

# Particle–Number Projected HFB Method with Skyrme Forces

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## Content

- PNP HFB
  - PNP, PAV, VAP
  - Application to Skyrme DFT
- Open problems
  - Shift Invariance
  - Energy Sum Rule
  - Deformation Energy Calculations

# HFB Method

Hamiltonian

$$H = \sum_{n_1 n_2} e_{n_1 n_2} c_{n_1}^\dagger c_{n_2} + \frac{1}{4} \sum_{n_1 n_2 n_3 n_4} \bar{v}_{n_1 n_2 n_3 n_4} c_{n_1}^\dagger c_{n_2}^\dagger c_{n_4} c_{n_3},$$

$$\bar{v}_{n_1 n_2 n_3 n_4} = \langle n_1 n_2 | V | n_3 n_4 - n_4 n_3 \rangle, \quad c_n | - \rangle = 0.$$

$$\begin{pmatrix} \alpha \\ \alpha^\dagger \end{pmatrix} = \begin{pmatrix} U^\dagger & V^\dagger \\ V^T & U^T \end{pmatrix} \begin{pmatrix} c \\ c^\dagger \end{pmatrix}$$

$$\alpha_k | \Phi \rangle = 0, \quad \hat{N} | \Phi \rangle \neq N | \Phi \rangle$$

$$\rho_{n'n} = \frac{\langle \Phi | c_n^\dagger c_{n'} | \Phi \rangle}{\langle \Phi | \Phi \rangle}, \quad \tilde{\rho}_{n'n} = \frac{\langle \Phi | s_n^* c_{\bar{n}} c_{n'} | \Phi \rangle}{\langle \Phi | \Phi \rangle},$$

$$\hat{T} \phi_n(\mathbf{r}, \sigma) = s_n \phi_{\bar{n}}(\mathbf{r}, \sigma), \quad s_n s_n^* = 1, \quad s_{\bar{n}} = -s_n$$

$$E[\rho, \tilde{\rho}] = \frac{\langle \Phi | H | \Phi \rangle}{\langle \Phi | \Phi \rangle}$$

$$\begin{pmatrix} h - \lambda & \tilde{h} \\ \tilde{h} & -h + \lambda \end{pmatrix} \begin{pmatrix} U \\ V \end{pmatrix}_k = E_k \begin{pmatrix} U \\ V \end{pmatrix}_k$$

$$h_{nn'} = \frac{\partial E[\rho, \tilde{\rho}]}{\partial \rho_{n'n}}, \quad \tilde{h}_{nn'} = \frac{\partial E[\rho, \tilde{\rho}]}{\partial \tilde{\rho}_{n'n}}$$

Bogoliubov  
Transformation

HFB Equations

# Particle Number Projected HFB Method

Projection After Variation (PAV)

HFB Energy

$$E_{HFB} = \frac{\langle \Phi | H | \Phi \rangle}{\langle \Phi | \Phi \rangle}, \quad \hat{N} | \Phi \rangle \neq N | \Phi \rangle$$

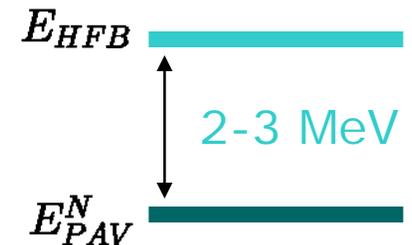
Particle Number  
Projection

$$P^N = \frac{1}{2\pi} \int d\phi e^{i\phi(\hat{N}-N)}$$

$$|\Psi^N\rangle = P^N |\Phi\rangle, \quad \hat{N} |\Psi^N\rangle = N |\Psi^N\rangle$$

PAV Energy

$$E_{PAV}^N = \frac{\langle \Psi^N | H | \Psi^N \rangle}{\langle \Psi^N | \Psi^N \rangle} = \frac{\langle \Phi | H P^N | \Phi \rangle}{\langle \Phi | P^N | \Phi \rangle}$$



# Particle Number Projected HFB Method

## Variation After Projection (VAP)

PNP HFB Energy

$$E^N[\rho, \tilde{\rho}] = \frac{\langle \Phi | H P^N | \Phi \rangle}{\langle \Phi | P^N | \Phi \rangle} = \frac{\int d\phi \langle \Phi | H e^{i\phi(\hat{N}-N)} | \Phi \rangle}{\int d\phi \langle \Phi | e^{i\phi(\hat{N}-N)} | \Phi \rangle}$$

$$\begin{pmatrix} \varepsilon^N + \Gamma^N + \Lambda^N & \Delta^N \\ -(\Delta^N)^* & -(\varepsilon^N + \Gamma^N + \Lambda^N)^* \end{pmatrix} \begin{pmatrix} U \\ V \end{pmatrix}_k = E_k^N \begin{pmatrix} U \\ V \end{pmatrix}_k$$

$$\varepsilon^N = \frac{1}{2} \int d\phi y(\phi) (Y(\phi) \text{Tr}[e\rho(\phi)] + [1 - 2ie^{-i\phi} \sin \phi \rho(\phi)] eC(\phi)) + h.c.$$

$$\Gamma^N = \frac{1}{4} \int d\phi y(\phi) (Y(\phi) \text{Tr}[\Gamma(\phi)\rho(\phi)] + 2[1 - 2ie^{-i\phi} \sin \phi \rho(\phi)] \Gamma(\phi)C(\phi)) + h.c.$$

$$\Lambda^N = -\frac{1}{4} \int d\phi y(\phi) (Y(\phi) \text{Tr}[\Delta(\phi)\bar{\kappa}^*(\phi)] - 4ie^{-i\phi} \sin \phi C(\phi)\Delta(\phi)\bar{\kappa}^*(\phi)) + h.c.$$

$$\Gamma_{n_1 n_3}(\phi) = \sum_{n_2 n_4} \bar{v}_{n_1 n_2 n_3 n_4} \rho_{n_4 n_2}(\phi),$$

$$\Delta^N = \frac{1}{2} \int d\phi y(\phi) e^{-2i\phi} C(\phi) \Delta(\phi) - (\dots)^T,$$

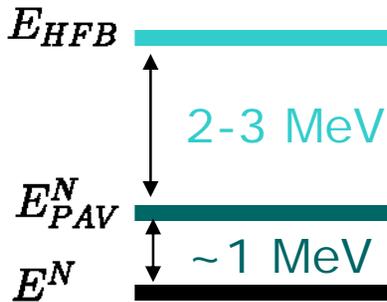
$$\Delta_{n_1 n_2}(\phi) = \frac{1}{2} \sum_{n_3 n_4} \bar{v}_{n_1 n_2 n_3 n_4} \kappa_{n_3 n_4}(\phi), \quad \bar{\Delta}_{n_3 n_4}^*(\phi) = \frac{1}{2} \sum_{n_1 n_2} \bar{\kappa}_{n_1 n_2}^*(\phi) \bar{v}_{n_1 n_2 n_3 n_4},$$

$$\rho(\phi) = C(\phi)\rho \quad \kappa(\phi) = (\phi)\kappa = \kappa C^T(\phi)$$

$$\bar{\kappa}(\phi) = e^{2i\phi} \kappa C^*(\phi) = e^{2i\phi} C^\dagger(\phi) \kappa$$

$$C(\phi) = e^{2i\phi} (1 + \rho(e^{2i\phi} - 1))^{-1},$$

$$x(\phi) = \frac{1}{2\pi} \frac{e^{-i\phi N} \det(e^{i\phi} I)}{\sqrt{\det C(\phi)}}, \quad y(\phi) = \frac{x(\phi)}{\int d\phi' x(\phi')}, \quad \int d\phi y(\phi) = 1,$$



# Skyrme HFB Method

Skyrme HFB  
Functional

$$\frac{\langle \Phi | H | \Phi \rangle}{\langle \Phi | \Phi \rangle} \iff E[\rho, \tilde{\rho}] = \int d\mathbf{r} [H(\mathbf{r}) + \tilde{H}(\mathbf{r})]$$

$H(\mathbf{r})$  and  $\tilde{H}(\mathbf{r})$  are normal and pairing energy densities, respectively, expressed in terms of particle and pairing local densities and currents

One-Body  
Density Matrices

$$\rho(\mathbf{r}\sigma, \mathbf{r}'\sigma') = \sum_{nn'} \rho_{nn'} \check{\psi}_n^*(\mathbf{r}', \sigma') \check{\psi}_n(\mathbf{r}, \sigma)$$

$$\tilde{\rho}(\mathbf{r}\sigma, \mathbf{r}'\sigma') = \sum_{nn'} \tilde{\rho}_{nn'} \check{\psi}_n^*(\mathbf{r}', \sigma') \check{\psi}_n(\mathbf{r}, \sigma)$$

Skyrme HFB  
Equations

$$h_{nn'} = \frac{\partial E[\rho, \tilde{\rho}]}{\partial \rho_{n'n}}, \quad \tilde{h}_{nn'} = \frac{\partial E[\rho, \tilde{\rho}]}{\partial \tilde{\rho}_{n'n}}$$

$$\begin{pmatrix} h - \lambda & \tilde{h} \\ \tilde{h} & -h + \lambda \end{pmatrix} \begin{pmatrix} U \\ V \end{pmatrix}_k = E_k \begin{pmatrix} U \\ V \end{pmatrix}_k$$

# PNP Skyrme HFB Method

## DFT Complication

HFB Method

$$E[\rho, \tilde{\rho}] \iff \langle \Phi | \hat{H} | \Phi \rangle$$

Densities  $\rho, \tilde{\rho}$  associated with a single state  $|\Phi\rangle$

PNP HFB Method

$$E^N[\rho, \tilde{\rho}] \iff \frac{\langle \Phi | H P^N | \Phi \rangle}{\langle \Phi | P^N | \Phi \rangle}$$

Requires knowledge of off-diagonal expectation values which are not automatically given by DFT

$$\langle \Phi(0) | \hat{H} | \Phi(\varphi) \rangle, \quad |\Phi(\varphi)\rangle = e^{i\varphi(\hat{N}-N)} |\Phi\rangle$$

'Mixed Densities'  
Prescription

Obviously certain extensions are necessary and they are not unique. Among various possibilities, the so-called 'mixed density' recipe is most frequently used in projection and other GCM calculations.

$$\rho(\varphi) = \langle \Phi(0) | \hat{\rho} | \Phi(\varphi) \rangle, \quad \tilde{\rho}(\varphi) = \langle \Phi(0) | \hat{\tilde{\rho}} | \Phi(\varphi) \rangle$$

# PNP Skyrme HFB Method

Energy Functional under 'Mixed densities' prescription

$$E^N[\rho, \tilde{\rho}] = \frac{\int d\varphi e^{-i\varphi N} \mathcal{I}(\varphi) E[\rho(\varphi), \tilde{\rho}(\varphi)]}{\int d\varphi e^{-i\varphi N} \mathcal{I}(\varphi)} = \int d\varphi y(\varphi) E[\rho(\varphi), \tilde{\rho}(\varphi)]$$

$$y(\varphi) = \frac{e^{-i\varphi N} \mathcal{I}(\varphi)}{\int d\varphi e^{-i\varphi N} \mathcal{I}(\varphi)}, \quad \int d\varphi y(\varphi) = 1, \quad \mathcal{I}(\varphi) = \langle \Phi(0) | \Phi(\varphi) \rangle$$

$$\rho(\varphi) = \langle \Phi(0) | \hat{\rho} | \Phi(\varphi) \rangle, \quad \tilde{\rho}(\varphi) = \langle \Phi(0) | \hat{\tilde{\rho}} | \Phi(\varphi) \rangle$$

$$E[\rho(\varphi), \tilde{\rho}(\varphi)] = \int d\mathbf{r} \mathcal{H}(\mathbf{r}, \phi)$$

$$\bullet \rho(\mathbf{r}), \tilde{\rho}(\mathbf{r}) \quad \Longrightarrow \quad \rho(\mathbf{r}, \phi), \tilde{\rho}(\mathbf{r}, \phi)$$

$$\bullet \tau(\mathbf{r}), \tilde{\tau}(\mathbf{r}) \quad \Longrightarrow \quad \tau(\mathbf{r}, \phi), \tilde{\tau}(\mathbf{r}, \phi)$$

$$\bullet \mathbf{J}_{ij}(\mathbf{r}), \tilde{\mathbf{J}}_{ij}(\mathbf{r}) \quad \Longrightarrow \quad \mathbf{J}_{ij}(\mathbf{r}, \phi), \tilde{\mathbf{J}}_{ij}(\mathbf{r}, \phi)$$

Energy  
Functional

Canonical Representation

Unprojected density  $\rho_n = v_n^2, \quad \tilde{\rho}_n = u_n v_n \quad \mathcal{I}(\varphi) = \prod_n (u_n^2 + v_n^2 e^{2i\varphi}).$

'Mixed' density  $\rho_n(\varphi) = \frac{v_n^2 e^{2i\varphi}}{u_n^2 + v_n^2 e^{2i\varphi}}, \quad \tilde{\rho}_n(\varphi) = \frac{u_n v_n e^{i\varphi}}{u_n^2 + v_n^2 e^{2i\varphi}}$

Projected density  $\rho_n^N = \int d\varphi y(\varphi) \rho_n(\varphi), \quad \tilde{\rho}_n^N = \int d\varphi y(\varphi) \tilde{\rho}_n(\varphi)$

# PNP Skyrme HFB Method

VAP under 'Mixed densities' prescription

$$\begin{pmatrix} h^N & \tilde{h}^N \\ \tilde{h}^N & -h^N \end{pmatrix} \begin{pmatrix} U \\ V \end{pmatrix}_k = E_k^N \begin{pmatrix} U \\ V \end{pmatrix}_k$$

$$h^N = \int d\phi y(\phi) Y(\phi) E(\phi) + \int d\phi y(\phi) e^{-2i\phi} C(\phi) h(\phi) C(\phi) - \left[ \int d\phi y(\phi) 2ie^{-i\phi} \sin(\phi) \tilde{\rho}(\phi) \tilde{h}(\phi) C(\phi) + h.c. \right],$$

$$\tilde{h}^N = \int d\phi y(\phi) e^{-i\phi} (\tilde{h}(\phi) C(\phi) + (\dots)^T)$$

$$h(\phi) = \frac{\partial E[\rho(\varphi), \tilde{\rho}(\varphi)]}{\partial \rho(\phi)}, \quad \tilde{h}(\phi) = \frac{\partial E[\rho(\varphi), \tilde{\rho}(\varphi)]}{\partial \tilde{\rho}(\phi)}$$

$$\rho_n = v_n^2, \quad \tilde{\rho}_n = u_n v_n, \quad \rho_n(\varphi) = \frac{v_n^2 e^{2i\varphi}}{u_n^2 + v_n^2 e^{2i\varphi}}, \quad \tilde{\rho}_n(\varphi) = \frac{u_n v_n e^{i\varphi}}{u_n^2 + v_n^2 e^{2i\varphi}}$$

$$C_\mu(\phi) = \frac{e^{2i\phi}}{u_\mu^2 + e^{2i\phi} v_\mu^2}, \quad y(\phi) = \frac{e^{-iN\phi} \prod_{\mu>0} (u_\mu^2 + e^{2i\phi} v_\mu^2)}{\sum_{l=0}^{L-1} e^{-iN\phi_l} \prod_{\mu>0} (u_\mu^2 + e^{2i\phi_l} v_\mu^2)}$$

$$Y_\mu(\phi) = ie^{-i\phi} \sin \phi C_\mu(\phi) - \sum_{l'=0}^{L-1} y(\phi_{l'}) ie^{-i\phi_{l'}} \sin \phi_{l'} C_\mu(\phi_{l'})$$

$$P^N = \frac{1}{2\pi} \int d\phi e^{i\phi(\hat{N}-N)} \implies P^N = \frac{1}{L} \sum_{l=1}^L e^{i\phi_l(\hat{N}-N)}, \quad \phi_l = \frac{2\pi}{L}(l-1)$$

PNP HFB  
Equations

Canonical  
Representation

Grid Points

# PNP Skyrme HFB Method

## Problems: Stability

Slow (even unstable) procedure

$$\begin{pmatrix} h^N & \tilde{h}^N \\ \tilde{h}^N & -h^N \end{pmatrix} \begin{pmatrix} U \\ V \end{pmatrix}_k = E_k^N \begin{pmatrix} U \\ V \end{pmatrix}_k$$

$$\text{Tr}\rho = \bar{N}, \quad \rho = V^*V^T$$

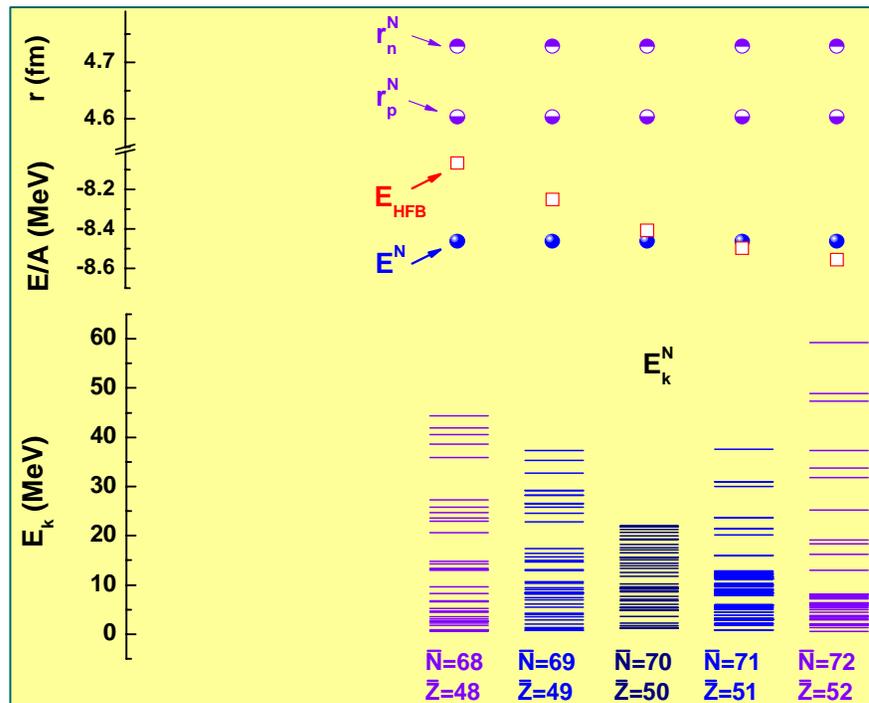
Stable procedure

$$\begin{pmatrix} h^N - \mu & \tilde{h}^N \\ \tilde{h}^N & -(h^N - \mu) \end{pmatrix} \begin{pmatrix} U \\ V \end{pmatrix}_k = E_k^N \begin{pmatrix} U \\ V \end{pmatrix}_k$$

$$\text{Tr}\rho = \bar{N}, \quad \rho = V^*V^T$$

$\mu$  is zero when the PNP solution is found

$$\text{Tr}\rho^N = N, \quad \rho^N = \int d\phi y(\phi) C(\phi) \rho$$

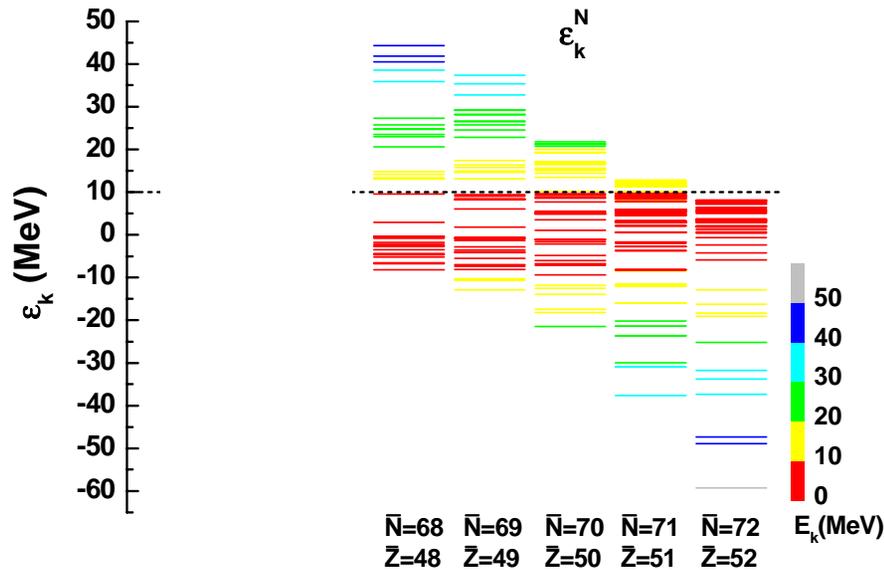
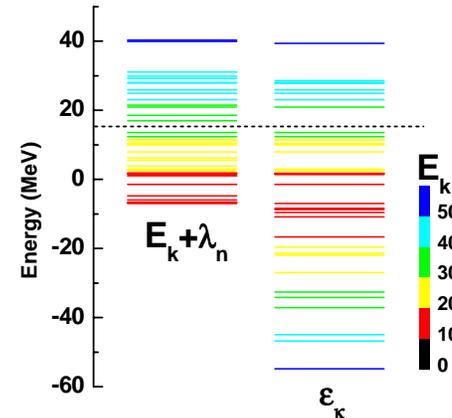


# PNP Skyrme HFB Method

Problems: Cut-off procedure for Delta pairing forces

$$\bar{e}_n = (1 - 2P_n)E_n + \lambda$$

$$\bar{\Delta}_n = 2E_n \sqrt{P_n(1 - P_n)}$$



Unprojected HFB  $h$  and  $\tilde{h}$   
appear in the PNP scheme  
at gauge angle  $\varphi = 0$

$$\tilde{E}_n = \begin{pmatrix} U \\ V \end{pmatrix}_n^\dagger \begin{pmatrix} h - \lambda & \tilde{h} \\ \tilde{h} & -h + \lambda \end{pmatrix} \begin{pmatrix} U \\ V \end{pmatrix}_n$$

# PNP Skyrme HFB Method

## Problems: Pairing Strength

In the standard Skyrme HFB method the pairing strength  $V_0$  is chosen in such a way that the HFB value of the average neutron gap  $\tilde{\Delta}_n$  at given cut-off energy  $\epsilon_{\text{cut}}$  reproduces the experimental value 1.245 MeV for the nucleus  $^{120}\text{Sn}$ .

The average neutron gap  $\tilde{\Delta}_n$  is no more defined within PNP HFB method. Therefore, the above procedure for adjusting the pairing strength is no more applicable.

A strict way of adjusting the pairing strength should be obtained by calculating mass differences and comparing with available experimental data.

We adjust the pairing strength to the total energy of some nucleus already calculated in PLN HFB approximation. This is rather crude approximation we are using just to analyze the quality of the PNP HFB treatment.

$$\tilde{H}(\mathbf{r}) = \frac{1}{2}V_0 \left[ 1 - V_1 \left( \frac{\rho}{\rho_0} \right)^\gamma \right] \sum_q \tilde{\rho}_q^2$$

$$V_1 = \begin{cases} 0 & \text{volume pairing} \\ 1 & \text{surface pairing} \end{cases}$$

$$\gamma = 1, \quad \rho_0 = 0.16 \text{ fm}^{-3}$$

HFB  
Method

PNP HFB  
Method

# PNP Skyrme HFB Method

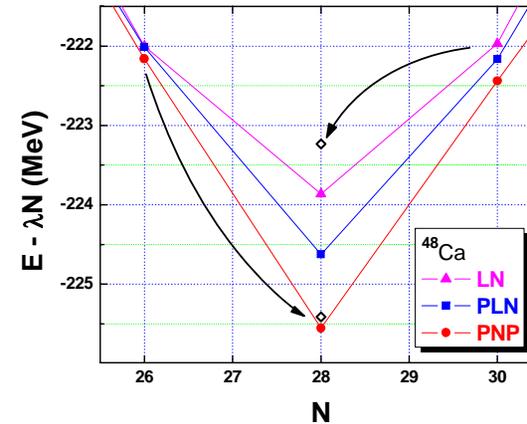
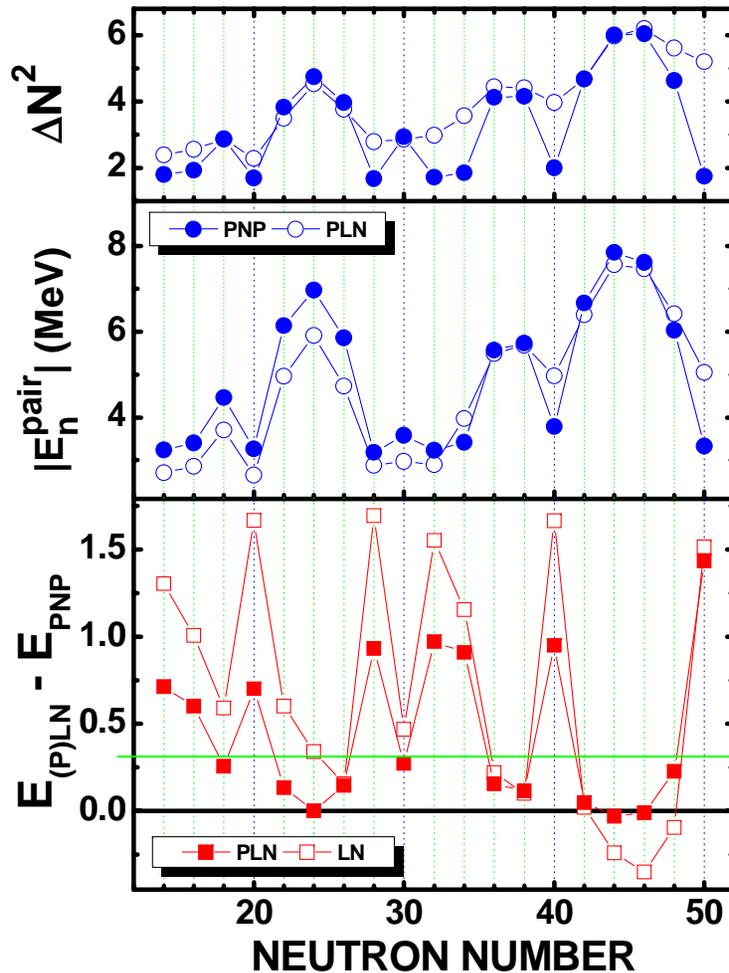
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## Ca Chain Calculations

- SLY4 + mixed delta pairing forces
- HFB within 20 major HO shells
- Complete Ca chain
- Comparison:
  - HFB+LN results (LN)
  - PAV HFB+LN results (PLN)
  - VAP PNP HFB results (PNP)
- PLN pairing strength fitted to  $\Delta_n$  @  $^{120}\text{Sn}$
- PNP pairing strength to PLN  $E_{\text{tot}}$  @  $^{44}\text{Ca}$
  
- **With L=9 gauge-angle points  
the code is just 9 times slower**

# PNP Skyrme HFB Method

## Some Results



- LN method should be avoided
  - One should use PLN instead
- PLN is a good approximation for open shell nuclei
  - total energy differences are less than 250 KeV
- PLN is wrong for closed shell nuclei
  - total energy differences could be more than 1 MeV
- One should try to correct PLN by projecting from neighboring nuclei

# PNP within DFT

## Well Known Singularity

$$\rho_n(\varphi) = \frac{v_n^2 e^{2i\varphi}}{u_n^2 + v_n^2 e^{2i\varphi}}, \quad \tilde{\rho}_n(\varphi) = \frac{u_n v_n e^{i\varphi}}{u_n^2 + v_n^2 e^{2i\varphi}}$$

$$u_n^2 = v_n^2 = \frac{1}{2}, \quad \varphi = \frac{\pi}{2}$$

We have found in mass table calculations that among all 5818 nuclei only about 50 of them have a neutron state with occupation 0.5 and other 48 nuclei with such a proton state. Therefore, we have about 100 questionable nuclei among 5818 which makes less than 2 percents. The situations however is much more serious when performing constrained HFB calculations.

# PNP HFB Method

## Shift invariance and Energy sum rule

$$|\Psi_N\rangle \equiv P^N |\Phi\rangle = \frac{1}{2\pi} \int_0^{2\pi} d\varphi e^{i\varphi(\hat{N}-N)} |\Phi\rangle$$

$$(P^N)^\dagger = P^N, \quad (P^N)^2 = P^N, \quad \sum_N P^N = 1$$

$$|\tilde{\Psi}_N\rangle = \frac{\Psi_N}{\sqrt{\langle\Psi_N|\Psi_N\rangle}}, \quad |\tilde{\Phi}\rangle = \frac{\Phi}{\sqrt{\langle\Phi|\Phi\rangle}}, \quad |\tilde{\Phi}\rangle = \sum_N b_N |\tilde{\Psi}_N\rangle,$$

$$b_N^2 = \frac{\langle\Phi|P_N|\Phi\rangle}{\langle\Phi|\Phi\rangle}, \quad \sum_N |b_N|^2 = 1$$

$$E_{HFB} = \frac{\langle\Phi|\hat{H}|\Phi\rangle}{\langle\Phi|\Phi\rangle}, \quad E^N = \frac{\langle\Psi_N|\hat{H}|\Psi_N\rangle}{\langle\Psi_N|\Psi_N\rangle}$$

Exact  
Relations

Shift  
Invariance

$$e^{i\eta(\hat{N}-N)} |\Psi_N\rangle = |\Psi_N\rangle, \quad \hat{N} |\Psi_N\rangle = N |\Psi_N\rangle \quad E^N = \frac{\langle\Phi|\hat{H}|\Psi_N\rangle}{\langle\Phi|\Psi_N\rangle} = \frac{\langle\Phi|\hat{H}\hat{S}_\eta|\Psi_N\rangle}{\langle\Phi|\hat{S}_\eta|\Psi_N\rangle}$$

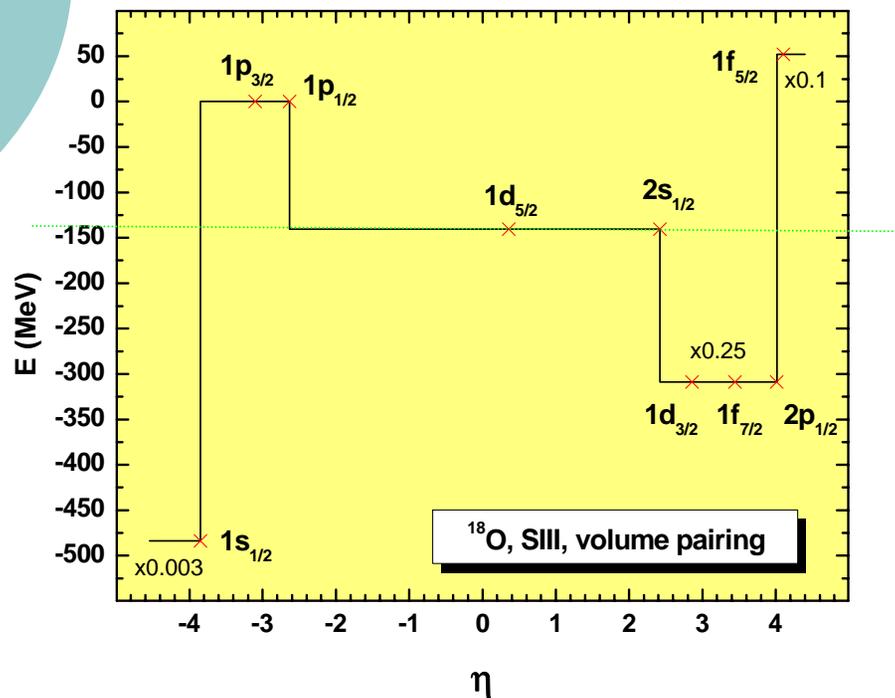
Energy  
Sum Rule

$$E_{HFB} = \sum_N |b_N|^2 E^N$$

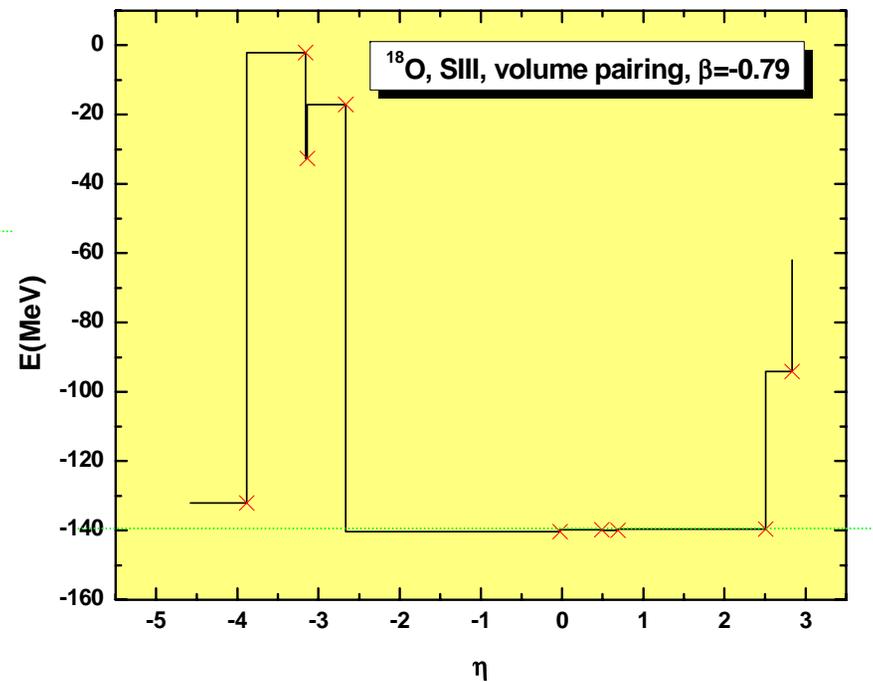
# PNP within DFT

## Broken Shift Invariance

### Spherical Nuclei



### Deformed Nuclei



# PNP within DFT

PNP Energy

$$E^N[\rho, \tilde{\rho}] = \frac{\int d\varphi e^{-\varphi N} \mathcal{I}(\varphi) E[\rho(\varphi), \tilde{\rho}(\varphi)]}{\int d\varphi e^{-\varphi N} \mathcal{I}(\varphi)}$$

$$\mathcal{I}(\varphi) = \langle \Phi(0) | \Phi(\varphi) \rangle = \prod_n (u_n^2 + v_n^2 e^{2i\varphi})$$

$$\rho_n(\varphi) = \frac{v_n^2 e^{2i\varphi}}{u_n^2 + v_n^2 e^{2i\varphi}}, \quad \tilde{\rho}_n(\varphi) = \frac{u_n v_n e^{i\varphi}}{u_n^2 + v_n^2 e^{2i\varphi}}$$

$$z = e^{i\varphi}, \quad d\varphi = \frac{dz}{iz}, \quad C_1(|z|=1) \quad \rho_n(z) = \frac{v_n^2 z^2}{u_n^2 + v_n^2 z^2}, \quad \tilde{\rho}_n(z) = \frac{u_n v_n z}{u_n^2 + v_n^2 z^2}$$

$$\mathcal{N}_N \equiv \int d\varphi e^{-\varphi N} \mathcal{I}(\varphi) E[\rho(\varphi), \tilde{\rho}(\varphi)] = -i \oint \frac{dz}{z^{N+1}} \prod_n (u_n^2 + v_n^2 z^2) E[\rho(z), \tilde{\rho}(z)]$$

$$\mathcal{D}_N \equiv \int d\varphi e^{-\varphi N} \mathcal{I}(\varphi) = \oint \frac{dz}{z^{N+1}} \prod_n (u_n^2 + v_n^2 z^2)$$

$$\oint_C dz f(z) = 2\pi i \sum_k \text{Rez}[f(z_k)] \quad z_k = \pm i |u_k/v_k|$$

$$\mathcal{N}_N = 2\pi i \sum_{|z_k| \leq 1} \text{Rez} \left[ \frac{1}{z_k^{N+1}} \prod_n (u_n^2 + v_n^2 z_k^2) E[z_k] \right]$$

$$\mathcal{D}_N = 2\pi i \text{Rez} \left[ \frac{1}{z_0^{N+1}} \prod_n (u_n^2 + v_n^2 z_0^2) \right]$$

New  
complex  
variable

Cauchy's  
residue  
theorem

# PNP within DFT

## PNP Energy

$$E_N = E_N[z_0] + \Delta E_N$$

$$z_0 = 0, \quad z_k = \pm i |u_k/v_k|$$

$$E_N[z_0] = \frac{\text{Rez} \left[ \frac{1}{z_0^{N+1}} \prod_n (u_n^2 + v_n^2 z_0^2) E[z_0] \right]}{\text{Rez} \left[ \frac{1}{z_0^{N+1}} \prod_n (u_n^2 + v_n^2 z_0^2) \right]} \quad \Delta E_N = \frac{2\pi i}{\mathcal{D}_N} \sum_{0 < |z_k| \leq 1} \text{Rez} \left[ \prod_n (u_n^2 + v_n^2 z_k^2) \frac{E[z_k]}{z_k^{N+1}} \right]$$

## PNP Energy – explicit pole dependence

- $E[\rho, \tilde{\rho}]$  leads to  $f_a(z)$  without pole at  $z_k$
- $a = d + p$ ,  $d$  power of  $\rho$ ,  $p$  power of  $\tilde{\rho}$
- $\nu_k$  is the degeneracy of the  $k$ -th canonical state
  - for deformed nuclei always  $\nu_k = 1$
  - for spherical nuclei  $\nu_k = (2j + 1)/2$
  - $\nu_k = 1$  for  $j = 1/2$ ,  $\nu_k = 2$  for  $j = 3/2$  ...

$$E_N = E_N[z_0] + \Delta E_N$$

$$\Delta E_N = \sum_{0 < |z_k| \leq 1} \sum_a \text{Rez} \left[ \frac{(u_k^2 + v_k^2 z_k^2)^{\nu_k} f_a(z_k)}{z_k^{N+1} (u_k^2 + v_k^2 z_k^2)^a} \right]$$

## Shifted PNP Energy

$$\underbrace{e^{\eta(\hat{N}-N)}}_{\hat{S}_\eta} |\Psi_N\rangle$$

$$C_1(|z| = 1) \Rightarrow C_\eta(|z| = e^{-\eta})$$

$$E_N(\eta) = E_N[z_0] + \Delta E_N(\eta)$$

$$\Delta E_N(\eta) = \sum_{|z_k| \leq e^{-\eta}} \sum_a \text{Rez} \left[ \frac{(u_k^2 + v_k^2 z_k^2)^{\nu_k} f_a(z_k)}{z_k^{N+1} (u_k^2 + v_k^2 z_k^2)^a} \right]$$

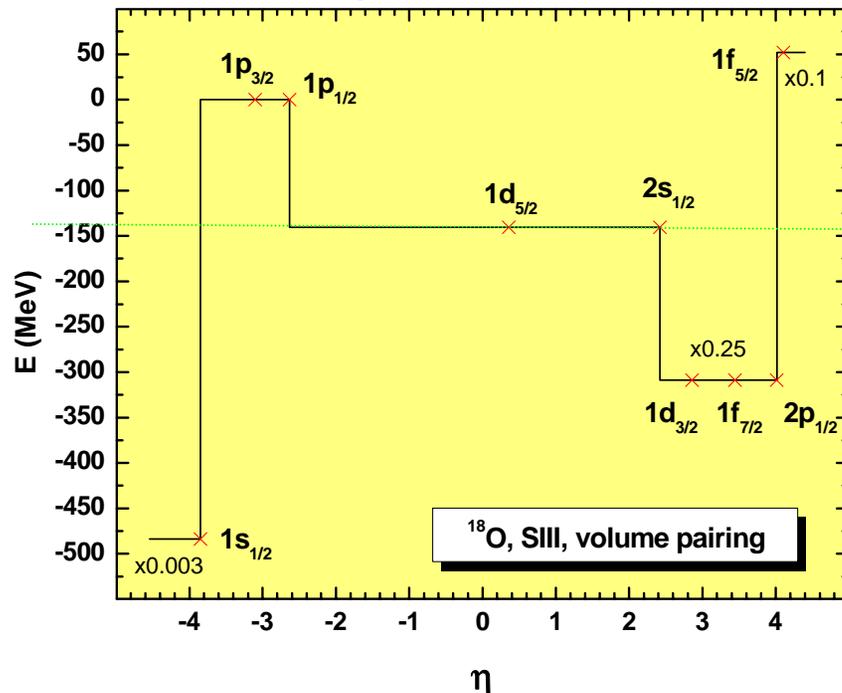
# PNP within DFT

## Local Shift Invariance

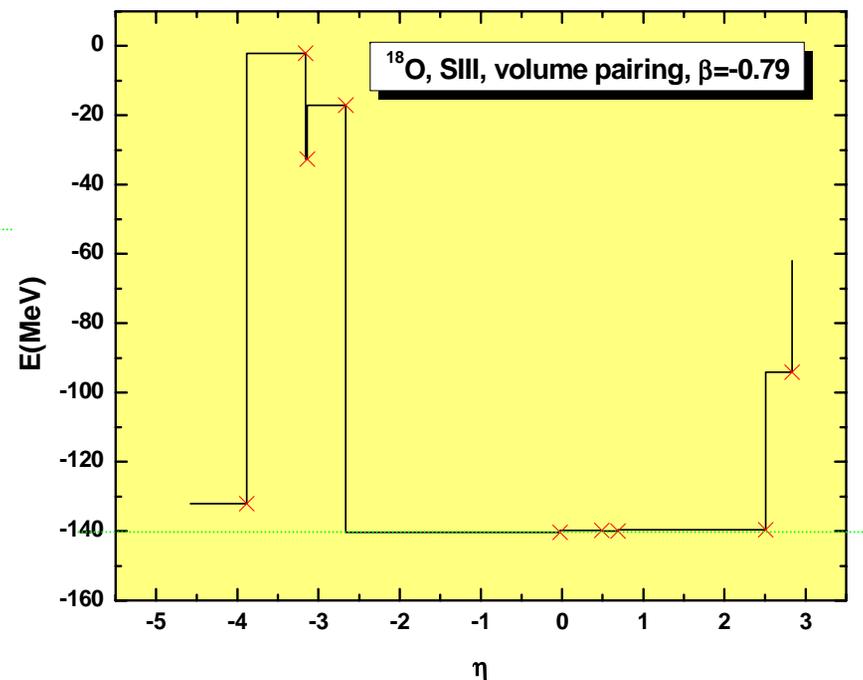
$$E_N(\eta) = E_N[z_0] + \Delta E_N(\eta) \quad z_0 = 0, \quad z_k = \pm i |u_k/v_k| \Leftrightarrow u_k^2 = v_k^2 = \frac{1}{2}, \quad \varphi = \frac{\pi}{2}$$

$$E_N[z_0] = \frac{\text{Rez} \left[ \frac{1}{z_0^{N+1}} \prod_n (u_n^2 + v_n^2 z_0^2) E[z_0] \right]}{\text{Rez} \left[ \frac{1}{z_0^{N+1}} \prod_n (u_n^2 + v_n^2 z_0^2) \right]} \quad \Delta E_N(\eta) = \sum_{|z_k| \leq e^{-\eta}} \sum_a \text{Rez} \left[ \frac{(u_k^2 + v_k^2 z_k^2)^{\nu_k} f_a(z_k)}{z_k^{N+1} (u_k^2 + v_k^2 z_k^2)^a} \right]$$

### Spherical Nuclei



### Deformed Nuclei



# PNP within DFT

## Exact versus Approximate DFT

In the ideal case when the functional  $E^N[\rho, \tilde{\rho}]$  is exactly equivalent to an expectation value of a given Hamiltonian  $H$  all residues from the poles  $z_k > z_0$  are strictly zero and the energy is defined only by the residue of the zero pole  $z_0 = 0$ .

*Kinetic energy term* as well as all linear terms in the energy functional correspond to a power  $a = 1$ . Then all residues of  $z_k > z_0$  will be zero since always one has  $\nu_k \geq 1$ .

*Coulomb energy* is a quadratic term,  $a = 2$ .

- The residue contribution of  $z_k > z_0$  is zero assuming one treats the exchange term exactly – the residue from the direct coulomb term exactly cancels with the residue of the exchange term.
- If one uses Slater approximation for the exchange term such a cancelation does not exist anymore.

### *pp and ph contributions*

ph channel:

$$\frac{t_0}{4}(1-x_0)\rho_n^2 \rightarrow \frac{t_0}{4}(1-x_0)\frac{v_k^4 z_k^4}{z_k^{N+1}(u_k^2 + v_k^2 z_k^2)^2}$$

pp-channel:

$$\frac{t_0}{4}(1-x_0)\tilde{\rho}_n^2 \rightarrow \frac{t_0}{4}(1-x_0)\frac{u_k^2 v_k^2 z_k^2}{z_k^{N+1}(u_k^2 + v_k^2 z_k^2)^2}$$

their sum cancels the pole contribution:

$$\frac{t_0}{4}(1-x_0)(\rho_n^2 + \tilde{\rho}_n^2) \rightarrow \frac{t_0}{4}(1-x_0)\frac{v_k^2 z_k^2}{z_k^{N+1}(u_k^2 + v_k^2 z_k^2)}$$

In the case of Skyrme forces for which contact pairing forces are used instead the original Skyrme force, one will see nonzero contributions from the poles  $z_k > z_0$  coming from both ph as well as pp terms.

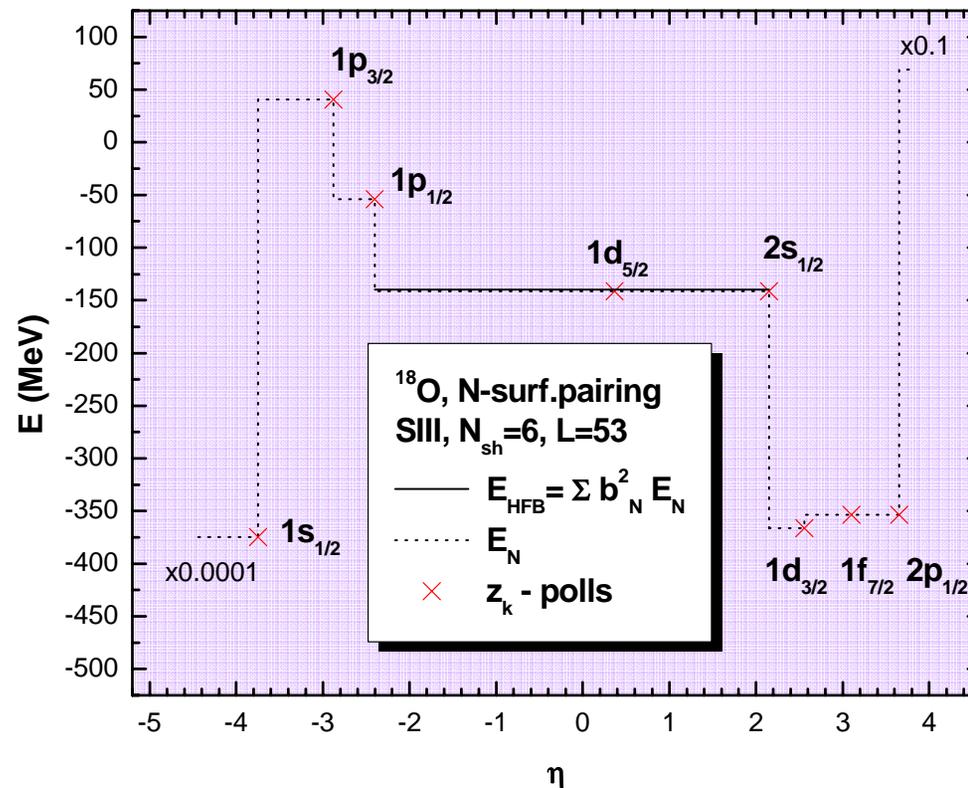
# PNP within DFT

Energy Sum Rule

$$\tilde{E}(\eta) = \sum_N |b_N|^2 E^N(\eta) \neq E_{HFB}$$

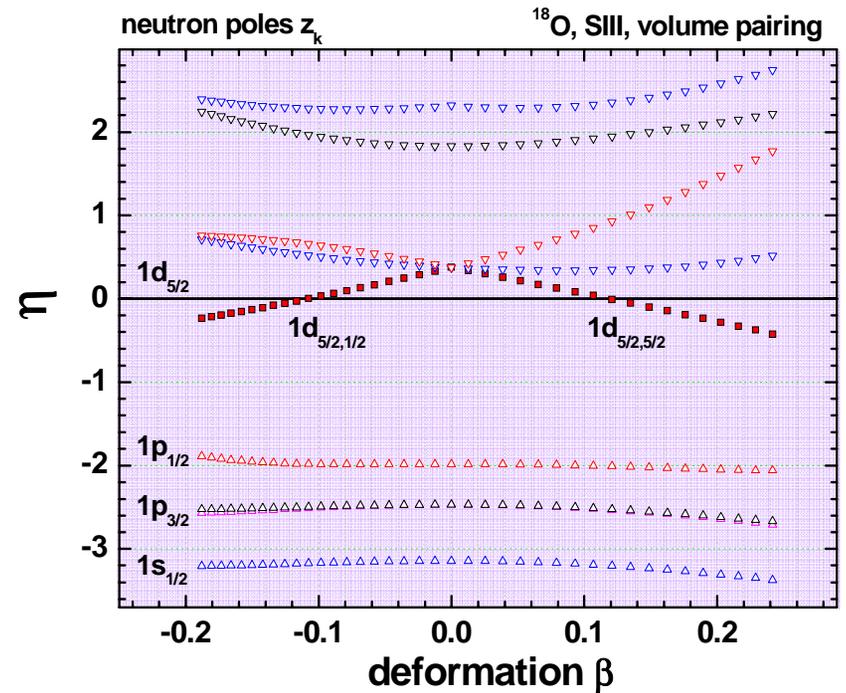
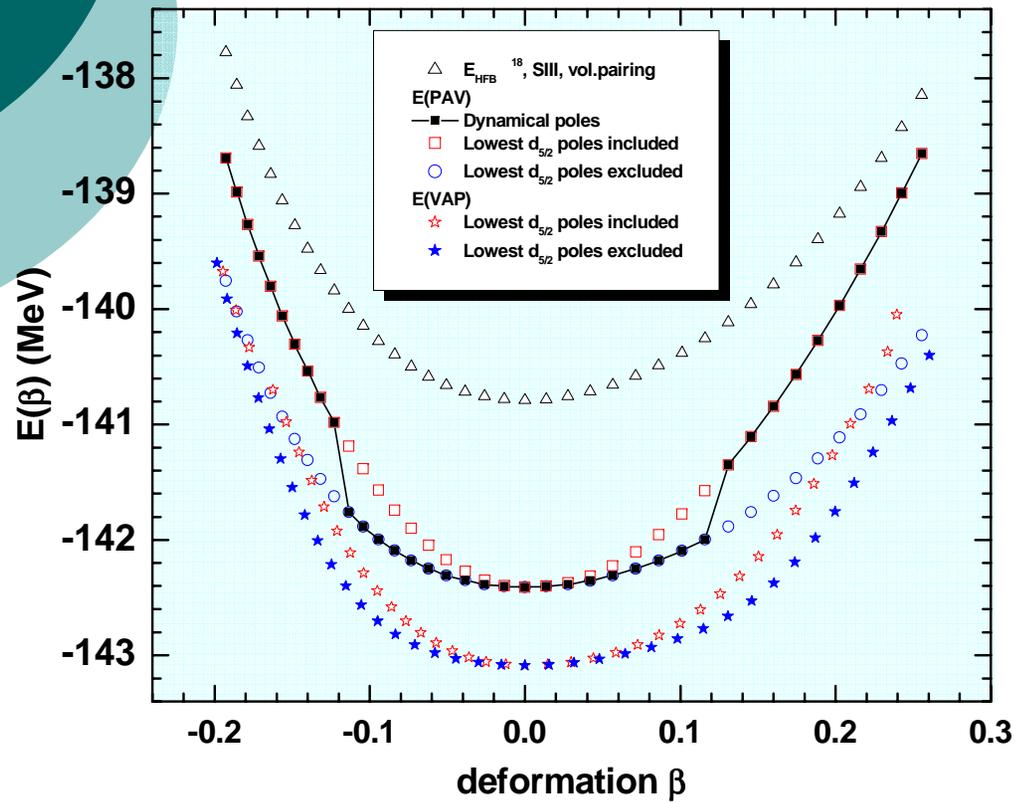
$$\tilde{E}(\eta = 0) = E_{HFB}$$

$$\tilde{E}(\eta = 0) = \sum_N |b_N|^2 E^N = \int d\varphi \sum_N e^{-i\varphi N} \mathcal{I}(\varphi) E[\rho(\varphi), \tilde{\rho}(\varphi)] = E[\rho(0), \tilde{\rho}(0)] = E_{HFB}$$



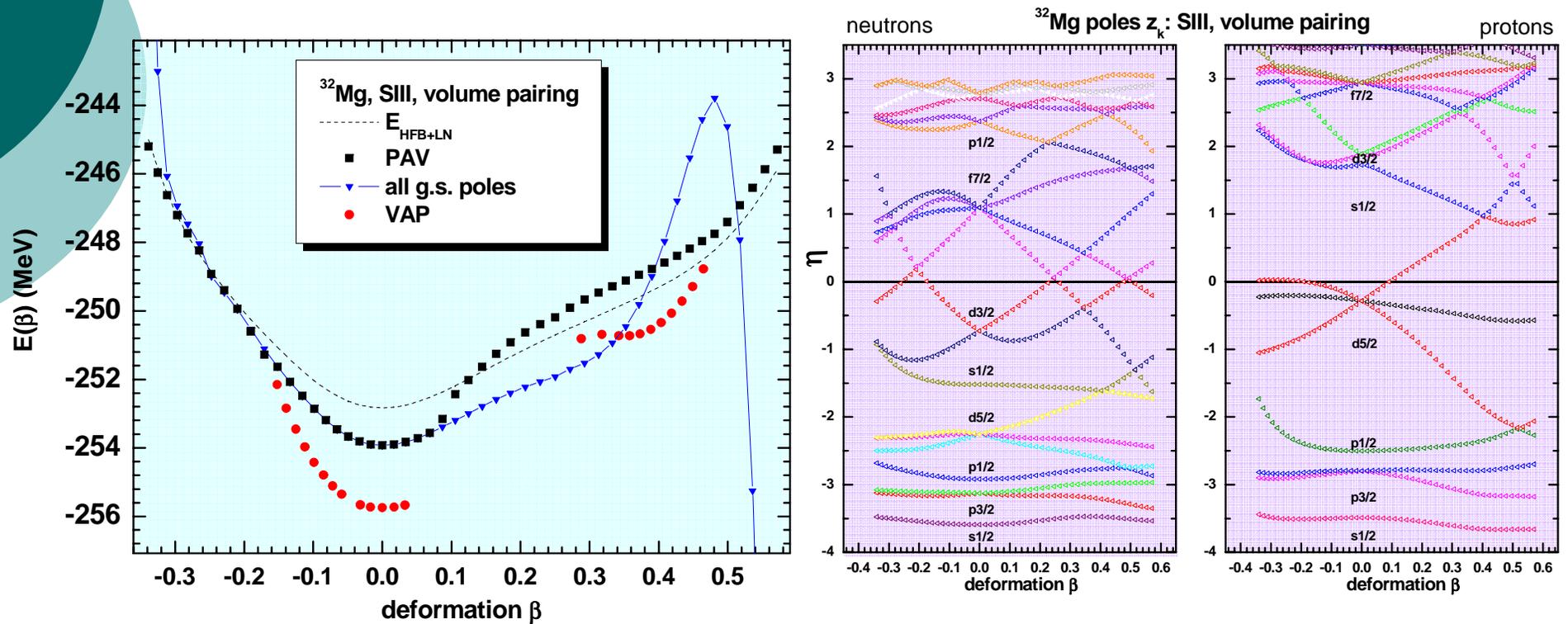
# PNP within DFT

## Deformation Energy Calculations



# PNP within DFT

## Deformation Energy Calculations



# PNP Skyrme HFB Method

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## Conclusions

When no singularity exists on the unit circle

- ⇒ LN method should be avoided – One should use PLN instead
- ⇒ PLN is a good for open shell nuclei – Error is less than 250 KeV
- ⇒ PLN is wrong for closed shell nuclei – Error could be more than 1 MeV
- ⇒ One should try to correct PLN – Projecting from neighboring nuclei

## PNP versus DFT

- ⇒ All singularities cancel if EDF is exact
- For an approximate functional:
- ⇒ Shift Invariance is broken – Locally it is satisfied
  - ⇒ Energy Sum rule is not satisfied – Satisfied on the unit circle only
  - ⇒ Density dependence is not analytical – Valid even for Gogny forces
  - ⇒ Instability in VAP – No solution at the moment