

# High precision determination of $w_0$

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for the BMW collaboration

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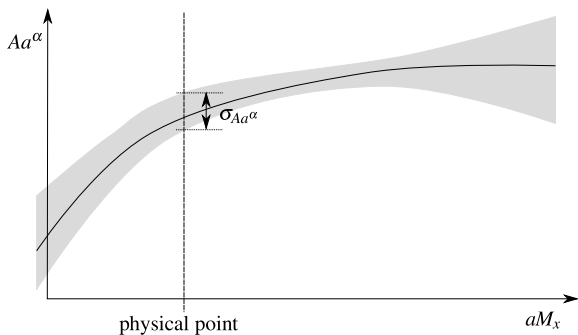


BMW  
collaboration

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# Scale setting

High precision scale setting is important for any high precision calculation on the lattice, especially for dimensionfull quantities.



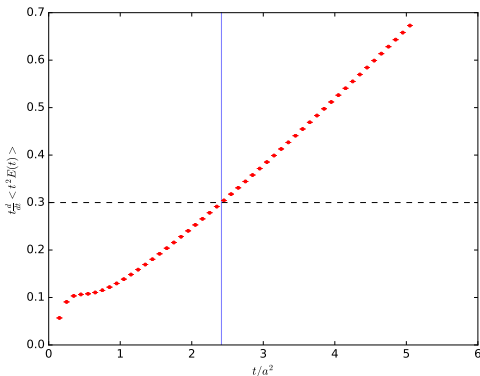
Scale setting uncertainty appears in several ways:

- Definition of the physical point
- Translation of the result in MeV.
- ...

[plot: L. Varnhorst]

$w_0$ 

Very promising for high precision scale setting:  $w_0$ .



[plot: L. Varnhorst]

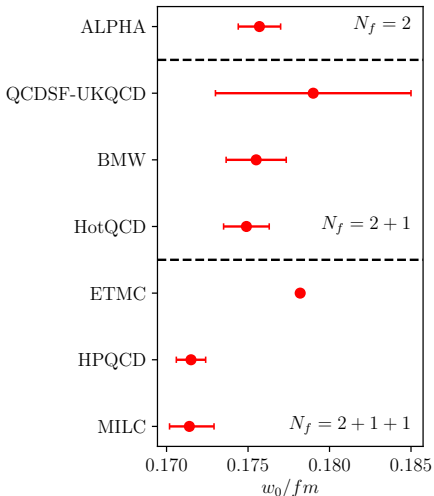
- Apply Wilson flow to “smooth out” the gauge fields and bring them closer to the classical solution. [Luscher:2010iy]
- Monitor the action density.
- Define  $w_0$  to be the value of  $\sqrt{t}$  where  $t \frac{d}{dt} \langle t^2 E(t) \rangle = 0.3$ . [Borsanyi:2012zs]
- Closely related to  $t_0$ . [Luscher:2010iy]

# Recent results

$w_0$  can not be determined experimentally.  $\Rightarrow$  It has to be determined once on the lattice.

In this talk I will present an ongoing effort to determine  $w_0$  with high precision in a blind analysis.

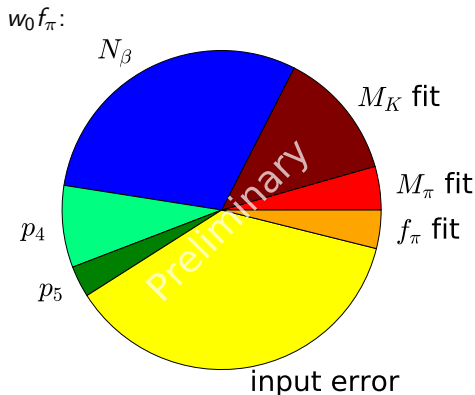
[ALPHA] M. Bruno *et al.* [ALPHA Collaboration], PoS LATTICE **2013** (2014) 321 [arXiv:1311.5585 [hep-lat]].  
 [QCDSF-UKQCD] R. Horsley *et al.* [QCDSF-UKQCD Collaboration], PoS LATTICE **2013** (2014) 249 [arXiv:1311.5010 [hep-lat]].  
 [BMW] S. Borsanyi *et al.*, "High-precision scale setting in lattice QCD," JHEP **1209** (2012) 010 [arXiv:1203.4469 [hep-lat]].  
 [HotQCD] A. Bazavov *et al.* [HotQCD Collaboration], Phys. Rev. D **90** (2014) 094503 [arXiv:1407.6387 [hep-lat]].  
 [ETMC] A. Deuzeman and U. Wenger, PoS LATTICE **2012** (2012) 162.  
 [HPQCD] R. J. Dowdall, C. T. H. Davies, G. P. Lepage and C. McNeile, Phys. Rev. D **88** (2013) 074504 [arXiv:1303.1670 [hep-lat]].  
 [MILC] A. Bazavov *et al.* [MILC Collaboration], Phys. Rev. D **93** (2016) no.9, 094510 [arXiv:1503.02769 [hep-lat]].



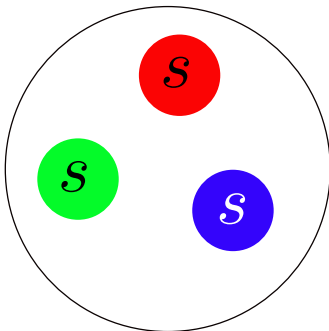
(Plot based on [MILC])

# Determination from $w_0 f_\pi$

- presented at LATTICE2017 by L. Varnhorst
- determined on 2+1+1 ensembles
- accuracy  $\approx 6\%$
- 2 challenges:
  - Input error of  $f_\pi$
  - What about QED?
- determination from  $w_0 M_\Omega$



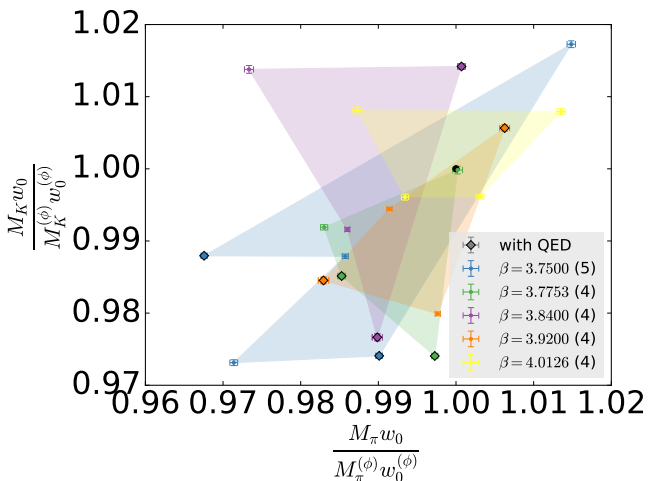
# $\Omega$ -mugshot



- hadron type: baryon
- valence quarks: sss
- charge: -
- mass:  $(1672.45 \pm 0.29)$  MeV,  $\Delta_{m_\Omega} = 0.17\text{‰}$  ( $\Delta_{f_\pi} = 1.5\text{‰}$ )

[PDG 2018]

# Lattice ensembles



The ensembles are generated with a staggered fermion action on stout-smearred gauge configurations. The gauge action is tree-level Symanzik improved.



## QED-Fit

$$w_0 M_\Omega = A + BM_{\pi^0}^2 w_0^2 + CM_{K_x}^2 w_0^2 + Ee_v^2 + Fe_v e_s + Ge_s^2$$

$$M_{K_x}^2 = M_{K^+}^2 + M_{K^0}^2 - M_{\pi^+}^2$$

$$A = A_0 + A_1 a^2$$

$$B = B_0 + B_1 a^2$$

$$C = C_0 + C_1 a^2$$

$$E = E_0 + E_1 a^2$$

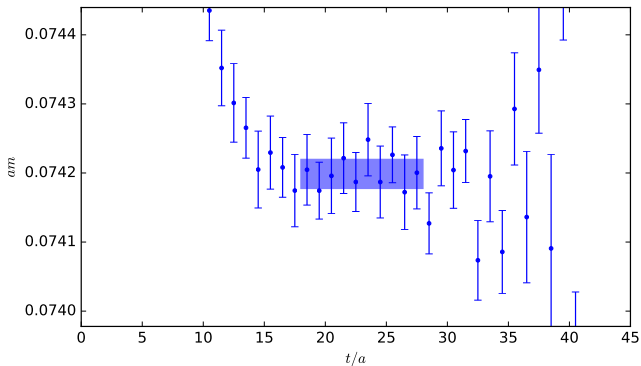
$$F = F_0$$

$$G = G_0$$

The scale is set by  $w_0$ . To determine the final result, the equation is solved for  $w_0$  in the continuum, at the physical point. In this way the determination is self consistent.

# Mass determination

$M_\pi$ ,  $M_K$  and  $M_\Omega$  are extracted by fitting appropriate correlators in two different fit ranges each.



$M_\Omega$  is also determined by solving the generalized eigen value problem (GEVP)  $\rightarrow$  talk by L. Varnhorst, 20.06.2019, 17:10

# QED-effects

$$w_0 M_\Omega = A + B M_{\pi^0}^2 w_0^2 + C M_{K_x}^2 w_0^2 + E e_V^2 + F e_V e_s + G e_s^2$$

$$[w_0 M_\Omega]_{\text{iso}} = A + B [M_{\pi^0}^2 w_0^2]_{\text{iso}} + C [M_{K_x}^2 w_0^2]_{\text{iso}}$$

$$\frac{\partial^2}{\partial e_V^2} [w_0 M_\Omega] = B \frac{\partial^2}{\partial e_V^2} [M_{\pi^0}^2 w_0^2] + C \frac{\partial^2}{\partial e_V^2} [M_{K_x}^2 w_0^2] + E$$

$$\frac{\partial}{\partial e_s e_V} [w_0 M_\Omega]_{q11} = B \frac{\partial}{\partial e_s e_V} [M_{\pi^0}^2 w_0^2] + C \frac{\partial}{\partial e_s e_V} [M_{K_x}^2 w_0^2] + F$$

$$\frac{\partial^2}{\partial e_s^2} [w_0 M_\Omega] = B \frac{\partial^2}{\partial e_s^2} [M_{\pi^0}^2 w_0^2] + C \frac{\partial^2}{\partial e_s^2} [M_{K_x}^2 w_0^2] + G$$

Perform a combined, fully correlated fit to all components, which takes x- and y-Errors into account

For more details on QED effects → talk by B. Toth, 18.06.2019, 17:10

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# Systematic error estimation

$$w_0 M_\Omega = A + BM_{\pi^0}^2 w_0^2 + CM_{K^x}^2 w_0^2 + Ee_v^2 + Fe_v e_s + Ge_s^2$$

$$M_{K^x}^2 = m_{K^+}^2 + M_{K^0}^2 - M_{\pi^+}^2$$

$$A = A_0 + A_1 a^2$$

$$B = B_0 + B_1 a^2$$

$$C = C_0 + C_1 a^2$$

$$E = E_0 + E_1 a^2$$

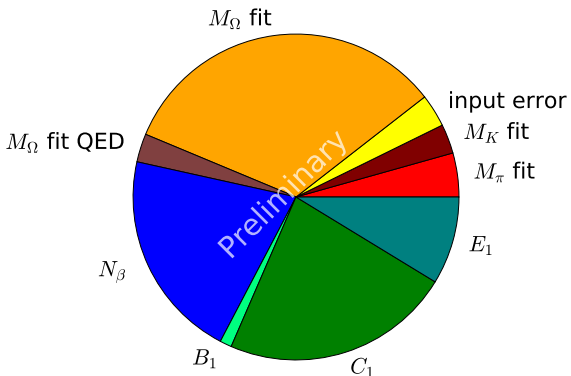
$$F = F_0$$

$$G = G_0$$

$$\beta = 3.7500, 3.7753, 3.8400, 3.9200, 4.0126$$

Everything colored in red is varied. If there are only three  $\beta$ s,  $E_1$  is excluded. Also the different results of the omega mass fits are varied. There is a total of 1920 fits.

# Final systematic error budget

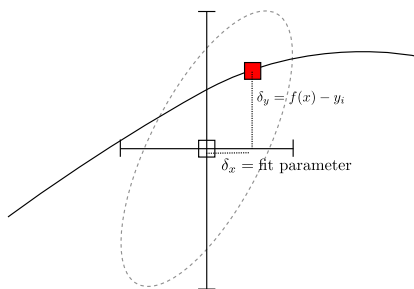


$$w_0 M_\Omega = A + BM_{\pi^0}^2 w_0^2 + CM_{K_X}^2 w_0^2 + Ee_v^2 + Fe_v e_s + Ge_s^2$$

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# Correlation and $\chi$ -errors

The statistic error is estimated by the Jackknife method.



[plot: L.Varnhorst]

For one quantity:

$$\vec{\delta} = \begin{pmatrix} f(x_i) - y_i \\ \delta_{x,i} \end{pmatrix} \quad \chi^2 = \sum_i \vec{\delta}^T C^{-1} \vec{\delta}$$

where  $C$  is the covariance matrix.

When generalized to several quantities, the number of fit parameters and the size of  $C$  increases. This can be numerically difficult.



# Statistic error estimation

Computationally easier, but leading to the same minimum:

Minimizing:

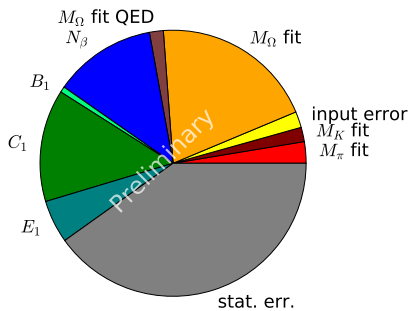
$$\chi^2 = \sum_i \frac{y_i^{(k)} - f^{(k)}(x_i)}{\sigma(y_i^{(k)} - f^{(k)}(x_i))} C_{kl}^{-1} \frac{y_i^{(l)} - f^{(l)}(x_i)}{\sigma(y_i^{(l)} - f^{(l)}(x_i))}$$

$$C_{lm} = \sum_k \left( y_l^{(k)} - f^{(k)}(x_l) - (y_l - f(x_l)) \right) \left( y_m^{(k)} - f^{(k)}(x_m) - (y_m - f(x_m)) \right)$$

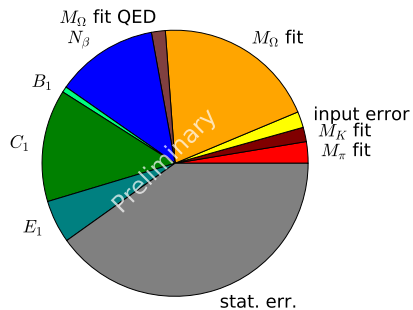
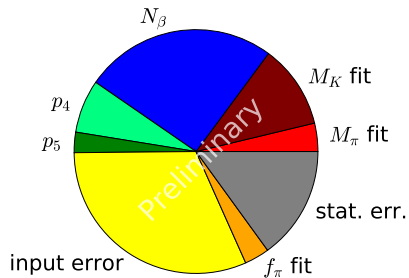
$\sigma$ : Jackknife error,  $f$ : fit function for  $w_0 M_\Omega$

# Full error

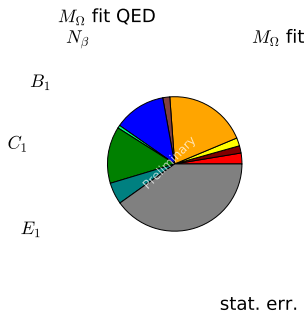
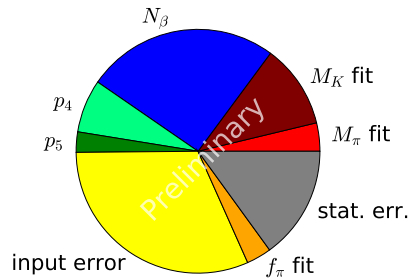
$w_0 M_\Omega$ :



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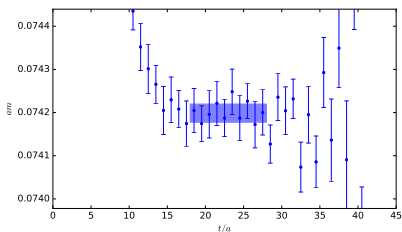
 $w_0 M_\Omega$ : $w_0 f_\pi$ :

## Full error

 $w_0 M_\Omega$ : $w_0 f_\pi$ :

# Analysis overview

- 1 Determine  $M_\Omega$ ,  $M_\pi$ , and  $M_K$  either by fitting the correlator or solving GEVP.



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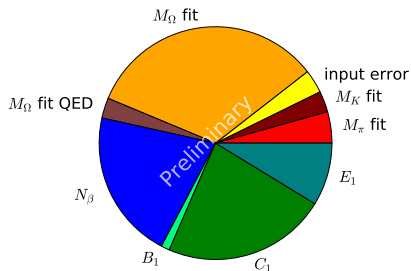
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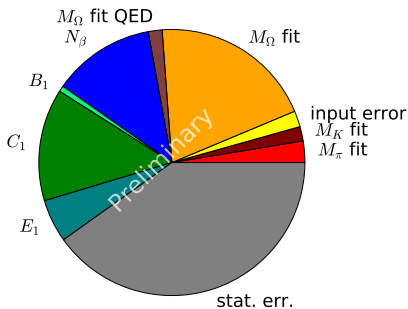
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- 3 Solve the fit function for  $w_0$  so the determination is self consistent.
- 4 Estimate the systematic error via the Histogram method by various analyses.
- 5 Determine the statistic error via the Jackknife method.



# Summary

- $w_0 M_\Omega$  is a suitable quantity to reduce the systematic error on  $w_0$ , due to the precise determination in experiment.
- The overall error can be reduced below 2‰

$w_0 M_\omega$ :

