

Dense \bar{K} nuclei and their excited states

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1. Introduction
2. New framework of Antisymmetrized Molecular Dynamics
 - K - p - \bar{K}^0 n mixing
 - J & T projections
3. Ground state of $ppnK^-$, $pppK^-$, $pppnK^-$, ${}^6\text{Be}K^-$, ${}^9\text{BK}^-$
and ${}^{11}\text{CK}^-$
4. Excited states of $ppnK^-$ and ${}^{11}\text{CK}^-$
5. Summary

Introduction

Studies of deeply bound kaonic nuclei by Akaishi & Yamazaki

- phenomenological $\bar{K}N$ potential

- free $\bar{K}N$ scattering data
- X-ray data of kaonic hydrogen atom
- binding energy and width of $\Lambda(1405)$

Strongly attractive.

$$V_{\bar{K}N} \\ I=0$$

- Analysis of ${}^8\text{Be}K^-$ with $\alpha\alpha K^-$ model \Rightarrow

*Binding energy of $K^- = 113 \text{ MeV}$
Central density = $5 \rho_0$*

Kaonic nucleus

K⁻ deeply bound?

Highly dense?

Peculiar structure?



What kind of structure does A -nucleons + K^- system favor?

Formalism

Kaonic nuclei ... $\left\{ \begin{array}{l} \bullet \text{ unknown structure} \\ \bullet \text{ peculiar structure ex) isovector deformation} \end{array} \right.$

\Rightarrow \star Method : **Antisymmetrized Molecular Dynamics**
no assumption on clusters, shape, etc.

$V_{I=0}^{\bar{K}N}$ plays an essential role in kaonic nuclei.

Improving

\bullet K^-p/\bar{K}^0n mixing

K^- -proton \longleftrightarrow $I=0 \bar{K}N$ int. \longleftrightarrow \bar{K}^0 -neutron

\bullet Angular momentum & isospin projections

Wave function

Essence of K^-p/\bar{K}^0n mixing

Nucleon's wave function

$$|\varphi_i\rangle = \sum_{\alpha=1}^{Nn} C_{\alpha}^i \exp\left[-\nu\left(\mathbf{r} - \frac{\mathbf{Z}_{\alpha}^i}{\sqrt{\nu}}\right)^2\right] |\sigma_{\alpha}^i \tau_{\alpha}^i\rangle$$

Superposition of Gaussian wave packets

$$|\tau_{\alpha}^i\rangle = \left(\frac{1}{2} + \gamma_{\alpha}^i\right) |p\rangle + \left(\frac{1}{2} - \gamma_{\alpha}^i\right) |n\rangle$$

p-n mixing

Anti-kaon's wave function

$$|\varphi_K\rangle = \sum_{\alpha=1}^{Nk} C_{\alpha}^K \exp\left[-\nu\left(\mathbf{r} - \frac{\mathbf{Z}_{\alpha}^K}{\sqrt{\nu}}\right)^2\right] |\tau_{\alpha}^K\rangle$$

Superposition of Gaussian wave packets

\bar{K}^0 - K^- mixing

$$|N\rangle = a|\text{proton}\rangle + b|\text{neutron}\rangle$$

$$|K\rangle = x|K^-\rangle + y|\bar{K}^0\rangle$$

Total wave function

$$|\Phi\rangle = \det[|\varphi_i\rangle] \otimes |\varphi_K\rangle$$

$$|\Phi^{\pm}\rangle = |\Phi\rangle \pm |P\Phi\rangle$$

Charge projection

$$d\alpha \exp\left[-i\alpha(\hat{T}_z - M)\right] |\Phi^{\pm}\rangle$$

Use $|P_M \Phi^{\pm}\rangle$ as a trial function

J & T projections (VBP)

$$|\Phi^\pm\rangle \longrightarrow |P_{MK}^J P_{TzTz'}^T \Phi^\pm\rangle \quad : \text{Eigen state of angular momentum J and isospin T}$$

$$|P_{MK}^J P_{TzTz'}^T \Phi^\pm\rangle = \int d\Omega_{Ang} D_{MK}^{J*}(\Omega_{Ang}) \hat{R}_{Ang}(\Omega_{Ang}) \times \int d\Omega_{iso} D_{TzTz'}^{T*}(\Omega_{iso}) \hat{R}_{iso}(\Omega_{iso}) |\Phi^\pm\rangle$$

J projection

T projection

$$\hat{R}_{Ang}(\Omega) = \exp[-i\alpha\hat{J}_z] \exp[-i\beta\hat{J}_y] \exp[-i\gamma\hat{J}_z]$$

$$\hat{R}_{iso}(\Omega) = \exp[-i\alpha\hat{T}_z] \exp[-i\beta\hat{T}_y] \exp[-i\gamma\hat{T}_z]$$

Calculate various expectation values with $|P_{MK}^J P_{TzTz'}^T \Phi^\pm\rangle$.

Formalism

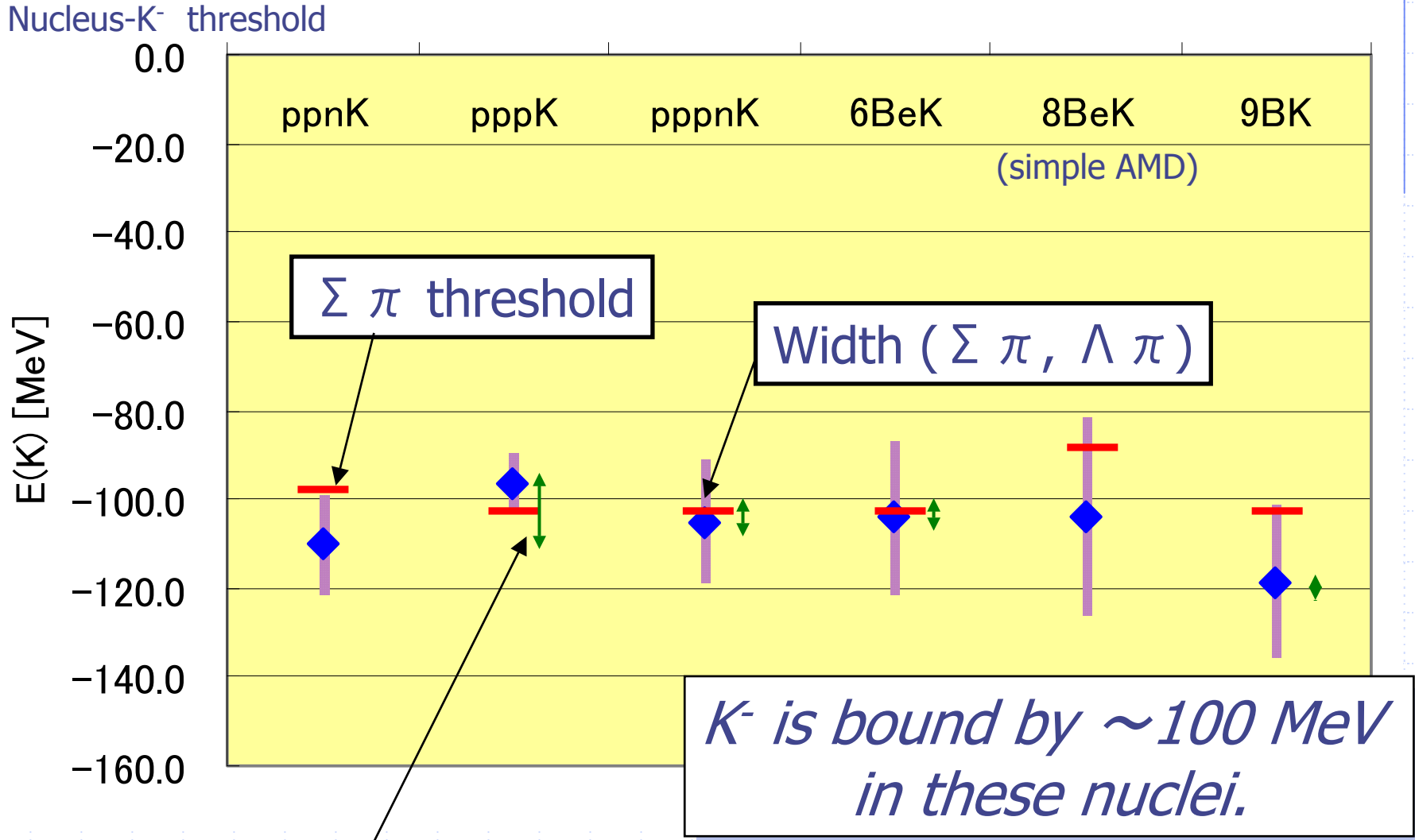
1. Hamiltonian $\hat{H} = \hat{T} + \hat{V}_{NN} + \hat{V}_{KN} + \hat{V}_{Coulomb} - \hat{T}_G$
2. Variational parameters $\{X_\alpha^i\} = \{C_\alpha^i, \mathbf{Z}_\alpha^i, \gamma_\alpha^i, C_\alpha^K, \mathbf{Z}_\alpha^K, \gamma_\alpha^K\}$
are determined by Frictional cooling eq. with constraint.
3. G-matrix method \Rightarrow Effective interaction $\hat{V}_{NN}, \hat{V}_{KN}$
bare NN int = Tamagaki potential (OPEG)
bare KN int = AY potential

given density and starting energy of \bar{K}
 \rightarrow G-matrix

AMD calculation
 \rightarrow density and starting energy of \bar{K}

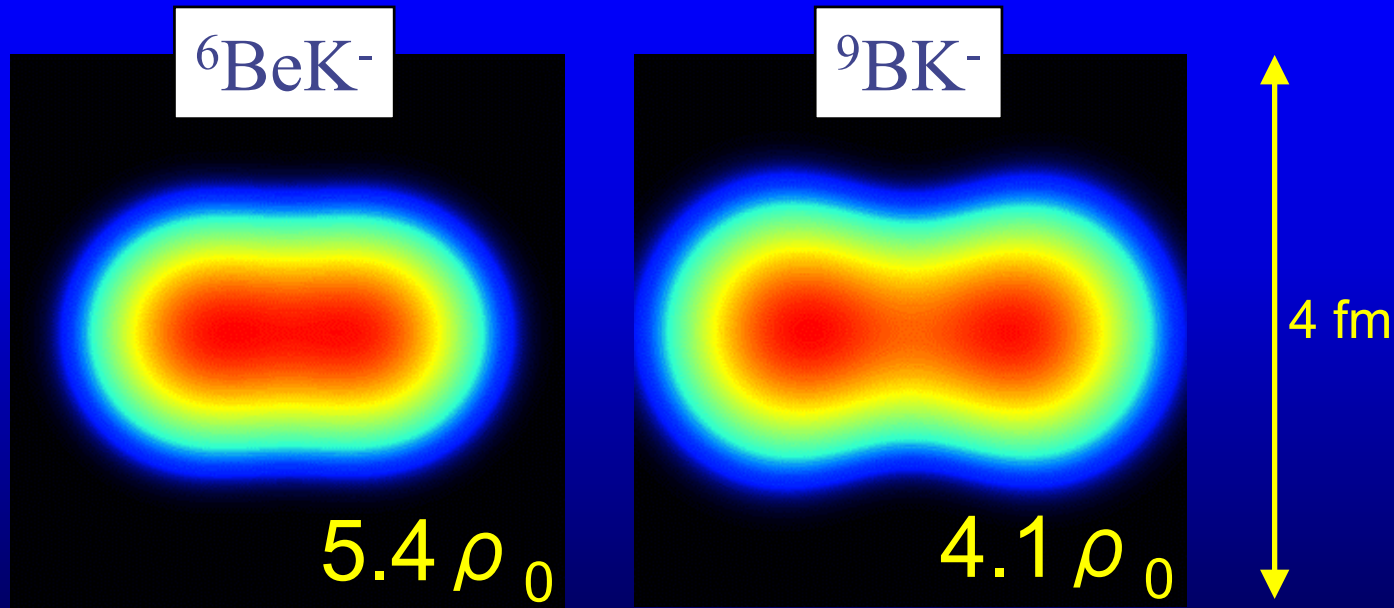
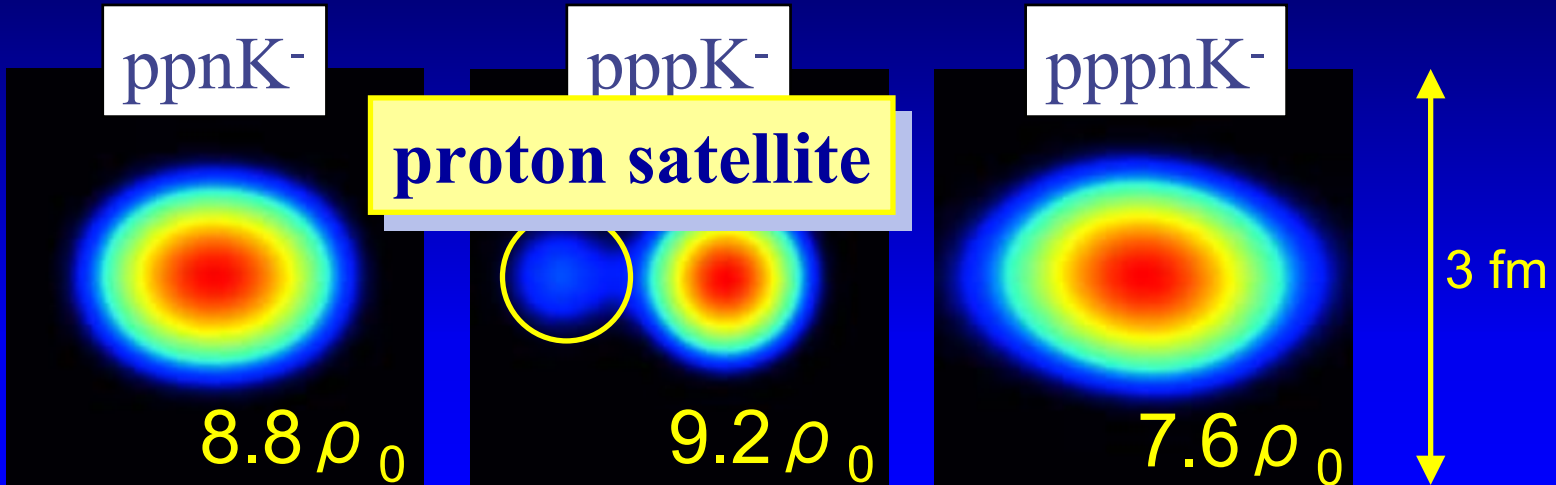
*Repeat until
getting consistency*

Binding energy of K^- and Decay-width

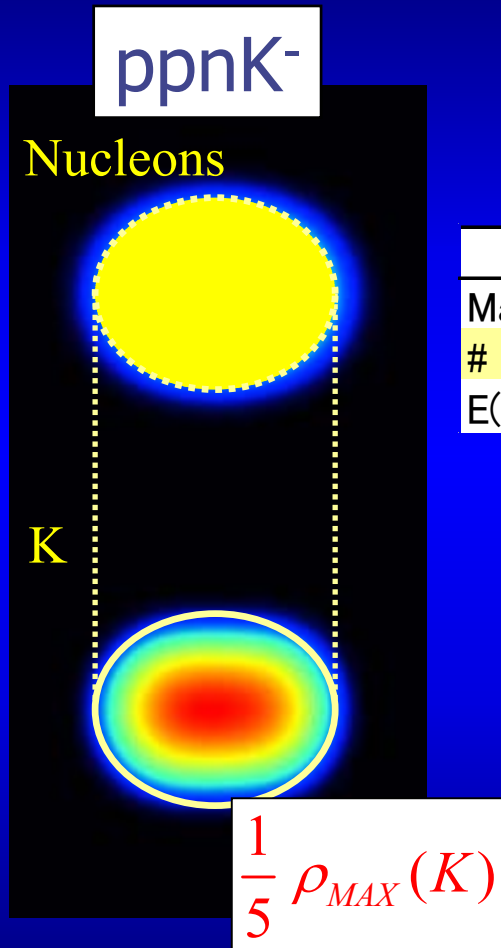


Uncertainty due to the inconsistency between the obtained results and the G-matrix used in the calculation

Nucleon density distribution



Number of nucleons near K⁻ meson



one center like

two center like

	ppnK	pppK	pppnK	6BeK	9BK
Max Kdens	0.67	1.48	0.62	0.34	0.35
# of nucleons	1.67	1.14	1.78	2.55	2.53
E(K)	110.3	96.7	105.0	104.2	117.0

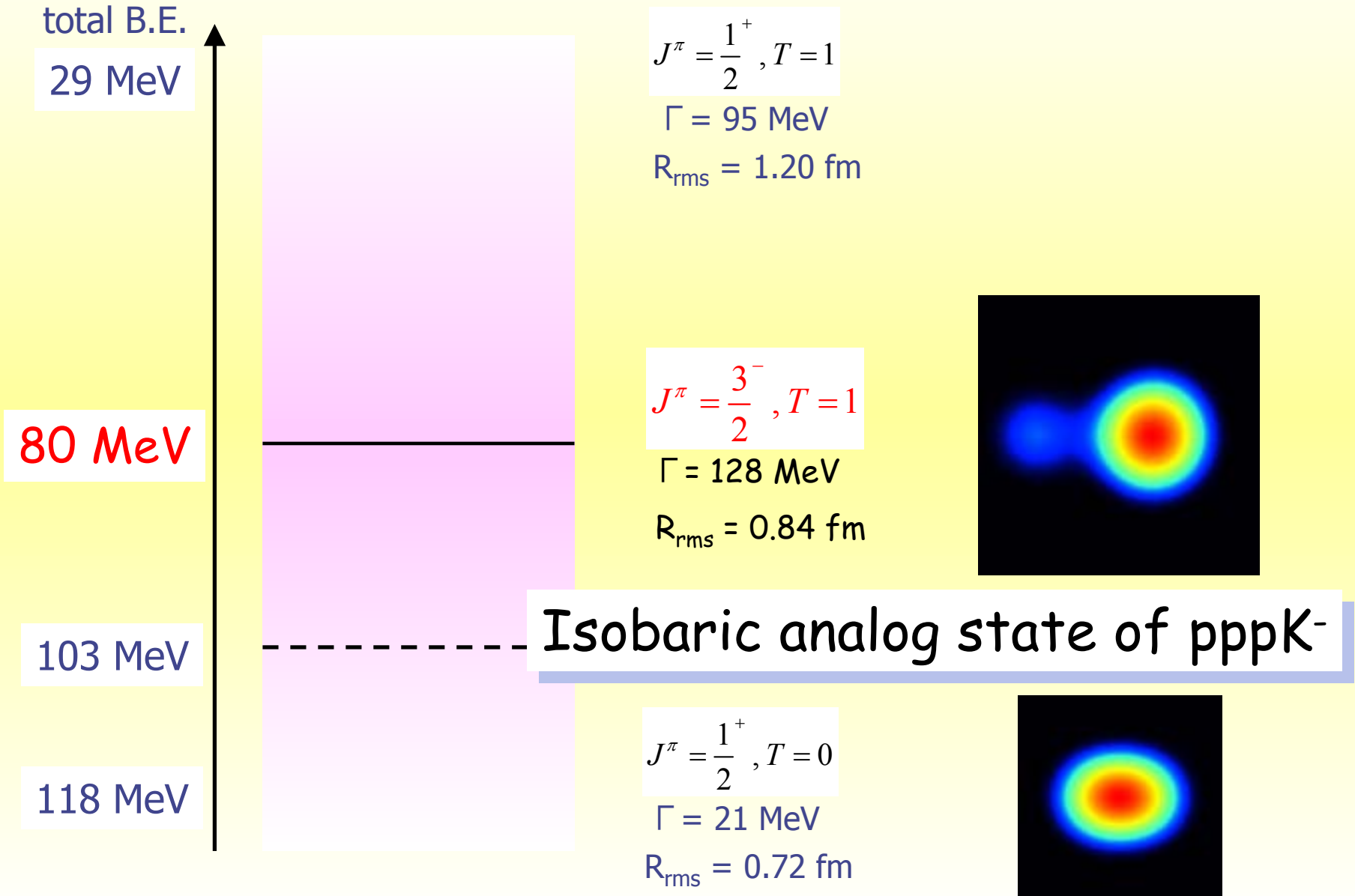
strange structure

Single K⁻ meson can interact with limited numbers of nucleons?



Saturation of E(K)

Excited state of ppnK⁻



Ground and Excited states of $^{11}\text{C}\text{K}^-$

total B.E. [MeV]

169 ----- $\Sigma \pi$ threshold

176 ----- $^9\text{BK}^- + \text{deuteron}$

180

Excited

$$J^\pi = \frac{1}{2}^+, T = 1$$

$$\Gamma = 49 \text{ MeV}$$

$$R_{\text{rms}} = 1.50 \text{ fm}$$

$$\rho_{\text{max}} = 4.8 \rho_0$$

182

$$J^\pi = \frac{1}{2}^+, T = 0$$

$$\Gamma = 46 \text{ MeV}$$

$$R_{\text{rms}} = 1.53 \text{ fm}$$

$$\rho_{\text{max}} = 4.8 \rho_0$$

191

Ground

$$J^\pi = \frac{3}{2}^-, T = 0$$

$$\Gamma = 46 \text{ MeV}$$

$$R_{\text{rms}} = 1.48 \text{ fm}$$

$$\rho_{\text{max}} = 4.8 \rho_0$$

$E(K) = 117 \text{ MeV}$

Excited states are also below the $\Sigma \pi$ -emission threshold.
Decay to the only $\Lambda \pi$ channel.

Structure of $^{11}\text{CK}^-$

• Ground
191 MeV

$$J^\pi = \frac{3^-}{2}, T = 0$$

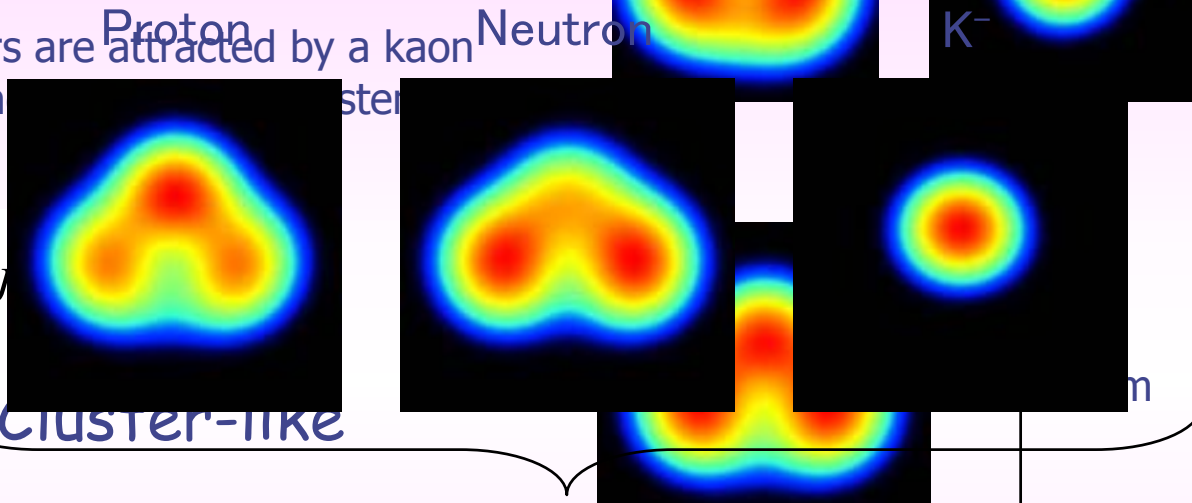
Three clusters are attracted by a kaon being at the center

Nucleons

Kaon

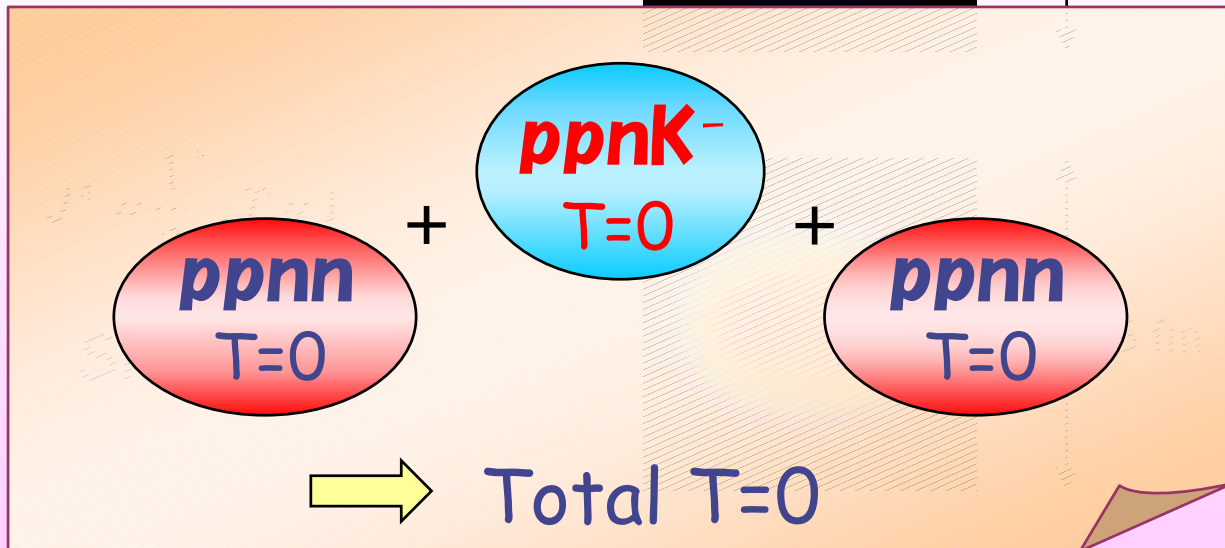
4 fm

• Excited
182 MeV



Cluster-like

• Excited
180 MeV



Summary

- We have improved AMD so that we can treat K^-p/\bar{K}^0n mixing and perform J & T projections.

- Results:

	E(K) [MeV]	width [MeV]	Max ρ [fm ⁻³]	Rrms [fm]
ppnK ⁻	110.3	21.2	1.50	0.72
pppK ⁻	96.7	12.5	1.56	0.81
pppnK ⁻	105.0	25.9	1.29	0.97
6BeK ⁻	104.2	33.3	0.91	1.17
9BK ⁻	118.5	33.0	0.71	1.45
11CK ⁻	117.4	46.0	0.81	1.48

In the ground state of kaonic nuclei, K^- is deeply bound by ~ 100 MeV and forms highly dense state.

Saturation of E(K) is related to the number of nucleons with which a K^- can interact.

- Excited states

- ppnK⁻ $J^\pi=3/2^-$, T=1 : Isobaric analog state of the ground state of pppK⁻ 38 MeV above the ground state, Γ is very large to be 128 MeV.
- ¹¹CK⁻ Two excited states as well as the ground state are below the $\Sigma \pi$ threshold.
 - $J^\pi=1/2^+$, T=1 : shell-like structure
 - $J^\pi=1/2^+$, T=0 : cluster-like structure

⁴He(T=0) + ppnK⁻(T=0) + ⁴He(T=0) configuration.