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**TWO-BODY TRANSITION DENSITIES IN NCSM:
MICROSCOPIC DESCRIPTION OF
CORE-SWELLING IN ${}^6\text{He}$**

SHELL MODEL CODE DEVELOPMENT

- ▶ Antoine - on-the-fly shell model code developed at Strasbourg.
 - ▶ Works in m-scheme.
 - ▶ Utilizes proton and neutron subspaces. $D \gg D_p, D_n$
- ▶ pAntoine - technical code developments at Chalmers (H.T. Johansson & B. D. Carlsson).
- ▶ Very large model spaces can be handled.
- ▶ Very efficient (runs on a single node).

${}^6\text{Li}, N_{\text{max}}=22$

Non-zero elements:

$N_{\text{non-zero}}=5 \times 10^{14} \sim 6 \text{ PB data}$

$2.5 \times 10^9 \text{ multi/sec/machine}$

NEW TRANSITION DENSITY CODE NEEDED

- Handle large model spaces
- Efficient, utilizes smaller subspace dimension
- Higher-rank operators

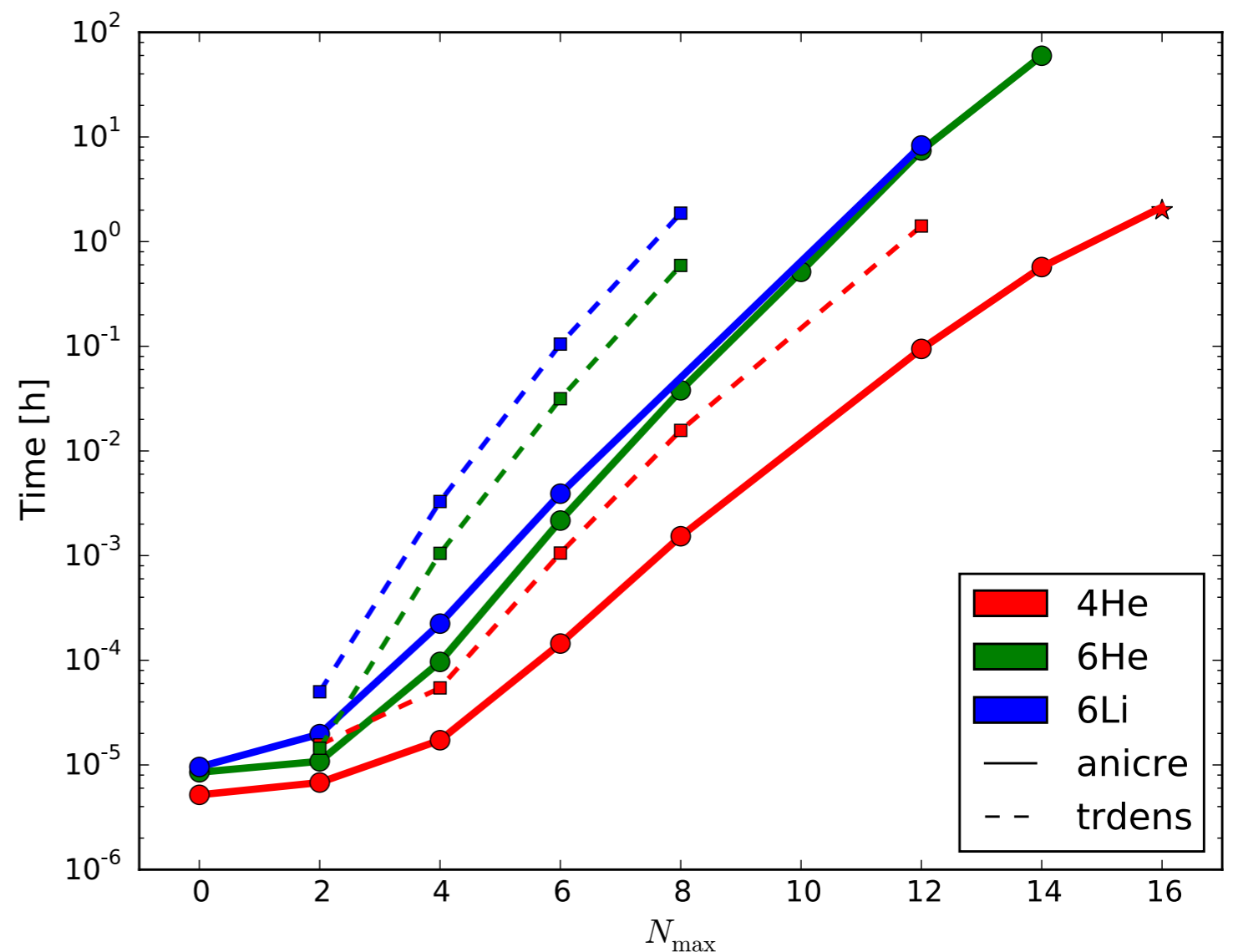
N_{max}	Dimension of model space: ${}^6\text{Li}$
16	7.9×10^8
18	2.7×10^9
20	8.6×10^9
22	2.5×10^{10}

ANICRE

$$(\xi_f J_f || T_\lambda^2 || \xi_i J_i) = \hat{\lambda}^{-1} \sum (a, b || T_\lambda^2 || c, d) (\xi_f J_f || [[a_a^\dagger a_b^\dagger][a_c a_d]]_\lambda || \xi_i J_i)$$

- Ongoing code development project
 - ▶ For the moment limited to single-threaded calculations
 - ▶ One- and two-body transition densities
- Efficient use of hash tables and indexing
- Efficient JM-ordering and spin-couplings

Two-body transition densities

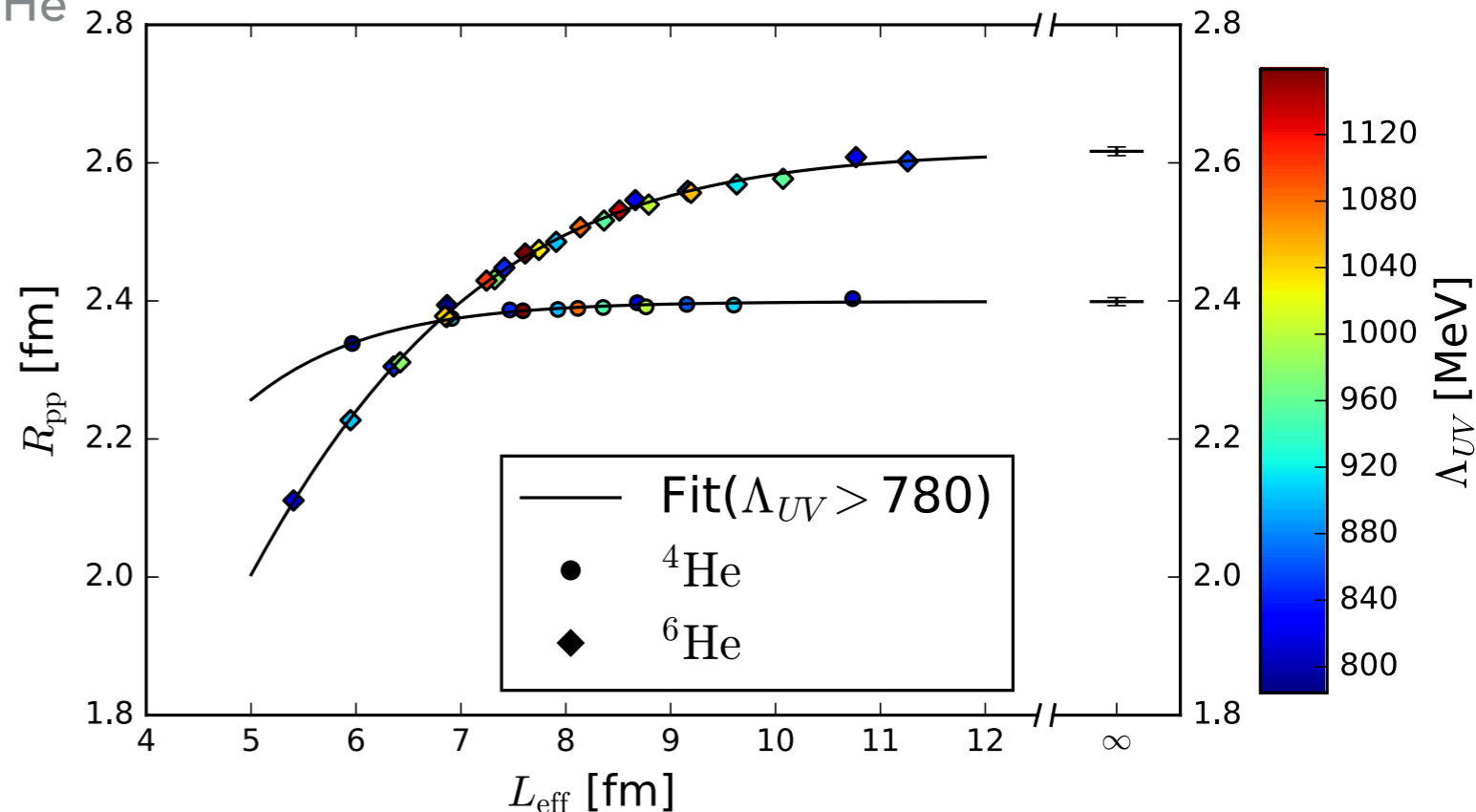
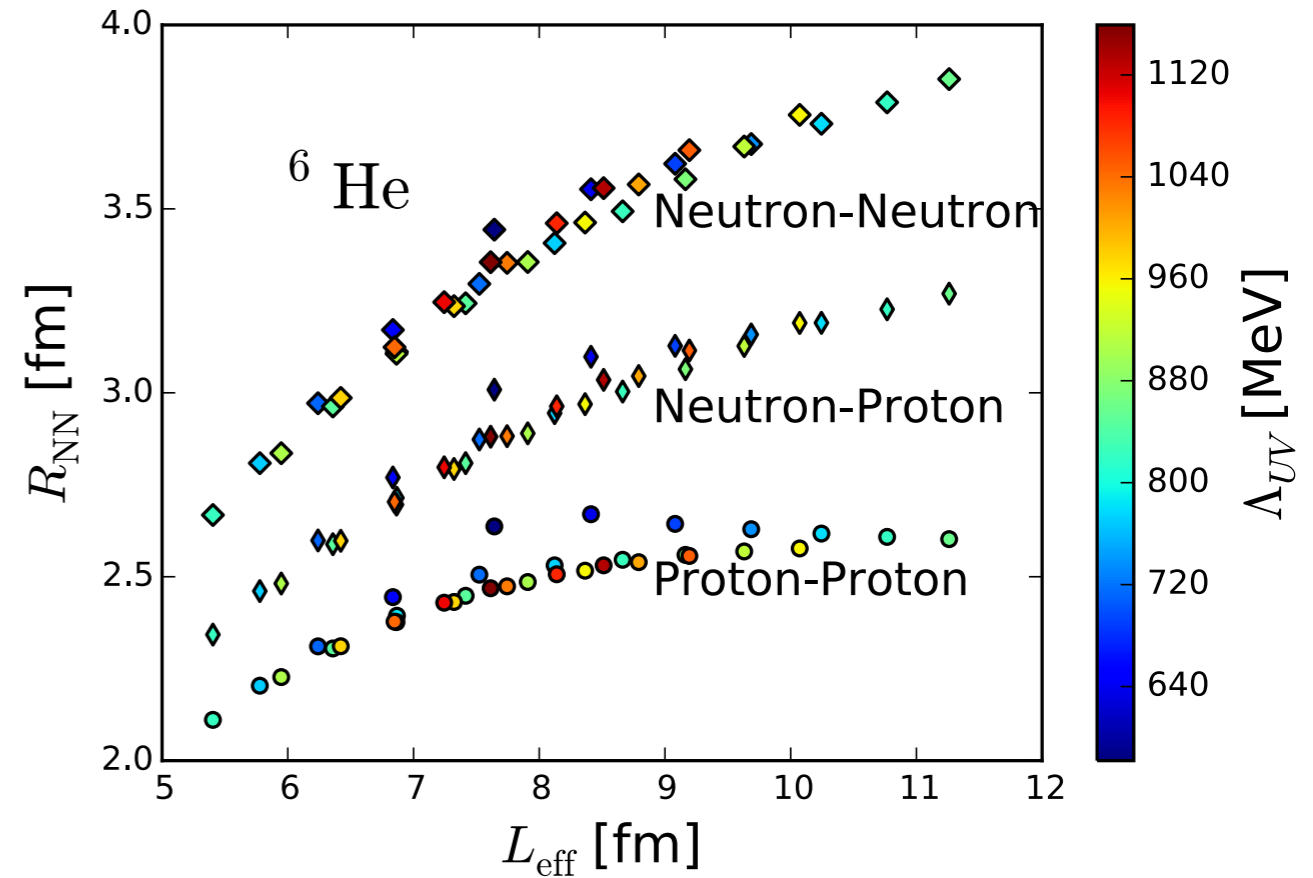


NUCLEON-NUCLEON DISTANCES

- Computed from two-body transition densities
- Test application of transition density code

CORE-SWELLING IN ${}^6\text{He}$

- Radius difference between the α -core in ${}^6\text{He}$ and the free α -particle.
- Important input in cluster models.
- A measure of core-swelling: increased p-p distance
- In our calculations: The core-swelling effect $\sim 9\%$ (Computed with NNLOopt)



NUCLEON-NUCLEON DISTANCES

Further information:
Come and see my poster.
Thank you!

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