Hyperasymptotic approximation to the top, bottom, and charm pole mass

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(Received 5 September 2019; accepted 13 January 2020; published 5 February 2020)

We construct hyperasymptotic expansions for the heavy quark pole mass regulated using the principal value (PV) prescription. We apply such hyperasymptotic expansions to the B/D meson masses, and $\bar{\Lambda}$ computed in the lattice. The issue of the uncertainty of the (top) pole mass is critically reexamined. The present theoretical uncertainty in the relation between \bar{m}_t , the $\overline{\text{MS}}$ top mass, and $m_{t,\text{PV}}$, the top pole mass regulated using the PV prescription, is numerically assessed to be $\delta m_{t,\text{PV}} = 28$ MeV for $\bar{m}_t = 163$ GeV.

DOI: 10.1103/PhysRevD.101.034002

I. INTRODUCTION

We construct hyperasymptotic expansions for the heavy quark pole mass regulated using the principal value (PV) prescription along the lines of [1]. We generalize the discussion of that reference by including possible ultraviolet renormalons. We then apply such expansions to various observables. The name hyperasymptotic we borrow from [2]. For a treatise of hyperasymptotic expansions in the context of ordinary differential equations see [3].

In [1] we studied observables characterized by having a large scale $Q \gg \Lambda_{\text{QCD}}$, and for which the operator product expansion (OPE) is believed to be a good approximation. We computed them within an hyperasymptotic expansion. More specifically, the perturbative part of the OPE was summed up using the PV prescription: S_{PV} . The difference between S_{PV} and the full nonperturbative (NP) result is assumed to exactly scale as the intrinsic NP terms of the OPE. In general terms:

$$\begin{aligned} \text{Observable}\left(\frac{Q}{\Lambda_{\text{QCD}}}\right) \\ &= S_{\text{PV}}(\alpha_X(Q)) + K_X^{(\text{PV})} \alpha_X^{\prime}(Q) \frac{\Lambda_X^d}{Q^d} (1 + \mathcal{O}(\alpha_X(Q))) \\ &+ \mathcal{O}\left(\frac{\Lambda_X^{d'}}{Q^{d'}}\right), \end{aligned} \tag{1}$$

where the last term refers to genuine higher order terms in the OPE (d' > d > 0). Then, since S_{PV} can not be computed exactly, we obtain it approximately along an hyperasymptotic expansion (a combination of (truncated) perturbative sums and of NP corrections). This is possible if enough terms of the perturbative expansion are known, and if the divergent structure of the leading renormalons of the observable is also known. This allows us to have a clear (parametric) control on the error of the computation. Two alternative methods were considered in [1] depending on how the truncation of the leading perturbative sum

$$S_T(\alpha) = \sum_{n=0}^{N} p_n^{(X)} \alpha_X^{n+1}(\mu)$$
 (2)

is made:

(1) N and $\mu \sim Q$ large but finite:

$$N = N_P \equiv |d| \frac{2\pi}{\beta_0 \alpha_X(\mu)} (1 - c\alpha_X(\mu)), \quad (3)$$

(2) $N \to \infty$ and $\mu \to \infty$ in a correlated way. We considered two options:

(A)
$$N+1 = N_S(\alpha) \equiv |d| \frac{2\pi}{\beta_0 \alpha_X(\mu)};$$

(B) $N = N_A(\alpha) \equiv |d| \frac{2\pi}{\beta_0 \alpha_X(\mu)} (1 - c' \alpha_X(Q))$ (4)

where c' > 0 but *c* is arbitrary otherwise. Note that in case (1), *c* can partially simulate changes on the scale or scheme of α_X . *d* is the dimension associated to a given renormalon.

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Note that in this paper d can be positive (infrared renormalons) or negative (ultraviolet renormalons), unlike in [1], where only positive d's were considered. Note also that genuine NP corrections are only associated to positive d's.

We will not study the modifications the inclusion of ultraviolet renormalons produce in case (2). In this paper we are mainly concerned in the scenario where the leading renormalon is of infrared nature and subleading renormalons can be ultraviolet and/or infrared. This is the case of the pole mass. In such scenario the precision we can obtain in case (2) is limited by the approximate knowledge of the leading infrared renormalon and we cannot add further to the discussion given in [1]. Different is the case (1), which we discuss in the next section.

The structure of the paper is as follows. In Sec. II we review the general case. Compared with [1] we include the possible effect of ultraviolet renormalons. In Sec. III we study the pole mass of a heavy quark in the large β_0 approximation. We use it as toy-model to test our methods. We then move to real QCD. In Sec. IV we study the B/D meson mass and lattice evaluations of $\overline{\Lambda}$. Finally, in Sec. V we do a dedicated study of the top mass.

In general we will avoid to make explicit the scheme (X) and scale (μ) dependence unless necessary.

II. GENERAL FORMULAS

A. *N* large and $\mu \sim Q \gg \Lambda_{\text{OCD}}$. Eq. (3). Case (1)

This case was already discussed at length in [1]. We now give the general expression after the inclusion of ultraviolet renormalons (for a more detailed discussion see [4]). It can be written in the following way

$$S_{\rm PV}(Q) = S_P + \sum_{\{|d|\}} S_{|d|} + \sum_{\{d>0\}} \Omega_d + \sum_{\{d<0\}} \Omega_d, \quad (5)$$

where

$$S_P \equiv \sum_{n=0}^{N_P(|d_{\min}|)} p_n \alpha^{n+1}(\mu) \equiv S_{|d|=0},$$
 (6)

and (|d| > 0)

$$S_{|d|} \equiv \sum_{n=N_P(|d|)+1}^{N_P(|d'|)} (p_n - p_n^{(as)}) \alpha^{n+1}(\mu),$$
(7)

where the asymptotic behavior associated to renormalons with dimensions $\leq |d|$ is included in $p_n^{(as)}$, and d' is the dimension of the closest renormalon to the origin in the Borel plane fulfilling that |d'| > |d|. Ω_d is a modification of the definition of terminant given in [5] that is more suitable to our case. Whereas in [5] $(p_N \alpha^N \times \text{the})$ terminant refers to the completion of the superasymptotic approximation to give the complete result, here Ω_d is the completation of the part of the perturbative series associated to the singularity located at $u \equiv \frac{\beta_0 t}{4\pi} = \frac{d}{2}$ in the Borel plane using the PV prescription. For the case of infrared renormalons (d > 0)the general analytic expression of Ω_d can be found in [1]. For a generic ultraviolet renormalon (d < 0) that produces the asymptotic behavior

$$p_n^{(as)} = Z_{O_d}^X \frac{\mu^d}{Q^d} \frac{\Gamma(n+b'+1)}{\Gamma(b'+1)} \left(\frac{\beta_0}{2\pi d}\right)^n \left\{ 1 + c_1 \frac{b'}{n+b'} + c_2 \frac{b'^2}{(n+b')(n+b'-1)} + \dots \right\},$$
(8)

 $\Omega_{d<0}$ reads

$$\Omega_{d<0} = \Delta \Omega_{\rm UV}(db) + c_1 \Delta \Omega_{\rm UV}(db-1) + \cdots, \quad (9)$$

where¹ (we define $\eta_c \equiv -b' + \frac{2\pi |d|c}{\beta_0} - 1$ where $b' = db - \gamma$)

$$\begin{split} \Delta\Omega_{\rm UV}(db) &= Z_{O_d}^{\chi} \frac{\mu^d}{Q^d} (-1)^{N_P+1} \frac{1}{\Gamma(b'+1)} \left(\frac{\beta_0}{2\pi |d|}\right)^{N_P+1} \alpha^{N_P+2} \int_0^\infty dx \frac{e^{-x} x^{N_P+1+b'}}{1 + \frac{x\beta_0 \alpha}{2\pi |d|}} \\ &= Z_{O_d}^{\chi} \frac{\mu^d}{Q^d} (-1)^{N_P+1} \frac{\pi}{\Gamma(b'+1)} \left(\frac{\beta_0}{|d|}\right)^{-b'-1/2} \alpha(\mu)^{1/2-b'} e^{\frac{-2\pi |d|}{\beta_0 \alpha(\mu)}} \left\{1 + \frac{\alpha(\mu)}{\pi} \frac{\beta_0}{12|d|} \left[-1 + 3\eta_c^2\right] \right. \\ &+ \frac{\alpha^2(\mu)}{\pi^2} \frac{\beta_0^2}{1152|d|^2} \left[13 - 48\eta_c - 60\eta_c^2 + 48\eta_c^3 + 36\eta_c^4\right] + \mathcal{O}(\alpha^3) \right\}. \end{split}$$

Joining all terms together we have

$$\Omega_{d<0} = \sqrt{\alpha(\mu)} K_X^{(P)} \frac{Q^{|d|}}{\mu^{|d|}} e^{\frac{-2\pi|d|}{\beta_0 \alpha(\mu)}} \left(\frac{\beta_0 \alpha(\mu)}{4\pi} \right)^{-b'} \{ 1 + \bar{K}_{X,1}^{(P)} \alpha(\mu) + \bar{K}_{X,2}^{(P)} \alpha^2(\mu) + \mathcal{O}(\alpha^3(\mu)) \},$$
(11)

¹An sketch of how these computations are done is given in the Appendix A.

where

$$K_X^{(P)} \equiv Z_{\mathcal{O}_d}^X(-1)^{N_p+1} \left(\frac{\beta_0}{\pi^2 |d|}\right)^{-1/2} \frac{1}{\Gamma(b'+1)} \left(\frac{2}{|d|}\right)^{-b'},$$
(12)

$$\bar{K}_{X,1}^{(P)} \equiv \left(\frac{2}{\pi}\right)^{1/2} \left(c_1 \frac{\beta_0 b'}{2\sqrt{2\pi}|d|} + \frac{\beta_0}{12|d|\sqrt{2\pi}}(-1+3\eta_c^2)\right),\tag{13}$$

$$\bar{K}_{X,2}^{(P)} \equiv \left(\frac{2}{\pi}\right)^{1/2} \left(c_2 \frac{b'^2 \beta_0^2}{4\sqrt{2}|d|^2 \pi^{3/2}} + c_1 \frac{b' \beta_0^2 (-1 + 3(\eta_c + 1)^2)}{24\sqrt{2}|d|^2 \pi^{3/2}} + \frac{\beta_0^2}{1152|d|^2 2^{1/2} \pi^{3/2}} [13 - 48\eta_c - 60\eta_c^2 + 48\eta_c^3 + 36\eta_c^4] \right).$$
(14)

Note that in this case μ is in the denominator. If we set the anomalous dimension to zero (b' = db), $\Omega_{d<0} \sim \sqrt{\alpha(\mu)} \frac{\Lambda_{\rm QCD}^{|d|}}{\mu^{2|d|}}$ (unlike for infrared renormalons where $\Omega_{d>0} \sim \sqrt{\alpha(\mu)} \frac{\Lambda_{\rm QCD}^{d}}{Q^d}$). If one takes μ very large this term will be quite small. In practice, if we take $\mu \sim Q$, we may need this term.

 $S_{\rm PV}$ will be computed truncating the hyperasymptotic expansion in a systematic way. This means truncating Eq. (5) as follows (note that we always define *D* to be positive):

$$S_{\rm PV}^{(D,N)}(Q) = \sum_{\{|d|\}} S_{|d| < D} + \sum_{\{|d| \le D\}} \Omega_d + \sum_{n=N_P(D)+1}^{N_P(D)+N} (p_n - p_n^{(as)}) \alpha^{n+1}(\mu).$$
(15)

For each value of the couple (D, N), we can state the parametric accuracy of $S_{PV}^{(D,N)}(Q)$. For instance for $S^{(0,N_P)}$ the error would be (up to a numerical and a $\sqrt{\alpha_X}$ factor)

$$\delta S^{(0,N_P)} \sim \mathcal{O}(e^{-|d_{\min}|_{\overline{\beta}_0 a_X(\mathcal{Q})}^{2\pi}}).$$
(16)

This is what is commonly named the superasymptotic approximation. For $S^{(|d_{\min}|,0)}$ the parametric form of the error reads (up to a numerical and a possible $\alpha_X^{3/2}$ factor):

$$\delta S^{(|d_{\min}|,0)} \sim \mathcal{O}(e^{-|d_{\min}|\frac{2\pi}{\beta_0 a_X(Q)}(1+\ln(|d/d_{\min}|))}), \qquad (17)$$

where *d* is the location of the next renormalon closest to the origin. This corresponds to the first term in the hyper-asymptotic approximation. The expression for the error in the general case $S_{PV}^{(D,N)}(Q)$ reads $(N \neq N_P \text{ but large})$

$$\delta S^{(D,N)} \sim \mathcal{O}(e^{-D\frac{2\pi}{\beta_0 \alpha_X(\mathcal{Q})}(1+\ln(|d/D|)}\alpha_X^N),$$
(18)

where d is the location of the next renormalon closest to the origin after D.

B. $m_{\rm PV}$, general formulas

For the case of the heavy quark mass, which we discuss at length in this paper, we have $[\bar{m} = m_{\overline{\text{MS}}}(m_{\overline{\text{MS}}})]$

$$m_{\rm PV}(\bar{m}) = m_P + \bar{m}\Omega_m + \sum_{n=N_P+1}^{2N_P} (r_n - r_n^{(as)})\alpha^{n+1}(\mu) + \bar{m}\Omega_2 + \bar{m}\Omega_{-2} + \mathcal{O}(e^{-2\frac{2\pi}{\beta_0 a}(1+\ln(3/2))}), \quad (19)$$

where

$$m_P \equiv \bar{m} + \sum_{n=0}^{N_P} r_n \alpha^{n+1}(\mu);$$
 (20)

the coefficients r_n for $n \leq 3$ were computed in [6–9];

$$\Omega_{m} = \sqrt{\alpha_{X}(\mu)} K_{X}^{(P)} \frac{\mu}{\bar{m}} e^{-\frac{2\pi}{\bar{\rho}_{0}\alpha_{X}(\mu)}} \left(\frac{\beta_{0}\alpha_{X}(\mu)}{4\pi}\right)^{-b} (1 + \bar{K}_{X,1}^{(P)}\alpha_{X}(\mu) + \bar{K}_{X,2}^{(P)}\alpha_{X}^{2}(\mu) + \mathcal{O}(\alpha_{X}^{3}(\mu))),$$
(21)

where now $K_X^{(P)}$ and $K_{X,i}^{(P)}$ read

$$K_X^{(P)} = -\frac{Z_m^X 2^{1-b} \pi}{\Gamma(1+b)} \beta_0^{-1/2} \left[-\eta_c + \frac{1}{3} \right], \qquad (22)$$

$$\bar{K}_{X,1}^{(P)} = \frac{\beta_0/(\pi)}{-\eta_c + \frac{1}{3}} \left[-b_1 b \left(\frac{1}{2} \eta_c + \frac{1}{3} \right) - \frac{1}{12} \eta_c^3 + \frac{1}{24} \eta_c - \frac{1}{1080} \right],$$
(23)

$$\bar{K}_{X,2}^{(P)} = \frac{\beta_0^2 / \pi^2}{-\eta_c + \frac{1}{3}} \left[-w_2(b-1)b\left(\frac{1}{4}\eta_c + \frac{5}{12}\right) + b_1 b\left(-\frac{1}{24}\eta_c^3 - \frac{1}{8}\eta_c^2 - \frac{5}{48}\eta_c - \frac{23}{1080}\right) - \frac{1}{160}\eta_c^5 - \frac{1}{96}\eta_c^4 + \frac{1}{144}\eta_c^3 + \frac{1}{96}\eta_c^2 - \frac{1}{640}\eta_c - \frac{25}{24192} \right],$$
(24)

where we have applied the general expression obtained in [1] to this case. In particular (*b* and s_n are defined in [1]),

$$\eta_c = -b + \frac{2\pi c}{\beta_0} - 1, \qquad b_1 = s_1, \qquad w_2 = \left(\frac{s_1^2}{2} - s_2\right) \frac{b}{b-1}.$$
(25)

Finally,

$$r_n^{(\rm as)}(\mu) = Z_m^X \mu (\frac{\beta_0}{2\pi})^n \sum_{k=0}^{\infty} c_k \frac{\Gamma(n+1+b-k)}{\Gamma(1+b-k)}.$$
 (26)

The coefficients c_k are pure functions of the β -function coefficients, as first shown in [10]. They can be found in [11–13]. At low orders they read ($c_0 = 1$)

$$c_{1} = s_{1}, \qquad c_{2} = \frac{1}{2} \frac{b}{b-1} (s_{1}^{2} - 2s_{2}),$$

$$c_{3} = \frac{1}{6} \frac{b^{2}}{(b-2)(b-1)} (s_{1}^{3} - 6s_{1}s_{2} + 6s_{3}). \qquad (27)$$

Note that

$$m_{\rm OS}^{(N)} = \bar{m} + \sum_{n=0}^{N} r_n \alpha^{n+1}(\mu).$$
 (28)

Therefore, m_P is nothing but the pole mass truncated to order $N = N_P$.

Our knowledge of the other terminants, Ω_2 and Ω_{-2} , is limited. We do not know the renormalization group structure of Ω_{-2} , except in the large β_0 . On the other hand, the renormalization group structure of Ω_2 is exactly known (provided the coefficients of the beta function are known to all orders). The reason is that it is linked to the kinetic operator of the HQET Lagrangian, the Wilson coefficient of which is protected by reparametrization invariance [14]. Therefore, it has no anomalous dimension and the Wilson coefficient is 1 in dimensional regularization to all orders in perturbation theory. Still, in the large β_0 approximation, the coefficient Z_2^X is equal to zero. If it is different from zero beyond the large β_0 approximations has been a matter of debate [15]. We will retake this discussion in the following sections.

We also give the formulas that apply to Eq. (4), i.e., to case (2): the limit $(N, \mu) \rightarrow \infty$. A general discussion can be found in [1]. It was argued that the limit (2A) was likely to be logarithmic divergent (see also [16]), and no formulas could be found that are valid beyond the large β_0 approximation. Therefore, we will not study this case further. For the limit (2B) formulas with NP exponential accuracy were found in [1] generalizing results from [17]. These formulas were valid beyond the large β_0 approximation. For the specific case of the pole mass they read

$$m_{\rm PV} = m_A + K_X^{(A)} \Lambda_X + \mathcal{O}(\alpha \Lambda_X), \qquad (29)$$

where

$$m_A = \bar{m} + \lim_{\mu \to \infty; 2B} \sum_{n=0}^{N_A} r_n \alpha^{n+1}(\mu),$$
 (30)

and

$$K_X^{(A)} = \frac{2\pi}{\beta_0} Z_m^X \left(\frac{\beta_0}{4\pi}\right)^b \int_{-c',\mathrm{PV}}^{\infty} dx e^{-\frac{2\pi}{\beta_0}x} \frac{1}{(-x)^{1+b}}.$$
 (31)

As discussed in [1], there is more than one way to take the $\mu \to \infty$; (2B) limit. One is to take Eq. (67) of [1] for N_A instead of limit (B) of Eq. (4). Both methods are general but require the knowledge of r_n and the beta function coefficients to all orders. This potentially limits their applicability in practice. Another option is to interpret the $\mu \to \infty$ limit as a change of scheme (where $\mu_0 \sim \bar{m}$):

$$\alpha_{X'}(\mu) = \frac{\alpha_X(\mu_0)}{1 + \frac{\beta_0}{2\pi} \alpha_X(\mu_0) \ln(\frac{\mu}{\mu_0})} \quad \text{and}$$
$$N_A(\alpha) \equiv \frac{2\pi}{\beta_0 \alpha_{X'}(\mu)} (1 - c' \alpha_X(\mu_0)). \quad (32)$$

This method still requires the knowledge of r_n to all orders. On the other hand, there is no need to know the β -function to all orders.

Irrespectively of which of the above methods we use to take the $\mu \rightarrow \infty$ limit we have

$$\lim_{\mu \to \infty; 2B)} \sum_{n=0}^{N_A} r_n \alpha^{n+1}(\mu) = \int_0^{\frac{4\pi}{\beta_0 \chi}} dt e^{-t/\alpha_X(\mu_0)} B[m_{\rm PV} - \bar{m}](t),$$
(33)

where $2/\chi = 1 - c'\alpha(\mu_0)$. The right-hand side of Eq. (33) can not be computed exactly. An approximated determination can be obtained by approximating the Borel transform to $(u \equiv \frac{\beta_0 t}{4\pi})$

$$B[m_{\rm PV} - \bar{m}](t) = \sum_{n=0}^{N_{\rm max}} \frac{(r_n - r_n^{(as)})}{n!} t^n + \frac{Z_m \mu}{(1 - 2u)^{1+b}} (1 + c_1(1 - 2u) + c_2(1 - 2u)^2 + \cdots),$$
(34)

where N_{max} is the number of perturbative coefficients that are known.

III. $m_{PV}(\bar{m})$ IN THE LARGE β_0 APPROXIMATION

Here the discussion runs parallel to the discussion for the static potential in the large β_0 approximation made in Sec. III of [1]. Nevertheless, we do not have the same analytic control as for the static potential. Note also that now we have ultraviolet renormalons. Moreover, the pole mass has the extra complication that it is ultraviolet divergent and needs renormalization. This makes the Borel transform more complicated and we do not have the exact μ factorization one has in the static potential. We take the Borel transform from [18–20]: $B[m_{\rm PV} - \bar{m}](u)$

$$=\bar{m}\frac{C_{F}}{4\pi}\left[\left(\frac{\bar{m}^{2}}{\mu^{2}}\right)^{-u}e^{-c_{\overline{\rm MS}}u}6(1-u)\frac{\Gamma(u)\Gamma(1-2u)}{\Gamma(3-u)}-\frac{3}{u}+R(u)\right],$$
(35)

where $c_{\overline{\text{MS}}} = -5/3$, $u = \frac{\beta_0}{4\pi}t$ and

$$R(u) = \sum_{n=1}^{\infty} \frac{1}{(n!)^2} \frac{d^n}{dz^n} G(z) \Big|_{z=0} u^{n-1} = -\frac{5}{2} + \frac{35}{24}u + \mathcal{O}(u^2),$$
(36)

$$G(u) = -\frac{1}{3}(3+2u)\frac{\Gamma(4+2u)}{\Gamma(1-u)\Gamma^2(2+u)\Gamma(3+u)}.$$
 (37)

This expression has been derived in the \overline{MS} scheme. Whereas the scheme dependence of the first term in Eq. (35) can be reabsorbed in changes of μ and $c_{\overline{\text{MS}}}$ (it would then be equivalent to a change of scale), controlling the scheme dependence of R(u) is more complicated. We will not care much, as R(u) has to do with the high energy behavior, and should only affect m_P , the finite sum. Therefore, when we change from the \overline{MS} to the lattice scheme, we will just change $c_{\overline{\text{MS}}} \rightarrow c_{\text{latt}}$ and leave R(u)unchanged. Strictly speaking then, the object we compute in the lattice scheme is not the pole mass, still it will have the same infrared behavior. The fact that we will obtain the same result after subtracting m_P from m_{PV} in both cases will be a nice confirmation that high-energy cancellation has effectively taken place and what is left is low energy.² The value of c_{latt} that we use is the same to the one used in [1]. To determine it we take the $n_f = 0$ number for a Wilson action of $d_1 = 5.88359$ [25] and use $c_{\text{latt}} = -2(\frac{5}{6} + \frac{2\pi d_1}{\beta_0})$. This is enough for our purposes, as we only use this scheme for checking the consistency between the results obtained with different schemes. Note that this yields two values of c_{latt} if we introduce the n_f dependence of β_0 : $c_{\text{latt}}(n_f = 0) = -8.38807$ and $c_{\text{latt}}(n_f = 3) = -9.88171$.

A. *N* large and $\mu \sim \bar{m} \gg \Lambda_{\text{OCD}}$. Eq. (3). Case (1)

We now take Eq. (19) in the large β_0 approximation and truncate it at different orders in the hyperasymptotic expansion. We then compare such truncations with the exact solution. We can study (even if in the large β_0 approximation) up to which values of \bar{m} the OPE is a good approximation of $m_{\rm PV}$. Remarkably enough we can actually check more than one term of the OPE (hyperasymptotic) expansion. Note that in the large β_0 approximation $\Omega_2 = 0$, but not Ω_{-2} , which in the large β_0 approximation reads $(\eta_c^{(\beta_0)} = \frac{2\pi c}{\beta_0} - 1)$

$$\Omega_{-2} = \sqrt{\alpha(\mu)} K_X^{(P)} \frac{\Lambda_X^2 \bar{m}^2}{\mu^4} \{ 1 + \bar{K}_{X,1}^{(P)} \alpha(\mu) + \bar{K}_{X,2}^{(P)} \alpha^2(\mu) + \mathcal{O}(\alpha^3(\mu)) \},$$
(38)

$$K_X^{(P)} \equiv Z_{-2}^X (-1)^{N_p + 1} \left(\frac{\beta_0}{2\pi^2}\right)^{-1/2}, \quad Z_{-2}^X = -\frac{C_F e^{c_X}}{\pi}, \quad (39)$$

$$\bar{K}_{X,1}^{(P)} \equiv \frac{\beta_0}{24\pi} \left(-1 + 3\eta_c^{(\beta_0)2} \right),\tag{40}$$

$$\bar{K}_{X,2}^{(P)} \equiv \frac{\beta_0^2}{4608\pi^2} [13 - 48\eta_c^{(\beta_0)} - 60\eta_c^{(\beta_0)2} + 48\eta_c^{(\beta_0)3} + 36\eta_c^{(\beta_0)4}].$$
(41)

We also explore the scheme dependence by performing the computation in the lattice and the \overline{MS} scheme. We will do these analyses for the cases with $n_f = 0$ and $n_f = 3$. The first in view of comparing with quenched lattice simulations, the second to simulate a more physical scenario, for which we can draw some conclusions that could be applied beyond the large- β_0 limit. In Figs. 1–3 we plot $m_{\rm PV} - \bar{m}, \quad m_{\rm PV} - m_P, \quad m_{\rm PV} - m_P - \bar{m}\Omega_m, \quad m_{\rm PV} - m_P - \bar{m}\Omega_m,$ $\bar{m}\Omega_m - \sum_{n=N_P+1}^{2N_P} (r_n - r_n^{(as)}) \alpha^{n+1}$, and $m_{\rm PV} - m_P - \bar{m}\Omega_m - m_P - \bar{m}\Omega_m - m_P - \bar{m}\Omega_m$ $\sum_{n=N_p+1}^{2N_p} (r_n - r_n^{(as)}) \alpha^{n+1} - \bar{m} \Omega_{-2} \text{ with } n_f = 0 \text{ light flavors.}$ In the counting of Eq. (15) this corresponds to (0,0), $(0, N_P)$, $(1,0), (1, N_P), (2,0)$ precision. We do such computation in the lattice (Fig. 1) and the \overline{MS} (Fig. 2) scheme. In Fig. 3 we compare the results in the lattice and \overline{MS} scheme. We observe a very nice convergent pattern in all cases down to surprisingly small scales. To visualize the dependence on c, we show the band generated by the smallest positive and negative possible values of c that yield integer values for N_P . The size of the band generated by the different values of c(the c dependence) decreases after introducing Ω_m to its associated sum. On the other hand Ω_{-2} (an ultraviolet renormalon) gives a very small contribution, in particular in the lattice scheme. This is consistent with interpreting the lattice scheme as the \overline{MS} scheme using a much higher renormalization scale μ for the scale of the strong coupling.

Let us discuss the results in more detail. We first observe that the \bar{m} dependence of $m_{\rm PV}$ is basically eliminated in $m_{\rm PV} - m_P$, as expected. This happens both in the lattice and $\overline{\rm MS}$ scheme. The latter shows a stronger *c* dependence. This is to be expected, as in the $\overline{\rm MS}$, we truncate at smaller orders in *N*. This makes the truncation error bigger. As we can see in the upper panel of Fig. 3, both schemes yield consistent predictions for $m_{\rm PV} - m_P$. We can draw some interesting observations out of this analysis. For $m_{\rm PV} - m_P$

²To make an analogy, the situation is similar to determinations of the infrared behavior of the energy of an static source in lattice perturbation theory. In [21-23] two different discretizations were used for the static quark propagators. This affected the ultraviolet, but let the infrared behavior unchanged, as it was nicely seen in those simulations. See also the discussion in [24].



FIG. 1. Upper panel: We plot $m_{\rm PV} - \bar{m}$ (black line) and the differences: (a) $m_{\rm PV} - m_P$ (cyan), (b) $m_{\rm PV} - m_P - \bar{m}\Omega_m$ (orange), (c) $m_{\rm PV} - m_P - \bar{m}\Omega_m - \sum_{n=N_P+1}^{2N_P} (r_n - r_n^{\rm (as)}) \alpha^{n+1}$ (green), and (d) $m_{\rm PV} - m_P - \bar{m}\Omega_m - \sum_{n=N_P+1}^{2N_P} (r_n - r_n^{\rm (as)}) \alpha^{n+1} - \bar{m}\Omega_{-2}$ (blue) in the large β_0 approximation using the lattice scheme with $n_f = 0$ light flavors. For each difference, the bands are generated by the difference of the prediction produced by the smallest positive or negative possible values of c that yields integer values for N_P . The (c) and (d) bands are one on top of the other. Lower panel: As in the upper panel but in a smaller range. $r_0^{-1} \approx 400$ MeV. The value of N_P depends on the scale 1/m we use. For instance for c positive, $N_P = 9$ for $1/m \in [0.003, 0.0045]$, $N_P = 8$ for $1/m \in [0.006, 0.0015]$, $N_P = 7$ for $1/m \in$ $[0.0165, 0.0435], N_P = 6 \text{ for } 1/m \in [0.045, 0.01185], N_P = 5$ for $1/m \in [0.12, 0.321]$, $N_P = 4$ for $1/m \in [0.3225, 0.876]$ and $N_P = 3$ for $1/m \in [0.8775, 1.299]$.

is better to choose a larger factorization scale, if we have enough coefficients of the perturbative expansion. This is particularly so at large distances: We can still get good results at very large distances in the lattice scheme.

We now turn to $m_{PV} - m_P - \bar{m}\Omega_m$. Adding the new correction brings much better agreement with expectations (which we remind is to get zero). After the introduction of $\bar{m}\Omega_m$, the \overline{MS} scheme yields more accurate results than the lattice scheme. This can already be seen in the upper panel of Fig. 3, and in greater detail in the lower panel of Fig. 3.

 $m_{\rm PV} - m_P - \bar{m}\Omega_m$ shows some dependence on \bar{m} , which is more pronounced in the lattice than in the $\overline{\rm MS}$ scheme. As in the large β_0 approximation the difference between both schemes is somewhat equivalent to a change of scale,



FIG. 2. As in Fig. 1 but in the $\overline{\text{MS}}$ scheme. The values of N_P for c positive are, for instance, $N_P = 6$ for 1/m = 0.003, $N_P = 5$ for $1/m \in [0.0045, 0.0105]$, $N_P = 4$ for $1/m \in [0.012, 0.03]$, $N_P = 3$ for $1/m \in [0.0315, 0.0825]$, $N_P = 2$ for $1/m \in [0.084, 0.2235]$, $N_P = 1$ for $1/m \in [0.225, 0.6105]$ and $N_P = 0$ for $1/m \in [0.612, 1.5]$.

these results point to that $\mu = \overline{m}$ in \overline{MS} scheme is close to the natural scale and minimize higher order corrections. Note that the lattice scheme computation is equivalent to the $\overline{\text{MS}}$ scheme choosing $\mu_{\text{latt}} = \mu_{\overline{\text{MS}}} e^{\frac{-c_{\text{latt}}}{2}} e^{\frac{c_{\overline{\text{MS}}}}{2}}$. This gives around a factor 30!. Once $\sum_{n=N_p+1}^{2N_p} (r_n - r_n^{(as)}) \alpha^{n+1}$ is incorporated in the prediction most of the difference between schemes disappears. The effect of introducing Ω_{-2} is very small, in particular in the lattice scheme. This is to be expected, since the lattice scheme corresponds to a larger renormalization scale μ . In any case, the difference between schemes gets smaller and smaller as we go to higher orders in the hyperasymptotic expansion, in particular at short distances. We also want to stress that this analysis opens the window to apply perturbation theory at rather large distances. Note that in the upper panel plots in Figs. 1, 2, and 3, we have gone to very large distances.

As some concluding remarks let us emphasize the following points. $m_{\rm PV} - m_P$ is more or less constant with relatively large uncertainties. This is to be expected, as the next correction in magnitude is $\bar{m}\Omega_m$ which is approximately constant (mildly modulated by $\sqrt{\alpha(\mu)}$). After introducing this term the error is much smaller and we can see more structure. In particular we are sensitive to



FIG. 3. Comparison of lattice and $\overline{\text{MS}}$ scheme results for $n_f = 0$ obtained in Figs. 1 and 2. Upper panel: We plot $m_{\text{PV}} - \bar{m}$ and the differences: (a) $m_{\text{PV}} - m_P$, and (b) $m_{\text{PV}} - m_P - \bar{m}\Omega_m$ in the lattice and $\overline{\text{MS}}$ scheme with $n_f = 0$ light flavors. Lower panel: Lower panel Figs. 1 and 2 combined.

 $\sum_{n=N_P+1}^{2N_P} (r_n - r_n^{(as)}) \alpha^{n+1}$. Here we find (at the level of precision we have now) a sizable difference between lattice and $\overline{\text{MS}}$. This can be expected: $\sum_{n=N_P+1}^{2N_P} (r_n - r_n^{(as)}) \alpha^{n+1}$ is the object we expect to be more sensitive to the scale.

Another interesting observation is that truncated sums behave better in the lattice scheme than in the $\overline{\text{MS}}$ scheme. Nevertheless, this could be misleading. The sums are truncated at the minimal term. Therefore, one needs more terms in the lattice scheme. If the number of terms is not an issue (which could be the case with dedicated numerical stochastic perturbation theory (NSPT) [26,27] computations in the lattice scheme) then the lattice scheme looks better. But as soon as Ω_m is introduced in the computation $\overline{\text{MS}}$ behaves better (at least in the large β_0 approximation).

Overall, we observe a very nice convergence pattern up to (surprisingly) rather large scales in the lattice and $\overline{\text{MS}}$ scheme. The agreement with the theoretical prediction (which is zero) is perfect at short distances. The estimated error is also expected to be small. It will be interesting to see if this also happens beyond the large β_0 .

We now turn to the $n_f = 3$ case. To easy the comparison with [1], we use the same value: $\Lambda_{\overline{\text{MS}}}(n_f = 3) = 174 \text{ MeV}$ (which yields $\alpha(M_{\tau}) \approx 0.3$). The general conclusions do not change if we fix $\Lambda_{\overline{\text{MS}}}$ (in the large β_0 approximation)



FIG. 4. As in Fig. 1 but with $n_f = 3$ light flavors. The value of N_P depends on the scale 1/m we use. For instance for *c* positive, $N_P = 1$ for 1/m = 0.005, $N_P = 10$ for $1/m \in [0.01, 0.015]$, $N_P = 9$ for $1/m \in [0.02, 0.04]$, $N_P = 8$ for $1/m \in [0.045, 0.115]$, $N_P = 7$ for $1/m \in [0.12, 0.315]$, $N_P = 6$ for $1/m \in [0.32, 0.865]$, and $N_P = 5$ for $1/m \in [0.87, 1.5]$.

using the world average value of α . We note that Λ_{OCD} for the physical case $(n_f = 3)$ is smaller than for the $n_f = 0$ case (if one sets the physical scale according to $r_0^{-1} \approx 400$ MeV). On top of that the running is less important. All this points to that the convergence should be even better than in the $n_f = 0$ case (and it was quite good already there). We show our results in Figs. 4-6 (these are the analogous of Figs. 1–3 but with $n_f = 3$). These plots confirm our expectations. Down to scales as low as 667 MeV we see no sign of breakdown of the OPE. This is so in both the lattice and the \overline{MS} schemes. Note that the precision we get is extremely high as we go to small scales: Using truncation (c): $m_P + \bar{m}\Omega_m +$ $\sum_{n=N_{P}+1}^{2N_{P}} (r_{n} - r_{n}^{(as)}) \alpha^{n+1}$, one gets m_{PV} in the \overline{MS} scheme with a precision below 1 MeV at scales of the order of the mass of the bottom, and in the lattice scheme with a precision below 2 MeV. Using truncation (d): $m_P + \bar{m}\Omega_m + \sum_{n=N_P+1}^{2N_P} (r_n - r_n^{(as)}) \alpha^{n+1} + \bar{m}\Omega_{-2}$, the precision does not significantly change, in particular in the lattice scheme. This reflects that ultraviolet renormalons play a minor role. The rest of the discussion follows parallel the one for $n_f = 0$.



FIG. 5. As in Fig. 1 but with $n_f = 3$ light flavors and in the $\overline{\text{MS}}$ scheme. The value of N_P depends on the scale 1/m we use. For instance for *c* positive, $N_P = 7$ for 1/m = 0.005, $N_P = 6$ for 1/m = 0.01, $N_P = 5$ for $1/m \in [0.015, 0.035]$, $N_P = 4$ for $1/m \in [0.04, 0.105]$, $N_P = 3$ for $1/m \in [0.11, 0.285]$, $N_P = 2$ for $1/m \in [0.29, 0.775]$, and $N_P = 1$ for $1/m \in [0.78, 1.5]$.

In the above numerics, we have used the exact expression for Ω_m and Ω_{-2} . In full QCD, we will not know the exact expression. Therefore, it makes sense to study how well the exact result is reproduced by its semiclassical expansion. We observed in [1] that Ω_m is very well saturated by the first terms of such expansion. Truncating the expansion produces differences much smaller than the typical precision of the different terms of the hyperasymptotic expansion. For Ω_{-2} , we compare in Tables I and II the exact result and the truncated semiclassical expansion for an illustrative set of values. We observe that the exact result is very well saturated by the first terms of the expansion computed in Eq. (38). Truncating the expansion produces differences much smaller than the typical precision of the different terms of the hyperasymptotic expansion. As expected $n_f = 3$ is better than $n_f = 0$. Note that in the large β_0 approximation we exactly have $\Lambda = \mu e^{-2\pi/(\beta_0 \alpha(\mu))}$.

An alternative, very effective, presentation of the above results can be done by plotting the relative accuracy of the prediction at each order in α , and at each order of the superasymptotic expansion. We note that we have one observable for each value of \bar{m} . Therefore, for illustration,



FIG. 6. Comparison of lattice and $\overline{\text{MS}}$ scheme results for $n_f = 3$. Upper panel: We plot m_{PV} and the differences: (a) $m_{\text{PV}} - m_P$, and (b) $m_{\text{PV}} - m_P - \bar{m}\Omega_m$ in the lattice and $\overline{\text{MS}}$ scheme with $n_f = 3$ light flavors. Lower panel: Figs. 4 and 5 combined.

we take two extreme cases. We use $m_{\rm PV}$ with $\bar{m} =$ 1.25 GeV and $\bar{m} = 163$ GeV. For the theoretical prediction we take the smallest positive value of c corresponding to lattice or $\overline{\text{MS}}$ scheme.³ We use the exact expressions for Ω_m and Ω_{-2} . Nevertheless, the NNLO truncated expression for Ω_m is precise enough to yield the same result. For Ω_{-2} we could truncate earlier with no visible effect. We show the results in Fig. 7. We stress that several terms of the hyperasymptotic expansion are included. First, we nicely see that, once reached the minimum, $N \sim N_P$, both schemes yield similar precision, but in the lattice scheme (bigger factorization scale μ) more terms of the perturbative expansions are needed to reach the same precision. We can see a gap when Ω_m is included, with significant better precision in the \overline{MS} scheme. One important lesson one may extrapolate from this exercise is that, if the number of perturbative coefficients is fixed, the smaller the renormalization scale μ , the better. One can obtain much better precision for an equal number of perturbative coefficients. Another observation is that the minimal term determined numerically need not to coincide with the minimal term

³Taking different values of c do not change the picture. The new points stand on top of the old ones where they overlap.

TABLE I. $\bar{m}\Omega_{-2}$ in the large β_0 approximation for $n_f = 0$ in r_0^{-1} units compared with Eq. (38) truncated at different powers of α . Upper panel computed in the $\overline{\text{MS}}$ scheme. Lower panel in the lattice scheme. Lattice seems to be better but both schemes yield very good results.

$\overline{\text{MS-Scheme}} \ (n_f = 0)$						
\bar{m} in r_0^{-1}	С	$\bar{m}\Omega$ Exact	$\left \frac{\Omega \text{LO}}{\Omega \text{Exact}} - 1\right \times 10^2$	$\left \frac{\Omega \text{NLO}}{\Omega \text{Exact}} - 1\right \times 10^3$	$\left \frac{\Omega \text{NNLO}}{\Omega \text{Exact}} - 1\right \times 10^4$	
0.6667	0.1786	0.2089	33.8725	147.64	3372.63	
0.8333	0.5693	0.0572	8.1940	93.6387	922.993	
1	0.8885	0.0362	14.1752	45.6019	16.0275	
1.25	1.2791	0.0260	5.7282	13.6703	130.93	
1.6667	0.0321	0.0199	12.0969	7.3723	12.2818	
2.5	0.7419	0.0094	4.9357	7.7804	6.5465	
5	0.2047	0.0042	2.9590	0.5254	4.8485	
10	1.4182	0.0018	0.2254	2.2970	2.3334	
100	0.1972	0.0001	1.2994	0.1190	0.3921	
		L	attice-Scheme $(n_f = 0)$			
\bar{m} in r_0^{-1}	С	$\bar{m}\Omega Exact \times 10^9$	$\left \frac{\Omega LO}{\Omega Exact} - 1\right \times 10^3$	$\left \frac{\Omega \text{NLO}}{\Omega \text{Exact}} - 1\right \times 10^4$	$\left \frac{\Omega \text{NNLO}}{\Omega \text{Exact}} - 1\right \times 10^5$	
0.6667	0.8101	0.33643	23.0068	12.0234	0.06620	
0.8333	1.2008	0.26320	13.94	5.2020	12.6513	
1	1.5200	0.21971	11.9588	13.8372	5.6692	
1.25	0.1599	0.17233	20.0735	0.3374	6.8322	
1.6667	0.6636	0.12061	15.0134	9.1010	3.3885	
2.5	1.3734	7.7980	1.2511	7.4891	5.2999	
5	0.8362	3.5950	14.9013	4.4610	0.4998	
10	0.2990	1.7262	4.2813	2.3906	2.5543	
100	0.8287	14.527	9.6569	1.8830	0.05590	

TABLE II. $\bar{m}\Omega_{-2}$ in the large β_0 approximation for $n_f = 3$ in GeV units compared with Eq. (38) truncated at different powers of α . Upper panel computed in the $\overline{\text{MS}}$ scheme. Lower panel in the lattice scheme. Lattice seems to be better but both schemes yield very good results.

$\overline{\text{MS}}$ -Scheme ($n_f = 3$)						
\bar{m} in GeV	С	$\bar{m}\Omega$ Exact	$\left \frac{\Omega \text{LO}}{\Omega \text{Exact}} - 1\right \times 10^2$	$\left \frac{\Omega \text{NLO}}{\Omega \text{Exact}} - 1\right \times 10^3$	$\left \frac{\Omega \text{NNLO}}{\Omega \text{Exact}} - 1\right \times 10^4$	
0.6667	0.4916	0.00375	3.6513	8.8280	16.5412	
0.8333	0.8113	0.00274	5.0007	2.9016	7.864	
1	1.0724	0.00223	1.6450	3.9632	11.1774	
1.25	1.3921	0.00183	6.2621	3.6008	0.7661	
1.6667	0.3717	0.00119	0.8631	2.7541	4.6880	
2.5	0.9525	0.00072	2.2110	0.5893	3.1281	
5	0.5130	0.00032	1.7608	1.5455	0.9363	
10	0.0735	0.00015	2.8072	0.2396	0.5520	
100	0.5069	0.00001	0.9405	0.4412	0.1384	
		La	attice-Scheme $(n_f = 3)$			
\bar{m} in GeV	С	$\bar{m}\Omega Exact \times 10^{11}$	$\left \frac{\Omega LO}{\Omega Exact} - 1\right \times 10^3$	$\left \frac{\Omega \text{NLO}}{\Omega \text{Exact}} - 1\right \times 10^4$	$\left \frac{\Omega \text{NNLO}}{\Omega \text{Exact}} - 1\right \times 10^5$	
0.6667	0.6457	0.13921	14.5539	5.0568	0.1286	
0.8333	0.9653	0.10969	9.6413	1.9386	3.4992	
1	1.2264	9.1458	6.7772	6.1070	1.8276	
1.25	0.1137	7.2422	15.151	0.1208	2.0634	
1.6667	0.5258	5.1755	9.9772	4.3766	1.2143	
2.5	1.1065	3.3731	1.7019	3.6030	1.9541	
5	0.6670	1.5918	10.8713	2.5255	0.06625	
10	0.2275	77.282	4.1604	1.2278	1.1874	
100	0.6609	674.45	7.7077	1.7587	4.5364	



FIG. 7. $|m_{PV} - m_{PV}^{Hyperasymptotic}|$ for $\bar{m} = 163$ GeV (upper panel) and $\bar{m} = 1.25$ GeV (lower panel). Blue points are $|m_{PV} - m_N|$. Orange points are $|m_{PV} - m_P - \bar{m}\Omega_m - \sum_{n=N_P+1}^{N} (r_n - r_n^{(as)})\alpha^{n+1}|$ with c = 1.21/1.39 and c = 1.36/0.11 (the smallest positive values that yield integer N_P) in the \overline{MS} and lattice scheme respectively for $\bar{m} = 163/1.25$ GeV. Green points are $|m_{PV} - m_P - \bar{m}\Omega_m - \sum_{n=N_P+1}^{2N_P} (r_n - r_n^{(as)})\alpha^{n+1} - \bar{m}\Omega_{-2} - \sum_{n=N_P+1}^{N} (r_n - r_n^{(as)})\alpha^{n+1}|$, where in the last sum the two first renormalons are subtracted. Change of color correspond to the inclusion of Ω_m and Ω_{-2} . Full points have been computed in the \overline{MS} scheme and empty points in the lattice scheme. We work with $n_f = 3$.

computed using $N = N_P$ (though it should not be much different). The difference reflects how much the exact coefficient is saturated by the asymptotic expression. The effect of Ω_{-2} is very small compared with the effect due to Ω_m . For the case of the top ($\bar{m} = 163 \text{ GeV}$) we can still see the sign alternating behavior of the perturbative series associated to the d = -2 renormalon in the \overline{MS} scheme. In the lattice scheme the effect is so small that it cannot be seen and the precision is set by the next renormalon located at d = 3. If one makes \bar{m} small, $\bar{m} = 1.25$ GeV, green and orange points mix in the MS scheme. This effect is more pronounced in the lattice scheme, where one can continuously move from the orange to the green points. The effect of the ultraviolet renormalon is very small and the precision is set by the u = 3/2renormalon.

B. $(N,\mu) \rightarrow \infty$. Eq. (4). Case (2)

We take Eq. (29) in the large β_0 limit by setting b = 0. As before, we have no analytic expressions to compare with (unlike the case of the static potential). Therefore, we directly focus on taking the limit (2B) and numerically check its convergence and how it compares with method (1).

Method (2B) has the pleasant feature that the generated $\mathcal{O}(\Lambda_{\text{OCD}})$ correction complies with the OPE. It also yields results that do not depend on N (and μ) anymore. Still, it has some errors and does not reach the precision of method (1). There is a residual scheme dependence associated to uncomputed terms of $\mathcal{O}(\alpha \Lambda_{\text{OCD}})$. Part of it can be estimated by the residual dependence in c'. In order to estimate it, we compute m_A for different values of c'. On the one hand c'cannot be very large, as $c'\alpha(\bar{m})$ should be relatively close to zero. On the other hand we cannot make $c'\alpha(\bar{m})$ to get arbitrary close to zero, as the $\mathcal{O}(\Lambda_{\text{OCD}})$ correction diverges logarithmically in c'. We also note that there is a value of $c' = c'_{\min}$ that makes that $K_X^{(A)} = 0$ so that the $\mathcal{O}(\Lambda_{\rm QCD})$ correction vanishes. Therefore, we compute m_A for different values of c'. For illustration we show some results in Fig. 8. We draw lines for $m_{\rm PV} - m_A - K_X^{(A)} \Lambda_X$ at c' = 1 and $c' = c_{\rm min}$ generating a band. We also explore the dependence on the scheme by comparing the results in the lattice and $\overline{\text{MS}}$ scheme. We stress again that, in the large β_0 approximation, lattice and \overline{MS} schemes basically correspond to a redefinition of μ , but quite large indeed. On the other hand the final result is μ independent. Nevertheless, the way the $\mu \to \infty$ limit is taken is fixed by N_A , as defined in Eq. (4), which is dependent on μ . This explains why different results are obtained.

In Fig. 8, we also compare with results obtained using method (1), more specifically we compare with $m_{\rm PV} - m_P - \bar{m}\Omega_m$, as they both have analogous power accuracy (though method (1) is parametrically more precise). For Ω_m we take the exact expression but using its approximated expression does not change the discussion, as the difference is very small. What we see is that the $\overline{\rm MS}$ scheme yields more precise predictions than the lattice scheme, and that method (1) yields considerable better results than method (2B).

Another issue specific of method (2B) is to determine how large we need to take N (and consequently μ) of the truncated sum such that it approximates well m_A . For illustrative purposes we show the convergence in Fig. 9 for $n_f = 3$ in the lattice and $\overline{\text{MS}}$ scheme. We find that we have to go to relatively large values of μ (and N) to get it precise. This can be a problem if one wants to go beyond the large β_0 . This problem would be less severe if one can use the asymptotic expression for the coefficients beyond certain n. Nicely enough, we find that the use of asymptotic expression for the coefficients for $n > N^*$ (~3 in the $\overline{\text{MS}}$ and ~8 in the lattice scheme) is very efficient and basically yields the same results as the exact result. Finally, we also remind





FIG. 8. Upper panel: We plot (a) $m_{\rm PV} - m_A - K_X^{(A)} \Lambda_X$ for $n_f = 0$ in the lattice and $\overline{\rm MS}$ scheme. For each case, we generate bands by computing m_A with c' = 1 and $c' = c'_{\rm min} = 0.652$. We also compare with (b) $m_{PV} - m_P - \bar{m}\Omega_m$ obtained with method (1) with the bands generated for Fig. 3. Lower panel: As the upper panel with $n_f = 3$, $c'_{\rm min} = 0.534$ and taking the bands obtained with method (1) for Fig. 6 for (b) $m_{\rm PV} - m_P - \bar{m}\Omega_m$.

that to approximate well m_A by the truncated sum is more costly for small values of c'.

IV. $\bar{\Lambda}_{PV}$ FROM LATTICE AND *B* PHYSICS

We now abandon the large- β_0 approximation. Our aim is to determine $\bar{\Lambda}_{PV}$. We will determine it first in gluedynamics ($n_f = 0$) in Sec. IV B. To study the scheme dependence of the result it will be useful to estimate the higher order coefficients of the β function in the Wilson action lattice scheme. We do so in the next section.

A. β -function coefficients in the Wilson action lattice scheme

In [22,23] it was shown that renormalon dominance allowed to give an accurate value for β_3^{latt} assuming that c_3 [see Eq. (44)] is already saturated by the renormalon in the $\overline{\text{MS}}$ scheme. We can estimate higher order terms of the β function in the lattice scheme (using the Wilson action) by also assuming that for n > 3 the coefficients c_n in the $\overline{\text{MS}}$ scheme are saturated by the renormalon. We show such



FIG. 9. Upper panel: We plot $m_{\rm PV} - m_A - K_X^{(A)} \Lambda_X$ for $n_f = 3$ in the lattice scheme with c = 1 versus the truncated sums $m_{\rm PV} - \sum_{n=0}^{N_A} r_n \alpha^{n+1}(\mu) - K_X^{(A)} \Lambda_X$, where μ is fixed using N_A defined in Eq. (4). Lower panel: As in the upper panel but in the $\overline{\rm MS}$ scheme.

estimates in Table III. These coefficients of the β function improve the agreement with the phenomenological parametrization of $\alpha_{\text{latt}}(1/a)$ obtained in [28] in the range $\beta \in$ (6, 6.8) (see Fig. 10). It is also worth mentioning that we observe a geometrical growth of the coefficients of the β function. Elucubrative, this would indicate that the beta function in this scheme has a finite radius of convergence, and one can take the ansatz

$$\beta^{\text{latt}}(\alpha) = \nu \frac{d}{d\nu} \alpha$$
$$\simeq -2\alpha \left\{ \sum_{n=0}^{3} \beta_n \left(\frac{\alpha}{4\pi}\right)^{n+1} - 1.4 \times 10^8 \left(\frac{\alpha}{4\pi}\right)^5 \frac{1}{1 - 10^2 \frac{\alpha}{4\pi}} \right\},\tag{42}$$

which would have a pole at around $\beta = 6/g^2 \simeq 3.8$.

B. $\bar{\Lambda}_{PV}$ from lattice

We determine $\bar{\Lambda}_{PV}$ in gluedynamics $(n_f = 0)$ from the energy of a meson made of a static quark and a light valence quark:

TABLE III. Estimates of the coefficients of the beta function for the bare coupling in the lattice scheme using renormalon dominance and $Z_m^{\overline{\text{MS}}} = 0.62$ [23]. The error quoted in the table gives the difference with the values of the beta coefficients obtained if one uses instead $Z_m^{\overline{\text{MS}}} = 0.6$ [13] (which yields more negative values), and it is only meant to illustrate the typical spread of values of the beta coefficients if one uses different values of $Z_m^{\overline{\text{MS}}}$.

β_3	eta_4	β_5	β_6
$-1.16(3) \times 10^{6}$	$-1.35(10) \times 10^{8}$	$-1.44(28) \times 10^{10}$	$-1.41(60) \times 10^{12}$

$$E_{\rm MC}(a) = \delta m_{\rm latt}^{\rm PV} + \bar{\Lambda}_{\rm PV} + \mathcal{O}(a\Lambda_{\rm QCD}^2). \tag{43}$$

 $\delta m_{\text{latt}}^{\text{PV}}$ has the following asymptotic series in powers of $\alpha = \alpha_{\text{latt}}(1/a)$:

$$\delta m_{\text{latt}}^{\text{PV}} \sim \sum_{n=0}^{\infty} \frac{1}{a} c_n \alpha^{n+1},$$
 (44)

where $\frac{r_n^{(as)}(\nu)}{\nu} = c_n^{(as)}$, since $m_{\rm PV}$ and $\delta m_{\rm latt}^{\rm PV}$ have the same leading infared renormalon (located at d = 1). The coefficients c_n are known from $n = 0 \div 19$ in the lattice scheme for a Wilson action [21–23]. We then adapt Eq. (19) to $\delta m_{\rm latt}^{\rm PV}$ to determine $\bar{\Lambda}_{\rm PV}$:

$$\bar{\Lambda}_{PV}(n_f = 0) = E_{MC}(a) - \delta m_{latt}^P - \frac{1}{a}\Omega_m - \sum_{N_P+1}^{N'=2N_P} \frac{1}{a} [c_n - c_n^{(as)}]\alpha^{n+1} + \mathcal{O}(a\Lambda_{QCD}^2),$$
(45)

where $\delta m_{\text{latt}}^P = \sum_{n=0}^{N_p} \frac{1}{a} c_n \alpha^{n+1}$. In the counting of Eq. (15) this corresponds to $(1, N_P)$ precision. The expression we use for Ω_m is Eq. (21) truncated to $\mathcal{O}(\alpha^3)$ here and in the rest of the paper. The error committed by this truncation is smaller than the error associated to Z_m . Therefore, we will neglect it in the following. The renormalon behavior associated to subleading renormalons of $E_{\text{MC}}(a)$ is not well known,



FIG. 10. Same caption as in Fig. 1 of [29] including more terms in the perturbative expansion using the β -function coefficients listed in Table III.

except that the next singularity in the Borel plane is expected to be at |u| = 1 (d = 2). Therefore, we stop the second perturbative expansion at $N' = 2N_P$ such that the reminder should be of $\mathcal{O}(a\Lambda_{\text{OCD}}^2)$. For the coefficients $c_n^{(as)}$ we use $Z_m^{\text{latt}}(n_f = 0) = 17.9(1.0)$ [23]. We also truncate the 1/n expansion in Eq. (26) to $\mathcal{O}(1/n^3)$. This means using the estimates for β_3 and β_4 listed in Table III. We take $E_{\rm MC}(a)$ from [30–32]. These points expand over the following energy range: $1/a \sim 2.93 r_0^{-1} \div 9.74 r_0^{-1}$. We show our results in Fig. 11. They follow the same logic than Figs. 1-6 in Sec. III. We observe that the subtraction of the perturbative expansion accounts for most of the 1/adependence. Still we have enough precision to be sensitive to $\mathcal{O}(a\Lambda_{\text{QCD}}^2)$ effects. A strict fit setting the $\mathcal{O}(a\Lambda_{\text{QCD}}^2)$ correction to zero gives a large $\chi^2_{red} \sim 6-7$. The inclusion of a pure *Ka* term to Eq. (45) gives a good fit.⁴ The statistical error is small and the $\chi^2_{\rm red} = 1.17/1.06$ (for the smallest |c|with positive/negative c value) is good. Overall, we obtain (using the smallest c positive, which means $N_P = 7$ except for $\beta = 5.7$ where we have $N_P = 6$)

$$\bar{\Lambda}_{\rm PV} = 1.42 r_0^{-1} (\text{stat.})^{-0.01}_{+0.04} (c)^{+0.05}_{-0.05} (Z_m)^{+0.16}_{-0.16}.$$
 (46)

This number is not very different from the number obtained in [29] using a superasymptotic approximation truncated at the minimal term determined numerically (typically this always gives slightly better results than truncating at the minimal term predicted by theory).

Let us now discuss the error budget in Eq. (46). The first error is the statistical error of the fit. The remaining errors are different ways to estimate the error produced by the approximate knowledge of the hyperasymptotic expansion. One possibility is to take the modulus of the difference with the evaluation using the *c* negative with the smallest possible modulus. This is the second error we quote in Eq. (46). The last error we include is due to the variation of $Z_m^{\text{latt}}(n_f = 0) = 17.9(1.0)$ [23] (correlated with the

⁴Unlike for the pole mass, it is not clear what is the operator of the OPE that would produce the NP correction and the associated u = 1 renormalon. Therefore, if for the pole mass we can be certain that the NP correction has the form Ka, without any anomalous dimensions nor any nontrivial $\ln(a)$ dependence, we can not exclude the possibility that this $\mathcal{O}(a)$ correction may have a non trivial anomalous dimension and/or $\ln(a)$ dependence.



FIG. 11. Upper panel: $E_{\rm MC}$ is the Monte Carlo lattice data [30–32]. The continuous lines are drawn to guide the eye. The other lines correspond to Eq. (45) truncated at different orders in the hyperasymptotic expansion. (a) $E_{\rm MC}(a) - \delta m_P(1/a)$, (b) $E_{\rm MC}(a) - \delta m_P(1/a) - \frac{1}{a}\Omega_m$, (c) $E_{\rm MC}(a) - \delta m_P(1/a) - \frac{1}{a}\Omega_m - \sum_{N_P+1}^{N'=2N_P} \frac{1}{a}[c_n - c_n^{(\rm as)}]a^{n+1}$ (in this last case we include the error of the lattice points in the middle of the band), (d) is the fit of the right-hand side of Eq. (45) to $\bar{\Lambda}_{\rm PV}(n_f = 0) - Ka$. For each difference and for the final fit, the bands are generated by the difference of the prediction produced by the smallest positive or negative possible values of *c* that yield integer values for N_P . Lower panel: As in the upper panel but in a smaller range. $r_0^{-1} \approx 400$ MeV.

error of c_n). The error it produces in Ω_m is small. Comparatively, most of the error associated to Z_m comes from the differences in $\sum_{N_p+1}^{N'=2N_p} \frac{1}{a} [c_n - c_n^{(as)}] \alpha^{n+1}$ evaluated at different Z_m . Whereas $\sum_{N_p+1}^{N'=2N_p} \frac{1}{a} [c_n - c_n^{(as)}] \alpha^{n+1}$ is

quite small for the central value of Z_m , it significantly changes after variation of Z_m . This variation is only partially compensated by the variation of the coefficients c_n , which have smaller errors, producing a significant change in $\sum_{N_p+1}^{N'=2N_p} \frac{1}{a} [c_n - c_n^{(as)}] \alpha^{n+1}$. We have also determined the central value in Eq. (46) not including the $\mathcal{O}(1/n^3)$ corrections in the asymptotic expressions for $c_n^{(as)}$. The difference we obtain is -0.08. This is significant, showing that the 1/n corrections are sizable in the lattice scheme. On the other hand, the difference is well inside the error associated to Z_m . Actually, the difference with evaluations including the $\mathcal{O}(1/n^4)$ corrections in the asymptotic expressions for $c_n^{(as)}$ is -0.03. This shows a convergent pattern, which we illustrate in Table IV. Overall, the largest source of uncertainty comes from the incomplete knowledge of $\sum_{N_p+1}^{N'=2N_p} \frac{1}{a} [c_n - c_n^{(as)}] \alpha^{n+1}$, which is closely linked to the incomplete knowledge of Z_m . This discussion points to that more accurate determinations of Z_m can be possible and, then, that the error of $\bar{\Lambda}_{PV}$ associated to Z_m could be made smaller. We believe these issues deserve further study that we leave for future work.

One error that we do not include here is the error associated to α . From the lattice point of view, we are talking of the relation between $\alpha(1/a)$ and r_0 . We use the phenomenological formula deduced in [28]. The error of this formula is claimed to be around 0.5-1% in the range $\beta \in (5.7, 6.92)$ [having a look to Fig. 10 a more conservative range could be (6,6.8)].

1. Scheme dependence

It is interesting to consider the scheme dependence of Eq. (46). In [29] relative large differences were found for fits to $\overline{\Lambda}$ after (approximated) scheme conversion to the $\overline{\text{MS}}$ scheme. The real problem is not transforming the coefficients c_n from the lattice to the $\overline{\text{MS}}$ scheme, but transforming α_{latt} to $\alpha_{\overline{\text{MS}}}$ with enough precision (in a way we need the relation between α_{latt} and $\alpha_{\overline{\text{MS}}}$ with NP, exponential, accuracy). This needs the coefficients of the β function in the lattice scheme to high orders. We show estimates in Table III. The inclusion of these coefficients of the β function makes that the determinations of $\overline{\Lambda}_{\text{PV}}$ in the $\overline{\text{MS}}$ and lattice scheme approach each other as we include more terms in the perturbative expansion of the relation between $\alpha_{\overline{\text{MS}}}$ and α_{latt} . We show the comparison in Table IV.

TABLE IV. Determinations of $\bar{\Lambda}_{PV}$ in the lattice and \overline{MS} scheme from fits of $\bar{\Lambda}_{PV} - Ka$ to the right-hand side of Eq. (45). The first three numbers show the impact in the fit of including the $\mathcal{O}(1/n^m)$ corrections for m = 2, 3, 4 in the asymptotic expressions for $c_n^{(as)}$ in the lattice scheme (in the \overline{MS} this effect is negligible). The other numbers are the fit of $\bar{\Lambda}_{PV}$ in the \overline{MS} scheme, using $\alpha_{\overline{MS}} = \alpha_{\text{latt}}(1 + \sum_{n=0}^{N_{tr}} d_n \alpha_{\text{latt}}^n)$ truncated at $N_{tr} = 4, 5, 6, 7$.

latt	$\mathcal{O}(\frac{1}{n^2})$	$\mathcal{O}(\frac{1}{n^3})$	$\mathcal{O}(\frac{1}{n^4})$	MS	$N_{tr} = 7$	$N_{tr} = 6$	$N_{tr} = 5$	$N_{tr} = 4$
$\bar{\Lambda}_{\rm PV}$	1.33	1.42	1.45	$\bar{\Lambda}_{PV}$	1.48	1.52	1.59	1.68

C. $\bar{\Lambda}_{pot}^{PV}$ from lattice

As an extra check of the method, we now consider the ground state energy of two static sources in the fundamental representation at a fixed distance r_0 computed in the lattice: $E_{\Sigma_g^+}(r_0; a)$. This object has the same renormalon as twice the pole mass. Following [33] we define the quantity

$$\bar{\Lambda}_{\text{pot}}(a) \equiv \frac{E_{\Sigma_g^+}(r_0; a)}{2} + \Delta, \tag{47}$$

where Δ is just a constant to fix the normalization at $r = r_0$. For $\bar{\Lambda}_{\text{pot}}(a)$ we perform an OPE assuming $r_0 \gg a$, and compute it using the PV prescription. We then have

$$\bar{\Lambda}_{\text{pot}}^{\text{PV}} \equiv \frac{E_{\Sigma_{g}^{+}}(r_{0}; a)}{2} + \Delta - \delta m_{\text{latt}}^{P} - \frac{1}{a} \Omega_{m} - \sum_{n=N_{P}+1}^{N'=2N_{P}} \frac{1}{a} [c_{n} - c_{n}^{(\text{as})}] \alpha^{n+1} + \mathcal{O}(a \Lambda_{\text{QCD}}^{2}). \quad (48)$$

We show the results in Fig. 12. The lattice data is taken from [34,35], as analyzed in [33]. A nicely flat curve appears. This object does not show $O(a\Lambda_{\rm QCD}^2)$ artifacts. This is consistent with the discussion in [28], though there the discussion was only made for energy differences. This has the potentially important consequence that potentials computed with different β 's can be related with perturbation theory with good accuracy. There is no need to subtract independent constants for each $\beta \equiv 6/g^2$.

D. $\bar{\Lambda}_{\rm PV}$ from *B* meson mass

We now move to the physical case with $n_f = 3$ light quarks. Using HQET we approximate the B/D meson mass by (we use spin averaged masses)

$$m_{B(D)} = m_{\rm PV} + \bar{\Lambda}_{\rm PV} + \mathcal{O}\left(\frac{1}{m_{\rm PV}}\right). \tag{49}$$

It is not the aim of this paper to determine \bar{m}_b (nor \bar{m}_c). We are rather interested to know the error associated to determinations of $m_{\rm PV}$ if \bar{m} is known, and vice versa. We will then later use this analysis for the top quark mass determination. For this purpose we use its hyperasymptotic approximation

$$m_{\rm PV}(\bar{m}_{b/c}) = m_{\rm P}(\bar{m}_{b/c}) + \bar{m}_{b/c}\Omega_m + \sum_{n=N_P+1}^{N'=2N_P} [r_n - r_n^{\rm (as)}]\alpha^{n+1} + \cdots.$$
(50)

To make the error analysis we use the bottom case, and take $\bar{m}_b = 4.186$ GeV from [36]. We obtain (we have added a -2 MeV to the relation between the $\overline{\text{MS}}$ bottom mass and the pole mass due to the charm quark [13])



FIG. 12. Upper panel: $\bar{\Lambda}_{pot}(a)$ is the Monte Carlo lattice data [34,35], as analyzed in [33]. The continuous lines are drawn to guide the eye. The other lines correspond to Eq. (48) truncated at different orders in the hyperasymptotic expansion: (a) $\bar{\Lambda}_{pot}(a) - \delta m_P(1/a)$, (b) $\bar{\Lambda}_{pot}(a) - \delta m_P(1/a) - \frac{1}{a}\Omega_m$, (c) $\bar{\Lambda}_{pot}(a) - \delta m_P(1/a) - \frac{1}{a}\Omega_m - \sum_{N_P+1}^{N'=2N_P} \frac{1}{a}[c_n - c_n^{(as)}]\alpha^{n+1}$ (in this last case we include the error of the lattice points in the middle of the band). For each difference the bands are generated by the difference of the prediction produced by the smallest positive or negative possible values of *c* that yield integer values for N_P . Lower panel: As in the upper panel but in a smaller range. $r_0^{-1} \approx 400$ MeV.

$$m_{b,\rm PV} = 4836(\mu)^{+8}_{-17}(Z_m)^{-11}_{+12}(\alpha)^{+8}_{-9}$$
 MeV. (51)

For the variation of μ we take the range $\mu \in (\bar{m}_b/2, 2\bar{m}_b)$. For Z_m we take $Z_m^{\overline{\text{MS}}}(n_f = 3) = 0.5626(260)$ from [13]. For the variation of α we take $\Lambda_{\overline{\text{MS}}}^{(n_f=3)} = 332 \pm 17$ MeV from [37]. The central value has been obtained with $N_P = 3$ (c = 0.3611). Therefore, the last term of Eq. (50) is set to zero, as we do not have more terms of the perturbative expansion. In the counting of Eq. (15) the precision is then (1,0). Within the hyperasymptotic counting the $\sum_{n=N_P+1}^{N'=2N_P} [r_n - r_n^{(as)}] \alpha^{n+1}$ term should roughly scale as (assuming the next renormalon is located at |u| = 1)

$$\sim e^{-\frac{2\pi}{\beta_0 a_X(\mu)}(1+\ln(2))}$$
. (52)

This is the expected scaling if $\mu \sim \bar{m}$. Nevertheless, the dependence on μ will be quite different depending on whether the next renormalon is ultraviolet ($\sim \mu^{-2}$) or infrared ($\sim \mu^2$). Actually, the magnitude is also expected to be different, being more important for an eventual infrared renormalon. As the situation is somewhat uncertain we do not dwell further in this issue. To roughly estimate the size of subleading terms we could compute with $N_P = 2$ (c = 1.7935). In the counting of Eq. (15) the precision is then $(1, N_{\text{max}} - N_P)$. The difference is below 1 MeV (after including $[r_3 - r_3^{(as)}]\alpha^4$, otherwise the difference is 7.5 MeV). Even computing with $N_P = 1$, which formally allows us to reach the next renormalon located at $2N_P = 2$ [i.e., $(1, N_P)$ precision in the counting of Eq. (15)], the difference is \sim 7 MeV. Alternatively, the remaining μ scale dependence of $m_{\rm P}(\bar{m}_{b/c}) - \bar{m}_{b/c}\Omega_m$ also gives a measure of the uncomputed $\sum_{N_P+1}^{N'=2N_P} [r_n - r_n^{(as)}] \alpha^{n+1}$ term, as such scale dependence should cancel in the total sum. We will then take it as the associated error. This is the first error quoted in Eq. (54). Actually, the error associated to Z_m is also a measure of the lack of knowledge of higher order terms in perturbation theory. Therefore, there is some degree of double counting by considering these two errors separately.

It is interesting to analyze the error of the superasymptotic approximation of $m_{\rm PV}$ (only computing m_P). If we vary μ in the range $\mu \in (\bar{m}_b/2, 2\bar{m}_b)$, we obtain $(N_P = 3, c = 0.3611)$

$$m_{b,P} = 5077(\mu)^{+134}_{-242}$$
 MeV. (53)

Nicely enough, it agrees with Eq. (51) within one sigma. This is also so if we take $N_P = 2$ (c = 1.7935): $m_{b,P} = 4922^{+107}_{-167}$ MeV. We find that the scale dependence of the superasymptotic approximation is large. Therefore the inclusion of $\bar{m}\Omega_m$ is crucial to make the result much more scale independent. On the other hand note that there is no error associated to Z_m at this order (which is, in any case, comparatively small).

Using Eq. (51) and Eq. (49) we can determine $\bar{\Lambda}_{PV}$. We work at leading order in 1/m. We obtain

$$\bar{\Lambda}_{\rm PV} = 477(\mu)^{-8}_{+17}(Z_m)^{+11}_{-12}(\alpha)^{-8}_{+9}(\mathcal{O}(1/m))^{+46}_{-46} \text{ MeV}, \quad (54)$$

where we have included an extra error source compared with Eq. (51). This extra error is associated to the O(1/m)corrections. The existence or not of genuine NP 1/mcorrections may introduce a significant error. In case they exist, if we take the hyperfine energy splitting as a measure of 1/m corrections, we find shifts from the central values of order ~46 MeV and ~140 MeV for *B* and *D* mesons, respectively. As Eq. (54) has been obtained from the *B* meson spin-average mass we conservatively estimate the error associated to genuine NP 1/m corrections to be of order ~46 MeV, as it is the most we can do from phenomenology and perturbation theory. Let us recall however that recent lattice simulations point to much smaller genuine NP 1/m corrections for the spin-independent average [38].

Earlier direct determinations of $\bar{\Lambda}_{PV}$ or m_{PV} can be found in [39,40]. The formulas are equivalent to those used here to one order less (using $N_P = N_{max} = 2$). They also include less terms in the sum in Eq. (26). More recently, a determination of $\bar{\Lambda}$ has been obtained in [38] using lattice data. In this case the formulas are equivalent to those used here since $N_P = 3$ (see Eq. (57) of Ref. [1]) except for the fact that the scale μ was always fixed equal to the heavy quark mass and that the mass was obtained in the MRS scheme [41]. In this reference is also given the relation between the PV and the MRS mass. Using it we obtain (where we combine quadratically the error of $Z_m^{\overline{MS}}$ and $\Lambda_{\overline{MS}}$)

$$\bar{\Lambda}_{PV} - \bar{\Lambda}_{MRS} = \cos(\pi b) \frac{4\pi\Gamma(-b)}{2^{1+b}\beta_0} Z_m^X \Lambda_X \bigg|_{n_f=3}$$
$$= -120(8) \text{ MeV}.$$
(55)

The prediction of [38] translates then to $\bar{\Lambda}_{PV} = 435(31)$, where we only include the error quoted in [38]. In particular, we do not include the error in Eq. (55). Note that Eq. (55) scales like $\mathcal{O}(\Lambda_{\text{OCD}})$, whereas $\bar{m}\Omega_m$ scales like $\mathcal{O}(\sqrt{\alpha}\Lambda_{\rm OCD})$. There is a 40 MeV difference with the number given in Eq. (54). 10 MeV can be understood because the value of \bar{m}_b used in [38] is around 10 MeV bigger. Another 10 MeV can be understood by the inclusion of 1/m nonperturbative effects. The remaining 20 MeV difference are more difficult to identify, though they are well inside uncertainties. Leaving aside the different α 's used, another source of difference is the value of Z_m . The value used in [38] comes from [42] (where the effect of scale variation was not included in the error analysis). This determination used a sum rule that is free of the leading pole mass renormalon. The possibility of using sum rules to determine the normalization of renormalons was first considered in [43]. For the determination of Z_m , sum rules were first used in [12]. Later sum rule analyses can be found in [44]. Alternatively one can use the ratio of the exact and asymptotic expression of the coefficients r_n to determine Z_m as in [13,21–23,45]. For an extra discussion on this issue see [46]. Finally, it is worth mentioning that Z_m can be determined either from the static potential or from the pole mass (and its relatives). The only value of Z_m that uses the static potential is from [13]. A preference for determinations of Z_m from the static potential can be theoretically motivated, as they are less affected by subleading renormalons. There are no ultraviolet renormalons and the next infrared renormalon is located at u = 3/2. On the other hand, the pole mass is expected to have renormalons at |u| = 1. Only in the event that there is no u = 1 renormalon and the effect of the u = -1 renormalon is subleading both determinations would be on equal footing on theoretical grounds. In any case, irrespectively of this discussion, consistent numbers are obtained between different analyses.

E. $(N,\mu) \rightarrow \infty$. Eq. (4). Case (2B)

All previous determinations of $\bar{\Lambda}_{PV}$ have been obtained using limit (1). For completeness we have also explored how limit (2B) performs for $\bar{\Lambda}_{PV}$, even though it is, in principle, less precise. We have considered the different methods to take the limit (2B) discussed at the end of Sec. II B, and compared with the numbers obtained above. We first consider the evaluation of m_{PV} using the right-hand side of Eq. (33) with Eq. (34). The central value is determined using $c'_{\min} = 1.076$, which is the value that makes $K_{\overline{\text{MS}}}^{(A)} = 0$, and $\mu_0 = \bar{m}_b = 4.186$ MeV. We obtain $\bar{\Lambda}_{PV} = 453$ MeV. The difference with Eq. (54) is 24 MeV, which is quite reasonable. We can also explore the μ dependence. Taking the variation $\mu_0 \in (\bar{m}_b/2, 2\bar{m}_b)$, we obtain $\bar{\Lambda}_{PV} = 453^{-36}_{+55}(\mu)$ MeV. Comparatively with Eq. (54) the μ scale dependence is much larger. We next consider the limit as taken in Eq. (32). This requires the knowledge of the coefficients r_n to all orders. For n > 3 we take the asymptotic expression. On the other hand the running of the beta function is only needed to one loop. This allows us to go to orders as high as $N_A = 3000$ (though it already converges at smaller values of N_A). Remarkably enough, we obtain the same result than before: 453 MeV. There is a residual dependence on c'. For illustration, if we take instead c' = 2, we obtain $\bar{\Lambda}_{PV} = 438$ MeV (the result using the right hand side of Eq. (33) with Eq. (34) yields the same number), and the scale dependence is larger: $\bar{\Lambda}_{\rm PV} = 438^{-55}_{+99}(\mu)$. The value of c' we have used to make the analysis can be an issue. As discussed in [1,17], taking $\chi - 2$ very small deteriorates the convergence and larger values of N_A are needed. This problem aminorates by taking larger values of c'. Since for the limit as taken in Eq. (32) we can go to very large N_A this is not a problem. We have also performed a similar analysis with $n_f = 0$ and r_0 units, relevant for the analyses performed in Sec. IV B. The discussion follows parallel to the one we just had with the difference that we now know 20 coefficients of the perturbative expansion. The value we obtain: $\bar{\Lambda}_{\rm PV} = 1.37 r_0^{-1}$ (using a quadratic fit) is indeed quite close to the value obtained in Sec. IV B, though less precise.

We have more problems with the other ways to take the $\mu \to \infty$ limit discussed at the end of Sec. II B. The direct use of N_A in Eq. (4) or of N_A in Eq. (67) in [1] requires, besides the coefficients r_n to all orders, the β -function coefficients to all orders as well. We do not have them. Instead we use truncated version of the β function. This makes the numerical calculation much more challenging,

since the running in μ is more complicated. Therefore, we had problems to go to very large N_A . For $N_A \ge 200$ we find instabilities is some cases. As mentioned before, the value of c' we use to make the analysis can be an issue. Taking $\chi - 2$ very small deteriorates the convergence. This problem aminorates by taking larger values of c'. In the lattice scheme determination of quenched $\bar{\Lambda}_{PV}$ we indeed observe convergence to the value obtained before using c' = 2. Using $c'_{min} = 1.076$ the convergence is less good. Determinations in the \overline{MS} scheme do not show convergence if we stop at $N_A \le 200$, though with an slight better behavior using c' = 2 rather than c'_{min} . Overall, as the precision we get with method (2B) is worse than with method (1), we will not study this limit in more detail.

V. TOP MASS

A. About the pole mass ambiguity

The top quark mass is one of the key parameters of the standard model. A lot of experimental work has been devoted to its determination (see for instance [47-49]). Whereas this is a matter of debate, it is typically assumed that the masses obtained from experiment correspond to the pole mass. Thus, there has been an ongoing discussion on the intrinsic uncertainty of these determinations (see for instance [45,50], and [51] for a more recent discussion). We believe that, without further qualifications, the question is ill posed, or may lead to confusion. It is well known that the pole mass is well defined (infrared finite and gauge independent) at finite (albeit arbitrary) order in perturbation theory [52]. It is also well known that such series is divergent.⁵ Therefore, no numerical value can be assigned to the infinite sum of the perturbative series of the pole mass. Truncated sums are well defined but depend on the order of truncation (a detailed discussion relevant for the analysis made in the present paper can be found in [1]). These truncated sums can be related with observables or with intermediate definitions of the heavy quark mass, like the PV mass (which regulates via Borel resummation the infinite sum), in a well-defined way.

In this context, the shortest answer to the above posed question is that the ambiguity (of a well-defined mass) is *zero*. As a matter of principle, m_{PV} (or m_P) can be *defined* with arbitrary accuracy (this also applies to any threshold mass), if one computes high enough orders of the perturbative series, and if \bar{m} is given. One can discuss (actually one can compute) the scheme/scale dependence (if they have) of them. In this respect, there is no much conceptual difference with respect to asking about the scheme/scale dependence of minimal subtraction schemes for the heavy quark masses.

⁵Actually this is only proven in the large β_0 approximation [19,20], and it is also supported by numerical analyses [21–23], but there is no analytic proof.

A quite a different question is to determine the typical difference (that not ambiguity) between (reasonable) different definitions of the pole mass. The short answer to this question is that the differences are (at most) of order Λ_{OCD} for (reasonable) different definitions of the pole mass. We emphasize that one can not be more precise unless stating the specific definition used for the pole mass. For instance, the difference between $m_{\rm PV}$ and m_P is of $\mathcal{O}(\sqrt{\alpha}\Lambda_{\rm OCD})$ with a known prefactor. Truncating the perturbative series at order N near N^* are also legitimate definitions of the pole mass. The typical difference between truncating at different N is of order Λ_{OCD} : see for instance Eq. (62) of [22]. One could even use M_B as a definition for the pole mass. Its difference with $m_{\rm PV}$ is of order $\Lambda_{\rm OCD}$. If one defines an imaginary mass by doing the Borel integral just above the positive real axis, the difference with $m_{\rm PV}$ is of $\mathcal{O}(i\Lambda_{\rm OCD})$. The authors of [45] choose to divide this number by π and take the modulus as their definition of the ambiguity. These examples illustrate that, even if the ambiguity is of $\mathcal{O}(\Lambda_{OCD})$, the coefficient multiplying Λ_{OCD} is arbitrary. Overall, it should be clear that no much more can be said, and we are indeed against of dwelling too much on this issue. Instead, we strongly advocate to avoid generic discussions about the pole mass, which is not well defined beyond perturbation theory, and restrict the discussion to the precision and errors of specific, NP well-defined, heavy quark masses the perturbative expansion of which can be related with the perturbative expansion of the pole mass.

Once working with NP well-defined heavy quark masses like $m_{t,PV}$ or $m_{t,P}$, we can address the more relevant question of determining the precision with which \bar{m}_t can be determined if $m_{t,PV}$ or $m_{t,P}$ is known (and vice versa, if \bar{m}_t is known what is the uncertainty of m_{PV}) with nowadays knowledge of the perturbative expansion. In other words, with which precision the theoretical expression is known. For reference we will take the value $\bar{m}_t = 163$ GeV in the following. We will see in the next section that indeed the precision is quite good and that the error is significantly smaller than typical numbers assigned for the ambiguity of the pole mass. We will not dwell in this paper on the precision with which $m_{t,PV}$ or $m_{t,P}$ can be *determined* from experiment as such discussion is observable dependent.

B. Decoupling and running

We now turn to an issue specific to the top quark (as compared with the bottom and charm quark). The top quark mass is much larger than Λ_{QCD} . The latter is the scale that characterizes renormalon associated effects and it is the precision we want to achieve. This obviously generates ratios of quite disparate scales. In the context of threshold masses with an explicit infrared cutoff ν_f , this calls for resummation of the large logarithms: $\ln \nu_f/m_t$. This is possible, and first done in [33] in the RS scheme (see also [46] for an extra discussion on this issue). Here, we approach the problem in a different way. We want to work with expansions for the perturbative series of the pole mass truncated at the minimal term: m_P , and to improve upon it using hyperasymptotic expansions. Nevertheless, at the scale of the top mass, we do not have enough terms to reach the asymptotic behavior of the perturbative expansion. We use instead that the top quark pole mass and the pole mass of a fictitious top quark with mass m'_t share the same leading infrared renormalon. Therefore, the leading infrared renormalon cancels in the difference. We can then decrease the top mass in a renormalon free way until we reach a top mass low enough that we can use the hyperasymptotic expansion. Such renormalon free running is determined by the following function (not compulsory to take $\mu = \bar{m}$ but it simplifies the computation)

$$\mathcal{F}(\bar{m}, n_f) \equiv \frac{d}{d\bar{m}} (m_{\rm PV}(\bar{m}) - \bar{m})$$

$$\simeq \frac{d}{d\bar{m}} \sum_{n=0} r_n^{(n_f)}(\bar{m}; \mu = \bar{m}) \alpha_{(n_f)}^{n+1}(\bar{m})$$

$$\equiv \sum_{n=1}^{N+1} f_n(\bar{m}) \left(\frac{\alpha_{(n_f)}(\bar{m})}{\pi}\right)^n.$$
(56)

This formula is correct up to $N \sim 2N_P$, since $\bar{m}\Omega_m$ and $r_n^{(as)}$ are independent of \bar{m} [see Eq. (19)], so that their derivative with respect to \bar{m} vanishes. The coefficients $r_n^{(n_f)}$ are evaluated for n_f massless particles. In the context of the MSR threshold mass the running is implemented in a similar way (see, for instance, [44]). Equation (56) makes explicit that such running is just a natural consequence of the relation between observables and their OPEs (for illustration, it follows from the fact that $M_B - M_D$, the *B* minus *D* meson mass difference is free of the leading infrared renormalon), and not linked to an specific threshold mass definition.

There is still another issue specific to the top quark: there are two heavy quarks (the bottom and charm), with masses much larger than $\Lambda_{\rm QCD}$, that generate extra corrections to the pole- $\overline{\rm MS}$ mass relation due to the finite mass of the bottom and charm quark. Therefore, we have for $\bar{m} \sim \bar{m}_t$

$$m_{\rm PV}(\bar{m}) = \bar{m} + \sum_{n=0}^{N_{\rm max}} r_n^{(n_f)}(\bar{m}; \mu = \bar{m}) \alpha_{(n_f)}^{n+1}(\bar{m}) + \delta m_b^{(n_f)}(\bar{m}) + \delta m_c^{(n_f)}(\bar{m}) + \delta m_{bc}^{(n_f)}(\bar{m}),$$
(57)

where it is implicit that N_{max} (the number of known terms of the perturbative expansion) is not large enough to see the decoupling of the bottom nor charm and certainly $N_{\text{max}} < N_P$. n_f stands for the number of active flavors. At the top mass scale we take $n_f = 5$. The $\mathcal{O}(\alpha^2)$ term of $\delta m_Q^{(n_f)}$ was computed in [53] and the $\mathcal{O}(\alpha^3)$ term in [54]. Note as well that at $\mathcal{O}(\alpha^3)$ there is a new contribution including a vacuum polarization of the bottom and charm at the same time. We name it $\delta m_{bc}^{(n_f)}$ and it has been computed in [50]. As we decrease the value of \bar{m}_t the bottom and charm

As we decrease the value of \bar{m}_t the bottom and charm quark will decouple. This decoupling will be absorbed in $\delta m_{b/c/bc}^{(n_f)}$, which are polynomials in powers of $\alpha^{(n_f)}$. In general this is not just changing n_f in the original expressions from $n_f = 5$ to $n_f = 4$ or 3. The explicit expressions can be found in the Appendix B. The renormalon is associated to scales smaller than the bottom and charm quark masses. Therefore, such scales should be decoupled before we talk about the hyperasymptotic expansion. As we have mentioned above we do such decoupling by varying the mass of the top till reaching a fictitious top with a mass small enough such that, first the bottom, and later the charm, decouple. Overall, our final formula is the following:

of the top it becomes much more effective to decouple, first

the bottom, and afterwards the charm quark. Once this is

done the corrections due to the bottom and charm masses to

$$m_{\rm PV}(\bar{m}_t) = \bar{m}_t + \int_{\mu_b}^{\bar{m}_t} d\bar{m} \left(\mathcal{F}(\bar{m}, 5) + \frac{d}{d\bar{m}} \left(\delta m_b^{(5)}(\bar{m}) + \delta m_c^{(5)}(\bar{m}) + \delta m_{(bc)}^{(5)}(\bar{m}) \right) \right) \\ + \int_{\mu_c}^{\mu_b} d\bar{m} \left(\mathcal{F}(\bar{m}, 4) + \frac{d}{d\bar{m}} \left(\delta m_b^{(4)}(\bar{m}) + \delta m_c^{(4)}(\bar{m}) + \delta m_{(bc)}^{(4)}(\bar{m}) \right) \right) + m_{\rm PV}(\mu_c) - \mu_c.$$
(58)

We emphasize that $\mathcal{F}(\bar{m}, n_f)$ is expanded in powers of α before integration. We take μ_b small enough such that the bottom decouples and μ_c small enough such that the bottom and charm decouple, and also such that we reach the asymptotic limit of the pole- $\overline{\text{MS}}$ mass perturbative expansion with the existing known coefficients. Therefore,

$$m_{\rm PV}(\mu_c) = m_P(\mu_c) + \mu_c \Omega_m + \delta m_b^{(3)}(\mu_c) + \delta m_c^{(3)}(\mu_c) + \delta m_{(bc)}^{(3)}(\mu_c) + \mathcal{O}\bigg(\mu_c e^{-\frac{2\pi}{\beta_0 \alpha(\mu_c)}(1 + \ln(2))}\bigg).$$
(59)

The $\mathcal{O}(\mu_c e^{\frac{2\pi}{\beta_0 a(\mu_c)}(1+\ln(2))})$ term stands for subleading corrections in the hyperasymptotic expansions, which are not known.

Let us now discuss in more detail the dependence on the bottom and charm quark, in particular the effects associated to the fact that they have masses much bigger than Λ_{OCD} (for the analysis we take $\bar{m}_b = 4.186 \text{ GeV}$ and $\bar{m}_c =$ 1.223 GeV [36] but the sensitivity to the specific values we use is very tiny). As already discussed in [19], the natural scale of a *n*-loop integral is not \bar{m}_t but $\bar{m}_t e^{-n}$. For the case of the bottom versus charm quark it was observed in [13]⁶ that the charm quark effectively decouples at order α^2/α^3 for the case of the charm quark effects in the bottom pole mass- $\overline{\text{MS}}$ mass relation. If we lower the mass of the top we can also observe at which scales it is more convenient to decouple the bottom and charm quark in the top pole mass- $\overline{\text{MS}}$ mass relation. This can be illustrated in Fig. 13, where we plot the corrections associated to the bottom and charm with and without decoupling in terms of the fictitious top mass (assuming a single heavy quark). Obviously for very large top masses it is not convenient to do the decoupling. Nevertheless, as we decrease the mass



FIG. 13. Upper panel: Plot of the correction to the PV mass of a top mass with varying \bar{m}_t mass due to a heavy quark with $\overline{\text{MS}}$ mass equal to 4.185 GeV (bottom) with and without decoupling (assuming a single heavy quark). Lower panel: As in the upper panel with a heavy quark with $\overline{\text{MS}}$ equal to 1.223 GeV (charm). We use Eqs. (B7) and (B10).

 $^{^{6}}$ In that reference MeV should read GeV instead from Eq. (8) to Eq. (12).

uncertainty associated to the $O(\alpha^4)$ corrections is negligible. Also the correction associated to the bottom and charm quark masses to Eq. (58) is, comparatively to the total running, very small. From this analysis we will take as central values $\mu_b = 20$ GeV and $\mu_c = 5$ GeV. For these values we obtain

$$\begin{split} \int_{\mu_{b}}^{\bar{m}_{t}} d\bar{m} \frac{d}{d\bar{m}} (\delta m_{b}^{(5)}(\bar{m}) + \delta m_{c}^{(5)}(\bar{m}) + \delta m_{(bc)}^{(5)}(\bar{m})) \\ &+ \int_{\mu_{c}}^{\mu_{b}} d\bar{m} \frac{d}{d\bar{m}} (\delta m_{b}^{(4)}(\bar{m}) + \delta m_{c}^{(4)}(\bar{m}) + \delta m_{(bc)}^{(4)}(\bar{m})) \\ &+ \delta m_{b}^{(3)}(\mu_{c}) + \delta m_{c}^{(3)}(\mu_{c}) + \delta m_{(bc)}^{(3)}(\mu_{c}) \\ &= -2.5|_{\mathcal{O}(a^{2})} + 0.8|_{\mathcal{O}(a^{3})} = -1.7 \text{ MeV}. \end{split}$$
(60)

The specific value depends on μ_b and μ_c but the good convergence and smallness of this correction holds true for other values of μ_b and μ_c . The implementation of the decoupling of the bottom and charm in [50] produces a much larger correction. An even larger effect is observed in the implementation performed in [45], where the perturbative expansion is always performed at the scale of the top mass (using renormalon based estimates for the higher order coefficients), decoupling the bottom, and later the charm, depending on the order of perturbation theory. Therefore, we take our numbers as optima, and the error negligible compared with other uncertainties.

We next explore the convergence pattern of the perturbative expansion. We first consider the perturbative expansion associated to \mathcal{F} . We find

$$\int_{\mu_b}^{\bar{m}_t} d\bar{m} \mathcal{F}(\bar{m}, 5) + \int_{\mu_c}^{\mu_b} d\bar{m} \mathcal{F}(\bar{m}, 4)$$

= 8445 + 837 + 53 - 43 = 9291(22) MeV. (61)

We observe a convergent pattern. For the last two terms the convergence deteriorates. On the other hand the perturbative expansion becomes sign alternating. This may indicate sensitivity to the u = -1 renormalon. We discuss this further in the next section. For sign-alternating asymptotic perturbative expansion the left-over is $\sim -1/2\times$ (the last computed term) (see [5]).⁷ Therefore, we take it as the error of the truncation of the perturbative expansion, which is the error we quote in Eq. (61). We also explore the dependence of Eq. (58) on μ_b and μ_c . The dependence is very small, as we can see in Fig. 14. For μ_c the variation is negligible, and for μ_b one gets variations of ~5 MeV for a central value of μ_b or around 20 GeV. Therefore, we will neglect it for the total error budget.

The other source of error is associated to the approximate determination of Eq. (59) [except for the δm_q terms, which



FIG. 14. Plots of Eq. (58) in terms of μ_b (upper panel) and μ_c (lower panel) truncating the perturbative expansion of $\mathcal{F}(\bar{m}, n_f)$ at different orders in α in Eq. (61). In the upper figure we set $\mu_c = 5$ GeV. In the lower figure we set $\mu_b = 20$ GeV.

have already been taken into account in Eq. (60)]. The error analysis is equal to the one in Eq. (54) adapted by changing $\bar{m}_b = 4.186 \text{ GeV} \rightarrow \mu_c = 5 \text{ GeV}$ [the error associated to α is only computed for the full Eq. (58)]

$$(m_P(\mu_c) + \mu_c \Omega_m)|_{\mu_c = 5 \text{ GeV}} = 5744(\mu)^{+7}_{-15}(Z_m)^{+9}_{-9} \text{ MeV}.$$
(62)

Finally, we also include the error associated to α . Combining all errors we obtain

$$m_{t,PV}(163 \text{ MeV}) = 173033(\text{h.o.})^{+22}_{-22}(\mu)^{+7}_{-15}(Z_m)^{+9}_{-9}(\alpha)^{+119}_{-123} \text{ MeV}.$$
(63)

By far the largest uncertainty is associated to α . For the purely theoretical error budget, the error is associated to higher order corrections in perturbation theory. They show up in different ways. One is the approximate knowledge of Z_m , which shows up in Ω_m . The other is the error in μ , which is a measure of the $\mathcal{O}(e^{\frac{2\pi}{\beta_0 a_X(\mu)}(1+\ln(2))})$ corrections to Eq. (59). h.o. stands for the error associated to higher order terms in perturbation theory of Eq. (61). All these errors

⁷We emphasize that these arguments do not apply to IR renormalons (and in particular to the u = 1/2 renormalon).

TABLE V. The coefficients f_n of $\mathcal{F}(\bar{m}, n_f)$. Note that $f_4(n_f = 0)$ has a 9% error from the determination in [9]. The $n_f = 10^{20}$ case is used as a test for comparison with the large β_0 . The last three (four) rows are the coefficients f_n in the large β_0 approximation.

$\mathcal{F}(\bar{m}, n_f)$	f_1	f_2	f_3	f_4	f_5
$\overline{n_f} = 0$	4/3	6.11	25.52	18.46	
$n_f = 3$	4/3	4.32	12.76	-63.37	
$n_f = 6$	4/3	2.53	-0.74	-105.70	
(Large β_0 /exact) $n_f = 10^{20}$	4/3	-5.97×10^{19}	-4.15×10^{38}	-2.54×10^{58}	-5.09×10^{77}
(Large β_0) $n_f = 0$	4/3	9.85	-11.31	114.33	-377.22
(Large β_0) $n_f = 3$	4/3	8.06	-7.57	62.62	-169.04
(Large β_0) $n_f = 6$	4/3	6.27	-4.58	29.46	-61.86

would profit from higher order perturbative computations. We have also explored other sources of uncertainty, and find them to be comparatively very small: the error (and the effect) associated to the finite mass of the bottom and charm quark is found to be very small, and similarly for variations in the values of μ_b and μ_c .

It is also useful to make the error estimate of the ratio of the PV and $\overline{\text{MS}}$ top mass. We obtain ($\bar{m}_t = 163 \text{ GeV}$)

$$\left[\frac{m_{t,\rm PV}}{\bar{m}_t} - 1\right] \times 10^5 = 6155(\rm h.o.)^{+13}_{-13}(\mu)^{+4}_{-9}(Z_m)^{+6}_{-6}(\alpha)^{+73}_{-75}.$$
(64)

Note that there is no ambiguity error associated to this number. Except for α all errors are associated to the lack of knowledge of higher order terms of the perturbative expansion. In comparison with [45] we find that our result is less sensitive to Z_m and to its associated error.

C. |u| = 1 renormalons

The perturbative expansion of $\mathcal{F}(\bar{m}, n_f)$ is free of the u = 1/2 renormalon. Therefore, it is the ideal object on which to study the subleading renormalons of the pole mass. In principle, these are located at u = 1 and u = -1. The existence of an infrared renormalon at u = 1 has been a matter of debate [15]. The existence of an ultraviolet renormalon at u = -1 can be established in the large β_0 approximation [19,20] but not beyond. With respect to this discussion some interesting observations can be drawn out of our analysis. The coefficients f_n show an interesting dependence in n_f (with changes of sign of different powers of n_f). In Table V we give the numbers of f_n for different values of n_f and also in the large β_0 approximation. We observe that for $n_f = 3$ the $\mathcal{O}(\alpha^4)$ flips sign. For $n_f = 6$, the $\mathcal{O}(\alpha^3)$ and $\mathcal{O}(\alpha^4)$ flip sign. The situation is somewhat puzzling. Let us first note that the sign of the coefficients would be interchanged compared with the large β_0 predictions (for $n_f = 3$). This could still be understood from a u = -1 renormalon if Z_{-2}^X flips sign from the large β_0 prediction to real QCD. This would indicate a large dependence of Z_{-2}^X on n_f compared with what has been seen for Z_m^X , where the large β_0 approximation gave the right sign and order of magnitude. For $n_f \to \infty$, the results agree with QED expectations (β_0 becomes negative and the perturbative series is nonsign-alternating). For $n_f = 6$ we observe that the last two terms are negative. One may then wonder if what we are seeing for $n_f = 6$ (and maybe also for $n_f = 3$) is that the u = -1 renormalon becomes effectively infrared. Obviously, we need higher order coefficients f_n to clarify this issue.⁸

It is usual lore that infrared renormalons dominate over ultraviolet ones (this is somewhat based on large β_0 analyses where ultraviolet renormalons are typically suppressed by the factor $\sim e^{d\frac{e_X}{2}}$ whereas infrared renormalons are enhanced by the factor $\sim e^{-d\frac{e_X}{2}}$). If we take this seriously, and also the numbers we obtain for f_n as an indication of the existence of the u = -1 renormalon, this may indicate that the u = 1renormalon is indeed zero. In this respect, it is worth mentioning the analysis of [38] where the NP correction associated to the u = 1 renormalon was found to be zero within errors. This is consistent with this discussion.

On the theoretical side it is also interesting to see where the u = 1 renormalon would show up in a perturbative computation of the heavy quarkonium mass. For the purposes of this discussion, the heavy quarkonium mass would read

$$M_{nl} = 2m_Q + \left\langle \frac{\mathbf{p}^2}{m_Q} \right\rangle_{nl} + \left\langle V_0 \right\rangle_{nl} + \left\langle \frac{V_1}{m_Q} \right\rangle_{nl} + \mathcal{O}\left(\frac{1}{m_Q^2}\right),$$
(65)

where V_0 is the static potential, and V_1 is the $1/m_Q$ potential. OPE analyses in the static limit show that V_0 does not have renormalon at u = 1. The virial theorem: $\langle \frac{\mathbf{p}^2}{m_Q} \rangle_{nl} = \langle rV'_0 \rangle_{nl}$, also guaranties that the kinetic term does not have such u = 1 renormalon. Therefore, any possible u = 1 infrared renormalon of the pole mass should cancel with the analogous infrared renormalon of the V_1/m_Q

⁸The coefficients of the perturbative expansion of the pole mass itself are also a polynomial in powers of n_f . The sign dependence of the different powers of n_f has been studied in [55,56].

potential. The fact that the latter can be written in a closed way in terms of Wilson loops [57] may open a venue on which to study this issue in further detail. This is postponed to future work.

VI. CONCLUSIONS

In this paper we have constructed hyperasymptotic expansions for the heavy quark pole mass (and for associated quantities) regulated using the PV prescription along the lines of [1]. We generalize the discussion of that reference by including possible ultraviolet renormalons. Such organization of the computation allows us to have a parametric control of the error committed when truncating the hyperasymptotic expansion.

In Sec. III the hyperasymptotic expansion of the pole mass of a heavy quark in the large β_0 is computed. We use it as a toy-model observable to test our methods. It works as expected. We can see the u = 1/2 infrared renormalon and the u = -1 ultraviolet renormalon. The next infrared renormalon is located at u = 3/2. Compared with the static potential case studied in [1] in the large β_0 approximation, infrared renormalons are located at the same points in the Borel plane. On the other hand, the pole mass has ultraviolet renormalons, whereas the static potential does not. In practice the main difference comes from the relevance of the u = -1 renormalon. In general, because of the u = -1 renormalon, it is necessary to stop the second perturbative expansion [see Eq. (19)] at $N \sim 2 \times \frac{2\pi}{\beta_0 \alpha^2}$ otherwise the perturbative series would start to diverge, as we can observe in Fig. 7 in the \overline{MS} scheme. Nevertheless, the importance of this renormalon heavily depends on the factorization scale μ one uses. If one takes μ high enough, one could indeed do perturbation theory until $N \sim 3 \times \frac{2\pi}{\beta_0 \alpha}$, where the u = 3/2 renormalon shows up. We can see the irrelevance of the u = -1 renormalon in the lattice scheme, which is equivalent to the \overline{MS} scheme with a much larger μ , in Fig. 7. One should keep in mind, though, that one needs perturbation theory to a much higher order in the lattice scheme to reach the same precision than in the $\overline{\text{MS}}$ scheme. We expect this qualitative behavior of ultraviolet renormalons to also hold true beyond the large β_0 approximation.

We next move to real QCD. We have performed determinations of $\bar{\Lambda}_{PV}$ using quenched lattice QCD. For these observables perturbative expansions to high orders are available [21–23]. This allows us to test the method and go beyond the superasymptotic and the leading term in the hyperasymptotic approximation. We observe $\mathcal{O}(a\Lambda_{QCD}^2)$ corrections for the *B* meson mass in the static approximation, but not for an analogous observable from the static potential. Nevertheless, we do not have enough precision to quantitatively study these effects. The limiting factor is the error of the normalization of the leading renormalon, and, related, the lack of knowledge of the higher order beta

function coefficients. The latter affects the O(1/n) corrections to the asymptotic formula of the perturbative series coefficients. These effects are sizable in the lattice scheme. On the other hand they are quite small in the $\overline{\text{MS}}$ scheme. On top of that the higher order coefficients of the perturbative expansion of δm_{latt} are not known with enough precision to disentangle the subleading renormalon (their error is strongly correlated with the error of Z_m). All these considerations forbid quantitative analyses beyond the leading term in the hyperasymptotic approximation. Further investigations are needed to improve on these issues, particularly on the error of Z_m , which also affects the discussion below.

We also determine $\bar{\Lambda}_{PV}$ from the physical *B* meson mass assuming that the \overline{MS} heavy quark mass is known. The result can be found in Eq. (54). In this analysis, we determine the error associated to the incomplete knowledge of the perturbative expansion in determinations of the heavy quark mass. We translate this result to the case of the top mass, which we study in detail in Sec. V. In this section the issue of the uncertainty of the (top) pole mass is critically reexamined. In particular, the bottom and charm quark finite mass effects are carefully incorporated. In our implementation we find these to be very small. We find the present uncertainty in the relation between \bar{m}_t and m_{PV} to be (for $\bar{m}_t = 163 \text{ GeV}$)

$$m_{t,\text{PV}}(163 \text{ MeV}) = 173033(\text{th})^{+25}_{-28}(\alpha)^{+119}_{-123} \text{ MeV},$$
 (66)

$$\left[\frac{m_{t,\rm PV}}{\bar{m}_t} - 1\right] \times 10^5 = 6155(\rm th)^{+15}_{-17}(\alpha)^{+73}_{-75},\tag{67}$$

where we have combined the theoretical errors quoted in Eqs. (63) and (64) in quadrature. There is no ambiguity associated to the renormalon in this number. The precision is systematically improvable the more terms of the perturbative expansion get to be known in the future. Interestingly enough, it seems we have found some evidence for the existence of the next renormalon at u = -1 but not of a possible renormalon at u = 1. We believe this makes very timely a quantitative determination of the renormalization group structure of the u = -1 renormalon, which to our knowledge is lacking. We leave this for future work.

ACKNOWLEDGMENTS

We thank M. Steinhauser for comments on the manuscript. C. A. thanks the IFAE group at Universitat Autònoma de Barcelona for warm hospitality during part of this work. This work was supported in part by the Spanish FPA2017-86989-P and SEV-2016-0588 grants from the ministerio de Ciencia, Innovación y Universidades, and the 2017SGR1069 grant from the Generalitat de Catalunya; and by the Chilean FONDECYT Postdoctoral Grant No. 3170116, and by FONDECYT Regular Grant No. 1180344.

APPENDIX A: EVALUATION OF Ω_d

We here briefly sketch how we compute the integrals that appear in Ω_d . For d < 0 we use the recursion formulas developed in [5] (see for instance Eq. (46) of Chapter XXI). For d > 0, we can use also such formulas (see for instance Eq. (47) of Chapter XXI). In this case, we can also alternatively perform the integration in the following way. For simplicity, we take the case b = 0and d = 1, as the method is similar for the more general case.

$$I = \int_{0,\text{PV}}^{\infty} du e^{-\frac{4\pi}{\beta_0 \alpha} u} \frac{(2u)^{N+1}}{1-2u} = -\frac{1}{2} e^{-\frac{2\pi}{\beta_0 \alpha}} \int_{-\frac{1}{2},\text{PV}}^{\infty} \frac{dy}{y} e^{-\frac{4\pi}{\beta_0 \alpha} y} e^{(\frac{2\pi}{\beta_0 \alpha} (1-c\alpha)+1) \ln(1+2y)}, \quad (A1)$$

where in the second equality we set $N = N_P$ according to Eq. (3). We also do the change of variables (where $K = c - \frac{\beta_0}{2\pi}$)

$$y = \frac{x}{2} \frac{1}{\sqrt{1 - K\alpha}} \sqrt{\frac{\beta_0 \alpha}{\pi}}.$$
 (A2)

We can then expand the exponent in powers of α and *x*:

$$I = -\frac{1}{2} e^{-\frac{2\pi}{\beta_0 a}} \int_{-\sqrt{\frac{\pi}{\beta_0 a}(1-Ka)}, \text{PV}}^{\infty} \frac{dx}{x} \times e^{-x^2} e^{-\frac{4\pi}{\beta_0 a}K_2^x} \sqrt{\frac{\beta_0 a}{\pi} \frac{1}{\sqrt{1-Ka}} + \frac{2}{3}x^3} \sqrt{\frac{\beta_0 a}{\pi} \frac{1}{\sqrt{1-Ka}} + \cdots}.$$
 (A3)

The inferior limit of the integral is then extended to $-\infty$. We then have

$$I \simeq -\frac{1}{2} e^{-\frac{2\pi}{\beta_0 a}} \int_{-\infty, \text{PV}}^{\infty} \frac{dx}{x} e^{-x^2} e^{-\frac{4\pi}{\beta_0 a} K_2^x} \sqrt{\frac{\beta_0 a}{\pi} \frac{1}{\sqrt{1-\kappa a}} + \frac{2}{3} x^3} \sqrt{\frac{\beta_0 a}{\pi} \frac{1}{\sqrt{1-\kappa a}}} + \cdots$$
$$\simeq e^{-\frac{2\pi}{\beta_0 a}} \frac{1}{6} \sqrt{\beta_0 a} \left(-1 + \frac{6\pi}{\beta_0} K \right) + \cdots .$$
(A4)

Irrespectively of considering d > 0 or d > 0, it is not clear to us what is the asymptotic structure of this expansion. This is something that we are investigating. In any case, at present, we have not seen evidence of asymptotic behavior of this expansion for all cases we have considered. This does not preclude however that if we go to higher orders we will find an asymptotic behavior for this perturbative expansion.

APPENDIX B: BOTTOM AND CHARM FINITE MASS CONTRIBUTIONS TO $m_{t,PV}$

We define

$$\begin{split} \delta m_q^{(1)} &\equiv \frac{\bar{m}}{3} \left(\left(1 - \frac{\bar{m}_q}{\bar{m}} \right) \left(1 - \frac{\bar{m}_q^3}{\bar{m}^3} \right) \left(\text{Li}_2 \left(\frac{\bar{m}_q}{\bar{m}} \right) - \frac{1}{2} \ln^2 \left(\frac{\bar{m}_q}{\bar{m}} \right) + \ln \left(1 - \frac{\bar{m}_q}{\bar{m}} \right) \ln \left(\frac{\bar{m}_q}{\bar{m}} \right) - \frac{\pi^2}{3} \right) \\ &+ \left(1 + \frac{\bar{m}_q}{\bar{m}} \right) \left(1 + \frac{\bar{m}_q^3}{\bar{m}^3} \right) \left(\text{Li}_2 \left(- \frac{\bar{m}_q}{\bar{m}} \right) - \frac{1}{2} \ln^2 \left(\frac{\bar{m}_q}{\bar{m}} \right) + \ln \left(1 + \frac{\bar{m}_q}{\bar{m}} \right) \ln \left(\frac{\bar{m}_q}{\bar{m}} \right) + \frac{\pi^2}{6} \right) \\ &- \frac{\bar{m}_q^2}{\bar{m}^2} \left(\ln \left(\frac{\bar{m}_q}{\bar{m}} \right) + \frac{3}{2} \right) + \ln^2 \left(\frac{\bar{m}_q}{\bar{m}} \right) + \frac{\pi^2}{6} \end{split}$$
(B1)

(note that this coefficient is n_f -independent),

$$\delta m_q^{(2,n_f)} = \frac{\bar{m}}{64} \left[h\left(\frac{\bar{m}_q}{\bar{m}}\right) + w\left(1,\frac{\bar{m}_q}{\bar{m}}\right) + n_f p\left(\frac{\bar{m}_q}{\bar{m}}\right) \right],\tag{B2}$$

where $n_f = 5$ for q = b and $n_f = 4$ for q = c, and we use the representation for the functions h(x), w(x, y) and p(x) given in Ref. [50], and

$$\delta m_{bc}^{(2)} = \frac{\bar{m}}{64} w \left(\frac{\bar{m}_b}{\bar{m}}, \frac{\bar{m}_c}{\bar{m}} \right). \tag{B3}$$

We then have

$$\delta m_{b/c}^{(5)} = \delta m_{b/c}^{(1)} \frac{\alpha_{(5)}^2(\bar{m})}{\pi^2} + \delta m_{b/c}^{(2,5/4)} \frac{\alpha_{(5)}^3(\bar{m})}{\pi^3}$$
(B4)

$$\delta m_{bc}^{(5)} = \delta m_{bc}^{(2)} \frac{\alpha_{(5)}^3(\bar{m})}{\pi^3} \tag{B5}$$

$$\delta m_b^{(4)} = \left[\delta m_b^{(1)} + \delta m_{b,dec}^{(1)}\right] \frac{\alpha_{(4)}^2(\bar{m})}{\pi^2} + \left[\delta m_b^{(2,5)} + \delta m_{b,dec}^{(2)}\right] \frac{\alpha_{(4)}^3(\bar{m})}{\pi^3} \tag{B6}$$

$$\delta m_c^{(4)} = \delta m_c^{(1)} \frac{\alpha_{(4)}^2(\bar{m})}{\pi^2} + \delta m_c^{(2,4)} \frac{\alpha_{(4)}^3(\bar{m})}{\pi^3} \tag{B7}$$

$$\delta m_{bc}^{(4)} = \left[\delta m_{bc}^{(2)} + \delta m_{bc,dec}^{(2)}\right] \frac{\alpha_{(4)}^3(\bar{m})}{\pi^3} \tag{B8}$$

$$\delta m_b^{(3)} = \left[\delta m_b^{(1)} + \delta m_{b,dec}^{(1)}\right] \frac{\alpha_{(3)}^2(\bar{m})}{\pi^2} + \left[\delta m_b^{(2,5)} + \delta m_{b,dec}^{(2)}\right] \frac{\alpha_{(3)}^3(\bar{m})}{\pi^3} \tag{B9}$$

$$\delta m_c^{(3)} = \left[\delta m_c^{(1)} + \delta m_{c,dec}^{(1)}\right] \frac{\alpha_{(3)}^2(\bar{m})}{\pi^2} + \left[\delta m_c^{(2,4)} + \delta m_{c,dec}^{(2)}\right] \frac{\alpha_{(3)}^3(\bar{m})}{\pi^3} \tag{B10}$$

$$\delta m_{bc}^{(3)} = \left[\delta m_{bc}^{(2)} + \delta m_{bc,dec}^{(2)} + \delta m_{cb,dec}^{(2)} \right] \frac{\alpha_{(3)}^3(\bar{m})}{\pi^3},\tag{B11}$$

where $\delta m^{(i)}_{(q,{\rm dec})}$ are generated by the decoupling and read

$$\delta m_{(q,\text{dec})}^{(1)} = -\frac{2}{9}\bar{m} \left(\frac{71}{32} + \ln\left(\frac{\bar{m}_q^2}{\bar{m}^2}\right) + \frac{\pi^2}{4} \right),\tag{B12}$$

$$\delta m_{(q,dec)}^{(2,n_f)} = \bar{m} \left\{ \left[\frac{2353}{11664} + \frac{7}{27} \zeta(3) + \frac{13\pi^2}{162} - \left(\frac{\pi^2}{54} + \frac{71}{432}\right) \ln\left(\frac{\bar{m}^2}{\bar{m}_q^2}\right) \right] n_f + \frac{8\mathrm{Li}_4(\frac{1}{2})}{27} - \frac{751}{216} \zeta(3) + \frac{61\pi^4}{1944} - \frac{113\pi^2}{72} - \frac{29869}{2916} + \frac{\ln^4(2)}{81} + \frac{2}{81} \pi^2 \ln^2(2) - \frac{11}{81} \pi^2 \ln(2) + \left(\frac{1225}{288} - \frac{1}{18} \zeta(3) + \frac{\pi^2}{9} + \frac{1}{27} \pi^2 \ln(2)\right) \ln\left(\frac{\bar{m}^2}{\bar{m}_q^2}\right) + \frac{1}{27} \ln^2\left(\frac{\bar{m}^2}{\bar{m}_q^2}\right) \right\} + \frac{1}{3} \ln\left(\frac{\bar{m}^2}{\bar{m}_q^2}\right) \delta m_q^{(1)}.$$
(B13)

Note that $\delta m_{(b,dec)}^{(2)} = \delta m_{(b,dec)}^{(2,5)}$ and $\delta m_{(c,dec)}^{(2)} = \delta m_{(b,dec)}^{(2,4)}$. This last expression indeed corresponds to Eq. (17) of [13] changing \bar{m}_b by \bar{m} .

Finally, we also have

$$\delta m_{bc,dec}^{(2)} = \frac{1}{3} \ln \left(\frac{\bar{m}^2}{\bar{m}_b^2} \right) \delta m_c^{(1)}, \tag{B14}$$

$$\delta m_{cb,dec}^{(2)} = \frac{1}{3} \ln \left(\frac{\bar{m}^2}{\bar{m}_c^2} \right) [\delta m_b^{(1)} + \delta m_{b,dec}^{(1)}]. \tag{B15}$$

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