Numerical analysis of the self-energy in covariant loop quantum gravity

Pietropaolo Frisoni[®],^{1,*} Francesco Gozzini[®],^{2,†} and Francesca Vidotto^{®1,3,‡}

¹Department of Physics & Astronomy, Western University, London, Ontario N6A 3K7, Canada

²Aix Marseille University, Université de Toulon, CNRS, CPT, 13288 Marseille, France ³Department of Philosophy and Rotman Institute, Western University, London, Ontario N6A 3K7, Canada

(Received 17 January 2022; accepted 12 April 2022; published 20 May 2022)

We study numerically the first order radiative corrections to the self-energy, in covariant loop quantum gravity. We employ the recently developed sl2cfoam-next spinfoam amplitudes library and some original numerical methods. We analyze the scaling of the divergence with the infrared cutoff, for which previous analytical estimates provided widely different lower and upper bounds. Our findings suggest that the divergence is approximately linear in the cutoff. We also investigate the role of the Barbero-Immirzi parameter in the asymptotic behavior, the dependence of the scaling on some boundary data, and the expectation values of boundary operators.

DOI: 10.1103/PhysRevD.105.106018

I. INTRODUCTION

The spinfoam, or covariant, formulation of loop quantum gravity (LQG) [1], is based on the definition of truncated amplitudes that define the background independent dynamics of quantum spacetime [2,3] in a way similar to the Feynman graphs' definition of the dynamics of quantum field theory. In a Feynman graph, radiative corrections appear as loops (set of propagators forming a circle) and may give rise to ultraviolet divergences. In a spinfoam, radiative corrections appear as bubbles (set of faces forming a sphere) and may give rise to infrared divergences when the cosmological constant is taken to be zero.¹ Here, we investigate the first order correction to a single edge of the spinfoam, which is analogous to the first order contribution to the self-energy in quantum field theory, and hence is referred to as a self-energy radiative correction. We use the version of the amplitude developed in Refs. [7-15] generally denoted as the EPRL amplitude. For reviews and notation, see Refs. [16–18].

The spinfoam infrared divergences have been studied analytically from a number of different perspectives [19– 28]. EPRL divergences have been studied in the Euclidean [29,30] and Lorentzian theories [31,32]. In particular, the last two references by Riello and Doná provide lower and upper bounds to the degree of divergence of the self-energy, resulting in a wide window. In Ref. [31], Riello gives a detailed analysis of the critical point that contributes to the

*pfrisoni@uwo.ca

amplitude in the large-spin scaling limit, based on the techniques introduced in Ref. [33], in the case of nondegenerate geometric configurations. He finds a logarithmic divergence in the cutoff K. He then estimates in a final appendix that the contribution of the degenerate sector might be linear in K, hence dominating the total divergence, postponing a detailed analysis of this sector to future studies. On the other hand, in Ref. [32], Doná disregards the interference between the different terms in the sum, and this allows him to estimate an upper bound on the divergence that scales as K^9 . He also proves that the upper bounds derived with his algorithm provide excellent estimates of different kinds of infrared divergences restricted to the three-dimensional model, suggesting that this might be also true in four dimensions, underlining the need for a complete numerical analysis for future studies.

These two results leave a window of possibilities which spans more than nine powers of the cutoff. In this work, we address the issue numerically. What we find is consistent with Riello's rough estimate: for small values of the Barbero-Immirzi parameter, the numerical analysis is compatible with a *linear* divergence in K.

The numerical calculation is possible thanks to the recently developed sl2cfoam-next library [34]. Our computations were performed on the clusters Cedar and Graham of Compute Canada (www.computecanada .ca) and on the Mesocentre of Aix-Marseille University (mesocentre.univ-amu.fr). The analysis and visualization of the computed data have been done with Julia [35] and *Mathematica* [36].

In Sec. II, we describe the structure of the self-energy amplitude. In Sec. III, we recall the calculation of the degree of divergence in topological SU(2) BF model, which can be found exactly. This allows us to compare

gozzini@cpt.univ-mrs.fr

[‡]fvidotto@uwo.ca

¹The general definition of the spinfoam amplitudes includes a nonzero cosmological constant [4–6] that plays the role of an infrared cutoff and makes the amplitudes finite.



FIG. 1. The two-complex C of the self-energy. All vertices and edges but only one internal and two external faces are depicted.

the numerical results with the analytical ones in the literature. In Sec. IV, we define the EPRL amplitude. In Sec. V, we describe the results of the numerical analysis of the EPRL self-energy amplitude. Finally, in the Appendixes, we report some further numerical studies and consistency checks.

II. SELF-ENERGY

Consider two 4-simplices joined by four tetrahedra. The boundary of this cellular complex is formed by two tetrahedra. In quantum gravity, this cellular complex can be associated to the transition from a single quantum of space to a single quantum of space, via a splitting into four virtual quanta, which then recombine [2,3]. Notice the similarity with the self-energy Feynman graph in a quantum field theory. This, however, is only an analogy; there is no notion of energy or radiation involved here. This graph should be understood as a higher order term in the spinfoam expansion.

The dual of this cellular complex is a 2-complex C. It is formed by two pentavalent vertices v_{\pm} joined by four internal edges e_a , with a = 1, 2, 3, 4. Each of the two vertices has one other edge e_{\pm} , ending on the boundary. The 2-complex has four external faces f_a that end on the boundary and six internal faces f_{ab} , dual to the surfaces that separate the four internal tetrahedra. See Fig. 1. The six faces f_{ab} , taken together, form the bubble we are interested in.

The boundary of the 2-complex C is the graph Γ formed by the two 4-valent nodes n_{\pm} where the edges e_{\pm} end (dual to the initial and final quanta of space), joined by four links l_a that bound the faces f_a (see Fig. 2). The kinematical LQG boundary Hilbert space is therefore $\mathcal{H}_{\Gamma} = L_2[SU(2)^4/SU(2)^2]_{\Gamma}$. The graph Γ is represented in Fig. 2.

A basis in \mathcal{H}_{Γ} is given by the spinnetwork states $|j_a, i_{\pm}\rangle$, where j_a are the spins attached to the links l_a . In the two intertwiner spaces, we choose the recoupling basis i_{\pm} defined by pairing the links j_1 and j_2 at the node n_{\pm} , i.e., the basis that diagonalizes the modulus square of the sum of the SU(2) generators in the SU(2) representation j_1 and j_2 [37–39]. The label *i* indicates the spin of the virtual link in this recoupling basis.



FIG. 2. The boundary graph.

We focus on the subspaces $\mathcal{H}_j \subset \mathcal{H}_{\Gamma}$ where the four spins are equal $j_a = j$ and denote the basis states as

$$|j, i_{\pm}\rangle = |j, i_{+}\rangle \otimes |j, i_{-}\rangle. \tag{1}$$

We are interested in the EPRL amplitude [16]

$$W(j, i_{\pm}) = \langle W_{\mathcal{C}} | j, i_{\pm} \rangle \tag{2}$$

relative to the 2-complex C. To compute this amplitude numerically, we need an infrared cutoff to bound the sum over the spins associated to the internal faces. We denote this cutoff K, and we define the cutoff amplitude as

$$W(j, i_{\pm}, K) = \langle W_{\mathcal{C}, K} | j, i_{\pm} \rangle.$$
(3)

We are interested in studying the dependence of this amplitude on K.

III. BF AMPLITUDE

As a first step, we consider the amplitude

$$W_{\rm BF}(j,i_{\pm}) = \langle W_{\mathcal{C}}^{\rm BF} | j, i_{\pm} \rangle \tag{4}$$

in the topological SU(2) BF theory, where the amplitude can be computed analytically [32]. The BF self-energy amplitude can be written as

$$W_{\rm BF}(j, i_{\pm}) = \sum_{j_{ab}, i_a} d_{\{j_{ab}\}} d_{\{i_a\}} \prod_{\pm} \{15j\}_{j_a, j_{ab}, i_{\pm}, i_a}, \quad (5)$$

where $d_{\{j_{ab}\}} = \prod_{(a,b)} (2j_{ab} + 1)$ and $d_{\{i_a\}} = \prod_a (2i_a + 1)$, with $a, b = 1, ...4, a \neq b$. The BF vertex amplitude $\{15j\}_{j_{AB},i_A}$ is the 15-j Wigner symbol defined contracting the five intertwiners i_A with the ten spins J_{AB} , with $A, B = 1, ...5, A \neq B$. See Appendix C for notation and conventions. A graphical representation of this amplitude that will be useful later on is the following:



Here, for completeness, we should include in the expression for the amplitude the dimensional factor of the boundary faces and intertwiners, but this is a constant overall factor that can be ignored in the analysis since it does not affect the scaling of the divergence; that is, what interests us. In (6), the six internal faces j_{ab} are highlighted in red. If we choose a recoupling basis for the internal intertwiners as



then this picture gives directly the expression of the two BF vertex amplitudes in terms of Wigner 3-j symbols: each intersection of three lines is one such symbols, and the lines specify the contractions' scheme

The multiple sum over the spins j_{ab} and the virtual spins i_a of the intertwiners are constrained by the Mandelstam identities that need to be satisfied for the Wigner symbols not to vanish. But it is clear from the last picture that the fixed values of the external spins j_a do not prevent the internal spins j_{ab} for growing arbitrarily large.

This is the origin of the divergence. To study it numerically, we introduce a cutoff $j_{max} = K$ on the six internal faces j_{ab} , therefore bounding all the sums by K. That is, we define



FIG. 3. Log-log plot of the BF self-energy amplitude as a function of the cutoff *K* on the internal spins j_{ab} . The continuous curve represents the function $4.8 \cdot K^9$.

$$W_{\mathrm{BF}}(j, i_{\pm}; K) = \sum_{j_{ab} \leq K \atop i_a} d_{\{j_{ab}\}} d_{\{i_a\}} \prod_{\pm} \{15j\}_{j_a, j_{ab}, i_{\pm}, i_a}.$$

We have computed this quantity numerically for a fixed value of the j_a and i_{\pm} , for different values of K. In the BF case, it is known analytically that the degree of divergence does not depend on j_a and i_{\pm} . The result is shown in Fig. 3, in which the continuous curve is $W_{\rm BF} = 4.8 \cdot K^9$.

This result is consistent with the analytical result [32], which gives

$$W_{\rm BF}(j, i_{\pm}; K) \propto K^9. \tag{8}$$

We briefly recall how to derive this result. Consider the simple case where all external spins are zero. The BF amplitude can be expressed in Fourier transform as a multiple integral over SU(2), given by one integration per link and one delta function per face. In this case, we have

$$W_{\rm BF} = \int \mathrm{d}g_a \prod_{a>b} \delta(g_a g_b^{-1}).$$

It is easy to see that one integration is redundant, and the other three can be used to eliminate three delta functions, leaving

$$W_{\rm BF} = \delta(1)^3. \tag{9}$$

To regularize this expression, one can write the delta function as a sum over characters χ^j of irreducible representations of SU(2) and introduce an infrared cutoff *K*, as follows:



FIG. 4. Log-log plot of the total number of different configurations of internal spins j_{ab} to be summed over for increasing values of the cutoff K. The power-law fit is ~1.7 · K^{5.6}.

$$\delta(U) = \lim_{K} \delta_K(U) = \lim_{K} \sum_{j=1}^{K} (2j+1)\chi^j(U).$$

Then, easily, $\delta_K(1) \approx K^3$, and (8) follows from (9).

The numerical analysis in Fig. 3 shows that the asymptotic scaling of the self-energy BF divergence (8) is reached for very low values of the infrared cutoff K. The computation involved in producing Fig. 3 is nowadays trivial and can be easily performed on a laptop thanks to very efficient routines for computing Wigner (3,6,9)-j symbols [40,41]. However, we note that the computational complexity grows for a large power of the cutoff index K, approximately equal to the number of internal faces of the foam. We plot in Fig. 4 the total number of combinations of internal spins that must be computed for each value of the cutoff K, till K = 20. We can see that the triangular inequalities reduce the complexity from K^6 to effectively $\sim K^{5.6}$. This still gives a large number of configurations to consider and provides a considerable numerical challenge in the EPRL case, whose vertex amplitude is much more computationally complex than the BF one, as we show in the next section.

IV. EPRL

Let us now come to the quantum gravity theory. The EPRL vertex amplitude is defined using a restriction of the unitary irreducible representations in the principal series of $SL(2, \mathbb{C})$ [12,13]. In order to evaluate it numerically, a more tractable formulation has been derived in Ref. [42] and discussed in Refs. [43,44]. In this formulation, the vertex is expressed as a sum over SU(2) 15-j symbols weighted by one "booster function" *B*4 per edge (see the Appendix C 2 for explicit formulas). The amplitude of a vertex bounded by spins j_f and intertwiners i_e can thus be written as a sum,

$$V(j_f, i_e, \gamma) = \sum_{\substack{l_f \ge j_f \\ k_e}} d_{\{k_e\}} \left(\prod_e B_4(j_f, l_f; i_e, k_e; \gamma) \right) \{15j\}_{l_f, k_e},$$
(10)

with virtual intertwiners k_e and virtual spins $l_f \ge j_f$, e = 2...5. One booster function must be replaced by the identity, corresponding to the suppression of a redundant SL(2, C) integration.² Notice that we keep the dependence on the Barbero-Immirzi parameter explicit. In this expression, the infinite sum over l_f , called the "shell expansion," can be shown analytically to converge, but in view of the numerical calculation, it is convenient to truncate it at a finite value defining

$$V(j_f, i_e, \gamma) = \lim_{\Delta l \to \infty} V(j_f, i_e, \gamma; \Delta l),$$
(11)

where



As before, we have neglected dimensional factors attached to the boundary intertwiners i_e and to the boundary spins j_f . In the following, we will refer to the cutoff Δl as the number of *shells*. Even if the convergence of the amplitudes a function of Δl is assured [43,44], it is not possible to have a unique prescription to set the optimal Δl to get an acceptable convergence, since it depends on the details of data such as the face spins j_f and the Barbero-Immirzi parameter. In addition, the convergence depends strongly on the structure of the 2-complex, and there is no known general procedure to estimate the error made in truncating the sum over the auxiliary spins. In the next section, we describe a technique to extrapolate the limit $\Delta l \rightarrow \infty$ for the self-energy.

²The 15-j symbol in (10) depends also on the four spins which label the gauge fixed edge, even if in the analytic notation it is not emphasized.

With the form of the vertex described above, we can write the self-energy amplitude explicitly, in a form suitable for the numerical analysis. This gives



with q = 1...4. Compared to the diagrammatic expressions used for the EPRL vertex (12), we have rearranged some links to emphasize the structure of the contractions between the vertices, which defines the spinfoam associated with the triangulation described in Sec. II.

The EPRL self-energy amplitude (13) is the main object of the analysis performed in this paper. The interested reader can derive the complete analytical expressions by comparing the graphical counterparts with the definitions given in Appendix C.

When the two boundary intertwiners i_+ and i_- have the same value *i*, the EPRL self-energy amplitude simplifies as

$$W(j, i, \gamma; K, \Delta l) = \sum_{\substack{j_{ab} \leq K\\i_a}} d_{\{j_{ab}\}} d_{\{i_a\}} V(j, j_{ab}; i, i_a; \gamma; \Delta l)^2$$
(14)

since the two vertex amplitudes entering the sum over the internal faces j_{ab} are identical. Except for the coherent state (Appendix A 1) and the boundary observable analysis (Appendix A 2), we shall always consider the two boundary intertwinwers i_{\pm} to have the same value. In the following, we use i = (0, 0). The results with different values are substantially identical, differing only by a slightly overall shift of the points, as we show in Appendix A 1.

Numerically, the vertex amplitude $V(j, j_{ab}; i, i_a; \gamma; \Delta l)$ in (14) can be efficiently computed using the recently developed library sl2cfoam-next [34] given the list of boundary spins $(j_i = j, j_{ab})$ and the number of shells Δl . The squared sum over the internal intertwiner indices i_a can be implemented by contracting the resulting *vertex tensor* with itself. All the configurations of internal spins are automatically distributed by the library across the

available cluster nodes, and each node parallelizes the computation of the various shells over the available local CPUs. The particular code employed for this work is available at the repository [45].

From (12), it is clear that the computation of a single EPRL vertex amplitude is much more complex than the corresponding BF vertex amplitude with the same boundary spins and intertwiner indices. In fact, a single EPRL amplitude requires the computation of roughly $(\Delta l + 1)^6$ different BF amplitudes and $(\Delta l + 1)^3$ booster functions, for each different set of intertwiners i_a that bound the vertex [whose number is $\sim (2j+1)^5$ if the boundary spins are of order $\sim j$]. In the present case, a different EPRL vertex tensor (i.e., a bundle of amplitudes at fixed boundary spins with running boundary intertwiner indices) must be computed for each configuration of the internal spins j_{ab} , whose number as a function of the cutoff K has been estimated in the previous section (see Fig. 4). From these simple estimates, we see that the numerical complexity of computing (14), using our methods, grows quickly when any of the parameters K, Δl , or j increases, roughly with a power law between five to six powers of the increasing parameter. We provide in the next section some data for the resources employed in a computation of (14) up to K = 10and $\Delta l = 20$.

V. DIVERGENCE ANALYSIS

We have studied the amplitude (14) numerically and considered the following questions:

- (1) What is the asymptotic scaling of the EPRL divergence?
- (2) What is the dependence on the Barbero-Immirzi parameter?
- (3) At fixed boundary spins and Barbero-Immirzi parameter, does the scaling change using boundary spinnetwork states or intrinsic coherent states?
- (4) What are the configurations which contribute most to the divergence itself?

The first two questions are addressed in the main text, the third is addressed in Appendix A 1, and the fourth is addressed in Appendix B. The main difficulty toward answering these questions is the limit $\Delta l \rightarrow \infty$ that cannot be taken numerically. To address this problem, we have relied on a property of the dependence of the amplitude on Δl at fixed *K*, that has emerged from the numerical analysis itself.

A. Extrapolation scheme

We show in Fig. 5 the amplitude (14) as a function of the cutoff *K*, for various values of the number of shells Δl and for $\gamma = 0.1$ and $\gamma = 1.0$. From a simple qualitative analysis, we infer that the convergence of the shell approximation is faster for a reduced bulk spins cutoff, while it becomes slower as *K* increases. The highest curve is a lower bound



FIG. 5. The EPRL divergence (14) as a function of the cutoff parameter *K*, for a number of shells $\Delta l = 0$ to 10. Darker colors correspond to larger Δl . A portion of the plot around K = 5 is highlighted to show the convergence properties of the shells approximation. *Left:* $\gamma = 0.1$. *Right:* $\gamma = 1.0$.

to the amplitude as a function of *K*, which is recovered in the limit $\Delta l \rightarrow \infty$. Convergence in the number of shells appears to depend on γ , with the case $\gamma = 1.0$ converging faster.

The convergence of (14) in the parameter Δl can be extrapolated from the data reported in Fig. 5. To this aim, we have computed the ratios of the differences between adjacent curves of Fig. 5, at fixed *K*, as a function of Δl . That is, we have studied the function

$$f(\gamma, K, \Delta l) \equiv \frac{W(\gamma, K, \Delta l+2) - W(\gamma, K, \Delta l+1)}{W(\gamma, K, \Delta l+1) - W(\gamma, K, \Delta l)}, \quad (15)$$

for each fixed value of *j* and *i*. The result is shown in Fig. 6 for $K \ge 5$ and $\gamma = 1.0$. It appears from Fig. 6 that the function $f(\gamma, K, \Delta l)$ tends to a constant for larger enough values of *K*, as a function of Δl (however, see also the more complete analysis for the case $\gamma = 0.1$, that is studied in a following section). Since (15) is the ratio between



FIG. 6. Function (15) for $\gamma = 1.0$. The curves appear to tend to a constant value $c_{K,\gamma}$.

successive terms of a series, this means that the sum over the parameter Δl can be well approximated by a geometric series when *K* is large. Defining

$$c_{K,\gamma} \equiv f(\gamma, K, N-2) \tag{16}$$

where *N* is the largest Δl parameter that has been reached numerically, our assumption implies for $\Delta l \ge N$

$$W(\gamma, K, \Delta l + 1) - W(\gamma, K, \Delta l) \approx (c_{K,\gamma})^{\Delta l}.$$
 (17)

We thus approximate the divergence in the limit $\Delta l \rightarrow \infty$ as

$$W(\gamma, K) \equiv \lim_{\Delta l \to \infty} W(\gamma, K, \Delta l)$$

$$\approx W(\gamma, K, N-1) + \frac{W(\gamma, K, N) - W(\gamma, K, N-1)}{1 - c_{K,\gamma}},$$
(18)

where we used the elementary limit of the geometric series. We note that the approximation improves when (i) more shells can be computed numerically, i.e., the index N increases, and (ii) the ratios (15) are closer to true constants in Δl . We applied this extrapolation scheme to the amplitudes computed numerically for different values of γ . In particular, in the case $\gamma = 0.1$, where the convergence in Δl is the slowest, hence the approximation is the least accurate, we pushed the Δl parameter to very large values in order to study the accuracy of our extrapolation scheme and obtained very precise results for a physically relevant case.

B. Case $\gamma = 0.1$

In Fig. 7, we show the amplitude (14) for the case $\gamma = 0.1$, with the computed data for $\Delta l \leq 20$ and the extrapolated curves from N = 10 and from N = 20.



FIG. 7. The complete analysis for the case $\gamma = 0.1$ till $N = \Delta l_{\text{max}} = 20$. Left: all the curves at various Δl with the extrapolation from $\Delta l = 20$. *Right:* The extrapolations from $\Delta l = 10$ and $\Delta l = 20$ with the corresponding linear fits. The slopes of the fits are 8.1×10^{-6} and 8.6×10^{-6} , respectively. The standard errors of the fits, obtained with Julia's LSQFIT package, are 2.3×10^{-8} and 4.6×10^{-9} , respectively.

From these plots, we can verify that the approximation gives results that are already accurate starting with low values of N, since the difference between the two extrapolations is small, around 5% for K = 10. Plotting the ratios (15) till very large values of Δl , as we did in Fig. 8, we can see that for the case $\gamma = 0.1$ the approximation $f(\gamma, K, \Delta l) \approx$ const is more accurate for larger values of K and less accurate for smaller cutoffs. This is not an issue since convergence in Δl is reached sooner for small cutoffs (since the ratios are smaller, the series converges faster). Also, Fig. 8 shows that the extrapolation is actually a lower bound to the true amplitude, although very close to the exact result.

Importantly, the qualitative behavior of the divergence, either extrapolated from N = 10 or N = 20, is clearly linear. In Fig. 7, we show the fit of the extrapolated curves with straight lines. It is clear that the extrapolated points fall



FIG. 8. Function (15) for $\gamma = 0.1$ plotted till $N = \Delta l_{\text{max}} = 20$. It is evident that the approximation $f \approx \text{const}$ for $\Delta l > N$ provides a lower bound to the exact limit $\Delta l \rightarrow \infty$.

exactly on a straight line for $K \gtrsim 5$ in both cases, with the better approximation at N = 20 being also closer to linear behavior than the one at N = 10. We conclude in this case that the overall divergence is very close to linear in the cutoff parameter *K*.

A few comments are in order. First, our results are in accordance with the rough analytical estimate of Riello [31]. It is important to notice that his main analysis focuses on what Riello calls the *nondegenerate* sector of the divergence and disregards the contributions from *degenerate* configurations when the 4-simplices dual to the two spinfoam vertices v_{\pm} have nonmaximal dimension. He found that this sector contributes with a leading factor of log *K* to the divergence. However, in Appendix C of Ref. [31], the remaining degenerate sector is estimated to contribute with a dominant factor of *K* to the overall divergence, which is what our numerical analysis shows. Our result also clearly shows that the interference effects neglected in Ref. [32] result in a strong suppression of the divergence, of approximately eight powers of *K*.

Second, our numerical analysis does not completely exclude the possibility that the behavior for very large values of K is different from what is inferred in the regime $K \leq 10$. There are, however, multiple arguments supporting our conclusion: (i) Our findings are compatible with past analytical estimates [31]. (ii) From Fig. 7, we see that improving the approximation also improves the matching with the linear fit, which appears perfect for as much as about 12 data points. (iii) We can compare the EPRL divergence with the BF case, where the asymptotic powerlaw divergence is already manifest for spins of order ~ 10 (see Fig. 3) and analytically it is seen to be independent from the details of the chosen boundary data. (iv) Recent numerical investigations [34,46] strongly support the hypothesis that the semiclassical (i.e., asymptotic) regime of theory is reached for relatively small spins if the



FIG. 9. Fit of the amplitude $W(\gamma, K)$ extrapolated via Eq. (18) with N = 10 for decreasing values of the Barbero-Immirzi parameter $\gamma \neq 0.1$ and with N = 20 for $\gamma = 0.1$. The fit is made cutting the values of $W(\gamma, K)$ for $K \leq 4$ and superimposing the resulting curve with all the extrapolated points.

Barbero-Immirzi parameter γ is small (e.g., around 0.1). This appears to be related to the slower frequency of asymptotic oscillations of the amplitude whenever γ is small [46]. Hence, we can assume that we are looking at the asymptotic divergence for $K \gtrsim 5$, in this case. This is also supported by our analysis for larger values of γ , which do not appear to converge to a fixed (γ -independent) asymptotic behavior in the considered range of the cutoff, as we show in the next paragraph.

As a final comment for the $\gamma = 0.1$ case, we report some technical details about the hardware resources that were employed. The computation of the divergence at $K \le 10$ and $\Delta l = 20$ ran on 640 CPUs for ~60 h, for a total of ~40, 000 CPU-h. The whole computation at $\gamma = 0.1$ from $\Delta l = 0$ to 20 took about 125,000 CPU-h. Recall that the computational complexity of the EPRL model scales as $(\Delta l + 1)^6$; thus, the plots at largest number of shells took most of the allocated CPU time. Symmetries in the internal spins j_{ab} peculiar to this case (j = 1/2, $i_{\pm} = 0$) have been implemented in the code to reduce the computational time of about a factor ~6.

C. Case $\gamma \gtrsim 0.1$

The extrapolation method outlined above allowed us to investigate also larger values of the Barbero-Immirzi

parameter γ without replicating the already substantial computational effort required for the case $\gamma = 0.1$. We have studied the scaling of the asymptotic divergence for nine values of γ , ranging between 0.1 and 10. The result is shown in Fig. 9, for which we used N = 10 in Eq. (18). Notice that the extrapolation scheme is more effective in the cases $\gamma > 0.1$ than in the case $\gamma = 0.1$, as the curves of the ratios given by the function (15) become closer to constants as γ increases (compare Fig. 6 with Fig. 8). The data of Fig. 9 are fitted with a function,

$$W(\gamma, K) = a + bK^c, \tag{19}$$

where *a*, *b*, *c* are real coefficients. The best fit for the values of these parameters are shown in Table I. Table I shows that the amplitude is suppressed for large γ . The divergence is increasingly well fitted by an approximately linear scaling as the Barbero-Immirzi parameter decreases, while for $\gamma \gtrsim 1$, the curve is approximated by a quadratic function in the range $K \in [0, 10]$. We interpret this apparent dependence of the scaling on γ as a sign that the asymptotic or semiclassical regime is not reached for $K \le 10$ when $\gamma \gtrsim 1$. A possible educated guess is that the nonlinear behavior for large γ could correspond to the initial nonlinear part of the curve seen also at small γ , namely, that the

TABLE I. TH COCHICICIIIS LAUR.	TABLE	I.	Fit	coefficients	table.
--------------------------------	-------	----	-----	--------------	--------

	$W(\gamma, K) = a + bK^c \ (j = 0.5, i = 0), \ K \in [0, 10]$					
γ	а	b	С			
10	-1.7×10^{-32}	9.2×10^{-33}	2.3			
7.5	-2.2×10^{-29}	8.2×10^{-30}	2.3			
5	-4.3×10^{-25}	1.1×10^{-25}	2.3			
3	-5.7×10^{-20}	1.1×10^{-20}	2.3			
1	-8.6×10^{-11}	2.6×10^{-11}	1.9			
0.75	-1.7×10^{-9}	1.1×10^{-9}	1.8			
0.5	-7.6×10^{-8}	6.7×10^{-8}	1.5			
0.25	-1.5×10^{-6}	2.0×10^{-6}	1.2			
0.1	-8.8×10^{-6}	8.6×10^{-6}	1			

dependence could become linear for any γ at sufficiently high *K*. This intriguing hypothesis could have interesting consequences for a renormalization procedure. We leave the accurate testing of the large- γ sector for future works. As the computational complexity dramatically increases when *K* increases, it is likely that different numerical methods or approximations would be needed in order to test this sector, which is interesting from a theoretical point of view but does not appear interesting physically, given the current status of the theory.

VI. CONCLUSIONS

In this work, we applied new computational techniques to the study of the infrared divergence of the spinfoam selfenergy graph in the EPRL model. We have computed the divergent amplitudes with a running cutoff under an approximation (the "shells approximation"), which we then lifted by introducing an extrapolation scheme suited for the case considered. We tested our assumptions by refining the approximation up to a large number of shells, using considerable computational resources, for the most interesting case with the Barbero-Immirzi parameter $\gamma = 0.1$. We also explored the divergence at other values of γ and investigated the dependence and reliability of our results by using different boundary data and by computing geometrical boundary observables.

Our findings considerably refine the upper and lower bounds given in the previous literature [31,32]. We have shown that the numerical evidence clearly points to a *linear* divergence in the infrared cutoff parameter. On one hand, our result shows that the divergence scaling is much less severe than expected by approximate power-counting arguments of BF and Euclidean theories [29,30] or upper-bound estimates to the Lorentzian theory [32]. On the other hand, our result also shows that the total divergence is of higher order than the analytical estimate of the so-called nondegenerate sector [31] and thus provides the first strong evidence that virtual "degenerate" (i.e., of nonmaximal dimension) configurations effectively dominate the first order quantum corrections to the bare spinfoam propagator. This is a point that was hinted at in previous works (see especially Appendix C of Ref. [31]) but left unverified.

As in ordinary quantum field theories, understanding quantum corrections to the bare spinfoam propagator is tightly related to the issue of renormalization, which is still an active area of research in the context of spinfoam and group field theories. Our work fills one of the very first steps toward this program by providing a precise estimate of the degree of divergence. Our results about the divergence at various values of γ also suggest that a possible renormalization scheme might involve the running of the Barbero-Immirzi parameter. The numerical testing of this hypothesis is left for future work, as the possible generalization of our extrapolation scheme to different divergent graphs.

This is one of the first works in the field of covariant LQG to specifically exploit high-performance codes on a computer cluster. The computations accomplished in this paper would not have been possible using a single or a few machines. We believe that complex computational projects will prove to be more and more fundamental to the advancement of the field in the coming years, and we hope that our work and our codes will provide a useful and encouraging ground for progressing in this direction. In this regard, the application of Markov chain Monte Carlo (MCMC) methods in covariant LQG to study the quantum regime in spinfoams composed of multiple vertex amplitudes glued together will appear soon [47].

ACKNOWLEDGMENTS

We thank Pietro Donà for numerous contributions, especially during the early stages of this work. We are grateful to Carlo Rovelli for discussing at length this manuscript and providing many suggestions that improved it. The Centre de Calcul Intensif d'Aix-Marseille and the Shared Hierarchical Academic Research Computing Network (SHARCNET) are acknowledged for granting access to their high-performance computing resources. We thank in particular the Compute/Calcul Canada staff for the constant support provided with the Cedar and Graham clusters. This work was supported by the Natural Science and Engineering Council of Canada (NSERC) through the Discovery Grant "Loop Quantum Gravity: from Computation to Phenomenology." We acknowledge support also from the QISS JFT Grant No. 61466. F. V.'s research is supported by the Canada Research Chairs Program.

APPENDIX A: CONSISTENCY CHECKS

1. Divergence with coherent boundary states

In this subsection, we test whether the results presented in Sec. V depend on the chosen boundary data. The basis chosen diagonalizes one of the two dihedral angles of the boundary tetrahedra, thus leaving the other maximally



FIG. 10. Self-energy amplitude (13) with fixed intertwiners (first two panels) vs coherent (third panel) boundary data (A1). Using coherent intertwiners, we find no difference in scaling by changing the orientation of the normals \vec{n}_a .

spread. Would the results change using a more semiclassical boundary state? To answer this question, we have repeated the computation using intrinsic coherent intertwiners [16] as boundary states.

The definition of the intrinsic semiclassical intertwiners is based on the relation between polyhedra and SU(2)invariants [48,49] and coherent intertwiners [8]. A 4-valent coherent intertwiner is defined by the group averaging,

$$||j_a, \vec{n}_a\rangle \equiv \int dg \bigotimes_{a=1}^4 g |j_a, \vec{n}_a\rangle = \sum_i c_i(\vec{n}_a) |j_a, i\rangle, \quad (A1)$$

where $|j_a, \vec{n}_a\rangle$ are SU(2) coherent states, $g \in SU(2)$. The vectors \vec{n}_a can be parametrized as $\vec{n} \equiv (\sin \Theta \cos \Phi, \sin \Theta \sin \Phi, \cos \Theta)$. The decomposition of a 4-valent coherent intertwiner, in the recoupling channel *i* with outgoing links, reads

$$c_{i}(\vec{n}_{a}) \equiv \langle j_{a}, i | j_{a}, \vec{n}_{a} \rangle$$

$$= \sum_{m_{a}} \begin{pmatrix} j_{1} & j_{2} & j_{3} & j_{4} \\ m_{1} & m_{2} & m_{3} & m_{4} \end{pmatrix}^{(i)} \prod_{a=1}^{4} D_{m_{a}, j_{a}}^{j_{a}}(\vec{n}_{a})$$

$$= \underbrace{\stackrel{\vec{n}_{1}}{j_{1}} \underbrace{\stackrel{\vec{n}_{2}}{j_{2}} & \vec{n}_{3}}{j_{4}}, \qquad (A2)$$

where $D_{m,j}^{j}(\vec{n}) = D_{m,j}^{j}(\phi, \Theta, -\phi)$ is the *SU*(2) Wigner matrix parametrized as

$$D^{j}_{m,j}(\phi,\Theta,-\phi) = e^{-im\phi}d^{j}_{m,j}(\Theta)e^{ij\phi},$$

with $d_{m,j}^{j}$ as a small Wigner matrix [50]. In order to be interpreted as quantum polyhedra, the boundary tetrahedra of the spinfoam must satisfy the closure condition

$$\sum_{f} \Lambda_f \vec{n}_f = 0, \tag{A3}$$

where Λ_f is the eigenvalue of the area operator associated with each face of the tetrahedron.

An example of the numerical analysis of the amplitude obtained by contracting the amplitude (13) with coherent states (A1), for increasing boundary spin values, is shown³ in Fig. 10, compared with boundary spinnetwork states. The scaling of the self-energy amplitude turns out to be unchanged whether using boundary spinnetwork states $|j, i_{\pm}\rangle$ or coherent states $||j, \vec{n}_{\pm f}\rangle$, apart for a global factor that multiplies all curves. Also, using coherent states, the scaling does not seem to depend on the orientation of the normals $\vec{n}_{\pm f}$.

2. Boundary observables

Another check of consistency of the geometrical picture and the numerical calculation can be obtained by looking at the geometry of the two boundary tetrahedra. Their geometrical properties are jointly fixed by choosing the boundary data and by the dynamics. Here, we show the resulting expectation values of some geometrical boundary operators, that is,

$$\langle O \rangle = \frac{\langle W | O | \Psi \rangle}{\langle W | \Psi \rangle}$$
 (A4)

where the bra W contains the propagator, namely, the dynamics, while the ket Ψ turns out to be the tensor product of the *in* and *out* states of the LQG Hilbert space. With the term "propagator," we refer to the square matrices (they are such since the self-energy triangulation has two boundary tetrahedra) in which the element *a*, *b* corresponds to the amplitude (13) with $a = i_+, b = i_-$. Computing the expectation values (A4) of geometric operators describing boundary tetrahedra is a significant numerical investigation, especially considering that computations carried out with the EPRL model are still in their primordial stages [44,51–55], and, as far as we know, at present time, there are no

³Normals \vec{n}_a satisfy condition (A3), and we considered the absolute value in order to obtain real numbers.



FIG. 11. Dynamic expectation value of the cosine of the external dihedral angle operator (A8). We find excellent agreement with the value $-0.\overline{3}$.

such numerical computations with spinfoams containing a bubble.

Since in the calculation of (A4) the dimensional factor of the boundary intertwiners is relevant, in this section, it is convenient to define a "normalized" amplitude as

$$W_N(j, i_{\pm}, K) \equiv W(j, i_{\pm}, K) \sqrt{d_{\{i_{\pm}\}}},$$
 (A5)

where $d_{\{i_{\pm}\}} = \prod_{\pm} (2i_{\pm} + 1) = (2i_{+} + 1)(2i_{-} + 1).$

a. Angles

The shape of the tetrahedra in twisted geometry is measured by the angle operator

$$A_{ab}|j,i_{\pm}\rangle = \cos(\theta_{ab})_{i_{\pm}}|j,i_{\pm}\rangle, \qquad (A6)$$

which is interpreted as the cosine of the external dihedral angle between the faces a and b of the tetrahedron defined on the node associated with the intertwiner i_{\pm} . The spinnetwork basis states diagonalize the dihedral angle θ_{ab} between faces a and b. The equation for the dihedral angle $\cos(\theta_{ab})$ in terms of intertwiner spin i_{\pm} reads

$$\cos(\theta_{ab})_{i_{\pm}} = \frac{i_{\pm}(i_{\pm}+1) - j_a(j_a+1) - j_b(j_b+1)}{2\sqrt{j_a(j_a+1)j_b(j_b+1)}}.$$
 (A7)

We consider the expectation value (A4) of the angle operator (A6) in any of the two (equal) boundary regular tetrahedra of the triangulation using the spinnetwork state. According to the recoupling basis (j_1, j_2) , we focus on the angle between faces 1 and 2. The expectation value on node + can be computed as

$$\frac{\langle W|A_{12}|W\rangle}{\langle W|W\rangle} = \frac{\sum_{i_{\pm}} [W_N(j, i_{\pm}, K)]^2 \cos(\theta_{12})_{i_{\pm}}}{\sum_{i_{\pm}} [W_N(j, i_{\pm}, K)]^2}, \quad (A8)$$



PHYS. REV. D 105, 106018 (2022)

FIG. 12. Dynamic expectation value of the volume operator (A10). We neglected all the constant factors in (A10) and (A11). The volumes scale as $j^{\frac{3}{2}}$.

where we considered the case in which all boundary spins are equal to *j*. Carrying out the numerical computation of (A8), we obtain a value that is in agreement with the geometric value of the external angle of a regular tetrahedron up to the tenth significant digit, as shown in Fig. 11. We repeated the calculation for increasing values of boundary spins *j* up to j = 5, also varying Δl , finding exactly the same value and the unaltered precision. This is consistent with the fact that we are looking at the only angle which is completely sharp, while the others turn out to be spread.

b. Volumes

The volume operator of a tetrahedron reads

$$V_{abc} = \frac{\sqrt{2}}{3} (8\pi G \hbar \gamma)^{\frac{3}{2}} \sqrt{|\vec{J}_a \cdot \vec{J}_b \times \vec{J}_c|}, \qquad (A9)$$

where J_a is the vector of SU(2) generators on link *a*. The operator (A9) is not diagonal in the spinnetwork states. The dynamic expectation value (see Ref. [16] for the derivation) on node + turns out to be

$$\frac{\langle W|V|W\rangle}{\langle W|W\rangle} = \frac{\sum_{i_{\pm},i'_{+}} \sqrt{d_{\{i_{+}\}}d_{\{i'_{+}\}}} W_{N}(j,i_{\pm},K)W_{N}(j,i'_{+},K)v(i_{+},i'_{+},j)}{\sum_{i_{\pm}} [W_{N}(j,i_{\pm},K)]^{2}} \tag{A10}$$

where $v(i_+, i'_+, j)$ is the matrix element

$$v(i_{+},i_{+}',j) = \frac{\sqrt{2}}{3} (8\pi G\hbar\gamma)^{\frac{3}{2}} \sum_{v} \langle j,i_{+}'|v\rangle \sqrt{|q_{v}|} \langle v|j,i_{+}\rangle,$$
(A11)

in which q_v are the eigenvalues of the volume, while $|v\rangle$ corresponds to the eigenvector relative to the eigenvalue q_v .



FIG. 13. The largest individual contributions to the divergent amplitude (14) come from "semiclassical" bulk spins configurations with $\bar{j} \sim j$ and low σ . Triangular inequalities result in a triangular shape in the plane (\bar{j}, σ).

In Fig. 12, we report the numerical values obtained from (A10). We see that the scaling of the expectation value of the boundary volume as a function of the spin *j* corresponds to that actually existing between the volume of a regular tetrahedron and the area of one of its faces, that is $V \propto j^{3/2}$. As for the angle operator, we repeated the calculation by varying the number of shells, finding the same trend.

Finally, for both the angle operator and for the volume, we repeated the calculation with boundary coherent states (A1) verifying that the expectation value is dominated by the classical contribution.

The picture that emerges is that the boundary observables have an expectation value that coincides with the classical one and are not affected by the presence of divergence in the bulk of the spinfoam. Namely, even if the amplitude (13) is divergent, the expectation values (A8), (A10) are finite and fully consistent with the geometry of a regular tetrahedron.

APPENDIX B: INTERNAL CONTRIBUTIONS

We add here a tentative analysis of the kind of configurations that contribute to the divergence. For this aim, it is convenient to rewrite Eq. (14) as

$$W(j, i, \gamma; K, \Delta l) = \sum_{j_{ab} \le K} \omega(j_{ab}; j, i, \gamma, \Delta l).$$
(B1)

Now, consider the average and the standard deviation of the sextuplet of internal spins $\{j_{ab}\}$,

$$\bar{j}(j_{ab}) = \frac{1}{6} \sum_{a \neq b} j_{ab}, \qquad \sigma(j_{ab}) = \sqrt{\frac{1}{6} \sum_{a \neq b} (j_{ab} - \bar{j}_{ab})^2},$$
(B2)

and define

$$w(\bar{j},\sigma;j,i,\gamma,K,\Delta l) = \sum_{\bar{j}(j_{ab})=\bar{j} \atop \sigma(j_{ab})=\sigma} \omega(j_{ab};j,i,\gamma,\Delta l).$$
(B3)

That is, we write the contributions to the divergence as a function of the average and standard deviation of the internal spins, adding the contributions with different sextuplets that have the same average and standard deviation. The function $100 - |\log w(\bar{j}, \sigma; j, i, \gamma, K, \Delta l)|$ is plotted for fixed $j, i, \gamma, \Delta l$ and various values of K in Fig. 13, where the factor 100 is only for visual purposes. Increasing Δl has the effect of increasing the value of each point, while increasing K also implies adding more configurations, as shown in Fig. 4. Figure 13 indicates that the largest individual contributions to the amplitude (14) come from semiclassical bulk spins configurations with $\overline{j} \sim j$ and low σ , as one might expect, while the divergence comes from the proliferation of smaller contribution with arbitrary high spin and deviations.

We also note that the divergence of the self-energy, manifestly shown in Fig. 9, occurs in the limit $\Delta l \rightarrow \infty$ only. In fact, in Ref. [32], it is shown that the sum over the internal spins of the amplitude (14) is convergent in the simplified model $\Delta l = 0$. Even if an analytical proof is not yet available, it seems likely that the amplitude has a finite limit $K \rightarrow \infty$ for each finite Δl , since the internal spins configurations in the limit become more and more suppressed. Since also the limit $\Delta l \rightarrow \infty$ at fixed *K* is finite, we can schematize the role of the two limits with the following four-node diagram:

APPENDIX C: CONVENTIONS AND BASIC FORMULAS

We collect here the basic definitions needed for the calculation in this paper.

1. SU(2) Wigner j-symbols

We use the definition of the Wigner's (3jm) symbol given in Ref. [50] with the orthogonality property

$$\sum_{m_1,\dots,m_2} \begin{pmatrix} j_1 & j_2 & j_3 \\ m_1 & m_2 & m_3 \end{pmatrix} \begin{pmatrix} j_1 & j_2 & j_3 \\ m_1 & m_2 & n_3 \end{pmatrix} = \frac{\delta_{j_3 l_3} \delta_{m_3 n_3}}{2j_3 + 1}.$$
(C1)

These are normalized and vanish if triangular inequalities are not satisfied. We define the (4jm) symbol as the contraction of two (3jm) symbol using an intertwiner k,

$$i_{m_1m_2m_3m_4}^{(k)} \equiv \begin{pmatrix} j_1 & j_2 & j_3 & j_4 \\ m_1 & m_2 & m_3 & m_4 \end{pmatrix}^{(k)} \\ \equiv \sum_{m_i} (-1)^{k-m_i} \begin{pmatrix} j_1 & j_2 & k \\ m_1 & m_2 & m_i \end{pmatrix} \begin{pmatrix} k & j_3 & j_4 \\ -m_i & m_3 & m_4 \end{pmatrix},$$
(C2)

with the orthogonality relations

$$\sum_{m_1,m_2,m_3} \begin{pmatrix} j_1 & j_2 & j_3 & j_4 \\ m_1 & m_2 & m_3 & m_4 \end{pmatrix}^{(k_1)} \begin{pmatrix} j_1 & j_2 & j_3 & j_4 \\ m_1 & m_2 & m_3 & m_4 \end{pmatrix}^{(k_2)} = \frac{\delta_{k_1k_2}}{2k_1+1} \frac{\delta_{j_4l_4}\delta_{m_4n_4}}{2j_4+1},$$
(C3)

where $\frac{\delta_{k_1k_2}}{2k_1+1}$ is the normalization factor. A synthetic notation for (4jm) symbols is also

$$\binom{j_f}{m_f}^{(k)} \equiv \binom{j_1 & j_2 & j_3 & j_4}{m_1 & m_2 & m_3 & m_4}^{(k)}.$$
 (C4)

With these definitions and the conventional definitions of Wigner's matrices for SU(2), $D_{mn}^{(j)}(g)$, we have

$$\int dg D_{m_1 n_1}^{j_1}(g) D_{m_2 n_2}^{j_2}(g) D_{m_3 n_3}^{j_3}(g) D_{m_4 n_4}^{j_4}(g) = \sum_k d_k \begin{pmatrix} j_1 & j_2 & j_3 & j_4 \\ m_1 & m_2 & m_3 & m_4 \end{pmatrix}^{(k)} \begin{pmatrix} j_1 & j_2 & j_3 & j_4 \\ n_1 & n_2 & n_3 & n_4 \end{pmatrix}^{(k)}.$$
(C5)

We denote the 4-valent intertwiner with the ket $|k\rangle$, defined as

$$|k\rangle \equiv i_{m_1m_2m_3m_4}^{(k)}|j_1,m_1\rangle|j_2,m_2\rangle|j_3,m_3\rangle|j_4,m_4\rangle \quad (C6)$$

where the (4jm) symbol is defined in (C2). The states (C6) constitute a basis in the intertwiner space of a 4-valent node,

$$\mathcal{I}_4 \equiv Inv[V^{j_1} \otimes V^{j_2} \otimes V^{j_3} \otimes V^{j_4}], \tag{C7}$$

where V^{j_i} is the irreducible representation of spin j_i . The resolution of the identity in the intertwiner space \mathcal{I}_4 and the orthogonality condition can be written as

$$\mathbf{1}^{(\mathcal{I}_4)} = \sum_{k=0}^{2j} (2k+1) |k\rangle \langle k|,$$
(C8)

$$\langle k|k'\rangle = \frac{\delta_{k,k'}}{(2k+1)}.$$
(C9)

The 15-j symbol we use in this work is the irreducible symbol of first type, following the convention of Ref. [56]. Its graphical representation and its definition in terms of Wigner's 6-j symbols is

$$= (-1)^{\sum_{i=1}^{5} j_i + l_i + k_i} \sum_{s} d_s \left\{ \begin{array}{cc} j_1 & k_1 & s \\ k_2 & j_2 & l_1 \end{array} \right\} \left\{ \begin{array}{cc} j_2 & k_2 & s \\ k_3 & j_3 & l_2 \end{array} \right\} \left\{ \begin{array}{cc} j_3 & k_3 & s \\ k_4 & j_4 & l_3 \end{array} \right\} \left\{ \begin{array}{cc} j_4 & k_4 & s \\ k_5 & j_5 & l_4 \end{array} \right\} \left\{ \begin{array}{cc} j_5 & k_5 & s \\ j_1 & k_1 & l_5 \end{array} \right\}$$

2. Booster functions

The booster functions [43,44], also known as B4 functions [42], are the noncompact remnants of the $SL(2, \mathbb{C})$ integrals. These functions turn out to encode all the details of the EPRL model, such as the *Y*-map. We define them as

$$B_4(j_f, l_f; i, k) \equiv \frac{1}{4\pi} \sum_{p_f} \left(\frac{j_f}{p_f} \right)^{(i)} \left(\int_0^\infty \mathrm{d}r \sinh^2 r \prod_{f=1}^4 d_{j_f l_f p_f}^{(\gamma j_f, j_f)}(r) \right) \left(\frac{l_f}{p_f} \right)^{(k)} = \underbrace{\frac{j_2 - i - j_3}{j_4 - j_4}}_{l_4 - j_4 - j$$

where $d^{(\rho,k)}(r)$ are the boost matrix elements for γ -simple irreducible representations of $SL(2, \mathbb{C})$ in the principal series, γ is the Barbero-Immirzi parameter, and the (4jm) symbols are defined in Appendix C 1.

The booster function $B_4(j_f, l_f; i, k)$ is interpreted as a quantum tetrahedron being boosted among adjacent frames: the two sets j_f and l_f describe the four areas of the tetrahedron in the two frames connected

by a boost, and the two intertwiners i and k describe the quantum intrinsic shape of the tetrahedron [42]. For a precise interpretation of the booster functions and their semiclassical limit, we refer to Ref. [57]. The explicit form of the boost matrix elements can be found in the literature. For the general form, see Refs. [42,58]. We just report here the case of simple irreducible representations:

$$d_{jlp}^{(\gamma j,j)}(r) = (-1)^{\frac{j-l}{2}} \frac{\Gamma(j+i\gamma j+1)}{|\Gamma(j+i\gamma j+1)|} \frac{\Gamma(l-i\gamma j+1)}{|\Gamma(l-i\gamma j+1)|} \frac{\sqrt{2j+1}\sqrt{2l+1}}{(j+l+1)!} \left[(2j)!(l+j)!(l-j)!\frac{(l+p)!(l-p)!}{(j+p)!(j-p)!} \right]^{1/2} \times e^{-(j-i\gamma j+p+1)r} \sum_{s} \frac{(-1)^{s}e^{-2sr}}{s!(l-j-s)!} {}_{2}F_{1}[l+1-i\gamma j,j+p+1+s,j+l+2,1-e^{-2r}].$$
(C12)

- [1] A. Perez, The spin-foam approach to quantum gravity, Living Rev. Relativity **16**, 3 (2013).
- [2] R. Oeckl, A 'general boundary' formulation for quantum mechanics and quantum gravity, Phys. Lett. B 575, 318 (2003).
- [3] C. Rovelli, D. Colosi, L. Doplicher, W. Fairbairn, L. Modesto, and K. Noui, Background independence in a nutshell, Classical Quantum Gravity 22, 2971 (2005).
- [4] W. J. Fairbairn and C. Meusburger, q-Deformation of Lorentzian spin foam models, Proc. Sci. QGQGS2011 (2011) 017.
- [5] M. Han, Cosmological constant in LQG vertex amplitude, Phys. Rev. D 84, 064010 (2011).
- [6] H. M. Haggard, M. Han, W. Kamiński, and A. Riello, Fourdimensional quantum gravity with a cosmological constant from three-dimensional holomorphic blocks, Phys. Lett. B 752, 258 (2016).
- [7] R. Pereira, Lorentzian LQG vertex amplitude, Classical Quantum Gravity 25, 085013 (2008).

- [8] E. R. Livine and S. Speziale, A new spinfoam vertex for quantum gravity, Phys. Rev. D 76, 084028 (2007).
- [9] E. R. Livine and S. Speziale, Consistently solving the simplicity constraints for spinfoam quantum gravity, Europhys. Lett. 81, 50004 (2008).
- [10] L. Freidel and K. Krasnov, A new spin foam model for 4d gravity, Classical Quantum Gravity 25, 125018 (2008).
- [11] J. Engle, R. Pereira, and C. Rovelli, Flipped spinfoam vertex and loop gravity, Nucl. Phys. B798, 251 (2008).
- [12] J. Engle, R. Pereira, and C. Rovelli, The Loop-Quantum-Gravity Vertex-Amplitude, Phys. Rev. Lett. 99, 161301 (2007).
- [13] J. Engle, E. Livine, R. Pereira, and C. Rovelli, LQG vertex with finite Immirzi parameter, Nucl. Phys. B799, 136 (2008).
- [14] W. Kaminski, M. Kisielowski, and J. Lewandowski, Spinfoams for all loop quantum gravity, Classical Quantum Gravity 27, 095006 (2010).

- [15] Y. Ding, M. Han, and C. Rovelli, Generalized spinfoams, Phys. Rev. D 83, 124020 (2011).
- [16] C. Rovelli and F. Vidotto, *Covariant Loop Quantum Gravity* (Cambridge University Press, Cambridge, England, 2014).
- [17] C. Rovelli, Zakopane lectures on loop gravity, Proc. Sci. QGQGS2011 (2011) 003.
- [18] P. Donà and S. Speziale, Introductory lectures to loop quantum gravity, in *TVC 79. Gravitation: Théorie et Expérience*, edited by A. Bounames and A. Makhlouf (Hermann, Paris, 2013).
- [19] B. Dittrich, The continuum limit of loop quantum gravity a framework for solving the theory, in *100 Years of General Relativity: Volume 4*, Loop Quantum Gravity: The First 30 Years, edited by A. Bounames and A. Makhlouf (World Scientific, Singapore, 2017).
- [20] B. Dittrich, S. Mizera, and S. Steinhaus, Decorated tensor network renormalization for lattice gauge theories and spin foam models, New J. Phys. 18, 053009 (2016).
- [21] B. Bahr and S. Steinhaus, Investigation of the spinfoam path integral with quantum cuboid intertwiners, Phys. Rev. D 93, 104029 (2016).
- [22] V. Bonzom, R. Gurau, and V. Rivasseau, Random tensor models in the largeNlimit: Uncoloring the colored tensor models, Phys. Rev. D 85, 084037 (2012).
- [23] D. Benedetti and R. Gurau, Phase transition in dually weighted colored tensor models, Nucl. Phys. B855, 420 (2012).
- [24] S. Carrozza, D. Oriti, and V. Rivasseau, Renormalization of a SU(2) tensorial group field theory in three dimensions, Commun. Math. Phys. 330, 581 (2014).
- [25] J. B. Geloun, T. A. Koslowski, D. Oriti, and A. D. Pereira, Functional renormalization group analysis of rank-3 tensorial group field theory: The full quartic invariant truncation, Phys. Rev. D 97, 126018 (2018).
- [26] V. Bonzom and M. Smerlak, Bubble divergences from twisted cohomology, Commun. Math. Phys. 312, 399 (2012).
- [27] V. Bonzom and M. Smerlak, Bubble divergences: Sorting out topology from cell structure, Ann. Henri Poincaré 13, 185 (2012).
- [28] A. Baratin, S. Carrozza, D. Oriti, J. Ryan, and M. Smerlak, Melonic phase transition in group field theory, Lett. Math. Phys. **104**, 1003 (2014).
- [29] C. Perini, C. Rovelli, and S. Speziale, Self-energy and vertex radiative corrections in LQG, Phys. Lett. B 682, 78 (2009).
- [30] T. Krajewski, J. Magnen, V. Rivasseau, A. Tanasa, and P. Vitale, Quantum corrections in the group field theory formulation of the EPRL/FK models, Phys. Rev. D 82, 124069 (2010).
- [31] A. Riello, Self-energy of the Lorentzian EPRL-FK spin foam model of quantum gravity, Phys. Rev. D 88, 24011 (2013).
- [32] P. Donà, Infrared divergences in the EPRL-FK spin foam model, Classical Quantum Gravity 35, 175019 (2018).
- [33] J. W. Barrett, R. J. Dowdall, W. J. Fairbairn, F. Hellmann, and R. Pereira, Lorentzian spin foam amplitudes: Graphical calculus and asymptotics, Classical Quantum Gravity 27, 165009 (2010).

- [34] F. Gozzini, A high-performance code for EPRL spin foam amplitudes, Classical Quantum Gravity 38, 225010 (2021).
- [35] J. Bezanson, A. Edelman, S. Karpinski, and V. B. Shah, Julia: A fresh approach to numerical computing, SIAM Rev. 59, 65 (2017).
- [36] Wolfram company, Mathematica, Ver. 12.2., Wolfram, 2020.
- [37] E. Bianchi and Y. Ding, Lorentzian spinfoam propagator, Phys. Rev. D 86, 104040 (2012).
- [38] E. Bianchi and H. M. Haggard, Discreteness of the Volume of Space from Bohr-Sommerfeld Quantization, Phys. Rev. Lett. **107**, 011301 (2011).
- [39] L. Freidel and S. Speziale, Twisted geometries: A geometric parametrisation of SU(2) phase space, Phys. Rev. D 82, 084040 (2010).
- [40] H. T. Johansson and C. Forssén, Fast and accurate evaluation of Wigner 3j, 6j, and 9j symbols using prime factorisation and multi-word integer arithmetic, SIAM J. Sci. Comput. 38, A376 (2016).
- [41] J. Rasch and A. C. H. Yu, Efficient storage scheme for precalculated Wigner 3j, 6j and Gaunt coefficients, SIAM J. Sci. Comput. 25, 1416 (2004).
- [42] S. Speziale, Boosting Wigner's nj-symbols, J. Math. Phys. (N.Y.) 58, 032501 (2017).
- [43] P. Donà and G. Sarno, Numerical methods for EPRL spin foam transition amplitudes and Lorentzian recoupling theory, Gen. Relativ. Gravit. **50**, 127 (2018).
- [44] P. Donà, M. Fanizza, G. Sarno, and S. Speziale, Numerical study of the Lorentzian Engle-Pereira-Rovelli-Livine spin foam amplitude, Phys. Rev. D 100, 106003 (2019).
- [45] github.com/PietropaoloFrisoni/Self-energy.
- [46] S. K. Asante, B. Dittrich, and H. M. Haggard, Effective Spin Foam Models for Four-Dimensional Quantum Gravity, Phys. Rev. Lett. 125, 231301 (2020).
- [47] F. V. Pietropaolo Frisoni, Francesco Gozzini, Markov chain Monte-Carlo methods for graph refinement in covariant loop quantum gravity (to be published).
- [48] J. W. Barrett, R. J. Dowdall, W. J. Fairbairn, H. Gomes, and F. Hellmann, Asymptotic analysis of the EPRL foursimplex amplitude, J. Math. Phys. (N.Y.) 50, 112504 (2009).
- [49] P. Donà, M. Fanizza, G. Sarno, and S. Speziale, SU(2) graph invariants, Regge actions and polytopes, Classical Quantum Gravity 35, 045011 (2018).
- [50] V. D. Aleksandroviic, A. N. Moskalev, and K. V. Kel'manoviic, *Quantum Theory of Angular Momentum: Irreducible Tensors, Spherical Harmonics, Vector Coupling Coefficients, 3nj Symbols* (World Scientific, Singapore, 1988).
- [51] P. Donà, M. Fanizza, G. Sarno, and S. Speziale, Numerical study of the Lorentzian Engle-Pereira-Rovelli-Livine spin foam amplitude, Phys. Rev. D 100, 106003 (2019).
- [52] F. Gozzini and F. Vidotto, Primordial fluctuations from quantum gravity, Front. Astron. Space Sci. 7, 629466 (2021).
- [53] M. Han, Z. Huang, H. Liu, D. Qu, and Y. Wan, Spinfoam on a Lefschetz thimble: Markov chain Monte Carlo

computation of a Lorentzian spinfoam propagator, Phys. Rev. D **103**, 084026 (2021).

- [54] M. Han, Z. Huang, H. Liu, and D. Qu, Complex critical points and curved geometries in four-dimensional Lorentzian spinfoam quantum gravity, arXiv:2110.10670.
- [55] P. Frisoni, Studying the EPRL spinfoam self-energy, in *Proceedings of The Sixteenth Marcel Grossmann Meeting* (World Scientific, 2021).
- [56] A. P. Yutsin, I. B. Levinson, and V. V. Vanagas, *Mathemati*cal Apparatus of the Theory of Angular Momentum (Israel Program for Scientific Translation, Jerusalem, Israel, 1962).
- [57] P. Donà and S. Speziale, Asymptotics of lowest unitary SL(2,C) invariants on graphs, Phys. Rev. D 102, 086016 (2020).
- [58] W. Ruhl, *The Lorentz Group and Harmonic Analysis* (Benjamin, New York, 1970).