

2N and 3N systems in three dimensional formalism - a compilation.

Few-nucleon fusion reactions in 3D formalism.

Kacper Topolnicki

Institute of Physics Jagiellonian University.

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- Classical QM - always start with the Schrödinger equation:

$$i\hbar\partial_t |\psi(t)\rangle = \hat{H} |\psi(t)\rangle$$

or

$$|\psi(t)\rangle = \exp(-i\hat{H}(t - t_0)) |\psi(t_0)\rangle.$$

- Hands on approach - the proton and the neutron are two states of the spin $\frac{1}{2}$ isospin $\frac{1}{2}$ nucleon. We use three dimensional states:

$$| \mathbf{k}_1 \mathbf{k}_2 \rangle \otimes |\uparrow\downarrow\rangle^{\text{isospin}} \otimes |\uparrow\uparrow\rangle^{\text{spin}}, | \mathbf{Kp} \rangle \otimes |\uparrow\downarrow\rangle^{\text{isospin}} \otimes |\uparrow\downarrow\rangle^{\text{spin}}$$

or

$$| \mathbf{k}_1 \mathbf{k}_2 \mathbf{k}_3 \rangle \otimes |\uparrow\downarrow\downarrow\rangle^{\text{isospin}} \otimes |\uparrow\downarrow\downarrow\rangle^{\text{spin}}, | \mathbf{Kpq} \rangle \otimes |\uparrow\downarrow\downarrow\rangle^{\text{isospin}} \otimes |\downarrow\uparrow\downarrow\rangle^{\text{spin}}.$$



- Classical non-relativistic QM, but calculations get quite complicated.
- A lot of pieces of the puzzle must fit together.
 - ▶ Analytical calculations are practically impossible. This is especially true for $3N$ systems.
 - ▶ Numerical code implementing analytical results must not contain errors. Typically thousands of lines of FORTRAN code. Each +, – must be in the proper place. This process must be automated.
 - ▶ Our solution - extensive use of symbolic programming within the Mathematica system.
- Exponential increase in efficiency. What used to take months now takes 15 min and a click of a button.



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- Chose a potential ($2N$, $3N$) and a problem ($2N$ bound state, $3N$ bound state, transition operator calculation).
- Calculate the hard part (analytical expressions and FORTRAN code) using Mathematica.
- Construct a FORTRAN implementation of linear operators (resulting directly from the Schrödinger equation with some additional constraints on the states of the system under consideration) from automatically generated code.
- Use Krylov subspace methods to reduce the size of the operators (this is especially needed for large $3N$ systems and requires the use of powerful computing clusters - JUQUEEN in FZJ).
- Solve the reduced (say 40×40 dimensional) linear (eigen) equation using classical methods. We use LAPACK or Mathematica linear solvers.
- Compare results with classical PWD calculations.

WHY 3D CALCULATIONS?



- The choice of partial wave channels is to some degree arbitrary. If we want more precise predictions, we take a larger number of channels but it is not always true that PWD states are organized in such a way that taking a large number of channels produces matrix elements of operators organized by their magnitude. 3D calculations utilize all partial waves.
- Using Krylov subspace methods and 3D representation automatically organizes matrix elements according to their size. This gives hope for better precision.
- Why not! Rare opportunity to gain direct insight into the nuclear processes.



- ϕ_1, ϕ_2 describe the 2N bound state.
- Linear operator (acting in the space of scalar functions ϕ).
- Expressed in terms of integrals but an explicit matrix representation is also available.

$$\begin{aligned}
 (\check{K}^d(E_d)\phi)_q(|\mathbf{p}|) = & \\
 \frac{1}{E_d - \frac{\mathbf{p}^2}{m}} \int d^3\mathbf{p}' \sum_{j=1}^6 v_j^{00}(\mathbf{p}, \mathbf{p}') \sum_{k''=1}^2 & \\
 \left(\sum_k (A^d(\mathbf{p}))_{qk}^{-1} B_{kjk''}^d(\mathbf{p}, \mathbf{p}') \right) \phi_{k''}(|\mathbf{p}'|) &
 \end{aligned}$$

Time independent Schrödinger equation

$\rightarrow (\check{K}^d(E)\phi)_q(|\mathbf{p}|) = \lambda\phi_q(|\mathbf{p}|)$ - find E such that $\lambda \approx 1$ ($E \approx E_d$).



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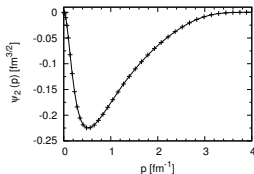
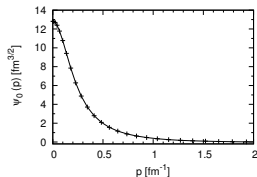
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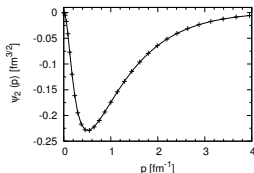
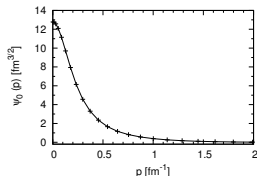
- $\phi_q(|\mathbf{p}|)$ - $2 \cdot 40 = 80$ dimensional vector.
- $\check{K}^d(E_d)$ - 80×80 matrix.
- $\left(\sum_k (A^d(\mathbf{p}))_{qk}^{-1} B_{kjk''}^d(\mathbf{p}, \mathbf{p}') \right)$ - calculated in Mathematica.
- $v_j^{00}(\mathbf{p}, \mathbf{p}')$ - 2N potential (decomposed).
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The s-wave (left) and d-wave (right) component of the deuteron wave function - chiral NNLO potential.



The same but for the Bonn B potential.

Two-nucleon systems in three dimensions
Phys. Rev. C 81 3 (2010)

Golak, J. and Glöckle, W. and Skibiński R. and Witała H. and Rospędzik D. and Topolnicki, K. and Fachruddin, I. and Elster, Ch. and Nogga, A.

TRANSITION OPERATOR (T MATRIX)



- Linear operators in the space of scalar functions t .
- The transition operator is fully determined by the set of 6 t functions.
- An explicit matrix representation is available.

$$\begin{aligned}
 (\check{B}t)_k^{\{\gamma\}} (\{E\}, |\mathbf{p}'|, \{|\mathbf{p}|\}, x') = \\
 \int_0^{+\infty} d|\mathbf{p}''| \int_{-1}^1 dx'' \int_0^{2\pi} d\phi'' \sum_{j=1}^6 \sum_{j'=1}^6 \frac{|\mathbf{p}''|^2}{\{E\} - \frac{|\mathbf{p}''|^2}{m} + i\epsilon} \\
 v_j^{\{\gamma\}} (|\mathbf{p}'|, |\mathbf{p}''|, \sqrt{1-x'^2}\sqrt{1-x''^2} \cos \phi'' + x'x'') \\
 \mathcal{B}_{kjj'} (|\mathbf{p}'|, \{|\mathbf{p}|\}, x', |\mathbf{p}''|, x'', \phi'') \\
 t_{j'}^{\{\gamma\}} (\{E\}, |\mathbf{p}''|, \{|\mathbf{p}|\}, x'')
 \end{aligned}$$

$$\begin{aligned}
 (\check{F}(|\mathbf{p}''|)t)_k^{\{\gamma\}} (\{E\}, |\mathbf{p}'|, \{|\mathbf{p}|\}, x') = \\
 m \int_{-1}^1 dx'' \int_0^{2\pi} d\phi'' \sum_{j=1}^6 \sum_{j'=1}^6 \\
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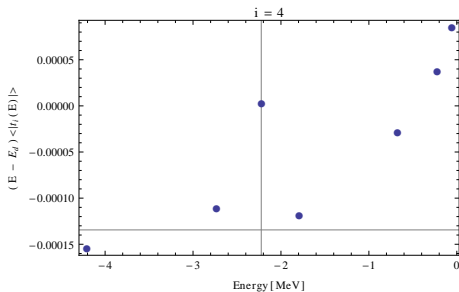
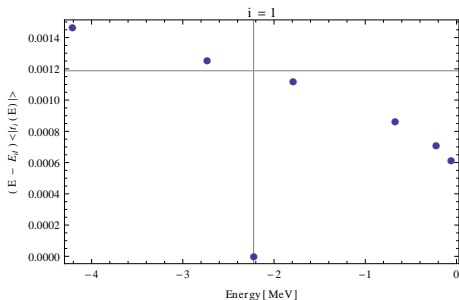


- $t_k^{\{\gamma\}}(\{E\}, |\mathbf{p}'|, \{|\mathbf{p}|\}, \mathbf{x}')$ - for each γ , E , $|\mathbf{p}'|$ - $6 \cdot 40 \cdot 40 = 9600$ dimensional vector.
- $(\check{B}t)_k^{\{\gamma\}}(\{E\}, |\mathbf{p}'|, \{|\mathbf{p}|\}, \mathbf{x}')$ - $4 \cdot 40 \cdot < \text{number of energies} >$ 9600×9600 dimensional independent problems.
- Cases with $E > 0$ and $E < 0$ need to be considered separately.
 - ▶ $E < 0$ - singularity around the deuteron binding energy, we need to substitute $\check{V} | \phi_d \rangle \frac{1}{E - E_b} \langle \phi_d | \check{V}$. All expressions simple to calculate with our tools.
 - ▶ $E > 0$ - problem with singularity in \check{B} (we introduce \check{f}).

TRANSITION OPERATOR $E < 0$



A slice through $t_i^{\{\gamma\}}(\{E\}, |\mathbf{p}'|, \{\mathbf{p}\}, \mathbf{x}')$ (the cross represents deuteron substitution):

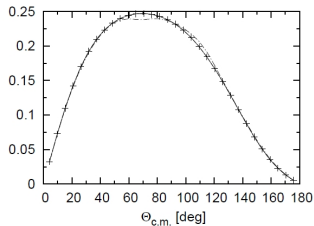
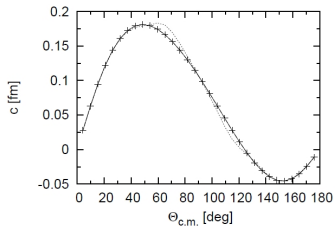
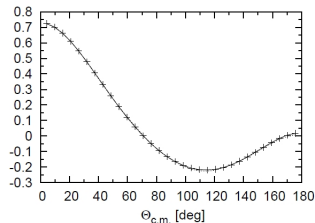
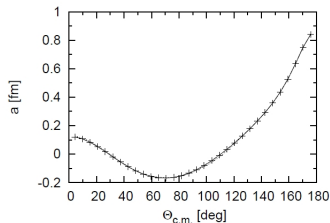


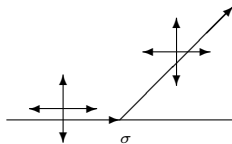
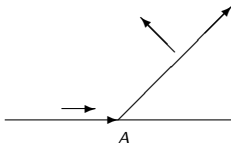
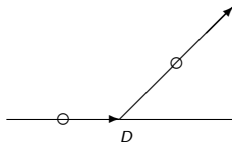
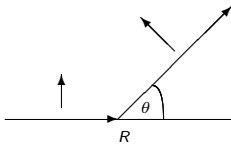
T OPERATOR ON SHELL - $E = 300$ MeV



*Different Methods for the
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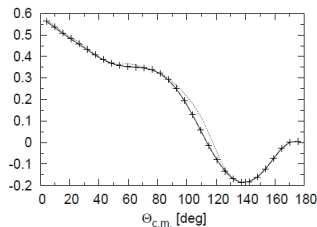
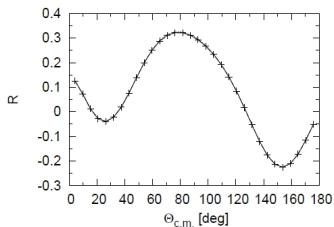
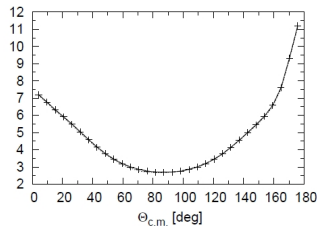
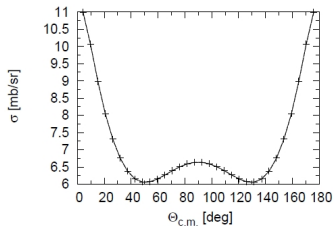


The Quantum Mechanical Few-Body Problem. Walter Glöckle (Springer-Verlag)

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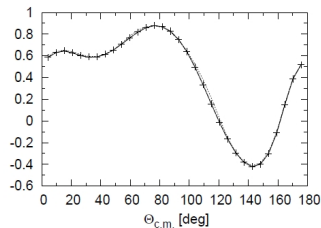
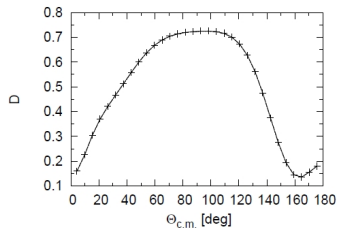
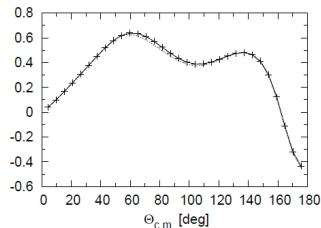
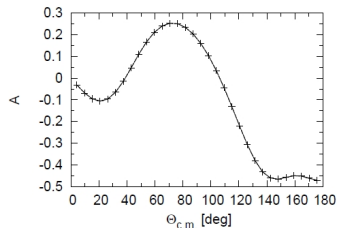
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3N BOUND STATE

- Linear operator in the space of β scalar functions.
- The 3N bound state is determined by the 8 β functions.
- Currently no explicit matrix representation is available - it is constructed from integrals.

$$\begin{aligned}
 & (\check{P}_{1223}^{\text{scalar}} \beta)_{t' T'}^{(k)} (|\mathbf{p}'|, |\mathbf{q}'|, \hat{\mathbf{p}}' \cdot \hat{\mathbf{q}}') = \\
 & \sum_{tT} \sum_{i=1}^8 \beta_{tT}^{(i)} (|\mathbf{P}^{2312}(\mathbf{p}', \mathbf{q}')|, |\mathbf{Q}^{2312}(\mathbf{p}', \mathbf{q}')|, \\
 & \hat{\mathbf{P}}^{2312}(\mathbf{p}', \mathbf{q}') \cdot \hat{\mathbf{Q}}^{2312}(\mathbf{p}', \mathbf{q}')) \\
 & C_{t' T' k; t T i}^{1223}(\mathbf{p}', \mathbf{q}')
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Schrödinger equation

$\rightarrow (\check{G}_0(E) (\check{V}^{\text{scalar}} + \check{V}^{(1)\text{scalar}}) (\check{\mathbf{1}} + \check{P}_{1223}^{\text{scalar}} + \check{P}_{1323}^{\text{scalar}})) \beta = \lambda \beta$ solve and find E such that $\lambda \approx 1$.

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 & (C^{-1}L)_{kji}(\mathbf{p}, \mathbf{q}; \mathbf{p}, \mathbf{p}'; \mathbf{p}', \mathbf{q})
 \end{aligned}$$

$$\begin{aligned}
 & (\check{V}^{(1)\text{scalar}} \beta)_{t'T'}^{(i)}(|\mathbf{p}|, |\mathbf{q}|, \hat{\mathbf{p}} \cdot \hat{\mathbf{q}}) = \\
 & \sum_{i=1}^8 \int d^3 \mathbf{p} \int d^3 \mathbf{q} \sum_{t'} \beta_{t'T'}^{(i)}(|\mathbf{p}|, |\mathbf{q}|, \hat{\mathbf{p}} \cdot \hat{\mathbf{q}}) \\
 & \sum_{r=1}^8 C_{kr}^{-1}(\mathbf{p}, \mathbf{q}) E_{ri}^{ttT}(\mathbf{p}\mathbf{q}; \mathbf{p}\mathbf{q}\mathbf{p}\mathbf{q}; \mathbf{p}\mathbf{q}) \equiv \\
 & \sum_{i=1}^8 \int d^3 \mathbf{p} \int d^3 \mathbf{q} \sum_{t'} \beta_{t'T'}^{(i)}(|\mathbf{p}|, |\mathbf{q}|, \hat{\mathbf{p}} \cdot \hat{\mathbf{q}}) (C^{-1}E)_{ri}^{ttT}(\mathbf{p}\mathbf{q}; \mathbf{p}\mathbf{q}\mathbf{p}\mathbf{q}; \mathbf{p}\mathbf{q})
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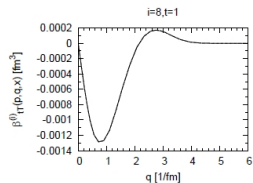
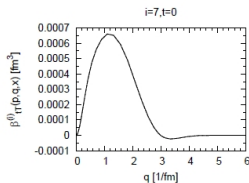
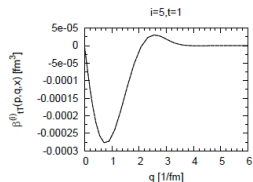
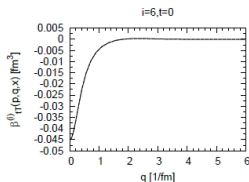
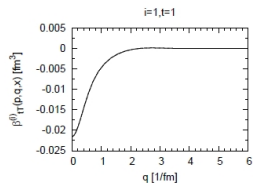
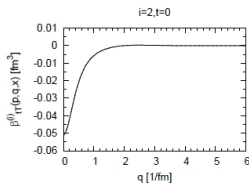


- $\beta_{t'T'}^{(k)}(|\mathbf{p}'|, |\mathbf{q}'|, \hat{\mathbf{p}}' \cdot \hat{\mathbf{q}}')$ - $3 \cdot 8 \cdot 40 \cdot 40 \cdot 40 = 1536000$ dimensional vectors.
- $\check{V}^{\text{scalar}}, \check{V}^{(1)\text{scalar}}, \check{P}_{1223}^{\text{scalar}}, \check{P}_{1323}^{\text{scalar}}$ - 1536000×1536000 dimensional operators.
- Large computational resources necessary - JUQUEEN in FZJ JUELICH.

*A Three-Dimensional Treatment of the Three-Nucleon Bound State
Few-Body Systems 2012*

Golak, J. and Topolnicki, K. and Skibiński, R. and Glöckle, W. and Kamada, H. and Nogga, A.

	PWD	3D
λ	1.0	0.99976
$\langle E_{\text{kin}} \rangle$	33.448	33.412
$\langle E_{\text{pot}}^{2N} \rangle$	-41.329	-41.273
$\langle E_{\text{pot}}^{3N} \rangle$	-0.765	- 0.770
total energy	-8.646	-8.631





- We developed a new framework for dealing with 2N and 3N systems.
- The results for the deuteron, t-matrix and 3N bound state have been verified and published in:
 - ▶ *Phys. Rev. C 81 3 (2010)*
 - ▶ *Few-Body Systems 2012 (53 237-252)*
 - ▶ *Few-Body Systems 2012 (1-20)*
- Current work is focused on compiling a collection of FORTRAN codes, Mathematica notebooks and packages that can, together with a comprehensive description (aka phd thesis), be used by anyone to reconstruct 2N and 3N calculations.
- Our tools can also be deployed in processes involving EM probes:
 - ▶ *Deuteron Disintegration in Three Dimensions, Few-Body Systems 2012*
- We start employing our three dimensional tools to study the decay of the muonic atom in $\mu^- + d \rightarrow \nu_\mu + n + n$ and other electro-weak processes.