

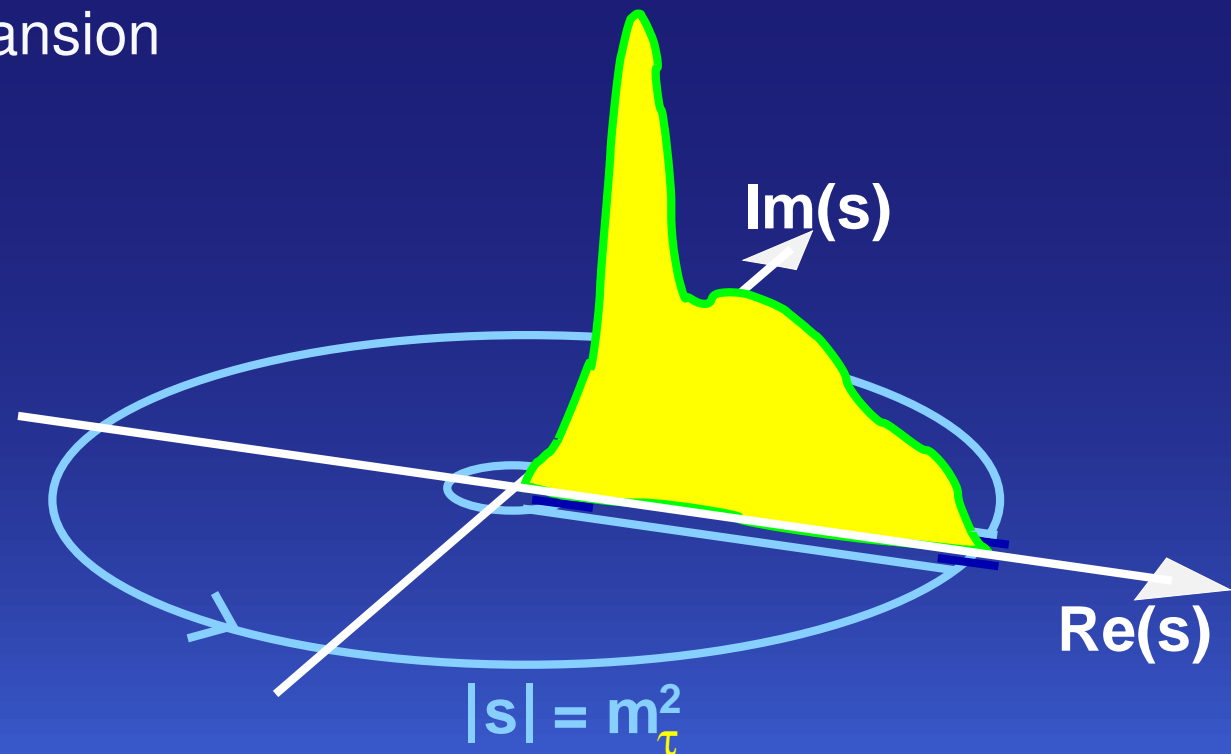
# FOPT and CIPT in $\tau$ Decays

$\alpha_s$  Workshop

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- ▶ Perturbative QCD Methods in  $\tau$  Decays
  - Fixed Order Perturbation Theory
  - Contour Improved Perturbation Theory
- ▶ Generalized FOPT
  - Modified Taylor Expansion
- ▶ Generalized RCPT
- ▶ Numerical Results
- ▶ Conclusions



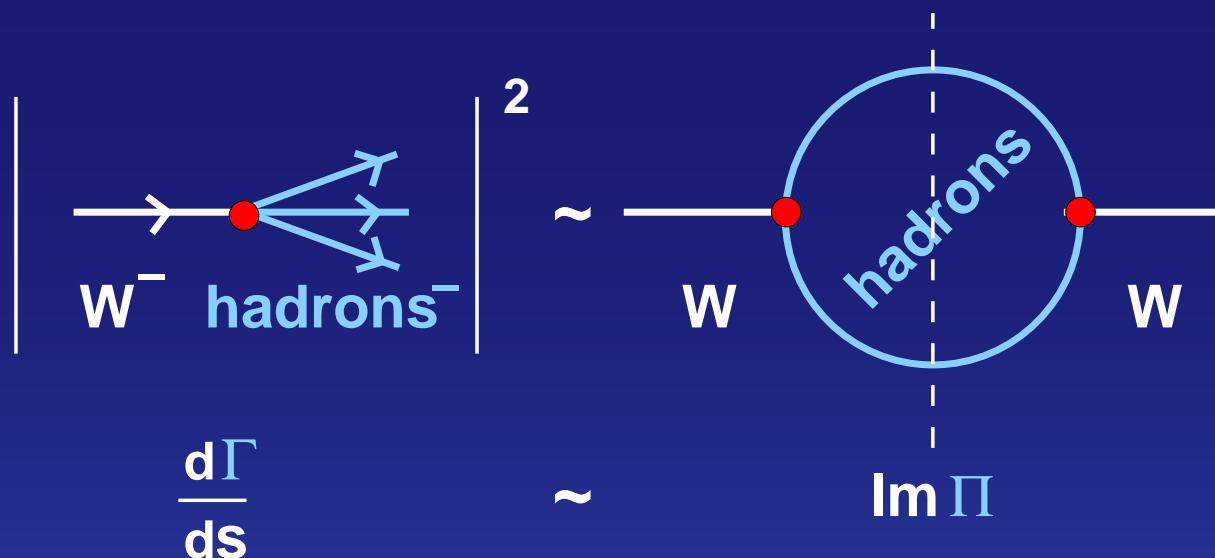
arXiv:0904.1796



# Perturbative QCD Methods in Hadronic $\tau$ Decays

► Use the optical theorem:

- The differential decay width  $d\Gamma/ds$  for the  $\tau$  going into hadrons is proportional to the imaginary part of the vacuum polarization  $\text{Im}\Pi$  (also called spectral function) of the  $W$  propagator
- this means it is enough to calculate the inclusive vacuum polarization of the  $W$  instead of every single final state of the hadronic  $\tau$  decays

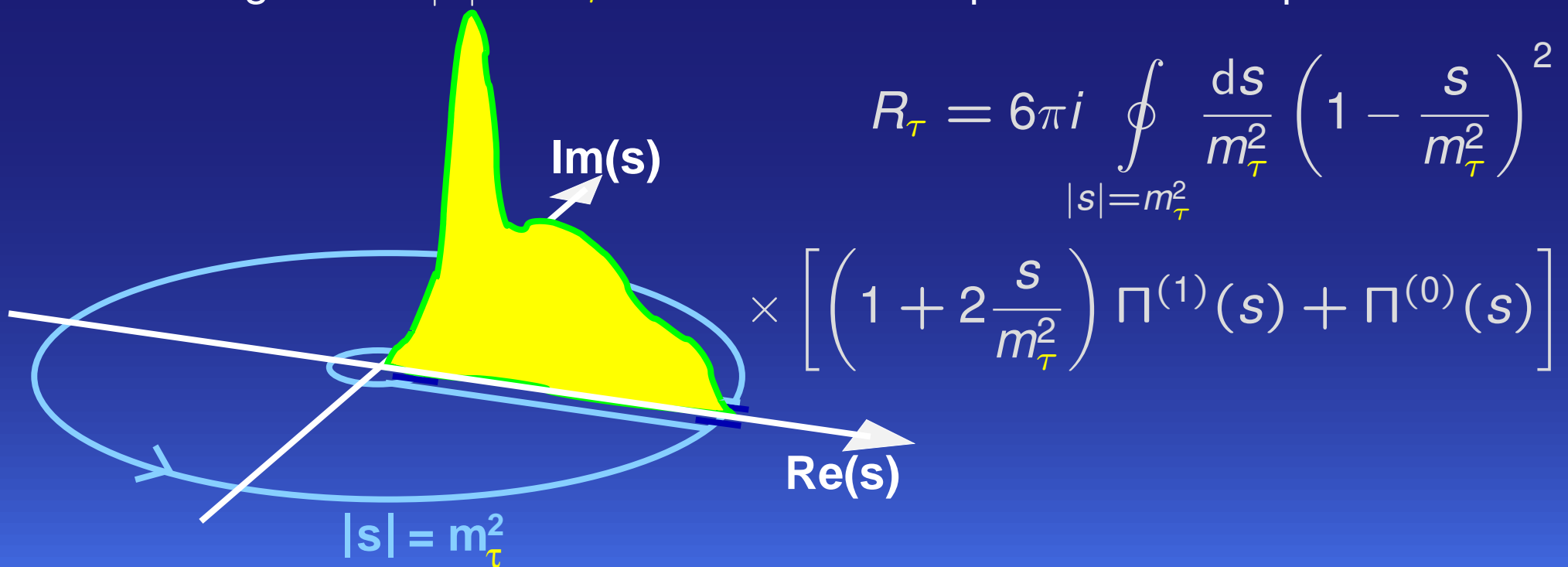


$$R_\tau = 12\pi \int_0^{m_\tau^2} \frac{ds}{m_\tau^2} \left(1 - \frac{s}{m_\tau^2}\right)^2 \left[ \left(1 + 2\frac{s}{m_\tau^2}\right) \text{Im}\Pi^{(1)}(s) + \text{Im}\Pi^{(0)}(s) \right]$$

# Perturbative QCD Methods in Hadronic $\tau$ Decays ► Cauchy

## ► Use the Cauchy theorem:

- since  $\Pi$  is analytic (except for the real positive  $s$ -axis where it might have poles) in the entire complex  $s$ -plane
- and due to the identity  $\text{Im}\Pi(s_+ + i\epsilon) = \frac{1}{2i}(\Pi(s_+ + i\epsilon) - \Pi(s_+ - i\epsilon))$
- the integral along the real positive  $s$ -axis can be expressed as a circular integral at  $|s| = m_\tau^2$
- this means all QCD calculations are done at a **well defined** rather high scale  $|s| = m_\tau^2$  over the vacuum polarization amplitude



## Perturbative QCD Methods in Hadronic $\tau$ Decays ► Adler-Function

- All the dynamics of the vacuum polarization amplitude is in its logarithmic derivative, the so-called Adler-function:  $D(s) = -s \frac{d\Pi(s)}{ds}$
- the Adler-function can be written as polynomial in the strong coupling constant  $\alpha_s$ :  $D(s) = \frac{1}{4\pi^2} \sum_{n=0}^{\infty} K_n \left( \frac{\alpha_s(-s)}{\pi} \right)^n$ , where the  $K_n$  are known up to  $n = 4$  and  $K_5$  has been estimated:

$$K_0 = K_1 = 1$$

$$K_2 = \frac{299}{24} - 9\zeta(3) = 1.63982 \dots$$

$$K_3 = \frac{58057}{288} - \frac{779}{4}\zeta(3) + \frac{75}{2}\zeta(5) = 6.37101 \dots$$

$$K_4 = \frac{78631453}{20736} + \frac{4185}{8}\zeta(3)^2 - \frac{1704247}{432}\zeta(3) \\ + \frac{34165}{96}\zeta(5) - \frac{1995}{16}\zeta(7) = 49.0757 \dots$$

$$K_5 \simeq 275, \quad \text{I'll use } 400 \pm 400$$

- convert the contour integral over  $\Pi(s)$  via partial integration in a contour integral over  $D(s)$

$$\begin{aligned} \oint_{|s|=m_\tau^2} ds g(s) \Pi(s) &= \oint_{|s|=m_\tau^2} \frac{ds}{s} \left[ G(s) - G(m_\tau^2) \right] (-s) \frac{d}{ds} \Pi(s) \\ &= \oint_{|s|=m_\tau^2} \frac{ds}{s} \left[ G(s) - G(m_\tau^2) \right] D(s) \end{aligned}$$

- where the polynomial  $g(s) = 6\pi i \frac{1}{m_\tau^2} \left(1 - \frac{s}{m_\tau^2}\right)^2 \left(1 + 2\frac{s}{m_\tau^2}\right)$
- has the antiderivative  $G(s) = 3\pi i \left(2\frac{s}{m_\tau^2} - 2\frac{s^3}{m_\tau^6} + \frac{s^4}{m_\tau^8}\right)$
- note that  $g(s)$  describes the  $J = 1$  part only since for massless quarks the scalar and pseudo-scalar parts vanish

$$R_\tau = 3 \sum_{n=0}^5 \frac{K_n}{2\pi i} \oint_{|s|=m_\tau^2} \frac{ds}{s} \left(1 - 2\frac{s}{m_\tau^2} + 2\frac{s^3}{m_\tau^6} - \frac{s^4}{m_\tau^8}\right) \left(\frac{\alpha_s(-s)}{\pi}\right)^n$$

- the perturbative descriptions of hadronic  $\tau$  decays start all with the same integral given on the previous slide
- they differ in the way the integral is calculated
- most interesting is the treatment of  $\alpha_s(-s)$  on the circle  $|s| = m_\tau^2$
- QCD does not tell us how large  $\alpha_s(\mu^2)$  at a given scale  $\mu$  is, but QCD does tell us what  $\alpha_s(\mu_1^2)$  at some scale  $\mu_1$  is if we know it at some other scale  $\mu_0$
- this prediction is made with the so-called  $\beta$ -function:

$$\beta(a_s) = \frac{da_s}{d \ln s} = -\beta_0 a_s^2 - \beta_1 a_s^3 - \beta_2 a_s^4 - \beta_3 a_s^5 + O(a_s^6), \text{ with } a_s = \frac{\alpha_s(\mu^2)}{4\pi}, \text{ and}$$

where the  $\beta_n$  are known up to  $n = 3$ :

$$\beta_0 = 11 - \frac{2}{3}n_f, \quad \beta_1 = 102 - \frac{38}{3}n_f, \quad \beta_2 = \frac{2857}{2} - \frac{5033}{18}n_f + \frac{325}{54}n_f^2,$$

$$\beta_3 = \frac{149753}{6} + 3564\zeta(3) - \left( \frac{1078361}{162} + \frac{6508}{27}\zeta(3) \right) n_f + \left( \frac{50065}{162} + \frac{6472}{81}\zeta(3) \right) n_f^2 + \frac{1093}{729}n_f^3.$$

- ▶ The Adler-function shown used  $K_n$  for a fixed choice of renormalization scale  $-\mu^2 = m_\tau^2$
- ▶ Physics should not depend on this choice
- ▶ The truncation of the perturbative series introduces a residual dependency:  $\mu^2 \frac{d}{d\mu^2} D(s, \mu^2) = 0 + O(\alpha_s^5(\mu^2))$
- ▶ This can be solved for each order with the  $\beta$ -function ( $\eta = \ln(-s/\mu^2)$ ):

$$K_0(\eta) = K_1(\eta) = K_0 = K_1 = 1$$

$$K_2(\eta) = K_2 - \frac{\beta_0}{4} \eta$$

$$K_3(\eta) = K_3 - \left( \frac{\beta_1}{16} + \frac{\beta_0}{2} K_2 \right) \eta + \frac{\beta_0^2}{16} \eta^2$$

$$K_4(\eta) = K_4 - \left( \frac{\beta_2}{64} + \frac{\beta_1}{8} K_2 + \frac{3\beta_0}{4} K_3 \right) \eta + \left( \frac{5\beta_1\beta_0}{128} + \frac{3\beta_0^2}{16} K_2 \right) \eta^2 - \frac{\beta_0^3}{64} \eta^3$$

$$K_5(\eta) = K_5 - \left( \frac{\beta_3}{256} + \frac{\beta_2}{32} K_2 + \frac{3\beta_1}{16} K_3 + \beta_0 K_4 \right) \eta + \left( \frac{3\beta_0^2}{8} K_3 + \frac{7\beta_1\beta_0}{64} K_2 + \frac{3\beta_2\beta_0}{256} + \frac{3\beta_1^2}{512} \right) \eta^2 + \left( \frac{\beta_0^3}{16} K_2 + \frac{13\beta_1\beta_0^2}{768} \right) \eta^3 + \frac{\beta_0^4}{256} \eta^4$$

## ► Fixed Order Perturbation Theory

- make Taylor-expansion of  $\alpha_s(-s)$  on the circle  $|s| = m_\tau^2$  around  $\alpha_s(m_\tau^2)$
- insert Taylor-expansion in the integral which becomes solvable in all orders
- order the result in powers of  $\alpha_s(m_\tau^2)$
- keep only the terms up to a fixed order in  $\alpha_s(m_\tau^2)$

$$\begin{aligned}
 R_\tau = & 3 \left( 1 + \frac{\alpha_s(m_\tau^2)}{\pi} + 5.2023 \dots \frac{\alpha_s^2(m_\tau^2)}{\pi^2} + 26.3659 \dots \frac{\alpha_s^3(m_\tau^2)}{\pi^3} \right. \\
 & + 127.0786 \dots \frac{\alpha_s^4(m_\tau^2)}{\pi^4} \\
 & \left. + (K_5 + 307.7825 \dots) \frac{\alpha_s^5(m_\tau^2)}{\pi^5} + O(\alpha_s^6) \right)
 \end{aligned}$$



## ► Contour Improved Perturbation Theory

- evolve  $\alpha_s(-s)$  in small steps numerically using the  $\beta$ -function on the circle  $|s| = m_\tau^2$
- insert the numerical values in the integral and solve it numerically too
- the result contains integrals over terms of the form

$$s^k \left[ \left( \frac{\alpha_s(m_\tau^2)}{\pi} \right)^l \ln^m \frac{-s}{m_\tau^2} \right]^n, \text{ with } k = 0, 1, 3, 4; n = 1, \dots, 5 \text{ (up to } K_5$$

for  $D(s)$ );  $l = 2, \dots, 5$  (up to  $\beta_3$  for the evolution of  $\alpha_s$ );  
 $m = 1, \dots, l - 1$

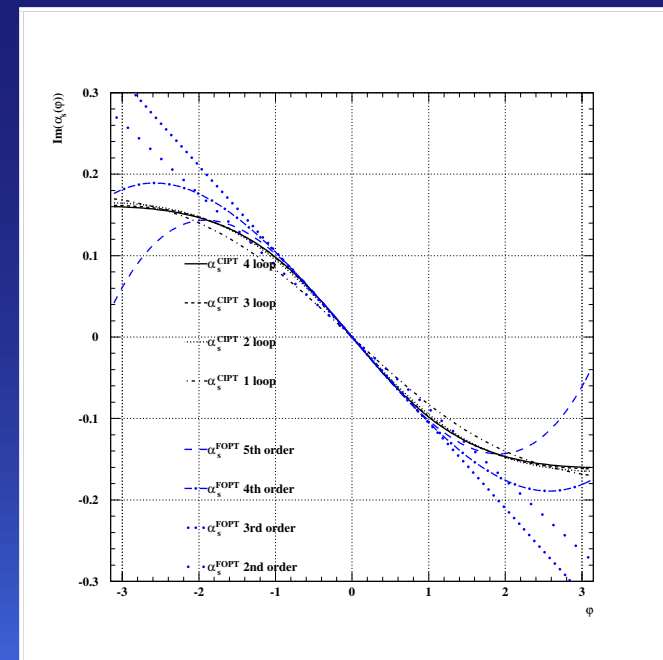
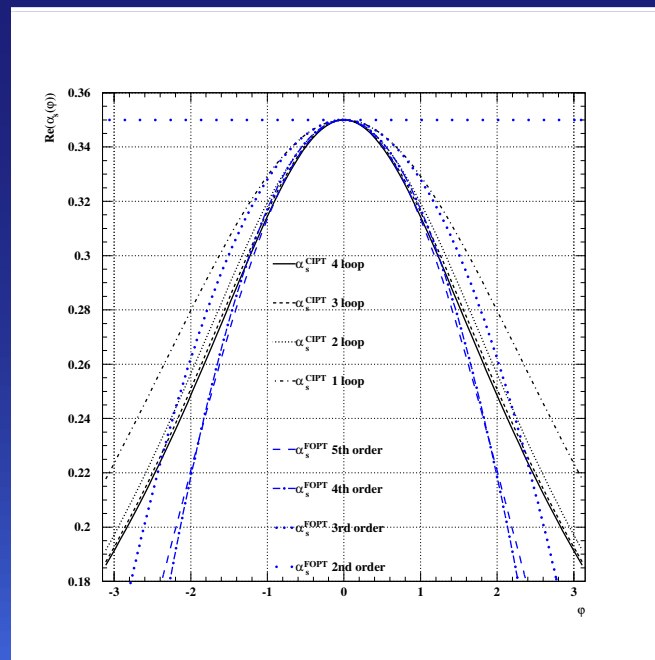
- the terms with total power in  $\alpha_s$  larger than the chosen order in FOPT are creating the **first** difference between CIPT and FOPT
- for up to the 5<sup>th</sup> order the neglected terms in FOPT are those with  $l \times n > 5$
- from the 50 logarithmic terms 39 are neglected in FOPT

- $\blacktriangleright$  let's have a closer look to the  $\alpha_s$  expansion used in FOPT
- $\blacktriangleright$  from the  $\beta$  function the following equation for  $\alpha_s(-s)$  as function of  $\alpha_s(s_0)$  follows:

$$\begin{aligned}
 \frac{\alpha_s(-s)}{\pi} &= \frac{\alpha_s(s_0)}{\pi} - \frac{1}{4} \beta_0 \ln \frac{-s}{s_0} \left( \frac{\alpha_s(s_0)}{\pi} \right)^2 + \\
 &\frac{1}{16} \left( \beta_0^2 \ln^2 \frac{-s}{s_0} - \beta_1 \ln \frac{-s}{s_0} \right) \left( \frac{\alpha_s(s_0)}{\pi} \right)^3 - \\
 &\frac{1}{128} \left( 2 \beta_0^3 \ln^3 \frac{-s}{s_0} - 5 \beta_0 \beta_1 \ln^2 \frac{-s}{s_0} + 2 \beta_2 \ln \frac{-s}{s_0} \right) \left( \frac{\alpha_s(s_0)}{\pi} \right)^4 + \\
 &\frac{1}{1536} \left( 6 \beta_0^4 \ln^4 \frac{-s}{s_0} - 26 \beta_0^2 \beta_1 \ln^3 \frac{-s}{s_0} + \right. \\
 &\quad \left. 9 \left( \beta_1^2 + 2 \beta_0 \beta_2 \right) \ln^2 \frac{-s}{s_0} - 6 \beta_3 \ln \frac{-s}{s_0} \right) \left( \frac{\alpha_s(s_0)}{\pi} \right)^5 + O(\alpha_s(s_0)^6)
 \end{aligned}$$

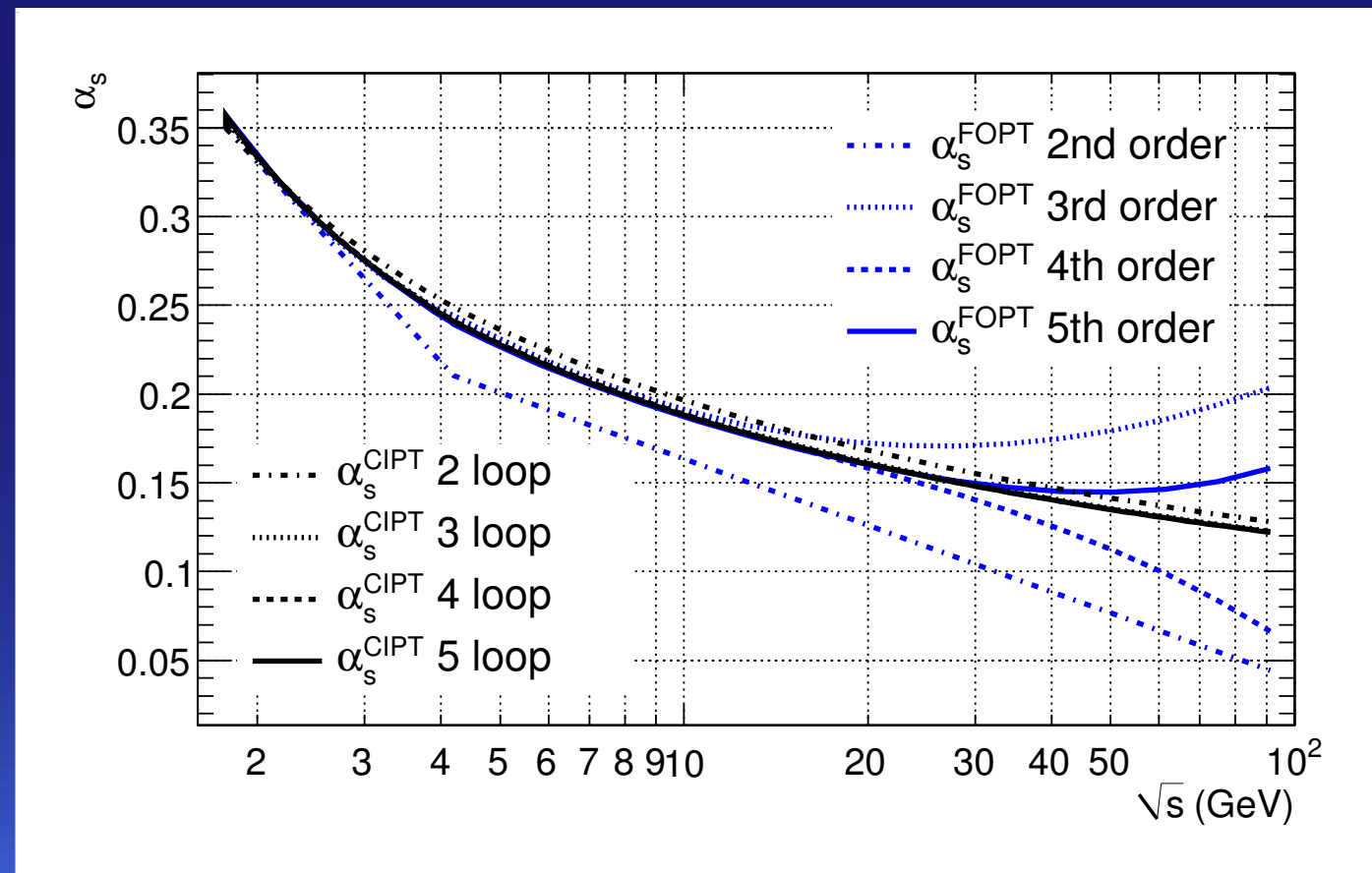
- $\blacktriangleright$  we can compare the numerically solved  $\alpha_s$  with the Taylor expansion from the prev. slide on the circle  $s = m_\tau^2$ .
- $\blacktriangleright$  note that we even keep all known orders (up to  $\alpha_s^5$ ) for the best
- $\blacktriangleright$  while in FOPT the term for  $\alpha_s^5$  enters with just first order (i.e. the constant value  $\alpha_s(s) = \alpha_s(s_0)$ )
- $\blacktriangleright$  the usage of the Taylor expansion over a wide distance from  $s$  to  $s_0$  creates the **second** difference between FOPT and CIPT

- $\blacktriangleright$  plots (real part: left, imaginary part: right) show Taylor-expanded  $\alpha_s(s = m_\tau^2 \exp(i\varphi))$  (blue) and numerical result (black) for various orders



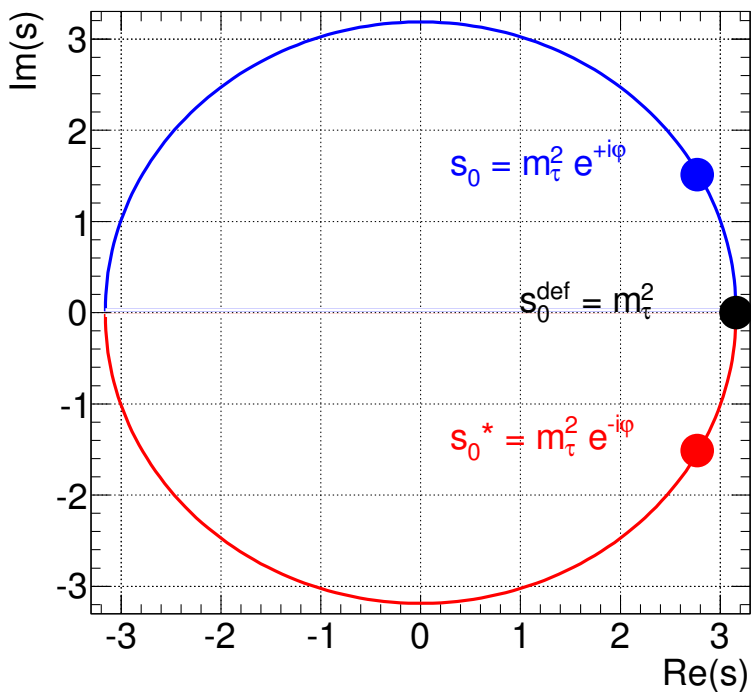
# Perturbative QCD Methods in Hadronic $\tau$ Decays $\blacktriangleright$ $\alpha_s$ expansion

- $\blacktriangleright$  another way of showing the differences between FOPT and CIPT is to apply the two methods to the evolution of  $\alpha_s$  from the  $\tau$  to the  $Z^0$
- $\blacktriangleright$  for CIPT 100 small logarithmic steps are used
- $\blacktriangleright$  for FOPT just 2 steps (at flavor thresholds) are used
- $\blacktriangleright$  beyond 2 loop the CIPT results become almost identical
- $\blacktriangleright$  but even at 5<sup>th</sup> order the deviation for FOPT is 30%
- $\blacktriangleright$  again the main difference comes not from higher orders but from the large step width



# Generalized FOPT

- ▶ How can the difference between FOPT and CIPT be absorbed into an uncertainty?
  - The Taylor expansion of  $\alpha_s$  on the circle  $|s| = m_\tau^2$  in FOPT is used from  $\alpha_s(m_\tau^2)$  to  $\alpha_s(-m_\tau^2 - i\epsilon)$  and from  $\alpha_s(m_\tau^2)$  to  $\alpha_s(-m_\tau^2 + i\epsilon)$



- The maximum angular distance on the circle is  $180^\circ$  on both semi-circles
- Splitting each semi-circle in 2 parts and developing  $\alpha_s$  around the splitting point ensures that again the angular distance is at most  $180^\circ$
- The splitting points should be symmetric on both semi-circles to keep the symmetry  $\alpha_s(s^*) = \alpha_s(s)^*$  which is important to keep integrals real
- If more and more equidistant splitting points would be added the result would converge towards CIPT

- ▶ Moving the splitting point on the semi-circles from  $\phi = 0^\circ$  (the usual FOPT) to  $\phi = 180^\circ$  (anti FOPT) shows the impact of this arbitrary choice

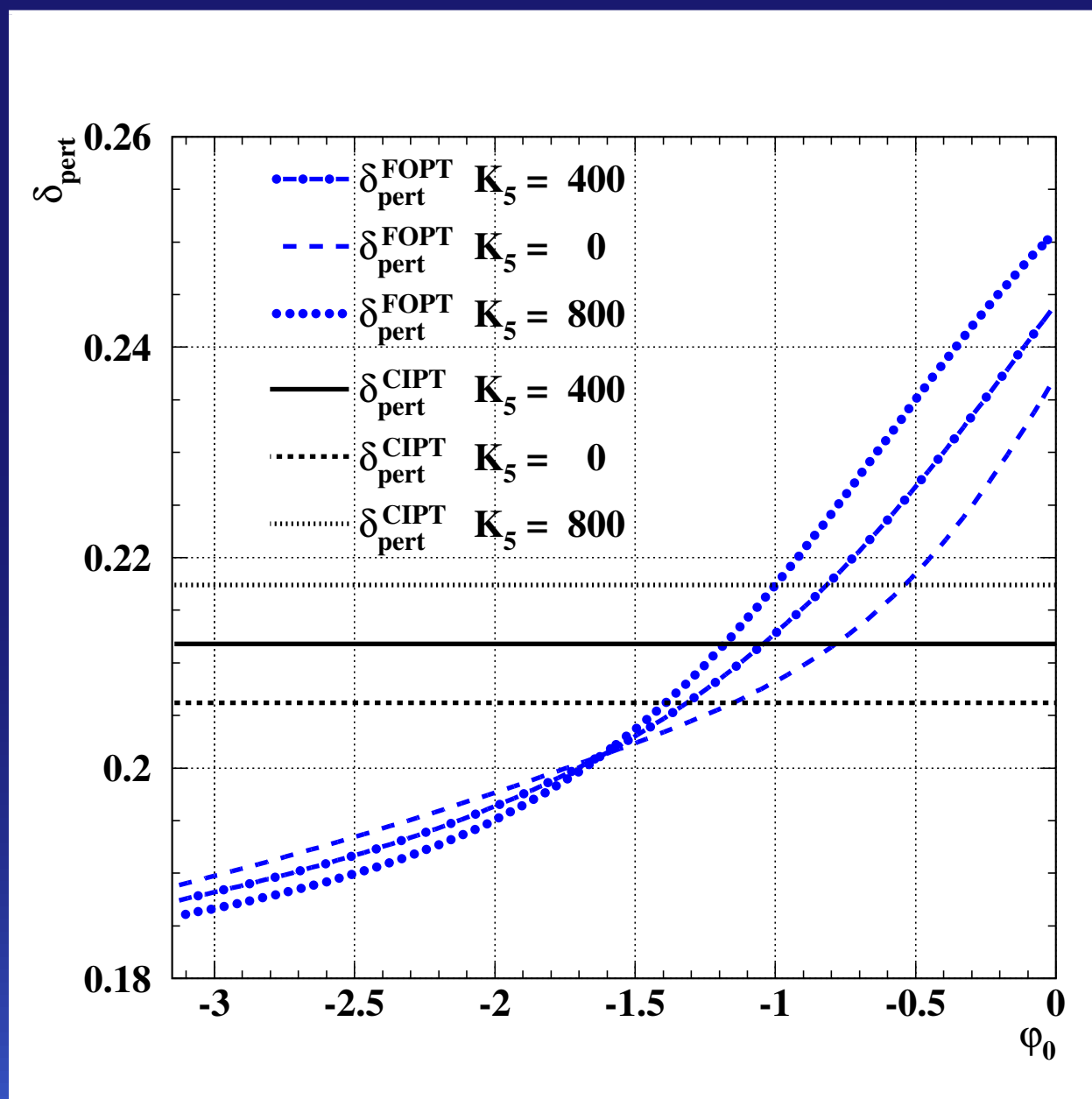
- ▶ The FOPT result around the new development points of  $\alpha_s(m_\tau^2 \exp(\pm i\varphi_0))/\pi = a \pm ib$  looks like this:

$$\begin{aligned}
 \delta_{\text{pert}} = & a + 0.8488 b + (-4.5 \varphi_0 + 0.8082) a b + \\
 & (1.9099 \varphi_0 + 5.2023) (a^2 - b^2) + \\
 & (-5.0625 \varphi_0^2 + 5.2138 \varphi_0 + 26.366) (a^3 - 3 a b^2) + \\
 & (4.2972 \varphi_0^2 + 27.410 \varphi_0 + 12.356) (b^3 - 3 a^2 b) + \\
 & (-9.6687 \varphi_0^3 - 101.51 \varphi_0^2 - 71.629 \varphi_0 + 127.08) (a^4 - 6 a^2 b^2 + b^4) + \\
 & (45.563 \varphi_0^3 - 100.94 \varphi_0^2 - 918.59 \varphi_0 - 521.11) (a^3 b - a b^3) + \\
 & (25.629 \varphi_0^4 - 92.897 \varphi_0^3 - 1220.5 \varphi_0^2 - 1272.5 \varphi_0 + \\
 & \quad K_5 + 307.78) (a^5 - 10 a^3 b^2 + 5 a b^4) + \\
 & (21.755 \varphi_0^4 + 324.78 \varphi_0^3 + 271.83 \varphi_0^2 - 1612.0 \varphi_0 + \\
 & \quad 0.8488 K_5 - 1413.5) (b^5 + 5 a^4 b - 10 a^2 b^3),
 \end{aligned}$$

- ▶  $\delta_{\text{pert}}$  remains real for all choices of  $\alpha_s$  and  $\varphi_0$
- ▶ the usual FOPT result is retained for  $\varphi_0 = 0$  and  $b = 0$
- ▶ re-inserting the Taylor expanded  $\alpha_s$  and Taylor expanding again also gives the usual FOPT result

# Generalized FOPT ► comparison to CIPT

- the plot shows  $\delta_{\text{pert}}$  for the generalized FOPT (blue) vs. CIPT (black) as function of  $\varphi_0$
- the 3 lines for each are different choices of  $K_5$  (0, 400, 800)
- CIPT remains constant for all choices of  $\varphi_0$
- variation for FOPT due to  $\varphi_0$  is 4 times larger than the variation due to  $K_5$



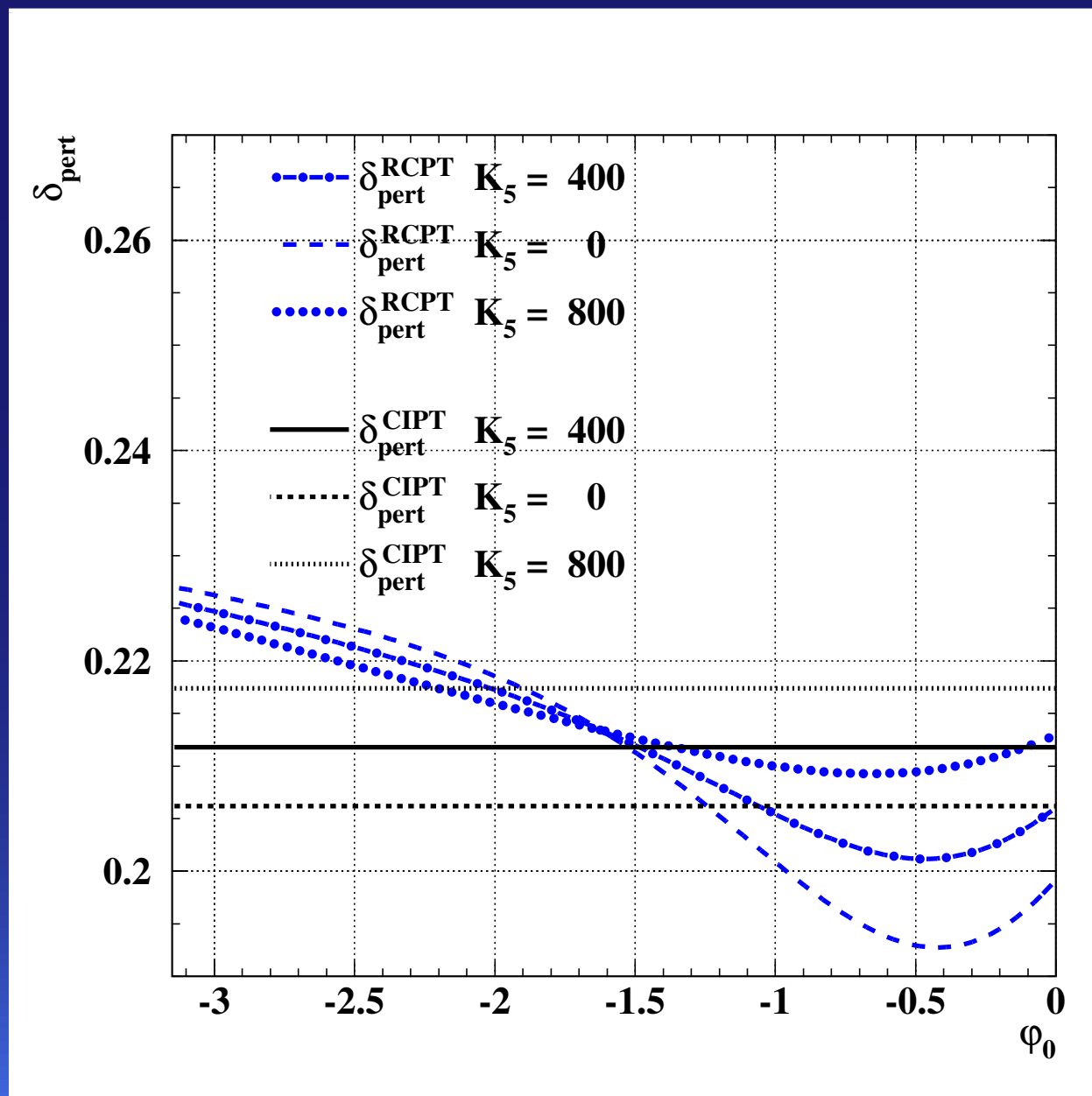
► For Renormalon Chain Perturbation Theory a similar generalization approach can be used

- usually the RCPT results are corrected up to the order used in FOPT with the known exact FOPT terms, while higher order terms are resummed to all orders in the large  $\beta_0$  approximation

►  $\delta_{\text{pert}}^{\text{RCPT}} =$

$$\delta_{\text{renormalon}} - \delta_{\text{large-}\beta_0}^{\text{FOPT}} + \delta_{\text{pert}}^{\text{FOPT}}$$

- in this equation the same arbitrary development point  $\varphi_0 = 0$  is usually used for  $-\delta_{\text{large-}\beta_0}^{\text{FOPT}} + \delta_{\text{pert}}^{\text{FOPT}}$
- therefore the same generalization idea to allow for other values of  $\varphi_0$  can be applied
- in addition I use 2-loop instead of 0-loop matching to go from the  $\overline{\text{MS}}$ -scheme to the V scheme
- the plot shows the comparison of generalized RCPT and CIPT like for generalized FOPT and CIPT on the slide before





# Numerical Results

- ▶ For numerical comparisons I'll use the  $\delta_{\text{pert}}$  from M. Davier et al. Eur.Phys.J.C56, 305 (2008), arXiv:0803.0979:

- ▶  $\delta_{\text{pert}} = 0.2042 \pm 0.0038_{\text{exp}} \pm 0.0033_{\text{non-pert}}$

$$\alpha_s^{\text{CIPT}}(m_\tau^2) = 0.3406 \pm 0.0047_{\text{exp}} \pm 0.0041_{\text{non-pert}} \pm 0.0066_{K_5}$$

$$\alpha_s^{\text{FOPT}}(m_\tau^2) = 0.3535 \pm 0.0061_{\text{exp}} \pm 0.0053_{\text{non-pert}} \pm 0.0208_{\varphi_0} \begin{matrix} +0.0005 \\ -0.0001 \end{matrix} K_5$$

$$\alpha_s^{\text{RCPT}}(m_\tau^2) = 0.3440 \pm 0.0030_{\text{exp}} \pm 0.0026_{\text{non-pert}} \pm 0.0061_{\varphi_0} \pm 0.0019_{K_5}$$

- ▶ All three results agree within errors due to the variation of  $\varphi_0$
- ▶ Averaging the RCPT and CIPT result only (since FOPT has a much larger uncertainty) yields:

$$\alpha_s(m_\tau^2) = 0.3423 \pm 0.005_{\text{exp}} \pm 0.007_{\Delta K_5} \pm 0.004_{\text{non-pert}} \begin{matrix} +0.005 \\ -0.001 \end{matrix} \mu$$

- ▶ the last error is due to the variation of the renormalization scale  $0.4 \leq \mu^2 / m_\tau^2 \leq 1.6$

- ▶ evolution to the  $Z^0$  gives finally:

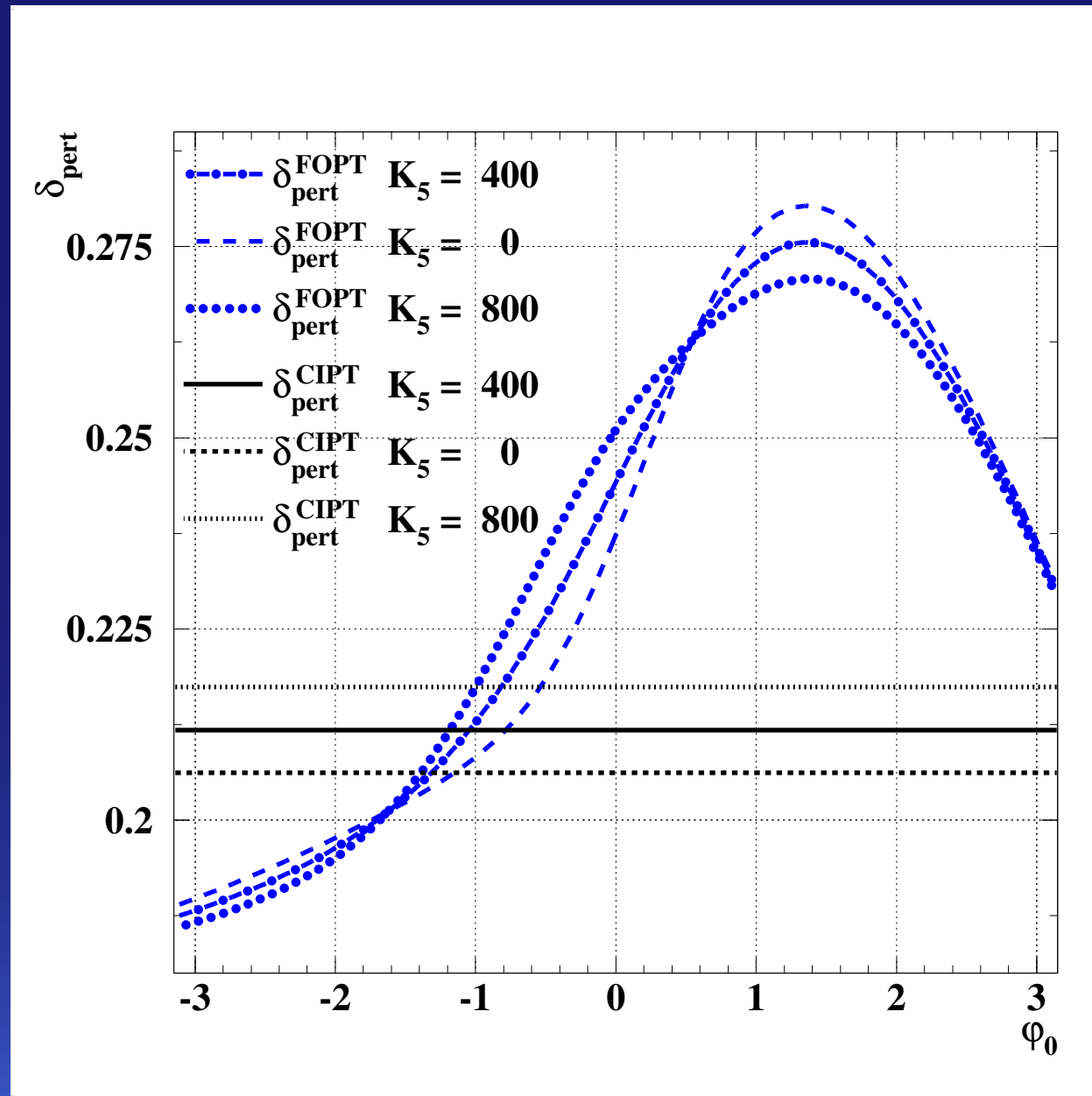
$$\alpha_s(m_{Z^0}^2) = 0.1213 \pm 0.0006_{\text{exp}} \pm 0.0008_{\Delta K_5} \pm 0.0004_{\text{non-pert}} \overset{+0.0005}{-0.0001}_{\mu} \pm 0.0002_{\text{ev}}$$

- ▶ where the last error is the evolution uncertainty due to the variation of the thresholds  $m_q < m_{\text{thresh}} < 2m_q$  and the quark masses itself within their respective errors

# Conclusions

- ▶ The differences between FOPT and CIPT can be mainly attributed to two reasons
  - truncation of the Taylor series of  $\alpha_s$
  - usage of the Taylor series over a long distance
- ▶ Uncertainties due to this can be made visible by allowing alternative development points of  $\alpha_s$  on the complex circle  $s_0 = m_\tau^2 \exp(i\varphi_0)$ 
  - these uncertainties are large:
    - ▶ about 4 times larger than uncertainties due to neglected higher orders
  - the usual chosen point  $s_0 = m_\tau^2$  sits at an extreme end making  $\alpha_s$  from FOPT appearing smaller than from CIPT
  - averaging over all possible choices gives a result much closer to CIPT – but still with larger error
- ▶ My proposal is to test FOPT type results always with the generalization procedure presented here in order to estimate the uncertainty and central value

- in some comments I received to the paper it was suggested to allow  $\varphi_0 > 0$  as well
- of course this is possible but one has to keep in mind that for all  $\varphi_0 > 0$  the angular distance in the Taylor expansion for  $\alpha_s$  is larger than  $180^\circ$  and thus larger than for default FOPT
- the plot shows the full range for possible  $\varphi_0$



- to illustrate the effect of more and more development points on the circle here is an example with 4 such points
- the points are arranged equidistantly on the semicircle and the angle  $\varphi_0$  is the angle relative to the quadrant
- the variation of  $\delta_{\text{pert}}$  with  $\varphi_0$  is already half of the value for 2 development points

