

Calculations of kaonic nuclei based on chiral meson-baryon coupled channel interaction models

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Introduction

- Self-consistent calculations of K^- -nuclear quasi-bound states using the following chiral meson-baryon interaction models:
 - **Prague** (P NLO)
(A. Cieply, J. Smejkal, Nucl. Phys. A 881 (2012) 115)
 - **Kyoto-Munich** (KM NLO)
(Y. Ikeda, T. Hyodo and W. Weise, Nucl. Phys. A 881 (2012) 98)
 - **Murcia** (M1 and M2)
(Z. H. Guo and J. A. Oller, Phys. Rev. C 87,(2013) 035202)
 - **Bonn** (B2 and B4)
(M. Mai and U.-G. Meißner, Nucl. Phys. A 900, (2013) 51)

Free-space $K^- p$ amplitudes

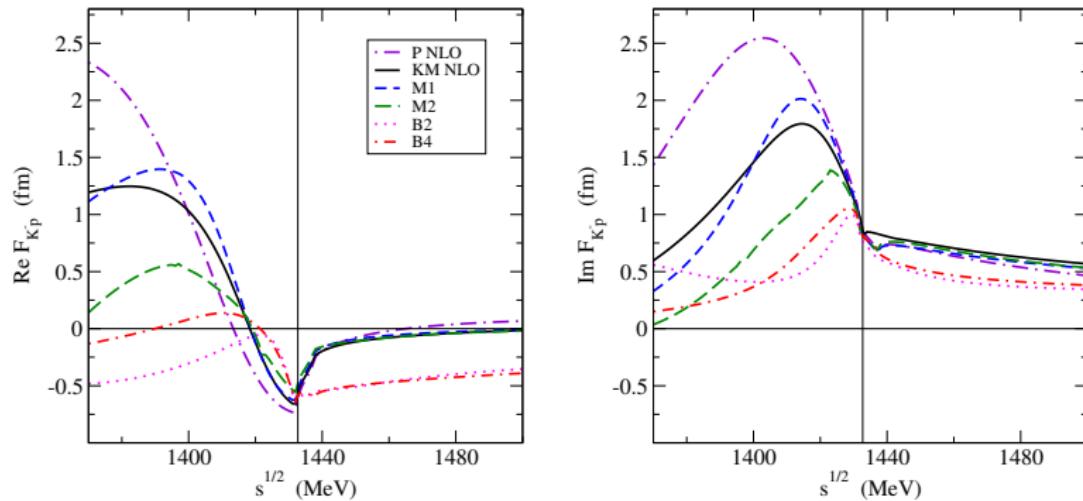


Fig.1: Energy dependence of real (left) and imaginary (right) parts of free-space $K^- p$ amplitudes in considered models.

Free-space K^-n amplitudes

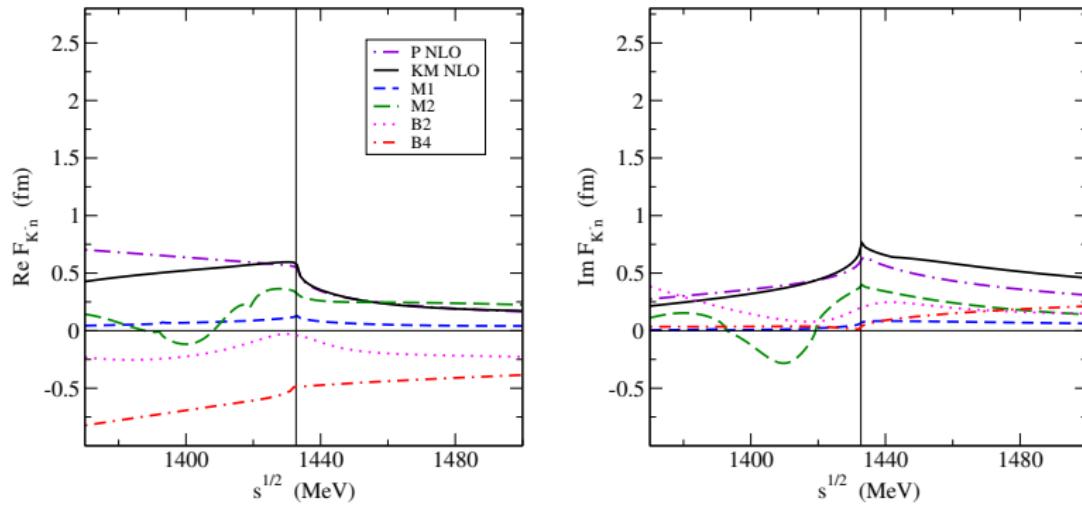


Fig.2: Energy dependence of real (left) and imaginary (right) parts of free-space K^-n amplitudes in considered models.

Model

- Klein-Gordon equation for K^-

$$\left[\omega_K^2 + \vec{\nabla}^2 - m_K^2 - \Pi_K(\vec{p}_K, \omega_K, \rho) \right] \phi_K = 0$$

complex energy $\omega_K = m_K - B_K - i\Gamma_K/2 - V_C = \tilde{\omega}_K - V_C$

- self-energy operator

$$\Pi_K = 2\text{Re}(\tilde{\omega}_{K-})V_{K-} = -4\pi \frac{\sqrt{s}}{m_N} \left(F_0 \frac{1}{2} \rho_p + F_1 \left(\frac{1}{2} \rho_p + \rho_n \right) \right),$$

F_0 and F_1 – isospin 0 and 1 scattering amplitudes

- Nucleus described within the RMF model

Model

- Free space amplitudes → in-medium amplitudes - WRW method
(T. Wass, M. Rho, W. Weise, Nucl. Phys. A 617 (1997) 449)

$$F_1 = \frac{F_{K^-n}(\sqrt{s})}{1 + \frac{1}{4}\xi_k \frac{\sqrt{s}}{m_N} F_{K^-n}(\sqrt{s})\rho}, \quad F_0 = \frac{[2F_{K^-p}(\sqrt{s}) - F_{K^-n}(\sqrt{s})]}{1 + \frac{1}{4}\xi_k \frac{\sqrt{s}}{m_N} [2F_{K^-p}(\sqrt{s}) - F_{K^-n}(\sqrt{s})]\rho}$$

where

$$\xi_k = \frac{9\pi}{p_f^2} 4 \int_0^\infty \frac{dt}{t} \exp(iqt) j_1^2(t), \quad q = \frac{1}{p_f} \sqrt{\tilde{\omega}_{K^-}^2 - m_{K^-}^2}.$$

- P NLO + Pauli + SE model
(A. Cieply, J. Smejkal, Nucl. Phys. A 881 (2012) 115)

$$F_{ij}(p, p'; \sqrt{s}) = -\frac{g_i(p)g_j(p')}{4\pi f_i f_j} \sqrt{\frac{M_i M_j}{s}} [(1 - C(\sqrt{s}) \cdot G(\sqrt{s})^{-1}) \cdot C(\sqrt{s})]_{ij},$$

$$G_i(\sqrt{s}; \rho) = \frac{1}{f_i^2} \frac{M_i}{\sqrt{s}} \int_{\Omega_i(\rho)} \frac{d^3 \vec{p}}{(2\pi)^3} \frac{g_i^2(p)}{p_i^2 - p^2 - \Pi_i(\sqrt{s}, \vec{p}; \rho) + i0}.$$

In-medium modified $\bar{K}N$ amplitudes

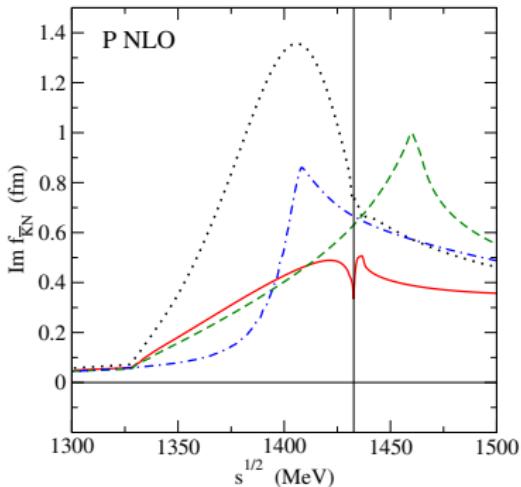
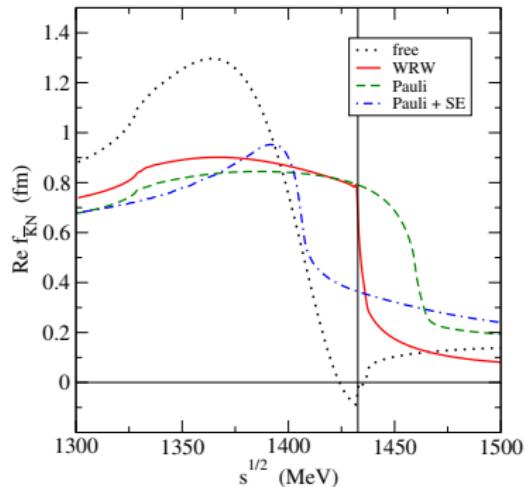


Fig.3: Energy dependence of reduced free-space (dotted line) $f_{\bar{K}N} = \frac{1}{2}(f_{K-p} + f_{K-n})$ amplitude compared with WRW modified amplitude (solid line), Pauli (dashed line), and Pauli + SE (dot-dashed line) modified amplitude for $\rho_0 = 0.17 \text{ fm}^{-3}$ in the P NLO model.

Energy dependence

- Amplitudes as a function of available energy \sqrt{s}

$$(s = (E_N + E_{K^-})^2 - (\vec{p}_N + \vec{p}_{K^-})^2)$$

- $\bar{K}N$ cms frame $\rightarrow K^-$ -nucleus frame $\vec{p}_N + \vec{p}_{K^-} \neq 0$

(A. Cieplý, E. Friedman, A. Gal, D. Gazda, J. Mareš, PLB 702, 402 (2011))

$$\sqrt{s} = m_N + m_{K^-} - 8.5 - \xi_N B_{K^-} + \xi_{K^-} \text{Re}V_{K^-}(r) - \xi_N T_N \left(\frac{\rho}{\rho_0}\right)^{2/3},$$

where $V_{K^-} = V_{K^-} + V_C$, $T_N = 23$ MeV, and $\xi_{N(K^-)} = m_{N(K^-)} / (m_N + m_{K^-})$.

K^- optical potential

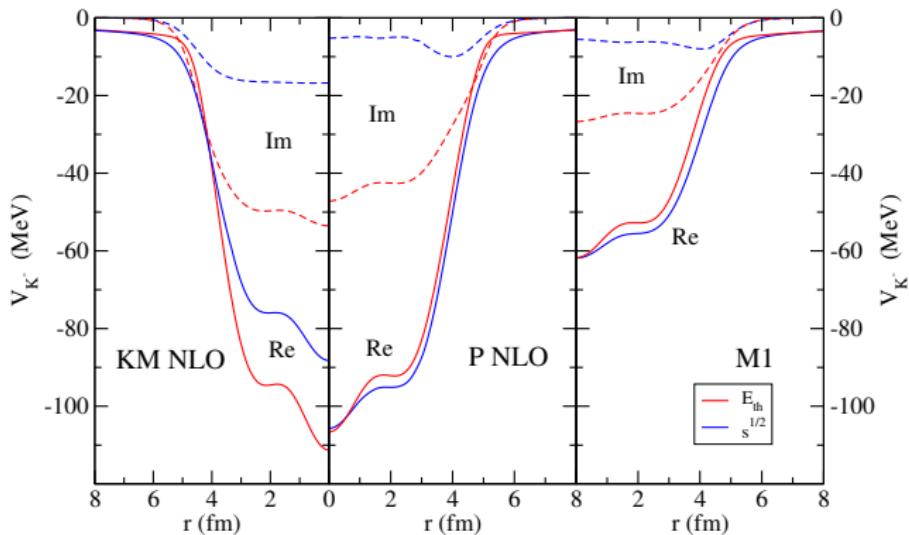


Fig.4: K^- optical potential in ${}^{40}\text{Ca}$ calculated for self-consistent \sqrt{s} and at threshold E_{th} in the KM NLO, P NLO, and M1 models.

K^- optical potential

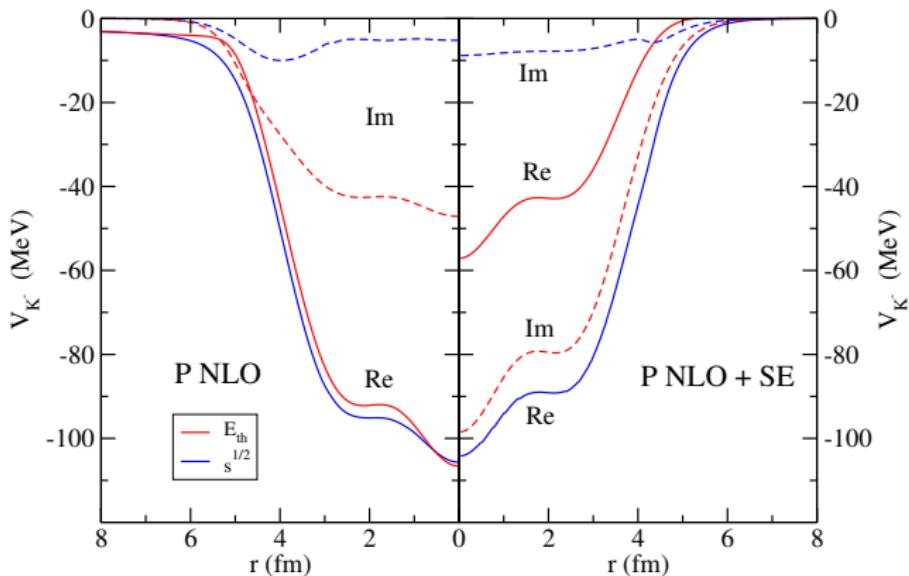


Fig.5: K^- optical potential in ${}^{40}\text{Ca}$ calculated self-consistently for in-medium \sqrt{s} and at threshold E_{th} using WRW modified amplitudes in the P NLO model compared with P NLO + Pauli + SE model.

K^- 1s binding energies and widths

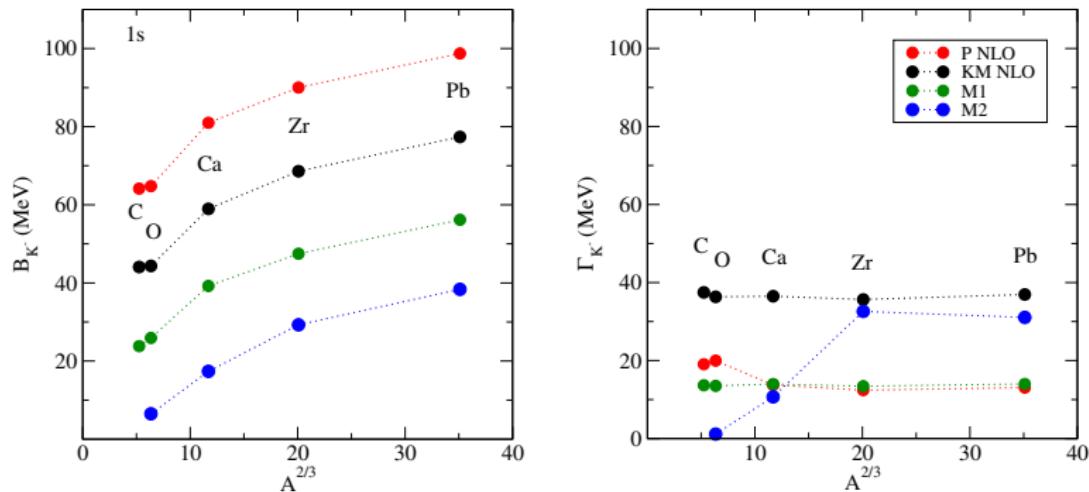


Fig.6: 1s K^- binding energies (left) and corresponding widths (right) in various nuclei calculated self-consistently in the P NLO, KM NLO, M1, and M2 models.

K^- 1s binding energies and widths

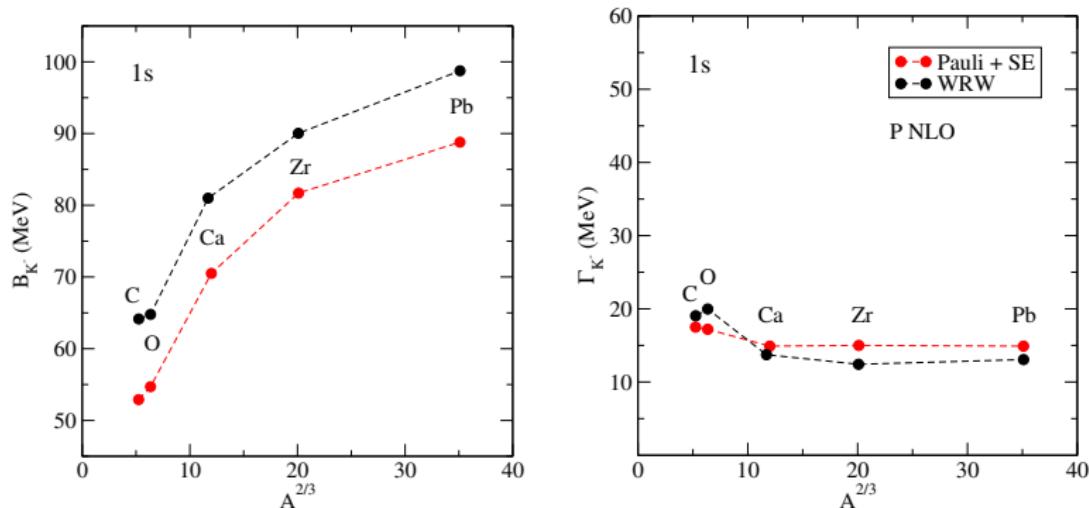


Fig.7: K^- binding energies (left) and widths (right) calculated using the WRW method in the P NLO model compared with P NLO + Pauli + SE model.

K^- spectrum in ^{40}Ca

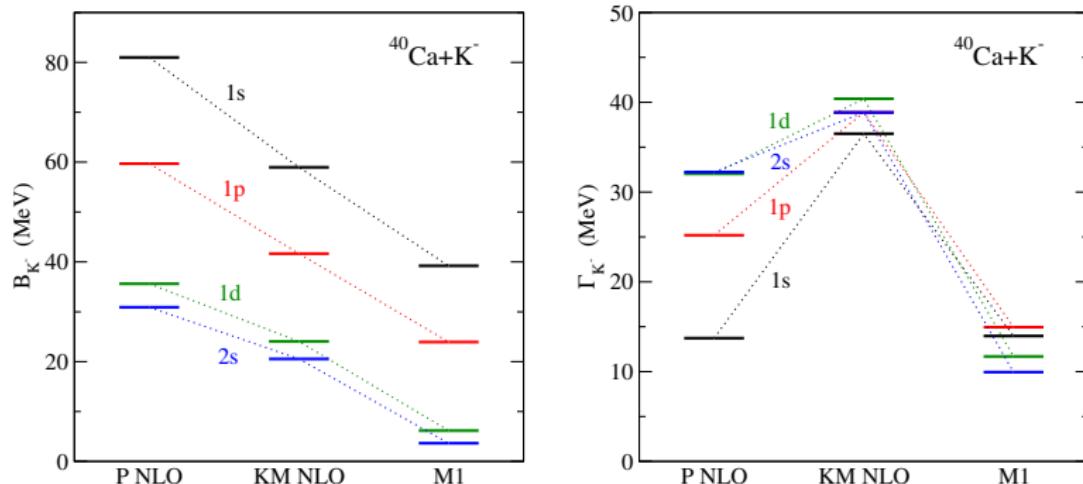


Fig.8: K^- binding energies (left) and widths (right) in s, p and d levels in ^{40}Ca calculated self-consistently in the P NLO, KM NLO, and M1 models.

Effects not included in the present calculations

- dynamic approach — core polarization effect up to $\simeq 5$ MeV in K^- binding energies
- P-wave interaction — weakly attractive $\simeq +5$ MeV in binding energy, $\simeq -1$ MeV in width
- $2N$ absorption $\bar{K}NN \rightarrow YN$

The effect of $2N$ absorption (CS30 model):

		C	O	Ca	Zr	Pb
$\sqrt{s} + \text{SE} + \text{dyn.}$	B_K	55.7	56.0	70.2	80.5	87.0
	Γ_K	12.3	12.1	10.8	10.9	10.8
$\sqrt{s} + \text{SE} + \text{dyn.} + 2N$	B_K	54.0	55.1	67.6	79.6	86.3
	Γ_K	44.9	53.3	65.3	48.7	47.3

(D. Gazda, J. Mareš, NPA 881, 2012, 159)

Conclusions

- Calculations of K^- -nuclear quasi-bound states based on chiral meson-baryon interaction models:
 - large model dependence of K^- binding energies
 - small K^- widths (10 - 20 MeV) in P NLO and M1 models,
twice larger widths (\simeq 40 MeV) in KM NLO model
 - 2N absorption important — adds up to 50 MeV to the widths
 - no bound states found for B2 and B4 models