

Chapter 3

Nuclear Force and Two-Nucleon Systems

3-1 The Deuteron

Binding Energy ; - The deuteron is a very unique nucleus in many respects.

- 1- Only loosely bound nucleus
- 2- having binding energy much less than the average value between a pair of nucleons in all the other stable nuclei.

$$E_B = \left\{ \left[\sum M_p + \sum M_n \right] - M_{\text{nucleus}} \right\} c^2 \quad (M_p \approx M_n)$$

$$E_B = \left\{ [938.783 + 939.566] - 1876.124 \right\} = 2.225 \text{ MeV}$$

$$\bar{E}_B = 2.22457312(22) \text{ MeV} \quad \text{more precise}$$

This value is obtained, using the radiative capture of a neutron by hydrogen.



$$\text{If } E_n \approx 0 \rightarrow E_\gamma = E_B$$

Partly because of small E_B , the deuteron has no excited state.

Ground state Property	Value
Binding energy E_B	2.22457312(22) MeV
Spin & Parity J^π	1^+
Isospin T	0
Magnetic dipole moment μ_d	0.857406(1) μ_N
Electric quadrupole Q_d	0.28590(30) $e \text{ fm}^2$
Radius r_d	1.963(4) fm

Parity, Spin and Isospin;

From experimental information $\rightarrow \Pi_d = +$

$$\Psi_d = \Psi_n \Psi_p \Psi_{\text{rel.}}$$

↑ \nearrow
 intrinsic \searrow
 (r-dep.) (l-dep)

$\Psi_n \Psi_p = +$ Since Ψ_n and Ψ_p are two different states of a single particle nucleon, their intrinsic wave func. have the same parity.

$\rightarrow \Psi_{\text{rel.}}$ determines the parity of deuteron

$$Y_{LM}(0,\varphi) \xrightarrow{\Pi} Y_{LM}(\pi-\theta, \pi+\varphi) = (-)^L Y_{LM}(0,\varphi)$$

$\left\{ \begin{array}{l} \bar{r} = \bar{r}_1 - \bar{r}_2 \xrightarrow{\Pi_{12}} -\bar{r} \\ (r, \theta, \varphi) \xrightarrow{\Pi_{12}} (r, \pi-\theta, \varphi+\pi) \end{array} \right.$

Since $\Pi_d = + \rightarrow L = \text{even}$

The ground state of deuteron has $J=1$

$$\bar{J} = \bar{L} + \bar{S} \quad S = 0, 1$$

$$\left\{ \begin{array}{l} J=1 \\ L=\text{even} \end{array} \right. \rightarrow \cancel{S=0} \quad \cancel{L=2}$$

$$\rightarrow (L, S) = (0, 1) \text{ and } (2, 1) \quad \text{for ground state}$$

$\overset{\circ}{P}$	$\overset{\dagger}{L}$
dominant part	small but significant amount admixture

$$\Psi_d = \Psi_r X_s \xi_t$$

\downarrow \downarrow
 sym sym

$\rightarrow \xi_t$ should be antisym.

$$\rightarrow T = 0 \quad \cancel{T \neq 0}$$

$$X_{ns}^S = \begin{cases} X_1' = \alpha_1 \alpha_2 \\ X_0' = \frac{1}{\sqrt{2}} \{ \alpha_1 \beta_2 + \alpha_2 \beta_1 \} \\ X_{-1}' = \beta_1 \beta_2 \end{cases} \quad \begin{aligned} \alpha &= |\frac{1}{2}, \frac{1}{2}\rangle_S \\ \beta &= |\frac{1}{2}, -\frac{1}{2}\rangle_S \end{aligned}$$

$$\xi_{t_2}^+ = \xi_0^+ = \frac{1}{\sqrt{2}} \{ \alpha'_1 \beta'_2 - \alpha'_2 \beta'_1 \} \quad \begin{aligned} \alpha' &= |\frac{1}{2}, \frac{1}{2}\rangle_T \\ \beta' &= |\frac{1}{2}, -\frac{1}{2}\rangle_T \end{aligned}$$

Alternatively, if

$$\xi_{t_2}^1 = \begin{cases} \xi_1^+ & (\text{P-P system}) \\ \xi_0^+ & (\text{n-p system}) \\ \xi_{-1}^+ & (\text{n-n system}) \end{cases} \quad (\text{if deuteron } \in \xi_0^+)$$

→ We would expect the other two states. However, no such states have been observed.

- i) The existence of P-P system may be rejected since the Coulomb repulsion is of order 1 Mev (large fraction of E_B)
 - ii) This limitation, however, does not apply for n-n system.
- Conclusion $T=1$ bound state: impossible

Therefore, possible states are:

$T=0, S=1, L=0 : {}^3S_1$ (triplet S-state)

$T=0, S=1, L=2 : {}^3D_1$ (, D-")

If L and S are good quantum number;

$$[H, L^2] = [H, S^2] = 0$$

→ deuteron would have to be in either one of these two states.

However, this is not the case.

3-2) Deuteron Magnetic Dipole Moment:

Mag. dipole operator:

Mag. dipole moment has two sources:

- { 1 - μ_n, μ_p coming from the intrinsic spin and orbital motion of quarks.
- { 2 - μ_p^{orb} coming from orbital motion of proton.

From electromagnetic theory,

$$\mu_i^{\text{orb.}} = \frac{e\hbar[c]}{2M_p c} l_i \quad \text{in CGS } [c] \text{ is ignored.}$$

$$\mu_i^{\text{orb.}} = g_e^{(i)} l_i \quad g_e^{(i)} = \begin{cases} 1 & \text{for } p \\ 0 & \text{for } n \end{cases}$$

$$\mu_i^{\text{spin}} = g_s^{(i)} S_i$$

$$\text{Since } S = \frac{1}{2} \rightarrow g_s^{(i)} = \begin{cases} g_p = 2 \mu_p = 5.585695 \mu_N & \text{for } p \\ g_n = 2 \mu_n = -3.826085 \mu_N & \text{for } n \end{cases}$$

for free nucleons

If we assume Nucleon structure $\underset{\text{bound}}{=}$ nucleon structure $\underset{\text{free}}$

We may use then values in nuclei as well.

$$\mu_j = g_p S_p + g_n S_n + \ell_p$$

Since $m_p \approx m_n \rightarrow \ell_p = \frac{1}{2} L$ (rel. angular mom.)
or angular mom. of deuteron

$$\mu_d = g_p S_p + g_n S_n + \frac{1}{2} L$$

Contribution from 3S_1 -state:

For this state $L=0$, we will see soon that;

$$\mu_j({}^3S_1) = \mu_p + \mu_n = 0.879805 \mu_N$$

There is a small difference with the observed value;

$$\mu_d(\text{obs.}) - \mu_d({}^3S_1) = 0.857406 - 0.879805 = -0.022399 \mu_N$$

This difference may be attributed to three possibilities:

1) Nucleon structure $\underset{\text{bound}}{\neq}$ nucleon structure $\underset{\text{free}}$

$$\rightarrow g_{p \text{ bound}} \neq g_{p \text{ free}} \quad g_{n \text{ bound}} \neq g_{n \text{ free}}$$

2) There are contributions from charged mesons exchanged between the proton and the neutron, and these have not been included in our calculations.

3) There are small admixtures of 3D_1 -state in the ground state of deuteron.

i) Item 1 is extremely unlikely, since the deuteron is a loosely bound system:

Because $E_B = 2.22 \text{ MeV} \ll E_{B(\text{quarks})} \sim \text{hundreds of MeV}$

→ $E_B = 2.22 \text{ MeV}$ is not expected to affect the motion of quarks.

ii) The effect of mesonic currents is possible.

iii) Item 3 is more likely to be major case for this discrepancy

Expectation value of the magnetic dipole operator:

By convention, the mag. moment, similar to other static electromagnetic moments is defined as the expectation value of the z -component, or $q=0$ component in spherical tensor notation of the operator in the salustate of Max. M, $M=J$.

$$\mu_d = \langle J, M=J | \mu_z | J, M=J \rangle$$

Now first we show;

$$\langle \alpha', j_m' | V_q | \alpha, j_m \rangle = \frac{\langle \alpha', j_m | J \cdot V | \alpha, j_m \rangle}{j(j+1)} \langle j_m' | J_q | j_m \rangle$$

$$J \cdot V = \sum_q (-1)^q J_{1q} V_{1,-q}$$

$$\begin{aligned} \langle \alpha', j_m | J \cdot V | \alpha, j_m \rangle &= \langle \alpha', j_m | (J_0 V_0 - J_1 V_{-1} - J_{-1} V_{+1}) | \alpha, j_m \rangle \\ &= m \langle \alpha', j_m | V_0 | \alpha, j_m \rangle + \frac{1}{\sqrt{2}} \sqrt{(j+m)(j-m+1)} \langle \alpha', j_{m+1} | V_{-1} | \alpha, j_m \rangle \\ &\quad - \frac{1}{\sqrt{2}} \sqrt{(j-m)(j+m+1)} \langle \alpha', j_{m+1} | V_{+1} | \alpha, j_m \rangle = C_{jm} \langle \alpha' j | V | \alpha j \rangle \end{aligned} \quad (1)$$

Remember $J_{\pm} = \mp \frac{1}{\sqrt{2}} (J_x \pm i J_y) = \mp \frac{1}{\sqrt{2}} J_{\pm}'$, $J_z = J_z'$

C_{jm} does not depend on α', α and V .

Also since $J \cdot V$ is a scalar op $\rightarrow C_{jm}$ is independent of m

$$\rightarrow C_{jm} \rightarrow C_j$$

If $V \rightarrow J$

$$\rightarrow \langle \alpha, j_m | J^2 | \alpha, j_m \rangle = C_j \langle \alpha j | J | \alpha j \rangle \quad (2)$$

By Wigner-Eckart theorem;

$$\langle \alpha', j_m' | T_k | \alpha, j_m \rangle = (J_K, m_q | j_K j_m') \langle \alpha' j | T_k | \alpha j \rangle$$

$$\frac{\langle \alpha', j_{m'} | V_q | \alpha, j_m \rangle}{\langle \alpha, j_{m'} | J_q | \alpha, j_m \rangle} = \frac{\langle \alpha' j | V | \alpha j \rangle}{\langle \alpha j | J | \alpha j \rangle}$$

$$(1) ad(\gamma) \rightarrow \frac{\langle \alpha' j | V | \alpha j \rangle}{\langle \alpha j | J | \alpha j \rangle} = \frac{\langle \alpha', j_m | J \cdot V | \alpha, j_m \rangle}{\langle \alpha, j_m | J^2 | \alpha, j_m \rangle}$$

$$\text{Since } \langle \alpha, j_m | J^2 | \alpha, j_m \rangle = j(j+1)$$

$$\rightarrow \langle \alpha', j_{m'} | V_q | \alpha, j_m \rangle = \frac{\langle \alpha', j_m | J \cdot V | \alpha, j_m \rangle}{j(j+1)} \langle j_{m'} | J_q | j_m \rangle$$

Now

$$\langle j_m | \mu_z | j_m \rangle = \frac{\langle j_m | \mu \cdot J | j_m \rangle}{j(j+1)} \langle j_m | J_z | j_m \rangle$$

$$= \frac{m}{j(j+1)} \langle j_m | \mu \cdot J | j_m \rangle$$

\downarrow
projection of μ on J

$$\mu_j = g_p s_p + g_n s_n + \frac{1}{2} L = \frac{1}{2} \left\{ (g_p + g_n) S + (g_p - g_n)(s_p - s_n) + L \right\}$$

But

$$\langle S m_s | (S_p - S_n) | S' m'_s \rangle = \begin{cases} 0 & s = s' \\ \pm 0 & s = 1, s' = 0 \\ \pm 0 & s = 0, s' = 1 \end{cases}$$

But we are interested in $s = s' = 1$ case. Therefore its contribution in our case is Zero.

This reduces M_d to a func. of L and S only (in our case).

$$\langle Jm | \mu_z | Jm \rangle = \frac{M}{J(J+1)} \langle Jm | \frac{1}{2} \{ (g_p + g_n)(S \cdot J) + (L \cdot J)^2 \} | Jm \rangle$$

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

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

$$M_d = \langle J_{m=j} | \mu_z | J_{m=j} \rangle = \frac{1}{4(J+1)} \left\{ (g_p + g_n) (J(J+1) - L(L+1) - S(S+1)) \right. \\ \left. + (J(J+1) + L(L+1) - S(S+1)) \right\}$$

For 3S_1 -state, $L=0$, $J=1$, $S=1$

$$\mu_d({}^3S_1) = \mu_p + \mu_n$$

For 3D_1 -state $L=2$, $J=1$, $S=1$

$$\mu_d({}^3D_1) = \frac{1}{8} \{ (g_p + g_n) (-2) + \zeta \} = 0.310 \mu_N$$

Since $\mu_d({}^3D_1) < \mu_d({}^3S_1)$

\rightarrow any admixture of 3D_1 -state will reduce the value of μ_d from that given by $\mu_d({}^3S_1)$.

$$Y_{LSJ}^M = \sum_{m_s m_e} (l S m_e m_s / JM) X_s^{m_s} Y_e^{m_e}$$

With the notation: ${}^b X_a \quad \begin{cases} X: l \\ a: J \\ b: 2S+1 \end{cases}$

$$^3S_1: \quad J=1 \quad M=1 \quad l=2 \quad S=1$$

$$Y_{011}^1 = (0101/11) Y_0^0 \alpha(1) \alpha(2) = Y_0^0 \alpha(1) \alpha(2)$$

$$^3D_1: \quad J=1 \quad M=1 \quad l=2 \quad S=1$$

$$Y_{211}^1 = \sum_{m_s} (21(1-m_s)m_s / 11) Y_2^{1-m_s} X_1^{m_s}$$

$$= (212-1/11) Y_2^2 X_1^{-1} + (2110/11) Y_2^1 X_1^0 + (2101/11) Y_2^0 X_1^1$$

$$= \sqrt{\frac{3}{5}} Y_2^2 \beta(1) \beta(2) - \sqrt{\frac{3}{10}} Y_2^1 \frac{1}{\sqrt{2}} [\alpha(1) \beta(2) + \beta(1) \alpha(2)] + \frac{1}{\sqrt{10}} Y_2^0 \alpha(1) \alpha(2)$$

when $\begin{cases} X_1^1 = \alpha(1) \alpha(2) \\ X_1^0 = \frac{1}{\sqrt{2}} [\alpha(1) \beta(2) + \beta(1) \alpha(2)] \\ X_1^{-1} = \beta(1) \beta(2) \end{cases}$

$$^3S_1: \quad \underline{\Phi}_{011}^1(r) = \frac{U_0(r)}{r} Y_{011}^1 = \frac{U_0(r)}{r} Y_0^0 X_1^1$$

$$^3D_1: \quad \underline{\Phi}_{211}^1(r) = \frac{U_2(r)}{r} Y_{211}^1$$

Admixture of 3D_1 -state:

$$|\Psi_1\rangle = a |{}^3S_1\rangle + b |{}^3D_1\rangle$$

experimental
↓

$$\left\{ \begin{array}{l} \langle \Psi_1 | \mu_z | \Psi_1 \rangle = a^2 \mu_s ({}^3S_1) + b^2 \mu_d ({}^3D_1) = 0.857 \mu_N \\ \quad (\text{no off-diagonal matrix element for } \mu_z) \\ \langle \Psi_1 | \Psi_1 \rangle = a^2 + b^2 = 1 \end{array} \right.$$

$$\rightarrow b^2 = 0.04 \rightarrow 4\% \text{ admixture of Hg } {}^3D_1$$

3-3 Deuteron Electric Quadrupole Moment:

Multipole Expansion of potential:

$$\Phi(x) = \sum_{l=0}^{\infty} \sum_{m=-l}^l \frac{4n}{2l+1} q_{lm} \frac{Y_{lm}(\theta, \phi)}{r^{l+1}}$$

$$\Phi(x) = \int \frac{f(x')}{|x-x'|} d^3x'$$

For the potential outside the charge distribution; $r_< = r'$, $r_> = r$.

$$\Phi(x) = 4n \sum_{lm} \frac{1}{2l+1} \left[\int Y_{lm}^*(\theta, \phi') r'^l f(x') d^3x' \right] \frac{Y_{lm}(\theta, \phi)}{r^{l+1}}$$

$$\rightarrow q_{lm} = \int Y_{lm}^*(\theta, \phi') r'^l f(x') d^3x'$$

$$q_{l,-m} = (-1)^m q_{lm}^*$$

$$q_{10} = \frac{1}{\sqrt{6\pi}} \int g(x') d^3x' = \frac{1}{\sqrt{6\pi}} q$$

$$q_{11} = -\sqrt{\frac{3}{8\pi}} \int (x' - iy') g(x') d^3x' = -\sqrt{\frac{3}{8\pi}} (P_x - iP_y)$$

$$q_{10} = \sqrt{\frac{3}{4\pi}} \int z' g(x') d^3x' = \sqrt{\frac{3}{4\pi}} P_z$$

$$q_{12} = \frac{1}{4} \sqrt{\frac{15}{2\pi}} \int (x' - iy')^2 g(x') d^3x' = \frac{1}{12} \sqrt{\frac{15}{2\pi}} (Q_{11} - 2iQ_{12} - Q_{22})$$

$$q_{21} = -\sqrt{\frac{15}{8\pi}} \int z' (x' - iy') g(x') d^3x' = -\frac{1}{3} \sqrt{\frac{15}{8\pi}} (Q_{13} - iQ_{23})$$

$$q_{20} = \frac{1}{2} \sqrt{\frac{5}{4\pi}} \int (3z'^2 - r'^2) g(x') d^3x' = \frac{1}{2} \sqrt{\frac{5}{4\pi}} Q_{33}$$

$$P = \int x' g(x') d^3x'$$

$$Q_{ij} = \int (3x_i x_j - r'^2 \delta_{ij}) g(x') d^3x'$$

$$\Phi(r) = \frac{q}{r} + \frac{P_r}{r^3} + \frac{1}{2} \sum_{ij} Q_{ij} \frac{x_i x_j}{r^5} + \dots$$

In a nucleus it is not possible to observe the charge distribution density directly, and the measured multipole moments may be used to infer something about the source.

Since $\underset{l,m}{Y(\theta, \phi)} \xrightarrow{D} (-1)^l Y_{lm}(\theta, \phi)$

All odd multipole electric operators give vanishing expectation values (for pure ground states).

→ The lowest non-vanishing multipole moment → quadrupole moment

$$Q_{pd} = 0.28590 e \text{ fm}^2 \text{ (measured)}$$

Quadrupole operator:

For a spherical nucleus;

$$\langle x^2 \rangle = \langle y^2 \rangle = \langle z^2 \rangle$$

As a result;

$$\begin{aligned}\langle r^2 \rangle &= \langle x^2 + y^2 + z^2 \rangle = \langle x^2 \rangle + \langle y^2 \rangle + \langle z^2 \rangle \\ &= 3 \langle z^2 \rangle\end{aligned}$$

The charge quadrupole op.
which measures the lowest
order departure from a
spherical symmetry is defined by;

$$Q_z = e(3z^2 - r^2)$$

For spherical nucleus (charge distribution);

$$\langle Q_z \rangle = 0$$

For deuteron: $\langle Q_z \rangle = +0.29$

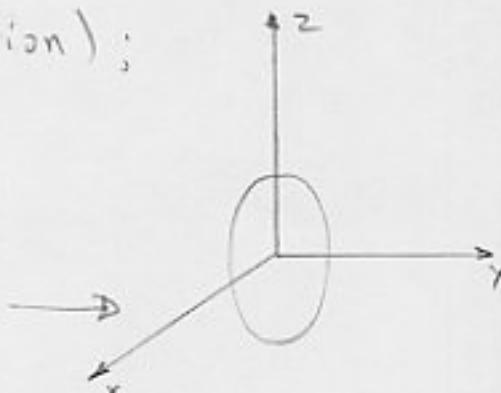
$$Q_z = e(3z^2 - r^2) = er^2(3G'_{20} - 1)$$

$$= \sqrt{\frac{16\pi}{5}} er^2 Y_{20}(\theta, \varphi)$$

spherical tensor
of rank 2

$$Q_A = \langle j | Q_z | j \rangle$$

Electric quadrupole moment
of a nucl. state



Remark :

$$\langle \alpha', j'm' | T_q^k | \alpha, jm \rangle = 0 \quad \text{unless } m' = q+m$$

$$\text{also } |j-k| \leq j' \leq j+k \quad (\text{triangle cond.})$$

Here $k=2$

$$\rightarrow \text{For } j < 1 \rightarrow Q_A = 0$$

At the same time since Q_2 operates on only in the coordinate space; it is independent of total spin S' .

$$\xrightarrow{\text{i.e.}} L \geq 1 \quad (\text{must be})$$

$$\rightarrow \langle {}^3S_1 | Q_2 | {}^3S_1 \rangle = 0$$

The nonvanishing $Q_d = 0.28590 \rightarrow$ Presence of 3D -component in the ground state

Expectation value of the Quadrupole operator:

Assumption: Deuteron ground state is a definite L.

$$|LS; JM\rangle = \sum_{M_L M_S} (LM_L, SM_S / JM) |LM_L\rangle |SM_S\rangle$$

$$Q_d(L) = \langle LS; JM | Q_2 | LS; JM \rangle$$

$$= \sum_{M_L M_S} \sum_{M'_L M'_S} (LM_L, SM_S / JM) (LM'_L, SM'_S / JM) \langle LM_L, SM_S | Q_2 | LM'_L, SM'_S \rangle$$

$$\text{where } |LM_L, SM_S\rangle \equiv |LM_L\rangle |SM_S\rangle$$

$$\text{Since } \langle SM_S | SM'_S \rangle = \delta_{M_S M'_S}$$

$$Q_d(L) = \sum_{M_L} (LM_L S(M-M_L) / JM)^2 \langle LM_L | Q_2 | LM_L \rangle$$

($M_L = M'_L$, because $Q_2 = 0$ for $M_L \neq M'_L$)

$$|LM_L\rangle = R_L(r) Y_{LM_L}(\theta, \varphi)$$

$$\int_0^\infty R_L^*(r) R_L(r) r^2 dr = 1$$

$$\begin{aligned} \langle LM_L | e(3z^2 - r^2) | LM_L \rangle &= \langle LM_L | \sqrt{\frac{16\pi}{5}} er^2 Y_{20}(\theta, \varphi) | LM_L \rangle \\ &= e \sqrt{\frac{16\pi}{5}} \int_0^\infty R_L^*(r) r^2 R_L(r) r^2 dr \int_{-1}^1 \int_0^1 Y_{LM_L}^*(\theta, \varphi) Y_{20}(\theta, \varphi) Y_{LM}(\theta, \varphi) d(C, \theta) d\varphi \end{aligned}$$

Since

$$\langle \alpha JM | T_q^k | \alpha' J'M' \rangle = (-)^{J-M} \begin{pmatrix} J & K & J' \\ -M & q & M' \end{pmatrix} \langle \alpha J || T^k || \alpha' J' \rangle$$

where

$$\langle \alpha J || T^k || \alpha' J' \rangle = (2k+1) \sum_{MM'} (-)^{J-M} \begin{pmatrix} J & K & J' \\ -M & q & M' \end{pmatrix} \langle \alpha JM | T_q^k | \alpha' J'M' \rangle$$

Here $T_q^k = Y_0^2$

$$\langle LM_L | Y_0^2 | LM_L \rangle = (-)^{L-M} \begin{pmatrix} L & 2 & L \\ -M_L & 0 & M_L \end{pmatrix} \langle L || Y^2 || L \rangle$$

But $\langle L || Y^k || L' \rangle = (-)^L \sqrt{\frac{(2\ell+1)(2\ell+1)(2\ell'+1)}{4\pi}} \begin{pmatrix} L & K & L' \\ 0 & 0 & 0 \end{pmatrix}$

$$\langle L || Y^2 || L \rangle = (-)^L (2L+1) \sqrt{\frac{5}{4\pi}} \begin{pmatrix} L & 2 & L \\ 0 & 0 & 0 \end{pmatrix}$$

$$\langle LM_L | Y_0^2 | LM_L \rangle = (-)^{M_L} \begin{pmatrix} L & 2 & L \\ -M_L & 0 & M_L \end{pmatrix} (2L+1) \sqrt{\frac{5}{4\pi}} \begin{pmatrix} L & 2 & L \\ 0 & 0 & 0 \end{pmatrix}$$

and where

$$\begin{pmatrix} J_1 & J_2 & J_3 \\ m_1 & m_2 & m_3 \end{pmatrix} = \frac{(-)^{J_1-J_2-m_3}}{\sqrt{2J_3+1}} (J_1 m_1, J_2 m_2 / J_1 J_2 J_3 - m_3)$$

$$\langle 2 2 1 | Y_0^2 | 2 2 \rangle = -\frac{1}{7} \sqrt{\frac{5}{\pi}}$$

$$\langle 2 1 1 | Y_0^2 | 2 1 \rangle = +\frac{1}{14} \sqrt{\frac{5}{\pi}}$$

$$\langle 2 0 1 | Y_0^2 | 2 0 \rangle = +\frac{1}{7} \sqrt{\frac{5}{\pi}}$$

The Clebsch-Gordon coeffs. for; S=1, L=2, M=S=1

$$(LM_L, S(M-M_L)/JM)^2 = (2 2 1 - 1 1 1)^2 = \frac{6}{10}$$

$$(2 1 1 0 1 1)^2 = \frac{3}{10}$$

$$(2 0 1 1 1 1)^2 = \frac{1}{10}$$

$$\begin{aligned}
 Q_d(L) &\equiv Q_d(2) \equiv Q_d(^3D_1) = \frac{6}{10} \langle L=2, M_L=2 | Q_2 | L=2, M_L=2 \rangle \\
 &+ \frac{3}{10} \langle L=2, M_L=1 | Q_2 | L=2, M_L=1 \rangle + \frac{1}{10} \langle L=2, M_L=0 | Q_2 | L=2, M_L=0 \rangle \\
 &= e\sqrt{\frac{16\pi}{5}} \left\{ \frac{6}{10} \left(-\frac{\sqrt{\frac{5}{\pi}}}{7} \right) + \frac{3}{10} \left(-\frac{\sqrt{\frac{5}{\pi}}}{14} \right) + \frac{1}{10} \left(\frac{\sqrt{\frac{5}{\pi}}}{7} \right) \right\} \langle r^2 \rangle_D
 \end{aligned}$$

where $\langle r^2 \rangle_D = \int_0^\infty R_D^*(r) r^2 R_D(r) r^2 dr$

An estimate:

If we take $\langle r^2 \rangle_D = \text{square of deuteron radius}$

$$\rightarrow Q_d(^3D_1) = -0.77 \text{ e fm}^2$$

but $Q_d = 0.28590 \text{ e fm}^2 \quad Q_d(^3D_1) \neq Q_d \text{ (experimentally)}$

\rightarrow The ground-state of deuteron is not entirely 3D_1 -state.

More realistic model:

$$|\Psi_d\rangle = a |{}^3S_1\rangle + b |{}^3D_1\rangle$$

$$\begin{aligned}
 Q_d &= a^2 \langle {}^3S_1, M=1 | Q_2 | {}^3S_1, M=1 \rangle + b^2 \langle {}^3D_1, M=1 | Q_2 | {}^3D_1, M=1 \rangle \\
 &+ 2ab \langle {}^3S_1, M=1 | Q_2 | {}^3D_1, M=1 \rangle
 \end{aligned}$$

with the notation:

$$|L S, JM\rangle \equiv |{}^3S_1, M=1\rangle$$

$\downarrow \downarrow \downarrow \downarrow$
 $0 1 1 1$

$$\langle {}^3S, M=1 | Q_2 | {}^3S, M=1 \rangle = 0$$

Since $L=0$

The main contribution comes from the last term (off-diagonal matrix)

$$\langle {}^3S, M=1 | Q_2 | {}^3D, M=1 \rangle \sim \int_0^\infty R_s^*(r) r^2 R_D(r) r^2 dr$$

$$|b|^2 = 4\% \text{ to } 7\%$$

(There are some parameters in the wave functions, R_s and R_D to be fixed)

3-4 Tensor Force and the Deuteron D-state

The analysis of $\mu_d \text{ad} Q_d \rightarrow |\Psi_d\rangle = a|{}^3S, \rangle + b|{}^3D, \rangle$

→ No definite L-value → L is not a good quantum number

$$\rightarrow [V \quad L^2] \neq 0$$

$$H = -\frac{\hbar^2}{2\mu} \nabla^2 + V$$

V : nucleon-nucleon int.
 μ : reduced mass

$$H|\Psi_d\rangle = E_d |\Psi_d\rangle$$

$|\Psi_d\rangle$: Ground-state wave func.
 eigenfunc. of H

From symmetry arguments;

the only contributions come from $L=0, L=2$

$$|\Psi_d\rangle = a|{}^3S, \rangle + b|{}^3D, \rangle$$

It may be convenient for us to think in terms of a matrix approach to the eigenvalue problem.

Since we are only interested in finding the amount of mixing between 3S_1 and 3D_1 states, we may use these two states as the basis to construct the Hamiltonian matrix;

$$\{H\} = \begin{pmatrix} H_{11} & H_{12} \\ H_{21} & H_{22} \end{pmatrix}$$

$$H_{11} = \langle {}^3S_1 | H | {}^3S_1 \rangle \quad H_{22} = \langle {}^3D_1 | H | {}^3D_1 \rangle \quad H_{12} = H_{21} = \langle {}^3D_1 | H | {}^3S_1 \rangle$$

In diagonalization, we find $\begin{cases} E_d \\ a, b \end{cases}$

But this is not of interest here;

We are interested to the type of H which can produce a mixing between 3S_1 and 3D_1 states.

If $H_{12} = H_{21} = 0 \rightarrow H$: diagonal in the basis states

The two eigenstates $|{}^3S_1\rangle$ and $|{}^3D_1\rangle$ without any admixtures between them.

The fact that ground state of deuteron is a mixture of these two states $\rightarrow \langle {}^3D_1 | H | {}^3S_1 \rangle \neq 0$

$$H = -\frac{\hbar^2}{2m} \nabla^2 + V$$

The kinetic energy term of H contributes only to the diagonal elements $\rightarrow \langle {}^3D, 1V | {}^1S, 1 \rangle \neq 0$

\rightarrow nuclear potential V is not diagonal in the basis span by states with definite orbital angular momentum and can therefore mix 3S , and 3P -states.

$$\langle {}^3D, 1V | {}^1S, 1 \rangle \neq 0$$

$$l=2 \quad \downarrow \quad l=0 \quad \uparrow \\ \text{must have } l=2$$

V must have a spatial part that is spherical tensor of rank 2, so as to be able to connect an S-state to a D-state.

$$\Psi_J = \Psi_r \Psi_{\text{spin}} \Psi_{\text{isospin}} \quad V = (r) (\text{spin}) (\text{isospin})$$

We ignore isospin part in the following discussion.

Since $[H, J^2] = 0$ (H conserves total angular momentum)

V must be a scalar in J

But we found that the spherical rank of spatial part of V is 2.

→ Intrinsic spin part must have rank 2, so that
a scalar product of these two rank-2 operators
can be constructed.

For spin $\frac{1}{2}$ system an arbitrary op may be expressed
as a linear combination of:

$$\alpha_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \quad \alpha_y = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix} \quad \alpha_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$$

$$I = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$$

In terms of spherical components;

$$\bar{\alpha} = \begin{cases} \alpha_{+1} = -\frac{1}{\sqrt{2}}(\alpha_x + i\alpha_y) = \sqrt{2} \begin{pmatrix} 0 & -1 \\ 0 & 0 \end{pmatrix} \\ \alpha_{-1} = +\frac{1}{\sqrt{2}}(\alpha_x - i\alpha_y) = \sqrt{2} \begin{pmatrix} 0 & 0 \\ 1 & 0 \end{pmatrix} \\ \alpha_0 = \alpha_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} \end{cases}$$

Its rank is $k=1$ (vector) and carries one unit of
angular momentum.

In general the product of two spherical tensors is not a spherical tensor.

$$[V^{k_1} \times U^{k_2}]_q^k \equiv T_q^k = \sum_{q_1, q_2} (k_1 q_1; k_2 q_2 / k q) V_{q_1}^{k_1} U_{q_2}^{k_2}$$

The product is a mixture of tensors with ranks;

$$|k_1 - k_2| \leq K \leq k_1 + k_2$$

If $V_{q_1}^{k_1}$ and $U_{q_2}^{k_2}$ are irreducible spherical tensors of rank k_1 and k_2 respectively, then T_q^k a spherical (irreducible) tensor of rank K .

For $K=2$:

$$T_0^0 = \frac{-V_0 U_0}{3} = \frac{1}{3} (V_+ U_{-1} + V_{-1} U_+ - V_0 U_0) \quad \left. \begin{array}{l} \\ \\ \end{array} \right\} \text{Irreducible}$$

$$T_1^1 = \frac{(V \times U)_q}{i\sqrt{2}} \quad \left. \begin{array}{l} \xrightarrow{\sqrt{3}} ? \\ \\ \end{array} \right\} "$$

$$T_{\pm 2}^2 = V_{\pm 1} U_{\pm 1}$$

$$T_{\pm 1}^2 = \frac{V_{\pm 1} U_0 + V_0 U_{\pm 1}}{\sqrt{2}} \quad \left. \begin{array}{l} \\ \\ \end{array} \right\} "$$

$$T_0^2 = \frac{V_+ U_- + 2 V_0 U_0 + V_- U_+}{\sqrt{6}}$$

$$\text{where } V_{\pm 1} = \mp \frac{1}{\sqrt{2}} (V_x \pm i V_y)$$

$$V_0 = V_z$$

Remark:

Cartesian tensor of rank 2 (dyadic):

$$T_{ij} = A_i B_j$$

$$A_i B_j = \underbrace{\frac{1}{3} (A \cdot B) \delta_{ij}}_{P_{ij}} + \underbrace{\frac{1}{2} (A_i B_j - A_j B_i)}_{Q_{ij}} + \underbrace{\left[\frac{1}{2} (A_i B_j + A_j B_i) - \frac{1}{3} (A \cdot B) \delta_{ij} \right]}_{R_{ij}}$$

$$AB = \begin{pmatrix} P_{11} & 0 & 0 \\ 0 & P_{22} & 0 \\ 0 & 0 & P_{33} \end{pmatrix} + \begin{pmatrix} 0 & Q_{12} & Q_{13} \\ -Q_{12} & 0 & Q_{23} \\ -Q_{13} & -Q_{23} & 0 \end{pmatrix} + \begin{pmatrix} R_{11} & R_{12} & R_{13} \\ R_{12} & R_{22} & R_{23} \\ R_{13} & R_{23} & R_{33} \end{pmatrix}$$

Diagonal

Anti-sym.

Sym.

$$T'_{ij} = \sum_{kl} R_{ik} R_{jl} T_{kl}$$

Under transformation:

Diagonal $\xrightarrow{\text{tr.}} \text{Diagonal}$

Anti-sym $\xrightarrow{\text{tr.}} \text{Anti-sym}$

Sym. $\xrightarrow{\text{tr.}} \text{Sym.}$

Diagonal matrix:

$\text{Tr}(P) = \text{const.}$

Any tr. leave it invariant. (scalar)

Anti-sym. matrix:

It has 3-independent elements - Q_{12}, Q_{13}, Q_{23} (vector)

Any tr. leave it antisymmetric.

Sym. matrix:

It has 5-independent elements:

$R_{11}, R_{22}, R_{33}, R_{12}, R_{13}, R_{23}$

with the cond. $\text{Tr}(R) = \text{const.}$

Any tr. leave it symmetric.

A two-body op in the nucleon intrinsic space may be constructed from the product of $\omega_{(1)}$ and $\omega_{(2)}$

$$[\omega_{(1)} \times \omega_{(2)}]_q^k$$

In Cartesian coord.:

$$\omega_{(1)} \cdot \omega_{(2)} = \omega_x(1) \omega_x(2) + \omega_y(1) \omega_y(2) + \omega_z(1) \omega_z(2)$$

In terms of α_{\pm} and α_0 :

$$\omega_{(1)} \cdot \omega_{(2)} = \omega_0(1) \omega_0(2) - \omega_{+1}(1) \omega_{-1}(2) - \omega_{-1}(1) \omega_{+1}(2)$$

which is in agreement with

$$J \cdot V = \sum_q (-1)^q J_{1q} V_{1-q}$$

$$T_0^\circ = [\omega_{(1)} \times \omega_{(2)}]_0^\circ = -\frac{1}{\sqrt{3}} (\omega_{(1)} \cdot \omega_{(2)})$$

$$T_0^\circ = \sum_{p,q} (1 p, 1 q / 100) \omega_p'(1) \omega_q'(2) \quad (p+q=0)$$

$$p, q = -1, 0, 1$$

Similarly

$$(\omega_{(1)} \times \omega_{(2)})_m^\dagger = \sum_{pq} (1 p, 1 q / 1 m) \omega_p'(1) \omega_q'(2) \quad (p+q=m)$$

Also;

$$[\alpha_{(1)} \times \alpha_{(2)}]_m^2 = \sum_{pq} (1P1g/2m) \alpha_p'(1) \alpha_q'(2)$$

$$[\alpha_{(1)} \times \alpha_{(2)}]_0^2 = \frac{1}{\sqrt{6}} \left\{ \alpha_1'(1) \alpha_1'(2) + \alpha_{-1}'(1) \alpha_1'(2) + 2 \alpha_0'(1) \alpha_0'(2) \right\}$$

$$[\alpha_{(1)} \times \alpha_{(2)}]_{\pm 1}^2 = \frac{1}{\sqrt{2}} \left\{ \alpha_{\pm 1}'(1) \alpha_0'(2) + \alpha_0'(1) \alpha_{\pm 1}'(2) \right\}$$

$$[\alpha_{(1)} \times \alpha_{(2)}]_{\pm 2}^2 = \alpha_{\pm 2}'(1) \alpha_{\pm 2}'(2)$$

Max. rank of $[\alpha_{(1)} \times \alpha_{(2)}] = 2$

To form a scalar product in J:

Rank of spin part $\stackrel{\text{must}}{=} \text{rank of orbital part}$

→ Max rank of orbital part = 2

This is adequate for our purposes;

We obtained before the orbital angular part of V must have rank 2 (P131)

Tensor operator:

An operator, formed by the scalar product of a second rank op. in intrinsic spin space and a similar one in coord. space, is called tensor operator.

The quantities we can use to form V being a scalar one:

$$\bar{r} = r\hat{n}, \quad S = \frac{1}{2}\alpha \quad \bar{\nabla} = \hat{i}\frac{\partial}{\partial x} + \hat{j}\frac{\partial}{\partial y} + \hat{k}\frac{\partial}{\partial z} \dots$$

$\bar{\nabla}$ is rejected, since it gives velocity-dep. forces (non-conservative)

i) $V = V(r) = V(\sqrt{r \cdot r})$ central

ii) $V = V(r)(\alpha_i \cdot \alpha_i)$ spin-dep. but central

iii) $V = V(r)(\alpha_i \cdot \hat{n})$

or $V = V(r)(\alpha_i \cdot \hat{n})$

or $V = V(r)(\alpha_i \times \alpha_j) \cdot \hat{n}$

are not scalars, they are pseudoscalars.

$$\text{Because } \hat{n} \xrightarrow{\pi} -\hat{n}$$

$$\alpha \xrightarrow{\pi} \alpha$$

There are two points:

- 1) Since \hat{n} is the only polar vector it must occur an even number of times in any acceptable product
- 2) Any polynomial in α can be reduced to an expression linear in α by means of spin identities.

For example $(\alpha \cdot \hat{n})^2 = I$, $(\alpha \cdot \hat{n})^3 = (\alpha \cdot \hat{n})$

→ An expression linear in α_1 and α_2 is the most general one we can use.

$$\text{iv) } V = V(r) (\alpha_1 \cdot \hat{n})(\alpha_2 \cdot \hat{n})$$

$$\text{or } V = V(r) (\alpha_1 \times \hat{n}) \cdot (\alpha_2 \times \hat{n})$$

The second one is a linear combination of the first and (α_1, α_2) by vector identities.

Hence, the combination $(\alpha_1 \cdot \hat{n})(\alpha_2 \cdot \hat{n})$ is essentially the only scalar leading to a non-central force.

It is useful to define the potential of the non-central force in such a way that its average over directions \hat{n} vanishes.

Since:

$$\overline{(\mathbf{A} \cdot \hat{\mathbf{n}})(\mathbf{B} \cdot \hat{\mathbf{n}})} = \frac{1}{3} (\mathbf{A} \cdot \mathbf{B})$$

all dir.

$$\hat{\mathbf{n}} = \hat{\mathbf{n}}(\theta, \varphi)$$

\downarrow
L-part

$$S_{12} = 3 (\alpha_1 \cdot \hat{\mathbf{n}}) (\alpha_2 \cdot \hat{\mathbf{n}}) - \alpha_1 \cdot \alpha_2$$

$$= \frac{3}{r^2} (\alpha_1 \cdot \vec{r}) (\alpha_2 \cdot \vec{r}) - \alpha_1 \cdot \alpha_2$$

$$S_1 = \frac{\hbar}{2} \alpha_1 \quad S_2 = \frac{\hbar}{2} \alpha_2 \quad S = S_1 + S_2 = \frac{\hbar}{2} (\alpha_1 + \alpha_2)$$

$$S_1 = \frac{1}{2} \quad S_2 = \frac{1}{2} \quad |S_1 - S_2| \leq S \leq S_1 + S_2$$

$$S = 0, 1$$

$$S^2 = \frac{\hbar^2}{4} (\alpha_1^2 + \alpha_2^2 + 2\alpha_1 \cdot \alpha_2) = \frac{\hbar^2}{4} (3I + 3[\alpha_1, \alpha_2])$$

$$S^2 = \frac{3}{2} \hbar^2 I + \frac{\hbar^2}{2} \alpha_1 \cdot \alpha_2 \quad \rightarrow \alpha_1 \cdot \alpha_2 = \frac{2}{\hbar^2} S^2 - 3I$$

$$S \cdot \hat{\mathbf{n}} = \frac{\hbar}{2} (\alpha_1 + \alpha_2) \cdot \hat{\mathbf{n}} = \frac{\hbar}{2} (\alpha_1 \cdot \hat{\mathbf{n}} + \alpha_2 \cdot \hat{\mathbf{n}})$$

$$(S \cdot \hat{\mathbf{n}})^2 = \frac{\hbar^2}{4} \left\{ (\alpha_1 \cdot \hat{\mathbf{n}})^2 + (\alpha_2 \cdot \hat{\mathbf{n}})^2 + 2(\alpha_1 \cdot \hat{\mathbf{n}})(\alpha_2 \cdot \hat{\mathbf{n}}) \right\}$$

$$\text{But } (\alpha_i \cdot \hat{\mathbf{n}})^2 = \alpha_i \cdot n_i \alpha_j \cdot n_j = \alpha_i \alpha_j r_i r_j = [\delta_{ij} I + i \epsilon_{ijk} \alpha_k] r_i r_j \\ = 0 + \hbar^2 I = I$$

$$(S \cdot \hat{\mathbf{n}}) = \frac{\hbar^2}{4} (I + I + 2(\alpha_1 \cdot \hat{\mathbf{n}})(\alpha_2 \cdot \hat{\mathbf{n}}))$$

$$\rightarrow (\hat{\alpha}_1 \cdot \hat{n})(\hat{\alpha}_2 \cdot \hat{n}) = \frac{?}{\hbar^2} (S \cdot \hat{n})^2 - I$$

$$S_{12} = \frac{6}{\hbar^2} (S \cdot \hat{n})^2 - 3I - \frac{2}{\hbar^2} S^2 + 3I$$

$$S_{12} = \frac{2}{\hbar^2} [3(S \cdot \hat{n})^2 - S^2]$$

It can be shown (Moshinsky 1958) :

$$S'_{12} = V(n) \left(\frac{32}{5}n\right)^{\frac{1}{2}} Y_2(\theta, \varphi) \cdot X_2$$

$Y_2(\theta, \varphi)$: Racah tensor with the components $Y_m^2(\theta, \varphi)$

$$X_2 : \quad " \quad " \quad \Rightarrow m=0 \text{ component } X_0^2 = \frac{1}{\sqrt{2}} (3S_z^2 - S^2)$$

3-5 Symmetry and Nuclear Force

One-body interaction: It acts only on a single body from an external source (centered force).

Ex The force acting on each electron in the atom from the nucleus (electrostatic force).

Two-body interaction: The interaction between a pair of particles at a time.

Ex.

The force between any pair of the electrons in the atom (electrostatic force).

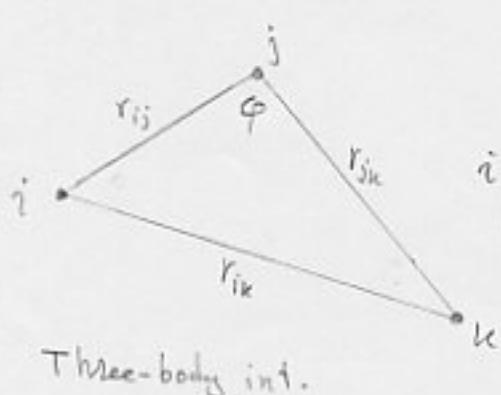
Three-body interaction: This interaction is present when at least there exist three particles.

Ex. The force between three nucleons in triton $\left\{ \begin{array}{l} {}^3\text{H} \\ {}^3\text{He} \end{array} \right.$ tritium system.

In a nucleus there is no fundamental one-body interaction.

The only one-body op. is Kinetic energy op.

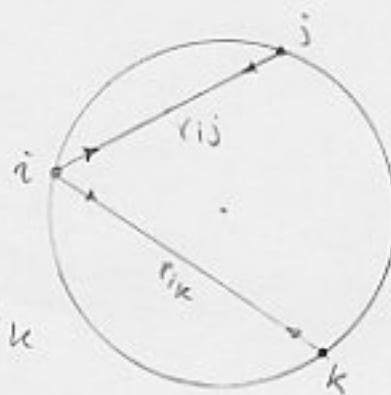
For convenience of carrying out calculations, one may, on occasion, construct an effective one-body term in the potential by taking an average of the two-body int. a particle experiences in the presence of all the other particles.



Three-body int.

$$V = V(r_{ij}, r_{jk}, r_{ik})$$

$$\text{or } V = V(r_{ij}, r_{jk}, \phi)$$



Two-body int

$$V = V(r_{ij})$$



one-body int.

$$V = V(r_i)$$

$$\rightarrow V = \sum_i V_i + \sum_{ij} V_{ij} + \sum_{ijk} V_{ijk}$$

i) With our present experimental equipment and theoretical knowledge

ii) And since $\sum_{ijk} V_{ijk} \ll \sum_{ij} V_{ij}$

→ it is unlikely to detect any three-body effect.

For this reason → we ignore any possible 3-body effect.

For two-body int. investigation we make use of deuteron,
but it is very limited system having only one bound state.

→ So, we will make use of scattering of one nucleon
off another one.

Charge Independence:

We assume the nuclear force is charge-independent.

i.e. the interaction of $n-n$, $p-p$ and $n-p$ are the same
if we remove the Coulomb interaction.

But still there are some symmetry-breaking effects, which
are small, but significant.

Ex.

Mass-difference of charged and neutral pions exchanged
between $p-n$ and $p-p$, $n-n$ causes a small difference
in their interactions.

The observed strengths of charge-symmetry breaking terms are except for a few highly specialized cases, smaller than or comparable to the accuracy we can achieve.

Charge independence implies that:

$$[H, T_z] = 0$$

$$\rightarrow T_z |\Psi\rangle = \frac{1}{2} (Z-N) |\Psi\rangle$$

$$\text{where } H |\Psi\rangle = E |\Psi\rangle$$

Charge symmetry breaking effects:

- 1) Coulomb
- 2) \rightarrow perturbation
- 3) Center of mass motion (The center of mass of nucleus (in particular P) does not coincide with the center of single particle potential well)
- 4) Finite size of Proton
- 5) \rightarrow neutron
- 6) Electromagnetic Spin-Orbit ($g_p \neq g_n$)
- 7) Vacuum polarization (virtual emission of e^-e^+ between P-P increases the repulsion)
- 8) p-n mass difference
- 9) Short range correlation
- 10) Isospin mixing in core

Isospin Invariance;

In addition to $[H, T_z] = 0$, for a nuclear Hamiltonian we have

$$[H, T^2] = 0$$

$$\rightarrow H|\Psi\rangle = E|\Psi\rangle$$

$$T^2|\Psi\rangle = T(T+1)|\Psi\rangle$$

Meaning:

$|\Psi\rangle$ remains unchanged if we replace some P with N or vice versa.

Ex.

$$A=16 \text{ system } T=1 \quad T_2 = \begin{cases} 1 & \left\{ \begin{array}{l} {}^{16}\text{O} \\ {}^{14}\text{N} \end{array} \right. \\ 0 & \left\{ {}^{15}\text{N} \right. \\ -1 & \left\{ {}^{14}\text{C} \right. \end{cases} \quad \begin{array}{ll} 8P-6N & \\ 7P-7N & \\ 6P-8N & \end{array}$$

${}^{16}\text{O} \xrightarrow[\text{into two } n]{\text{changing two } p} {}^{14}\text{C}$ members of isobar

Mathematically, the only difference in $|\Psi\rangle$ between the members is in their M_{Tz} -quanta number.

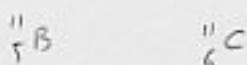
$$|{}^{16}\text{O}\rangle = T_+ |{}^{14}\text{N}\rangle = T_+ T_+ |{}^{14}\text{C}\rangle$$

$$|{}^{14}\text{C}\rangle = T_- |{}^{14}\text{N}\rangle \dots$$

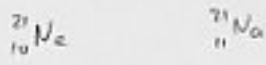
Isobaric Analogue states (IAS): A group of states related by a rotation in the isospin space.

Among the light nuclei there are many groups with the property $[H, T^2] = 0$

$3/2^+$	-	$3/2^-$
$5/2^+$	-	$5/2^-$
$7/2^+$	-	$7/2^-$
$9/2^+$	-	$9/2^-$
$1/2^+$	-	$1/2^-$
$3/2^+$	-	$3/2^-$
$5/2^+$	-	$5/2^-$
$7/2^+$	-	$7/2^-$
$9/2^+$	-	$9/2^-$



$1/2^+$	-	$3/2^+$
$3/2^+$	-	$1/2^+$
$5/2^+$	-	$3/2^+$
$7/2^+$	-	$5/2^+$
$9/2^+$	-	$7/2^+$



In heavy nuclei $F_{\text{Coulomb}} = 8t_{\text{strong}}$ (because of large number of p)

since; $[H^{\text{em}}, T^2] \neq 0$

$$\rightarrow T^2 |\Psi\rangle \neq T(T+1) |\Psi\rangle$$

$$\text{where } H|\Psi\rangle = E|\Psi\rangle$$

In contrast with the previous situation;

$$T^2 |\Psi_i\rangle = \sum_j \alpha_j |\Psi_j\rangle \quad (\langle \Psi_i |: \text{eigenstate of nuclear Hamiltonian})$$

In spite of the difficulty caused by electromagnetic interactions, the evidence for isospin invariance of the nuclear force itself is quite strong.

We will see;

Isospin symmetry limits the possible forms of nuclear potential.

Isospin Operators:

$$\mathcal{C}_1 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \quad \mathcal{C}_2 = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix} \quad \mathcal{C}_3 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$$

$$I = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$$

All other single-particle isospin operators can be expressed as a linear combination of I and \mathcal{C} .

For a system consisting of A particles;

$$T = \frac{1}{2} \sum_{i=1}^A \mathcal{C}_i$$

For two-particle system

$$T = \frac{1}{2} (\mathcal{C}_1 + \mathcal{C}_2)$$

Nuc. int. is $\left\{ \begin{array}{l} \text{two-body} \\ \text{scalar} \end{array} \right.$

T is one-body op., since it operates on a particle at a time.
Furthermore it is a vector.

$$T^2 = \frac{1}{4} (\mathcal{C}_1^2 + \mathcal{C}_2^2 + 2 \mathcal{C}_1 \cdot \mathcal{C}_2) = \frac{1}{4} (3I + 3I + 2 \mathcal{C}_1 \cdot \mathcal{C}_2)$$

$$\mathcal{C}_1 \cdot \mathcal{C}_2 = 2T^2 - 3I$$

$$\langle T | \mathcal{C}_1 \cdot \mathcal{C}_2 | T \rangle = \begin{cases} -3 & \text{for } T=0 \\ 1 & , \quad T=1 \end{cases}$$

mixed particle rank

$\rightarrow \gamma_1, \gamma_2$ may be used to express the isospin dependence of the nucl. force. (There are only two linearly indep. scalar two-body ops., I and γ_1, γ_2)

Ex.

We have already seen, a bound state is found for $T=0$, the deuteron, but not for $T=1$.

Remark: The property of isospin invariance is quite different from isospin dependence of the nucl. force (above example).

Other symmetries and General Form of Nucl. Potential:

- 1) The nucl. force between two nucleons must be invariant under a translation of two nucleons.

i.e. $V = V(|\vec{r}|=|\vec{r}_1 - \vec{r}_2|)$

- 2) Galilean invariance:

The potential may have \vec{p}_1 ad \vec{p}_2 dependence. But since $\vec{P} = \vec{p}_1 + \vec{p}_2$ corresponds to the center of mass momentum, it can not appear in V .

The only possibility is

$$\bar{P} = \frac{1}{2}(\bar{P}_1 - \bar{P}_2) \quad (\text{Galilean invariance})$$

- 3) U may depend on isospin.
- 4) V invariant under space rotation
- 5) V : time-reversal
- 6) " : space reflection (Parity)
- 7) " : permutation between two nucleons.

$$V = V(r, \alpha_1, \alpha_2, \gamma_1, \gamma_2, P)$$

Taking into account the symmetry requirements of nuclear potential, only a very limited number of linearly dependent set of single nucleon operators can be constructed.

Ex.

\vec{l} is not indep. variable, since $\vec{l} = \vec{r} \times \vec{p}$

It can be shown (Okubo and Marshall Ann. Phys. 4 (1958) 166) the most general two-body potential takes the form:

$$V(r; \alpha_1, \alpha_2, \gamma_1, \gamma_2) =$$

$$\begin{aligned}
 & V_o(r) + V_{\alpha}(r) \alpha_1 \alpha_2 + V_{\gamma}(r) \gamma_1 \gamma_2 + V_{\alpha\gamma}(r) (\alpha_1 \cdot \alpha_2)(\gamma_1 \cdot \gamma_2) \\
 & + V_{LS}(r) L \cdot S + V_{LS\gamma}(r) (L \cdot S)(\gamma_1 \cdot \gamma_2) \\
 & + V_T(r) S_{12} + V_{T\gamma}(r) S_{12} \gamma_1 \cdot \gamma_2 \\
 & + V_Q(r) Q_{12} + V_{Q\gamma}(r) Q_{12} \gamma_1 \cdot \gamma_2 \\
 & + V_{PP}(r) (\alpha_1 \cdot P)(\alpha_2 \cdot P) + V_{PP\gamma}(r) (\alpha_1 \cdot P)(\alpha_2 \cdot P)(\gamma_1 \cdot \gamma_2)
 \end{aligned}$$

Where $L \cdot S = \frac{1}{2} (l_1 \cdot l_2)(\alpha_1 \cdot \alpha_2)$

$$Q_{12} = \frac{1}{2} \left\{ (\alpha_1 \cdot L)(\alpha_2 \cdot L) + (\alpha_2 \cdot L)(\alpha_1 \cdot L) \right\}$$

To determine $V_o(r)$, $V_{\alpha}(r)$, ... we need information in addition to that generated from symmetry arguments above.

Procedure:

- { i) For example, we may make use of our knowledge of the basic nature of the nucle. force such as the meson-exchange picture of Yukawa.
- { ii) Fitting some assumed form of the radial dependence to experimental data.

The 12-terms in the potential may be divided into 5-groups:

i) $V_{\text{central}} = V_0(r) + V_\omega(r) \alpha_1 \alpha_2 + V_\chi(r) \chi_1 \chi_2 + V_{\text{cr}}(\alpha_1 \alpha_2)(\chi_1 \chi_2)$

Since tensorial ranks of the spatial parts of all 4-operators are zero, ($K=0$).

1) $V_0(r)$ depends only on r and is invariant under a rotation of coord. system.

2) The spatial dependence of the second term is also on r only, but it depends on intrinsic spin through the op. $\alpha_1 \alpha_2$.

$$\langle S | \alpha_1 \alpha_2 | S \rangle = \begin{cases} -3 & \text{for } S=0 \\ 1 & \text{, } S=1 \end{cases}$$

$S=0$ singlet state (scalar nucleon pair)

$S=1$ triplet = (vector " ")

There are two linearly indep. scalar two-body op:

$$I \text{ ad } \alpha_1 \alpha_2$$

3) Similar to case (2) but $\alpha_1 \alpha_2$ is replaced by $\chi_1 \chi_2$.

4) It has both spin and isospin dependence.

Since all ℓ -terms are scalars in intrinsic spin and orbital angular momentum

$$\rightarrow [V_{\text{central}} \mathbf{S}^2] = [V_{\text{central}} \mathbf{L}^2] = [V_{\text{central}} \mathbf{J}^2] = 0$$

ii) $V_{\text{spin-orbit}} = V_{LS}(r) \mathbf{L} \cdot \mathbf{S} + V_{LSc}(r) (\mathbf{L} \cdot \mathbf{S}) (\mathbf{\tau}_1 \cdot \mathbf{\tau}_2)$ $\left. \begin{array}{l} \text{non-central} \\ \mathbf{L} = \mathbf{L}(0, q) \end{array} \right\}$

$$[V_{S-O}, \mathbf{L}] \neq 0 \quad [V_{S-O}, \mathbf{S}] \neq 0 \quad [V_{S-O}, \mathbf{J}] = 0$$

i) The first term involves only $\mathbf{L} \cdot \mathbf{S}$ -term.

ii) The second term has additional op. $(\mathbf{\tau}_1 \cdot \mathbf{\tau}_2)$, isospin-dep. -

For example different mesons exchanged between nucleons, may or may not have an isospin dependent int.

The spatial part in both cases involves \mathbf{L} .

$$\mathbf{L} \xrightarrow{\square} \mathbf{L} \quad \text{axial vector}$$

To maintain parity invariance as well as rotational invariance of V ; we need another axial vector to obtain a scalar product.

$\mathbf{L} \cdot \mathbf{L}$ is not useful, since $[\mathbf{L}^2, \mathbf{L}] = [\mathbf{L}^2, \mathbf{S}] = 0$

\rightarrow which is a part of central force.

The only possibility: L · S

However, $\langle LS | L \cdot S | L' \cdot S' \rangle = 0$ for $L \neq L'$

i) Because rank of $L = 1 \rightarrow \langle \cdot \rangle = 0$ if $|L' - L| > 1$

→ For non-zero $\langle \cdot \rangle \quad L' = L \text{ or } L' = L \pm 1$

ii) On the other hand:

The parts of $\langle LS | = (-)^L$
" " $|L'S' \rangle = (-)^{L'}$

$$\begin{array}{ccc} S & \xrightarrow{n} & S \\ L & \xrightarrow{n} & L \end{array}$$

$$\Rightarrow \langle LS | L \cdot S | L' \cdot S' \rangle \xrightarrow{n} -\langle LS | L \cdot S | L' \cdot S' \rangle \quad \text{if } L' = L \pm 1$$

→ and must vanish.

iii) $V_{\text{Tensor}} = V_T(r) S_{12} + V_{T2}(r) S_{12} C_1 \cdot C_2$

iv) $V_{\text{quadratic}} = V_Q(r) Q_{12} + V_{Q2}(r) Q_{12} C_1 \cdot C_2$

The quadratic spin-orbit term enters only when there is
momentum dependence in potential.

$$V_{\text{inelastic}} = V_{pp}(r)(\alpha_1, p)(\alpha_2, p) + V_{pp\pi}(r)(\alpha_1, p)(\alpha_2, p)(\alpha_3, \vec{\alpha}_3)$$

gives no contributions in elastic interaction.

3-6 Yukawa Theory of Nuclear Interaction:

Beyond the symmetry arguments in the previous section, meson-exchange idea by Yukawa (1934) is useful starting point for the examination of nucleon-nucleon int.

A proper derivation of a boson exchange potential requires a relativistic quantum field theory treatment.



However, the essence may be obtained by drawing an analogy to classical electrodynamics.

$\Phi(r)$: electrostatic pot.

is a sol of $\nabla^2\Phi(r) = 0$ (1) (in source-free region)
Laplace's eqn.

In presence of a point source at origin:

$$\nabla^2\Phi(r) = -\left[\frac{1}{4\pi\epsilon_0}\right] 4\pi q\delta(r) \quad (2)$$

Poisson's eqn.

with the familiar sol. of Coulomb pot.

$$\Phi(r) = \left[\frac{1}{4\pi\epsilon_0}\right] \frac{q}{r}$$

When the electromagnetic field is quantized:

photons; emerge as the field quanta

charge; the source of the field

Our aim is to find an equ similar to (2) and its analogue in quantum field theory for a short-range nucleon-pot.

The equ. must be invariant under a Lorentz tr. so that to be correct in the relativistic limit as well.

→ The Schrödinger equ is ruled out

The field quanta exchanged between nucleons must be a boson, (because only bosons can be created and annihilated singly.)

A Fermion, on the other hand must be created and annihilated together with its antiparticle.)

The Dirac equ. is unsuitable (an equ for spin $\frac{1}{2}$ particles)

→ The prime candidate: Klein-Gordon equ.

$$E^2 = p^2 c^2 + m^2 c^4$$

We quantize this eqn. by:

$$\bar{E} \rightarrow i\hbar \frac{\partial}{\partial t}$$

$$\bar{p} \rightarrow -i\hbar \vec{\nabla}$$

(in the same way as in nonrelativistic quantum mechanics).

$$\rightarrow -\hbar^2 \frac{\partial^2}{\partial t^2} \varphi(r) = (-\hbar^2 c^2 \nabla^2 + m^2 c^4) \varphi(r)$$

m : mass of the field quanta.

$$\rightarrow \left(\nabla^2 - \frac{1}{c^2} \frac{\partial^2}{\partial t^2} \right) \varphi(r) = \frac{m^2 c^2}{\hbar^2} \varphi(r)$$

$$\rightarrow (\square + \frac{m^2 c^2}{\hbar^2}) \varphi(r) = 0$$

$$\text{where } \square = \partial_\mu \partial^\mu = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} - \vec{\nabla}^2$$

This is only analogous to (1), since it does not yet contain a source term for field quanta.

This point may be further demonstrated by:

$$\left\{ \begin{array}{l} m \rightarrow 0 \\ \text{and ignoring the time-dependence} \end{array} \right.$$

The result is quantized version of (1). (through commutating relations)

Including a source:

For simplicity, we consider only the static limit, and ignore terms involving time derivatives.

For point source with strength g located at origin:

$$\nabla^2 \varphi(\vec{r}) = \frac{m^2 c^2}{\hbar^2} \varphi(\vec{r}) - g \delta(\vec{r})$$

with sol. $\varphi(r) = \frac{g}{4\pi} \frac{e^{-\frac{mc}{\hbar r} r}}{r}$ Yukawa Pot.

In the limit $m \rightarrow 0$ and letting $g = [\frac{1}{4\pi\epsilon_0}]^{1/2} q$ it reduces to sol. of Coulomb potential.

If $m \neq 0$ the strength of the pot. drops by $\frac{1}{e}$ at $r_0 = \frac{\hbar}{mc}$: range of the force mediated by boson

For pions ($m \sim 140 \text{ MeV}$) $\rightarrow r_0 \sim 1.4 \text{ fm}$

We shall see that the exchange of a single pion gives a good representation of the long-range part of the nuclear pot.

3-7 Nucleon-Nucleon Scattering Phase Shifts:

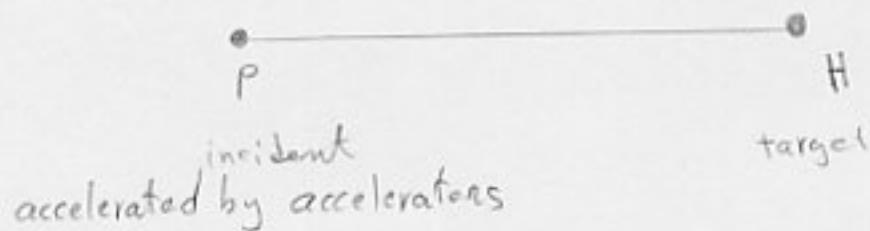
The general form of V in section 3-5 was obtained using the properties of the deuteron and symmetries of a two-nucleon system.

To make further progress, we need additional experimental information, and that is provided by scattering.

Nucleon-Nucleon Scattering:

There are 4-types of scattering:

i) p-p scattering;



ii) n-n scattering;

For incident; a) At low energies neutrons are used from nuclear reactions.

b) At higher energies, neutrons are used from proton bombardment

For example, through (p, n) reaction on a ^7Li target.

However, both { the intensity
energy resolution } of neutron beams obtained in
this way are much more limited than those of proton beams.

iii) The scattering of neutrons off proton target ($n-p$ scattering);
is an important source of information.

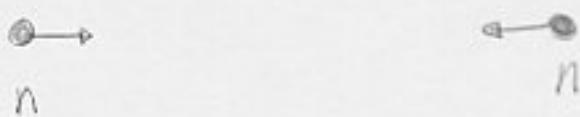
$p-p$ system or $n-n$ system can only be coupled to $T=1$,
whereas a $n-p$ system can be either $T=0$ or $T=1$.

iv) $p-n$ scattering.

Since neutron is unstable (half-life $\sim 10\text{ min.}$), it is not
possible to construct a fixed neutron target as it is with
protons.

Getting round of this limitation:

i) To carry out a colliding experiment.



Such an experiment requires high intensities in both beams
(which are not easily available).

2) Using deuterium.

Since the deuteron is a loosely bound system of n and p ;
 p_n or nn scattering results may be obtained by
 p_d or nd scattering.

$$p_n \sim p_d - (pp)$$

$$nn \sim nd - (np)$$

The procedure is valid provided:

- 1 - The accuracy in pp or np scattering data is comparable or better than p_d or nd data. The effect of deuteron binding energy can be corrected in a satisfactory manner.
- 2 - Three-body effects are negligible.

The information obtained from p_n and nn -scattering may not be any different from that obtained from np and pp -scattering.

For example the difference between np and p_n scattering is that the targets and incidents are replaced.

Under time-reversal invariance \rightarrow then two arguments should give identical results.

Isospin $\left\{ \begin{matrix} \text{up} \\ \text{down} \end{matrix} \right. \left. \begin{matrix} \xrightarrow{\text{time reversal}} \\ \text{up} \end{matrix} \right\} \text{down}$

Both nn and pp are $T=1$ systems.

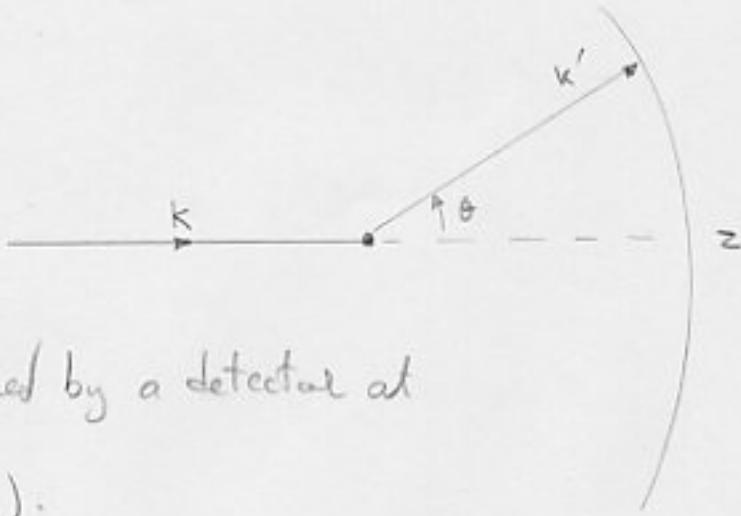
If nucl. force is charge-indep.:

$$\rightarrow \text{nn-scattering result} = (\text{pp-scattering result}) - \text{Coulomb int.}$$

But the accuracy in nn-scattering is still inadequate.

Scattering Cross Section:

The quantity measured in scattering experiment is:



The number of counts registered by a detector at a certain fixed angle (θ, φ) .

$$\text{The counting rate} \approx \Delta R, I, N, \frac{d\sigma}{d\Omega}$$

I: intensity of the incident beam

N: number of target nuclei

ΔR : The solid angle subtended by the detector.

$$\frac{d\sigma}{d\Omega} = \frac{d\sigma}{d\Omega}(k, \theta, \varphi)$$

In nonrelativistic limit; the scattering wave-func. is th sol.
of schrödinger equ. ;

$$-\frac{\hbar^2}{2\mu} \nabla^2 \Psi + (V - E) \Psi = 0 \quad \mu: \text{reduced mass}$$

For short-range nucl. forces $\rightarrow V=0$ except for a
small region
scattering region

$$\Psi(r, \theta, \varphi) \xrightarrow[r \rightarrow \infty]{} e^{ikz} + f(\theta, \varphi) \frac{e^{ikr}}{r}$$

We assume elastic scattering; $|k| = |k'|$

$$\frac{d\sigma}{dr}(\theta, \varphi) = |f(\theta, \varphi)|^2$$

For unpolarized incident beam ad target (non aligned spins),
the scattering is invariant with respect a rotation around z-axis,
(φ -indep.)

$$f(\theta, \varphi) \rightarrow f(\theta)$$

Partial wave analysis:

For a central pot. ℓ : conserved quantity in the reaction
(relative ang. momentum)

under such condns. \rightarrow it is useful to expand the wave func.
in the following way:

$$\Psi(r, \theta) = \sum_{\ell=0}^{\infty} a_{\ell} Y_{\ell 0}(\theta) R_{\ell}(kr)$$

Remark: In the absence of polarization $Y_{lm}(\theta, \varphi)$ contribute
with only $m=0$.

For free particle:

$$R_{\ell}(kr) \xrightarrow{\text{free}} J_{\ell}(kr) \xrightarrow{r \rightarrow \infty} \frac{1}{kr} \sin(kr - \frac{\ell\pi}{2})$$

$$k = \sqrt{\frac{2ME}{\hbar^2}}$$

For elastic scattering:

The probability current density = conserved
in each partial wave channel

$$R_{\ell}(kr) \xrightarrow[r \rightarrow \infty]{\text{scatt.}} \frac{1}{kr} \sin(kr - \frac{\ell\pi}{2} + \delta_{\ell})$$

$$f(\theta) = \frac{\sqrt{4\pi}}{k} \sum_{\ell=0}^{\infty} \sqrt{2\ell+1} e^{i\delta_{\ell}} \sin \delta_{\ell} Y_{\ell 0}(\theta)$$

$$\frac{d\sigma}{d\Omega} = \frac{4\pi}{k^2} \left| \sum_{\ell=0}^{\infty} \sqrt{2\ell+1} e^{i\delta_{\ell}} \sin \delta_{\ell} Y_{\ell 0}(\theta) \right|^2$$

$$\sigma = \int \frac{d\sigma}{dR} dR = \frac{4\pi}{k^2} \sum_{l=0}^{\infty} (2l+1) S^2 \delta_l(k)$$

Nucleon-nucleon Scattering phase shifts:

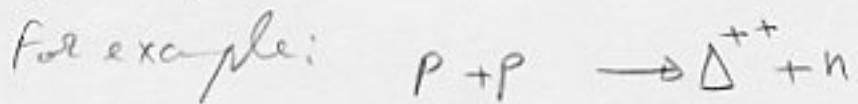
In realistic N-N scattering

1) Nucl. pot. in general depends on total intrinsic spin (S')

$\rightarrow J=l+s$ is conserved rather than l . (in scatt.)

For two nucleons $S=\{0, 1\}$ \rightarrow To determine S' we need to detect the orientation (polarization) of spins of the nucleons involved.

2) With sufficient energy, scattering can excite the internal degree of freedom of nucleons;



These are inelastic scatt.

We are dealing with identical particles:

$$\rightarrow \Psi_{(1,2)} = -\Psi_{(2,1)}$$

i) For PP-scatt. ; $T=1$ (sym), if $S=0$ (antisym)

\rightarrow Relative orbital wave func must be symmetric

$\rightarrow l: \text{even}$

For $S=0 \rightarrow J=l$

\rightarrow Partial waves for the lowest two orders are $\left\{ {}^1S_0, {}^1D_2 \right\}$

By the same token; for $S=1 \rightarrow l: \text{odd}$

The lowest order $l=1$ P-wave

$$S=1, l=1 \rightarrow J=0, 1, 2 \rightarrow \left\{ {}^3P_0, {}^3P_1, {}^3P_2 \right\} \quad \begin{array}{l} (\text{three-triplet state}) \\ l=1 \\ m=-1, 0, 1 \\ S=1 \\ m_s=-1, 0, 1 \end{array}$$

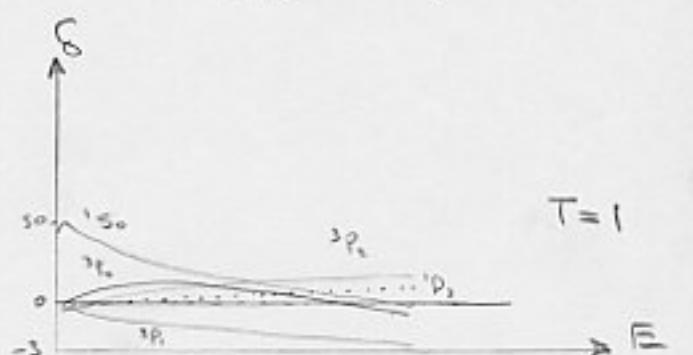
Contribution from
inelastic scatt. ≈ 0 for $E_{\text{scatt}} < 300 \text{ MeV}$
(imaginary part of δ)
in the lab frame

There is no admixture between

the two $J=0$ states $\left\{ {}^1S_0, {}^3P_0 \right\}$,

since they are of different

parity. $\rightarrow \left\{ \ell \right\}$ good quantum numbers



PP-scatt $E_{\text{scatt}} < 300 \text{ MeV}_{\text{lab}}$

(Coulomb effect removed)

ii) For np-scattering; $T=0$ or $T=1$

For $T=0$ (antisym), $S=0$ (antisym) $\rightarrow l: \text{odd}$

The lowest order $l=1$ ($l=3$ contribution in general ≈ 0)

For p-wave ($l=1$) np-scatt to be in $S=1 \rightarrow T=1$

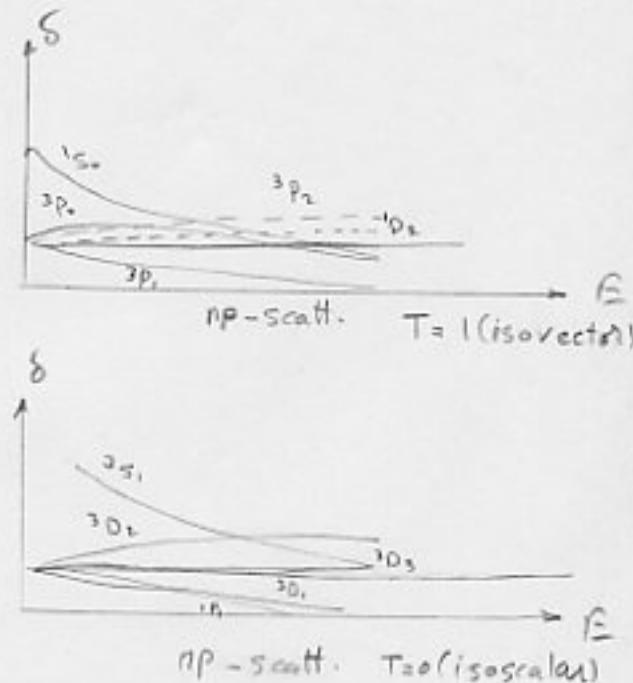
Then, the phase shifts in this case $(^1P_0, ^3P_1, ^3P_2)_{np}$ should be identical to those in pp-scatt. $(^3P_0, ^3P_1, ^3P_2)$; if:

- { 1 - Nucl. force is charge-indep.
- { 2 - Coulomb effects are removed.

The results are in good agreement.

But still there is a small difference, which is attributed to both:

- { 1 - experiment errors
- { 2 - charge dep. effects



For $T=0$ (antisym) $S=1$ (sym) $\rightarrow l: \text{even}$

$l = \text{even}$ \rightarrow makes it possible to have mixing of different J -partial waves (with the same J)

Up to now; each phase shift has been characterized by a definite l -value (as well as J -and S -values).
(even though the l is not fundamentally a good quantum number)

In general parity and other symmetry requirements restrict the mixing of different l -partial waves.

As in the case of deuteron;

Tensor force $\xrightarrow{\text{can cause}}$ admixture between two triplet states of the same J but different in l by two units ($l = J \pm 1$)

Now; for a given J , the scattering is now specified by two phase shifts;

δ_J for $l = J - 1$

δ_{J+1} for $l = J + 1$

and ϵ_J : which indicates the amount of mixing.

$$\epsilon = \delta(k) \quad \epsilon_J = \epsilon_J(k)$$

Using the def. of δ_J (convention) given in PhysRev. 105 (1957) 302,
the scattering matrix for a given J is written in the form;

$$\{S\} = \begin{pmatrix} e^{2i\delta_{J>}} G_{>2\epsilon_J} & i e^{i(\delta_{J>} + \delta_{J<})} S_{>2\epsilon_J} \\ i e^{i(\delta_{J>} + \delta_{J<})} S_{>2\epsilon_J} & e^{2i\delta_{J<}} G_{<2\epsilon_J} \end{pmatrix}$$

For the scattering from $l=J+1$ channel to $l=J+1$ channel the connection

is by $e^{2i\delta_J} = e^{2i\delta_{J>}} G_{>2\epsilon_J}$

From $l=J-1$ channel to $l=J-1$ channel:

$$e^{2i\delta_J} = e^{2i\delta_{J<}} G_{<2\epsilon_J}$$

On the other hand; the scattering from $l=J-1$ to $l=J+1$ and vice versa
is given by $e^{2i\delta_J} = e^{i(\delta_{J>} + \delta_{J<})} S_{>2\epsilon_J}$

These are generally referred to as the nuclear bar phase shifts.

Spin polarization in nucleon-nucleon scattering.

Scattering pot. may be $V = V(S)$

$$\rightarrow \alpha = \alpha(S) \quad \rightarrow \alpha(S=0) \neq \alpha(S=1)$$

It is not easy to observe the value of S directly in the experiment.

The detected quantity in the experiment is the orientation of the nucleon.

There are 4-possibilities in the initial state and in the final state.

$$|\frac{1}{2}, \frac{1}{2}\rangle, |\frac{1}{2}, -\frac{1}{2}\rangle, |\frac{1}{2}, -\frac{1}{2}\rangle, |-\frac{1}{2}, -\frac{1}{2}\rangle$$

If $V = V(S) \rightarrow$ the orientation of the nucleons may be changed as a result of scattering.

$\begin{cases} 4 \text{-different initial states,} \\ 4 \text{-final } \end{cases} \rightarrow 4 \times 4 \text{ different polarization measurements (in principle)}$

$$\{S\} = \begin{pmatrix} + & + & + & + \\ + & - & + & - \\ - & + & - & + \\ - & - & + & - \end{pmatrix}$$

16-quantities are not indep.

Time-reversal and other symmetries inherit in the scattering; leave only s -indep. matrix elements.

$$f_1 = f_{++,++} = f_{--,--} \quad f_2 = f_{++,--} = f_{--,++}$$

$$f_3 = f_{+-,+-} = f_{-+,-+} \quad f_4 = f_{+-,-+} = f_{-+,-+}$$

$$f_5 = f_{++,-+} = f_{-+,-} = f_{--,+-} = f_{-+,-+}$$

$$= f_{--,++} = f_{+-,++} = f_{++,-+} = f_{+-,--}$$

By the use of Transition op., T:

$$\langle K' | V | \Psi^+ \rangle = \langle K' | T | K \rangle$$

$$f(K, \theta) = f_{K,K'} = -\frac{\mu}{2\pi t^2} \langle K | V | \Psi^+ \rangle = -\frac{\mu}{2\pi t^2} \langle K' | T | K \rangle$$

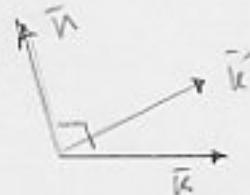
In place of f_i to f_5 , the T matrix element for nucleon-nucleon interaction is often written as a func. of five-coeffs., A, .. F in the following form:

$$T_{K,K}(1,2) = A + B \alpha_n(1) \alpha_n(2) + C \{ \alpha_n(1) + \alpha_n(2) \}$$

$$+ D \alpha_q(1) \alpha_q(2) + E \alpha_p(1) \alpha_p(2)$$

$\begin{cases} \vec{n} \\ \vec{p} \\ \vec{q} \end{cases}$ three vectors along which the nucleon spin components are taken.

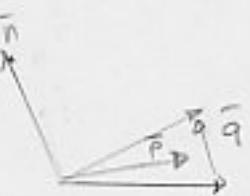
$$\hat{n} = \frac{\mathbf{k} \times \mathbf{k}'}{|\mathbf{k} \times \mathbf{k}'|} \quad \hat{q} = \frac{|\mathbf{k}' - \mathbf{k}|}{|\mathbf{k}' - \mathbf{k}|} \quad \hat{P} = \hat{q} \times \hat{n}$$



Ref.: Bystricky, Lehar, and Winteritz

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The amount of indep. information obtained from scattering is greatly increased with Polarization measurements.



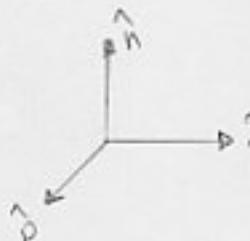
It is not difficult to obtain analyzing power (A_y) measurement, whereby a polarized beam is scattered from an unpolarized target (and the polarization of the particles in the final state is not detected).

At a fixed θ :

$$\sigma(S_{\text{incident}} \text{ parallel to } \hat{n}) - \sigma(S_{\text{incident}} \text{ anti-parallel to } \hat{n}) \neq 0 \quad \text{in general}$$

Such a difference is characterized by the analyzing power.

This supplies one of the independent quantities in the scattering.



The sum of two differential cross section supplies the other.

For additional information the polarization of the scattered particle must be measured.

Valuable information can also be obtained using polarized targets.

However this requires low temperature techniques to freeze the spin orientation of the nucleons in the target.

Inelastic Scattering:

If $K_{cm} >$ Required energy for meson production

$\left\{ \begin{array}{l} \text{Kinetic} \\ \text{energy of} \\ \text{incident in} \\ \text{cm.} \end{array} \right.$

→ The inelastic scattering become possible.

For the lightest mesons;

$$m_{\pi^\pm} \approx 140 \text{ Mev}$$

We expect pion production once the bombarding energy is above the threshold.

As the energy increases → $\left\{ \begin{array}{l} 1 - \text{Excitation of the internal deg. of freedom of the nucleon} \\ 2 - \text{Particle production} \end{array} \right.$
becomes more probable.

Inelastic scattering:

There is a loss in I (flux of the incident)

→ Probability is not conserved.

→ V : Complex

→ $\begin{cases} \text{Scattering amplitude} \\ \text{phase shifts} \end{cases}$: complex

$$f(\theta) = \sum_{\ell=0}^{\infty} \sqrt{4\pi(\ell\ell+1)} f_{\ell} Y_{\ell 0}(\theta)$$

$$\text{Since } S_{\ell}(k) \equiv 1 + 2ik f_{\ell}(k)$$

$$\rightarrow f_{\ell} = \frac{S_{\ell} - 1}{2ik}$$

For purely elastic scattering:

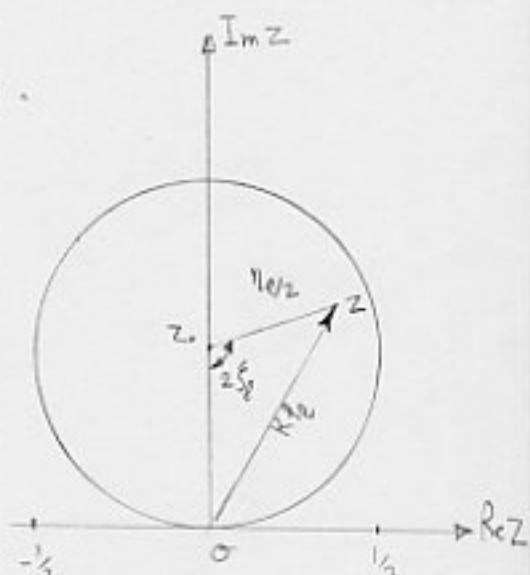
$$S_{\ell} = e^{i\delta_{\ell}} \quad |S_{\ell}| = 1$$

be real

$$f_{\ell} = \frac{e^{i\delta_{\ell}} - 1}{2ik} = \frac{e^{i\delta_{\ell}} - 1}{k} = \frac{1}{k \cot \delta_{\ell} - ik}$$

$$f(\theta) = \sum_{\ell=0}^{\infty} (2\ell+1) \left(\frac{e^{i\delta_{\ell}} - 1}{2ik} \right) P_{\ell}(\cos \theta)$$

$$= \frac{1}{k} \sum_{\ell=0}^{\infty} (2\ell+1) e^{i\delta_{\ell}} S_{\ell} P_{\ell}(\cos \theta)$$



Argand diagram

$$z_0(0, \frac{1}{2})$$

$$z = kf_{\ell} = \frac{1}{2i} (e^{i\delta_{\ell}} - 1)$$

$$Y_{\ell 0}(\theta) = \frac{P_{\ell}(\cos \theta)}{\sqrt{4\pi}}$$

In inelastic scattering:

$$S_e(k) \equiv n_e(k) e^{2i\delta_e(k)} \quad |S_e(k)| \leq 1$$

$$0 \leq n_e(k) \leq 1$$

$$f_e(k) = \frac{S_e(k)-1}{2ik} = \frac{n_e(k)e^{2i\delta_e(k)} - 1}{2ik} = \frac{n_e \Sigma \delta_e}{2k} + i \frac{1 - n_e \Sigma \delta_e}{2k}$$

$$\tilde{\rho}_e = 4\pi \sum_{l=0}^{\infty} (2l+1) |f_e(k)|^2 = 4\pi \sum_l (2l+1) \frac{1+n_e^2 - 2n_e \cos 2\delta_e}{4k^2}$$

We remember:

$$\Psi(\vec{r}) \rightarrow \frac{i}{2k} \sum_{l=0}^{\infty} (2l+1) i^l \underbrace{\left[\frac{e^{-i(kr - \frac{pl}{2})}}{r} - S_l(k) \frac{e^{i(kr - \frac{pl}{2})}}{r} \right]}_{\Psi_{\text{in}}} \underbrace{P_l(\cos\theta)}_{\Psi_{\text{out}}} \quad (1)$$

$$\Psi_{\text{in}} \sim \frac{i}{2k} \frac{e^{-ikr}}{r} P_l(\cos\theta)$$

$$J = \frac{i}{2im} (\Psi^* \nabla \Psi - \Psi \nabla \Psi^*) \quad \text{flux}$$

$$\text{Total flux} = \int J(\vec{r}) \cdot (d\vec{r}/dr) = \frac{i}{2im} \int dr (\Psi^* \frac{\partial \Psi}{\partial r} - \frac{\partial \Psi^*}{\partial r} \Psi)$$

$$\text{Total radial flux}_{\text{(in)}} = \left(\frac{\hbar k}{m} \right) \left(\frac{4\pi}{(2k)^2} \right) \quad \text{(inward)}$$

} Remark:
inward ≠ incident

$$\text{Total radial flux}_{\text{(out)}} = \left(\frac{\hbar k}{m} \right) |S_e(k)|^2 \frac{4\pi}{4k^2} \quad \text{for each } l\text{-value}$$

(outward)

$$\rightarrow \text{The net flux lost} = TRL_{\text{(in)}} - TRL_{\text{(out)}} \\ = \frac{\hbar k}{m} \left(\frac{n}{k^2} \right) \left(1 - n_e^2(k) \right) \quad \text{for each } l\text{-value}$$

Remark: Note that equ. (1) with the help of:

$$e^{ikr} \rightarrow \underbrace{\frac{i}{2k} \sum_{l=0}^{\infty} (2l+1) i^l}_{\text{incident}} \left[\underbrace{\frac{e^{-i(kr-\frac{pl}{2})}}{r}}_{\text{incoming}} - \underbrace{\frac{e^{i(kr-\frac{pl}{2})}}{r}}_{\text{outgoing}} \right] P_l(G_\theta)$$

can be cast into:

$$\Psi(r) \rightarrow e^{ikr} + \left[\underbrace{\sum_{l=0}^{\infty} (2l+1) \frac{S_l(k)-1}{2ik}}_{\text{outgoing}} P_l(G_\theta) \right] \frac{e^{ikr}}{r}$$

$$\text{Incident flux} = \frac{\hbar k}{m} \quad (\text{only in one dir.})$$

$$\text{Note that } \bar{J} = \frac{k}{2im} (\Psi^* \nabla \Psi - \Psi \nabla \Psi^*) = \frac{\hbar \bar{k}}{m} \quad \text{for } \Psi = e^{ikr}$$

$$\frac{\text{The net flux lost}}{\text{Incident flux}} = \frac{n}{k^2} \left(1 - n_e^2(k) \right) \quad \text{for each } l\text{-value}$$

$$\rightarrow \eta_{\text{inel}} = \frac{n}{k^2} \sum_l (2l+1) [1 - n_e^2(k)]$$

$$\begin{aligned}\omega_{\text{tot}} &= \omega_{\text{el}} - \omega_{\text{inel}} = \frac{n}{k^2} \sum_{\ell} (2\ell+1) \left(1 + \eta_{\ell}^2 - 2\eta_{\ell} \cos 2\beta_{\ell} + 1 - \eta_{\ell}^2 \right) \\ &= \frac{2n}{k^2} \sum_{\ell} (2\ell+1) (1 - \eta_{\ell} \cos 2\beta_{\ell}) \quad (1)\end{aligned}$$

Optical Theorem:

$$f(\theta) = \sum_{\ell=0}^{\infty} (2\ell+1) f_{\ell}(k) P_{\ell}(\cos \theta)$$

$$f(0) = \sum_{\ell=0}^{\infty} (2\ell+1) f_{\ell}(k)$$

$$\operatorname{Im} f(0) = \sum_{\ell=0}^{\infty} (2\ell+1) \operatorname{Im} f_{\ell}(k)$$

$$\begin{aligned}\operatorname{Im} f_{\ell}(k) &= \operatorname{Im} \left(\frac{\eta_{\ell} \sum 2\beta_{\ell}}{2k} + i \frac{1 - \eta_{\ell} \cos 2\beta_{\ell}}{2k} \right) \\ &= \frac{1 - \eta_{\ell} \cos 2\beta_{\ell}}{2k}\end{aligned}$$

$$\operatorname{Im} f(0) = \sum_{\ell} (2\ell+1) \frac{1 - \eta_{\ell} \cos 2\beta_{\ell}}{2k} \quad (2)$$

$$(1)(2) \rightarrow \operatorname{Im} f(0) = \frac{k}{4n} \omega_{\text{tot}}$$

So the optical tho. is satisfied.

i) If $\eta_{\ell}(k) = 1 \rightarrow$ no absorption $\omega_{\text{inel}} = 0$ purely elastic scatt.

ii) If $\eta_{\ell}(k) = 0 \rightarrow$ total absorption, however $\omega_{\text{el}} \neq 0$

Consider a black disc;

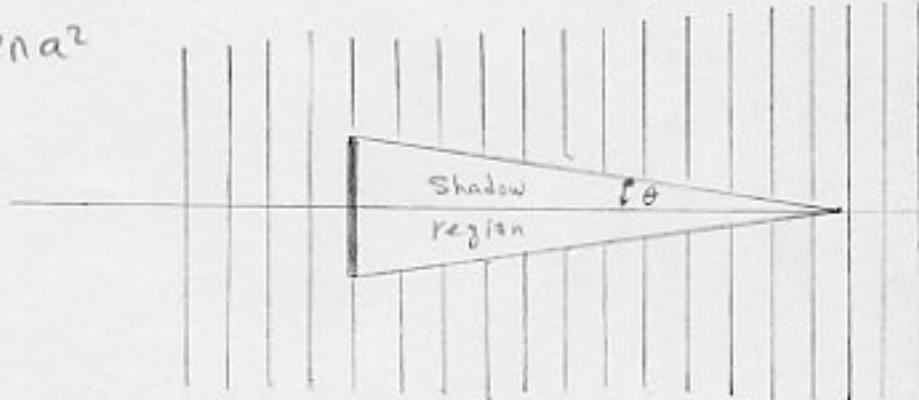
Black disc $\left\{ \begin{array}{l} 1 - \text{It has well defined edge} \\ 2 - \text{it is totally absorbing} \end{array} \right.$

Since $\left\{ \begin{array}{l} L = \sqrt{\ell(\ell+1)} h \approx \ell h \\ L = a\rho = a\ell k \end{array} \right.$ $\rightarrow \ell = \frac{a}{k}$ a : the radius of disc

$$\eta_e(k) = 0 \quad (\text{total absorption}) \rightarrow \sigma_{inel} = \frac{\pi}{k^2} \sum_{\ell=0}^{ka} (2\ell+1) = \frac{\pi}{k^2} (ka)^2 = \pi a^2$$

$$\eta_e(k) = 0 \quad \rightarrow \sigma_{el} = 4\pi \sum_{\ell=0}^{ka} (2\ell+1) \frac{1}{4k^2} = \pi a^2$$

$$\sigma_{tot} = \sigma_{el} + \sigma_{inel} = 2\pi a^2$$



Classically: One expects no elastic scatt. when there is total absorption.

Quantum mechanically:

Disc takes flux $\sim \pi a^2$ out of the incident beam.

However; far away, shadow gets filled (by diffraction of incident wave at the edge of disc $\sim \pi a^2$, which is elastically).

This elastic scattering is also called shadow scattering

It is strongly peaked up forward.

The angle to which it is confined can be estimated from the uncertainty principle:

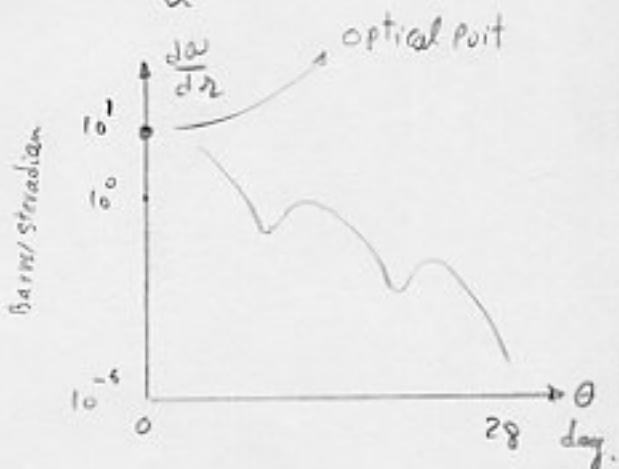
$$\Delta x \Delta p \sim \hbar$$

An uncertainty in the lateral dir of magnitude a

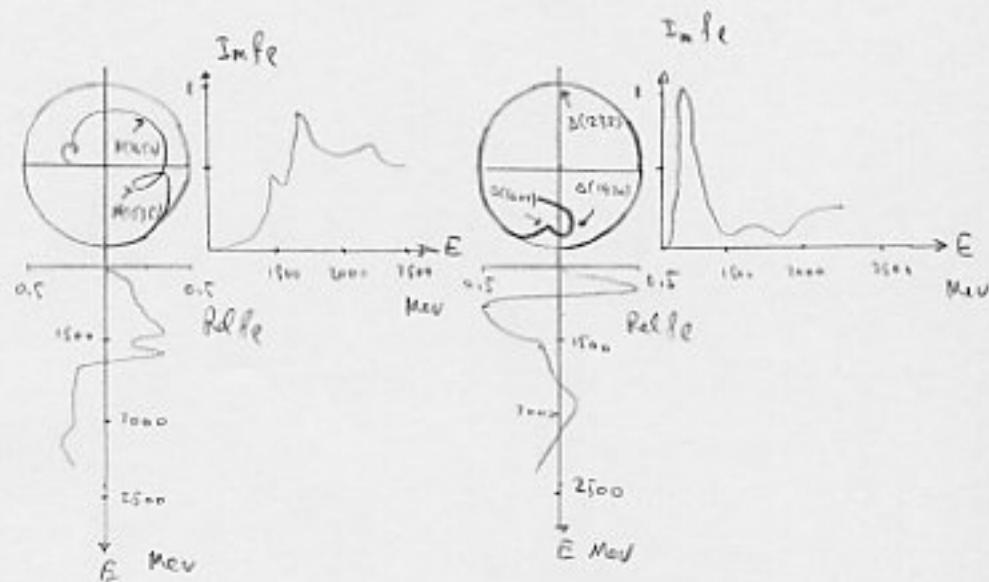
will be accompanied by \rightarrow an uncontrolled lateral momentum $P_{\perp} \sim \frac{\hbar}{a}$

$$P_{\perp} \approx P\theta \quad \text{small } \theta \quad \rightarrow \theta \sim \frac{\hbar}{ap} \sim \frac{1}{aK}$$

This agree with the optical result $\theta \sim \frac{\lambda}{a}$



scatt.; proton of ^{16}O nuclei
1000 Mev (1BeV)



$$\ell=0, J=\frac{1}{2}, T=\frac{1}{2}$$

$$\ell=1, J=\frac{3}{2}, T=\frac{3}{2}$$

(strong P_{33} resonance)

pion-nucleon scatt.

phase shifts \approx Real for $E_{\text{scatt.}} < 300 \text{ Mev}$ in the lab frame
 $\approx 150 \text{ Mev} \approx \text{CM} \approx$

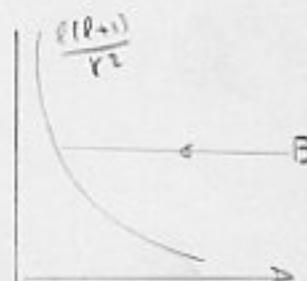
3-8) Low Energy Scattering Parameters:

$$-\frac{\hbar^2}{2\mu} \nabla^2 \psi + (V - E)\psi = 0$$

$$\psi(r, \theta) = \sum_{\ell=0}^{\infty} a_{\ell} Y_{\ell 0}(\theta) R_{\ell}(kr)$$

$$-\frac{\hbar^2}{2\mu} \left\{ \frac{1}{r^2} \frac{d}{dr} r^2 \frac{d}{dr} - \frac{\ell(\ell+1)}{r^2} \right\} R_{\ell}(kr) + V(r) R_{\ell}(kr) = E R_{\ell}(kr)$$

$$\hat{V}(r) = V(r) + \frac{\hbar^2}{2\mu} \frac{\ell(\ell+1)}{r^2} \quad \text{effective pot.}$$



$$U_e(k, r) = V R e(k, r)$$

$$\frac{1}{r^2} \frac{\partial}{\partial r} r^2 \frac{\partial}{\partial r} = \frac{d^2}{dr^2} + \frac{2}{r} \frac{d}{dr}$$

$$\rightarrow \frac{d^2}{dr^2} U_e + \left(k^2 - \frac{2M}{\hbar^2} V(r) - \frac{l(l+1)}{r^2} \right) U_e = 0$$

At low energies ($E < 10 meV$), only $l=0$ (s-wave) has important contribution.

$$\text{Def. : } \lim_{E \rightarrow 0} \alpha = 4\pi a^2 \quad a: \text{scattering length}$$

\uparrow
remains finite

In low energy scatt. α is often used instead of δ_0 .

They are related by

$$\alpha = \lim_{K \rightarrow 0} \text{Re} \left\{ - \frac{e^{i\delta_0} S(\delta_0)}{K} \right\} \quad \text{where } K^2 = \frac{2ME}{\hbar^2}$$

(convention here: for attractive V , $\rightarrow \alpha > 0$ for bound state)

The energy dependence of δ_0 at low energy is given by effective range parameter, r_e , defined by:

$$K \cot \delta_0 = -\frac{1}{\alpha} + \frac{1}{2} r_e K^2$$

$\{\alpha, r_e$ provide a useful way to parametrize information on low energy scatt..

Effective Range:

The energy dep. of scatt. at low energies is given by effective range.

For $\ell = 0$

$$\frac{d^2 U_0(k, r)}{dr^2} - \left\{ \frac{2\mu}{\hbar^2} V(r) - k^2 \right\} U_0(k, r) = 0 \quad (1)$$

$$\rightarrow \begin{cases} \ddot{U}_0(k_1, r) - \left\{ \frac{2\mu}{\hbar^2} V(r) - k_1^2 \right\} U_0(k_1, r) = 0 & E_1 = \frac{\hbar^2 k_1^2}{2\mu} \\ \ddot{U}_0(k_2, r) - \left\{ \frac{2\mu}{\hbar^2} V(r) - k_2^2 \right\} U_0(k_2, r) = 0 & E_2 = \frac{\hbar^2 k_2^2}{2\mu} \end{cases}$$

$$\rightarrow \int_0^\infty \left\{ U_0(k_2, r) \dot{U}_0(k_1, r) - U_0(k_1, r) \dot{U}_0(k_2, r) \right\} dr + (k_1^2 - k_2^2) \int_0^\infty U_0(k_1, r) U_0(k_2, r) dr = 0$$

Integration by parts:

$$\begin{aligned} & \left\{ U_0(k_2, r) \dot{U}_0(k_1, r) - U_0(k_1, r) \dot{U}_0(k_2, r) \right\} \Big|_0^\infty \\ &= (k_2^2 - k_1^2) \int_0^\infty U_0(k_1, r) U_0(k_2, r) dr \end{aligned} \quad (2)$$

This is true for arbitrary $V(r)$ including $V(r) = 0$

Now consider $V(r) = 0$ with the sol. $V_0(k, r)$

$$\frac{d^2 V_0(k, r)}{dr^2} + k^2 V_0(k, r) = 0 \quad (3)$$

Analogously:

$$\left\{ V_0(k_2, r) V'_0(k_1, r) - V_0(k_1, r) V'_0(k_2, r) \right\} \Big|_0^\infty \\ = (k_2^2 - k_1^2) \int_0^\infty V_0(k_1, r) V_0(k_2, r) dr \quad (4)$$

If $V(r)$ in eqn (1) is short in range; (1) and (3) are identical to each other in the asymptotic region.

$\rightarrow V_0(k, r) = U_0(k, r) = A \text{Si}(kr + \delta_0)$ Remark:
δ₀ is due to V(r)
for U(kr),
but for V(kr)
is just a change in
ref. point.

$\left\{ \begin{array}{l} U_0(k, r) \xrightarrow[r \rightarrow 0]{} 0 \\ U_0(k, r) \xrightarrow[r \rightarrow \infty]{} V_0(k, r) \end{array} \right.$

$$\rightarrow \left\{ U_0(k_2, r) U'_0(k_1, r) - U_0(k_1, r) U'_0(k_2, r) \right\} \Big|_0^\infty \\ = \lim_{r \rightarrow \infty} \left\{ V_0(k_2, r) V'_0(k_1, r) - V_0(k_1, r) V'_0(k_2, r) \right\} \quad (5)$$

(2), (4), (5) \rightarrow

$$V_o(k_1, 0) V'_o(k_2, 0) - V_o(k_2, 0) V'_o(k_1, 0)$$
$$= (k_1^2 - k_2^2) \int_0^\infty \{ V_o(k_1, r) V_o(k_2, r) - U_o(k_1, r) U_o(k_2, r) \} dr$$

(6)

Note that $V_o(k, 0) \neq 0$

Assuming ; $V_o(k, 0) = 1 \rightarrow V_o(k, r) = \frac{\sin(kr + \delta_o)}{\sin \delta_o}$

$$(6) \rightarrow V'_o(k_2, 0) - V'_o(k_1, 0)$$

$$= (k_1^2 - k_2^2) \int_0^\infty \{ V_o(k_1, r) V_o(k_2, r) - U_o(k_1, r) U_o(k_2, r) \} dr$$

$$\frac{k_2 \cot \delta_o(k_2) - k_1 \cot \delta_o(k_1)}{k_2^2 - k_1^2} = \int_0^\infty \{ V_o(k_1, r) V_o(k_2, r) - U_o(k_1, r) U_o(k_2, r) \} dr$$

If we choose $E_1 \approx E_2 \approx E$

$$\frac{d}{d(k^2)} k \cot \delta_o = \int_0^\infty \{ V_o^2(k, r) - U_o^2(k, r) \} dr \quad (7)$$

$$r_e \equiv 2 \int_0^\infty \{ V_o^2(k, r) - U_o^2(k, r) \} dr \quad \text{effective range.}$$

$$(7) \rightarrow K \cot \delta_0 = \int d(k) \frac{v_e}{2} + C$$

$$K \cot \delta_0(k) = (K \cot \delta_0)_{k=0} + \frac{1}{2} v_e k^2 + \dots$$

$v_e = v_e(k)$

Now using this def.;

$$a = \lim_{k \rightarrow 0} \operatorname{Re} \left\{ -\frac{1}{k} e^{i\delta_0} \right\} = \lim_{k \rightarrow 0} \operatorname{Re} \left\{ \frac{-1}{K \cot \delta_0(k) - ik} \right\}$$

$$a = \frac{1}{K \cot \delta_0(k) - \frac{1}{2} v_e k^2} \quad \rightarrow \quad K \cot \delta_0(k) = -\frac{1}{a} + \frac{1}{2} v_e k^2$$

$$\omega_0 = \frac{4\pi}{k^2} \sin^2 \delta_0(k) = \frac{4\pi}{k^2 + \left\{ \frac{1}{2} v_e k^2 - \frac{1}{a} \right\}^2}$$

$$\omega_0 \rightarrow 4\pi a^2$$

$k \rightarrow 0$

Neutron Scattering off hydrogen molecules:

In H-H molecule, (homonuclear molecule);

$$d_{\text{H-H}} = 7.8 \times 10^{-11} \text{ m} \gg \text{nuc. short range}$$

→ No. nuc. int between two protons in H₂.

$$\Psi_{\text{H}_2} = \Psi(r) \chi_s \text{ antisym.}$$

For this reason → Spin orientations of two protons are correlated with their relative orbital angular momentum.

→ Such a correlation may be exploited for neutron-proton scatt. length measurements.

There are two low-lying states for hydrogen molecule.

lower one: para-hydrogen $\Psi_{p-p}(r) = \text{sym.} \rightarrow S_z = 0$

higher one: Ortho - $\Psi_{p,p}(r) = \text{antisym.} \rightarrow S_z = 1$

$$S=1 \left\{ \begin{array}{l} S_2=1 \\ S_2=0 \\ S_2=-1 \end{array} \right. \quad S=0 \left\{ \begin{array}{l} S_2=0 \end{array} \right.$$

→ Statistical weight of (S=1) = 3 · (statistical weight of (S=0)=1)
(in a sample in equilibrium at room temp.)

But at low temperatures H_2 tends to go into the lowest energy state

→ Almost completely in a para-hydrogen state

→ At temp. T , we have $\alpha(T)(H_2)_{\text{para}} + \beta(T)(H_2)_{\text{ortho}}$

Intense neutron flux is available at low energies;

→ high precision measurements of neutron scatt. from H_2 .

By lowering the energy; $\lambda_n = \text{long} > d \sim d$

so that → scatt. off the two protons in a hydrogen molecule
is a coherent one (interference can be seen)

If neutrons are low in energy → little energy is transferred
to hydrogen target.

Remark: Energy received by target may cause transitions
from para $\xrightarrow{\text{to}}$ ortho-hydrogen states, and
this decreases the accuracy.

For this reasons; the neutron energy is kept low $\sim 0 \text{ mev} (10^{-3} \text{ eV})$

Corresponding to $T = 100^\circ \text{K}$

At this limit:

$$K \text{ and } \delta_0 \approx -\frac{1}{a} + \left(\frac{1}{2} r_0 K^2\right)$$

And the scatt. is characterized by two np-scatt. length alone.

For $\ell = 0$, NP-system is:

either $\begin{cases} S=0 & (\text{singlet state}) \\ T=1 \end{cases}$ or $\begin{cases} S=1 & (\text{triplet state}) \\ T=0 \end{cases}$

$$S = \frac{1}{2} (\alpha_n \cdot \alpha_p)$$

$$\text{Eigenvalue of } \alpha_n \cdot \alpha_p = \begin{cases} 1 & \text{for } S=1 \\ -3 & \Rightarrow S=0 \end{cases}$$

Define the projection operators:

$$P_t = \frac{1}{4} (3 + \alpha_n \cdot \alpha_p) \quad P_s = \frac{1}{4} (1 - \alpha_n \cdot \alpha_p)$$

$$\begin{cases} P_t |S=1\rangle = 1 |S=1\rangle \\ P_t |S=0\rangle = 0 |S=0\rangle \end{cases} \quad \begin{cases} P_s |S=1\rangle = 0 |S=1\rangle \\ P_s |S=0\rangle = 1 |S=0\rangle \end{cases}$$

The scatt. length with spin-dep of the nuc. forces:

$$\alpha = \alpha_t P_t + \alpha_s P_s = \frac{1}{4} (3\alpha_t + \alpha_s) + \frac{1}{4} (\alpha_t - \alpha_s) \alpha_n \cdot \alpha_p$$

For slow neutrons; ($\lambda \gg r_0$), we can take zero-range approx. (contact int., Breit 1947, S-func.)

$$V = \sum_i \left(\frac{2\pi t^2}{\mu} \right) a^{(i)} \delta(r_n - r_i) \quad (1)$$

i: the i-th proton

When we take the effects of E_b in Hz; different powers of $\frac{a}{d}$ or $\frac{a}{\lambda}$ show in V;

$\frac{a}{d}, (\frac{a}{d})^2, \dots$ corresponds to different orders of perturbations
or $\frac{a}{\lambda}, (\frac{a}{\lambda})^2, \dots$ in (1)

The cond. for applicability of (1) is;

$$a \ll d, a \ll \lambda$$

Moreover, in order to disregard finite-range of nucl. forces
(i.e. $V \rightarrow \delta(r)$) we must have theconds.;

$$r_0 \ll d \quad \text{and} \quad r_0 \ll \lambda$$

r_0 : effective range of nucl. force.

$$V = \frac{n\hbar^2}{2\mu} \left[(3\alpha_t + \alpha_s) + (\alpha_t - \alpha_s) \tilde{\omega}_n \cdot \tilde{\omega}_1 \right] \delta(r_n - r_1) \\ + \frac{n\hbar^2}{2\mu} \left[(3\alpha_t + \alpha_s) + (\alpha_t - \alpha_s) \tilde{\omega}_n \cdot \tilde{\omega}_2 \right] \delta(r_n - r_2) \quad (2)$$

$$\rightarrow V = V_S + V_A$$

$$V_S = \left(\frac{n\hbar^2}{2\mu} \right) \left[(3\alpha_t + \alpha_s) + (\alpha_t - \alpha_s) \tilde{\omega}_n \cdot S_H \right] [\delta(r_n - r_1) + \delta(r_n - r_2)] \quad (3)$$

$$V_A = \left(\frac{n\hbar^2}{2\mu} \right) [(\alpha_t - \alpha_s) \tilde{\omega}_n \cdot (\tilde{\omega}_1 - \tilde{\omega}_2)] [\delta(r_n - r_1) - \delta(r_n - r_2)]$$

$$(S_H = \frac{1}{2}(\tilde{\omega}_1 + \tilde{\omega}_2))$$

V_S : responsible for transitions from $\begin{cases} \text{Para-H to Para-H} \\ \text{or} \\ \text{Ortho-H to Ortho-H} \end{cases}$

(spin-sym of H_2 doesn't change)

V_A : responsible for transitions from $\begin{cases} \text{Para-H to Ortho-H} \\ \text{and} \\ \text{Vice Versa} \end{cases}$

For slow neutrons $\lambda \gg d \rightarrow$ Two protons can be considered at a single position $r_1 \approx r_2 \equiv r'$

$$\rightarrow [\delta(r_n - r_1) - \delta(r_n - r_2)] \approx 0$$

$$\rightarrow V_A \approx 0$$

Having V_S , only elastic scatt. of neutrons is possible.

(because of orthogonality of wave fns ad since $V_S \neq \text{dep}(r_1 - r_2)$)

↑
Protons

Comparing (1) and (2) \rightarrow

$$\alpha_H = \frac{1}{2} [(3\alpha_t + \alpha_s) + (\alpha_t - \alpha_s) \alpha_n \cdot S_H]$$

$$\langle \alpha_H^2 \rangle = \frac{1}{4} \left\{ (3\alpha_t + \alpha_s)^2 + 2(3\alpha_t + \alpha_s)(\alpha_t - \alpha_s) \langle \alpha_n \cdot S_H \rangle \right. \\ \left. + (\alpha_t - \alpha_s)^2 \langle (\alpha_n \cdot S_H)^2 \rangle \right\}$$

$$(\alpha_n \cdot S_H)^2 = \alpha_{nx}^2 S_{Hx}^2 + \alpha_{ny}^2 S_{Hy}^2 + \alpha_{nz}^2 S_{Hz}^2 \\ + 2 \alpha_{nx} \alpha_{ny} S_{Hx} S_{Hy} + 2 \alpha_{ny} \alpha_{nz} S_{Hy} S_{Hz} + 2 \alpha_{nz} \alpha_{nx} S_{Hz} S_{Hx}$$

For unpolarized reaction beam; $|N\rangle = \alpha \begin{pmatrix} 1 \\ 0 \end{pmatrix}_s + \beta \begin{pmatrix} 0 \\ 1 \end{pmatrix}_s \quad (\alpha = \beta)$

$$\langle \text{last 3-terms} \rangle = 0$$

$$\text{Since } \alpha_i^2 = 1$$

$$\langle \alpha_{nx}^2 S_{Hx}^2 + \alpha_{ny}^2 S_{Hy}^2 + \alpha_{nz}^2 S_{Hz}^2 \rangle = \langle S_{Hx}^2 + S_{Hy}^2 + S_{Hz}^2 \rangle \\ = \langle S_H^2 \rangle = S_H(S_H+1)$$

Also

$$\langle \alpha_n \cdot S_H \rangle = 0 \quad \text{for unpolarized beam.}$$

$$\langle \alpha_H^2 \rangle = \frac{1}{4} \left\{ (3\alpha_t + \alpha_s)^2 + (\alpha_t - \alpha_s)^2 S_H(S_H+1) \right\}$$

$$\alpha_H = 4/7 \langle \alpha_H^2 \rangle = 1/7 \left\{ (3\alpha_t + \alpha_s)^2 + (\alpha_t - \alpha_s)^2 S_H(S_H+1) \right\}$$

For para-H; $S_{\text{II}} = 0$

$$\alpha_{\text{para}} = \pi (3a_t + a_s)^2$$

For ortho-H; $S_{\text{II}} = 1$

$$\alpha_{\text{ortho}} = \pi (3a_t + a_s)^2 + 2\pi (a_t - a_s)^2$$

Using the experimental data for α_{para} and α_{ortho} one may obtain,

$|3a_t + a_s|$ and $|a_t - a_s| \rightarrow |a_t|, |a_s|$ and the sign of $\frac{a_t}{a_s}$

Neutron-Proton Scattering length:

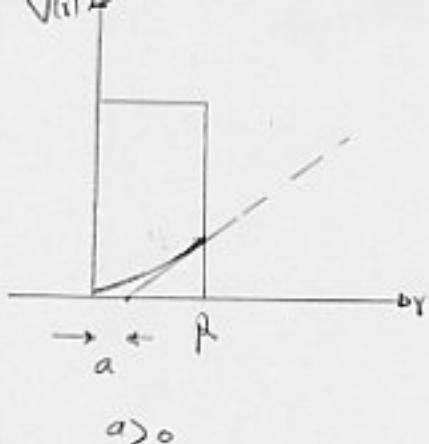
np system $\in T=1$

$$T=1 \quad \begin{cases} T_2=1 \text{ PP-system} \\ T_2=0 \text{ NP-} \\ T_2=-1 \text{ NN-} \end{cases}$$

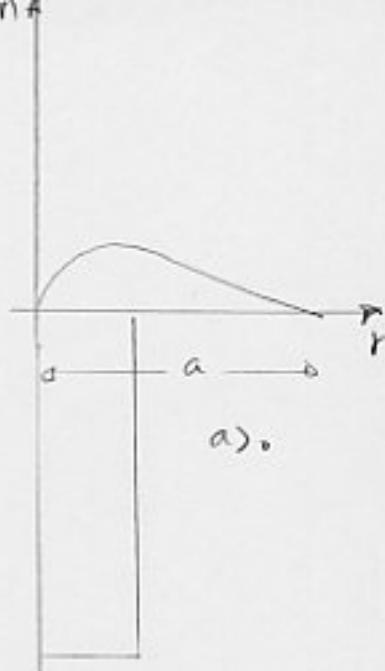
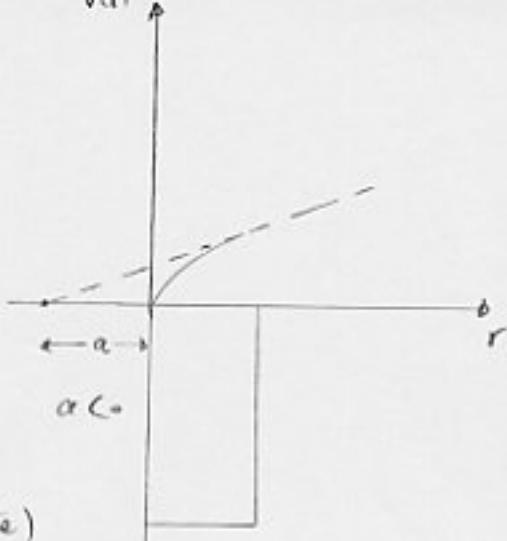
	$S=0, T=1$	$S=1, T=1$
a_{pp}	$-17.1 \pm 0.2 \text{ fm}$	-
r_e	$2.794 \pm 0.015 \text{ fm}$	-
a_{nn}	$-16.6 \pm 0.6 \text{ fm}$	-
r_e	$2.84 \pm 0.03 \text{ fm}$	-
a_{np}	$-23.71 (\pm 0.015) \text{ fm}$	$5.423 \pm 0.005 \text{ fm}$
r_e	$2.73 \pm 0.03 \text{ fm}$	$1.73 \pm 0.02 \text{ fm}$

Then we can compare a_{np} with a_{nn} and a_{pp} .

$$a = \lim_{k \rightarrow 0} \operatorname{Re} \left\{ -\frac{1}{k} e^{i\delta_0} \sin \delta_0 \right\}$$



(convention: $\alpha < 0$ bound state)



a : the intercept of outside wave-function with r -axis.

The sign of all three scatt. lengths are negative (in $T=1$)

→ No bound state (repulsive force) as we faced earlier.

a_{pp} is easily measured from low energy p-scatt. off H-target

However; Since,

$$\left(\frac{d\omega}{dr} \right)_{\text{Rutherford}} = \left\{ \left[\frac{1}{4\pi e^2} \right] \frac{zZ e^2}{4T z^2(\xi)} \right\}^2 = \left\{ \frac{4\pi c z Z}{4T} \frac{1}{z^2(\xi)} \right\}^2$$

$$\left(\frac{d\omega}{dr} \right)_{\text{Coulomb}} \sim \frac{1}{T^2}$$

pp-scattering at low energies is dominated by electromagnetic effects.

→ The accuracy is limited

$$\alpha_{\text{nucl.}} = \alpha_{\text{total}} - \alpha_{\text{Coulomb}}$$

↓
small ↓
large

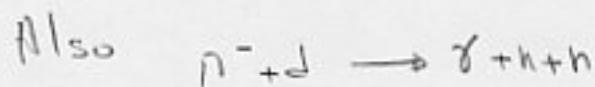
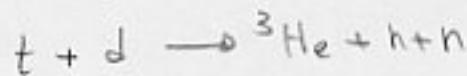
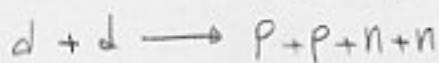
$$\alpha_{pp} = -7.823 \pm 0.01 \text{ fm} \quad \alpha_{pp} = -17.1 \pm 0.2 \text{ fm}$$

total nucl.

Measurements of α_{nn} are complicated by the absence of fixed
reaction targets.



(t: triton, d: deuterons)



due to the availability of good quality pion beams in the recent years
is used to reduce the uncertainty of α_{nn} .

α_{nn} is deduced from the int. of two n in the final state.

Comparison of α -values in $T=1$ states:

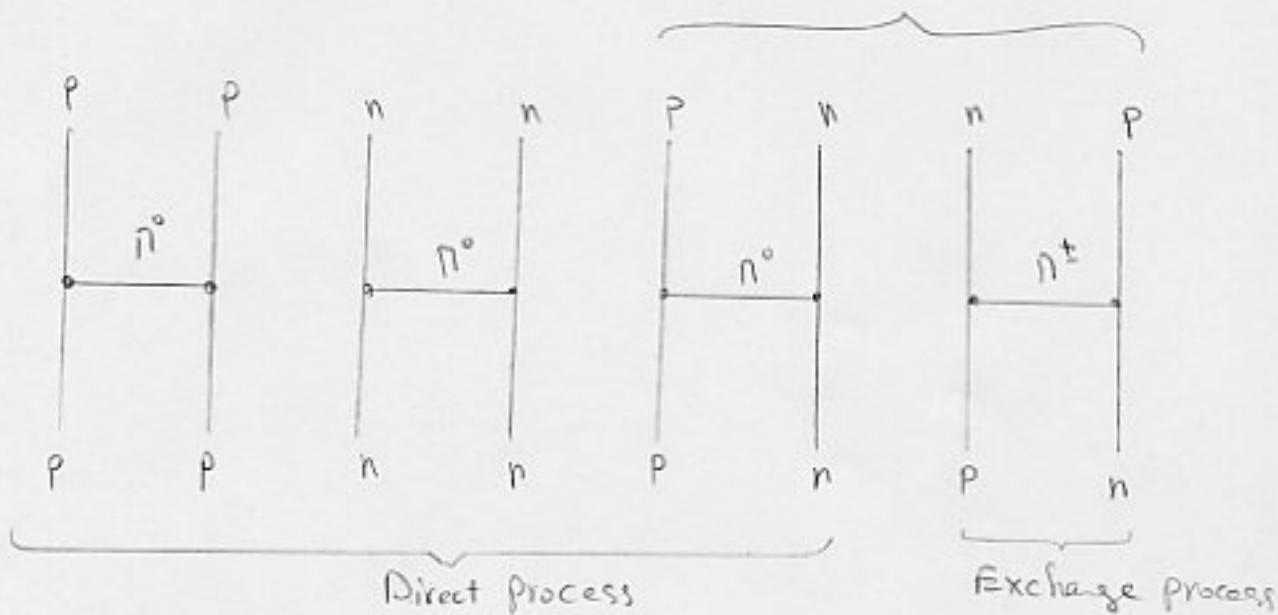
$$\alpha_{np} = -23.715 \pm 0.015 \text{ fm}$$

There is a big difference between α_{np} and the others.

This is due to:

- 1 - Charge dependent effects.
- 2 - But most of the difference is due to the mass difference of π^\pm and π^0 .

$$m_{\pi^\pm} - m_{\pi^0} = 4.6 \text{ MeV}$$



$\left\{ \begin{array}{l} a_{pp} = -17.1 \pm 0.2 \text{ fm} \\ a_{nn} = -166 \pm 0.6 \text{ fm} \end{array} \right.$ are equal within experimental errors.

But the last measured for a_{nn} is $-18.6 \pm 0.5 \text{ fm}$ from the reaction $\pi^- + d \rightarrow \gamma + n + n$

indicating the charge-dep. of nucl. force.

Remark: At low energies the two nucleons can not come very close together, and the long range of nucl. forces is dominated by the exchange of single pion.

For $T=0 \rightarrow S=1$

$$a_{np} = 5.423 \pm 0.005 \text{ fm} \quad a_{np} > 0 \rightarrow \text{bound state}$$

$$a_{np_{T=1}} \neq a_{np_{T=0}} \rightarrow \text{Isospin-dep. of nucleon forces.}$$

Comparing r_c shows the isospin dependence.

But these values don't give any indication of charge-dependence within experimental errors.

3.9 - The Nucl. Potential:

One Pion Exchange Pt.: (OPEP)

Yukawa's idea of a single one-pion exchange pol. can fit experimental data only for inter-nucleon distances $> 2 \text{ fm}$.

In retrospect this is not a surprise, since $m_\pi \sim 160 \text{ MeV}$
corresponds to a range of 1.6 fm

The experimental data from scatt. experiment show that the sources other than one-pion exchange are responsible.

For example, the values of δ_0 (Page 167) are $\left\{ \begin{array}{l} \text{large} \\ \text{positive} \end{array} \right.$ at low energies

→ indicating that the force is an attractive one -

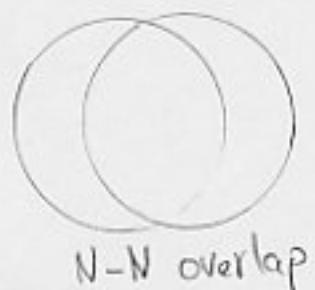
As the energy is increased { above 250 Mev (lab) , 1S_0 becomes negative
{ , 300 Mev (π) , 3S_1 " "

Showing → the forces now are repulsive (hard core evidence)

$$d_{NN} \sim 1 \text{ fm}$$

This strong repulsion is expected:

Three quarks (fermions) in each nucleon must occupy one of the three lowest available state.



In a collision six quarks of two nucleons are no longer two separate groups of



three quarks each (large overlap) .

→ Acc to the Pauli exclusion principle , 3 of the 6-quarks must go $\xrightarrow{\text{to}}$ higher states

A large amount of energy is required to make this transition.

→ This phenomena prevents the nucleons come very close to each other .

Justification of existence of hard core in terms of quarks
does not give a fair quantitative prediction.

OPEP explains the long-range behaviour of nuclear int.

But it has difficulties in, relating

$$g_{NN} \xrightarrow{\text{to}} g_{\pi N} \quad g: \text{strength of int.}$$

$g_{NN} = g_{\pi N}$ (the probability of emitting and absorbing virtual π)

Such probabilities = $f(g_{\pi N})$

In OPEP the $g_{\pi N}$ can not be used directly to calculate g_{NN} .

i.e. There are some other processes effective in th int. other than OPEP.

→ g_{NN} is often treated as a parameter adjusted to fit the N-N data.

One-Boson Exchange Pot. (OBEP):

Nucl. force may be divided into three parts:

i) Long-range part $r > 2\text{ fm}$

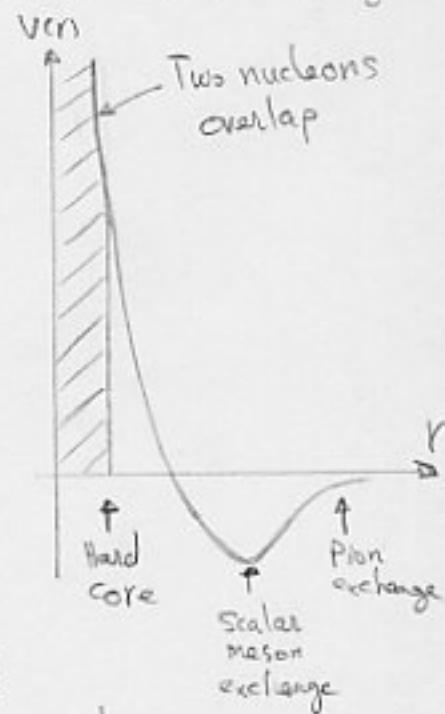
Dominated by one-pion exchange

ii) Intermediate-range part $1\text{ fm} < r < 2\text{ fm}$

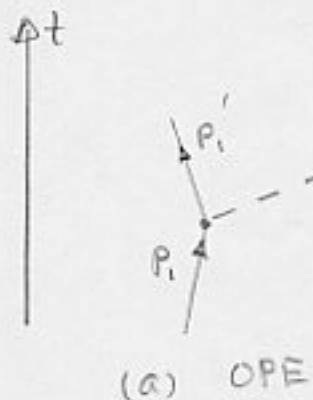
Dominated by two pions and heavy mesons

iii) Hard core int. $r \lesssim 1\text{ fm}$

Dominated by, heavy meson exchange, multi-pion exchange as well as QCD effects.

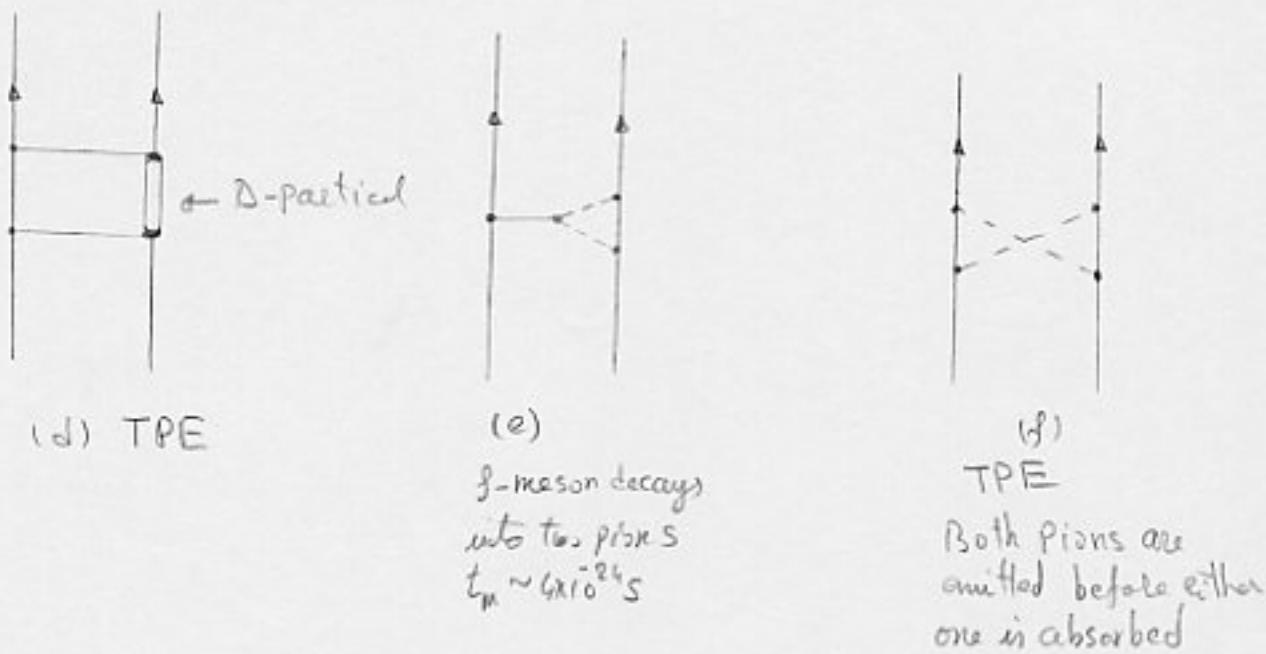


Feynman Diagrams:



Two-pion exchange
are emitted and absorbed one after another

For simplicity (a) is often shown in (b) form.



Nucleon-Nucleon Pot.:

There are two general approaches to constructing a pot. - that
that is the correct for $\left\{ \begin{array}{l} \text{long} \\ \text{intermediate ranges} \\ \text{short} \end{array} \right.$

i) Phenomenological pot.

Generalization of OPE picture $\xrightarrow{\text{to}}$ OBC picture

For simplification only a single boson is considered.
The multimeson exchange is compensated, by letting the
strength for each type of meson exchange to be determined
by fitting from NN scatt. data (as a parameter).

Hard core term is put by hand without any ref. to its source.

Some difficulties:

- a) Lifetimes of many of the mesons are sufficiently short
→ in the exchange process their decay must be taken into account.
- b) In order to fit experimental data, the range of each OBE term and consequently → their mass become adjustable parameters with little or no relation to real masses.
- ii) To make use of our knowledge of hadrons as much as possible and treat phenomenologically only those aspects, mainly short-range ints., of which we have incomplete knowledge.

Nucleon-Nucleon interaction for bound nucleons:

One of the reasons for constructing a correct form of
nuc. pot. is:

{ To investigate the nuclear structure
{ and nuclear reaction

We are faced with many-body system can always be expressed \rightarrow
in terms of two-body int.

i) However effective int. between two-nucleons inside the
nucleus is different with that of bare (under the influence
of all other nucleons).

ii) The second difficulty:

Can a nuclear pot. be specified completely within a
two-body system?

In two-particle system:

$$E = \text{const} , \quad \vec{P} = \text{conserved}$$

$$\frac{P_1^2}{2\mu_1} + \frac{P_2^2}{2\mu_2} = E$$

$$\frac{P^2}{2\mu} = E \rightarrow P_x^2 + P_y^2 + P_z^2 = 2\mu E \quad \sqrt{P_x^2 + P_y^2 + P_z^2} = \sqrt{2\mu E}$$

i.e. The sum of momenta of two nucleons is confined to lie on a spherical shell in momentum space with the radius of $\sqrt{2\mu E}$

E : energy in the center of mass

The two-body int (T-matrix elements) is said to be, on the energy shell, (on shell matrix elements).

In a nucleus with more than two nucleons;

{ Energy conservation applies to the nucleus as a whole
 { Momentum

and the momenta of a pair of nucleons are no longer restricted

$$\text{by } \frac{P_1^2}{2\mu_1} + \frac{P_2^2}{2\mu_2} = E$$

→ off the energy shell (off shell).

Relation quark-quark int.:

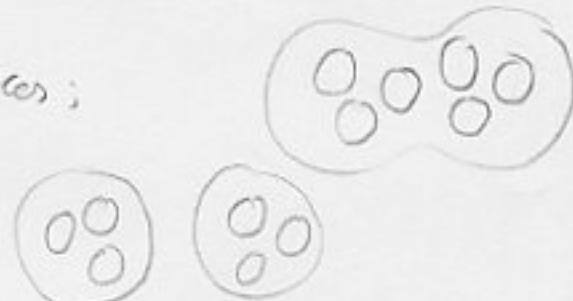
Although, it is accepted the force between nucleons is a facet of q-q strong int., a quantitative connection between nucleon force and q-q int. is still lacking.

The difficulty arises from QCD calculations at the low energies where nucleon phys. operates.

In order to study nucleon int. in quark model, we need a system of at least 6-quarks.

q-q force must give confinement cond. :

{ At close distances ;
otherwise ;



tightly clustered
(bag)

At large distances (compared with q-q distances) ;

The force between these
two bags

\approx
must be
consistent

meson exchange

At intermediate distances :

the force = must be attractive and consistent with
several pions and heavier mesons exchange.

Qualitatively ; we can see how nucle. force may arise from q-q int. by making an analogy with the force between chemical molecules.

Molecules with spherical charge distribution except no force between each other .

But since there is a force between them \rightarrow There must be a residual force present ;

\rightarrow This is Known as the Van der Waals force .

Suppose , a neutral molecule acquires an electric dipole moment \vec{P} (say, as a result of fluctuation in its shape and consequently , its charge distribution)

$$\Phi(r) = \left[\frac{1}{4\pi\epsilon_0} \right] \frac{\hat{r} \cdot \vec{P}}{r^2} = \left[\frac{1}{4\pi\epsilon_0} \right] \frac{PC_0\theta}{r^2} \quad \text{due to dipole}$$

$$E(r,\theta) = -\nabla \Phi = -\left[\frac{1}{4\pi\epsilon_0}\right] \left\{ \frac{\bar{P}}{r^3} - 3\left(\frac{PC_6\theta}{r^4}\right)\bar{r} \right\} \quad \text{due to dipole}$$

This electric field induces a dipole moment \bar{P}' in another molecule a ;

$$\bar{P}' = \chi \bar{E} \quad \chi: \text{polarizability of molecule}$$

The dipole-dipole int. between these two molecules

$$V(r) = -\bar{P}' \cdot \bar{E} = -\left[\frac{1}{4\pi\epsilon_0}\right]^2 \chi (1 + 3C_6\theta) \frac{P^2}{r^6}$$

Interaction energy is always negative, regardless of the orientation of the first dipole.

$$\rightarrow F \sim \frac{1}{r^7}$$

For spherically symmetric molecule;

$$\langle \bar{P} \rangle = 0 \quad (\bar{P} \text{ on average is zero})$$

However, because of fluctuation, the instantaneous value of \bar{P} may be $\neq 0$

$$\text{i.e. } \rightarrow \langle P^2 \rangle \neq 0$$

\longrightarrow An attractive force between two molecules results from this fluctuation.
Analogously:

Electrostatic force in molecules $\xrightarrow{\sim}$ Color force between quarks
 Van der Waals force between molecules $\xrightarrow{\sim}$ Color Van der Waals force between
 - 206 - two nucleons