

COUPLED-CHANNEL FORMALISM

As discussed in [1]: Main tension between $\pi\pi$ data sets is absorbed in inelastic parameters

$\Rightarrow$ those parameters belong to exclusive channels – what happens if we build them in explicitly?

Idea: Do different $\pi\pi$ data sets agree better with the other exclusive channels?

We construct three different channel structures:

- pseudoscalar + pseudoscalar ($\pi\pi, K\bar{K}$) [PP]
- vector + pseudoscalar ($\rho\eta, \omega\pi, K\bar{K}^*$) [VP]
- axial vector + pseudoscalar ($a_1\pi$) [AP]

COMPONENTS [2]

The $\mathcal{S}$ -, $\mathcal{M}$ -, and $\mathcal{T}$ -matrices:

$${}_{\text{out}}\langle q_1 q_2, b | \mathcal{S} - 1 | p_1 p_2, a \rangle_{\text{in}} = i(2\pi)^4 \delta^4(p_1 + p_2 - q_1 - q_2) \mathcal{M}_{ba},$$

$$\mathcal{M}(s)_{ab} = \xi_a(s) \mathcal{T}(s)_{ab} \xi_b(s),$$

with

$$\begin{aligned} \rho(s) &= \text{phase space factor,} \\ \xi(s) &= \beta(s) B(s) = \text{centrifugal barrier factor,} \\ \beta(s) &= \text{interaction structure function,} \\ B(s) &= \text{taming factor.} \end{aligned}$$

We are using the two-potential formalism and divide $\mathcal{T}$ into resonant and non-resonant (background) parts

$$\mathcal{T}_{ab} = \mathcal{T}_{ab}^{\text{B}} + \mathcal{T}_{ab}^{\text{R}}$$

BACKGROUND POTENTIAL

As background we use a contact interaction term

$$V_{\text{B}}(s) = f_0 \quad \Rightarrow \quad \mathcal{T}_{\text{B}} = (1 - V_{\text{B}}\Pi)^{-1} V_{\text{B}}$$

The loop integral of two non-interacting particles is dispersively defined as

$$\begin{aligned} \Pi(s) &= \Pi(s_0) + \frac{s - s_0}{2\pi i} \int_{s_{\text{thr}}}^{\infty} \frac{ds'}{s' - s_0} \frac{\text{disc}(\Pi(s'))}{s' - s} \\ \text{with } \text{disc}(\Pi(s)) &= 2i\xi(s)\rho(s)\xi(s). \end{aligned}$$

Two particle interaction is dressed by the background interaction

$$\begin{aligned} \gamma_{\text{in}}^{\dagger} &= 1 + \Pi\mathcal{T}_{\text{B}}, \quad \gamma_{\text{out}} = 1 + \mathcal{T}_{\text{B}}\Pi, \\ \text{with } \text{disc}(\gamma(s)) &= 2i\mathcal{M}_{\text{B}}^*(s)\rho(s)\gamma(s). \end{aligned}$$

RESONANCE POTENTIAL

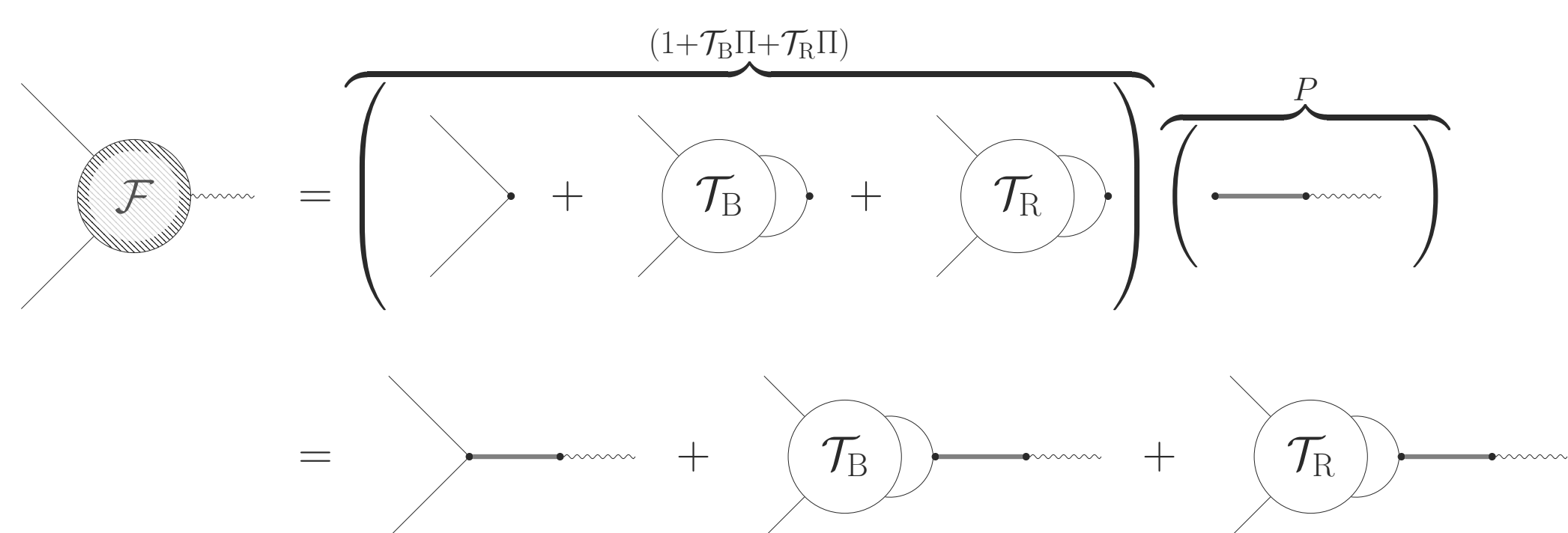
The resonance interaction is defined via

$$V_{\text{R}}(s) = -g^T G_{\text{R}}(s) g, \quad G_{\text{R}}^{kl}(s) = \frac{\delta_{kl}}{s - m_k^2} \Rightarrow \quad \mathcal{T}_{\text{R}} = \gamma_{\text{out}} (1 - V_{\text{R}}\Sigma)^{-1} V_{\text{R}} \gamma_{\text{in}}^{\dagger}$$

$$\begin{aligned} \Sigma(s)_{ij} &= \Sigma(s_0)_{ij} + \frac{s - s_0}{2\pi i} \int_s^{\infty} \frac{ds'}{s' - s_0} \frac{\text{disc}(\Sigma(s'))_{ij}}{s' - s} \\ \text{with } \text{disc}(\Sigma(s)) &= 2i\xi(s)\gamma(s)^* \rho(s)\xi(s)\gamma(s). \end{aligned}$$

FORM FACTOR AND OBSERVABLES

$$F = \gamma_{\text{out}} (1 - V_{\text{R}}\Sigma) P \quad \text{with} \quad P_i = \sum_R \frac{\alpha^R g_i^R}{s - m_R^2}$$



CHANNELS

PP channel:

$$\begin{aligned} \langle P(q_1) \bar{P}(q_2) | j^{\mu}(0) | 0 \rangle &= (q_1 - q_2)^{\mu} F_P^V(s), \\ \beta_{PP}(s)^2 &= \frac{1}{3} (q_1 - q_2)^{\mu} g_{\mu\nu} (q_1 - q_2)^{\nu}, \\ \beta_{PP}(s) &= \sqrt{\frac{s - 4M_P^2}{3}}. \end{aligned}$$

VP channel:

$$\begin{aligned} \langle P(q_0) V(q_V, \lambda) | j^{\mu}(0) | 0 \rangle &= \epsilon^{\mu\nu\alpha\beta} n_{\nu}^{\lambda} q_{0\alpha} k_{\beta} f_{VP}(s), \\ \beta_{VP}(s)^2 &= \frac{1}{3} \sum_{\lambda} (\epsilon^{\mu\nu\alpha\beta} n_{\nu}^{\lambda} q_{0\alpha} k_{\beta})^* g_{\mu\rho} (\epsilon^{\rho\gamma\eta\delta} n_{\gamma}^{\lambda} q_{0\eta} k_{\delta}), \\ \beta_{VP}(s) &= \sqrt{\frac{(s - (M_P + M_V)^2)(s - (M_P - M_V)^2)}{6}}. \end{aligned}$$

AP channel:

$$\langle a_1(q_1, \lambda) \pi(q_2) | j^{\mu}(0) | 0 \rangle = n_{\nu}^{\lambda} (T_1^{\mu\nu} F_1(s) + T_2^{\mu\nu} F_2(s)) \quad [3]$$

Definition of structure function more involved

Resonances in loop

$\rho\eta$, $K^*\bar{K}$ and $a_1\pi$ channel have resonances (R) with non-negligible width. In construction of Π_i we use spectral density to smear out discontinuity

$$\Pi_{RP}(s) = \frac{s - s_0}{\pi} \int_{(\sqrt{s_{\text{decay}}} + m_P)^2}^{\infty} \frac{\int_{s_{\text{decay}}}^{(\sqrt{s'} - m_P)^2} \rho(s', m_R^2, m_P^2) \xi_{RP}^2(s', m_R^2, m_P^2) \sigma_R(m_R^2) dm_R^2}{(s' - s)(s' - s_0)} ds'$$

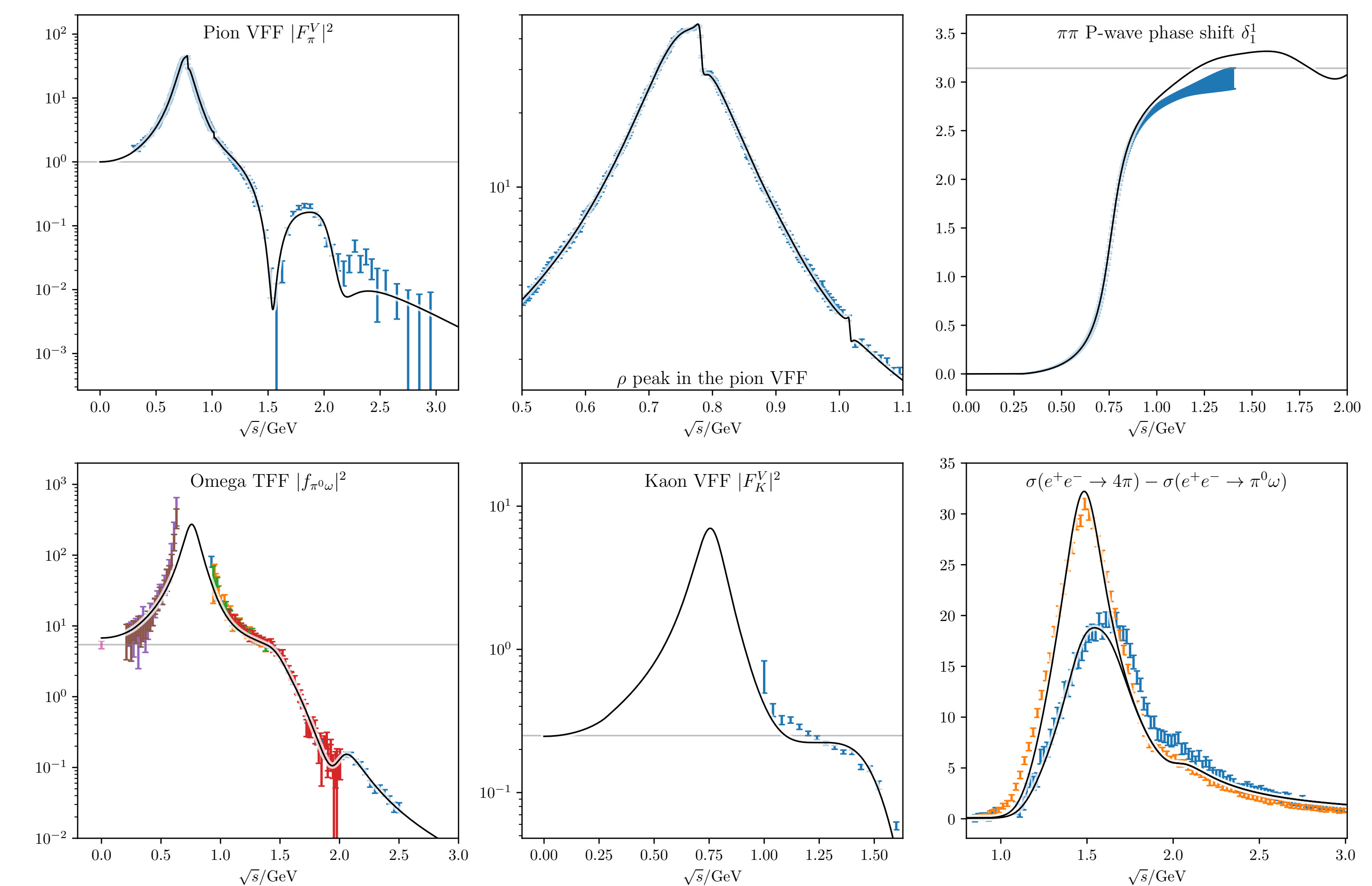
Loops have complicated analytic structure [4]

Effective 4π

We use the decay chain $a_1\pi \rightarrow \rho\pi\pi \rightarrow 4\pi$

- decouple structures into two different channels via change of basis
- have to take care of symmetry factors and charge states
 $\rightarrow$ simulated different interference effects in charged and partially charged final states
- smearing out complicated: pole position of a_1 not well-known
 $\rightarrow$ use τ -decay data for fit of spectral density of a_1

PRELIMINARY RESULTS



References

- [1] P. Stoffer, G. Colangelo, and M. Hoferichter, JINST **18**, C10021 (2023) [arXiv:2308.04217 [hep-ph]].
- [2] L. A. Heuser, G. Chanturia, F. K. Guo, C. Hanhart, M. Hoferichter, and B. Kubis [arXiv:2403.15539 [hep-ph]].
- [3] G. Colangelo, M. Hoferichter, M. Procura, and P. Stoffer, JHEP **09**, 074 (2015) [arXiv:1506.01386 [hep-ph]].
- [4] M. Döring, C. Hanhart, F. Huang, S. Krewald, and U.-G. Meissner, Nucl. Phys. A **829**, 170 (2009) [arXiv:0903.4337 [nucl-th]].

