_R(t) \left( 1 - 2 e_R(t) \sin \Phi_R(t) + \frac{904}{15} \nu e_R(t) R_R^5(t) \Omega_R^5(t) \cos \Phi_R(t) - \frac{5}{2} e_R^2(t) \cos 2\Phi_R(t) + O(v_R^{15}) \right) \] (5.29)

\[
\phi(t) = \Phi_R(t) + 2 e_R(t) \cos \Phi_R(t) + \frac{504}{5} \nu e_R(t) R_R^5(t) \Omega_R^5(t) \sin \Phi_R(t) - \frac{5}{4} e_R^2(t) \sin 2\Phi_R(t) + O(v_R^{15}) \right) \] (5.30)

where we have included the error terms, which can be derived as discussed in Sec. IV D, and have written \( A_R = e_R R_R \). The expressions for the two-loop renormalized initial conditions are given in (4.42)–(4.44) and (5.25).

It is straightforward to show that these resummed perturbative solutions satisfy the equations of motion through \( O(v^{10}) \) and that \( dr(t)/dt = o(t) \) to the same order. In the case where the initial data is fine-tuned so as to yield a quasicircular inspiral [i.e., by setting \( A_R(t_i) = e_R(t_i) = 0 \)], the resummed solutions become

\[
r_{qc}(t) = R_R(t) - \frac{29696}{75} \nu^2 R_R^{10} (t) \Omega_R^{10}(t)
\]

\[
o_{qc}(t) = \Omega_R(t)
\]

\[
\phi_{qc}(t) = \Phi_R(t).
\] (5.31)

Figure 2 shows the fractional errors between the numerical solution of (4.2) and the one-loop (blue) and two-loop (orange) resummed solutions for the orbital radius (top panel) and phase (bottom panel). We observe a marked global improvement in the two-loop resummed solution for \( r(t) \), providing at least an order of magnitude better...
VI. BEYOND LEADING ORDER RADIATION REACTION

In this formalism, the inclusion of higher-order radiation reaction forces is straightforward. The equations of motion through the 1PN correction to radiation reaction forces [5,6] are given by

\[
\ddot{r} - ro^2 = -\frac{M}{r^2} + \frac{64M^3\nu}{15r^4}\dot{r} + \frac{16M^2\nu}{5r^3}(\dot{r}^2 + \dot{r}^2\omega^2)
- \frac{8M^2\nu}{105r^5}(821 + 210\nu)r
+ \frac{8M^2\nu}{105r^4}\dot{r}((-362 + 245\nu)r^2\omega^2 - 775r^2)
- \frac{4M^2\nu}{35r^5}\dot{r}((-65 + 84\nu)r^4\omega^4
+ (59 + 84\nu)r^2\omega^2r^2 + 54\dot{r}^4) \tag{6.1}
\]

and

\[
\ddot{\omega} + 2\dot{\omega} = -\frac{24M^3\nu}{5r^3}\omega - \frac{8M^2\nu}{5r^2}\omega(\dot{r}^2 + \dot{r}^2\omega^2)
+ \frac{4M^4\nu}{105r^4}(1325 + 546\nu)\omega
- \frac{2M^3\nu}{105r^5}(2(205 + 777\nu)r^2\omega^2
- (1025 + 1414\nu)r^2)
+ \frac{2M^2\nu}{35r^6}\omega((313 + 42\nu)r^4\omega^4
- (1747 - 42\nu)r^2\omega^2r^2 + 40\dot{r}^4). \tag{6.2}
\]

We are interested in demonstrating how to handle higher PN-order secular terms in DRG, so we do not include the 1PN or higher potentials here, which do not (directly) generate secularly diverging perturbations. Of course, a fully consistent orbital solution should include all potentials that contribute to a given PN order.

As done in the previous section, we expand the solution around the background including perturbations up to order \(v_B^7\). Following (5.1), where now \(\delta^t \sim v_B^7R_R\) and \(\delta^\rho \sim v_B^5\Omega_R \sim v_B^7/R_B\), we find that the perturbed radial and angular frequency solutions contain the following contributions at this order:

\[
r(t) \supset -\frac{4\nu}{105}R_B^7\Omega_R^8(336\nu - 3179)(t - t_0) \tag{6.3}
\]

\[
\omega(t) \supset \frac{2\nu}{35}R_B^7\Omega_R^8(336\nu - 3179)(t - t_0) \tag{6.4}
\]

\[
\phi(t) \supset \frac{1}{35}\nu R_B^7\Omega_R^8(336\nu - 3179)(t - t_0)^2. \tag{6.5}
\]

At this order, there is no mixing between the subleading corrections (i.e., \(\delta^\rho\delta^t\)) and the leading \(v^5\) pieces (\(\delta^t\) and \(\delta^\omega\)). As such, there are no quadratic divergences in \(r(t)\) and \(\omega(t)\).

The associated counterterms then lead to the following RG equations:

\[
\frac{dR_R}{dt} (\tau) = -\frac{64\nu}{5}R_B^5\Omega_R^6
- \frac{4\nu}{105}(336\nu - 3179)R_B^3\Omega_R^6 \tag{6.6}
\]

\[
\frac{d\Omega_R}{dt} (\tau) = \frac{96\nu}{5}R_B^3\Omega_R^4
+ \frac{2\nu}{35}(336\nu - 3179)R_B^1\Omega_R^4 \tag{6.7}
\]

\[
\frac{d\Phi_R}{dt} (\tau) = \Omega_R. \tag{6.8}
\]

The exact solutions to the frequency and phase RG equations are

\[
\Omega_R (t) = \Omega_R (t_i) \frac{R_R (t)}{R_R (t_i)} ^ {1/2} = \frac{M^{1/2}}{R_R (t)} \tag{6.9}
\]

\[
-\frac{32\nu}{5}M^{5/2} (\Phi_R (t) - \Phi_R (t_i))
= \frac{1}{5} \left( R_B^{5/2} (t) - R_B^{5/2} (t_i) \right) + \frac{1}{3} \alpha M (R_B^{3/2} (t) - R_B^{3/2} (t_i))
+ \frac{1}{3} M^2 (R_B^{1/2} (t) - R_B^{1/2} (t_i))
- \alpha^{5/2} M^{5/2} \left( \tanh^{-1} \sqrt{\frac{R_B (t)}{\alpha M}} - \tanh^{-1} \sqrt{\frac{R_B (t_i)}{\alpha M}} \right). \tag{6.10}
\]

where

\[
\alpha = \frac{3179}{336} - \nu \approx 9.5 - \nu \tag{6.11}
\]

and we have used the fact that the combination \(R_B^3\Omega_R^6 = M\) is an RG invariant. Here and below, we choose the RG scale \(\tau\) to be the observation time \(t\).

The solution to the radial RG equation is found by first writing it as

\[
\frac{R_R^4}{R_R - \alpha M} dR_R = -\frac{64}{5}M^3\nu dt. \tag{6.12}
\]

Integrating both sides gives the exact but implicit relation,
\[-\frac{64\nu}{5} M^3 (t - t_i) \]
\[= \frac{1}{4} (R^\nu_R(t) - R^\nu_R(t_i)) + \frac{1}{3} \alpha M (R^3_R(t) - R^3_R(t_i)) \]
\[+ \frac{1}{2} \alpha^2 M^2 (R^2_R(t) - R^2_R(t_i)) + \alpha^2 M^3 (R^1_R(t) - R^1_R(t_i)) \]
\[+ \alpha^4 M^4 \log \left( \frac{R_R(t) - \alpha M}{R_R(t_i) - \alpha M} \right) \]  
(6.13)

Note that setting $\alpha = 0$ in these RG solutions recovers the one-loop 0PN results derived in the previous sections. The RG solution for the one-loop oscillation amplitude is also given exactly at this order by

$$A_R(t) = A_R(t_i) = \text{constant}. \quad (6.14)$$

A two-loop calculation would induce a nontrivial RG flow for the eccentricity as in the 0PN example in the previous section.

**VII. CONCLUSION**

In this work, we have shown how to utilize the dynamical renormalization group formalism to solve for the long-time behavior for binary inspirals by systematically resumming secularly growing perturbations. By utilizing this formalism, one can avoid the ambiguities intrinsic to using the adiabatic approximation and orbit averaging [20,21]. We generated an analytic form for the trajectory of an inspiral at second order in the leading (2.5PN) radiation reaction force. At this order, there exist highly nontrivial consistency checks of the formalism. In particular, it must be that all secular divergences have the right functional form to be absorbable into the initial conditions for the orbit. This attribute is called renormalizability in the context of field theory. Since this formalism solves the equations of motion directly (i.e., without appealing to any kind of averaging procedure), then to go beyond 1PN accuracy, one would require the 2PN correction to the radiation reaction force.

Perhaps the most fertile ground for this formalism is in spin dynamics, where finding closed form solutions becomes a significant challenge. In nearly all studies of spin effects on compact binary inspiral evolutions, the equations describing the components of the spin vectors are orbit averaged and, more recently, precession averaged [13,14]. However, the dynamical renormalization group does not require averaging over short time scales in the problem to render the problem more amenable for solving, whether analytically or numerically. Instead, a naive perturbative solution of the full, nonaveraged equations of motion for the binary’s orbital coordinates and spin vectors is the starting point for the dynamical renormalization group method. Incorporating spin effects will be the subject of a companion paper [30].

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