

3. NUCLEAR INTERACTIONS

(1)

Generalities

- Relevant degrees of freedom: structureless nucleons
- Non-relativistic framework (+ relativistic corrections)
- Nucleon dynamics in nuclei is to a large extent non-relativistic

$$\frac{|\vec{p}|}{m} \approx \frac{200 \text{ MeV}}{939 \text{ MeV}} \Rightarrow \left(\frac{v}{c}\right)^2 < 0.1$$

- Interaction is instantaneous, potential is time independent
- Strong + weak + EM accounted for
- Here we consider the ab initio Hamiltonian (as opposed to effective)

AB INITIO

in vacuum
nucleus independent
links nuclei to QCD

$$H |\Psi^A\rangle = E^A |\Psi^A\rangle$$

EFFECTIVE

in medium
nucleus dependent
focuses on phenomenological description

$$H_{\text{eff}}^A |\Psi_{\text{eff}}^A\rangle = E^A |\Psi_{\text{eff}}^A\rangle$$

General form

$$H = \sum_{i=1}^A \frac{p_i^2}{2M} + \frac{1}{2} \sum_{i \neq j=1}^A V(i,j) + \frac{1}{3!} \sum_{i \neq j \neq k=1}^A W(i,j,k) + \dots$$

which is the form of these operators?

how many do we need to include?

- Over the years, three different approaches have been developed to model nuclear interactions

- i) Symmetry-driven
- ii) Physics-driven
- iii) Effective field theory

Symmetry-driven approach

• Most general form of two-nucleon interaction

$$V(1,2) = V(\vec{r}_1, \vec{p}_1, \vec{s}_1, \vec{t}_1, \vec{r}_2, \vec{p}_2, \vec{s}_2, \vec{t}_2)$$

idea: use known symmetries to constrain the operatorial form

• Nuclear interactions invariant under

- exchange of two nucleons
- translations in space
- translations in time
- rotations
- Galilean boost
- parity
- time-reversal
- isospin (approximate)

→ Potential can be written as a part that is invariant under isospin rotations

$$\hat{V}(p,p) = \tilde{V}(p,n) = \tilde{V}(n,n)$$

$$V = \tilde{V} + V^{ISB}$$

+ small isospin-symmetry-breaking corrections

$$V^{ISB}(p,p) \neq V^{ISB}(p,n) \neq V^{ISB}(n,n)$$

The most general operator structure consistent with all symmetry constraints can be then derived.

Two examples

i) isospin operators

Potential must be a scalar with respect to isospin rotations

⇒ can depend only on scalar product

$$\begin{aligned} \Rightarrow V(\dots, \vec{t}_1, \vec{t}_2) &= F(\dots, \vec{t}_1 \cdot \vec{t}_2) \\ &= F_0(\dots) + F_1(\dots) \vec{t}_1 \cdot \vec{t}_2 + F_2(\dots) (\vec{t}_1 \cdot \vec{t}_2)^2 + \dots \end{aligned}$$

One can prove that

$$(\vec{t}_1 \cdot \vec{t}_2)^2 |T M_T\rangle = (3 - 2\vec{t}_1 \cdot \vec{t}_2) |T M_T\rangle \quad \rightarrow \text{Exercise}$$

$$\Rightarrow F_0(\dots) + F_1(\dots) \vec{t}_1 \cdot \vec{t}_2$$

ii) linear dependence on spin operators

Since the potential is a scalar under spatial rotations, terms linear in \vec{S} must appear under the form of a scalar product

$$V(\dots, \vec{S}_1, \vec{S}_2) = F(\dots, \underline{\vec{a} \cdot \vec{S}_1 + \vec{b} \cdot \vec{S}_2})$$

invariance under permutations $1 \leftrightarrow 2$ implies $\vec{a} = \vec{b}$

$$\Rightarrow \vec{a} \cdot (\vec{S}_1 + \vec{S}_2) = \vec{a} \cdot \vec{S}$$

possibilities: $\vec{r}, \vec{p}, \vec{L}, \vec{S}$ would be trivial (a constant)
 wouldn't be invariant under parity / only non-trivial possibility

At the end, the most general operator structure reads

(4)

$$V(1,2) = V^0 + V^S \vec{s}_1 \cdot \vec{s}_2 + V^T \vec{t}_1 \cdot \vec{t}_2 + V^{St} (\vec{s}_1 \cdot \vec{s}_2) (\vec{t}_1 \cdot \vec{t}_2)$$

with

$$V^i = \sum_{k=1}^5 f_k^i (\vec{r}_1^2, \vec{p}_1^2, \vec{L}^2) O_k \quad (i=0, s, t, st)$$

and

$$O_k = \begin{cases} 1 & \\ \vec{L} \cdot \vec{S} & \text{spin-orbit} \\ S_{12}^r \equiv 3(\vec{s}_1 \cdot \vec{r})(\vec{s}_2 \cdot \vec{r}) - \vec{s}_1 \cdot \vec{s}_2 & \left. \begin{array}{l} \\ \\ \end{array} \right\} \text{tensor} \\ S_{12}^p \equiv 3(\vec{s}_1 \cdot \vec{p})(\vec{s}_2 \cdot \vec{p}) - \vec{s}_1 \cdot \vec{s}_2 & \\ Q_{12} \equiv \frac{1}{2} \left[(\vec{s}_1 \cdot \vec{L})(\vec{s}_2 \cdot \vec{L}) + (\vec{s}_2 \cdot \vec{L})(\vec{s}_1 \cdot \vec{L}) \right] & \text{quadratic spin-orbit} \end{cases}$$

where $\vec{x} \equiv \frac{\vec{x}}{|\vec{x}|}$

- Tensors are defined with an additional $\vec{s}_1 \cdot \vec{s}_2$ such that the average over the angles of \vec{r} and \vec{p} gives zero

- Q_{12} is called quadratic spin-orbit because it can be reexpressed as a function of $(\vec{L} \cdot \vec{S})^2$, $\vec{L} \cdot \vec{S}$ and \vec{L}^2

→ This approach was followed to develop the Argonne-V18 potential model (AV18)

→ Three-nucleon forces? One could follow the same strategy, but it gets very complicated

Physics-driven approach

(5)

First models date back to the 1930's

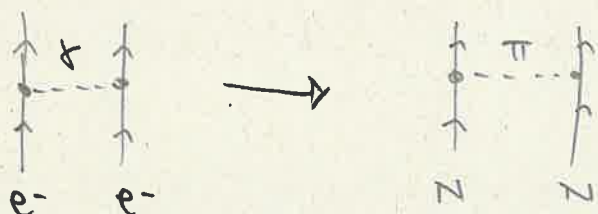
What was known at the time?

- Coulomb interaction between charged particles (infinite range)
- Nuclear interaction is short-ranged ($\sim 2 \text{ fm}$)

- All nuclei are compact
- Nuclear force negligible for atoms
- First scattering experiments

Yukawa 1935

Idea: nuclear force mediated by massive spin-0 boson



"Mesotron"

(later \rightarrow pion)

range $\sim 2 \text{ fm} \Rightarrow m \sim 100 \text{ MeV}$

\sim Compton wavelength of exchanged boson

Yukawa potential

$$V(r) \propto \frac{e^{-mr}}{r}$$

- One-pion exchange describes long-range attraction
- Generates tensor and $\vec{t} \cdot \vec{t}$ structures
- However, shorter-range part unaccounted for
 - \rightarrow Multi-pion exchange? Doesn't work.
 - \rightarrow 1960s: discovery of heavier mesons

\downarrow
Can model ranges smaller than $\frac{1}{m_\pi}$

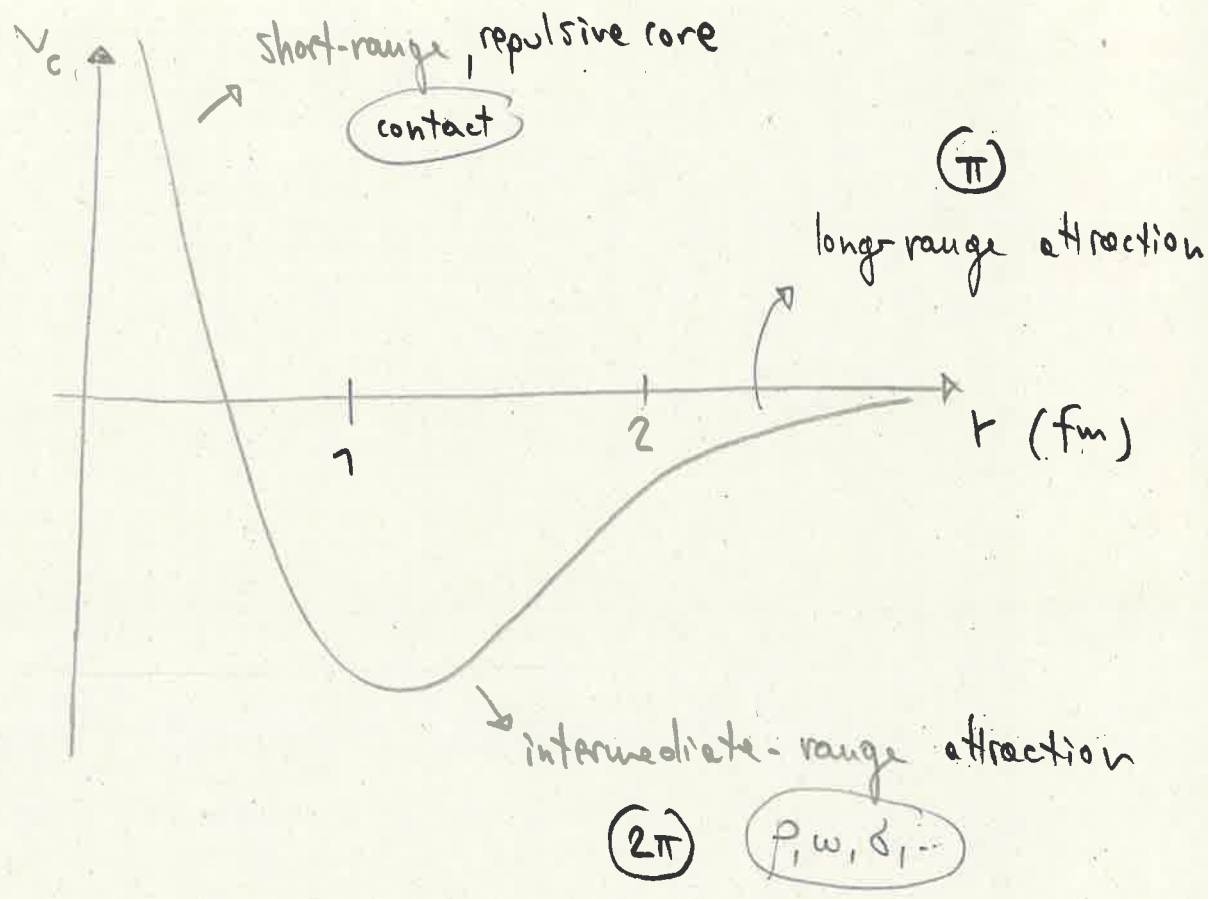
- ρ, ω, σ mesons generate different operator structures. (6)

- Very precise NN potentials constructed in this way

1980's $\rightarrow \chi^2 \sim 2$

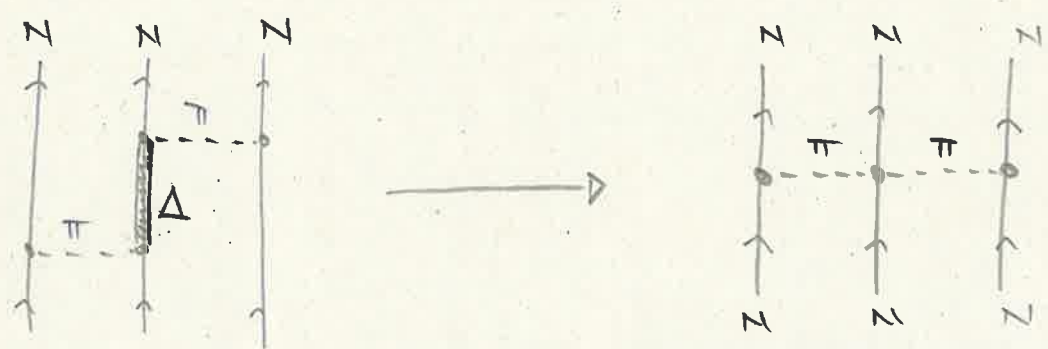
1990's $\rightarrow \chi^2 \sim 1$

One-boson exchange (OBE) potentials



- Three-nucleon forces?

Certain processes e.g. involving nucleonic excitations can not be described in terms of NN interactions



Fujita-Miyazawa 1957

Effective field theory approach (EFT)

(7)

- Symmetry-driven and physics-driven potentials very successful, but phenomenological models

no systematic way to improve

On the contrary, effective (field) theories provide a more systematic framework

- EFT introduced in nuclear physics in the 1990's (Weinberg, van Kolck, ...) → today: most used approach

- EFT principles

1) Separation of scales

→ Use separation of energy scales to define d.o.f. and expansion parameter ϵ

$$\epsilon = \frac{Q}{M}$$

typical scales at play in the considered physical system

high-energy scale

2) Write all possible interaction terms (Lagrangian) between d.o.f. allowed by symmetries of underlying theory

3) Order by size all these terms

→ Scaling $\epsilon \sim$ determine \sim ⇒ "Power counting"

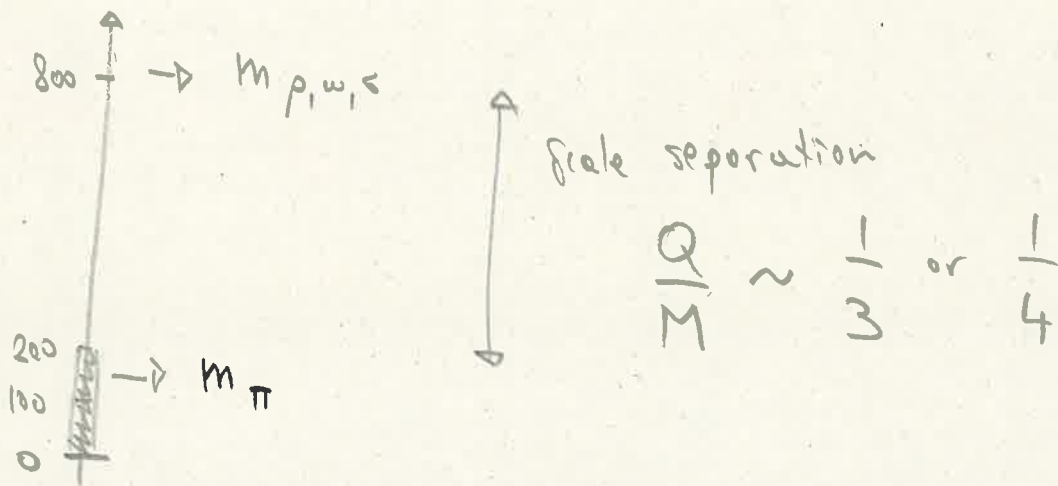
4) Truncate the expansion at a given order and compute observables

(8)

- Unknown parameters (coupling constants) have to be fixed (underlying theory or experiment)
- Order-by-order convergence has to be verified

- Application to nuclear forces

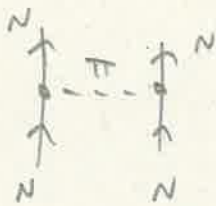
1) → Typical nucleon momenta $Q \sim 100-200 \text{ MeV}$



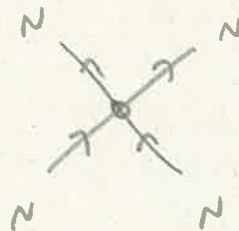
⇒ Degrees of freedom explicit in theory: N, π

2) → Interactions of two types

pion exchanges



contact interactions



3) → Most used ordering is Weinberg power counting

General equation for k-nucleon diagram

(9)

$$V = 2K - 4 + 2L + \sum_i \Delta_i$$


vertices


where

$$\begin{cases} L = \text{number of loops} \\ \Delta_i = d_i + \frac{h_i}{2} - 2 \end{cases}$$

derivatives nucleon fields

4) → At lowest order (= "leading order" (LO)) two terms

pure contact  $V_{ct}^{(0)}(p, p') = C_S + C_T \vec{S}_1 \cdot \vec{S}_2$ — central

one-pion exchange  $V_{\pi}^{(0)}(p, p') = -\frac{g_A^2}{4f_\pi^2} \vec{t}_1 \cdot \vec{t}_2 \frac{\vec{S}_1 \cdot \vec{q} \vec{S}_2 \cdot \vec{q}}{q^2 + m_\pi^2}$ — tensor

↓
Higher orders and many-nucleon sectors (→ see slides)

Experimental constraints: NN scattering

Reaction types:

- np scattering is the easiest: neutron beams on hydrogen targets

• pp scattering: technically easy, but EM interaction needs to be subtracted (beyond Coulomb, radiative corrections, ...)

• nn scattering: only indirect information (no n targets)

- nd scattering (then subtract np component)
- two neutrons in final state, e.g. $n+d \rightarrow n+n+p$
- comparison between different reactions



For energies beyond the pion production threshold ($\sim 2 m_\pi$ in lab frame) an inelastic component appears in the potential.

Cross section and partial-wave expansion

Relevant quantities

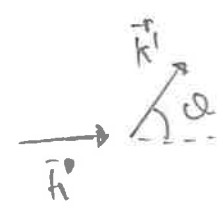
- differential cross section

$\frac{d\sigma}{d\Omega}(E, \theta)$

usually independent of φ

- total cross section

$\sigma(E)$



\rightarrow elastic scattering $\rightarrow |\vec{k}| = |\vec{k}'|$ and $E = \frac{k^2}{2\mu}$ with $\mu = \frac{M}{2}$ reduced mass

Wave function at $r \rightarrow \infty$

$\Psi_{\vec{k}}(\vec{r}) \xrightarrow{r \rightarrow \infty} e^{i\vec{k}\cdot\vec{r}} + f(k, \theta) \frac{e^{ikr}}{r}$

/
/
 incoming outgoing

$f(k, \theta)$ is the scattering amplitude $\Rightarrow \frac{d\sigma(k, \theta)}{d\Omega} = |f(k, \theta)|^2$ (11)

By expanding $\Psi_E(\vec{r})$ in partial waves one obtains

$$\Psi_k(\vec{r}) \xrightarrow{r \rightarrow \infty} \sum_{l=0}^{\infty} (2l+1) P_l(\cos\theta) \frac{(-1)^{l+1} e^{-ikr} + S_l(k) e^{ikr}}{2ikr}$$

incoming
outgoing

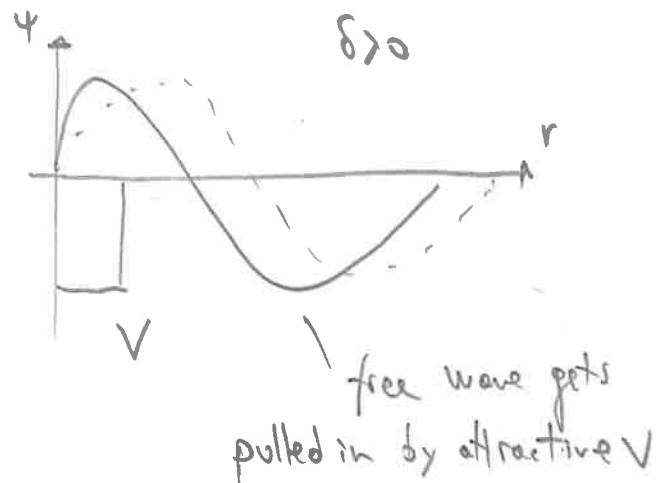
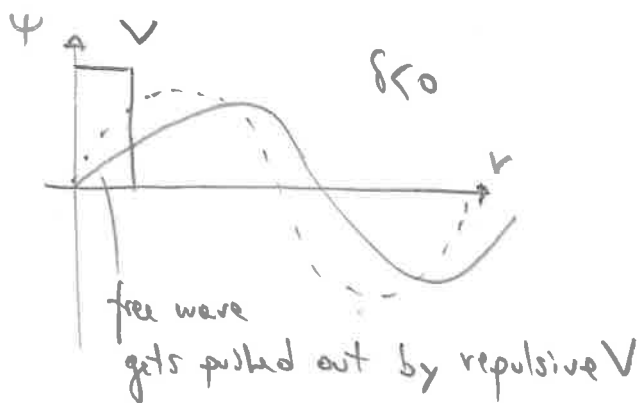
$S_l(k) = 1 + 2ik f_l(k)$ is the scattering matrix, usually parametrised as

$$S_l(k) = e^{2i\delta_l(k)} \quad (\text{S-matrix is unitary})$$

Then

$$f_l(k) = \frac{S_l(k) - 1}{2ik} = \frac{e^{i\delta_l(k)} \sin \delta_l(k)}{k} = \frac{1}{k \cot \delta_l(k) - ik}$$

Outside the range of the interaction ($r \rightarrow \infty$) the potential leads to a change in the phase of the outgoing wave



$$\Rightarrow \sigma_{\text{tot}}(E) = \int d\Omega \frac{d\sigma(k, \theta)}{d\Omega} = \int d\Omega |f(k, \theta)|^2 = \frac{4\pi}{k^2} \sum_{l=0}^{\infty} (2l+1) \sin^2 \delta_l(k)$$

Low energy scattering

Schwinger showed that $k^{2l+1} \cot \delta_l(k)$ has an analytic expansion

$$l=0 \rightarrow k \cot \delta_0(k) = -\frac{1}{a} + \frac{1}{2} r_e k^2 - P^2 r_e^3 k^4 + \mathcal{O}(k^6)$$

$$l=1 \rightarrow k^3 \cot \delta_1(k) = -\frac{3}{a_p^3} + \mathcal{O}(k^2)$$

Effective range expansion

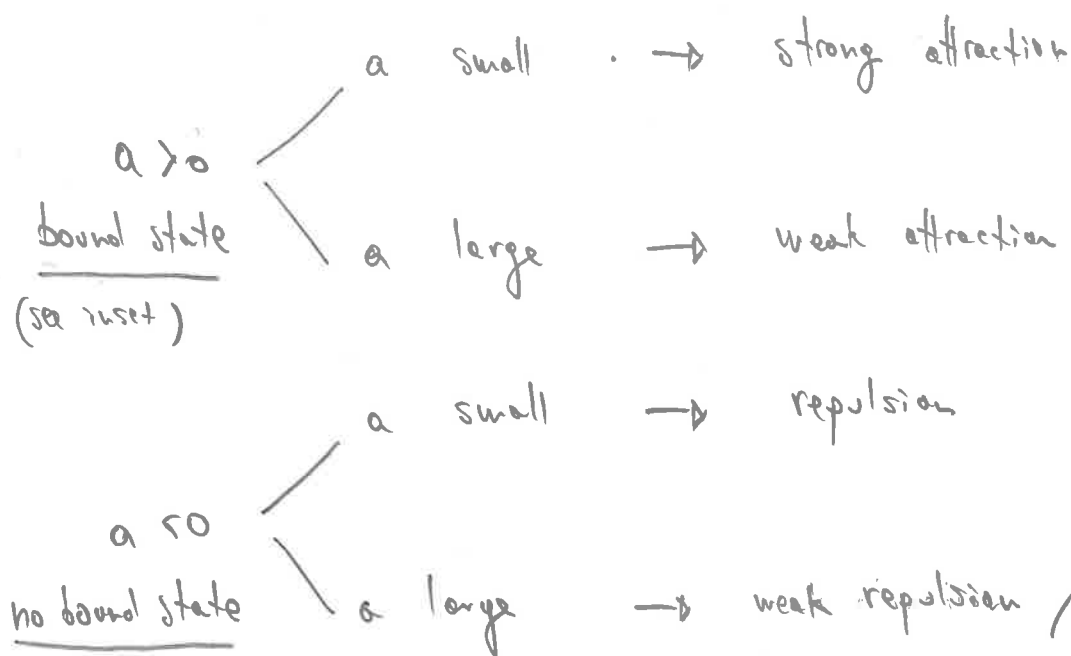
In the limit $k \rightarrow 0$ only $l=0$ scattering is present and

$$f_0(k) \rightarrow -a \Rightarrow \frac{d\sigma}{d\Omega} = a^2 \Rightarrow \sigma_{\text{tot}} = 4\pi a^2$$

in this limit a is the radius of an absorption disc equivalent to $\sigma_{\text{tot}}(0)$

\rightarrow effective range represents first correction \sim range of the interaction

Scattering length \rightarrow useful to expand $\cot \delta_0(k) \Rightarrow$ get $\delta_0(k) = -ka$



$$f_0(k) \xrightarrow{k \rightarrow 0} \frac{1}{-\frac{1}{a_0} - ik}$$

pole at $k = \frac{i}{a}$

$$\Rightarrow e^{ikr} \rightarrow e^{-\frac{r}{a}}$$

Case of nucleon-nucleon scattering

(13)

→ At low energies, only $L=0 \Rightarrow$ two possible channels:

spin singlet 1S_0 (s) and spin triplet 3S_1 (t)

$a_s(np) \approx -23.7 \text{ fm}$	$r_s(np) \approx 2.73 \text{ fm}$
$a_t(np) \approx +5.4 \text{ fm}$	$r_t(np) \approx 1.75 \text{ fm}$
$a_s(pp) \approx -17.1 \text{ fm}$ (after correcting for Coulomb)	$r_s(pp) \approx 2.84 \text{ fm}$
$a_s(nn) \approx -16.6 \text{ fm}$	$r_s(nn) \approx 2.66 \text{ fm}$

- Charge symmetry realised to a very good extent
- Charge independence more approximate
- Bound state in 3S_1 channel (n-p) \Rightarrow deuteron
- Scattering length large also in 1S_0 channel
 \Rightarrow almost-bound state (resonance) of two neutrons
(two protons excluded because of Coulomb) di-neutron

Phase shifts and high-energy scattering

14

Phase shifts are used to parametrise scattering data based on a partial wave analysis

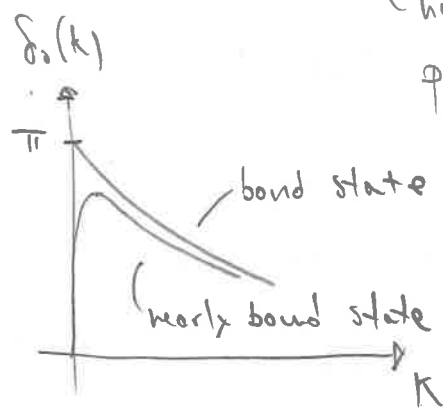
→ see e.g. Nijmegen partial wave analysis nn-online.org/NN

Discussion of $T=1$ and $T=0$ phase shifts (see figures)

- Presence of a bound state in 3S_1 channel (consistent with scattering length) and a nearly-bound state in 1S_0

Levinson's theorem $\delta_l(k=0) = n_l \pi$

↳ number of bound states in partial wave l



- S-waves become repulsive at high momentum \leftrightarrow short distances \Rightarrow short-range repulsion of NN interaction

• Consider 3P waves

1) Central component \sim average of 3P_0 , 3P_1 , and 3P_2 is small at small energies \Rightarrow there must be something else \Rightarrow spin-orbit

11) Consider spin-orbit operator

$$\vec{L} \cdot \vec{S} = \frac{1}{2} (\vec{J}^2 - \vec{L}^2 - \vec{S}^2)$$

$$= \frac{1}{2} [J(J+1) - L(L+1) - S(S+1)]$$

for 3P $\frac{1}{2} J(J+1) - 2$

- -2 for $^3P_0 \Rightarrow$ repulsive
- -1 for $^3P_1 \Rightarrow$ repulsive
- $+1$ for $^3P_2 \Rightarrow$ attractive

?? \Rightarrow there must be something else
 $\longleftrightarrow \delta < 0$
 $\longleftrightarrow \delta > 0$
 ↓
 tensor term

$V_{LS} \vec{L} \cdot \vec{S}$ where $V_{LS} < 0$

Experimental constraints: the deuteron

The deuteron is the only bound system of two nucleons.

Main properties

- Binding energy $\bar{E} = 2.2245 \text{ MeV}$
- Total angular momentum $J = 1$
- Magnetic dipole moment $\mu = 0.8574 \mu_N$ with $\mu_N = \frac{e}{2m_p}$
- Electric quadrupole moment $Q = 0.2859 \text{ efm}^2$
- R.m.s. radius $r = 1.97 \text{ fm}$

Proton-neutron system can exist in $T=0$ or $T=1$. Given $J=1$, there are four possibilities $^3S_1, ^1P_1, ^3P_1, ^3D_1$
 $+ \quad - \quad - \quad + \quad \leftarrow \text{parity} = (-1)^L$

Since parity has to be conserved, it can be at most a combination of the two states of the same parity.

Quadrupole moment = 0 for a spherical nucleus

⇒ cannot be pure L=0 state

If it were a superposition of the P states, it would mix T=0 and T=1. Not impossible, but not very probable given that we expect isospin symmetry (specially since there is no Coulomb).

Further hints come from the magnetic moment.

$$\mu = \langle \Psi_{JM} | M_z | \Psi_{JM} \rangle$$

for a nucleus in the state $|\Psi_{JM}\rangle$.

M_z is the third component of the magnetic moment operator

$$\vec{M} = \mu_N \left[\sum_{i=1}^Z (\vec{l}_i + g_p \vec{s}_i) + \sum_{i=1}^N g_n \vec{s}_i \right]$$

proton orbital angular momenta

nucleon spin g-factors

$$g_p = 5.585 \quad g_n = -3.826$$

For the deuteron one has

$$\mu = \mu_N \langle \Psi_{11} | (\vec{l}_p + g_p \vec{s}_p + g_n \vec{s}_n) | \Psi_{11} \rangle$$

and one finds

• for a superposition $|\Psi_{11}\rangle = \cos \omega_0 |^3S_1\rangle + \sin \omega_0 |^3D_1\rangle$

$$\mu = \mu_N (0.820 - 0.570 \sin^2 \omega_0)$$

• for a superposition $|\Psi_{11}\rangle = \cos \omega_1 |^1P_1\rangle + \sin \omega_1 |^3P_1\rangle$

$$\mu = \mu_N (0.500 + 0.190 \sin^2 \omega_1)$$

⇒ it has to be $^3S_1 + ^3D_1$ (cf. experimental value $\mu = 0.8574 \mu_N$)

The only term mixing different L components is the tensor
in the nuclear interaction

only operator that does not
commute with L^2

(17)