

How to identify hadronic molecules

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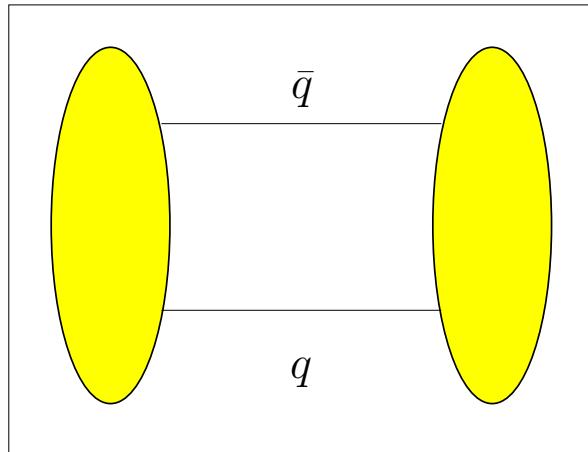
In collaboration with

M. Cleven, F.-K. Guo, Y. Kalashnikova, S. Krewald, A. Kudryavtsev, U.-G. Meißner, A. Nefediev

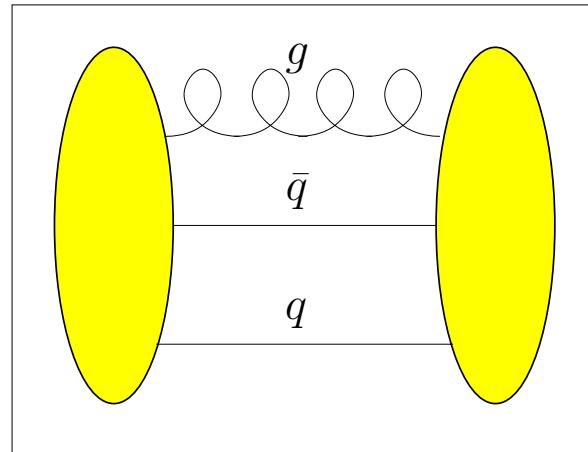
Key references:

- C. H., Yu. S. Kalashnikova, A. E. Kudryavtsev and A. V. Nefediev, Phys. Rev. D **75** (2007) 074015.
- F.-K. Guo, C.H., S. Krewald, U.-G. Meißner, Phys. Lett. **B666** (2008)251.
- M. Cleven, F.-K. Guo, C.H., U.-G. Meißner, Eur. Phys. J **A** in print; arXiv:1009.3804 [hep-ph].

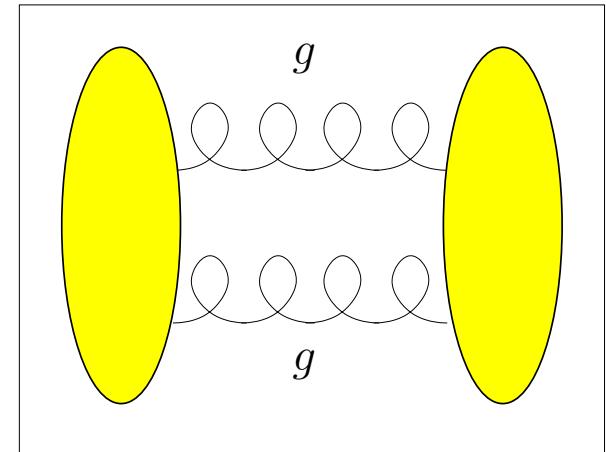
$\bar{q}q$



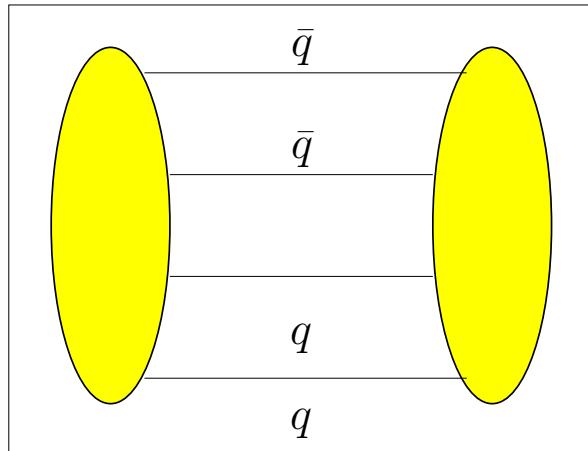
hybrid



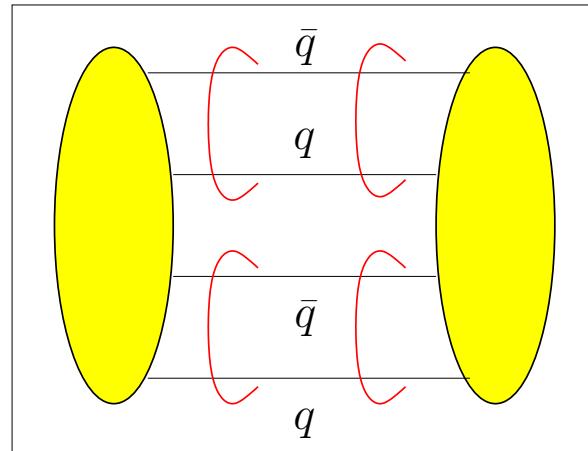
glueball



tetraquark



molecule



Only hadrons can
 go **on-shell** →
Unique analytic
 properties of
molecular amplitude

Definition using non-rel. QM

Landau (1960), Weinberg (1963), Baru et al. (2004)

Expand in terms of **non-interacting** quark and meson states

$$|\Psi\rangle = \begin{pmatrix} \lambda|\psi_0\rangle \\ \chi(\mathbf{p})|h_1 h_2\rangle \end{pmatrix},$$

here $|\psi_0\rangle$ = elementary state and $|h_1 h_2\rangle$ = two-hadron cont., then λ^2 equals probability to find the bare state in the physical state
 $\rightarrow \lambda^2$ is the quantity of interest!

The Schrödinger equation reads

$$\hat{\mathcal{H}}|\Psi\rangle = E|\Psi\rangle, \quad \hat{\mathcal{H}} = \begin{pmatrix} \hat{H}_c & \hat{V} \\ \hat{V} & \hat{H}_{hh}^0 \end{pmatrix} \rightarrow \chi(p) = \lambda \frac{f(p^2)}{E - p^2/(2\mu)}$$

introducing the transition form factor $\langle\psi_0|\hat{V}|hh\rangle = f(p^2)$,
Note: \hat{H}_{hh}^0 contains only meson kinetic terms!

Effective Coupling

Therefore

$$|\Psi\rangle = \lambda \begin{pmatrix} |\psi_0\rangle \\ -\frac{f(p^2)}{\epsilon + p^2/(2\mu)} |h_1 h_2\rangle \end{pmatrix},$$

For the normalization of the physical state we get

$$1 = \langle \Psi | \Psi \rangle = \lambda^2 \left(1 + \int \frac{f^2(p^2)}{(\epsilon + p^2/(2\mu))^2} d^3 p \right)$$

using

$$\int \frac{f^2(p^2) d^3 p}{(p^2/(2\mu) + \epsilon)^2} = \frac{4\pi^2 \mu^2 f(0)^2}{\sqrt{2\mu\epsilon}} + \mathcal{O}\left(\frac{\sqrt{\epsilon\mu}}{\beta}\right)$$

for **s-waves** with $\beta = \text{range of forces}$. Using $8\pi^2 \mu f(0)^2 = g$

$$1 = \lambda^2 \left(1 + \frac{\mu g / 2}{\sqrt{2\mu\epsilon}} + \mathcal{O}\left(\frac{\sqrt{\epsilon\mu}}{\beta}\right) \right)$$

Thus...

using for residue $g_{\text{eff}}^2/4\pi = \lambda^2 2(m_1 + m_2)^2 g$

$$\frac{g_{\text{eff}}^2}{4\pi} = 4(m_1 + m_2)^2(1 - \lambda^2)\sqrt{2\epsilon/\mu} \leq 4(m_1 + m_2)^2\sqrt{2\epsilon/\mu}$$

$(1 - \lambda^2)$ = Quantifies molecular component in physical state

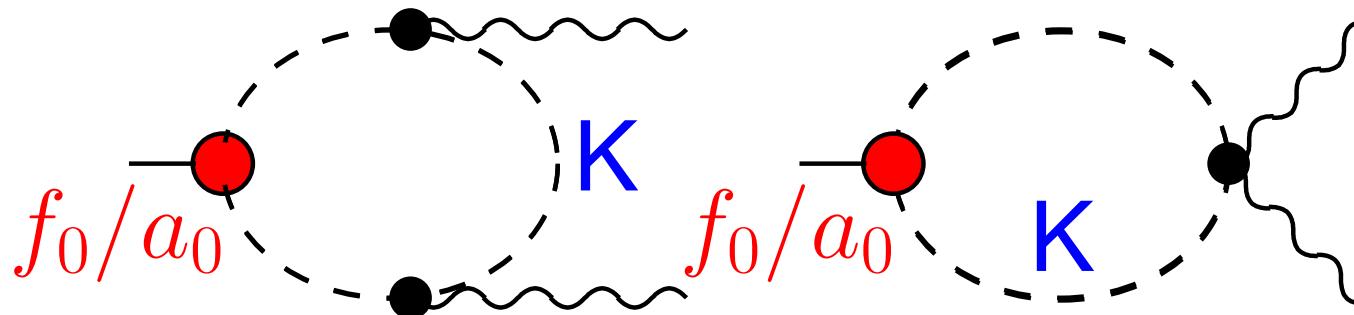
The **structure information** is hidden in the
effective coupling, extracted from experiment,
 independent of the phenomenology
 used to introduce the pole(s)

Formula also heavily used by Tuebingen group!

Example I: light scalar mesons

Two-photon decays of $f_0(980)/a_0(980)$

C.H. et al. (2007)



$$\Gamma_{\gamma\gamma} = g_{\text{eff}}^2 \frac{\alpha^2}{256\pi^3} \frac{1}{x} \left(\frac{1}{x^2} \arcsin(x)^2 - 1 \right)^2$$

where $x = 1 - \varepsilon/(2m)$. For the f_0/a_0 : $\varepsilon = 10$ MeV; $m = m_K$, s.t.

$$\Gamma_{\gamma\gamma}^{\text{theo.}} = (1 - \lambda^2)(0.22 \pm 0.07) \text{ keV} + \dots \text{ vs. } \Gamma_{\gamma\gamma}^{\text{exp. } f_0} = 0.22 \pm 0.02 \text{ keV}$$

Exp. value from R. Garcia-Martin, B. Moussallam (2010), analysing Belle data

Leading range corrections scale as $m\varepsilon/\beta^2 \simeq 1\%$

Example II: $D_s(2317)/D_s^*(2460)$

Experimental facts:

- Masses well below quark model predictions
- $M_D + M_K - M(D_s(2317)) = M_{D^*} + M_K - M(D_s^*(2460))$
 - $D_s(2317)/D_s^*(2460)$ as $K D^{(*)}$ bound states?
van Beeveren, Rupp; Oset, Gamermann; Lutz, Soyeur; ...

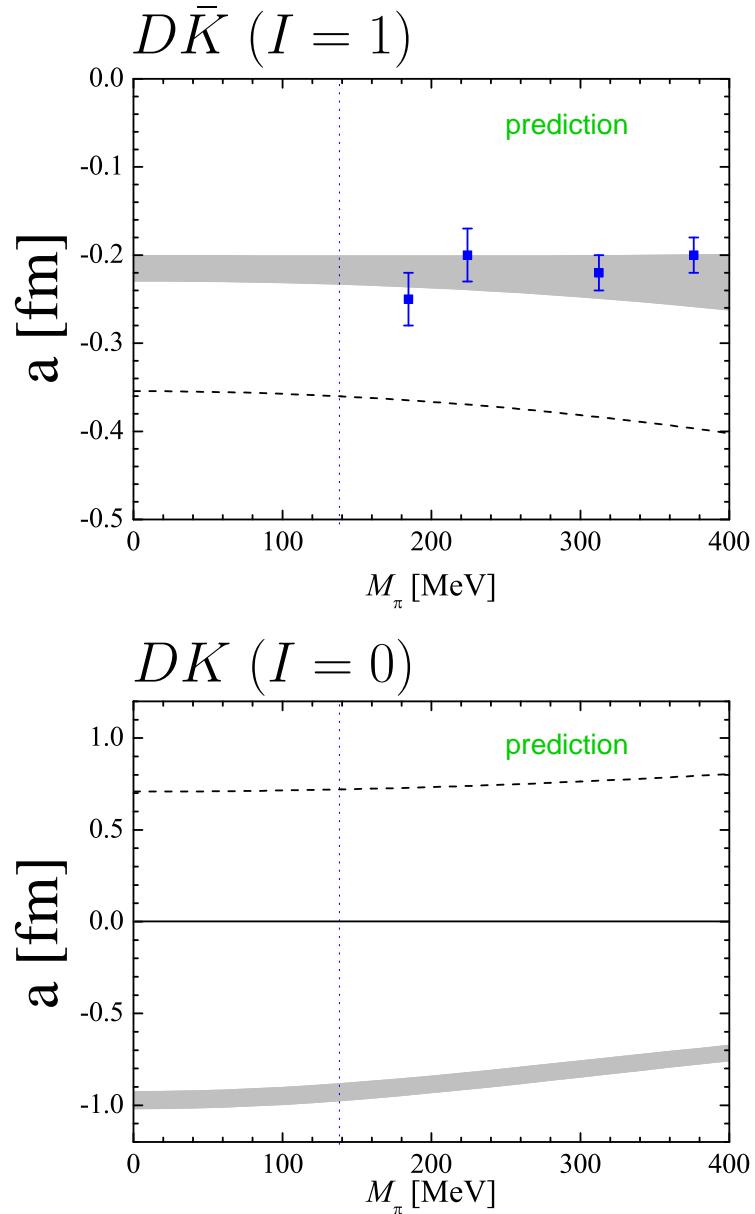
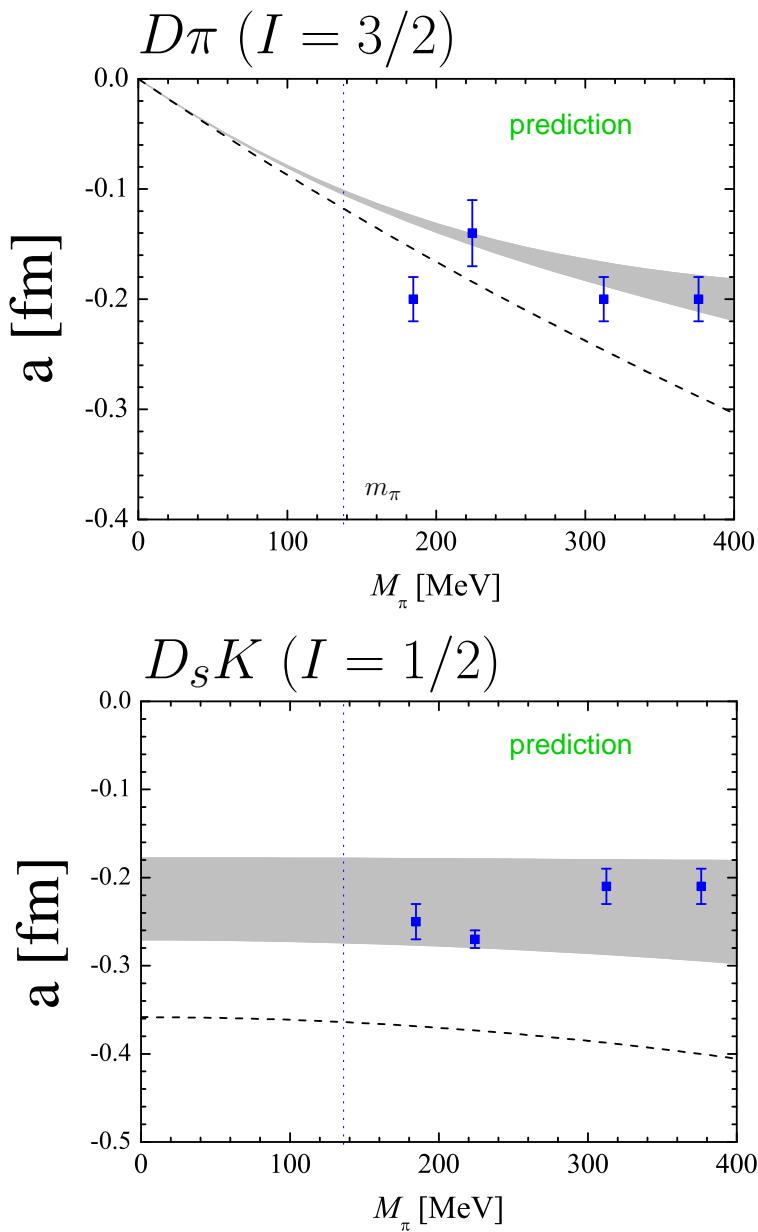
Here: Goldstone-Boson– $D^{(*)}$ -Meson scattering

- ChPT → controlled quark mass dependence
 - up to potential quark mass dep. of regulator*
- with unitarization → dynamical generation of poles
- use LEC to fix pole position of $D_s(2317)$

From residues: $D_s(2317)$ ($D_s^*(2460)$) as $K D(K D^*)$ molecules

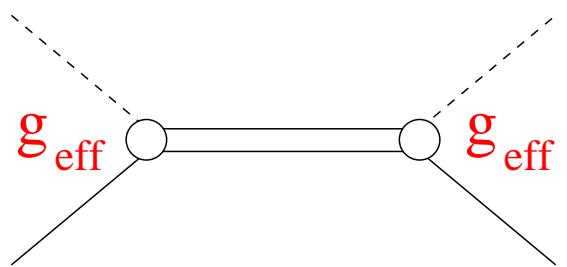
Prediction of other observables; Sensitivity to molecular nature?

Chiral extrapolation

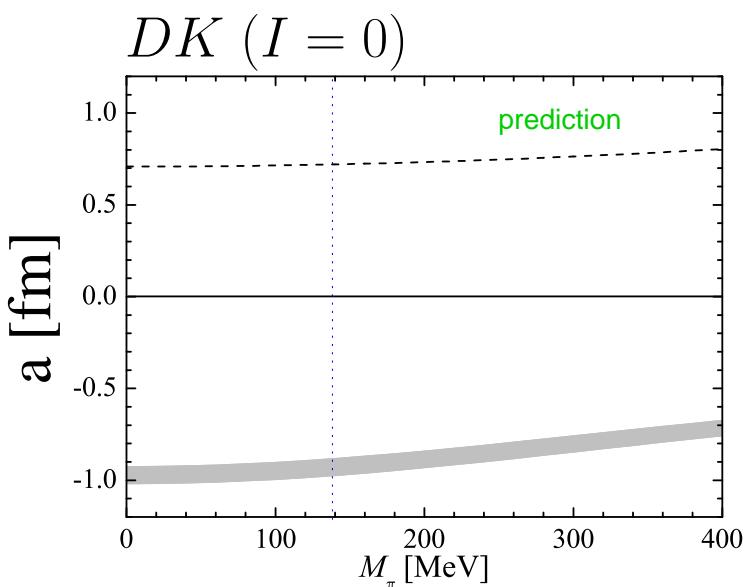
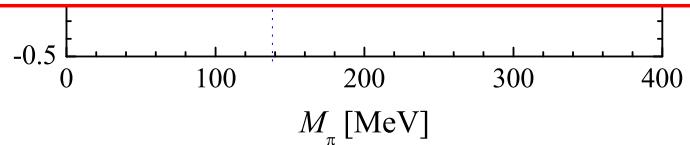
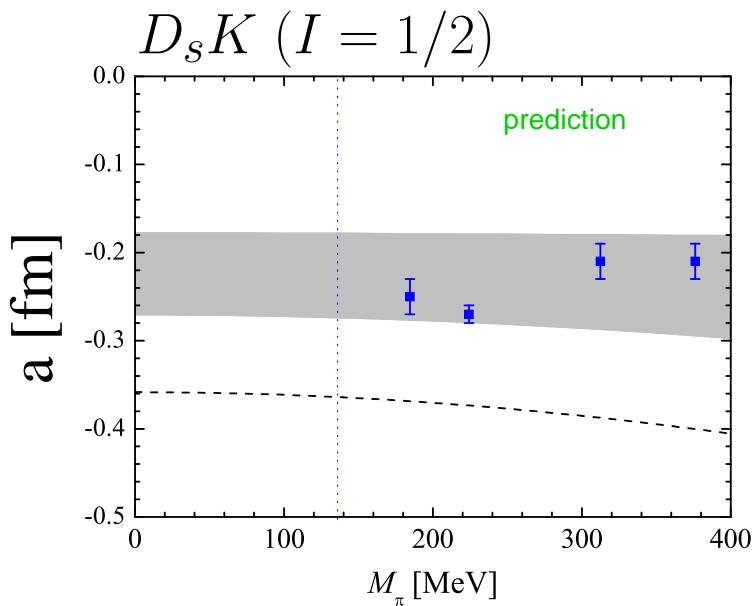
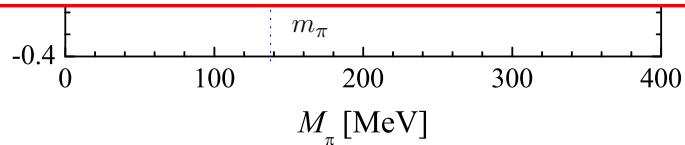


Lattice: Liu, Lin, Orginos (2008); UChPT: Guo et al. (2009)

Chiral extrapolation

$$D_S(2317): a = g_{\text{eff}} + \mathcal{O}(1/\beta) \simeq \left(\frac{2(1-\lambda^2)}{2-\lambda^2} \right) \frac{-1}{\sqrt{2m_K \epsilon}}$$


$a = 1 \text{ fm}$ for molecule ($\lambda^2 = 0$); smaller otherwise



Lattice: Liu, Lin, Orginos (2008); UChPT: Guo et al. (2009)

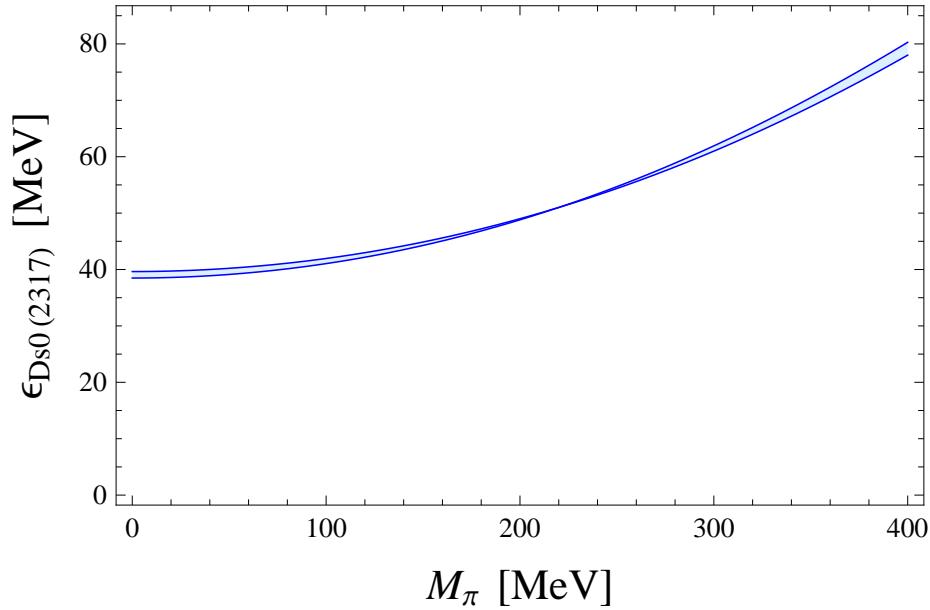
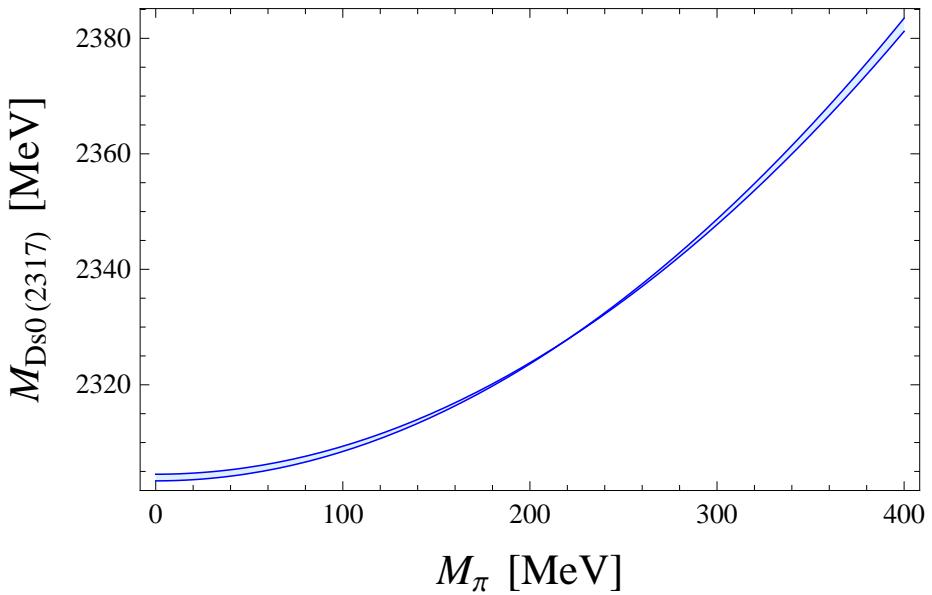
π mass dependence of poles

M. Cleven et al. (2010)

If $D_s^*(2317)$ as quark state: $c\bar{s}$

expect very weak light quark mass dependence

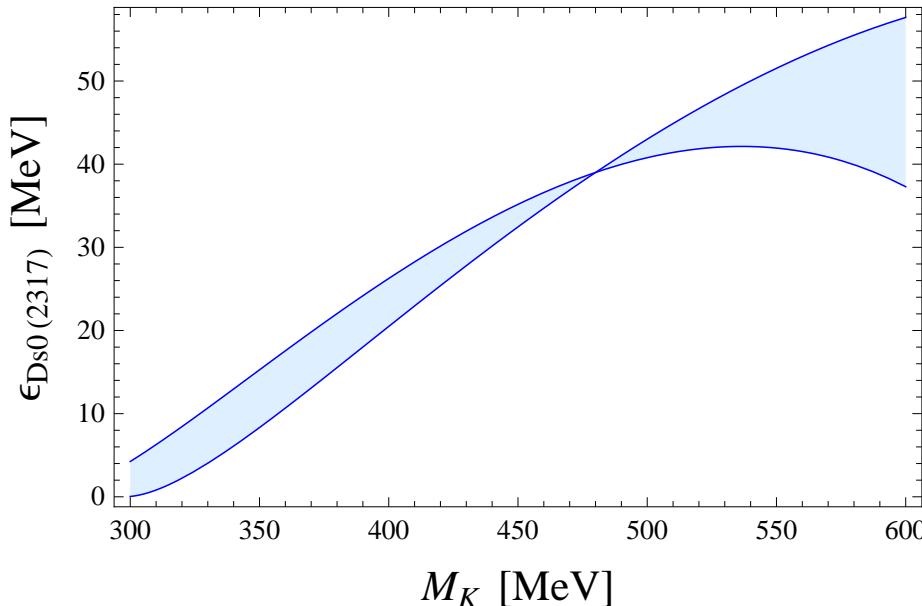
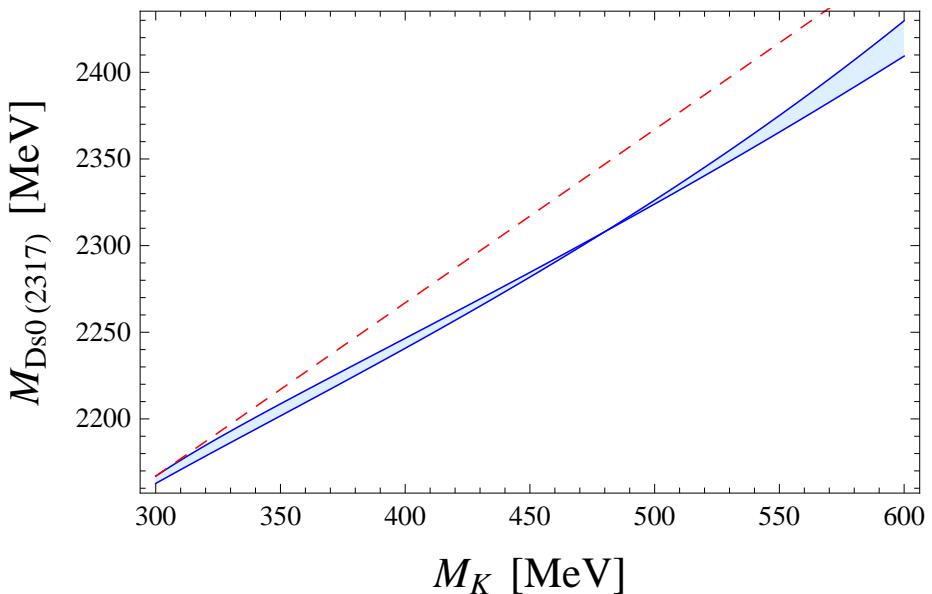
We find



Molecule shows relatively strong light quark mass dependence
since prominent component KD contains light quarks!

K mass dependence of poles

M. Cleven et al. (2010)



- very pronounced linear K –mass dependence
- state follows threshold
- for K –mass smaller than 300 MeV turns into resonance
note: interaction strength propto. M_K

Isospin breaking in QCD and EFT through quark mass and charge differences

The same effective operators lead to

→ mass differences, e.g.

$$\begin{aligned} \triangleright m_{D^+} - m_{D^0} &= \Delta m^{\text{strong}} + \Delta m^{\text{e.m.}} \\ &= ((2.5 \pm 0.2) + (2.3 \pm 0.6)) \text{ MeV} \end{aligned}$$

▷ $\pi^0 - \eta$ mixing → parameters fixed

→ Isospin breaking scattering amplitude

▷ e.g. $K D \rightarrow \pi^0 D_s$ predicted; with this

$$\Gamma(D_s(2317) \rightarrow D_s \pi^0) = (180 \pm 110) \text{ keV}$$

Lutz, Soyeur (2007); complete to NLO+uncertainty estimate: Guo et al. (2008)

much smaller in quark model → direct measurement (PANDA)

Conclusion

Progress through interplay of
controlled theory,
experiments of high quality and
numerical simulations with super computers