

Strange nuclear structures with high density formed by single/double K^- meson

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1. Introduction
2. New framework of Antisymmetrized Molecular Dynamics
 - $K^-p-\bar{K}^0n$ mixing
 - J & T projections
3. Results
ppn K^- , ppp K^- , pppn K^- , ${}^6\text{Be}K^-$, ${}^9\text{B}K^-$ and ${}^{11}\text{C}K^-$
4. Results = extremely proton-rich nuclei =
pppp K^- , ppp K^-K^- and pppp K^-K^-
5. Summary & Future plan

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.

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

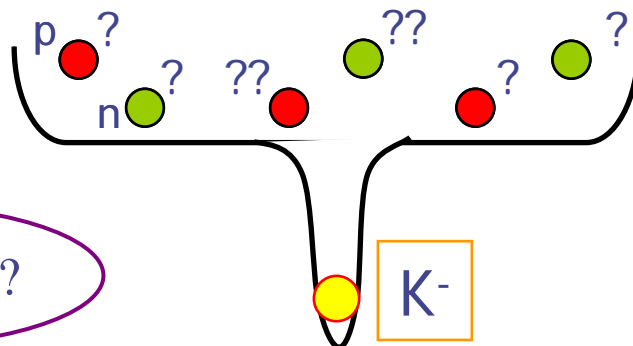
- Analysis of ${}^8\text{Be}K^-$ with $\alpha\alpha K^-$ model \Rightarrow

*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 ...

- unknown structure
- peculiar structure ex) isovector deformation

⇒ ☆ Method : **Antisymmetrized Molecular Dynamics**

no assumption on clusters, shape, etc.

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

Improving

• $\bar{K}^-p - \bar{K}^0n$ mixing



• Angular momentum & isospin projections

Wave function

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

Nucleon's wave function

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

$$|\sigma_{\alpha}^i\rangle = |\uparrow\rangle \text{ or } |\downarrow\rangle$$

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

p-n mixing

Anti-kaon's wave function

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

$$|\tau_{\alpha}^K\rangle = \left(\frac{1}{2} + \gamma_{\alpha}^K\right) |\bar{K}^0\rangle + \left(\frac{1}{2} - \gamma_{\alpha}^K\right) |K^-\rangle$$

\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

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

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

J & T projections (VBP)

After obtaining Φ^\pm by Tz-projected AMD cooling, it is projected onto the eigen state of the angular momentum J and the isospin T.

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

J projection

T projection

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

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

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

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*

Results — ppnK⁻ —

Model space

2 Gaussian / nucleon
5 Gaussian / kaon

Assumption of G-matrix

$E(K)=112.8\text{MeV}$, central density= 1.50fm^{-3}

Results

Projecting onto $J=1/2$ and $T=0$. Parity is positive.

	F(K)	Width	$\rho(0)$	Rrms
JT projection	110.3	21.2	1.50	0.72
simple AMD	105.2	23.7	1.39	0.72
BHF	108	20		0.97

G-matrix consistency : OK !

Quantum numbers

	After	Before
$J(J+1)$	0.75	1.36
\parallel	0.78	1.22
$\frac{1}{2} \cdot \left(\frac{1}{2} + 1\right)$	0.75	1.22
J2 (total sys.)	0.75	1.36
J2 (N sys.)	0.78	1.22
L2 (N)	0.03	0.14
S2 (N)	0.03	0.14
L2 (kaon)	0.03	0.14
T2	0.00	0.02
Tz	0.00	0.00

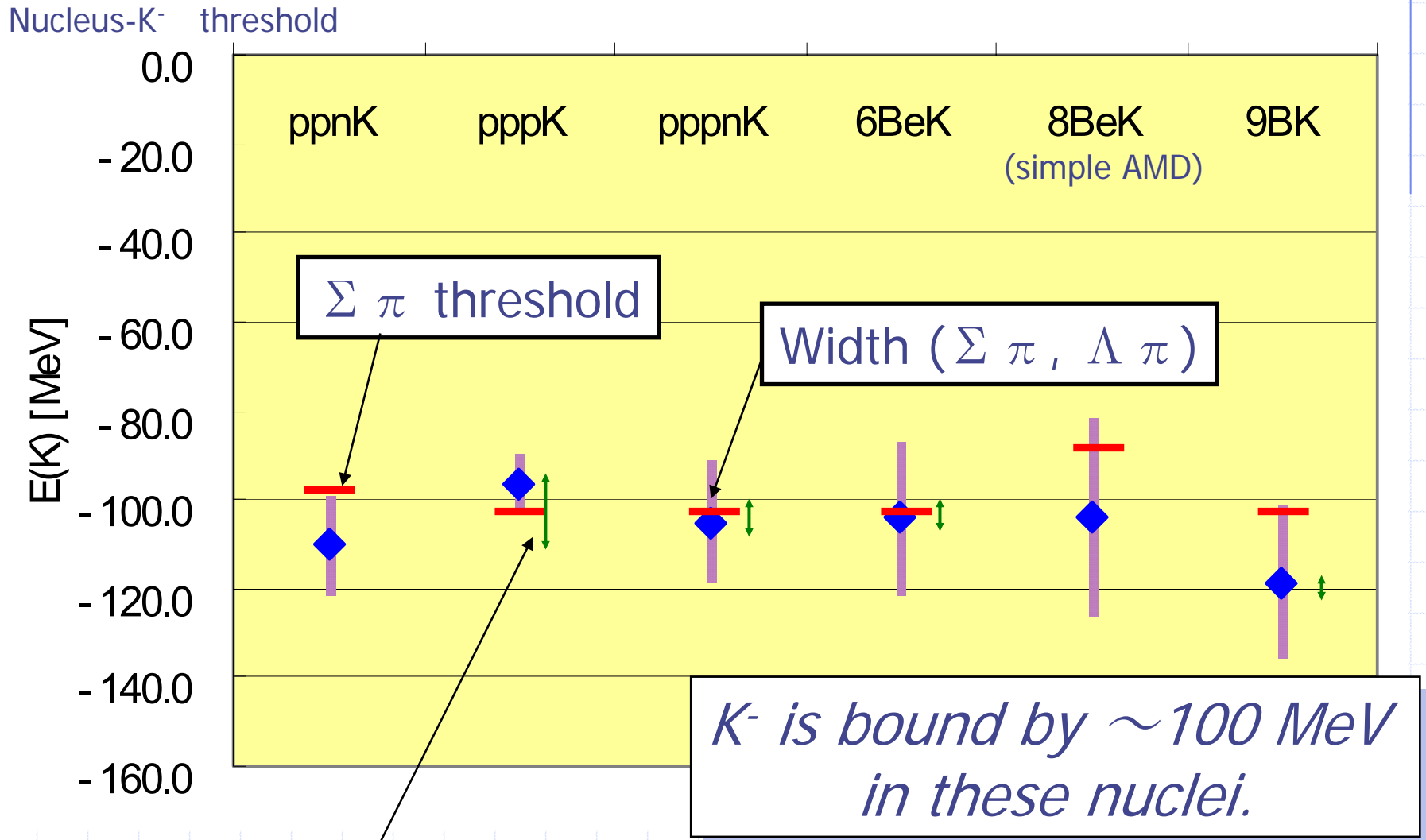
JT projection : OK !

Particle numbers

Proton	1.51
Neutron	1.49
K-	0.51
K0bar	0.49

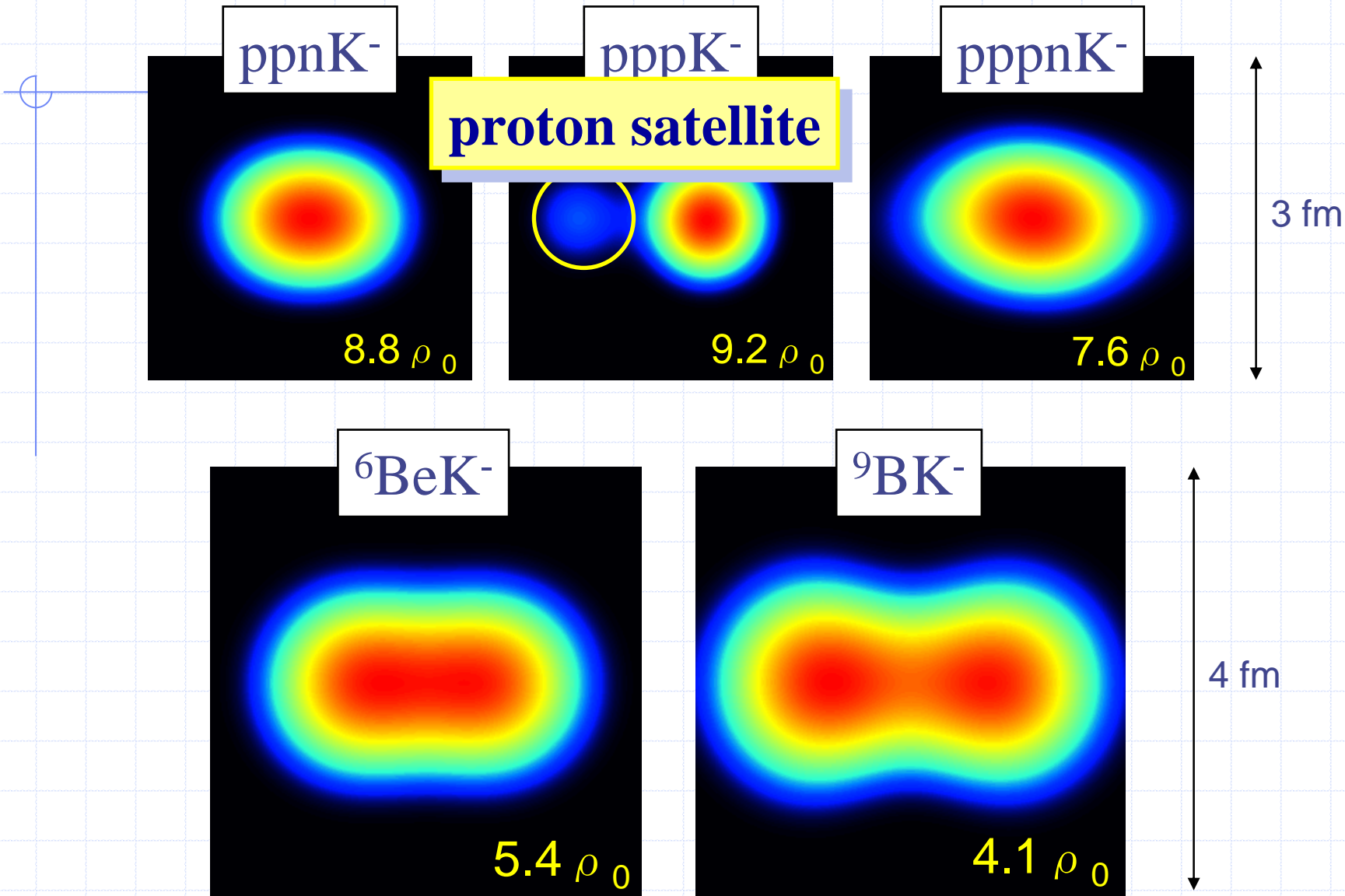
$\text{ppnK}^- : \text{pnn}\overline{\text{K}}^0$
= 1:1

Binding energy of K^- and width

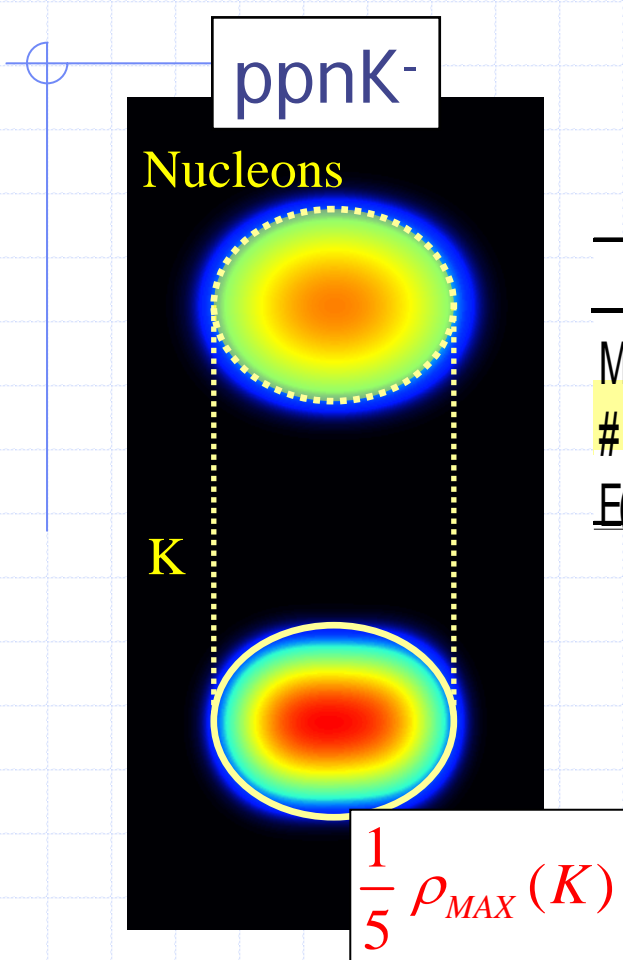


Uncertainty due to the inconsistency between the obtained results and the G-matrix used in the calculation

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

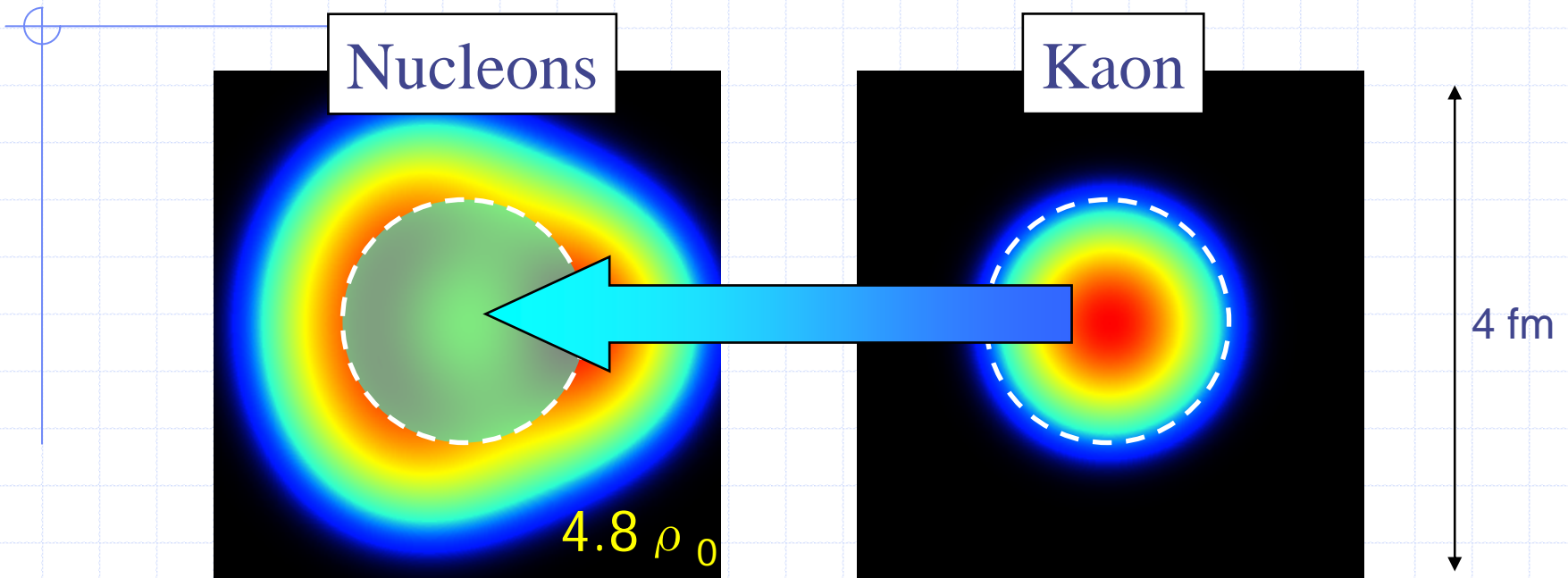
strange structure

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



Saturation of E(K)

Larger nucleus case = $^{11}\text{CK}^-$ =



Results

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

$$\Gamma = 37 \text{ MeV}$$

$$R_{\text{rms}} = 1.49 \text{ fm}$$

LS force : off

A K^- meson stays at the center of the system.

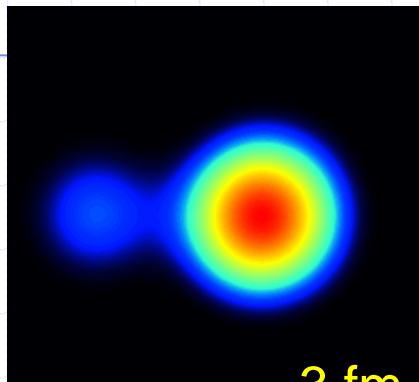


It attracts three clusters.

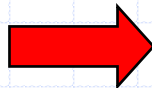
Results of extremely proton-rich kaonic nuclei

pppK⁻

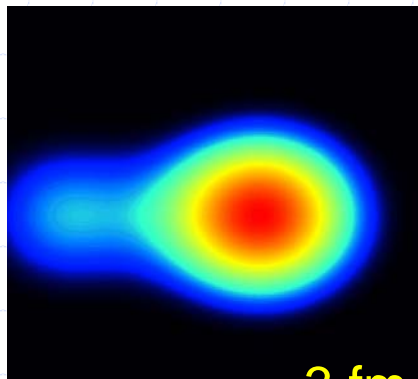
total B.E. 96.7 MeV
 E(K) 96.7 MeV
 Γ 12.5 MeV
 Rrms 0.81 fm



Proton



added



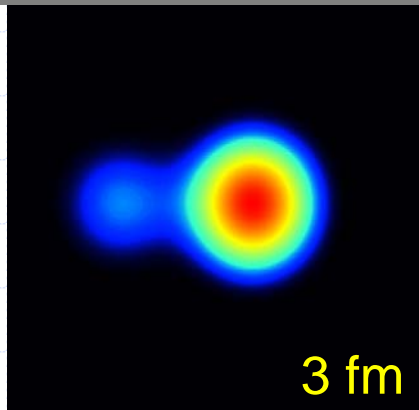
ppppK⁻

total B.E. 75.1 MeV
 E(K) 75.1 MeV
 Γ 161.8 MeV
 Rrms 0.92 fm

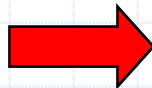
*These kaonic nuclei have proton-satellite structure.
 But not so deeply bound.*

ppppK⁻K⁻

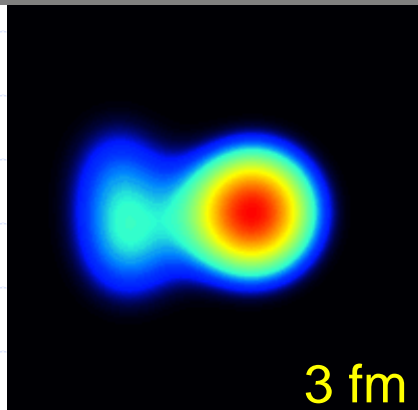
total B.E. 205.7 MeV
 E(K) 102.8 MeV
 Γ 51.9 MeV
 Rrms 0.65 fm
 ρ_{\max} 21.6 ρ_0



Proton



added



pppppK⁻K⁻

total B.E. 217.3 MeV
 E(K) 108.7 MeV
 Γ 41.2 MeV
 Rrms 0.73 fm
 ρ_{\max} 14.1 ρ_0

tentative

Summary & Future plan

- We have improved AMD so that we can treat K^-p/\bar{K}^0n mixing and perform J & T projections.
- We have calculated various kaonic nuclei – $ppnK^-$, $pppK^-$, $pppnK^-$, ${}^6\text{Be}K^-$, ${}^9\text{Be}K^-$ and ${}^{11}\text{C}K^-$.
Our results are ↓

	E(K) [MeV]	width [MeV]	ρ (0) [fm ⁻³]	Rrms [fm]	β	γ [deg]
$ppnK^-$	110.3	21.2	1.50	0.72	0.22	9.2
$pppK^-$	96.7	12.5	1.56	0.81	0.70	11.8
$pppnK^-$	105.0	25.9	1.29	0.97	0.54	3.8
${}^6\text{Be}K^-$	104.2	33.3	0.91	1.17	0.44	0.3
${}^9\text{Be}K^-$	118.5	33.0	0.71	1.45	0.46	20.8
${}^{11}\text{C}K^-$	117.0	37.0	0.82	1.49	0.36	46.4

*K^- is very deeply bound
and forms very highly dense state.*

- Saturation of E(K) is related to the number of nucleons with which a K^- can interact?
- Even if the KN interaction is very attractive, extremely proton-rich nuclei are not always deeply bound.

- ⊕ Excited states ?
- ⊕ KK interaction ?

Introduction

Remarks

$$v_{KN}^I(r) = v_D^I \exp\left[-(r/0.66 \text{ fm})^2\right],$$

$$v_{KN,\pi\Sigma}^I(r) = v_{C_1}^I \exp\left[-(r/0.66 \text{ fm})^2\right],$$

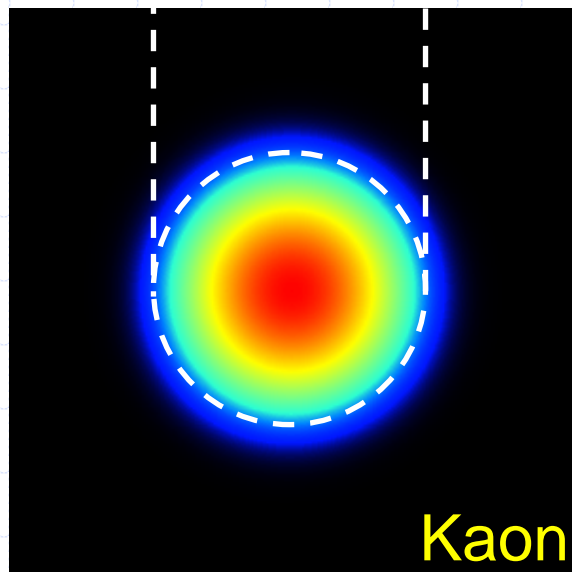
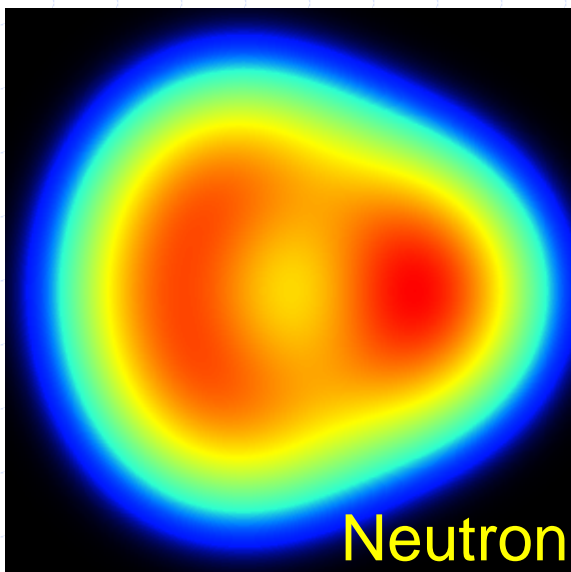
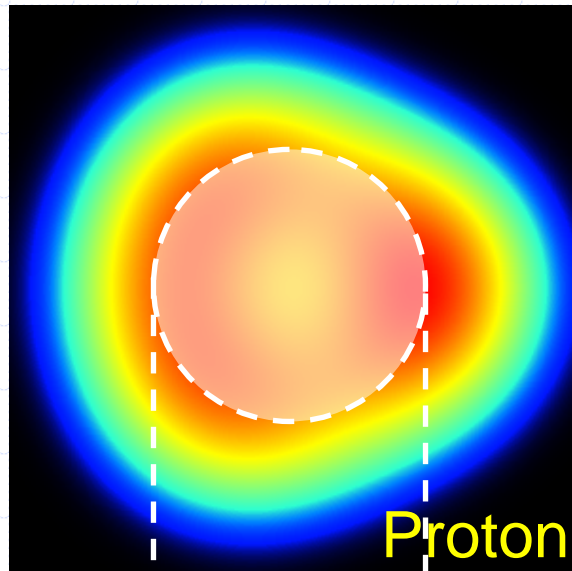
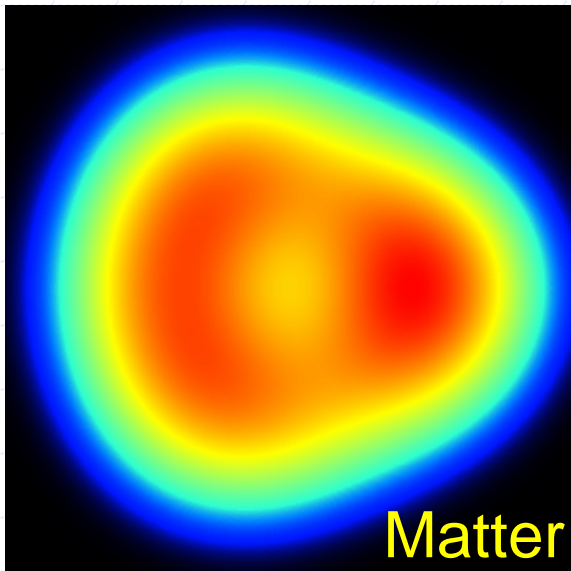
$$v_{KN,\pi\Lambda}^I(r) = v_{C_2}^I \exp\left[-(r/0.66 \text{ fm})^2\right]$$

$$v_D^{I=0} = -436 \text{ MeV}, \quad v_{C_1}^{I=0} = -412 \text{ MeV}, \quad v_{C_2}^{I=0} = \text{none}, \\ v_D^{I=1} = -62 \text{ MeV}, \quad v_{C_1}^{I=1} = -285 \text{ MeV}, \quad v_{C_2}^{I=1} = -285 \text{ MeV}.$$

Y. Akaishi and T. Yamazaki, PRC 65 (2002) 044005

- Not single channel, but **coupled** channel.
- Same property as KN interaction derived from Chiral theory.

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



$$\begin{aligned} E(K) &= 117 \text{ MeV} \\ \Gamma &= 37 \text{ MeV} \\ \rho_{\text{MAX}} &= 0.82 \text{ fm}^{-3} \\ R_{\text{rms}} &= 1.49 \text{ fm} \end{aligned}$$

LS force : off

size : 4fm x 4fm

Introduction

Studies of deeply bound kaonic nuclei by Akaishi & Yamazaki

- phenomenological $\bar{K}N$ potential

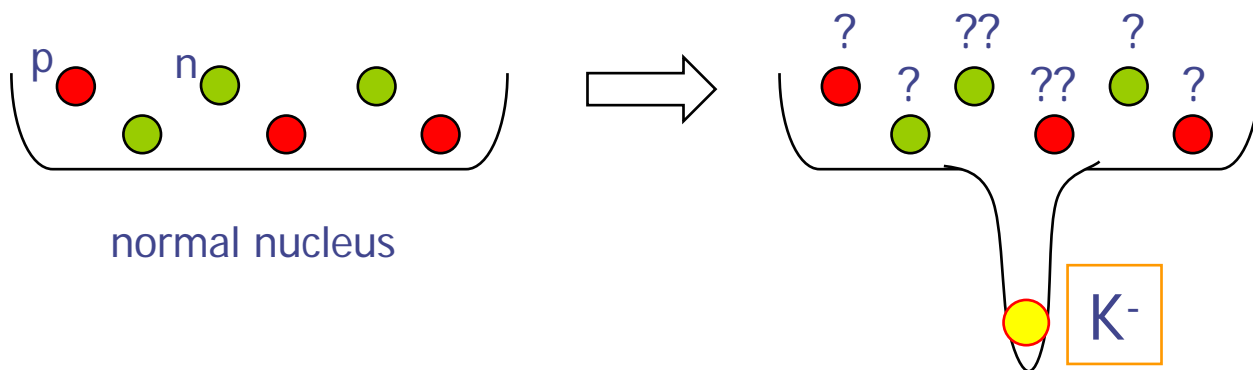
- free $\bar{K}N$ scattering data
- X-ray data of kaonic hydrogen atom
- binding energy and width of $\Lambda(1405)$

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 = $5 \rho_0$*

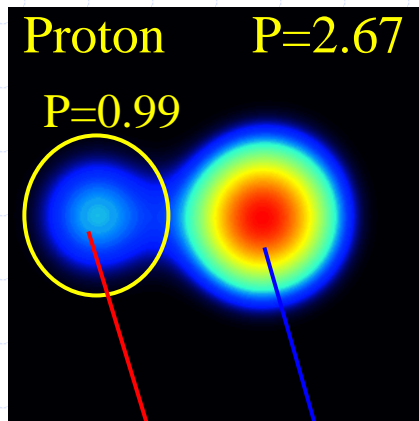


What kind of structure does A -nucleons + K^- system favor?

K^- deeply bound? highly dense? peculiar structure?

Proton satellite in pppK⁻

pppK⁻



$$\text{proton : } 2.66 = 1 + 1.66 = 1 + \frac{2}{3} \times 2 + \frac{1}{3} \times 1$$

$$\text{neutron : } 0.33 = \frac{1}{3} \times 1$$

$$\sqrt{\frac{2}{3}} |pp \otimes K^- \rangle + \sqrt{\frac{1}{3}} |pn \otimes \bar{K}^0 \rangle$$

$$|{}^3_{\bar{K}}\text{He}\rangle = |p\rangle \otimes |{}^2_{\bar{K}}\text{H}\rangle$$

Isospin

$$|1,1\rangle = |1/2, 1/2\rangle \otimes |1/2, 1/2\rangle$$

T, Tz

$$\sqrt{\frac{2}{3}} |1,1\rangle \otimes |1/2, -1/2\rangle + \sqrt{\frac{1}{3}} |1,0\rangle \otimes |1/2, +1/2\rangle$$

Clebsch-Gordan coefficient

Introduction

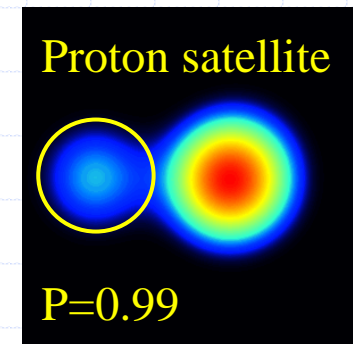
According to our study of various kaonic nuclei ...

- K^- is deeply bound in light nuclei.
- Since the main decay channel, $\Sigma \pi$, is closed, the deeply bound state is a discrete state.
- Due to the strong attraction from a K^- , nuclei are drastically shrunk to form dense state.
- Structures peculiar to kaonic nuclei are found.

$$E(K) \doteq 100 \text{ MeV.}$$

$$\rho = 4 \sim 8 \rho_0$$

1. Isovector deformation in ${}^8\text{Be}K^-$
2. Proton satellite in $\text{ppp}K^-$



Question

The strong attraction from K^- can make **extremely proton-rich nuclei** bound?
How are their structures?

Kaonic nucleiに期待できること

1、ハドロン物理の観点

K⁻が核子を引き寄せ、非常に高密度な状態が核内に。
高密度だが、割と冷たい。

- 高密度核物質、カイラル対称性の回復
- 高密度下でのNN/KN相互作用
- K凝縮、strange matter

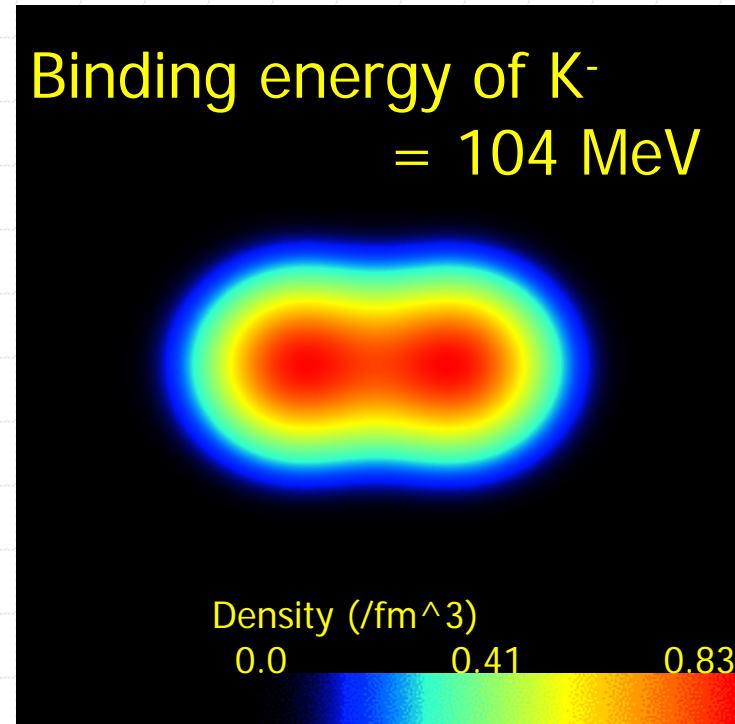
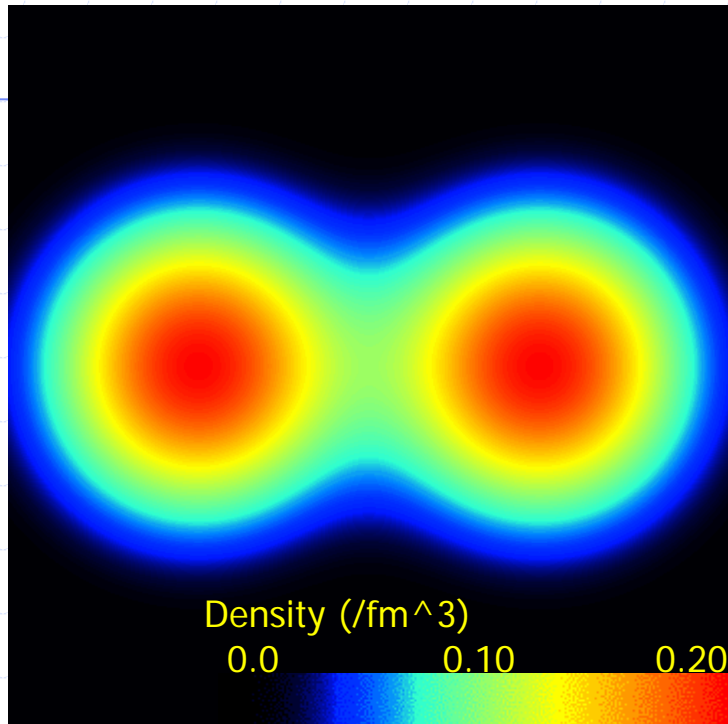
2、核構造の観点

K⁻からのアイソスピンに依存した強い引力

- 従来のクラスター構造が変化
 - • • 原子核がshrink
- isovector deformation
 - • • K⁻は陽子をより強くひきつける
- extreme proton-rich nuclei
 - • • p p K、p p p K、p p p p K、p p p K K、p p p p K Kなど

今までの原子核にない新しい構造様式？
多彩な構造？

Results // ^8Be and $^8\text{BeK}^-$ //



^8Be

rmsR = 2.46 fm
 β = 0.63
 Central density = 0.10

$/\text{fm}^3$

force: MV1 case 3

$^8\text{BeK}^-$

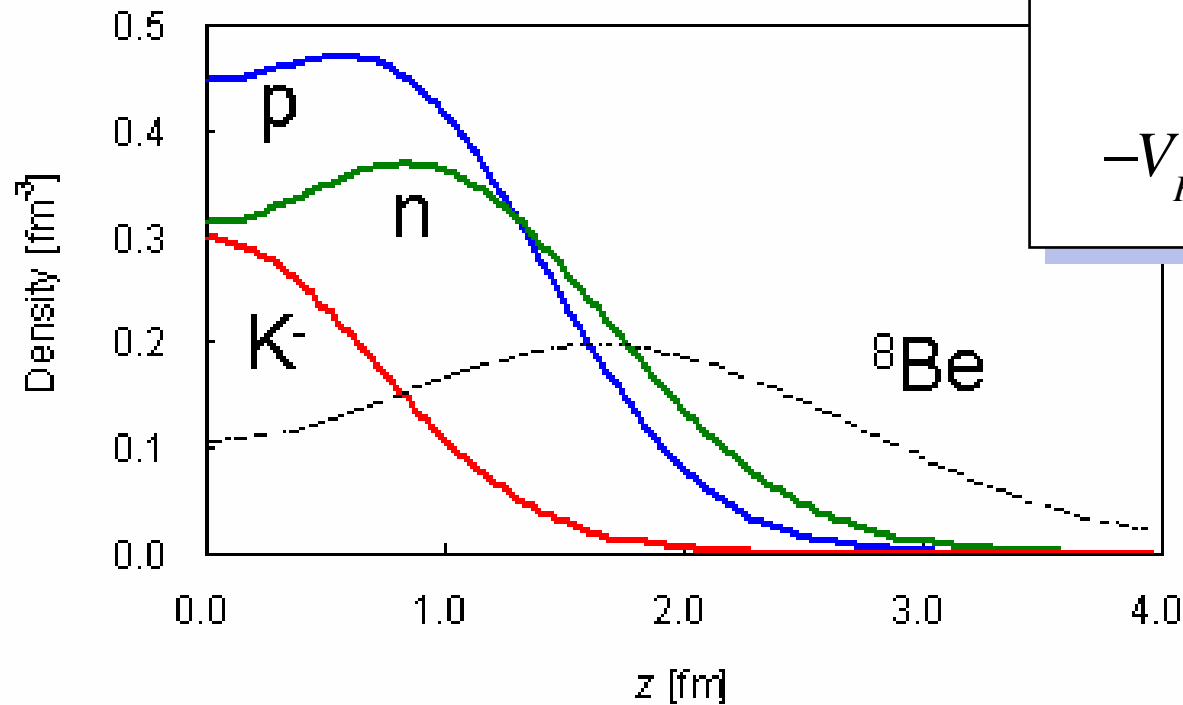
rmsR = 1.42 fm
 β = 0.55
 Central density = 0.76

$/\text{fm}^3$

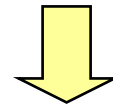
4.5 normal density

Isovector deformation in ${}^8\text{Be}K^-$

density distributions along Z-axis



$$-V_{I=0}^{\bar{K}N} \gg -V_{I=1}^{\bar{K}N}$$



$$-V_{K^-p} \gg -V_{K^-n}$$

New type of neutron skin

Neutrons are peripherally distributed, yet with quite high density.

Double kaonic nuclei

Motivation

- We have a simple question:
How do kaonic nuclei behave, if extra-one K^- is added?
- Multi K^- system is related to K^- condensation, strange quark matter, and so on.
Double kaonic nuclei, which contain two K^- 's, are the simplest case of Multi K^- systems.

Formalism

Wave func.: $|\Phi\rangle = \det[NN\cdots N] \otimes S[KK]$

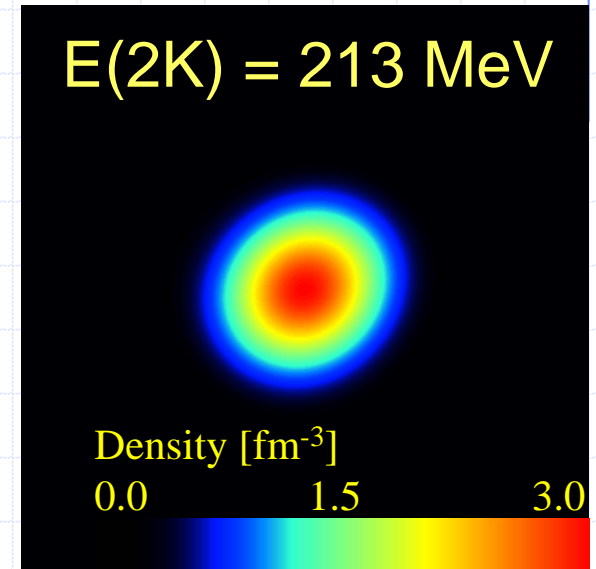
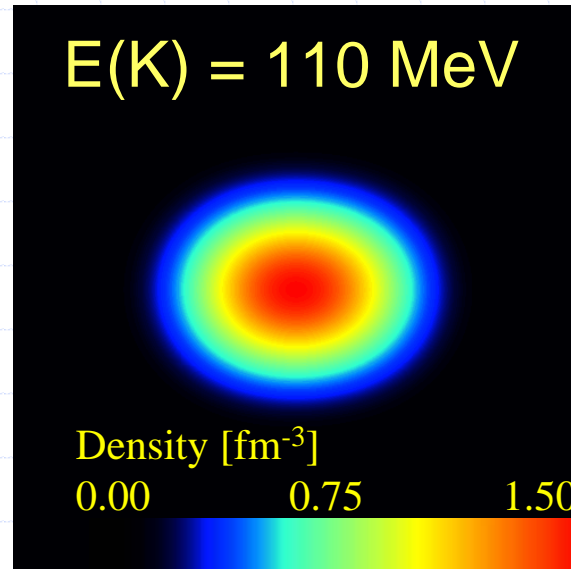
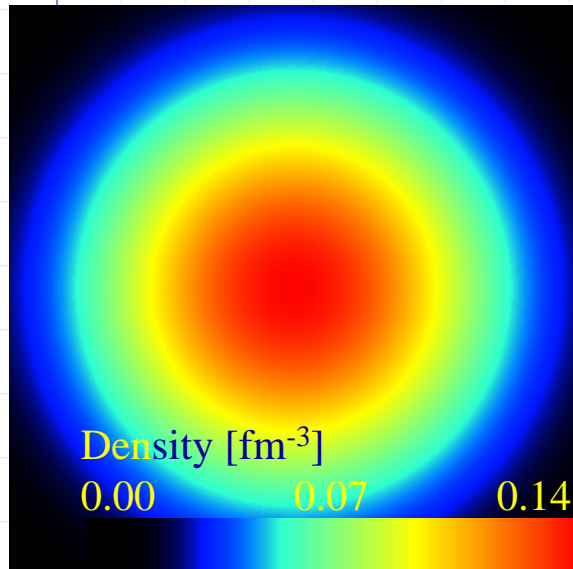
Symmetrized

calculate in quite the same way as single kaonic nuclei.

$V_{KK} = 0$ since we have no information in the present stage.

Double kaonic nucleus // ppnK-K- //

← 4 fm → ← 4 fm → ← 4 fm →



ppn

total B.E. = 6.0 MeV
central density = 0.14 fm^{-3}
 $R_{\text{rms}} = 1.59 \text{ fm}$

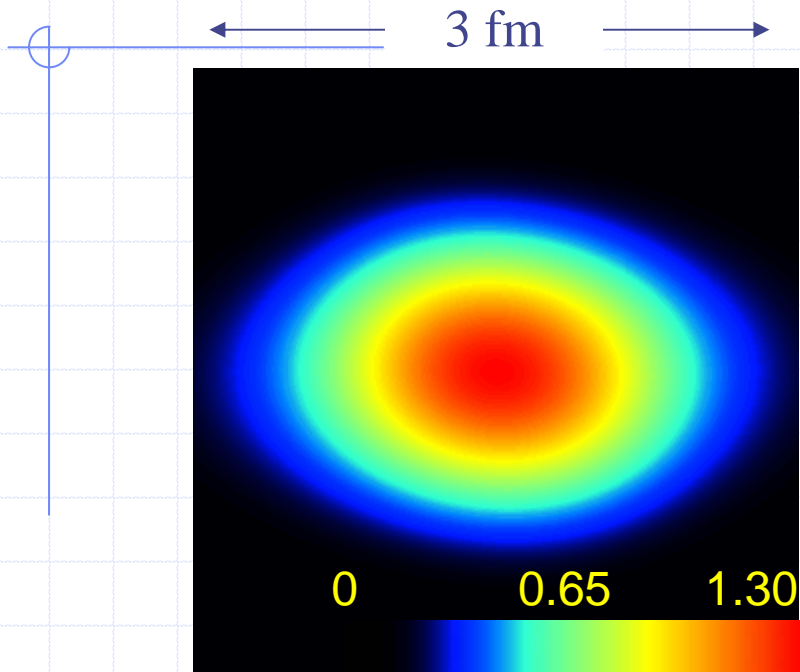
ppnK⁻

total B.E. = 118 MeV
central density = 1.50 fm^{-3}
 $R_{\text{rms}} = 0.72 \text{ fm}$

ppnK⁻K⁻

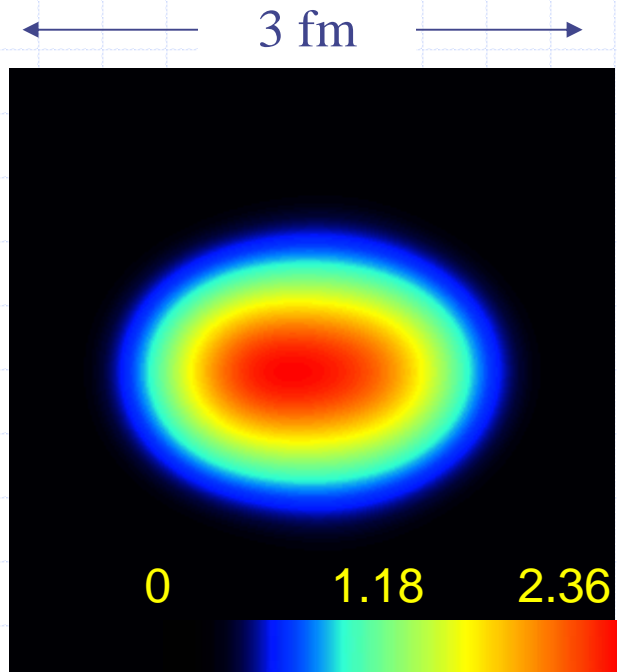
total B.E. = 221 MeV
central density = 3.01 fm^{-3}
 $R_{\text{rms}} = 0.69 \text{ fm}$

Double kaonic nucleus // pppnK-K- //



pppnK-

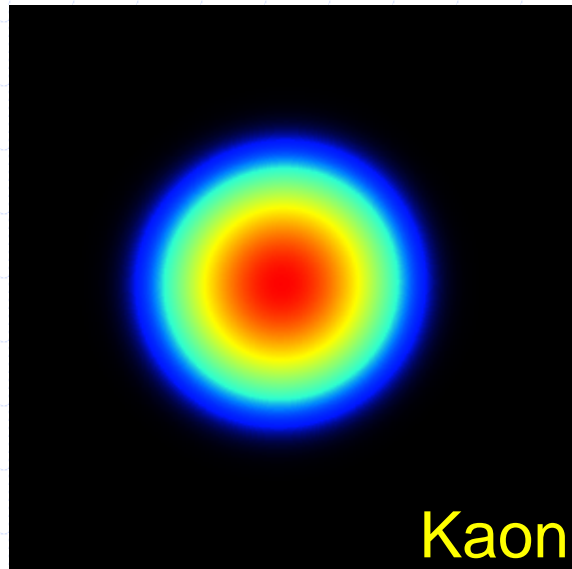
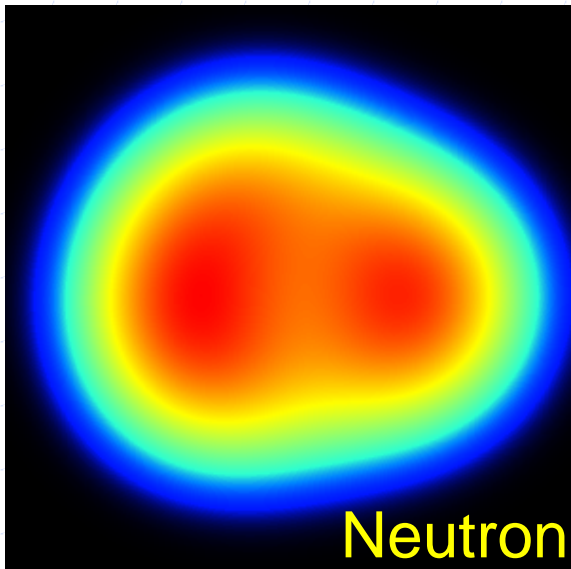
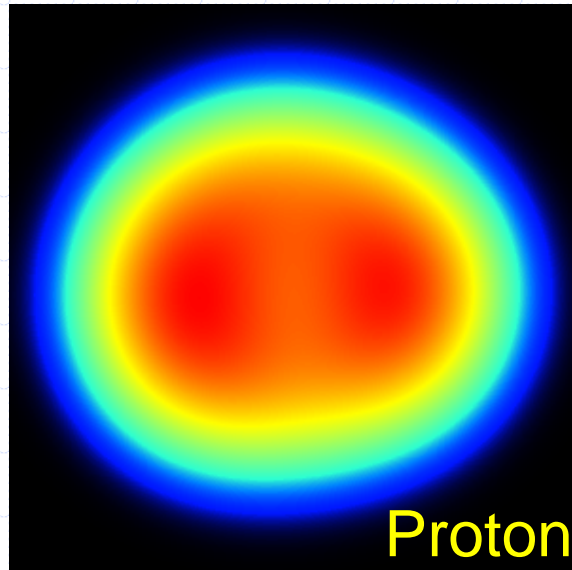
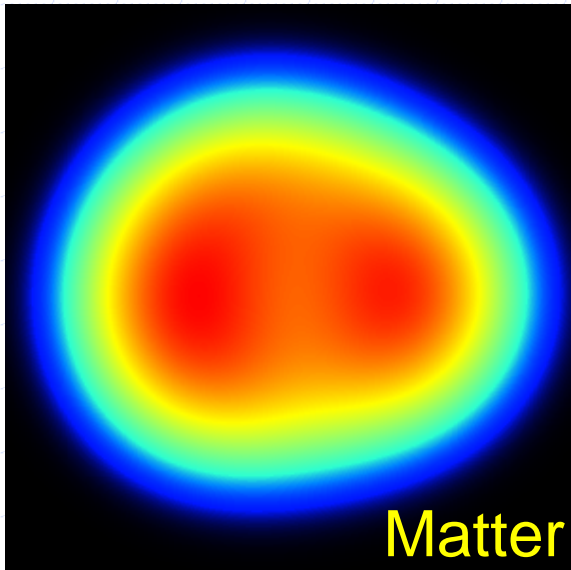
$E(K) = 105 \text{ MeV}$
central density = 1.29 fm^{-3}
 $R_{\text{rms}} = 0.97 \text{ fm}$



pppnK-K-

$E(2K) = 222 \text{ MeV}$
central density = 2.33 fm^{-3}
 $R_{\text{rms}} = 0.73 \text{ fm}$

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



$$\begin{aligned} E(K) &= 137 \text{ MeV} \\ \Gamma &= 33 \text{ MeV} \\ \rho_{\text{MAX}} &= 0.80 \text{ fm}^{-3} \\ R_{\text{rms}} &= 1.43 \text{ fm} \end{aligned}$$

LS force : on

size : 4fm x 4fm

$$\begin{aligned}
|{}^3_{\text{K}}\text{H}\rangle &= P_{T_Z=0} |\Phi_N \Phi_K\rangle \\
&= P_{T_Z=0} \left[\left(\sum_{m=-3/2}^{+3/2} P_{T_Z^N=m} \right) |\Phi_N\rangle \otimes \left(P_{T_Z^K=+1/2} + P_{T_Z^K=-1/2} \right) |\Phi_K\rangle \right] \\
&= \underbrace{P_{T_Z^N=+1/2} |\Phi_N\rangle \otimes P_{T_Z^K=-1/2} |\Phi_K\rangle}_{|ppn\text{K}^-\rangle} + \underbrace{P_{T_Z^N=-1/2} |\Phi_N\rangle \otimes P_{T_Z^K=+1/2} |\Phi_K\rangle}_{|pnn\bar{\text{K}}^0\rangle}
\end{aligned}$$

$$|ppn\text{K}^-\rangle$$

$$|pnn\bar{\text{K}}^0\rangle$$

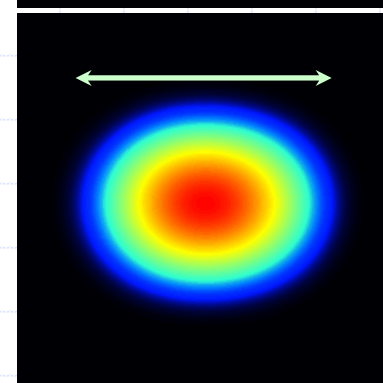
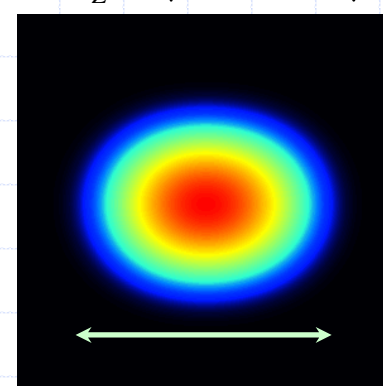
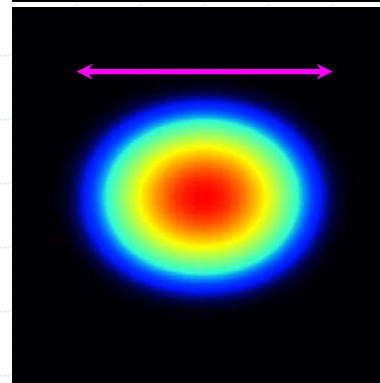
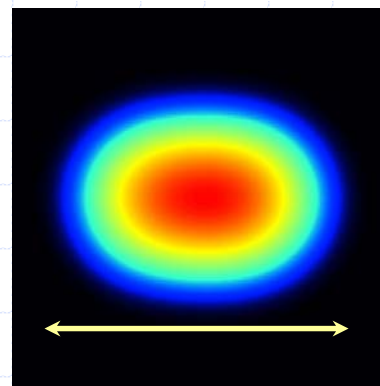
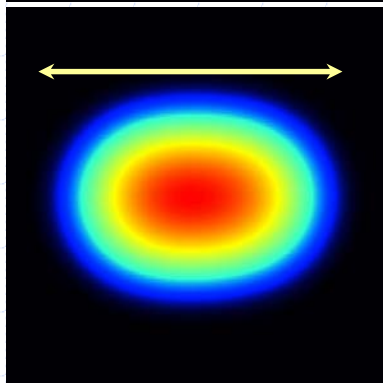
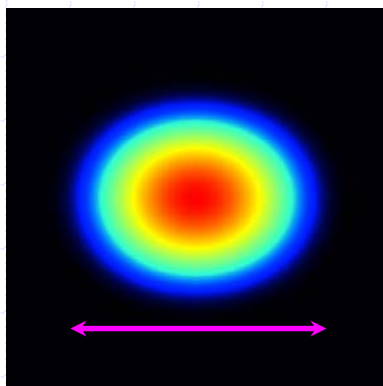
$$P_{T_Z^N=+1/2} |\Phi_N\rangle$$

$$P_{T_Z^N=-1/2} |\Phi_N\rangle$$

$$P_{T_Z=0} |\Phi_N \Phi_K\rangle$$

Proton

Neutron



$$\begin{aligned} \left| {}^3_{\text{K}}\text{H} \right\rangle &= P_{T_z=0} \left| \Phi_N \Phi_K \right\rangle ; \text{AMD wave function} \\ &= \left| \text{ppnK}^- \right\rangle + \left| \text{pnn}\bar{\text{K}}^0 \right\rangle \end{aligned}$$

Here,

$$\begin{aligned} {}_N \langle \text{ppnK}^- | {}^3_{\text{K}}\text{H} \rangle_N &= 0.71 \approx \frac{1}{\sqrt{2}} \\ {}_N \langle \text{pnn}\bar{\text{K}}^0 | {}^3_{\text{K}}\text{H} \rangle_N &= 0.70 \approx \frac{1}{\sqrt{2}} \end{aligned} \quad \Rightarrow \quad \left| {}^3_{\text{K}}\text{H} \right\rangle_N = \frac{1}{\sqrt{2}} \left| \text{ppnK}^- \right\rangle_N + \frac{1}{\sqrt{2}} \left| \text{pnn}\bar{\text{K}}^0 \right\rangle_N$$

normalized wave function

In addition,

$${}_N \langle \text{pnn}\bar{\text{K}}^0 | e^{i\pi\hat{T}_y} | \text{ppnK}^- \rangle_N = 0.999 \quad \Rightarrow \quad \left| \text{pnn}\bar{\text{K}}^0 \right\rangle_N = e^{i\pi\hat{T}_y} \left| \text{ppnK}^- \right\rangle_N$$

Therefore,

Intrinsic state in isospin space

$$\left| {}^3_{\text{K}}\text{H} \right\rangle_N = \sum_{\theta=0,\pi} \frac{1}{\sqrt{2}} e^{i\theta\hat{T}_y} \left| \text{ppnK}^- \right\rangle_N$$

holds good in AMD calculation of ${}^3_{\text{K}}\text{H}(T=0)$.

Treatment of K^- / \bar{K}^0 mixing

Usual practice = Coupled channel

$$\text{ex) } ppnK^- \quad |\Phi\rangle = \sum_a C_a |ppnK^-\rangle_a + \sum_b D_b |pnn\bar{K}^0\rangle_b$$

↔ Multi-Slater determinants

But some problems ...

1. common $\{\mathbf{Z}_i\}$ for all slater det., ${}_A C_Z + {}_A C_{Z-1}$ slater determinants !
2. different $\{\mathbf{Z}_i^{(a)}\}$ for each slater det., how many slater determinants ?
3. How is the effective frictional cooling for multi-slater det. ?
4. Calculation of non-diagonal matrix elements is somewhat tedious. $\langle ppnK^- | \hat{V} | pnn\bar{K}^0 \rangle$

Single slater determinat with charge-mixed s.p. wave func.

*In the single-particle state,
p and n are mixed, and
K- and K0bar are mixed.*

$$|N\rangle = a|p\rangle + b|n\rangle$$

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

$$|\Phi\rangle = |NNNK\rangle \text{ contains } ppnK^- \text{ and } pnn\bar{K}^0$$

But total Tz is restored with the Tz-projection.

Treatment of K^- / \bar{K}^0 mixing

Usual practice = Coupled channel

$$\text{ex) } ppnK^- |\Phi\rangle = C |ppnK^-\rangle + D |pnn\bar{K}^0\rangle$$

But in AMD treatment **non-diagonal matrix element** can't be calculated.

Calculation of $\langle ppnK^- | \hat{V} | pnn\bar{K}^0 \rangle$ needs the inverse matrix of the overlap matrix $B_{ij} = \langle \varphi_i | \varphi_j \rangle$

$$B = \begin{matrix} & \begin{matrix} p & p & n \end{matrix} \\ \begin{pmatrix} \alpha & \beta & 0 \\ 0 & \mathbf{0} & \gamma \\ 0 & 0 & \delta \end{pmatrix} & \begin{matrix} p \\ n \\ n \end{matrix} \end{matrix} \Rightarrow B^{-1} \text{ does not exist !}$$

In the Slater determinant $|\Phi\rangle = |NNNK\rangle$, $ppnK^-$ and $pnn\bar{K}^0$ are mixed.

*In the single-particle state,
p and n are mixed, and
K- and K0bar are mixed.*

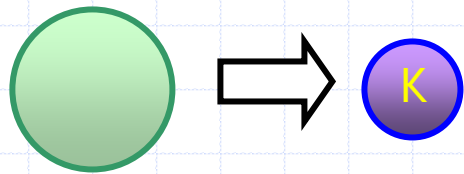
$$\begin{aligned} |N\rangle &= a|p\rangle + b|n\rangle \\ |K\rangle &= x|K^-\rangle + y|\bar{K}^0\rangle \end{aligned}$$

But total Tz is restored with the Tz-projection.

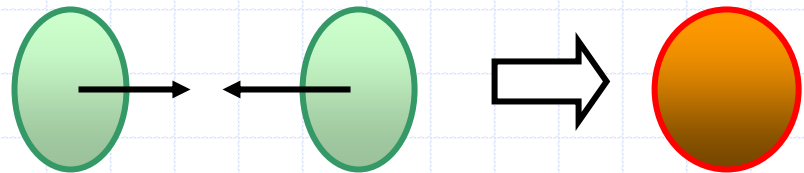
Introduction

If $\bar{K}N$ interaction is very strongly attractive...

- We can obtain a highly dense state by implanting K- into normal nuclei.

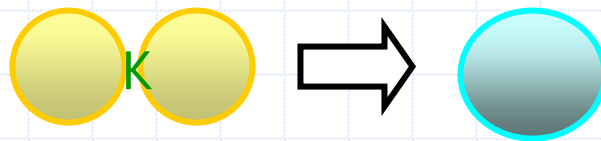


low temperature, high density

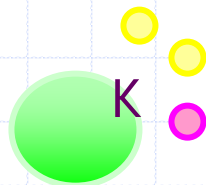


high temperature, high density

- Change of nuclear structure
: well-developed clustering structure vanishes ??

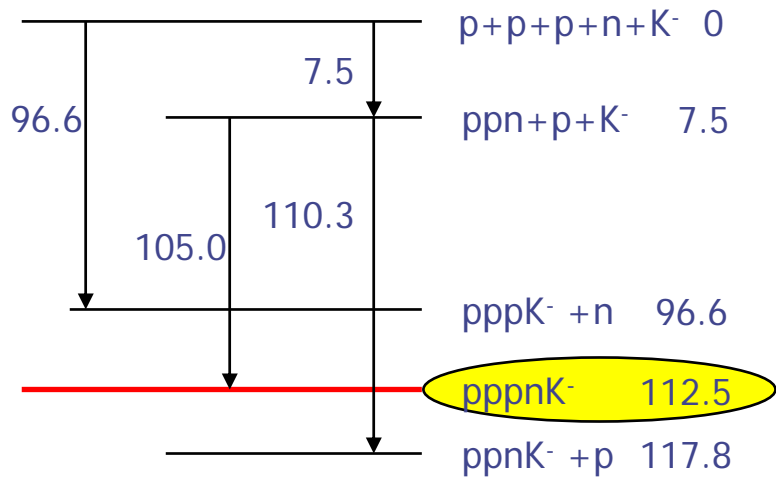


- Extending the drip line of unstable nuclei ???



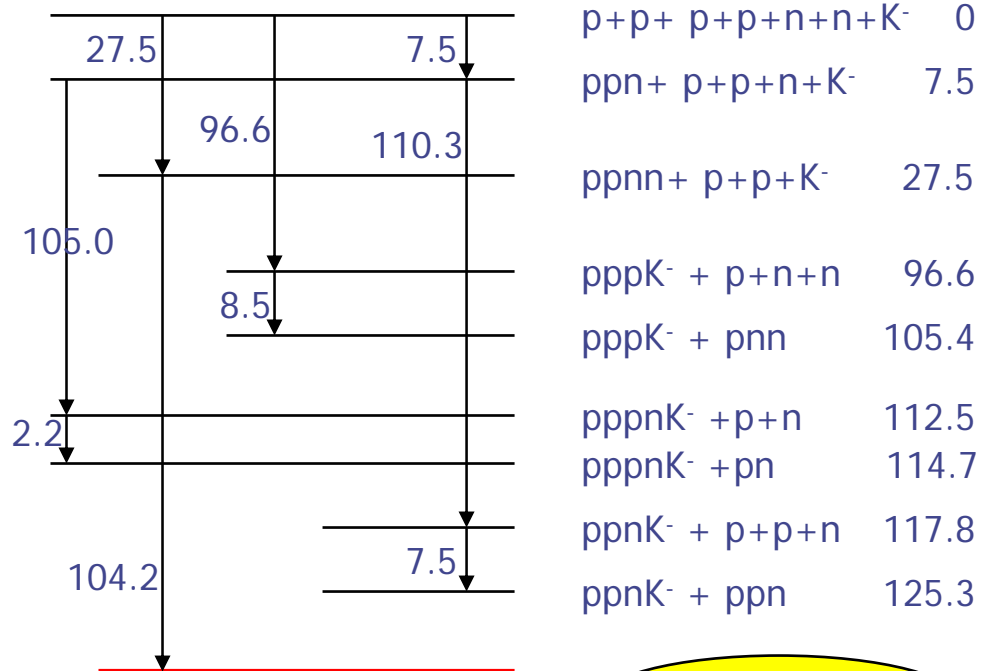
Stability for strong interaction

pppnK⁻



pppnK⁻ is unstable for strong interaction.

⁶BeK⁻



⁶BeK⁻ 131.7

⁶BeK⁻ is stable for strong interaction.

Formalism

- Hamiltonian

$$\hat{H} = \hat{T} + \hat{V}_C + \hat{V}_{Coul.} + \hat{V}_{KN} - \hat{T}_G$$

In kinetic energy part, taking the mass difference between a nucleon and K into consideration.

- Frictional cooling eq. with constraint

$$\dot{X}_\alpha^i = (\lambda + i\mu) \frac{1}{i\hbar} \left[\frac{\partial H}{\partial X_\alpha^{i*}} + \eta \frac{\partial W_G}{\partial X_\alpha^{i*}} \right] \text{ and c.c.}$$

negative

$$W_G = \left| \langle \mathbf{R}_G \rangle \right|^2 + \left| \langle \mathbf{P}_G \rangle \right|^2$$

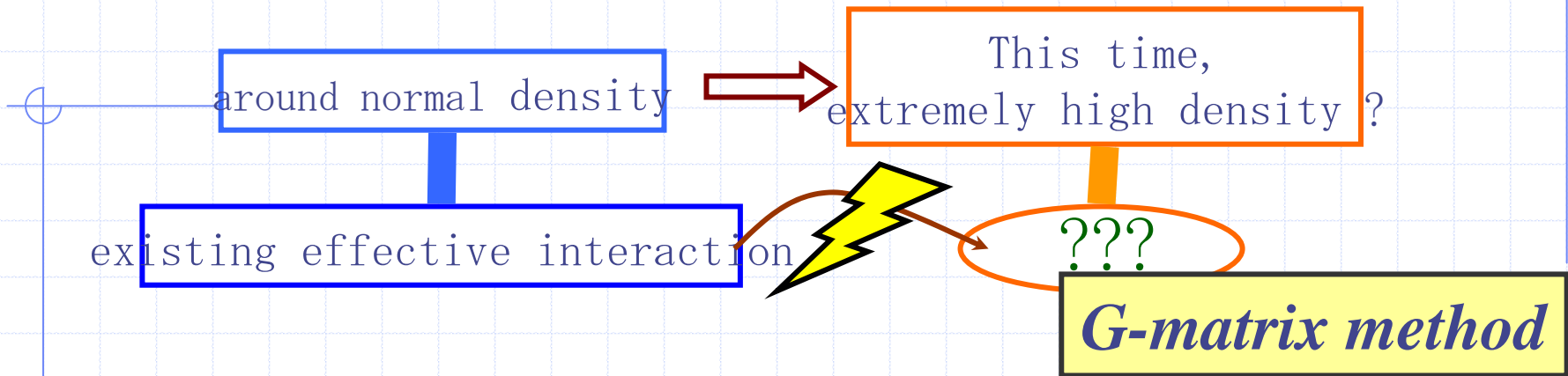
Constraint condition
The center of mass of the system is fixed to the origin.

$$\left\{ X_\alpha^i \right\} = \left\{ C_\alpha^i, \mathbf{Z}_\alpha^i, C_\alpha^K, \mathbf{Z}_\alpha^K \right\}$$

complex numbers

$$\mathbf{R}_G = \frac{\sum_{i=1}^A m_N \mathbf{r}_i + m_K \mathbf{r}_K}{Am_N + m_K}, \quad \mathbf{P}_G = \sum_{i=1}^A \mathbf{p}_i + \mathbf{p}_K$$

G-matrix calculation



- bare interaction and model space

NN int. : Tamagaki potential (OPEG)

KN int. : Akaishi-Yamazaki potential

model space : 2 Gaussian wave packets for one nucleon,
5 Gaussian wave packets for one anti-kaon.

- procedure

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

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

*Repeat
until getting
a consistent result.*

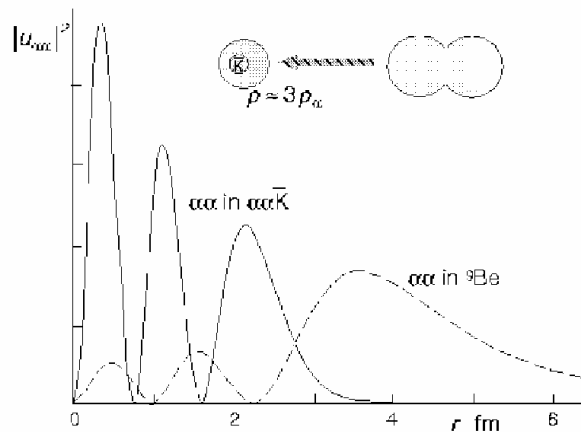
Introduction

Changing the matter density of ^8Be

Y.Akaishi solved $\alpha\alpha K^-$ system with ATMS method.

NN interaction = Hasegawa-Nagata No.1 which has 1.6 GeV repulsive core.

KN interaction = Y.A. & T.Y. force.



The distribution $|u_{\alpha\alpha}(r)|^2$ of the $\alpha\alpha$ relative motion in $\alpha\alpha K^-$ system.

Binding energy = 113 MeV
 $\Gamma_{\Delta+\pi}$ = 38 MeV
 Central density = $5 \rho_0$

K^- causes ^8Be to be shrunk.
 ^8Be implanted K^- loses two α structure.

Akaishi-Yamazaki $\bar{K}N$ interaction

1, free $\bar{K}N$ scattering data

A.D.Martin, Nucl.Phys.B179(1981)33

$$a^{l=0} = -1.76 + i 0.46 \text{ fm}, a^{l=1} = 0.37 + i 0.60 \text{ fm}$$

cf) A.D.Martin :

$$a^{l=0} = -1.70 + i 0.68 \text{ fm}, a^{l=1} = 0.37 + i 0.60 \text{ fm}$$

2, X-ray data of kaonic hydrogen atom

M.Iwasaki et al. Phys.Rev.Lett.78(1997)3067

T.M.Ito et al. Phys.Rev.C58(1998)2366

$$a_{K-p} = (a^{l=0} + a^{l=1}) / 2 = -0.70 + i 0.53 \text{ fm}$$

cf) T.M.Ito et al. :

$$a_{K-p} = (-0.78 \pm 0.15 \pm 0.03) + i (0.49 \pm 0.25 \pm 0.12) \text{ fm}$$

3, Binding energy and decay width of $\Lambda(1405)$

B.E. = -29.7 MeV, $\Gamma = 40$ MeV cf) $\Lambda(1405)$ exists 27 MeV below $K^- + p$ threshold

regarding $\Lambda(1405)$ as $l=0$ bound state of $\bar{K}N$

$V_{I=0}^{KN}$ is much more attractive than $V_{I=1}^{KN}$!!

り方

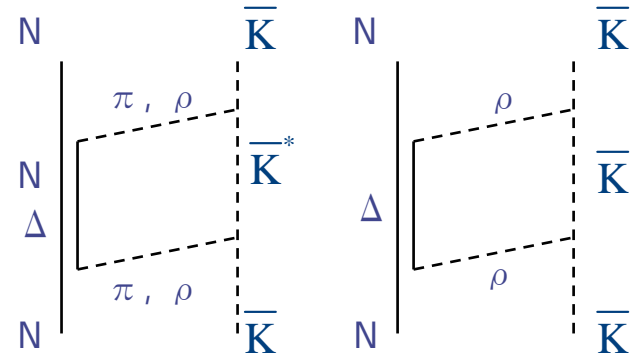
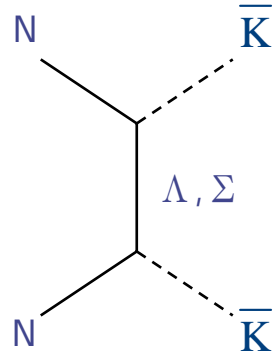
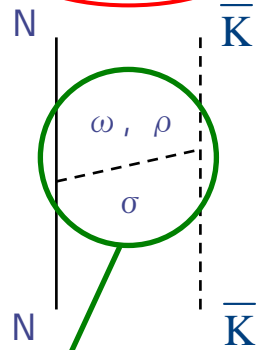
	W.Weise	Julichグループ (K. Holinde)	赤石・山崎
基礎理論	Chiral SU(3) effective Lagrangian	meson exchange	-----
目的	$\bar{K}N$ 系全体を合わす	NN,YN,KN, $\bar{K}N$ 全てを合わす	$\bar{K}N$ のbound stateにのみ着目
Channel	$pK^-, n\bar{K}^0, \Lambda\pi^0, \Sigma^0\pi^0, \Sigma^+\pi^-, \Sigma^-\pi^+$ KN threshold 下に $\Lambda(1405)$ と	(同左) + $N\bar{K}^*, \Delta\bar{K}, \Delta\bar{K}^*$	$pK^-, n\bar{K}^0, \Lambda\pi^0, \Sigma^0\pi^0, \Sigma^+\pi^-, \Sigma^-\pi^+$
ポテンシャルの導入	<p>いうresonance。カイラル摂動論で処理しようとする、ある種のダイヤグラムを無限オーダーとらねばならない。</p> <p style="text-align: center;">↓</p> <p>ポテンシャルで処理 q2まで取ったカイラルラグランジアンからBorn近似で求めたs波散乱長と等しい散乱長を持つようなポテンシャル。</p> <p>但しポテンシャルの形は仮定</p> <p>local : Yukawa型 separable : $\frac{\alpha_i^2}{\alpha_i^2 + k_i^2} \cdot \frac{\alpha_j^2}{\alpha_j^2 + k_j^2}$</p>	<p>相互作用ラグランジアン $W = -\int d^3x L_{int}$ から</p> <p>W : 2次 one boson exchange, resonance diagram W : 4次 box diagram のポテンシャル</p> <p>KN相互作用からG-parity変換で決まるものはそれを用いる。その他はfree parameter。</p> <p>vertexのform factorは仮定。</p>	<p>とにかくポテンシャルの形はGauss型としてしまえ！</p> <p>全チャンネルで $v_x^I = v_{cx}^I \exp[-(r/0.66)^2]$</p>
再現	<ul style="list-style-type: none"> $\Lambda(1405)$ の位置と幅 $\bar{K}N$低エネルギー散乱の全断面積 (K: 60~300 MeV/c in LAB) Branching ratio 	<ul style="list-style-type: none"> $\Lambda(1405)$ の位置と幅 $\bar{K}N$低エネルギー散乱の全断面積 (K: 60~300 MeV/c in LAB) 	<ul style="list-style-type: none"> $\Lambda(1405)$ の位置と幅 $\bar{K}N$低エネルギー散乱の散乱長 kaonic hydrogen atomのスペクトル <p>核物質中での散乱振幅の振る舞いはWeiseとconsistent</p>
論文	Nucl.Phys.A594(1995)325	Nucl.Phys.A513(1990)557	Phys.Rev.C65(2002)044005

Julich $\bar{K}N$ Quasi-potential

$$p_{\bar{K}}^{\text{LAB}} = 60 \sim 300 \text{ MeV}/c$$

Dominant

Minor



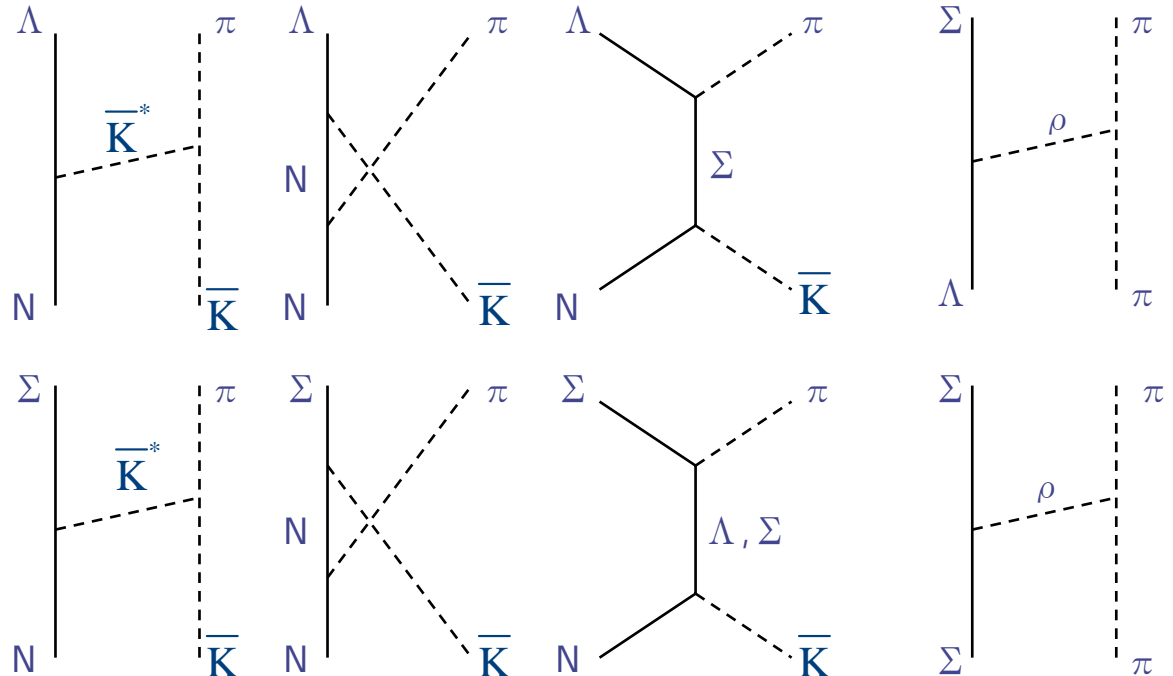
$\bar{K}N$ の場合、
全てのメソンが引力的に働く。



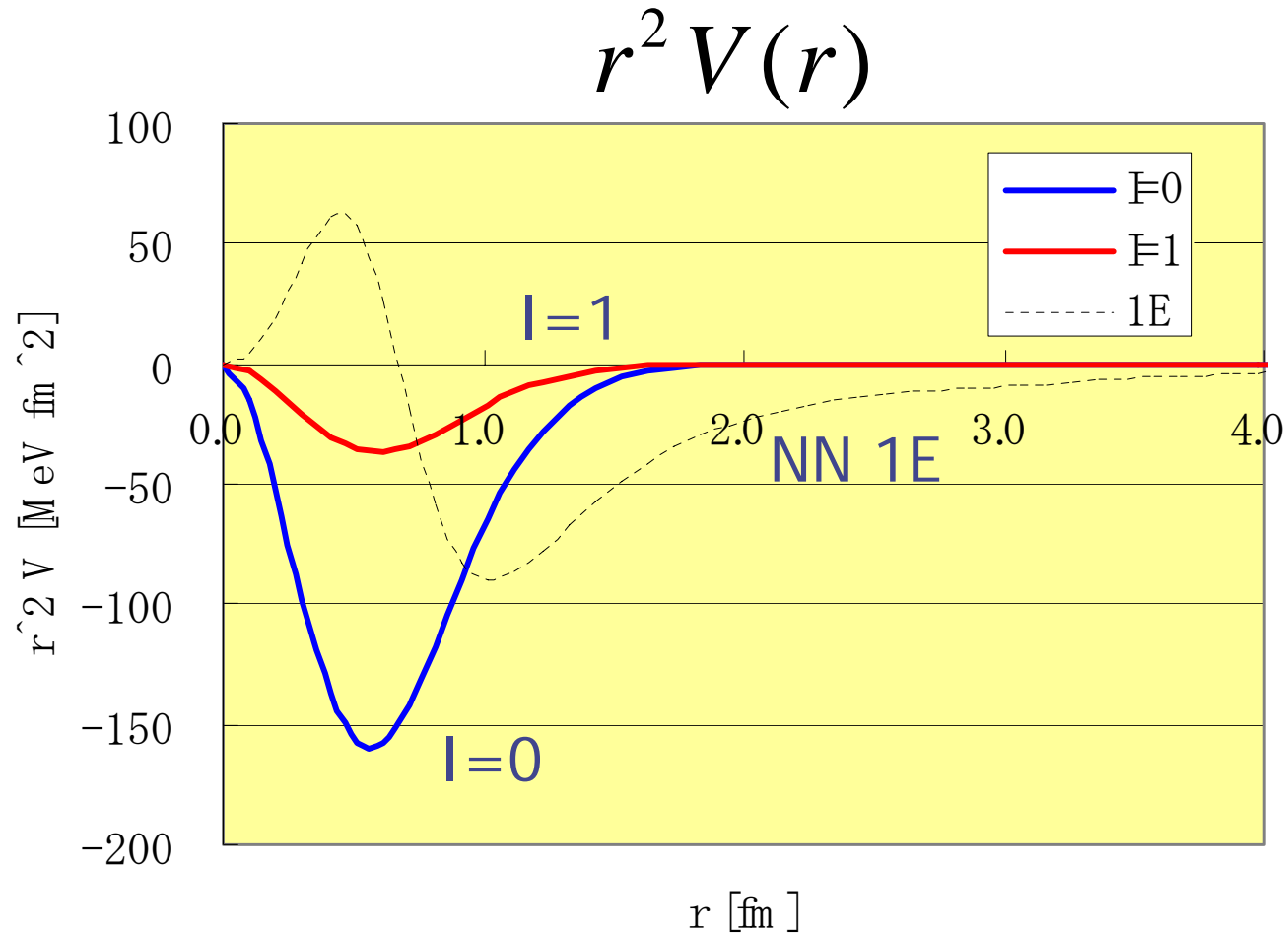
$\Lambda(1405)$ の形成

Weise とも consistent

local potentialを用いた場合、
レンジパラメータはこれらの
メソン質量に対応。
 $\Lambda(1405)$ を作る。

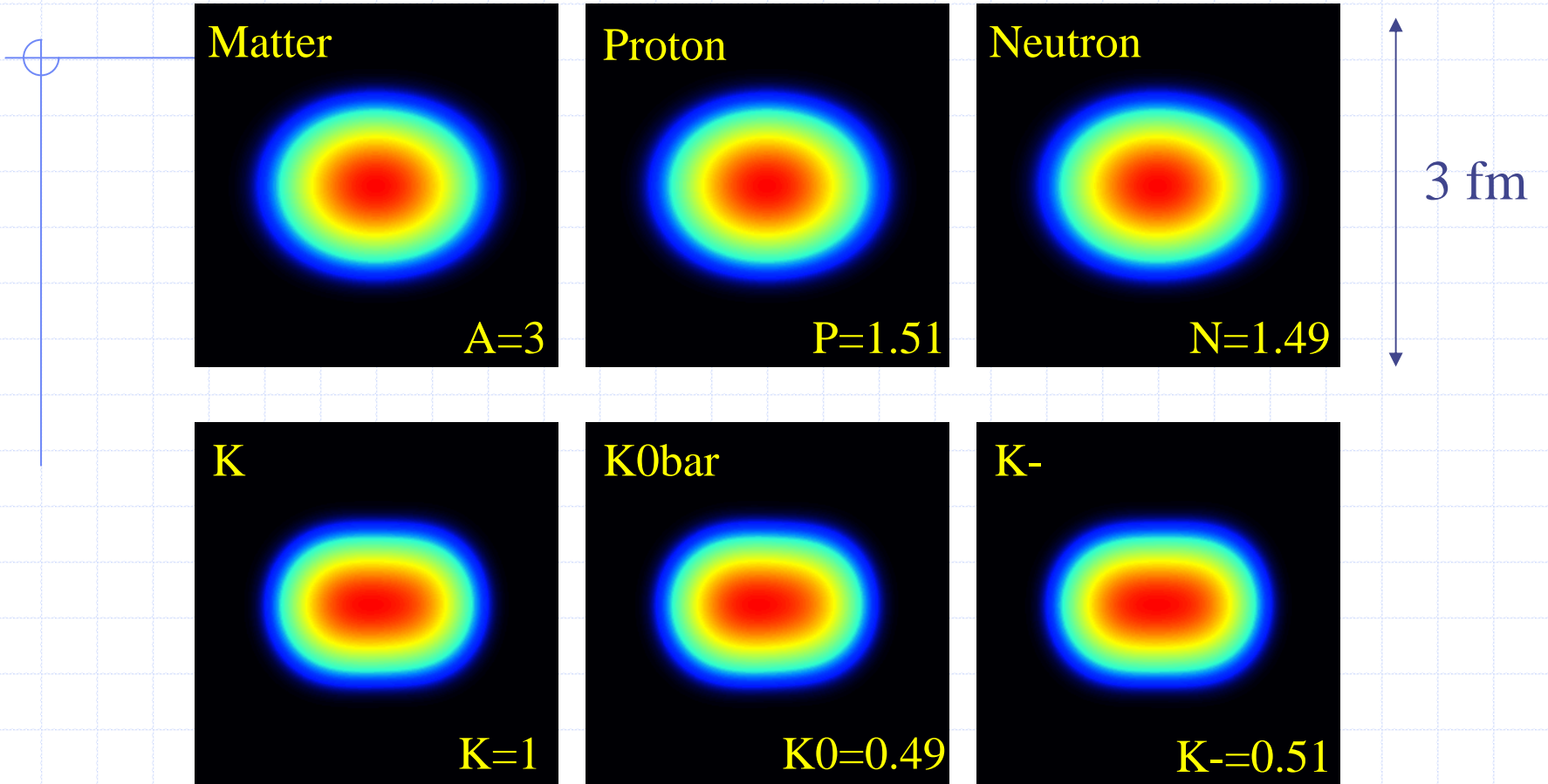


KN potential (G-matrix) for ${}^6\text{BeK}^-$



ppnK⁻

Density distributions of $|P_{Tz} \Phi\rangle$ in ZX plane.



$\rho(0) = 1.50 \text{ fm}^{-3}$... central density

$$\text{ppnK}^- : \text{pnn}\overline{\text{K}}^0 = 1 : 1$$

model space

2 Gaussian / nucleon
5 Gaussian / kaon

force

$E(K)=94.5\text{MeV}$, central density= 1.49fm^{-3}

Results

Projecting onto $J=3/2$ and $T=1$. Parity is negative.

	B F	width	dens0	rmsR	beta	gamma
JT projection	96.68	12.45	1.56	0.81	0.70	11.78

G-matrix consistency is a little violated.

Quantum numbers

	After	Before
J2 (total sys.)	3.73	4.20
J2 (N sys.)	3.69	3.96
L2 (N sys.)	1.98	3.15
S2 (N sys.)	0.75	0.81
L2 (kaon)	0.14	0.24
T2	2.00	2.00
Tz	1.00	1.00

Particle numbers

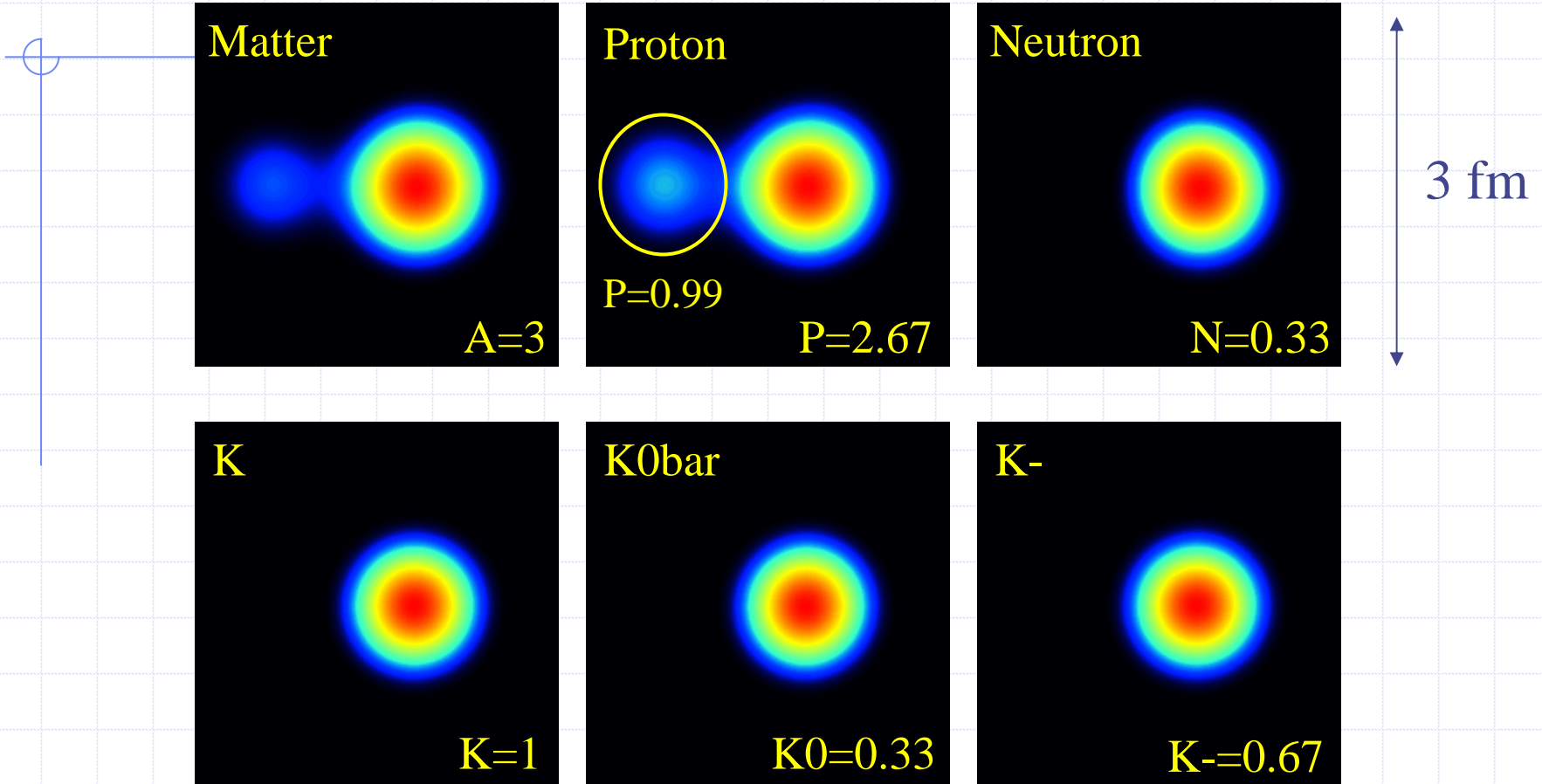
Proton	2.67
Neutron	0.33
K-	0.67
K0bar	0.33

Even if we start AMD cooling from various initial value γ which is concerned to the isospin, most of all solutions are converged to this result.

JT projection is correctly working.

pppK⁻

Density distributions of $|P_{Tz} \Phi\rangle$ in ZX plane.



$$|\frac{3}{K}\text{He}\rangle = |p\rangle \otimes |\frac{2}{K}\text{H}\rangle$$

$\rho(0) = 1.57\text{fm}^{-3}$... central density

$$\sqrt{\frac{2}{3}} |pp \otimes K^-\rangle - \sqrt{\frac{1}{3}} |pn \otimes \bar{K}^0\rangle$$

model space

2 Gaussian / nucleon
5 Gaussian / kaon

force

$E(K)=99.6\text{MeV}$, central density= 1.31fm^{-3}

Results

Projecting onto $J=1$ and $T=1/2$. Parity is negative.

	B.F	width	dens0	rmsR	beta	gamma
JT projection	105.01	25.85	1.29	0.97	0.54	3.80

G-matrix consistency is a little violated.

Quantum numbers

	After	Before
J2 (total sys.)	1.97	5.49
J2 (N sys.)	2.01	5.31
L2 (N sys.)	2.04	3.26
S2 (N sys.)	2.02	2.05
L2 (kaon)	0.05	0.18
T2	0.75	0.79
Tz	0.50	0.50

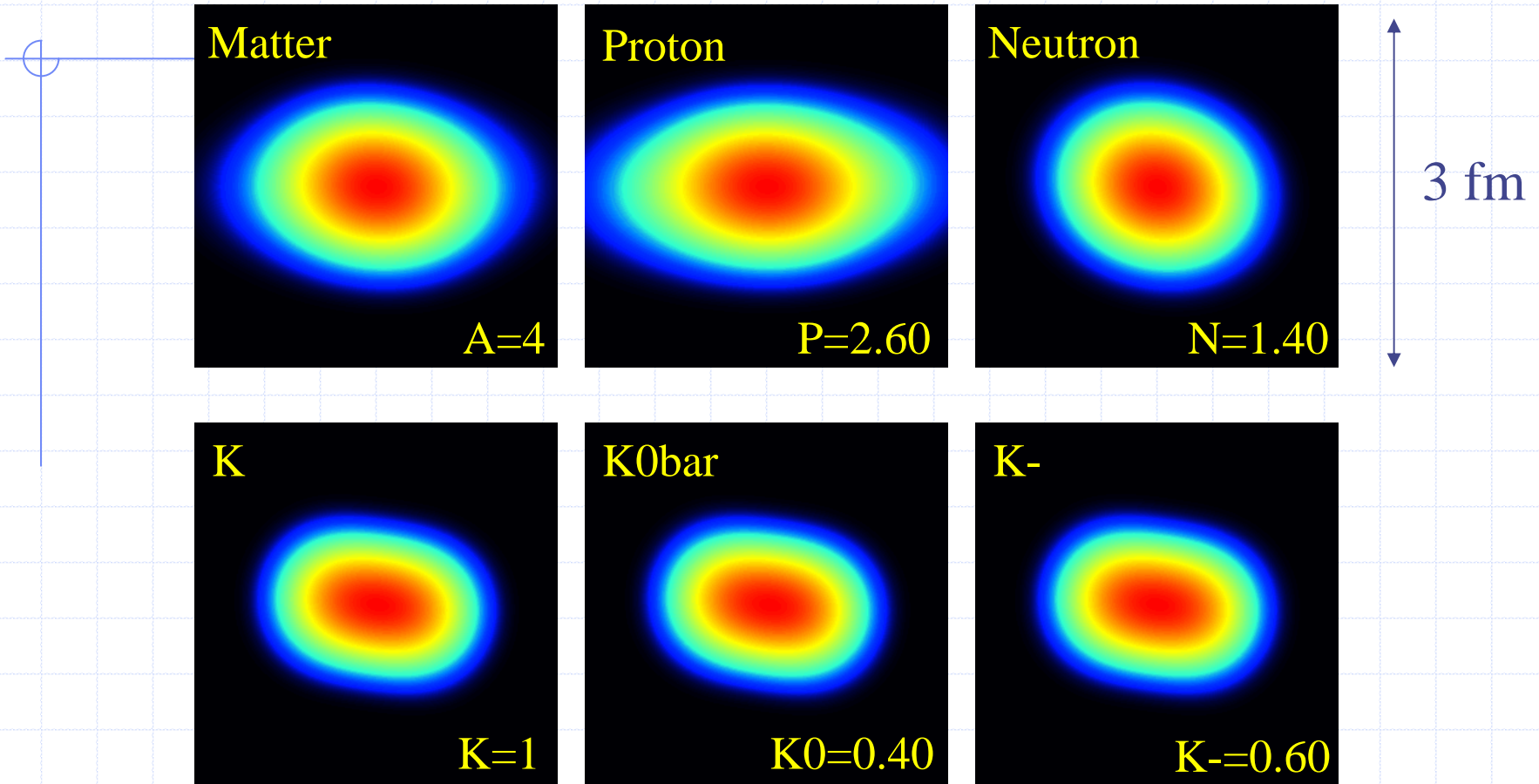
Particle numbers

Proton	2.60
Neutron	1.40
K-	0.60
K0bar	0.40

JT projection is correctly working.

pppnK-

Density distributions of $|P_{Tz}\Phi\rangle$ in ZX plane.



$\rho(\mathbf{0}) = 1.29\text{fm}^{-3}$... central density

Results — ${}^6\text{BeK}^-$ —

model space

2 Gaussian / nucleon
5 Gaussian / kaon

force

$E(K)=106\text{MeV}$, central density= 0.96fm^{-3}

Results

Projecting onto $J=0$ and $T=1/2$. Parity is negative.

	B.F	width	dens0	rmsR	beta	gamma
JT projection	104.24	33.40	0.91	1.17	0.44	0.30

G-matrix consistency is accomplished.

Quantum numbers

	After	Before
J2 (total sys.)	-0.06	4.86
J2 (N sys.)	0.03	4.60
L2 (N sys.)	0.03	4.58
S2 (N sys.)	0.00	0.02
L2 (kaon)	0.09	0.27
T2	0.75	0.97
Tz	0.50	0.50

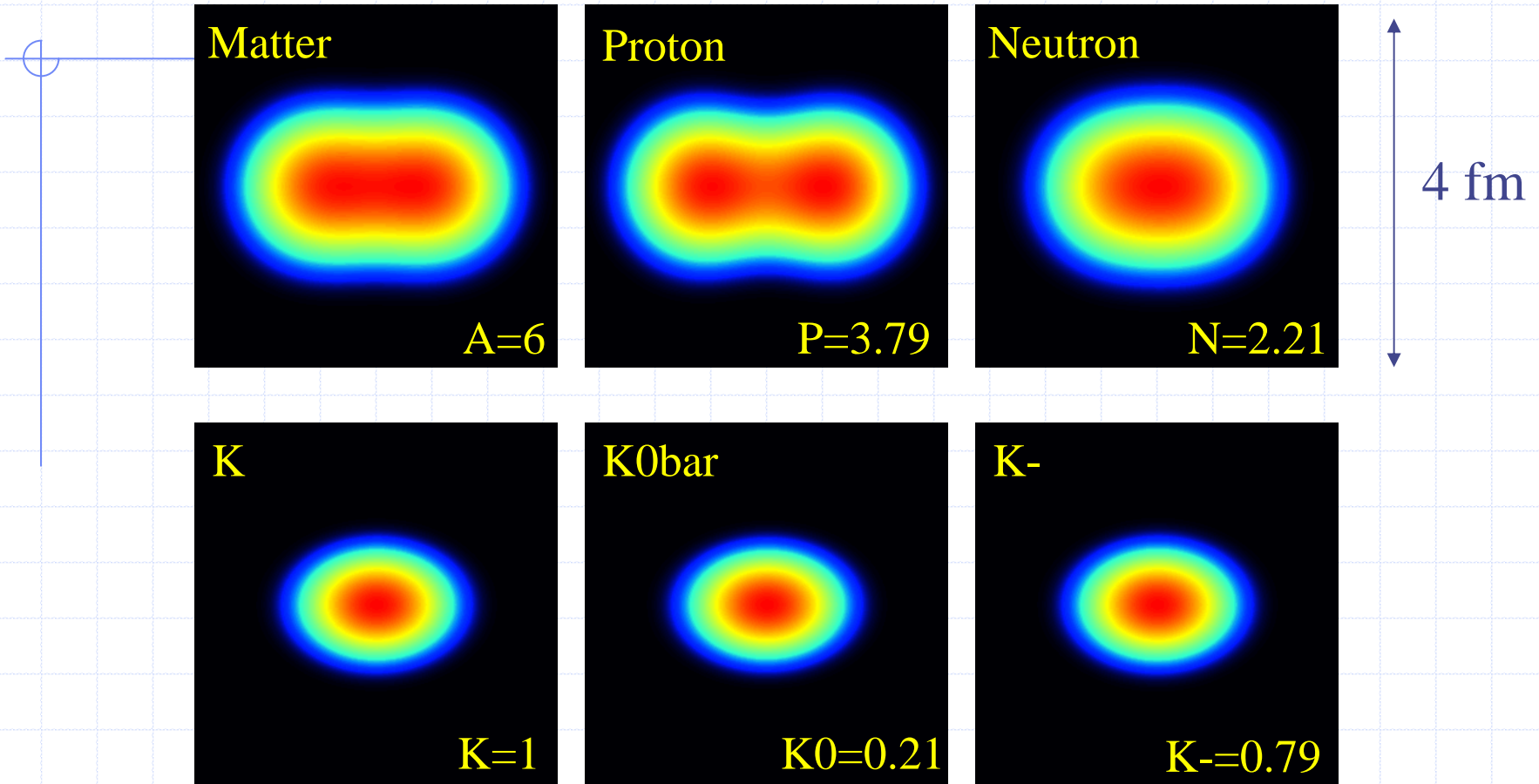
Particle numbers

Proton	3.79
Neutron	2.21
K-	0.79
K0bar	0.21

JT projection is correctly working.

${}^6\text{BeK}^-$

Density distributions of $|P_{Tz}\Phi\rangle$ in ZX plane.



$\rho(\mathbf{0}) = 0.91\text{fm}^{-3}$... central density

${}^9\text{BK}^-$

model space

2 Gaussian / nucleon
5 Gaussian / kaon

force

$E(K)=106\text{MeV}$, central density= 0.96fm^{-3}
for ${}^6\text{BeK}^-$

Results

Projecting onto $J=3/2$ and $T=0$. Parity is negative.

	$E(K)$	width	dens0	rmsR	beta	gamma
JT projection	99.6	35.5	0.67	1.46	0.46	16

G-matrix consistency is violated.

Quantum numbers

	After	Before
J2 (total sys.)	3.68	9.74
J2 (N sys.)	3.71	9.55
L2 (N sys.)	2.77	8.75
S2 (N sys.)	0.76	0.78
L2 (kaon)	0.08	0.18
T2	0.00	0.24
Tz	0.00	0.00

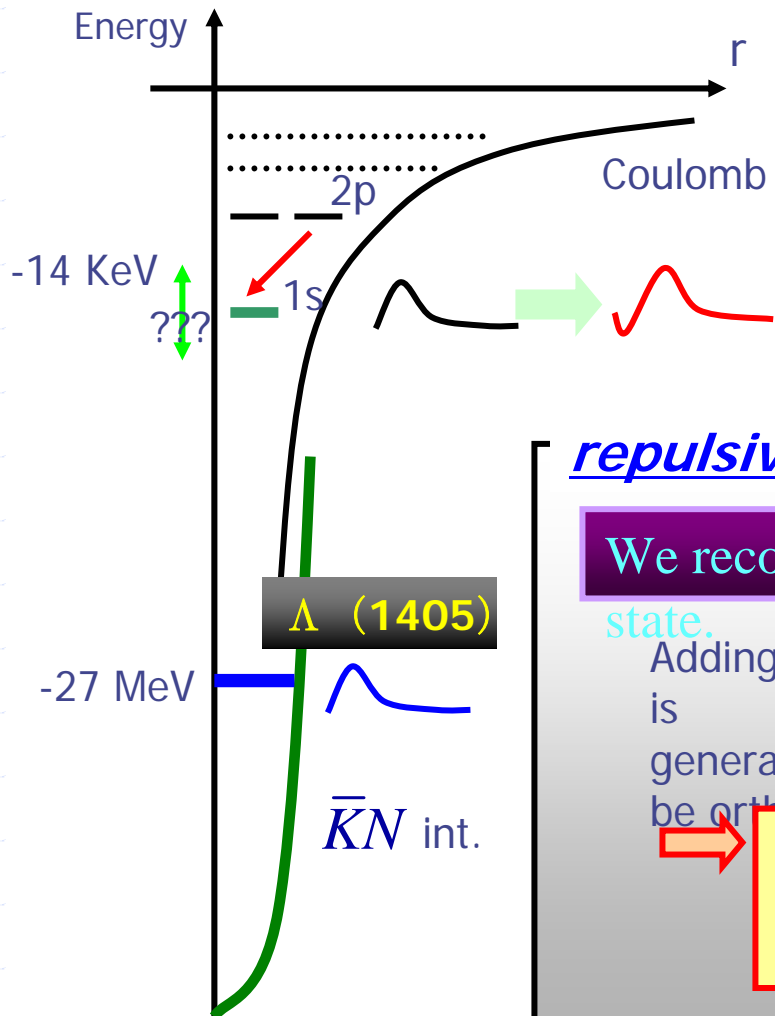
Particle numbers

Proton	4.69
Neutron	4.31
K-	0.69
K0bar	0.31

JT projection is almost correctly working.

KN interaction

Is $\bar{K}N$ interaction very strongly attractive ?



“kaonic hydrogen puzzle”

- atomic 1s shift caused by KN interaction
- atomic 2p \rightarrow 1s X-ray measurement : attractive
- analysis of low energy data : repulsive

Prof. Iwasaki investigated very precisely \rightarrow “repulsive” !!

*: M.Iwasaki et.al. Phys.Rev.Lett.78(1997)3067

repulsive \equiv attractive ???

We recognize $\Lambda(1405)$ is K^-p bound state.

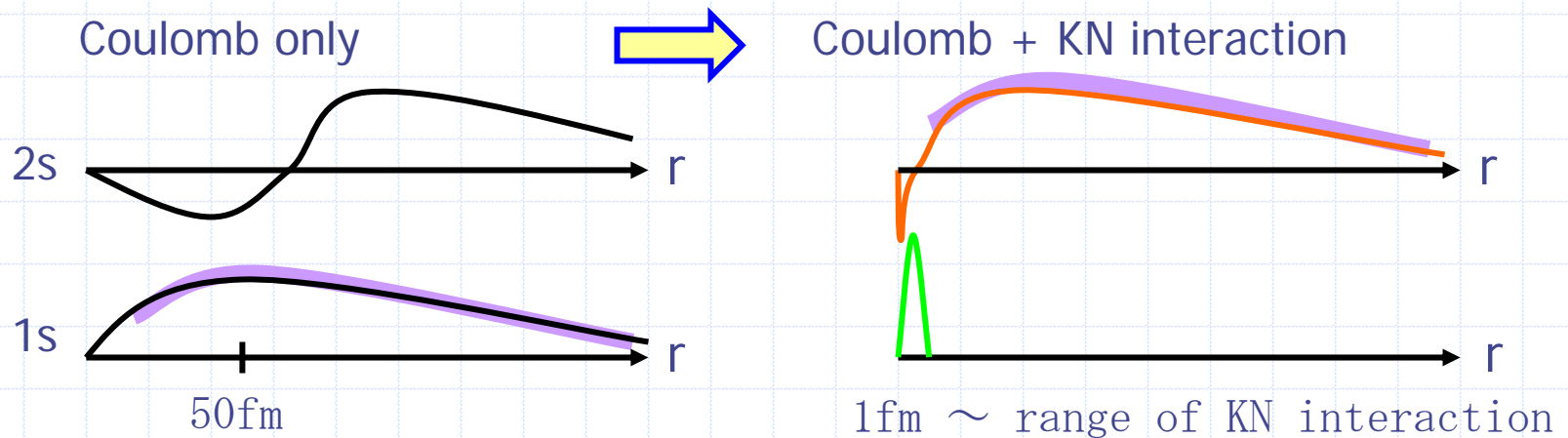
$\Lambda(1405) =$
3 quark state ??

Adding $\bar{K}N$ int., a new bound state (= $\Lambda(1405)$) is generated below the atomic 1s state. Then it must be orthogonal to the new bound state.

\rightarrow The original atomic 1s state has a node and its kinetic energy increases. As a result it shifts upward. = repulsive-like

kaonic hydrogen atom

about atomic 1s state shifted by KN interaction



2s state lowered by K^-N interaction is very similar to the original atomic 1s state. If ignoring a node, this lowered 2s state seems to be the solution obtained by changing the boundary condition a little --- $\psi = 0$ at $r=0$ fm \rightarrow $\psi = 0$ at $r=1$ fm --- So lowered 2s state appears energetically near the original 1s state.

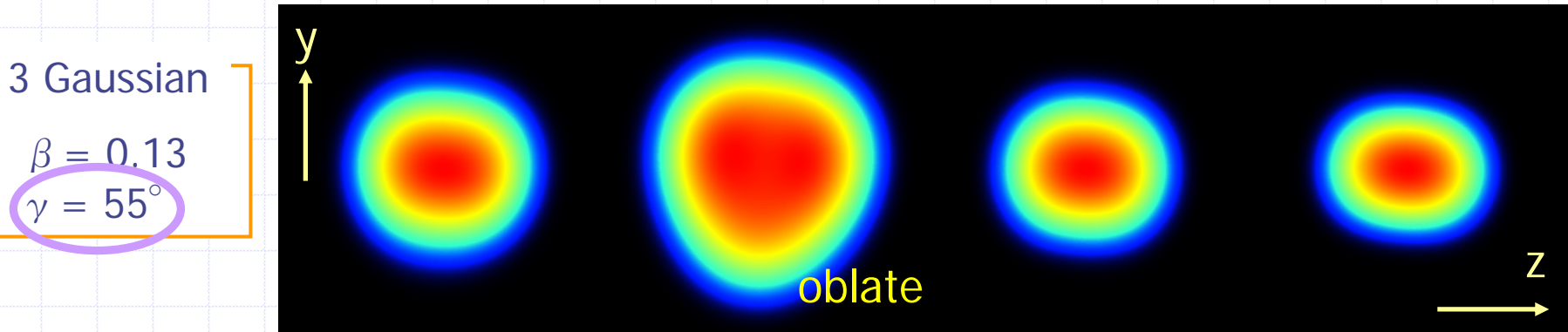
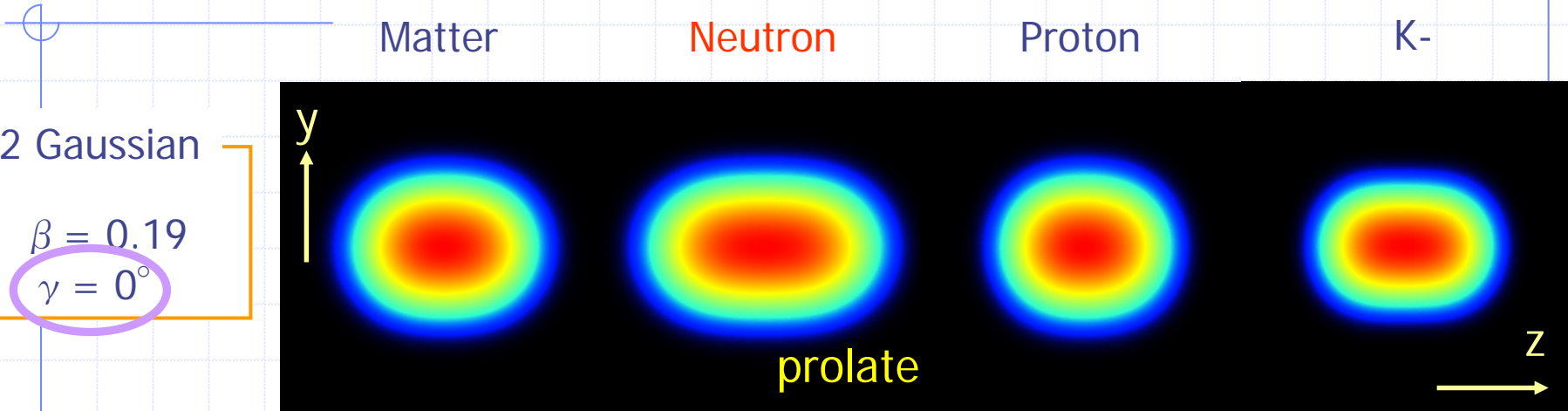
lowered 2s = 1s' (1s もどき !!)

Seeing with nuclear scale, it has a node. So it is 2s state.

Seeing with atomic scale, the shape of its wave function is almost that of original 1s state.

ppnK- // dependency on the number of wave packets //

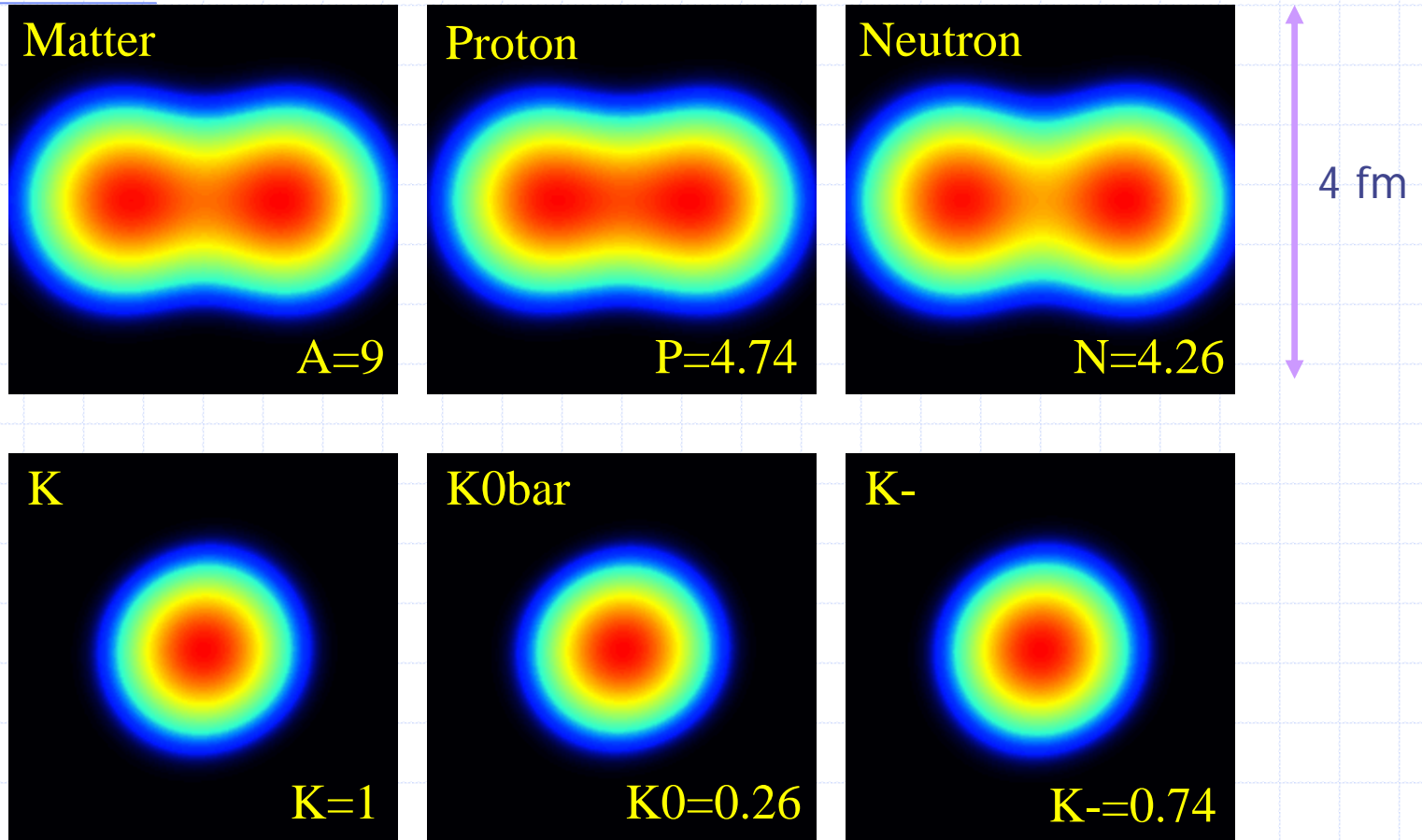
Density distribution in YZ plane



As expected, the neutron distribution is strongly dependent on the number of wave packets. As a result, the shape of the system is largely changed.

${}^9\text{B}K^-$ (tentative)

Density distributions of $|P_{Tz}\Phi\rangle$ in ZX plane.



Similar structure to ${}^8\text{Be}K^-$