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The strong coupling from the revised ALEPH data for hadronic τ decays

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ABSTRACT

We apply an analysis method previously developed for the extraction of the strong coupling from the OPAL data to the recently revised ALEPH data for non-strange hadronic τ decays. Our analysis yields the values $\alpha_s(m_\tau^2) = 0.296 \pm 0.010$ using fixed-order perturbation theory, and $\alpha_s(m_\tau^2) = 0.310 \pm 0.014$ using contour-improved perturbation theory. Averaging these values with our previously obtained values from the OPAL data, we find $\alpha_s(m_\tau^2) = 0.303 \pm 0.009$, respectively, $\alpha_s(m_\tau^2) = 0.319 \pm 0.012$. We present a critique of the analysis method employed previously, for example in analyses by the ALEPH and OPAL collaborations, and compare it with our own approach. Our conclusion is that non-perturbative effects limit the accuracy with which the strong coupling, an inherently perturbative quantity, can be extracted at energies as low as the τ mass. Our results further indicate that systematic errors on the determination of the strong coupling from analyses of hadronic τ -decay data have been underestimated in much of the existing literature.

I. INTRODUCTION

Recently, Ref. [1], for the ALEPH collaboration, updated and revised previous ALEPH results for the non-strange vector (V) and axial vector (A) spectral distributions obtained from measurements of hadronic τ decays. In particular, Ref. [1] corrects a problem in the publicly posted 2005 and 2008 versions of the correlations between different energy bins uncovered in Ref. [2].¹ The corrected data supersede those originally published by the ALEPH collaboration [3, 4].

One of the hadronic quantities of interest that can be extracted from these data is the strong coupling $\alpha_s(m_\tau^2)$ at the τ mass, through the use of Finite-Energy Sum Rules (FESRs) [5], as advocated long ago [6, 7]. Both the ALEPH and OPAL [8] collaborations have done so by applying an analysis strategy, developed in Refs. [7, 9], in which small, but non-negligible non-perturbative effects were estimated using a truncated form of the operator product expansion (OPE). A feature of the particular truncation scheme employed is that it assumes that, in addition to contributions which violate quark-hadron duality, also OPE contributions of dimension $D > 8$ unsuppressed by non-leading powers of α_s can be safely neglected. Given the goal of extracting $\alpha_s(m_\tau^2)$ with the best possible accuracy, these features of what we will refer to as the “standard analysis” have been questioned, starting with the work of Refs. [10, 11]. In these works, it was argued that both the OPE truncation to terms with $D \leq 8$ and the neglect of violations of quark-hadron duality lead to additional numerically non-negligible systematic uncertainties not included in the errors obtained on $\alpha_s(m_\tau^2)$ and the OPE condensates from the standard-analysis approach. In order to remedy this situation, in Refs. [12, 13], we developed a new analysis strategy designed to take both OPE and duality-violating (DV) non-perturbative effects consistently into account. This strategy was then successfully applied to the OPAL data [12, 13]. In the present article, we apply this analysis strategy to the corrected ALEPH data, and compare our results to those obtained from the OPAL data in Ref. [13] as well as to those of the recent re-analysis presented in Ref. [1].

The calculation of the order- α_s^4 term [14] in the perturbative expansion of the Adler function in 2008 led to a renewed interest in the determination of the strong coupling from hadronic τ decays, with many attempts to use this new information on the theory side of the relevant FESRs in order to sharpen the extraction of $\alpha_s(m_\tau^2)$ from the data [1, 4, 10, 12–19]. Since the perturbative series converges rather slowly, different partial resummation schemes have been considered, leading to variations in the obtained results. The majority of these post-2007 updates (Refs. [1, 4, 14–19]), however, were carried out assuming that the standard-analysis treatment of non-perturbative effects was essentially correct, with none of the references in this subset, with the exception of Refs. [1, 4], redoing the analysis starting from the underlying experimental data (the emphasis, instead, being on the merits of different resummation schemes for the perturbative expansion). Reference [10], which did revisit the determination of the higher- D OPE contributions, and performed a more careful treatment of these contributions, did not, however, include DV contributions in its analysis framework. While its results were tested for self-consistency, the absence of a representation of DV effects meant no estimate of the residual systematic error associated with their neglect was possible. The only articles to incorporate both the improved treatment of higher- D OPE contributions and an implementation of a physically motivated representation of DV effects

¹ The updated and corrected data can be found at <http://aleph.web.lal.in2p3.fr/tau/specfun13.html>.

were those of Refs. [12, 13], which, due to the problem with the then-existing ALEPH covariance matrices, were restricted to analyzing OPAL data. Our goal in this article is to reconsider the treatment of non-perturbative effects employing the newly released ALEPH data, which have significantly smaller errors than the OPAL data. We will present results for the two most popular resummation schemes for the perturbative (*i.e.*, $D = 0$ OPE) series: fixed-order perturbation theory (FOPT) and contour-improved perturbation theory (CIPT) [20], without trying to resolve the discrepancies that arise between them (for an overview of the two methods, see Ref. [21]).

This article is organized as follows. In Sec. II we give a brief overview of the necessary theory, referring to Ref. [12] for more details. In Sec. III we discuss the new ALEPH data set, and check explicitly that the current publicly posted version of the correlation matrices pass the test that led to the identification of the problem with the previous version [2]. We also show the comparison of the experimental ALEPH and OPAL non-strange spectral functions. Section IV summarizes our fitting strategy, developed in Refs. [12, 13]. Sections V and VI present the details of the fits, and the results we obtain from them for $\alpha_s(m_\tau^2)$ and dimension 6 and 8 OPE coefficients in the V and A channels. We explore the χ^2 landscape using the Markov-chain Monte Carlo code `hrothgar` [22], which in the case of the OPAL data proved useful in uncovering potential ambiguities. Also included is an estimate for the total non-perturbative contribution to the ratio of non-strange hadronic and electronic τ branching fractions. In Sec. VI we check how well the two Weinberg sum rules [23] and the sum rule for the electro-magnetic pion mass difference [24] are satisfied by our results. Finally, in Sec. VII, we present a critical discussion of the standard analysis employed in Refs. [1, 3, 4, 8], focusing on the most recent of these, described in Ref. [1]. We demonstrate explicitly the inconsistency of this analysis with regard to the treatment of non-perturbative effects, and conclude that, while the standard analysis approach was a reasonable one to attempt in the past, it must be abandoned in current or future determinations of $\alpha_s(m_\tau^2)$ from hadronic τ decay data. In our concluding section, Sec. VIII, we compare our approach with the standard-analysis method, highlighting and juxtaposing the assumptions underlying each, and summarize our results.

II. THEORY OVERVIEW

The sum-rule analysis starts from the correlation functions

$$\begin{aligned}\Pi_{\mu\nu}(q) &= i \int d^4x e^{iqx} \langle 0 | T \{ J_\mu(x) J_\nu^\dagger(0) \} | 0 \rangle \\ &= (q_\mu q_\nu - q^2 g_{\mu\nu}) \Pi^{(1)}(q^2) + q_\mu q_\nu \Pi^{(0)}(q^2) \\ &= (q_\mu q_\nu - q^2 g_{\mu\nu}) \Pi^{(1+0)}(q^2) + q^2 g_{\mu\nu} \Pi^{(0)}(q^2) ,\end{aligned}\tag{2.1}$$

where J_μ stands for the non-strange V or A current, $\bar{u}\gamma_\mu d$ or $\bar{u}\gamma_\mu\gamma_5 d$, while the superscripts (0) and (1) label spin. The decomposition in the third line employs the combinations $\Pi^{(1+0)}(q^2)$ and $q^2\Pi^{(0)}(q^2)$, which are free of kinematic singularities. Defining $s = q^2 = -Q^2$ and the spectral function

$$\rho^{(1+0)}(s) = \frac{1}{\pi} \text{Im} \Pi^{(1+0)}(s) ,\tag{2.2}$$

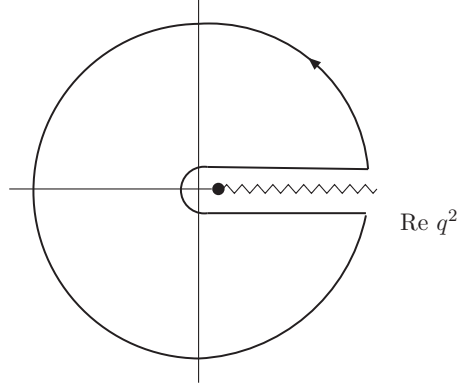


FIG. 1: Analytic structure of $\Pi^{(1+0)}(q^2)$ in the complex $s = q^2$ plane. There is a cut on the positive real axis starting at $s = q^2 = 4m_\pi^2$ (a pole at $s = q^2 = m_\pi^2$ and a cut starting at $s = 9m_\pi^2$) for the V (A) case. The solid curve shows the contour used in Eq. (2.3).

Cauchy's theorem and the analytical properties of $\Pi^{(1+0)}(s)$, applied to the contour in Fig. 1, imply the FESR

$$I_{V/A}^{(w)}(s_0) \equiv \frac{1}{s_0} \int_0^{s_0} ds w(s) \rho_{V/A}^{(1+0)}(s) = -\frac{1}{2\pi i s_0} \oint_{|s|=s_0} ds w(s) \Pi_{V/A}^{(1+0)}(s), \quad (2.3)$$

valid for any $s_0 > 0$ and any weight $w(s)$ analytic inside and on the contour [5].

The flavor ud V and A spectral functions can be experimentally determined from the differential versions of the ratios,

$$R_{V/A;ud} = \frac{\Gamma[\tau \rightarrow (\text{hadrons})_{V/A;ud} \nu_\tau(\gamma)]}{\Gamma[\tau \rightarrow e \bar{\nu}_e \nu_\tau(\gamma)]}, \quad (2.4)$$

of the width for hadronic decays induced by the relevant current to that for the electron mode. Explicitly [25],

$$\frac{dR_{V/A;ud}(s)}{ds} = 12\pi^2 |V_{ud}|^2 S_{EW} \frac{1}{m_\tau^2} \left[w_T(s; m_\tau^2) \rho_{V/A;ud}^{(1+0)}(s) - w_L(s; m_\tau^2) \rho_{V/A;ud}^{(0)}(s) \right], \quad (2.5)$$

where S_{EW} is a short-distance electroweak correction and $w_T(s; s_0) = (1 - s/s_0)^2(1 + 2s/s_0)$, $w_L(s; s_0) = 2(s/s_0)(1 - s/s_0)^2$. Apart from the pion-pole contribution, which is not chirally suppressed, $\rho_{V/A;ud}^{(0)}(s) = O[(m_d \mp m_u)^2]$, and the continuum part of $\rho_{V/A}^{(0)}(s)$ is thus numerically negligible. As a result, the spectral functions $\rho_{V/A;ud}^{(1+0)}(s)$ can be determined directly from $dR_{V/A;ud}(s)/ds$. The FESR (2.3) can thus be studied for arbitrary s_0 and arbitrary analytic weight $w(s)$. From now on, we will denote the experimental version of the spectral integral on the left-hand side of Eq. (2.3) by $I_{V/A;\text{ex}}^{(w)}(s_0)$ (generically, $I_{\text{ex}}^{(w)}(s_0)$) and the theoretical representation of the contour integral on the right-hand side by $I_{V/A;\text{th}}^{(w)}(s_0)$ (generically, $I_{\text{th}}^{(w)}(s_0)$).

For large enough $|s| = s_0$, away from the positive real axis, $\Pi^{(1+0)}(s)$ can be approximated by the OPE

$$\Pi_{\text{OPE}}^{(1+0)}(s) = \sum_{k=0}^{\infty} \frac{C_{2k}(s)}{(-s)^k} , \quad (2.6)$$

with the OPE coefficients C_{2k} logarithmically dependent on s through perturbative corrections. The term with $k = 0$ corresponds to the purely perturbative, mass-independent contributions, which have been calculated to order α_s^4 in Ref. [14], and are the same for the V and A channels. The C_{2k} with $k \geq 1$ are different for the V and A channels, and, for $k > 1$, contain non-perturbative $D = 2k$ condensate contributions. As in Refs. [12, 13], we will neglect purely perturbative quark-mass contributions to C_2 and C_4 , as they are numerically very small for the non-strange FERSs we consider in this article. For the same reason, we will neglect the s -dependence of the coefficients C_{2k} for $k > 1$. For the perturbative contribution, C_0 , we will use the result of Ref. [14] and extract $\alpha_s(m_\tau^2)$ in the $\overline{\text{MS}}$ scheme. Since the coefficient c_{51} of the order- α_s^5 term has not been calculated we will use the estimate $c_{51} = 283$ of Ref. [15] with a 100% uncertainty. We will also employ both FOPT and CIPT resummation schemes in evaluating the truncated perturbative series. For more details on the treatment of the $D > 0$ OPE contributions, we refer the reader to Ref. [12].

Perturbation theory, and in general the OPE, breaks down near the positive real $s = q^2$ axis [26]. We account for this by replacing the right-hand side of Eq. (2.3) by

$$-\frac{1}{2\pi i s_0} \oint_{|s|=s_0} ds w(s) \left(\Pi_{\text{OPE}}^{(1+0)}(s) + \Delta(s) \right) , \quad (2.7)$$

with

$$\Delta(s) \equiv \Pi^{(1+0)}(s) - \Pi_{\text{OPE}}^{(1+0)}(s) , \quad (2.8)$$

where the difference $\Delta(s)$ accounts, by definition, for the quark-hadron duality violating contribution to $\Pi^{(1+0)}(s)$. As shown in Ref. [11], Eq. (2.7) can be rewritten as

$$I_{\text{th}}^{(w)}(s_0) = -\frac{1}{2\pi i s_0} \oint_{|s|=s_0} ds w(s) \Pi_{\text{OPE}}^{(1+0)}(s) - \frac{1}{s_0} \int_{s_0}^{\infty} ds w(s) \frac{1}{\pi} \text{Im} \Delta(s) , \quad (2.9)$$

if $\Delta(s)$ is assumed to decay fast enough as $s \rightarrow \infty$. The imaginary parts $\frac{1}{\pi} \text{Im} \Delta_{V/A}(s)$ can be interpreted as the DV parts, $\rho_{V/A}^{\text{DV}}(s)$, of the V/A spectral functions.

The functional form of $\Delta(s)$ is not known, even for large s , and we thus need to resort to a model in order to account for DVs. Following Refs. [11, 27, 28],² we use a model based on large- N_c and Regge considerations, choosing to parametrize $\rho_{V/A}^{\text{DV}}(s)$ as³

$$\rho_{V/A}^{\text{DV}}(s) = e^{-\delta_{V/A} - \gamma_{V/A} s} \sin(\alpha_{V/A} + \beta_{V/A} s) . \quad (2.10)$$

This introduces, in addition to α_s and the $D \geq 4$ OPE condensates, four new parameters in each channel. As in Refs. [12, 13], we will assume that Eq. (2.10) holds for $s \geq s_{\text{min}}$, with s_{min} to be determined from fits to the data. This, in turn, assumes that we can take s_{min} significantly smaller than m_τ^2 , *i.e.*, that both the OPE and the *ansatz* (2.10) can be used in some interval below m_τ^2 .

² See also Refs. [29, 30].

³ In Ref. [12] we used $\kappa_{V/A} \equiv e^{-\delta_{V/A}}$; in Ref. [13] we switched to $\delta_{V/A}$.

Let us pause at this point to revisit the basic ideas underlying the DV *ansatz* (2.10). Since there exists, as yet, no theory of DVs starting from first principles in QCD, the *ansatz* (2.10) represents simply our best, physically motivated, guess as to an appropriate form of DV contributions to the V and A spectral functions. The damped oscillatory form employed is, however, far from arbitrary. First, it reflects the fact that DVs are expected to produce almost harmonic oscillations around the perturbative continuum, in line with expectations from Regge theory, in which resonances occur with equal squared-mass spacings on the relevant daughter trajectories. Second, the exponential damping factor in the *ansatz* reflects the understanding that the OPE is (at best) an asymptotic, and not a convergent, expansion. It is certainly the case that the OPE representation is more successful for euclidean $Q^2 \sim 2 \text{ GeV}^2$ than for comparable Minkowski scales, $q^2 \sim 2 \text{ GeV}^2$, where DV contributions are clearly visible in the spectral functions. Once DVs are identified as representing the irreducible error present in this asymptotic expansion, it is natural to assume that their contribution should exhibit an exponentially suppressed dependence on $s = q^2$, as in our *ansatz* (2.10). These qualitative expectations are also reflected in the explicit Regge- and large- N_c -motivated model discussed in much more detail in Refs. [11, 27–30]. These plausibility arguments aside, we will use the precise ALEPH data to subject the parametrization (2.10) to non-trivial tests described in detail in Sec. VII.

Several considerations underlie our choice of weight functions $w(s)$. First, we will choose weight functions which are likely to be well-behaved in perturbation theory, based on the findings of Ref. [31]. In particular, we will exclude weight functions with a term linear in s , and require the ones we use to include a constant term (which we will normalize to one). Second, because it is not known at which order the OPE might start to diverge (for the values of s_0 of interest), we wish to avoid terms in Eq. (2.6) with $D > 8$, about which essentially nothing is known. That means that if we do not want to arbitrarily set the coefficients C_D with $D > 8$ equal to zero, our weight functions are restricted to polynomials with degree not larger than three. Combining these constraints, we are left with the form

$$w(s; s_0) = 1 + a(s/s_0)^2 + b(s/s_0)^3 . \quad (2.11)$$

This allows us at most three independent weight functions, and limits the extent to which we can use sufficiently pinched weights, *i.e.*, weights with a (multiple) zero at $s = s_0$, which help to suppress DVs [32, 33]. The upshot is that, if we want to exploit the s_0 dependence of the data (instead of fitting only at $s_0 = m_\tau^2$, as was done in Refs. [1, 3, 4, 8]) and treat the OPE consistently, modeling DVs is unavoidable [12]. We emphasize that the s_0 dependence of fit results provides a crucial test of the validity of FESR fits to the data, as we will see below. As in Refs. [12, 13], we choose to consider the weight functions

$$\begin{aligned} \hat{w}_0(x) &= 1 , \\ \hat{w}_2(x) &= 1 - x^2 , \\ \hat{w}_3(x) &= (1 - x)^2(1 + 2x) = 1 - 3x^2 + 2x^3 = w_T(s; s_0) , \\ x &\equiv s/s_0 . \end{aligned} \quad (2.12)$$

The first choice, \hat{w}_0 , is predicated on the fact that pinching is known to suppress DV contributions and we need at least one weight which is sufficiently sensitive to DV contributions to fix the DV parameters. The remaining two weights \hat{w}_2 and \hat{w}_3 are singly and doubly pinched, respectively. For a more detailed discussion of our choices, we refer to Ref. [12]. An important observation is that these choices for what goes into the parametrization of

$I_{\text{th}}^{(w)}(s_0)$ did remarkably well in the analysis of the OPAL data. It therefore makes sense to see what happens if we apply the same strategy to the ALEPH data.

III. THE ALEPH DATA

In this section, we discuss the revised ALEPH data, which are available from Ref. [34]. First, we perform a minor rescaling, in order to account for more precise values of some “external” quantities (*i.e.*, quantities not directly measured by ALEPH, but used in their analysis of the data); this is discussed in Sec. III A, where we also specify our other inputs. Then, in Sec. III B we apply to the corrected covariance matrices the test of Ref. [2] that led us to uncover the problem with the previously published versions, and verify that the revised covariances pass this test. Finally, we compare the V and A spectral functions obtained from the ALEPH data with those from the OPAL data.

A. Data and normalization

We will use the following input values in our analysis:

$$\begin{aligned} m_\tau &= 1.77682(16) \text{ GeV} , \\ B_e &= 0.17827(40) , \\ V_{ud} &= 0.97425(22) , \\ S_{EW} &= 1.0201(3) , \\ m_\pi &= 139.57018(35) \text{ MeV} , \\ f_\pi &= 92.21(14) \text{ MeV} . \end{aligned} \tag{3.1}$$

Here B_e is the branching fraction for the decay $\tau \rightarrow e\bar{\nu}_e\nu_\tau$ and we have used the result of an HFAG fit of the τ branching fractions which incorporates $\pi_{\mu 2}$ and $K_{\mu 2}$ data and Standard Model expectations based on these data for the π and K branching fractions [35]; f_π is the π decay constant. The value for V_{ud} is from Ref. [36], that for S_{EW} from Ref. [37], and the values for m_τ , m_π and f_π are from the Particle Data Group [38]. Only the error on B_e has a significant effect in our analysis; errors on the other input quantities are too small to affect the final analysis errors in any significant way.

To the best of our knowledge, Ref. [1] uses the values $B_e = 0.17818(32)$ and $S_{EW} = 1.0198$. This value for B_e we infer from the ALEPH values for $R_V = 1.782(9)$ [1] and the corresponding branching fraction $B_V = 0.31747$ [34] specified in the publicly posted V data file (no error quoted). The continuum (pion-less) axial branching fraction $B_{A,\text{cont}} = 0.19369$ with $B_e = 0.17818$ translates into $R_{A,\text{cont}} = 1.08705$. From these values, and the quoted value $R_{ud} = 3.475(11)$ [1], it follows that the ALEPH value for R_π , the pion pole contribution to R_{ud} , is $R_\pi = 0.606$. However, if one employs the very precisely known value of f_π quoted above, obtained from $\pi_{\mu 2}$ decays, together with the quoted values for S_{EW} and V_{ud} , one finds instead the more precisely determined expectation $R_\pi = 0.6101$. Using this latter value as well as the ALEPH value $R_{ud} = 3.475(11)$ leads to $R_V + R_{A,\text{cont}} = 2.865$, instead of the ALEPH value $(B_V + B_{A,\text{cont}})/B_e = (0.31747 + 0.19369)/0.17818 = 2.8688$. We employ the more precise $\pi_{\mu 2}$ expectation for the important A channel pion-pole contribution, and take this difference into account by rescaling the V and continuum A non-strange spectral

functions by the common factor $2.865/2.8688 = 0.9987$, since we have no information on whether this rescaling should affect the V and A channels asymmetrically. Our rescaling is thus imperfect, but it is to be noted that the effect of this rescaling lowers our value for $\alpha_s(m_\tau^2)$ by less than one percent, a much smaller shift than that allowed by the total error, see Sec. VI.

The new ALEPH data use a variable bin width, with the highest bin, number 80, centered at $\text{sbin}(80) = 3.3375 \text{ GeV}^2$, which is above $m_\tau^2 = 3.1571 \text{ GeV}^2$. The next-highest bin, number 79, is centered at $\text{sbin}(79) = 3.0875 \text{ GeV}^2$, with a width $\text{dsbin}(79) = 0.1750 \text{ GeV}^2$, so that also $\text{sbin}(79) + \text{dsbin}(79)/2 > m_\tau^2$. In order to avoid using values of s larger than m_τ^2 , we will modify these values to

$$\begin{aligned}\text{sbin}(79) &= 3.07854 \text{ GeV}^2, \\ \text{dsbin}(79) &= 0.157089 \text{ GeV}^2,\end{aligned}\tag{3.2}$$

so that $\text{sbin}(79) + \text{dsbin}(79)/2 = m_\tau^2$.

Finally, ALEPH provides binned spectral data for $\text{sfm2}(\text{sbin})$, which are related to the spectral functions by

$$\text{sfm2}(\text{sbin}) = 100 \times \frac{12\pi^2 |V_{ud}|^2 S_{EW} B_e}{m_\tau^2} \Delta w^T(\text{sbin}; m_\tau^2) \rho^{(1+0)}(\text{sbin}),\tag{3.3}$$

in which

$$\Delta w^T(\text{sbin}; m_\tau^2) = \int_{\text{sbin}-\text{dsbin}/2}^{\text{sbin}+\text{dsbin}/2} ds w^T(s; m_\tau^2).\tag{3.4}$$

For infinitesimal $\text{dsbin} = ds$ one has $\Delta w^T(s; m_\tau^2) = w^T(s; m_\tau^2) ds$, but for finite bin width we have to make a choice in how we construct moments with other weights from the spectral functions obtained from Eq. (3.3). We choose to use the definition

$$I_{\text{ex}}^{(w)}(s_0) = \sum_{\text{sbin} \leq s_0} \left(\int_{\text{sbin}-\text{dsbin}/2}^{\text{sbin}+\text{dsbin}/2} ds w(s; s_0) \right) \rho^{(1+0)}(\text{sbin})\tag{3.5}$$

for all moments considered in this article.

B. Correlations

As shown in Ref. [2], there was a problem with the publicly posted 2005 and 2008 versions of the ALEPH covariance matrices. This problem, since corrected in Ref. [1], turns out to have resulted from an inadvertent omission of contributions to the correlations between different bins induced by the unfolding procedure. The problem was discovered by producing fake data sets from a multivariate gaussian distribution based on the posted ALEPH data and covariance matrices, and then comparing the resulting fake data to the actual ALEPH data. The result of this test is shown in Fig. 2, which is the same as Fig. 3 of Ref. [2]. The top panel shows the experimental data taken from Ref. [4], the bottom panel a typical fake data set produced using the corresponding covariance matrix. The absence of the strong correlations seen in the actual data from the corresponding fake data is what signals the existence of the problem with the previous version of the ALEPH covariance matrix. Figure 3 shows the result of performing the same test on the updated and corrected results

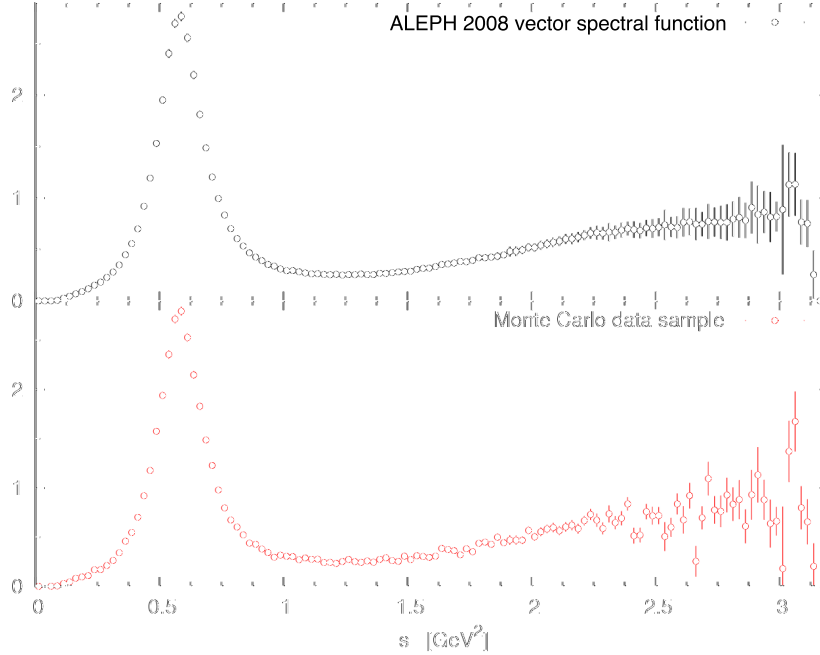


FIG. 2: *Vector spectral function times $2\pi^2$. Top panel: ALEPH data from 2008 [4]; bottom panel: Monte Carlo sample with 2008 covariance matrix.*

reported in Ref. [1], the top panel again showing the actual ALEPH data and the bottom panel a typical fake data set. The fake data (red points) obviously behave much more like the corresponding real data than was the case previously.⁴ We have examined many such fake data sets with the same conclusion. A similar exercise was, of course, carried out for the A -channel case.

C. Comparison with OPAL data

In Fig. 4 we show the vector and axial spectral functions as measured by ALEPH [1, 34] and OPAL [8]. The normalizations of the spectral functions for both experiments have been updated to take into account modern values for relevant branching fractions; for the normalization of ALEPH data, see Sec. III A above, and for the normalization of OPAL data, see Sec. III of Ref. [13].

While there is in general good agreement between the ALEPH and OPAL spectral functions, a detailed inspection reveals some tension between the two, given the size of the errors, for instance in the regions below 0.5 GeV^2 and around 2 GeV^2 in the vector channel, with possibly anti-correlated tensions in the same regions in the axial channel. The presence of a large $D = 0$, 1-loop α_s -independent contribution in the weighted OPE integrals enhances the impact of such small discrepancies on the output $\alpha_s(m_\tau^2)$. We quantify the impact of these differences below, showing that they lead to some tension between the values for $\alpha_s(m_\tau^2)$

⁴ Even though the new wider binning near the kinematic endpoint makes it somewhat harder to see such differences in this region.

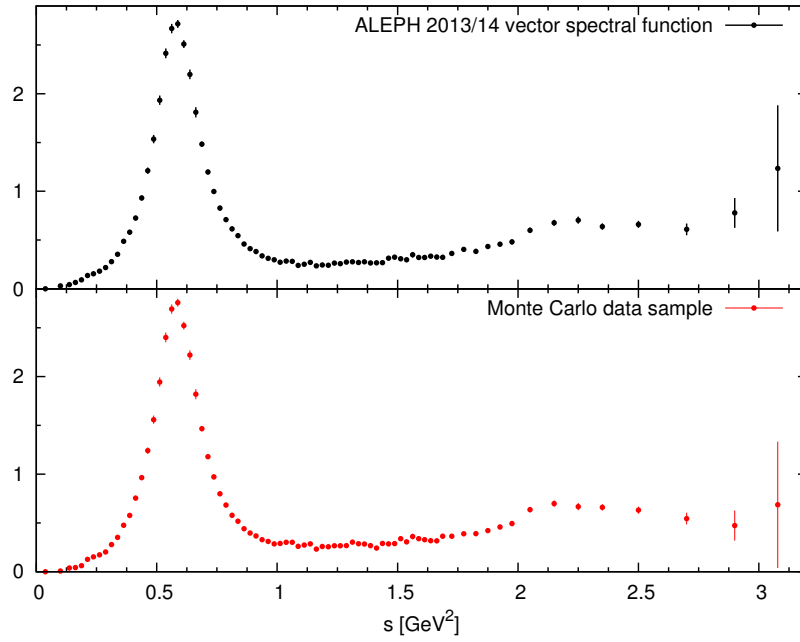


FIG. 3: *Vector spectral function times $2\pi^2$. Top panel: ALEPH data from 2013 [1]; bottom panel: Monte Carlo sample with 2013 covariance matrix.*

obtained from the two data sets, though the results turn out to agree within total estimated errors.

IV. FITTING STRATEGY

As already explained in Sec. II, and in more detail in Refs. [12, 13], non-pinned weights are needed in order to get a handle on the DV parameters of Eq. (2.10). The simplest and most robust choice of weight allowing us to extract these parameters is the weight $\hat{w}_0(x) = 1$. In order to check the stability of these simple fits, we also perform simultaneous fits of the weights \hat{w}_0 and \hat{w}_2 , and of \hat{w}_0 , \hat{w}_2 and \hat{w}_3 , as in Ref. [12, 13]. This gives us access to the $D = 6$ and $D = 8$ terms in the OPE, but also allows us to test for the consistency of the values of $\alpha_s(m_\tau^2)$ and the DV parameters between our different fits.

The values we obtain for $I_{\text{ex}}^{(w)}(s_0)$ from the ALEPH data are highly correlated, both between different values of s_0 , and between different weight functions. If we consider only fits using $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ for a range of s_0 values, it turns out that fully correlated χ^2 fits are possible, but if we also include $I_{\text{ex}}^{(\hat{w}_2)}(s_0)$ and $I_{\text{ex}}^{(\hat{w}_3)}(s_0)$ in the fits, the complete correlation matrices become too singular. For fits with multiple weights, we will follow Refs. [12, 13], using instead the block-diagonal “fit quality”

$$\mathcal{Q}^2 = \sum_w \sum_{s_0^i, s_0^j} \left(I_{\text{ex}}^{(w)}(s_0^i) - I_{\text{th}}^{(w)}(s_0^i; \vec{p}) \right) \left(C^{(w)} \right)_{ij}^{-1} \left(I_{\text{ex}}^{(w)}(s_0^j) - I_{\text{th}}^{(w)}(s_0^j; \vec{p}) \right), \quad (4.1)$$

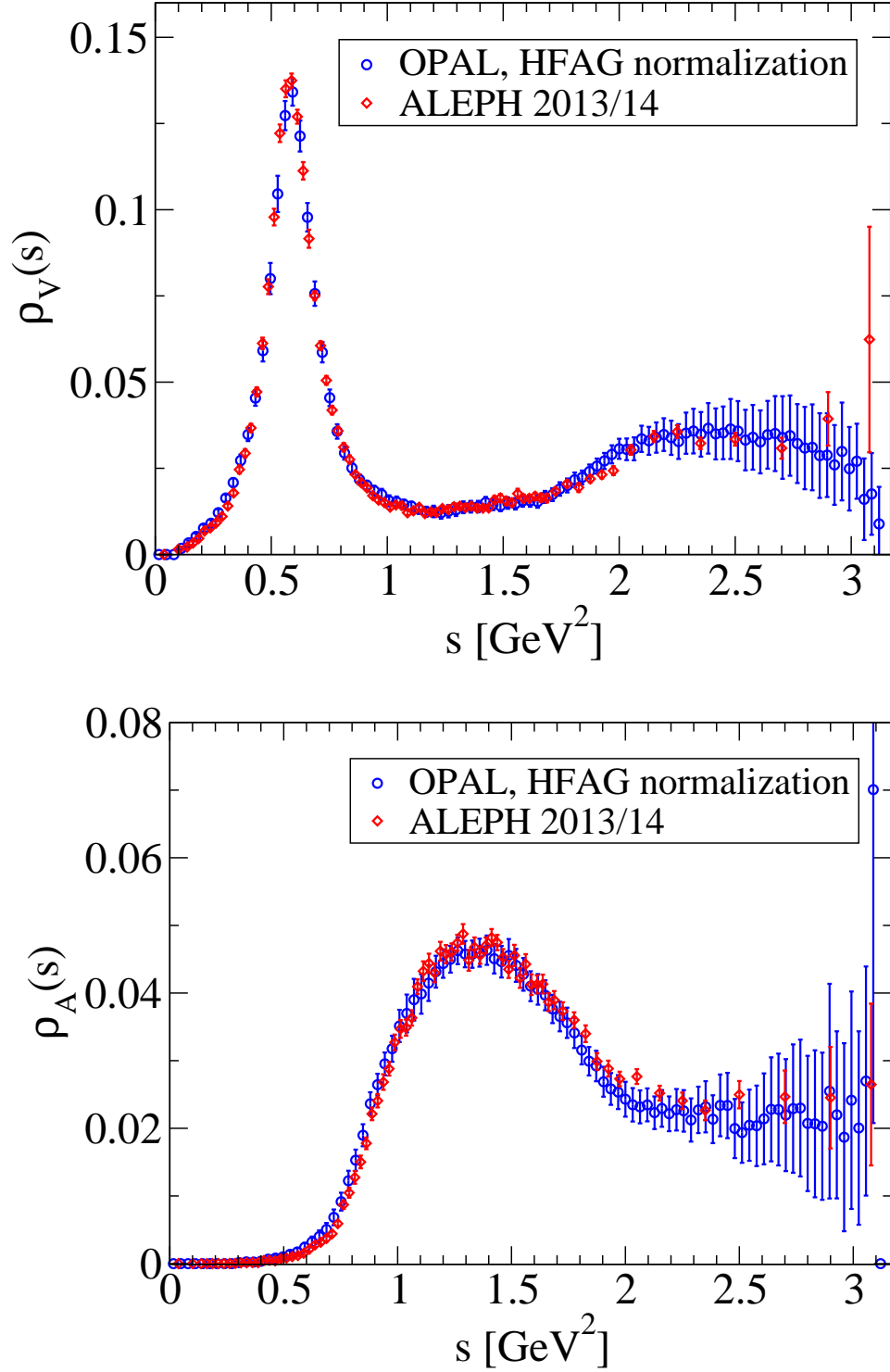


FIG. 4: Comparison of ALEPH and OPAL data for the spectral functions. Top panel: $I = 1$ vector channel; bottom panel: $I = 1$ continuum (pion-pole subtracted) axial channel.

where we have made the dependence of $I_{\text{th}}^{(w)}$ on the fit parameters \vec{p} explicit. The matrix $C^{(w)}$ is the (block-diagonal) covariance matrix of the set of moments with fixed weight w and s_0 running over the chosen fit window range. The sums over s_0^i and s_0^j are over bins i and j , and the sum over w is over \hat{w}_0 and \hat{w}_2 , or over \hat{w}_0 , \hat{w}_2 and \hat{w}_3 .⁵ The motivation for this choice is that the cross-correlations between two moments arise mainly because the weight functions used in multiple-moment fits appear to be close to being linearly dependent in practice (even though, as a set of polynomials, of course they are not). This near-linear dependence is possibly caused by the relatively large errors on the data for values of s toward m_τ^2 , because it is primarily in this region that the weights \hat{w}_0 , \hat{w}_2 and \hat{w}_3 differ from each other.

An important observation is that we can freely choose our fit quality \mathcal{Q}^2 , as long as errors are propagated taking the full data correlation matrix into account. In our case, we choose to estimate fit errors for fits using Eq. (4.1) by propagating the data covariance matrix through a small fluctuation analysis; for details on how this is done, we refer to the appendix of Ref. [12]. We note that the fit quality \mathcal{Q}^2 does not follow a standard χ^2 distribution, so that no absolute meaning can be attached to the minimum value obtained in a fit of this type.

The theoretical moments $I_{\text{th}}^{(w)}(s_0; \vec{p})$ are non-linear functions of (some of) the fit parameters \vec{p} , and it is thus not obvious what the probability distribution of the model parameters looks like. As in Ref. [13], we will therefore also explore the posterior probability distribution of the model parameters, assuming that the input data follow a multivariate gaussian distribution. In order to map out this probability distribution, we use the same Markov-chain Monte Carlo code `hrothgar` [22] as was used in Ref. [13], to which we refer for more details. The distribution generated by `hrothgar` is proportional to $\exp[-\mathcal{Q}^2(\vec{p})/2]$ on the space of parameters, given the data.

V. FITS

In this section we present our fits, leaving the discussion of α_s and other parameters obtained from these fits to Sec. VI. We first present fits to moments constructed from the V spectral function only, followed by fits using both the V and A spectral moments. We have considered χ^2 fits to $I_{\text{ex}}^{(\hat{w}_0)}$ and combined fits using fit qualities of the form (4.1) to $I_{\text{ex}}^{(\hat{w}_0)}$, $I_{\text{ex}}^{(\hat{w}_2)}$, and $I_{\text{ex}}^{(\hat{w}_3)}$. Below we will show only the χ^2 fits to $I_{\text{ex}}^{(\hat{w}_0)}$ and the \mathcal{Q}^2 fits to all three moments. The results from \mathcal{Q}^2 fits to the two moments $I_{\text{ex}}^{(\hat{w}_0)}$ and $I_{\text{ex}}^{(\hat{w}_2)}$ are completely consistent with these, and we therefore omit them below in the interest of brevity.

As reviewed above, and discussed in much more detail in Refs. [12, 13], the necessity to fit not only OPE parameters, but also DV parameters, makes it impossible to fit spectral moments for the sum of the V and A spectral functions. Already for $I_{\text{ex}}^{(\hat{w}_0)}$ this would entail a 9-parameter fit, and with the existing data such fits turn out to be unstable. Reference [1] did perform fits to moments of the $V + A$ spectral function at the price of neglecting duality violations and contributions from $D > 8$ terms in the OPE; we will compare our fits with those of Ref. [1] in detail in Sec. VII below.

From Fig. 4, we see that the only “feature” in the A channel is the peak corresponding to the a_1 meson. In contrast, the V channel data indicate the existence of more resonance-like

⁵ If only one weight is included in the sum, \mathcal{Q}^2 reverts to the standard χ^2 .

features than just the ρ meson peak around $s = 0.6 \text{ GeV}^2$, even though the resolution is not good enough to resolve multiple resonances beyond the ρ . If we wish to avoid making the assumption that already the lowest peak in each channel is in the asymptotic regime in which the *ansatz* (2.10) is valid, we should limit ourselves to fits to the V channel only. However, we will present also fits to the combined V and A channels below, and see that the results are consistent with those from fits to only the V channel.

In all cases, we find it necessary to include the moment \hat{w}_0 in our fits in order to determine both $\alpha_s(m_\tau^2)$ and the DV parameters. While one might consider fits to the spectral function itself, such fits are found to be insufficiently sensitive to the parameter $\alpha_s(m_\tau^2)$, and hence have not been pursued.⁶ Fits involving only pinched moments such as \hat{w}_2 and \hat{w}_3 , on the other hand, are insufficiently sensitive to the DV parameters. All our fits will thus include the spectral moments $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$, either in the V channel alone, or in the combined V and A channels. In the latter case, there is a separate set of DV parameters for each of these channels, but the fit parameter $\alpha_s(m_\tau^2)$ is, of course, common to both.⁷

⁶ It is important to distinguish fits to the spectral function itself from fits to *the moments of the spectral function*; they are quite different. Even in the case of the \hat{w}_0 moment, the integral $I_{V;\text{ex}}^{(\hat{w}_0)}(s_0)$ contains all the data from threshold to s_0 and always includes, in particular, the ρ peak. On the other hand, a fit of the DV *ansatz* (2.10) to the vector spectrum would probably only include data for s_0 between s_{min} and m_τ^2 , and, since one needs to choose $s_{\text{min}} \gg m_\rho^2$, the ρ peak is clearly excluded. The change in $I_{V/A;\text{ex}}^{(\hat{w}_0)}(s_0)$ as s_0 is increased from the upper edge of bin k to the upper edge of bin $k+1$ is, of course, equal to the average value of the relevant spectral function, $\rho_{V/A}$, in bin $k+1$. As such, in fits which employ all possible $s_0 \geq s_{\text{min}}$, the fact that the s_0 -dependence of $I_{V/A;\text{ex}}^{(\hat{w}_0)}(s_0)$ is one of the key elements entering the fit means that spectral function values in the interval $s_{\text{min}} \leq s \leq m_\tau^2$ are part of the input, but clearly not the *only* input.

Let us be even more specific. First, as already noted, even for single-weight \hat{w}_0 fits, the integral of the experimental spectral function over the region from threshold to s_{min} enters the \hat{w}_0 moment for all s_0 . While this is a region in which the OPE and the DV *ansatz* are not valid, this additional input turns out to be crucial; fits for both $\alpha_s(m_\tau^2)$ and the DV parameters are not possible without including it. Second, as seen in our previous analysis employing the OPAL data, fit results are not changed if, rather than using integrated data for all available $s_0 > s_{\text{min}}$, one instead employs a winnowed set thereof in the analysis. For such a winnowed set, it is only the sums of the experimental spectral function values over the bins lying between adjacent winnowed s_0 , and not the full set of spectral function values in all bins in those intervals, that determine the s_0 variation entering the fit. Finally, all of the multi-weight fits we employ involve weights, $w(x = s/s_0)$, which are themselves s_0 dependent. This means that the s_0 dependence of the DV part of the corresponding theory moments results not just from the values of $\rho^{\text{DV}}(s)$ in the interval $s_0 \leq s \leq m_\tau^2$ (where experimental constraints exist), but also involve s_0 - and $w(x)$ -dependent weighted integrals of the DV *ansatz* form in the interval from m_τ^2 to ∞ . It would thus be incorrect to characterize the moment-based fit analysis we employ as in any way representing simply a fit to the experimental spectral functions.

⁷ The $D > 2$ OPE coefficients are also generally different between the V and A channels [7]. In the case of C_4 (which, due to the absence of terms linear in x , does not enter for the weights we employ, in the approximation of dropping contributions higher-than-leading order in α_s) the full gluon condensate and leading-order quark condensate contributions are the same for the V and A channels. For polynomial weights with a term linear in x , $D = 4$ contributions would be present, and one could impose the resulting near-equality of C_4 in the V and A channels. This was done in the version of the analysis performed by

s_{\min} (GeV ²)	χ^2/dof	$p\text{-value}$ (%)	α_s	δ_V	γ_V	α_V	β_V
1.425	33.0/21	5	0.312(11)	3.36(36)	0.66(22)	-0.33(61)	3.27(33)
1.475	29.5/19	6	0.304(11)	3.32(41)	0.70(25)	-1.21(73)	3.72(39)
1.500	29.5/18	4	0.304(11)	3.32(41)	0.70(25)	-1.19(87)	3.71(45)
1.525	29.0/17	3	0.302(11)	3.37(43)	0.68(26)	-1.49(94)	3.86(48)
1.550	24.5/16	8	0.295(10)	3.50(50)	0.62(29)	-2.43(94)	4.32(48)
1.575	23.5/15	8	0.298(11)	3.50(47)	0.62(28)	-2.1(1.0)	4.15(53)
1.600	23.4/14	5	0.297(12)	3.50(48)	0.62(28)	-2.1(1.1)	4.16(56)
1.625	23.4/13	4	0.298(13)	3.47(50)	0.63(28)	-2.0(1.2)	4.12(62)
1.675	23.1/11	2	0.301(15)	3.35(60)	0.68(31)	-1.7(1.4)	3.96(70)
1.425	33.2/21	4	0.331(15)	3.20(34)	0.74(21)	-0.30(61)	3.24(33)
1.475	29.5/19	6	0.320(14)	3.16(40)	0.78(24)	-1.20(73)	3.70(39)
1.500	29.5/18	4	0.320(15)	3.16(40)	0.78(24)	-1.19(87)	3.69(45)
1.525	28.9/17	4	0.317(14)	3.22(42)	0.75(25)	-1.51(93)	3.85(48)
1.550	24.3/16	8	0.308(13)	3.36(49)	0.69(28)	-2.48(93)	4.33(48)
1.575	23.3/15	8	0.311(14)	3.35(46)	0.69(27)	-2.2(1.0)	4.17(52)
1.600	23.3/14	6	0.311(15)	3.36(47)	0.69(27)	-2.2(1.1)	4.19(56)
1.625	23.2/13	4	0.312(16)	3.33(49)	0.70(28)	-2.1(1.2)	4.15(62)
1.675	23.0/11	2	0.314(19)	3.23(58)	0.74(30)	-1.8(1.5)	4.02(74)

TABLE 1: V channel fits to $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ from $s_0 = s_{\min}$ to $s_0 = m_\tau^2$. FOPT results are shown above the double line, CIPT below; no $D > 0$ OPE terms included in the fit. γ_V and β_V in units of GeV⁻².

A. Fits to vector channel data

We begin with fits to the single moment $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$, as a function of s_{\min} , with s_{\min} defined to be the minimum value of s_0 included in the fit. Since these are χ^2 fits, one may estimate the p -values for these fits; they are shown in the third column of Tab. 1. We note that the p -values are not large, but they are not small enough to exclude the validity of our fit function based on the ALEPH data. Judged by p -value, the fits with $s_{\min} = 1.55$ and 1.575 GeV² are the best fits, and we thus take the average value of the central values for the fit parameters from these two fits as our best value, with a statistical error that is the larger of the two (noting that these are essentially equal in size). For the strong coupling, we find

$$\begin{aligned}
\alpha_s(m_\tau^2) &= 0.296(11) , & (\text{FOPT}) , \\
&= 0.310(14) , & (\text{CIPT}) .
\end{aligned}
\tag{5.1}$$

OPAL but not in the analyses of the ALEPH collaboration, including Ref. [1]. The fact that the fitted value of the gluon condensate obtained from independent V and A channel fits in Ref. [1] are not close to agreeing within errors is, in fact, a clear sign of the unphysical nature of these fits, see Sec. VII below.

	α_s	δ_V	γ_V	α_V	β_V
α_s	1	0.600	-0.606	0.689	-0.653
δ_V	0.600	1	-0.994	0.310	-0.297
γ_V	-0.606	-0.994	1	-0.330	0.315
α_V	0.689	0.310	-0.330	1	-0.996
β_V	-0.653	-0.297	0.315	-0.996	1

TABLE 2: *Parameter correlation matrix for the V channel \hat{w}_0 FOPT fit with $s_{\min} = 1.55 \text{ GeV}^2$ shown in Tab. 1.*

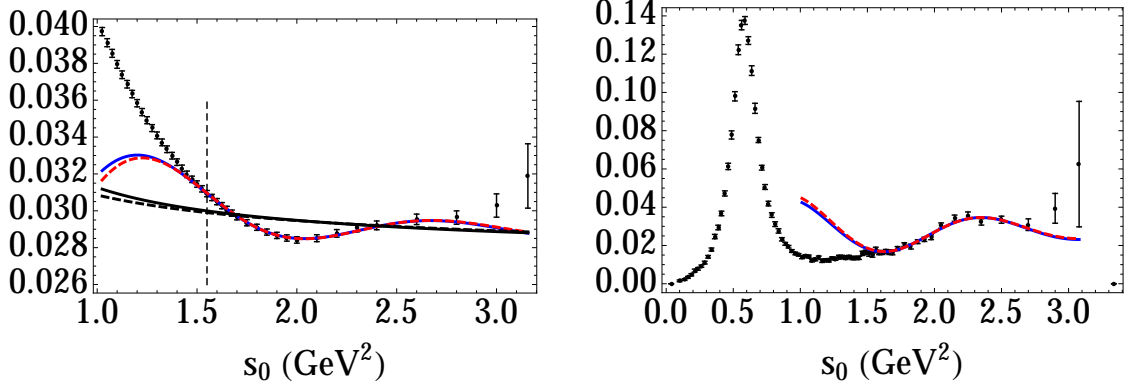


FIG. 5: *Left panel: comparison of $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ and $I_{\text{th}}^{(\hat{w}_0)}(s_0)$ for the $s_{\min} = 1.55 \text{ GeV}^2$ V channel fits of Tab. 1. Right panel: comparison of the theoretical spectral function resulting from this fit with the experimental results. CIPT fits are shown in red (dashed) and FOPT in blue (solid). The (much flatter) black curves in the left panel represent the OPE parts of the fits, i.e., the fit results with the DV parts removed. The vertical dashed line indicates the location of s_{\min} .*

The difference between the FOPT and CIPT results reflects the well-known fact that the two prescriptions show no sign of converging to one another as the truncation order is increased [14, 15]. We observe that the p -value starts to decrease again from $s_{\min} = 1.6 \text{ GeV}^2$, indicating that the data become too sparse for an optimal fit. We investigated the sensitivity of these fits to omitting the data in up to four bins with the largest values of s , and found no significant difference. This is no surprise, given the errors shown in Fig. 5. For illustration, we show the parameter correlation matrix for the FOPT fit with $s_{\min} = 1.55 \text{ GeV}^2$ in Tab. 2.

In Fig. 5 we show the results of CIPT and FOPT fits to $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ for $s_{\min} = 1.55 \text{ GeV}^2$. The left panel shows the results of the fits for the moment, the right-hand panel the OPE+DV versions of the spectral functions resulting from these fits.

As in Ref. [13], we studied the posterior probability distribution, using the same Markov-chain Monte Carlo code, `hrothgar` [22]. We remind the reader that it is not obvious what this distribution looks like, even if we assume that the data errors follow a multivariate gaussian distribution. For the fits of Tab. 1, this code generates points in the 5-dimensional parameter space, and computes the χ^2 value associated with each of these points. These

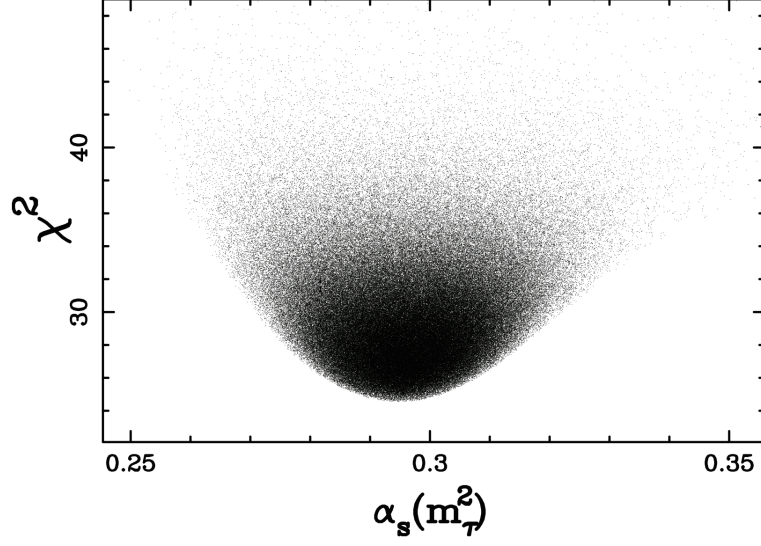


FIG. 6: χ^2 versus $\alpha_s(m_\tau^2)$, FOPT, $s_{\min} = 1.55 \text{ GeV}^2$, 1250000 points.

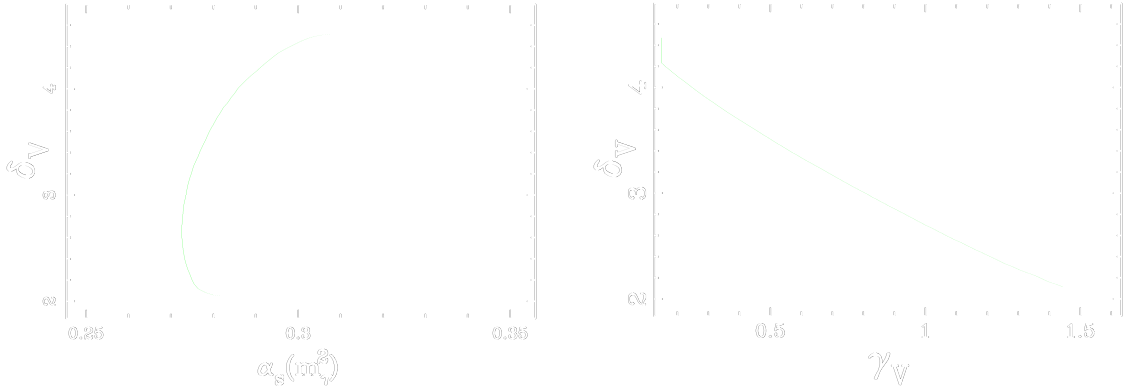


FIG. 7: Two-dimensional contour plots showing δ_V versus $\alpha_s(m_\tau^2)$ (left) and δ_V versus γ_V (right) for the V channel \hat{w}_0 FOPT, $s_{\min} = 1.55 \text{ GeV}^2$ fit. Blue (darker) areas and green (lighter) areas contain 68%, respectively, 95% of the distribution. γ_V in units of GeV^{-2}

points are distributed as $\exp[-\chi^2(\vec{p})]$, with \vec{p} the parameter vector, and χ^2 evaluated on the ALEPH data (including the full covariance matrix) and the values of the parameters at these points.

In Fig. 6 we show χ^2 as a function of $\alpha_s(m_\tau^2)$, choosing the FOPT fit with $s_{\min} = 1.55 \text{ GeV}^2$. Since for each $\alpha_s(m_\tau^2)$ points with many different values for the other four parameters are generated stochastically, the distribution appears as the cloud shown in the figure. This distribution shows a unique minimum for the value of χ^2 , at approximately $\alpha_s(m_\tau^2) = 0.295$, consistent with Tab. 1. The width of the distribution is also roughly consistent with the error of ± 0.010 , but we see that the distribution of points is not entirely symmetric around the minimum. There is no alternative (local) minimum, as was the case

s_{\min} (GeV ²)	\mathcal{Q}^2/dof	α_s	δ_V	γ_V	α_V	β_V	$10^2 C_{6V}$	$10^2 C_{8V}$
1.425	106.0/71=1.49	0.305(10)	3.02(38)	0.87(24)	-0.68(56)	3.43(31)	-0.59(17)	0.94(29)
1.475	93.3/65=1.43	0.302(10)	3.07(44)	0.85(27)	-1.41(68)	3.81(36)	-0.71(16)	1.19(28)
1.500	93.2/62=1.50	0.302(10)	3.08(45)	0.85(27)	-1.40(77)	3.80(40)	-0.71(18)	1.19(30)
1.525	85.6/59=1.45	0.298(10)	3.21(49)	0.78(29)	-1.96(78)	4.08(41)	-0.79(16)	1.36(27)
1.550	76.3/56=1.36	0.295(10)	3.30(52)	0.74(30)	-2.48(81)	4.33(41)	-0.86(14)	1.50(24)
1.575	74.5/53=1.41	0.297(10)	3.29(51)	0.74(29)	-2.25(87)	4.22(44)	-0.83(16)	1.43(27)
1.600	74.2/50=1.48	0.297(11)	3.31(51)	0.73(30)	-2.27(92)	4.23(47)	-0.83(16)	1.44(29)
1.625	73.8/47=1.57	0.298(11)	3.28(54)	0.74(31)	-2.16(99)	4.18(50)	-0.81(18)	1.40(32)
1.675	72.0/41=1.76	0.299(12)	3.28(63)	0.74(34)	-2.1(1.1)	4.13(57)	-0.80(21)	1.37(39)
1.425	98.6/71=1.39	0.328(16)	3.17(39)	0.77(25)	-0.43(61)	3.30(32)	-0.60(19)	0.83(35)
1.475	89.5/65=1.38	0.319(14)	3.11(44)	0.81(27)	-1.24(71)	3.72(37)	-0.76(16)	1.18(31)
1.500	89.4/62=1.44	0.319(15)	3.11(44)	0.81(27)	-1.20(81)	3.70(42)	-0.76(18)	1.16(34)
1.525	82.1/59=1.39	0.314(14)	3.22(48)	0.77(28)	-1.81(80)	4.00(42)	-0.85(15)	1.37(28)
1.550	73.7/56=1.32	0.309(13)	3.28(51)	0.74(30)	-2.39(82)	4.28(42)	-0.93(13)	1.53(25)
1.575	71.8/53=1.35	0.311(14)	3.28(50)	0.74(29)	-2.12(89)	4.15(45)	-0.89(15)	1.45(28)
1.600	71.7/50=1.43	0.311(14)	3.28(51)	0.74(29)	-2.16(94)	4.17(48)	-0.90(15)	1.46(29)
1.625	71.5/47=1.52	0.312(15)	3.24(53)	0.75(30)	-2.0(1.0)	4.11(51)	-0.88(17)	1.42(34)
1.675	69.8/41=1.70	0.313(16)	3.22(63)	0.76(33)	-1.9(1.2)	4.04(59)	-0.86(20)	1.38(42)

TABLE 3: V channel fits to $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$, $I_{\text{ex}}^{(\hat{w}_2)}(s_0)$ and $I_{\text{ex}}^{(\hat{w}_0)}(s_3)$ from $s_0 = s_{\min}$ to $s_0 = m_\tau^2$, FOPT results are shown above the double line, CIPT below; $D = 6, 8$ OPE terms included in the fit. γ_V and β_V in units of GeV⁻², C_{6V} in units of GeV⁶ and C_{8V} in units of GeV⁸.

with the OPAL data [13].

We also find the parameters δ_V and γ_V to be much better constrained than was the case for the corresponding fits to the OPAL data in Ref. [13]. The distributions in the δ_V - $\alpha_s(m_\tau^2)$ and δ_V - γ_V planes are shown in the left and right panels of Fig. 7.⁸ Since for all other fits presented in the rest of this article the conclusions about the posterior probability distribution found with **hrothgar** are similar, we will refrain from showing the analogues of Figs. 6 and 7 for those fits.

Next, we consider simultaneous fits to the moments $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$, $I_{\text{ex}}^{(\hat{w}_2)}(s_0)$ and $I_{\text{ex}}^{(\hat{w}_0)}(s_3)$; results for the same values of s_{\min} as before are given in Tab. 3. These fits are performed by minimizing \mathcal{Q}^2 as defined in Eq. (4.1), with correlations between different moments omitted. However, the full correlation matrix, including correlations between different moments, has been taken into account in the parameter fit error estimates shown in the table. These errors were determined by linear propagation of the full data covariance matrix; for a detailed explanation of the method, we refer to the appendix of Ref. [12].

Judging by the values of \mathcal{Q}^2/dof ,⁹ again the two fits for $s_{\min} = 1.55$ and 1.575 GeV² are

⁸ Note that the vertical axis covers the interval $\delta_V \in [2, 5]$, to be compared with the significantly larger interval $\delta_V \in [-2, 5]$ in Fig. 2 of Ref. [13].

⁹ Which, given the fact that \mathcal{Q}^2 is not equal to χ^2 for these fits, cannot easily be translated into p -values.

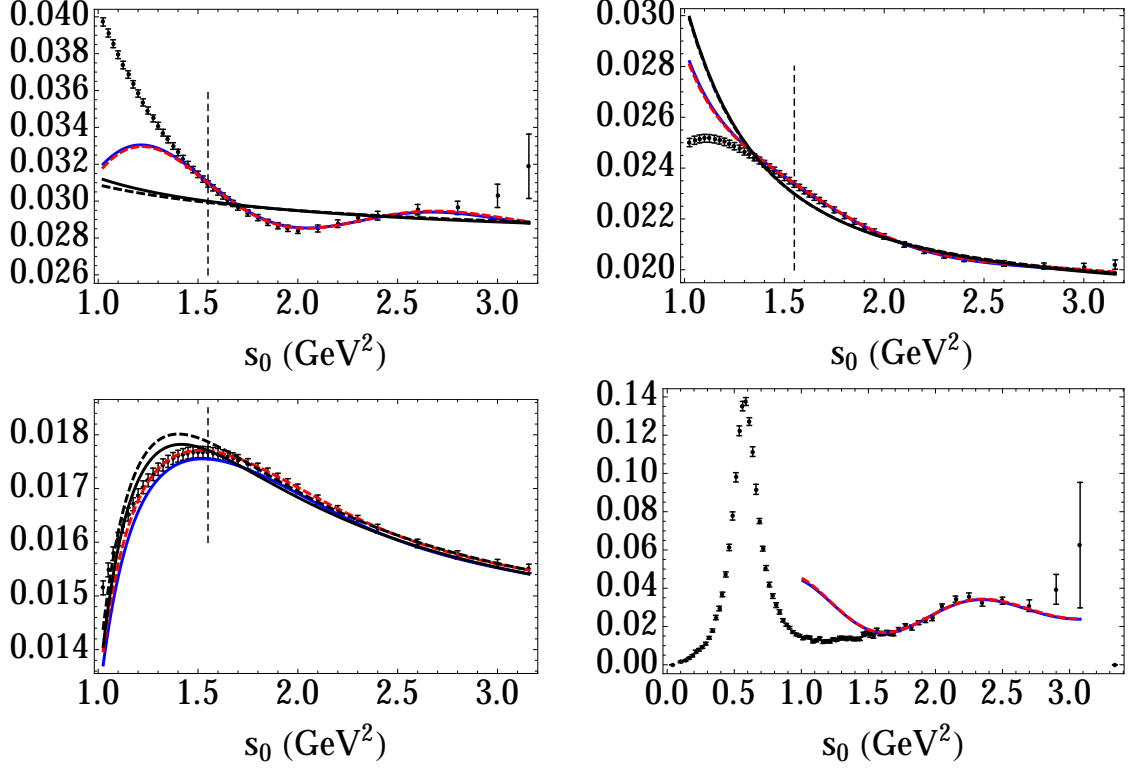


FIG. 8: Upper left panel: comparison of $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ and $I_{\text{th}}^{(\hat{w}_0)}(s_0)$ for the $s_{\text{min}} = 1.55 \text{ GeV}^2$ V channel fits of Tab. 3. Lower left panel and upper right panel: analogous comparisons for $I^{(\hat{w}_2)}(s_0)$ (upper right panel) and $I^{(\hat{w}_3)}(s_0)$ (lower left panel). CIPT fits are shown in red (dashed) and FOPT in blue (solid). Lower right panel: comparison of the theoretical spectral function resulting from this fit with the experimental results. The black curves (which are much flatter for the \hat{w}_0 case) represent the OPE parts of the fits. The vertical dashed line indicates the location of s_{min} .

the optimal ones. Averaging parameter values between these two fits, we find

$$\begin{aligned} \alpha_s(m_\tau^2) &= 0.296(10) , & (\text{FOPT}) , \\ &= 0.310(14) , & (\text{CIPT}) , \end{aligned} \quad (5.2)$$

in excellent agreement with Eq. (5.1). We have also considered fits involving only the two moments $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ and $I_{\text{ex}}^{(\hat{w}_2)}(s_0)$, and find results very similar those contained in Tabs. 1 and 3. In Fig. 8 we show the quality of the fits of Tab. 3 for $s_{\text{min}} = 1.55 \text{ GeV}^2$.

We end this subsection with several comments. First, we see that pinching indeed serves to suppress the role of DV contributions. The upper right panel in Fig. 8 shows the singly pinched \hat{w}_2 case and the lower left panel shows the doubly pinched \hat{w}_3 case. There is also a significant difference between the colored and black curves in all panels, though with the onset of this difference shifting to lower s_0 as the degree of pinching is increased. The existence of these differences implies that, with the errors on the ALEPH data, the presence of duality violations is evident for all three moments. This, in turn, implies that omitting duality violations from the theory side of the corresponding FESRs has the potential to produce a significant additional systematic error on $\alpha_s(m_\tau^2)$ (and the higher D OPE

s_{\min} (GeV ²)	χ^2/dof	$p\text{-value}$ (%)	α_s	δ_V δ_A	γ_V γ_A	α_V α_A	β_V β_A
1.500	49.8/37	8	0.310(14)	3.45(40) 1.85(38)	0.62(24) 1.38(20)	-1.0(1.0) 4.5(1.2)	3.60(53) 2.46(59)
1.525	48.6/35	6	0.309(15)	3.53(42) 1.99(40)	0.59(25) 1.31(20)	-1.2(1.2) 4.4(1.2)	3.71(60) 2.49(62)
1.550	40.0/33	19	0.297(11)	3.57(48) 1.56(49)	0.58(28) 1.44(22)	-2.33(97) 5.43(89)	4.27(50) 1.99(46)
1.575	38.7/31	16	0.300(12)	3.57(45) 1.67(51)	0.58(26) 1.41(23)	-1.9(1.1) 5.22(94)	4.08(55) 2.10(48)
1.600	37.2/298	14	0.300(12)	3.56(46) 1.41(57)	0.59(27) 1.52(25)	-2.0(1.2) 5.4(1.0)	4.10(59) 2.01(52)
1.625	35.4/27	13	0.300(13)	3.50(48) 0.90(72)	0.62(27) 1.73(29)	-1.9(1.3) 5.8(1.2)	4.07(64) 1.82(60)
1.500	49.7/37	8	0.327(18)	3.29(39) 1.92(39)	0.70(24) 1.35(20)	-1.0(1.0) 4.5(1.1)	3.59(53) 2.50(60)
1.525	48.5/35	6	0.326(19)	3.37(40) 2.06(41)	0.66(24) 1.28(21)	-1.2(1.2) 4.4(1.2)	3.70(60) 2.54(62)
1.550	39.7/33	20	0.311(13)	3.43(47) 1.61(49)	0.65(27) 1.43(22)	-2.38(96) 5.36(87)	4.28(49) 2.04(45)
1.575	38.4/31	17	0.315(15)	3.42(44) 1.72(52)	0.65(26) 1.39(24)	-2.0(1.1) 5.15(92)	4.10(56) 2.14(48)
1.600	36.9/29	15	0.314(15)	3.41(45) 1.46(58)	0.66(26) 1.50(25)	-2.1(1.2) 5.33(98)	4.13(59) 2.06(51)
1.625	35.1/27	14	0.314(16)	3.36(48) 0.96(72)	0.68(27) 1.71(29)	-2.0(1.3) 5.7(1.1)	4.11(64) 1.87(58)

TABLE 4: Combined V and A channel fits to $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ from $s_0 = s_{\min}$ to $s_0 = m_\pi^2$. FOPT results are shown above the double line, CIPT below; no $D > 0$ OPE terms are included in the fit. $\gamma_{V,A}$ and $\beta_{V,A}$ in units of GeV⁻².

coefficients) that cannot be estimated if only fits without DV parameters are attempted. We will return to this point in Sec. VII below. Second, we note that the spectral function itself below $s = s_{\min}$ is not very well described by the curves obtained from the fits. While the form of Eq. (2.10) constitutes a reasonable assumption for asymptotically large s , we do not know *a priori* what a reasonable value of s_{\min} should be. It is clear, however, that our *ansatz* works reasonably well for $s \gtrsim 1.5$ GeV², but that the asymptotic regime definitely does not include the region around the ρ peak.

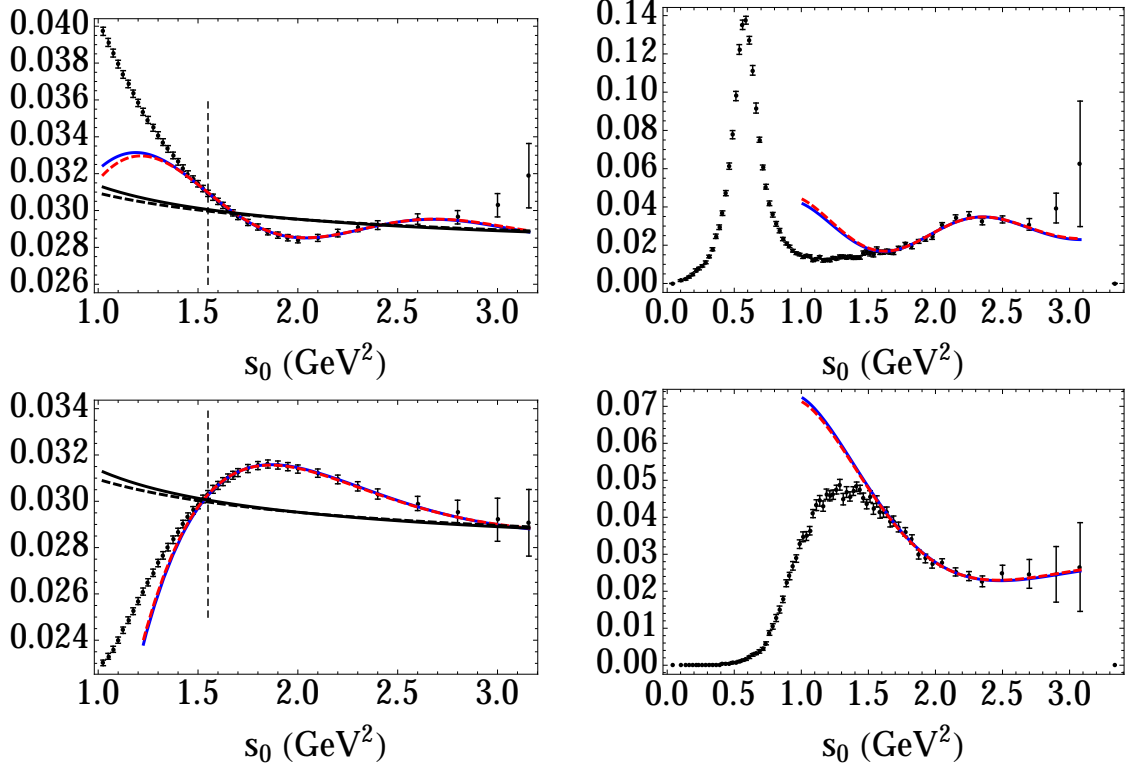


FIG. 9: *Left panels: comparison of $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ and $I_{\text{th}}^{(\hat{w}_0)}(s_0)$ for the $s_{\text{min}} = 1.55 \text{ GeV}^2$ combined V and A channel fits of Tab. 4 (V top, A bottom). Right panels: comparison of the theoretical spectral function resulting from this fit with the experimental results (V top, A bottom). CIPT fits are shown in red (dashed) and FOPT in blue (solid). The (much flatter) black curves on the left represent the OPE parts of the fits. The vertical dashed line indicates the location of s_{min} .*

B. Combined fits to vector and axial channel data

We now consider fits analogous to those of the preceding subsection, involving simultaneous fitting of the V and A spectral moments as a function of s_{min} . The fit parameter $\alpha_s(m_\tau^2)$ is common to the two channels, while the $D > 0$ OPE and DV parameters are distinct for each. Fits to $I_{\text{ex},V}^{(\hat{w}_0)}(s_0)$ and $I_{\text{ex},A}^{(\hat{w}_0)}(s_0)$ are shown in Tab. 4; we displayed fewer values of s_{min} for the sake of brevity.

Fits with $s_{\text{min}} = 1.55$ and 1.575 GeV^2 have the highest p -values, as before. Averaging the parameter values for these fits, we find

$$\begin{aligned} \alpha_s(m_\tau^2) &= 0.299(12) , & (\text{FOPT}) , \\ &= 0.313(15) , & (\text{CIPT}) , \end{aligned} \quad (5.3)$$

slightly higher values than those of Eqs. (5.1) and (5.2), but consistent within errors. The errors are χ^2 errors, since all correlations were taken into account in the fit; they are slightly larger than those found in the V -channel fits.

For $s_{\text{min}} = 1.55 \text{ GeV}^2$ we show the quality of the fits in the left panels of Fig. 9 and the V and A spectral-function comparisons obtained using parameter values from the fit in the

s_{\min} (GeV ²)	\mathcal{Q}^2/dof	α_s	$\delta_{V,A}$	$\gamma_{V,A}$	$\alpha_{V,A}$	$\beta_{V,A}$	$10^2 C_{6V,A}$	$10^2 C_{8V,A}$
1.475	182/131=1.39	0.297(7)	2.90(42)	0.95(26)	-1.61(65)	3.91(35)	-0.78(13)	1.31(23)
			2.26(35)	1.13(18)	4.92(58)	2.25(30)	-0.08(35)	1.12(96)
1.500	160/125=1.28	0.297(8)	2.92(43)	0.94(26)	-1.62(73)	3.91(39)	-0.78(14)	1.31(25)
			1.90(44)	1.29(21)	5.26(69)	2.08(36)	-0.26(44)	1.8(1.4)
1.525	149/119=1.25	0.294(8)	3.08(48)	0.86(28)	-2.16(75)	4.18(40)	-0.85(13)	1.46(23)
			1.86(48)	1.30(22)	5.38(72)	2.02(37)	-0.38(49)	2.1(1.6)
1.550	126/113=1.11	0.292(9)	3.19(51)	0.80(30)	-2.65(79)	4.42(41)	-0.90(13)	1.57(22)
			1.53(56)	1.42(24)	5.73(84)	1.84(43)	-0.63(61)	3.0(2.2)
1.575	124/107=1.16	0.293(9)	3.18(51)	0.81(29)	-2.47(84)	4.33(43)	-0.88(14)	1.52(24)
			1.57(61)	1.41(26)	5.67(86)	1.87(44)	-0.57(61)	2.8(2.2)
1.600	116/101=1.15	0.293(9)	3.20(52)	0.80(30)	-2.51(89)	4.35(46)	-0.89(14)	1.53(25)
			1.14(74)	1.59(29)	6.0(1.0)	1.72(53)	-0.73(72)	3.6(2.7)
1.625	112/95=1.18	0.294(10)	3.20(55)	0.79(31)	-2.43(95)	4.31(48)	-0.87(15)	1.50(28)
			0.85(92)	1.71(34)	6.2(1.2)	1.61(63)	-0.80(80)	4.0(3.2)
1.475	159/131=1.21	0.338(13)	3.45(32)	0.61(20)	-0.63(67)	3.42(35)	-0.58(16)	0.83(31)
			2.23(33)	1.25(21)	3.45(81)	3.02(42)	0.59(25)	-0.64(58)
1.500	146/125=1.17	0.328(15)	3.26(39)	0.72(24)	-0.92(79)	3.56(41)	-0.67(18)	1.00(35)
			1.96(41)	1.34(22)	4.41(89)	2.53(46)	0.25(40)	0.3(1.0)
1.525	136/119=1.14	0.320(13)	3.35(44)	0.69(26)	-1.59(79)	3.90(41)	-0.80(15)	1.26(29)
			1.93(46)	1.32(23)	4.76(83)	2.35(43)	0.05(43)	0.78(12)
1.550	118/113=1.04	0.312(13)	3.35(49)	0.70(29)	-2.28(81)	4.23(42)	-0.90(13)	1.48(25)
			1.59(55)	1.44(25)	5.37(89)	2.03(46)	-0.33(56)	2.0(1.8)
1.575	115/107=1.07	0.315(13)	3.35(48)	0.70(28)	-1.98(88)	4.09(45)	-0.86(15)	1.39(29)
			1.65(59)	1.42(27)	5.23(92)	2.11(47)	-0.22(55)	1.6(1.7)
1.600	108/101=1.07	0.314(14)	3.33(49)	0.71(29)	-2.04(93)	4.12(47)	-0.87(15)	1.41(30)
			1.23(70)	1.60(30)	5.6(1.1)	1.95(55)	-0.37(64)	2.2(2.2)
1.625	105/95=1.10	0.315(15)	3.28(53)	0.73(30)	-1.9(1.0)	4.06(51)	-0.85(17)	1.37(34)
			0.96(85)	1.71(35)	5.7(1.2)	1.87(63)	-0.42(71)	2.4(2.5)

TABLE 5: Combined V and A channel fits to $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$, $I_{\text{ex}}^{(\hat{w}_2)}(s_0)$ and $I_{\text{ex}}^{(\hat{w}_3)}(s_0)$ from $s_0 = s_{\min}$ to $s_0 = m_\tau^2$. FOPT results are shown above the double line, CIPT below; $D = 6, 8$ OPE terms included in the fit. $\gamma_{V,A}$ and $\beta_{V,A}$ in units of GeV⁻², $C_{6V,A}$ in units of GeV⁶ and $C_{8V,A}$ in units of GeV⁸.

corresponding right-hand panels. We note that the fit curves in the axial case are essentially determined by the shoulder of the a_1 resonance, in contrast to what happens in the vector case, where the ρ peak is well away from the region relevant for the shape of the fit curves.

Tab. 5 shows the results of the combined V and A channel fits to the three moments $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$, $I_{\text{ex}}^{(\hat{w}_2)}(s_0)$ and $I_{\text{ex}}^{(\hat{w}_3)}(s_3)$. Judging by the values of \mathcal{Q}^2/dof , the best fits are again those with $s_{\min} = 1.55$ and 1.575 GeV², leading to

$$\alpha_s(m_\tau^2) = 0.293(9) , \quad (\text{FOPT}) , \quad (5.4)$$

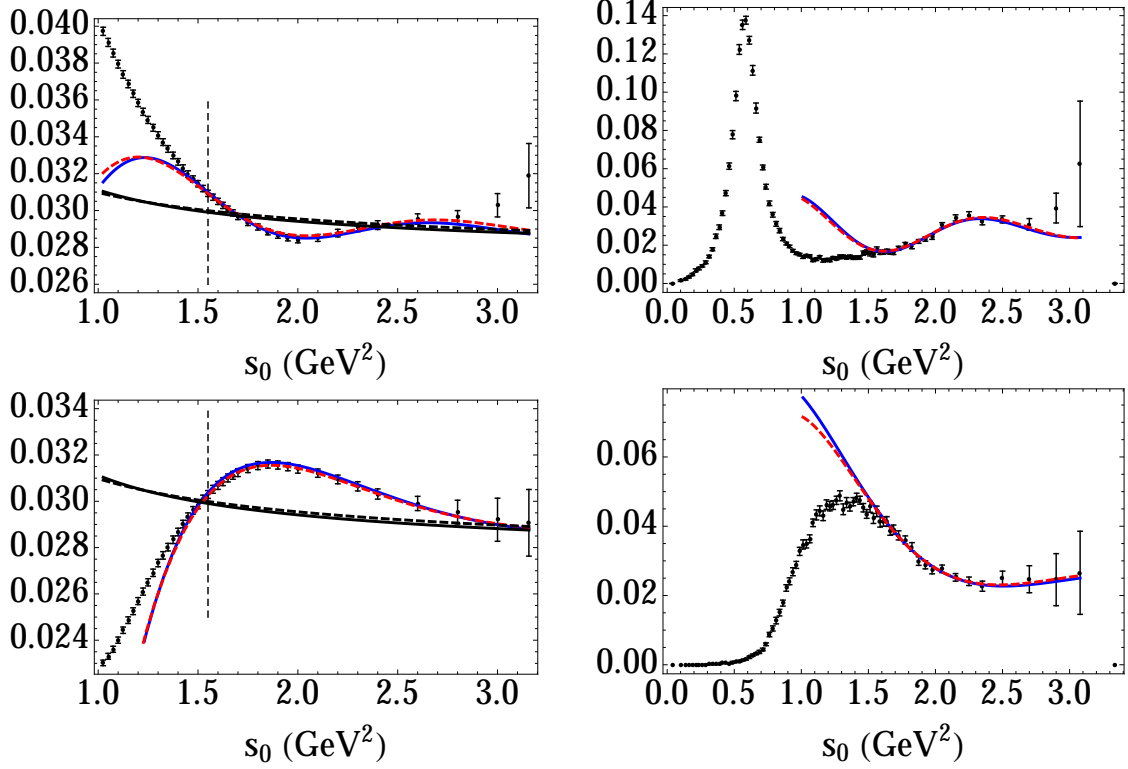


FIG. 10: *Left panels: comparison of $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ and $I_{\text{th}}^{(\hat{w}_0)}(s_0)$ for the $s_{\text{min}} = 1.55 \text{ GeV}^2$ combined V and A channel fits of Tab. 5 (V top, A bottom). Right panels: comparison of the theoretical spectral function resulting from this fit with the experimental results (V top, A bottom). CIPT fits are shown in red (dashed) and FOPT in blue (solid). The (much flatter) black curves on the left represent the OPE parts of the fits. The vertical dashed line indicates the location of s_{min} .*

$$= 0.313(13) , \quad (\text{CIPT}) .$$

These values are in good agreement with those of the other fits reported above. As before, fits to just the pair of moments $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ and $I_{\text{ex}}^{(\hat{w}_2)}(s_0)$ do not lead to any surprises. We show the quality of the fits of Tab. 5 for the moments $I_{\text{ex}}^{(\hat{w}_0)}(s_0)$ and the comparison of the resulting spectral functions to the experimental ones for both channels in Fig. 10. The fits for the other two moments look very similar to those in Fig. 8 for the V channel, and show a similar quality in the A channel.

VI. TESTS AND RESULTS

There are a number of consistency checks that can be applied once values for $\alpha_s(m_\tau^2)$ as well as the $D > 0$ OPE and DV parameters have been obtained from a fit. We will present some of these in Sec. VIA. Then, in Sec. VIB, we will present our final number for $\alpha_s(m_\tau^2)$, following this in Sec. VIC by a determination of the non-perturbative contribution to $R_{V+A;ud}$ and a comparison of the $D = 6$ OPE coefficients with the results of estimates

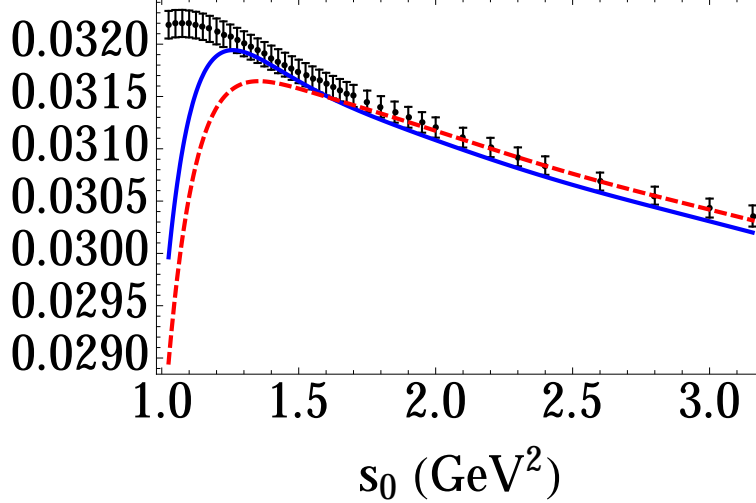


FIG. 11: The rescaled version of $R_{V+A;ud}(s_0)$ (the RHS of eq. (6.1)) as a function of s_0 . Theory curves from $s_{\min} = 1.55 \text{ GeV}^2$ entries of Tab. 5; CIPT (red, dashed) and FOPT (blue, solid).

based on the vacuum saturation approximation (VSA). In Sec. VID we will compare the present results with those from our fits to the OPAL data.

A. Tests

We consider first the comparison of the experimental value of

$$I_{ex,V}^{(\hat{w}_3)}(s_0) + I_{ex,A}^{(\hat{w}_3)}(s_0) = \frac{m_\tau^2}{12\pi^2|V_{ud}|^2 S_{EW}} R_{V+A;ud}(s_0) \quad (6.1)$$

with the function obtained from the fit. In Fig. 11 we show this comparison, using the parameter values for $s_{\min} = 1.55 \text{ GeV}^2$ from Tab. 5. The fitted curves are in good agreement everywhere above $s_0 \approx 1.3 \text{ GeV}^2$ ($s_0 \approx 1.5 \text{ GeV}^2$) for the FOPT (CIPT) fits.¹⁰ We include this test because (in rescaled form) it was originally advocated as an important confirmation of the analysis of Ref. [3]. One can see that our fits satisfy this test at least as well (see *e.g.* Fig. 73 of Ref. [3]). In other words, this test is not able to discriminate between the results of our analysis and those Refs. [1, 3, 4]. For more discussion on the comparison between our analysis and that of Refs. [1, 3, 4] we refer to Sec. VII.

As in Ref. [12], we may also consider the first and second Weinberg sum rules (WSRs) [23], as well as the DGMLY sum rule for the pion electro-magnetic mass splitting [24]. These sum rules can be written as

$$\int_0^\infty ds \left(\rho_V^{(1+0)}(s) - \rho_A^{(1+0)}(s) \right) = \int_0^\infty ds \left(\rho_V^{(1)}(s) - \rho_A^{(1)}(s) \right) - 2f_\pi^2 = 0, \quad (6.2)$$

¹⁰ We recall that even though correlations between different spectral moments are not included in the fit quality \mathcal{Q}^2 , those between bins within one spectral moment are included, making these fits strongly correlated.

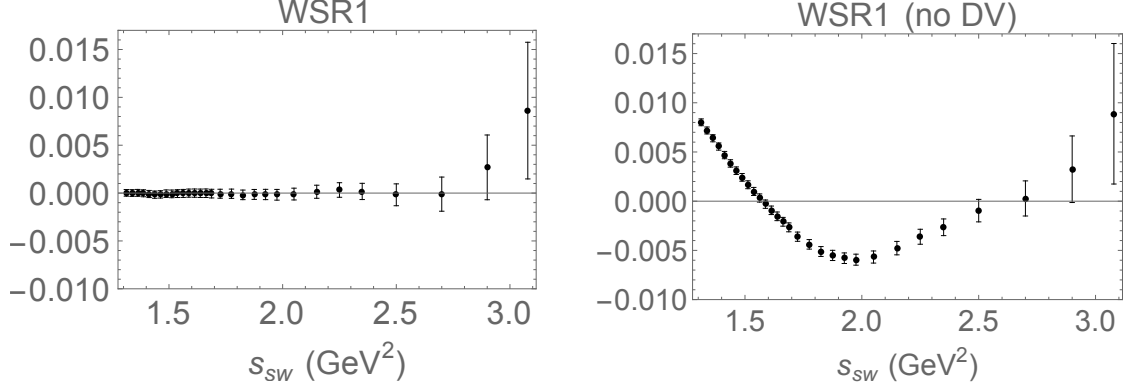


FIG. 12: The first Weinberg sum rule, with DVs (left panel) and without DVs (right panel), both in GeV^2 . Data have been used for $s < s_{\text{SW}}$, while the DV ansatz (2.10) with parameter values obtained from the $s_{\text{min}} = 1.55 \text{ GeV}^2$ fit has been used for $s > s_{\text{SW}}$. The figures shown use CIPT fits.

$$\int_0^\infty ds s \left(\rho_V^{(1+0)}(s) - \rho_A^{(1+0)}(s) \right) = \int_0^\infty ds s \left(\rho_V^{(1)}(s) - \rho_A^{(1)}(s) \right) - 2m_\pi^2 f_\pi^2 = 0 ,$$

$$\int_0^\infty ds s \log(s/\mu^2) \left(\rho_V^{(1)}(s) - \rho_A^{(1)}(s) \right) = \frac{8\pi f_0^2}{3\alpha} \left(m_{\pi^\pm}^2 - m_{\pi^0}^2 \right) ,$$

where f_0 is the pion decay constant in the chiral limit, and α is the fine-structure constant. For the second WSR we assume that terms of order $m_i m_j$, $i, j = u, d$ can be neglected. Without this assumption, the integral is linearly divergent, forcing us to cut it off. If we cut off the integral at s_0 , there would be an extra contribution proportional to $m_i m_j \alpha_s^2 s_0$ in this sum rule. This contribution is still very small at $s_0 = m_\tau^2$ (of order a few percent of the contribution $2m_\pi^2 f_\pi^2$), allowing us to assume that we are effectively in the chiral limit with regard to the second WSR. Even the term $2m_\pi^2 f_\pi^2$, while dominating the term proportional to $m_i m_j \alpha_s^2 s_0$, vanishes in the chiral limit, and itself turns out to be numerically negligible within errors. Also the DGMLY sum rule holds only in the chiral limit, and in that limit the integral on the left-hand side is independent of μ because of the second WSR.

In Fig. 12 we show the first integral in Eq. (6.2) as a function of the “switch” point s_{SW} below which we use the experimental data, and above which we use the DV ansatz (2.10) with parameters from the CIPT fit with $s_{\text{min}} = 1.55 \text{ GeV}^2$ of Tab. 5 in order to evaluate the integral. Using parameter values from Tab. 4 or FOPT fits leads to almost identical figures.¹¹ The figure on the left includes the contribution from Eq. (2.10), while the figure on the right omits such contributions. The latter is equivalent to the upper right panel of Fig. 8 in the first paper in Ref. [3]. Clearly, the first WSR is very well satisfied by our fits, but only if duality violations are taken into account. We do not show similar figures for the second WSR and the DGMLY sum rule, because our conclusions for these sum rules are very similar. Just as in Ref. [12, 13], these sum rules are satisfied within errors, but only if duality violations are taken into account. In particular, within errors, one may assume that our representation of the spectral functions is in the chiral limit, for the purpose of these

¹¹ The contribution from OPE terms to the spectral functions $\rho_{V,A}$ is suppressed by an extra power of α_s , and small enough to be negligible [11, 12].

three sum rules.

B. The strong coupling

The presence of duality violations forces us to make several assumptions in order to extract a value for $\alpha_s(m_\tau^2)$. These assumptions have been checked against the data, *cf.* Figs. 5 and 8–12. First, we need to assume that Eq. (2.10) provides a satisfactory description of duality violations for asymptotically large s . Second, we need to assume that $s \gtrsim 1.5 \text{ GeV}^2$ is already in the asymptotic region. And, finally, if we wish to also use the axial data, we need to assume that this is true both in the V and A channels. As already discussed above, this would amount to the assumption that the upper shoulder of the a_1 resonance is already more or less in the asymptotic region. Using only the V -channel fits, we avoid having to make this latter assumption, and doing so we find, from the results quoted in Eq. (5.2),

$$\begin{aligned} \alpha_s(m_\tau^2) &= 0.296(10)(1)(2) = 0.296 \pm 0.010, & (\overline{\text{MS}}, n_f = 3, \text{FOPT}), \\ &= 0.310(14)(1)(1) = 0.310 \pm 0.014, & (\overline{\text{MS}}, n_f = 3, \text{CIPT}), \end{aligned} \quad (6.3)$$

where the first error is the statistical fit error already given in Eq. (5.2), while the second represents half the difference between the $s_{\min} = 1.55$ and 1.575 GeV^2 results of Tab. 3 from which the average is derived. The third error represents the change induced by varying the estimated 6-loop $D = 0$ coefficient $c_{51} = 283$ [15] by the assumed 100% uncertainty about its central value, as in Ref. [12, 13]. The error from this latter uncertainty would be about ± 0.004 for both FOPT and CIPT if it were estimated from fits using only the moment with weight \hat{w}_0 ; this would raise both final errors by 0.001. We observe that the final errors we find are of the same order of magnitude as the difference between the FOPT and CIPT values of $\alpha_s(m_\tau^2)$. We also note that in all tables the value of $\alpha_s(m_\tau^2)$ is very stable as a function of s_{\min} for all values of s_{\min} included in these tables, except for possibly the lowest s_{\min} shown.

Equation (6.3) constitutes our final result for $\alpha_s(m_\tau^2)$ from the revised ALEPH data. Converting these results into values for α_s at the Z mass using the standard self-consistent combination of 4-loop running with 3-loop matching at the flavor thresholds [39], we find

$$\begin{aligned} \alpha_s(m_Z^2) &= 0.1155 \pm 0.0014, & (\overline{\text{MS}}, n_f = 5, \text{FOPT}), \\ &= 0.1174 \pm 0.0019, & (\overline{\text{MS}}, n_f = 5, \text{CIPT}). \end{aligned} \quad (6.4)$$

C. Non-perturbative quantities

As in Ref. [13], we would like to estimate the relative deviation of the aggregate dimension-6 condensates $C_{6,V/A}$ from the values given by the VSA. We express these condensates in terms of the VSA-violating parameters ρ_1 and ρ_5 by [7]

$$C_{6,V/A} = \frac{32}{81} \pi \alpha_s(m_\tau^2) \langle \bar{q}q \rangle^2 \begin{pmatrix} 2\rho_1 - 9\rho_5 \\ 11\rho_1 \end{pmatrix}, \quad (6.5)$$

with VSA results for $C_{6,V/A}$ corresponding to $\rho_1 = \rho_5 = 1$. Using $\langle \bar{q}q(m_\tau^2) \rangle = (-272 \text{ MeV})^3$ [40], and the averages of the results for $C_{6,V}$ and $C_{6,A}$ from the $s_{\min} = 1.55$ and 1.575 GeV^2

fits of Tab. 5, we find¹²

$$\begin{aligned}\rho_1 &= -4 \pm 4, & \rho_5 &= 5.9 \pm 0.9 & (\text{FOPT}) , \\ \rho_1 &= -2 \pm 3, & \rho_5 &= 5.9 \pm 0.8 & (\text{CIPT}) .\end{aligned}\tag{6.6}$$

While no conclusion can be drawn about the accuracy of the VSA for ρ_1 , it is clear that the VSA is a poor approximation for ρ_5 . The value for ρ_5 is consistent with the one we found from OPAL data in Ref. [13].

It is conventional to characterize the size of non-perturbative contributions to the ratio $R_{V+A;ud} = R_{V;ud} + R_{A;ud}$ of the total non-strange hadronic decay width to the electron decay width, where $R_{V/A;ud}$ have been defined in Eq. (2.5), by the parametrization

$$R_{V+A;ud} = N_c S_{\text{EW}} |V_{ud}|^2 (1 + \delta_P + \delta_{NP}) ,\tag{6.7}$$

where δ_P stands for the perturbative, and δ_{NP} for the non-perturbative contributions beyond the parton model. If one knows δ_{NP} , the quantity δ_P , and hence $\alpha_s(m_\tau^2)$ can be determined from the experimental value of $R_{V+A;ud}$. In such an approach, the error on $\alpha_s(m_\tau^2)$ is thus directly correlated with that on δ_{NP} . As in Ref. [13], our fits give access to the values of δ_{NP} , as well as those of $\delta^{(6)}$, $\delta^{(8)}$, and δ^{DV} , the contributions to δ_{NP} from the $D = 6$ and $D = 8$ terms in the OPE as well as the DV term. From the $s_{\text{min}} = 1.55 \text{ GeV}^2$ fits of Tab. 5, we find

$$\begin{aligned}\delta^{(6)} &= 0.058 \pm 0.026 , & \delta^{(8)} &= -0.036 \pm 0.017 , \\ \delta_{\text{DV}} &= -0.0016 \pm 0.0011 & (\text{FOPT}) , \\ \delta^{(6)} &= 0.040 \pm 0.024 , & \delta^{(8)} &= -0.024 \pm 0.015 , \\ \delta_{\text{DV}} &= -0.0009 \pm 0.0009 & (\text{CIPT}) .\end{aligned}\tag{6.8}$$

The FOPT and CIPT estimates for these quantities are consistent with each other. There is a strong correlation between $\delta^{(6)}$ and $\delta^{(8)}$, about -0.97 in the FOPT case.

The values for δ^{NP} derived from these results are

$$\begin{aligned}\delta^{NP} &= 0.020 \pm 0.009 & (\text{FOPT}) , \\ \delta^{NP} &= 0.016 \pm 0.010 & (\text{CIPT}) ,\end{aligned}\tag{6.9}$$

which differ by 1.6σ , respectively, 1.2σ from the values found using the the OPAL data in Ref. [13]. With the value $R_{V+A;ud} = 3.475(11)$ quoted in Ref. [1], one finds $\delta_P \approx 0.18$, an order of magnitude larger than δ_{NP} , indicating that $R_{V+A;ud}$ is a dominantly perturbative quantity. However, as in Ref. [13], we find an error on δ_{NP} much larger than that reported by standard analyses in the literature, almost an order of magnitude so, for example, when compared to Ref. [1]. The result is that the error on $\alpha_s(m_\tau^2)$ is underestimated in the standard analysis; for further discussion, we again refer to Sec. VII below.

¹² We neglected the smaller errors on α_s and $\langle \bar{q}q \rangle$.

D. Comparison with the fits of Ref. [13] to OPAL data

A particularly interesting check is to look for consistency of the results from our fits to the ALEPH data with those we obtained by fitting the OPAL data in Ref. [13]. For the strong coupling, our results from OPAL data were

$$\begin{aligned}\alpha_s(m_\tau^2) &= 0.325 \pm 0.018, & (\overline{\text{MS}}, n_f = 3, \text{FOPT, OPAL, Ref. [13]}) , \\ &= 0.347 \pm 0.025, & (\overline{\text{MS}}, n_f = 3, \text{CIPT, OPAL, Ref. [13]}) .\end{aligned}\quad (6.10)$$

The values (6.3) we find from the ALEPH data are 1.4, respectively, 1.3 σ lower than the OPAL values, assuming that the errors on the ALEPH and OPAL values are independent. We also note that the fits in Ref. [13] were not entirely unambiguous; a choice about the preferred range for δ_V had to be made. The fact that the difference between our central ALEPH- and OPAL-based values, as well as that between our central CIPT- and FOPT-based results, is, in each case, comparable in size to the error obtained in any of these analyses supports the notion that any improvement in the precision with which $\alpha_s(m_\tau^2)$ can be determined from hadronic τ decays will require significant improvements to the data. Of course, this assumes that the fit *ansatz* employed is valid in the region of s_0 larger than about 1.5 GeV². We will return to this point in Sec. VII below, as well as in the Conclusion.

The coupling $\alpha_s(m_\tau^2)$ is, of course, not the only fit parameter. One may for instance compare the values of the OPE and DV parameters between Tab. 3 above and Tab. 4 of Ref. [13] for $s_{\min} \approx 1.5$ GeV², and conclude that they agree between the ALEPH and OPAL fits within (sometimes fairly large) errors. However, comparing Tab. 5 above with Tab. 5 of Ref. [13], one observes that the OPE and DV parameters for the axial channel agree less well between the ALEPH and OPAL fits. This may be an indication that it is safer to restrict our fits to the vector channel. Results for $\alpha_s(m_\tau^2)$ are, nevertheless, found to be consistent between pure- V and combined V and A fits, both in this article and in Ref. [13].

E. Final results for the strong coupling from ALEPH and OPAL data

To conclude this section, we present our best values for the strong coupling at the τ mass extracted from the ALEPH and OPAL data for hadronic τ decays, and based on the assumptions that underlie our analysis. The FOPT and CIPT averages, weighted according to the errors in Eqs. (6.3) and (6.10), are

$$\begin{aligned}\alpha_s(m_\tau^2) &= 0.303 \pm 0.009, & (\overline{\text{MS}}, n_f = 3, \text{FOPT, ALEPH \& OPAL}) , \\ &= 0.319 \pm 0.012, & (\overline{\text{MS}}, n_f = 3, \text{CIPT, ALEPH \& OPAL}) .\end{aligned}\quad (6.11)$$

These convert to the values

$$\begin{aligned}\alpha_s(m_Z^2) &= 0.1165 \pm 0.0012, & (\overline{\text{MS}}, n_f = 5, \text{FOPT, ALEPH \& OPAL}) , \\ &= 0.1185 \pm 0.0015, & (\overline{\text{MS}}, n_f = 5, \text{CIPT, ALEPH \& OPAL}) .\end{aligned}\quad (6.12)$$

VII. THE ANALYSIS OF REF. [1]

We now turn to a discussion of what we have referred to as the standard analysis, which was used in Refs. [1, 3, 4, 8], and is based on Ref. [9]. We begin with a brief overview of

what is done in this approach. One considers spectral moments with the weights

$$w_{k\ell}(x) = (1-x)^2(1+2x)(1-x)^k x^\ell, \quad (7.1)$$

$$x = s/s_0,$$

choosing $(k, \ell) \in \{(0, 0), (1, 0), (1, 1), (1, 2), (1, 3)\}$, and evaluating these moments at $s_0 = m_\tau^2$ only. Ignoring logarithms,¹³ terms in the OPE contribute to these weights up to $D = 16$. The five $s_0 = m_\tau^2$ moment values are, of course, insufficient to determine the eight OPE parameters $\alpha_s(m_\tau^2)$, $\langle \frac{\alpha_s}{\pi} GG \rangle$, C_6 , C_8 , C_{10} , C_{12} , C_{14} and C_{16} , so some truncation is necessary. The standard analysis approach to this problem is to assume the OPE coefficients $C_{D=2k}$ for $D > 8$ are small enough that they may all be safely neglected in all of the FESRs under consideration, despite numerical enhancements of their contributions via larger coefficients in some of the higher degree weights. Duality violations are, similarly, assumed to be small enough that $\Delta(s)$ in Eq. (2.8) can be ignored as well, at least for s_0 close to m_τ^2 . With these assumptions, the remaining OPE parameters $\alpha_s(m_\tau^2)$, $\langle \frac{\alpha_s}{\pi} GG \rangle$, C_6 and C_8 are fitted using the $s_0 = m_\tau^2$ values of the five $w_{k\ell}$ spectral moments noted above, for each of the channels V , A , and $V + A$. The central values and errors for $\alpha_s(m_\tau^2)$ are taken from the fits (FOPT and CIPT) to the $V + A$ channel, based on the VSA-motivated expectation of significant $D = 6$ cancellation and the hope of similar strong DV cancellations in the $V + A$ sum. However, as we have seen in Eq. (6.6), VSA is a rather poor approximation. Furthermore, the fact that the spectral function for the $V + A$ combination is flatter in the region between 2 and 3 GeV² than is the case for the V or A channels separately may mislead one into believing that DVs are already negligible at these scales for the $V + A$ combination. In actual fact, however, though somewhat reduced in the $V + A$ sum, DV oscillations are still evident in the ALEPH $V + A$ distribution. In addition, since we have a good representation of the individual V and A channels, we also have a good representation of their sum. The fact that our fits yield results for γ_A significantly larger than those for γ_V implies that the level of reduction of DV contributions in going from the separate V and A channels to the $V + A$ sum is accidental in the window between 2 and 3 GeV², and does not persist to higher s , where the stronger exponential damping in the A channel would drive the result for the $V + A$ sum towards that for the V channel alone.

These assumptions should be compared with those that have to be made in order to carry out the analysis presented in this article (as well as in the OPAL-based analyses of Refs. [12, 13]). DVs are unambiguously present in the spectral functions, as can be seen, for example, in the relevant panels of Figs. 5, 8, 9 and 10. In the standard analysis, the hope is that the double or triple pinching of the weights in Eq. (7.1) is sufficient to allow DVs to be ignored altogether, and indeed, for example Fig. 8, shows that pinching significantly reduces the role of DV contributions, especially near $s_0 = m_\tau^2$. However, if, as in the standard analysis, one restricts one's attention to $s_0 = m_\tau^2$, and wishes to employ only weights which are at least doubly pinched, the number of OPE parameters to be fit will necessarily exceed the number of weights employed, making additional assumptions, such as the truncation in dimension of the OPE described above, unavoidable.¹⁴ With the standard-analysis choice of the set of weights of Eq. (7.1), one finds that the OPE must be truncated at dimension $D = 8$ in order to leave at least one residual degree of freedom in the fits. In our analysis, in

¹³ Which appear in subleading terms in α_s at each order in the OPE.

¹⁴ For a detailed discussion of this point, see Ref. [12].

contrast, we choose not to ignore DVs *a priori*. This requires us to model their contribution to the spectral functions (as we did through Eq. (2.10)), and to use not just the single value $s_0 = m_\tau^2$, but rather a range of s_0 extending down from m_τ^2 . The one assumption we *do* have to make is that the *ansatz* (2.10) provides a sufficiently accurate description of DVs for values of s_0 between approximately 1.5 GeV² and m_τ^2 .

Clearly, whatever choice is made, it needs to be tested. For our analysis framework, we have presented detailed tests already above. In this section we consider primarily the standard analysis, most recently used in Ref. [1]. Our conclusion, from what follows below, is that the assumptions made in this framework do not hold up to quantitative scrutiny, and hence that the standard analysis approach should no longer be employed in future analyses.¹⁵

The results presented in Tab. 4 of Ref. [1] already indicate that there are problems with the standard analysis. Let us consider the values obtained for the gluon condensate, $\langle \frac{\alpha_s}{\pi} GG \rangle$, in the different channels, together with the χ^2 value for each fit (recall that for each of these fits there is only one degree of freedom):

$$\begin{aligned} \langle \frac{\alpha_s}{\pi} GG \rangle &= (-0.5 \pm 0.3) \times 10^{-2} \text{ GeV}^4, & \chi^2 &= 0.43 & V, \\ &(-3.4 \pm 0.4) \times 10^{-2} \text{ GeV}^4, & \chi^2 &= 3.4 & A, \\ &(-2.0 \pm 0.3) \times 10^{-2} \text{ GeV}^4, & \chi^2 &= 1.1 & V + A. \end{aligned} \quad (7.2)$$

The χ^2 values correspond to p -values of 51%, 7%, and 29%, respectively, indicating that all fits are acceptable. For these fits to be taken as meaningful, however, their results should satisfy known physical constraints. One such constraint is that there is only one effective gluon condensate, whose values should therefore come out the same in all of the V , A and $V + A$ channels. This is rather far from the case for the results quoted in Eq. (7.2), where, for example, the V and $V + A$ channel fit values differ very significantly. It is, moreover, problematic to accept the $V + A$ channel value and ignore the V channel one when the p -value of the V -channel fit is, in fact, larger than that of the $V + A$ channel.

There can be several reasons for the inconsistencies in the results of Ref. [1]. One possibility is that some of the weights (7.1) have theoretical problems already in perturbation theory, as argued in Ref. [31]. Another possibility is that the assumptions underlying the standard analysis do not hold. Whatever the reason, the discrepant gluon condensate values point to a serious problem with the standard analysis framework.¹⁶

We now turn to quantitative tests of the OPE fit results reported in Ref. [1]. We focus on the $V + A$ channel, where DVs and $D > 4$ OPE contributions were expected to play a reduced role, and on the CIPT $D = 0$ treatment, since this is the only case for which the OPE fit parameter values are quoted in Ref. [1]. The tests consist of comparing the weighted OPE and spectral integrals for the weights $w_{k\ell}$ employed in the analysis of Ref. [1], not just at $s_0 = m_\tau^2$, but over an interval of s_0 extending below m_τ^2 . If the assumptions

¹⁵ We point out that the inadequacy of the standard analysis framework was already demonstrated in Refs. [10–13], but it appears important to re-emphasize this point in view of the continued use of this framework in the literature, in particular in the updated analysis of Ref. [1].

¹⁶ This problem already existed in earlier ALEPH analyses [3, 4], but in principle it might have been due to the problem with the data itself. Note that OPAL enforced equality of the gluon condensate between various channels, and were able to obtain reasonable fits as judged by the χ^2 , possibly because of the larger data errors.

made about $D > 8$ OPE and DV contributions being negligible are valid at $s_0 = m_\tau^2$ they should also be valid in some interval below this point. A good match between the weighted spectral integrals and the corresponding OPE integrals, evaluated using the results for the OPE parameters quoted in Ref. [1], should thus be found over an interval of s_0 . If, on the other hand, these assumptions are not valid, then the fit parameter values will contain contaminations from DV contributions and/or contributions with higher D , both of which scale differently with s_0 than do the $D = 0, 4, 6$ and 8 contributions appearing in the truncated OPE form. Such contamination will show up as a disagreement between the s_0 dependence of the fitted OPE representations and the experimental spectral integrals.

It is worth expanding somewhat on this latter point since the agreement of the OPE and spectral integrals at $s_0 = m_\tau^2$ for the weights $w_{k\ell}$ employed in the standard analysis is sometimes mistakenly interpreted as suggesting the validity of the assumptions underlying the standard analysis at $s_0 = m_\tau^2$. However, while the agreement is certainly a necessary condition for the validity of these assumptions, it is not in general, a sufficient one. This caution is particularly relevant since four parameters are being fit using only five data points, making it relatively easy for the effects of neglected, but in fact non-negligible, higher- D and/or DV contributions to be absorbed, *at a fixed s_0* , into the values of the four fitted lower- D parameters. That this is a realistic possibility is demonstrated by the alternate set of OPE fit parameters obtained in the analysis of Ref. [10], which neglected DV contributions, but not OPE contributions with $D > 8$. The results of this fit, including non-zero C_D with $D > 8$ and an $\alpha_s(m_\tau^2)$ significantly different from that obtained via the standard analysis of the same data [4], produced equally good agreement between the $s_0 = m_\tau^2$ OPE and spectral integral results for all the $w_{k\ell}$ employed in the standard analysis fit of Ref. [4], conclusively demonstrating that this agreement does not establish the validity of the standard analysis assumptions. So long as one works at fixed $s_0 = m_\tau^2$, there is no way to determine whether the results of the standard analysis are, in fact, contaminated by neglected higher- D OPE and/or DV effects or not. One may, however, take advantage of the fact that different contributions to the theory sides of the various FESR scale differently with s_0 , with integrated DV contributions oscillatory in s_0 and integrated $D = 2k$ OPE contributions scaling as $1/s_0^k$. If the $D = 0, 4, 6$ and 8 parameters obtained from the fixed- $s_0 = m_\tau^2$ standard analysis fit have, in fact, absorbed the effects of $D > 8$ and/or DV contributions, the fact that the nominal lower- D s_0 -scaling does not properly match that of the higher- D and/or DV contaminations will be exposed when one considers the same FESR, with the same standard analysis OPE fit parameter values, at lower s_0 . A breakdown of the standard analysis assumptions will thus be demonstrated by a failure of the agreement of the OPE and spectral integrals observed at $s_0 = m_\tau^2$ to persist over a range of s_0 below m_τ^2 . Such s_0 -dependence tests represent important self-consistency checks for all FESR analyses.

Before carrying out these self-consistency tests on the results of the standard analysis, it is useful to make explicit the relative roles of the various different D contributions entering the $s_0 = m_\tau^2$ results for the $w_{k\ell}$ -weighted OPE integrals employed in the $V + A$ CIPT fit of Ref. [1]. For the $D = 0$ contributions, it is important to remember that the leading one-loop contribution is independent of both s_0 and α_s . It is thus the difference of the full $D = 0$ contribution and this leading term which determines the α_s dependence of the $D = 0$ contributions, and which is relevant to the determination of $\alpha_s(m_\tau^2)$. Tab. 6 shows the $s_0 = m_\tau^2$ results for (i) the α_s -dependent $D = 0$ contributions and (ii) the $D = 4, 6$, and 8 contributions corresponding to the CIPT fit results of Tab. 4 of Ref. [1], for each of the $w_{k\ell}$ employed in that analysis. The sum of the $D = 6$ and 8 contributions, which is $\sim 1 - 2\%$

of the α_s -dependent $D = 0$ contribution for w_{00} and w_{10} , is, in contrast, $\sim 10 - 25\%$ of the corresponding $D = 0$ contributions for the w_{11} , w_{12} and w_{13} cases. Furthermore, for w_{11} , the $D = 4$ contribution is essentially the same size as the α_s -dependent $D = 0$ one.

(k, ℓ)	α_s -dependent $D = 0$	$D = 4$	$D = 6$	$D = 8$
(0, 0)	0.005173	-0.000008	-0.000117	0.000033
(1, 0)	0.004399	-0.000361	-0.000117	0.000082
(1, 1)	0.000365	0.000350	-0.000039	-0.000049
(1, 2)	0.000208	0.000002	0.000039	-0.000016
(1, 3)	0.000081	0.000000	0.000000	0.000016

TABLE 6: The $D = 4, 6$ and 8 and α_s -dependent $D = 0$ contributions to the $s_0 = m_\tau^2$, $V + A$, $w_{k\ell}$ moments corresponding to the $V + A$ OPE fit parameter results of Tab. 4 of Ref. [1].

It is clear from these observations that it is the w_{11} , w_{12} and w_{13} moments which dominate the determinations of the $D = 4, 6$ and 8 OPE parameters in the analysis of Ref. [1]. Bearing in mind the very slow variation with s_0 of the $D = 0$ contributions to the dimensionless OPE integrals and the $1/s_0^k$ scaling of the $D = 2k$ contributions, it is, moreover, clear that the relative roles of the non-perturbative contributions will grow significantly relative to the α_s -dependent $D = 0$ ones as s_0 is decreased. Studying the s_0 dependence of the match of the OPE to the corresponding spectral integrals for the w_{11} , w_{12} and w_{13} spectral weights thus provides a particularly powerful test of the reliability of the values for the $D = 4, 6$ and 8 parameters obtained in the fits of Ref. [1].

The results of these tests are shown in Fig. 13. It is clear that the s_0 -dependence of the experimental spectral integrals and fitted OPE integrals is very different, demonstrating conclusively the unreliability of the $D = 4, 6$ and 8 fit parameter values obtained in Ref. [1]. Changes in the values of the $D = 6$ and 8 parameters, which enter the w_{00} FESR, would of course also force a change in the $\alpha_s(m_\tau^2)$ required to produce a match between the $s_0 = m_\tau^2$ w_{00} -weighted OPE and spectral integrals.

It is worth expanding somewhat on these observations for the w_{13} case, where the source of the problem with the fit of Ref. [1] becomes particularly obvious. Because of the x^3 factor present in $w_{13}(x)$, the $D = 2$ and 4 contributions to the OPE part are completely negligible numerically, leaving the standard analysis version of the w_{13} -weighted OPE integral entirely determined by the parameters $\alpha_s(m_\tau^2)$ and $C_{8,V+A}$. With the results and errors for these quantities from Tabs. 4 and 5 of Ref. [1], one finds that, as s_0 is decreased from m_τ^2 to *e.g.* 2 GeV^2 , the α_s -dependent $D = 0$ contribution *decreases* by $0.000001(0)$, while the $D = 8$ contribution *increases* by $0.000086(20)$. This is to be compared to the *increase* in the corresponding spectral integral, which is $0.000028(8)$. Evidently the disagreement between the w_{13} -weighted OPE and spectral integral results seen in Fig. 13 results from a problem with the fit value for $C_{8,V+A}$. Trying to fix the problem with the w_{13} FESR through a change in $C_{8,V+A}$ alone turns out to exacerbate the problem with the w_{12} FESR. Working backward, one finds that attempting to change $C_{4,V+A}$, $C_{6,V+A}$ and $C_{8,V+A}$ so as to improve the match between the s_0 dependences of the OPE and spectral integrals for the w_{11} , w_{12} and w_{13} FESRs without any change in $\alpha_s(m_\tau^2)$ produces changes in the $D \geq 4$ contributions to the w_{10} and w_{00} FESRs that can only be compensated for by a decrease in $\alpha_s(m_\tau^2)$. The

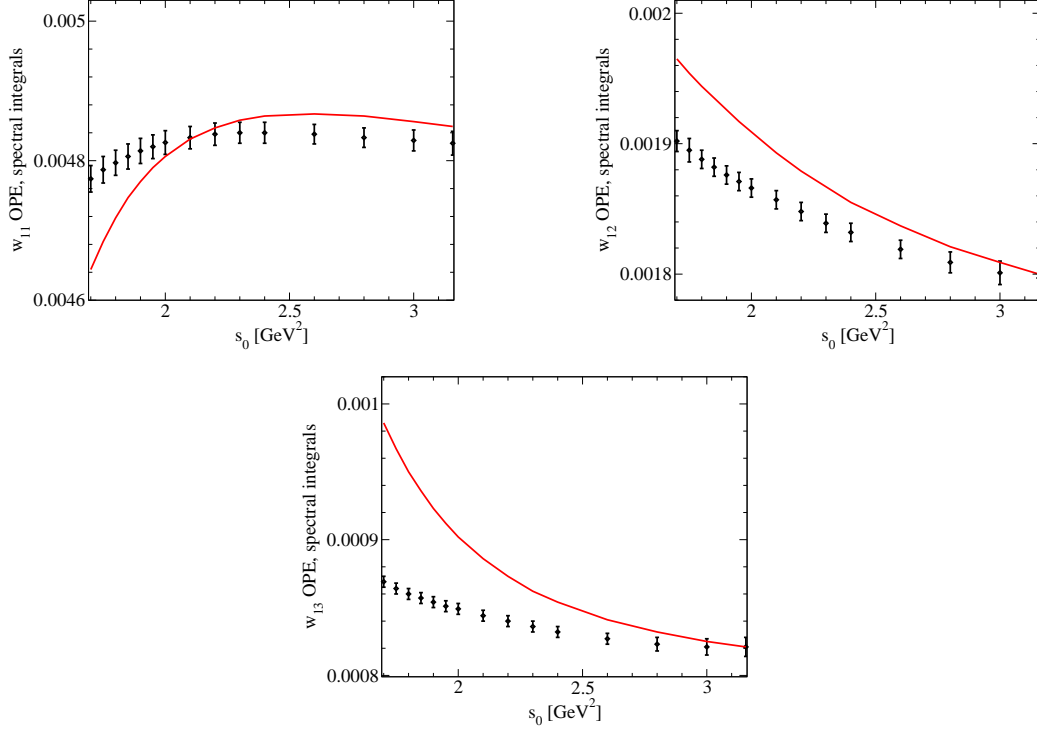


FIG. 13: Comparisons of the s_0 dependence of the $w_{k\ell} = w_{11}$, w_{12} and w_{13} $V + A$ spectral integrals to that of the corresponding OPE integrals evaluated employing as input the results of the CIPT fit for the OPE parameters from Tab. 4 of Ref. [1].

problem of the discrepancies between the s_0 -dependences of the OPE and spectral integrals in the w_{11} , w_{12} and w_{13} FESR parts of the standard analysis can thus not be resolved simply through shifts in $C_{4,V+A}$, $C_{6,V+A}$ and $C_{6,V+A}$ which leave the target of the analysis, namely the output $\alpha_s(m_\tau^2)$ value, unchanged.

A natural question, given the discussion above, is whether our approach produces a better match between experiment and theory for the higher spectral weights. The answer, as we will see below, is yes. Before embarking on this investigation, however, it is important to emphasize the non-optimal nature of the FESRs with weights w_{10} , w_{11} , w_{12} , and w_{13} . First, all of these weights contain a term linear in the variable x , a fact which, according to the arguments of Ref. [31], should make standard methods of estimating the uncertainty associated with truncating the integrated perturbative series for these weights much less reliable than is the case for the weights employed in our analysis. Second, the values of the C_D with $D > 8$ obtained from the fits reported in Ref. [10] were found to produce very strong cancellations amongst higher- D OPE contributions when employed in the higher (k, ℓ) $w_{k\ell}$ FESRs, making these FESRs particularly sensitive to any shortcomings in the treatment of higher- D OPE contributions, as well as a poor choice for use in attempting to fit the values of C_D with $D > 8$. The strong cancellation amongst higher- D OPE contributions for the higher- (k, ℓ) $w_{k\ell}$ moments turns out to be also a feature of the results of our extended analysis below, and hence not attributable to the neglect of DV contributions in Ref. [10]. Because of these strong cancellations, the use of the higher- (k, ℓ) $w_{k\ell}$ should be avoided

in future analyses, and we consider them below only for the sake of comparison with the results of the analysis of Ref. [1]. In making this comparison, we will focus on the CIPT resummation of perturbation theory, with the CIPT version of the standard analysis being the only one for which quantitative fit results are reported in Ref. [1].

To evaluate the OPE contributions to the w_{10} , w_{11} , w_{12} and w_{13} FESRs requires knowledge of five new quantities, $C_{4,V+A}$, $C_{10,V+A}$, $C_{12,V+A}$, $C_{14,V+A}$ and $C_{16,V+A}$, in addition to the OPE and DV parameters already obtained in our analysis. We estimate these using the $w(s) = (s/s_0)^{k-1}$ versions of the FESR Eq. (2.9), neglecting, as before, sub-leading contributions at each order $D > 2$ in the OPE. This yields, for $D = 2k > 2$,

$$(-1)^{k+1}C_{2k,V+A} = 2f_\pi^2 m_\pi^{2(k-1)} + \int_0^{s_0} ds s^{k-1} \rho_{V+A}^{(1)}(s) \\ + \int_{s_0}^\infty ds s^{k-1} \rho_{V+A}^{\text{DV}}(s) + \frac{1}{2\pi i} \oint_{|z|=s_0} dz z^{k-1} \Pi_{V+A}^{PT}(z), \quad (7.3)$$

where Π^{PT} is the perturbative contribution to $\Pi(z)$, corresponding to the $D = 0$ term in Eq. (2.6). The choices $k = 2, \dots, 8$ yield C_4, \dots, C_{16} , respectively. With $\alpha_s(m_\tau^2)$ and the V and A channel DV parameters from the $s_{\min} = 1.55 \text{ GeV}^2$ combined V and A CIPT fit of Tab. 5, we find, for the central values,

$$\begin{aligned} C_{4,V+A} &= 0.00268 \text{ GeV}^4, \\ C_{6,V+A} &= -0.0125 \text{ GeV}^6, \\ C_{8,V+A} &= 0.0349 \text{ GeV}^8, \\ C_{10,V+A} &= -0.0832 \text{ GeV}^{10}, \\ C_{12,V+A} &= 0.161 \text{ GeV}^{12}, \\ C_{14,V+A} &= -0.191 \text{ GeV}^{14}, \\ C_{16,V+A} &= -0.233 \text{ GeV}^{16}. \end{aligned} \quad (7.4)$$

For $C_{6,V+A}$ and $C_{8,V+A}$ the agreement with the values in Tab. 5 is excellent. With such values of the C_D , $D > 8$ contributions are far from negligible compared to the $D = 6$ and 8 ones for the $w_{k\ell}$ spectral weights with degree higher than three; the maximum scale, m_τ^2 , accessible in hadronic τ decays is not, it turns out, high enough to ensure that the OPE series is rapidly converging in dimension.

The theory parts $I_{\text{th}}^{(w_{k\ell})}(s_0)$ of the w_{10} , w_{11} , w_{12} and w_{13} FESRs produced by the results of Eq. (7.4) and Tab. 5 are compared to the corresponding spectral integrals in Fig. 14 as a function of s_0 . The agreement is obviously excellent, and far superior to that obtained from the standard analysis of Ref. [1]. This excellent agreement, over the whole range of s_0 shown, is completely destroyed if one removes the $D > 8$ contributions from the theory sides of the w_{10} , w_{11} , w_{12} and w_{13} FESRs. We emphasize again that the aim here is not a reliable determination of the OPE coefficients C_{4-16} , but a proof of existence of a set of values which, combined with our values for $\alpha_s(m_\tau^2)$ and the DV parameters, give an excellent representation of the s_0 dependence of the moments with the weights w_{10} , w_{11} , w_{12} and w_{13} (in addition, of course, to the weights included in our fits, in particular $w_{00} = \hat{w}_3$).

The problems demonstrated above with the standard analysis results of Ref. [1] could be a consequence of the neglect of non-negligible DVs, the breakdown of the assumption that $D > 8$ OPE contributions are negligible for all of the $w_{k\ell}$ employed, or both. In an attempt to clarify the situation, it is useful to consider a fit in which the potentially dangerous

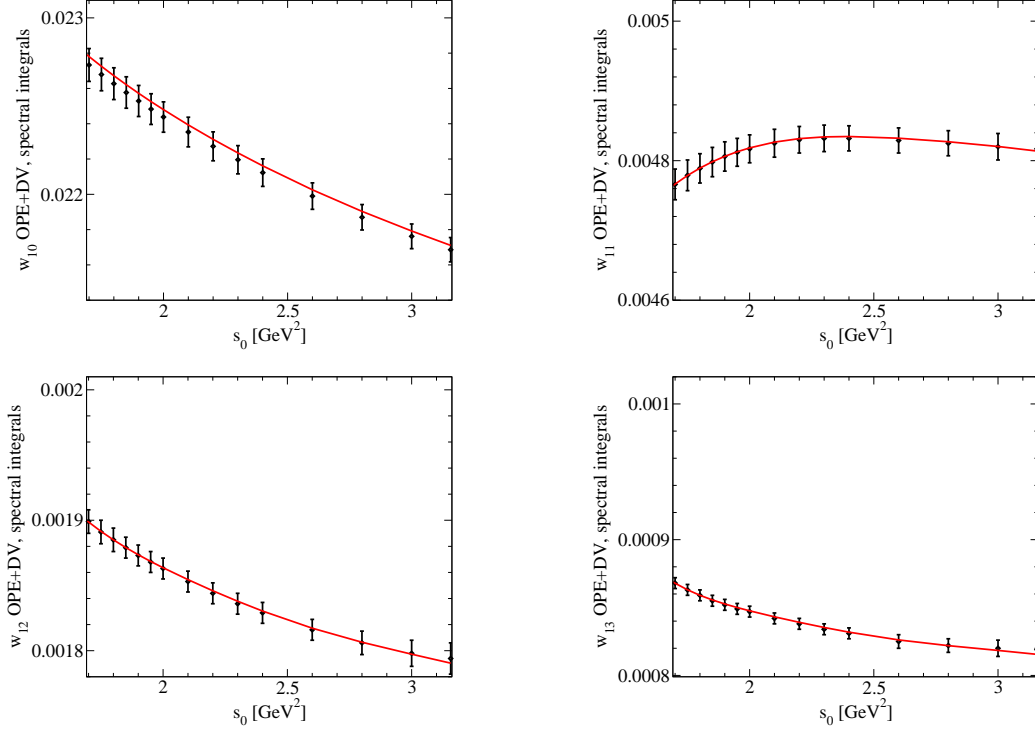


FIG. 14: Comparison of $I_{\text{th}}^{(w_{k\ell})}(s_0)$ (cf. Eq. (2.9) for $w_{k\ell} = w_{10}, w_{11}, w_{12}$ and w_{13} with $I_{\text{ex}}^{(w_{k\ell})}(s_0)$ for the $V + A$ channel using the results of the $s_{\min} = 1.55 \text{ GeV}^2$, combined V and A CIPT fit of Tab. 5 and Eq. (7.4). Top left panel: the w_{10} case; top right panel: the w_{11} case, bottom left panel: the w_{12} case; bottom right panel: the w_{13} case.

assumption about $D > 8$ OPE contributions is avoided. As an example, we consider a fit to the doubly pinched $\hat{w}_3 = w_{00}$ FESR in the $V + A$ channel ignoring DV contributions. Since the weight is doubly pinched, one expects DV contributions to be significantly suppressed, though the actual amount of suppression is not clear *a priori*. Since the OPE integrals still depend on three parameters, $\alpha_s(m_\tau^2)$, $C_{6,V+A}$ and $C_{8,V+A}$, it is, of course, necessary to consider the fit over a range of s_0 . To be specific, we focus on fits employing the FOPT resummation of perturbation theory. This exercise results in apparently perfectly acceptable fits, with p -values 10% and higher for $s_{\min} \geq 1.95 \text{ GeV}^2$. The fit quality drops dramatically as s_0 is lowered beyond this point, with p -values already at the 0.2% level for $s_{\min} = 1.90 \text{ GeV}^2$. The highest p -value, 57%, occurs for $s_{\min} = 2.2 \text{ GeV}^2$, and corresponds to

$$\begin{aligned} \alpha_s(m_\tau^2) &= 0.330 \pm 0.006, \\ C_{6,V+A} &= 0.0070 \pm 0.0022 \text{ GeV}^6, \\ C_{8,V+A} &= -0.0088 \pm 0.0042 \text{ GeV}^8. \end{aligned} \quad (7.5)$$

The quality of the resulting match between the fitted OPE and spectral integrals for $s_{\min} = 2.2 \text{ GeV}^2$, shown in the left panel of Fig. 15, is excellent. Despite this good quality match, the results of Eq. (7.5) are incomplete, in the sense that, in addition to the fit error induced by the covariances of the $V + A$ spectral data, there is an unspecified (and

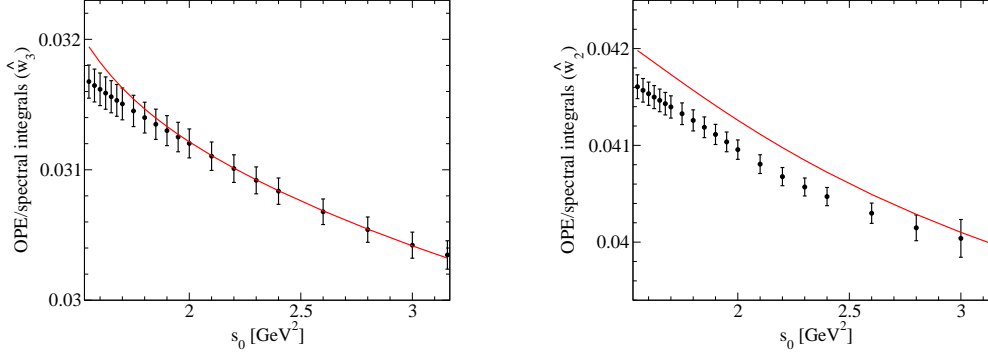


FIG. 15: Comparison of the $\hat{w}_3 = w_{00}$ -weighted spectral integrals (left panel) and \hat{w}_2 -weighted spectral integrals (right panel) with the corresponding OPE integrals evaluated using the results of the no-DV fit given in Eq. (7.5).

hence unquantified) systematic error associated with the neglect of DV contributions in the fit. Since the DV contribution to the FESR (2.9) involves the weighted integral of the DV component of the spectral function in the interval $s \geq s_0$, neglecting this systematic error would be reasonable if the $V + A$ spectral distribution showed no signs of DVs in the region $s > 2.2 \text{ GeV}^2$. This is, however, rather far from being the case, making the absence of an estimate for the residual systematic error associated with neglecting DV contributions problematic. One internally consistent way to test whether DV contributions are sufficiently small to be neglected for the \hat{w}_3 FESR is to demonstrate that they are already small for the singly pinched \hat{w}_2 FESR. Whether or not this is the case can be investigated by comparing the \hat{w}_2 -weighted OPE and spectral integrals, in the same s_0 range, using parameters obtained from the no-DV fit to \hat{w}_3 , Eq. (7.5). The results of this test are shown in Fig. 15 (right panel). The agreement between the OPE and spectral integrals is clearly not good, indicating the presence of significant DV contributions in the \hat{w}_2 FESR. This, together with the rapid deterioration of the \hat{w}_3 no-DV fit quality for $s_{\min} \leq 1.95 \text{ GeV}^2$, suggests that neglecting DV contributions to the \hat{w}_3 FESR is also dangerous.

The hope underlying existing FESR analyses which ignore DV effects is that the double pinching of the weight $w_{00} = \hat{w}_3$ is sufficient to make the residual DV contributions very small. While the arguments above make this possibility unlikely, it is still logically possible that, although DVs cannot be ignored in the singly-pinched \hat{w}_2 FESR, they can be ignored in the doubly-pinched \hat{w}_3 FESR. Let us therefore consider again the FOPT version of the \hat{w}_3 FESR in the $V + A$ channel, but now, rather than ignoring DVs, taking as external input the results for the DV parameters from the $s_{\min} = 1.55 \text{ GeV}^2$ FOPT fit of Tab. 5 and fitting the remaining OPE parameters $\alpha_s(m_\tau^2)$, $C_{6,V+A}$, and $C_{8,V+A}$, to the \hat{w}_3 weighted spectral integral in the $V + A$ channel in the presence of this estimate of the DV contributions. The results of this exercise, which are to be compared with Eq. (7.5), are

$$\begin{aligned} \alpha_s(m_\tau^2) &= 0.301 \pm 0.006 \pm 0.009, \\ C_{6,V+A} &= -0.0127 \pm 0.0020 \pm 0.0066 \text{ GeV}^6, \\ C_{8,V+A} &= 0.0399 \pm 0.0040 \pm 0.021 \text{ GeV}^8, \end{aligned} \tag{7.6}$$

where the first error is statistical and the second is that induced by the correlated uncertain-

ties of the external input DV parameters. The inclusion of the DV contributions induces a significant decrease in the value of $\alpha_s(m_\tau^2)$ and significant changes in the results for $C_{6,V+A}$ and $C_{8,V+A}$ (including changes in sign for both) as compared to the no-DV fit results of Eq. (7.5). The fit parameters are all changed in the direction of the results of the more detailed combined V and A fits discussed in Sec. V. This exercise clearly demonstrates that the effects of DVs on the parameters obtained from the $V + A \hat{w}_3$ FESR analysis are much larger than the nominal errors obtained on those parameters from the no-DV fit. This provides a further indication of the necessity of modeling DV effects in analyses attempting to extract $\alpha_s(m_\tau^2)$ from hadronic τ -decay data.

VIII. CONCLUSION

In this article, we reanalyzed the recently revised ALEPH data [1] for non-strange hadronic τ decays, with as primary goal the extraction of the strong coupling α_s at the scale m_τ . The rather low value of m_τ raises the question of to what extent the determination of a perturbative quantity like α_s in such an analysis might be “contaminated” by non-perturbative effects. Our specific aim was to take all known non-perturbative effects into account and arrive at a realistic estimate of the systematic error on the value of α_s extracted using hadronic τ data. This is important for three reasons. First, the value of α_s from τ decays, evolved to the Z mass, has long been claimed to be one of the most precise values available. Second, because the τ mass is so much smaller than other scales at which the strong coupling has been determined, $\alpha_s(m_\tau^2)$ provides a powerful test of the QCD running of the strong coupling, with the corresponding β function known to four-loop order. Finally, there continues to be some tension between the values of the $n_f = 5$ coupling $\alpha_s(M_Z^2)$ obtained from different sources. While lattice determinations involving analyses of small-size Wilson loops [41, 42], $c\bar{c}$ pseudoscalar correlators [43], the relevant combination of ghost and gluon two-point functions [44, 45], and employing the Schrödinger functional scheme [46] yield values, 0.1183(8) [41], 0.1192(11) [42], 0.1186(5) [43], 0.1196(11) [45], and 0.1205(20) [46], compatible both amongst one another and with the central value of the global electroweak fit result, $\alpha_s(M_Z^2) = 0.1196(30)$ [47], lower values have been obtained in a number of other analyses, *e.g.*, 0.1174(12) from lattice analyses of f_π/Λ_{QCD} [48], 0.1166(12) from an analysis of the static quark energy [49], 0.1118(17) from the recently revised JLQCD lattice determination from current-current two-point functions [50], and values in the range 0.1130 – 0.1160 from analyses of DIS data and shape observables in e^+e^- [51].

We have employed our analysis method previously [12, 13], using the OPAL data [8], but the revised ALEPH data have significantly smaller errors, and thus provide a more stringent test of our analysis method.

The fact that at such low scales non-perturbative effects are not negligible has of course been long known, and has been taken into account in the analysis of hadronic τ decays through the inclusion of higher-dimension condensate terms in the OPE. However, the experimental data are provided in the form of spectral functions, *i.e.*, as functions of $s = q^2$ with q denoting momentum in Minkowski space. Such values of q^2 , viewed as a complex variable, are outside the domain of validity of the OPE. While this is well known, it can also easily be inferred from the form of the vector spectral function in Fig. 4, which clearly shows oscillations that cannot be represented by the OPE. These oscillations lead unavoidably to the conclusion that violations of quark-hadron duality are, in general, significant at the scales accessible through experimental hadronic τ decay data.

It follows that in order to investigate the effect of duality violations on the extraction of α_s from τ decay data, they need to be taken into account. Unfortunately, a model is needed in order to parametrize the oscillations in the spectral functions, and this modeling necessitates making some assumptions on which to base the analysis. This is, however, true for any such analysis: the assumption that duality violations can be ignored in a given analysis amounts to assuming a model as well; in terms of the *ansatz* (2.10) it corresponds to taking the parameters $\delta_{V,A}$ to ∞ . We have, instead, assumed that this *ansatz* (with finite δ) provides a reasonable model of the resonance features present in the spectral functions for values of s in some region below m_τ^2 in which perturbation theory is still meaningful.¹⁷ As much as our aim is to find the most accurate value of $\alpha_s(m_\tau^2)$ possible given the data, an equally important goal was to test the validity of our approach, with the increased precision of the ALEPH data as compared to the OPAL data being particularly useful in this regard. This increased precision is, moreover, found to produce unique fit minima in the **hrothgar** studies of the multi-dimensional fit parameter space, improving the situation found for the corresponding fits to the OPAL data, and confirming that the precision of the ALEPH data is more than good enough to support fits incorporating an explicit representation of DV contributions.

Despite the recent resurgence of interest in this problem, triggered by the completion of the five-loop calculation of the Adler function in Ref. [14], very few investigations have carried out a complete analysis starting from the data. In essence, only two methods have been proposed through which to investigate non-perturbative effects, with the first being the method based on Refs. [7, 9], which was employed by Refs. [1, 3, 4, 8], and the second being the method we employed in this article, applying and extending ideas proposed in earlier work [10, 12, 13, 31]. In the absence of a detailed theoretical understanding of duality violations, it is important to test for the self-consistency of either analysis method using the data employed in the analysis.

In Sec. VII we demonstrated that the first method, used in Ref. [1], does not pass such tests. Indications supporting this conclusion have been published in earlier work, but now that the revised data are available, and in view of our critique in Sec. VII, we conclude that this method suffers from numerically significant systematic uncertainties not quantifiable within the analysis framework employed in Ref. [1], and hence must be discarded. The second method, employed in this article, does a much better job in fully describing the data, as we have shown in great detail in Secs. V, VI and VII above. However, there are some signs that also the limits of this method maybe in view. Fit qualities are typically larger than in the case of our analysis of the OPAL data [13], and a comparison of results based on ALEPH and OPAL data also shows some tension, even though errors are too large to say anything more conclusive. While these tensions may be caused by imperfections in the data (for instance slight discrepancies in the spectral function data visible in Fig. 4), it is by no means excluded that they point to shortcomings of the theory description as well.

We briefly reviewed, in Sec. II, why we consider the DV parametrization in Eq. (2.10) a physically sensible one. However, it remains relevant to test this form more quantitatively using experimental data. In this regard, we would like to stress that the exercise involving the x^N FESRs leading to the results of Eq. (7.4) represents a highly non-trivial test of this type. This follows from the fact that DV contributions to the x^N FESRs are generally not small, and oscillate with s_0 . The $D = 0$ OPE and DV contributions to the theory side

¹⁷ Up to the order considered [14].

of the x^N FESR for each N are, in this exercise, fixed by the results of the earlier fits involving the *ansatz* (2.10), leaving only a $D = 2N + 2$ OPE contribution controlled by C_{2N+2} to complete the theory side of the FESR. The different x^N considered provide very different weightings on the interval from s_0 to ∞ , and the different s_0 considered represent integration over different portions of the oscillations in the experimentally accessible region. Therefore, a problem with the DV *ansatz* would be expected to show up as an inability to successfully fit, with the single parameter C_{2N+2} , the s_0 -dependent difference between the experimental spectral integrals and the sum of the previously fixed $D = 0$ OPE and DV theory integral contributions. In fact, as we have seen, a set of C_{2N+2} exist which produce excellent matches to the experimental spectral integrals over a sizeable range of s_0 for all N ($N = 1, \dots, 7$) required to generate the results, shown in Fig. 14, for the weights $w_{k\ell}$ employed in Ref. [1]. The fact that the form (2.10) conforms to the qualitative features expected of the contribution representing the residual error of an asymptotic series, and the success of the detailed self-consistency tests just described, confirms that the *ansatz* (2.10) provides a good representation of DV effects in the channels of interest. Possible residual inaccuracies in this representation should, in any case, not be turned into an argument to not include DVs at all, since that strategy would lead to the presence of unquantifiable systematic errors which use of our *ansatz* strongly indicates are unlikely to be small.

It is interesting to compare the values of $\alpha_s(m_\tau^2)$ from the various analyses. First, the half differences between our ALEPH- and OPAL-based values are 0.015 (FOPT) and 0.019 (CIPT), while the average (between FOPT and CIPT) fit errors is about 0.012 for fits to ALEPH data (*cf.* Eq. (6.3)), and about double that for fits to OPAL data. Finally, the difference between the FOPT and CIPT values is 0.014 for the ALEPH-based values, and 0.022 for the OPAL-based values. These differences and errors are all comparable in size, and it appears reasonable to conclude that they reflect both the data and theory limitations on the accuracy with which $\alpha_s(m_\tau^2)$ can be obtained from analyses of hadronic τ decay, at least at present. We do not believe that it is meaningful to condense these results in the form of one central value and one aggregate error for $\alpha_s(m_\tau^2)$. Clearly, our ALEPH-based values are not in agreement with the value obtained in Ref. [1], despite using the same data. Averaging the values of Eq. (6.3) and adding half the difference between the two values as an error estimate for the CIPT/FOPT perturbative uncertainty, we would find a value $\alpha_s(m_\tau^2) = 0.303 \pm 0.014$, to be compared with the value 0.332 ± 0.012 quoted in Ref. [1]. It should be emphasized again that the error in the latter value does not include a component accounting for the systematic problems identified in Sec. VII.

One may ask whether one can do better. First, it would be interesting to apply our analysis method to data with better statistics, and such data are in principle available from the BaBar and Belle experiments. Such data would allow us to scrutinize our theoretical understanding in more detail and would, as can be seen from Fig. 4, be especially useful in the upper part of the spectrum. However, to date the analyses required to produce inclusive hadronic spectral functions from these data are not complete, and thus such an investigation must be postponed until they become available. Second, it would be nice to develop a deeper insight into the theory itself, or, lacking that, to develop new tools for testing any given model for duality violations. A recent idea in this direction based on functional analysis can be found in Ref. [52]. Finally, we note that the difference between the results for $\alpha_s(m_\tau^2)$ obtained using the FOPT and CIPT resummation schemes represents, at present, an important limitation on the accuracy with which α_s can be obtained at a scale as low as m_τ^2 ; further progress will require an improved understanding of this issue.

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