

# *Pionless EFT for Few-Body Systems*

Betzalel Bazak

Physique Theorique

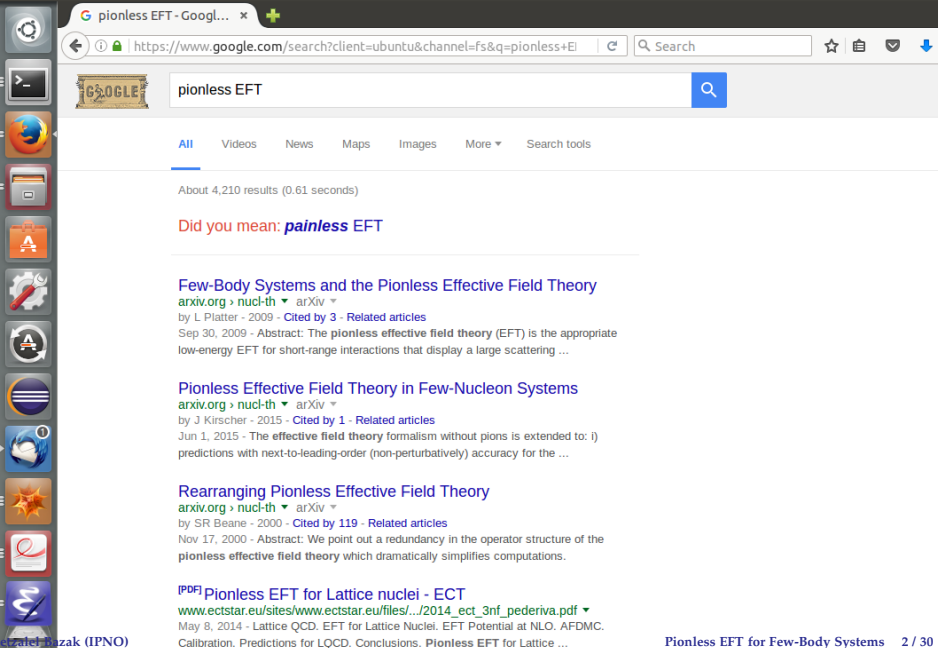
Institut de Physique Nucleaire d'Orsay

Nuclear Physics from Lattice QCD

Institute for Nuclear Theory

April 7, 2016

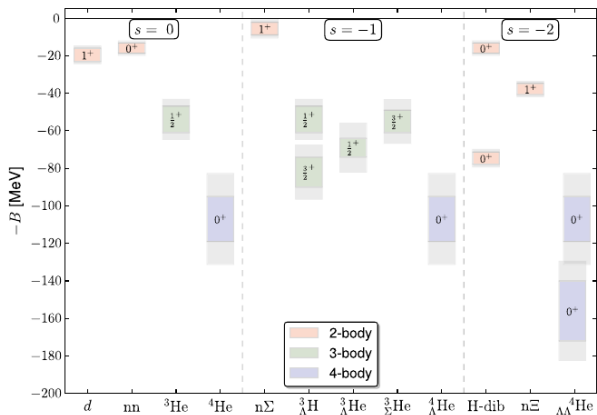
Seattle



Google Search results for "pionless EFT".

Search results for "pionless EFT":

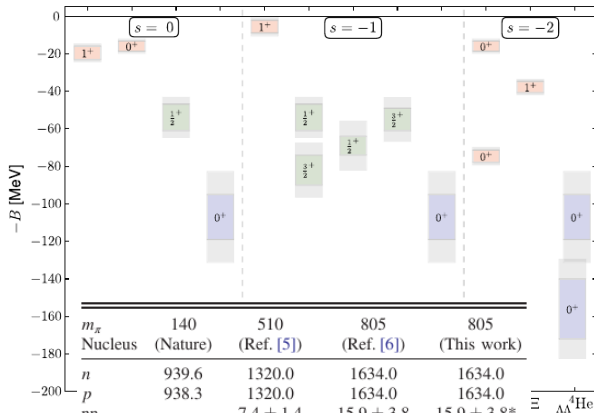
- About 4,210 results (0.61 seconds)
- Did you mean: [painless EFT](#)
- [Few-Body Systems and the Pionless Effective Field Theory](#)  
arxiv.org > nucl-th > arXiv  
by L Platter - 2009 - Cited by 3 - Related articles  
Sep 30, 2009 - Abstract: The pionless effective field theory (EFT) is the appropriate low-energy EFT for short-range interactions that display a large scattering ...
- [Pionless Effective Field Theory in Few-Nucleon Systems](#)  
arxiv.org > nucl-th > arXiv  
by J Kirscher - 2015 - Cited by 1 - Related articles  
Jun 1, 2015 - The effective field theory formalism without pions is extended to: i) predictions with next-to-leading-order (non-perturbatively) accuracy for the ...
- [Rearranging Pionless Effective Field Theory](#)  
arxiv.org > nucl-th > arXiv  
by SR Beane - 2000 - Cited by 119 - Related articles  
Nov 17, 2000 - Abstract: We point out a redundancy in the operator structure of the pionless effective field theory which dramatically simplifies computations.
- [Pionless EFT for Lattice nuclei - ECT](#)  
www.ectstar.eu/sites/www.ectstar.eu/files/.../2014\_ect\_3nf\_pederiva.pdf  
May 8, 2014 - Lattice QCD. EFT for Lattice Nuclei. EFT Potential at NLO. AFDMC. Calibration. Predictions for LQCD. Conclusions. Pionless EFT for Lattice ...



S. R. Beane *et al.* (NPLQCD Collaboration), Phys. Rev. D **87**, 034506 (2013)

N. Barnea *et al.*, Phys. Rev. Lett. **114**, 052501 (2015);

J. Kirscher *et al.*, Phys. Rev. C **92**, 054002 (2015)



S. R. Beane *et al.*

$m_x$	140	510	805	805
Nucleus	(Nature)	(Ref. [5])	(Ref. [6])	(This work)
$n$	939.6	1320.0	1634.0	1634.0
$p$	938.3	1320.0	1634.0	1634.0
$nn$	...	$7.4 \pm 1.4$	$15.9 \pm 3.8$	$15.9 \pm 3.8^*$
$D$	2.224	$11.5 \pm 1.3$	$19.5 \pm 4.8$	$19.5 \pm 4.8^*$
${}^3n$	...	...	...	...
${}^3\text{H}$	8.482	$20.3 \pm 4.5$	$53.9 \pm 10.7$	$53.9 \pm 10.7^*$
${}^3\text{He}$	7.718	$20.3 \pm 4.5$	$53.9 \pm 10.7$	$53.9 \pm 10.7$
${}^4\text{He}$	28.30	$43.0 \pm 14.4$	$107.0 \pm 24.2$	$89 \pm 36$
${}^5\text{He}$	27.50	...	...	$98 \pm 39$
${}^5\text{Li}$	26.61	...	...	$98 \pm 39$
${}^6\text{Li}$	32.00	...	...	$122 \pm 50$

$\Xi$   
 $\Lambda\Lambda$   ${}^4\text{He}$   
34506 (2013)

N. Barnea *et al.*, Phys. Rev. Lett. **114**, 052501 (2015);

J. Kirscher *et al.*, Phys. Rev. C **92**, 054002 (2015)

# Universality

- Consider particles interacting through 2-body potential with range  $R$ .
- Classically, the particles 'feel' each other only within the potential range.
- But, in the case of resonant interaction, the wave function can have much larger extent.
- At low energies, the 2-body physics is completely governed by the scattering length,  $a$ .

$$\lim_{k \rightarrow 0} k \cot \delta(k) = -\frac{1}{a} + \frac{1}{2} r_{\text{eff}} k^2$$

From Sakurai's book

- When  $|a| \gg R$  the potential details has no influence: *Universality*.

# Universality

- Consider particles interacting through 2-body potential with range  $R$ .
- Classically, the particles 'feel' each other only within the potential range.
- But, in the case of resonant interaction, the wave function can have much larger extent.
- At low energies, the 2-body physics is completely governed by the scattering length,  $a$ .

$$\lim_{k \rightarrow 0} k \cot \delta(k) = -\frac{1}{a} + \frac{1}{2} r_{\text{eff}} k^2$$

From Sakurai's book

- When  $|a| \gg R$  the potential details has no influence: *Universality*.

# Universality

- Consider particles interacting through 2-body potential with range  $R$ .
- Classically, the particles 'feel' each other only within the potential range.
- But, in the case of resonant interaction, the wave function can have much larger extent.
- At low energies, the 2-body physics is completely governed by the scattering length,  $a$ .

$$\lim_{k \rightarrow 0} k \cot \delta(k) = -\frac{1}{a} + \frac{1}{2} r_{\text{eff}} k^2$$

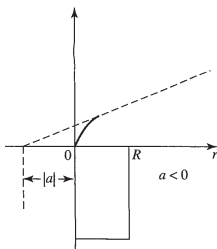
From Sakurai's book

- When  $|a| \gg R$  the potential details has no influence: *Universality*.

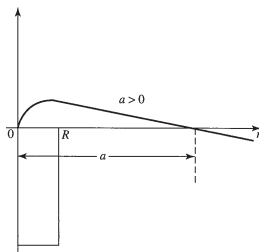
# Universality

- Consider particles interacting through 2-body potential with range  $R$ .
- Classically, the particles 'feel' each other only within the potential range.
- But, in the case of resonant interaction, the wave function can have much larger extent.
- At low energies, the 2-body physics is completely govern by the scattering length,  $a$ .

$$\lim_{k \rightarrow 0} k \cot \delta(k) = -\frac{1}{a} + \frac{1}{2} r_{\text{eff}} k^2$$



From Sakurai's book



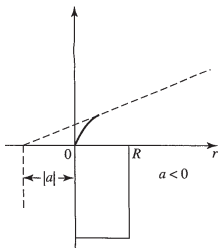
- When  $|a| \gg R$  the potential details has no influence: *Universality*.



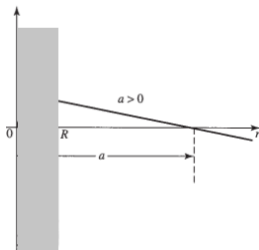
# Universality

- Consider particles interacting through 2-body potential with range  $R$ .
- Classically, the particles 'feel' each other only within the potential range.
- But, in the case of resonant interaction, the wave function can have much larger extent.
- At low energies, the 2-body physics is completely govern by the scattering length,  $a$ .

$$\lim_{k \rightarrow 0} k \cot \delta(k) = -\frac{1}{a} + \frac{1}{2} r_{\text{eff}} k^2$$



From Sakurai's book



- When  $|a| \gg R$  the potential details has no influence: *Universality*.

# Universality

- Generally,  $a \approx r_{\text{eff}} \approx R$ .  
Universal systems are fine-tuned to get  $a \gg r_{\text{eff}}, R$ .
- Corrections to universal theory are of order of  $r_{\text{eff}}/a$  and  $R/a$ .
- For  $a > 0$ , we have universal dimer with energy  $E = -\hbar^2/ma^2$ .
- $^4\text{He}$  Atoms:  $a \approx 170.9a_0$ , ( $a_0$  = the Bohr radius), is much larger than its van der Waals radius,  $r_{\text{vdW}} \approx 9.5a_0$ .
- Nucleus:  $a_s \approx -23.4$  fm,  $a_t \approx 5.42$  fm,  $R = \hbar/m_\pi c \approx 1.4$  fm.  
Deuteron binding energy, 2.22 MeV, is close to  $\hbar/ma_t^2 \approx 1.4$  MeV.
- Ultracold atoms near a Feshbach resonance,

$$a(B) = a_{bg} \left( 1 + \frac{\Delta}{B - B_0} \right)$$

# Universality

- Generally,  $a \approx r_{\text{eff}} \approx R$ .  
Universal systems are fine-tuned to get  $a \gg r_{\text{eff}}, R$ .
- Corrections to universal theory are of order of  $r_{\text{eff}}/a$  and  $R/a$ .
- For  $a > 0$ , we have universal dimer with energy  $E = -\hbar^2/ma^2$ .
- **$^4\text{He}$  Atoms:**  $a \approx 170.9a_0$ , ( $a_0$  = the Bohr radius), is much larger than its van der Waals radius,  $r_{\text{vdW}} \approx 9.5a_0$ .
- **Nucleus:**  $a_s \approx -23.4$  fm,  $a_t \approx 5.42$  fm,  $R = \hbar/m_\pi c \approx 1.4$  fm.  
Deuteron binding energy, 2.22 MeV, is close to  $\hbar/ma_t^2 \approx 1.4$  MeV.
- **Ultracold atoms** near a Feshbach resonance,

$$a(B) = a_{bg} \left( 1 + \frac{\Delta}{B - B_0} \right)$$

# Universality

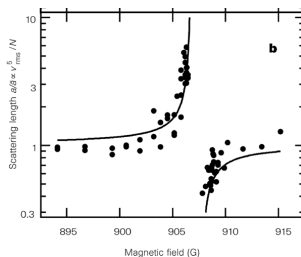
- Generally,  $a \approx r_{\text{eff}} \approx R$ .  
Universal systems are fine-tuned to get  $a \gg r_{\text{eff}}, R$ .
- Corrections to universal theory are of order of  $r_{\text{eff}}/a$  and  $R/a$ .
- For  $a > 0$ , we have universal dimer with energy  $E = -\hbar^2/ma^2$ .
- **$^4\text{He}$  Atoms:**  $a \approx 170.9a_0$ , ( $a_0$  = the Bohr radius), is much larger than its van der Waals radius,  $r_{\text{vdW}} \approx 9.5a_0$ .
- **Nucleus:**  $a_s \approx -23.4$  fm,  $a_t \approx 5.42$  fm,  $R = \hbar/m_\pi c \approx 1.4$  fm.  
Deuteron binding energy, 2.22 MeV, is close to  $\hbar/ma_t^2 \approx 1.4$  MeV.
- **Ultracold atoms** near a Feshbach resonance,

$$a(B) = a_{bg} \left( 1 + \frac{\Delta}{B - B_0} \right)$$

# Universality

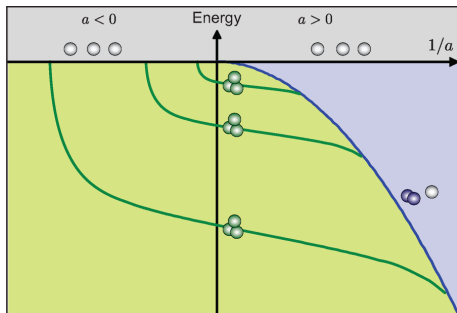
- Generally,  $a \approx r_{\text{eff}} \approx R$ .  
Universal systems are fine-tuned to get  $a \gg r_{\text{eff}}, R$ .
- Corrections to universal theory are of order of  $r_{\text{eff}}/a$  and  $R/a$ .
- For  $a > 0$ , we have universal dimer with energy  $E = -\hbar^2/ma^2$ .
- **$^4\text{He}$  Atoms:**  $a \approx 170.9a_0$ , ( $a_0$  = the Bohr radius), is much larger than its van der Waals radius,  $r_{\text{vdW}} \approx 9.5a_0$ .
- **Nucleus:**  $a_s \approx -23.4$  fm,  $a_t \approx 5.42$  fm,  $R = \hbar/m_\pi c \approx 1.4$  fm.  
Deuteron binding energy, 2.22 MeV, is close to  $\hbar/ma_t^2 \approx 1.4$  MeV.
- **Ultracold atoms** near a Feshbach resonance,

$$a(B) = a_{bg} \left( 1 + \frac{\Delta}{B - B_0} \right)$$



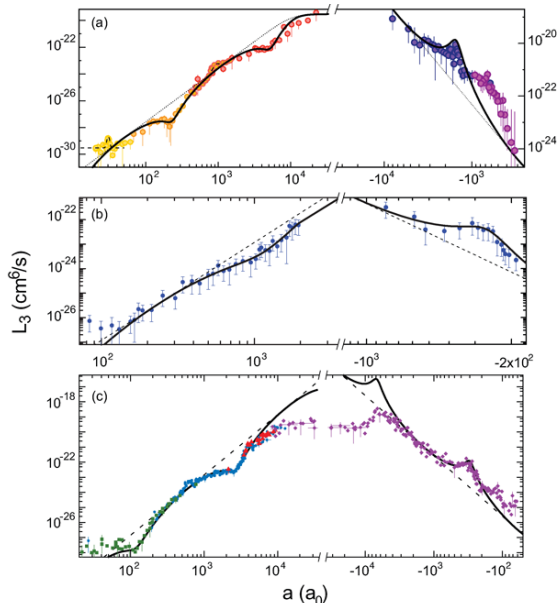
# Efimov Physics

- *The unitary limit:*  $E_2 = 0$ ,  $a \rightarrow \infty$ .
- In 1970 V. Efimov found out that if  $E_2 = 0$  the 3-body system will have an **infinite** number of bound states.
- The 3-body spectrum is  $E_n = E_0 e^{-2\pi n/s_0}$  with  $s_0 = 1.00624$ .



F. Ferlino and R. Grimm, *Physics* **3**, 9 (2010)

# Efimov Physics in Ultracold Atoms



●  $^{39}\text{K}$   
M. Zaccanti *et al.*,  
Nature Phys. **5**, 586 (2009).

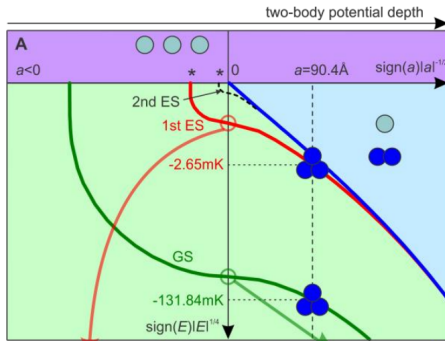
●  $^7\text{Li}$   
N. Gross, Z. Shotan, S.  
Kokkelmans, and L.  
Khaykovich,  
Phys. Rev. Lett. **103**, 163202  
(2009).

●  $^7\text{Li}$   
S.E. Pollack, D. Dries, and  
R.G. Hulet,  
Science **326**, 1683 (2009)

F. Ferlaino and R. Grimm, Physics **3**, 9 (2010)

# $^4\text{He}$ Atoms

- The system of  $^4\text{He}$  atoms is known to be a natural candidate for universal physics, since  $a \approx 170.9a_0 \gg r_{vdW} \approx 9.5a_0$ .
- The  $^4\text{He}$  dimer is bound by 1.62 mK ( $\hbar^2/ma^2 \approx 1.48$  mK).
- $E_3 \approx 131.84$  mK,  $E_3^* \approx 2.6502$  mK, giving a ratio of  $E_3^*/E_3 \approx 49.7$ .
- Recently this excited state was observed experimentally.



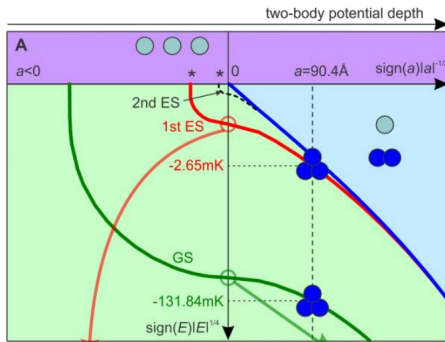
Theory: E. Hiyama and M. Kamimura, *Phys Rev A*. 85, 062505 (2012);

Experiment: M. Kunitski *et al*, *Science* 348 551 (2015).



# $^4\text{He}$ Atoms

- The system of  $^4\text{He}$  atoms is known to be a natural candidate for universal physics, since  $a \approx 170.9a_0 \gg r_{vdW} \approx 9.5a_0$ .
- The  $^4\text{He}$  dimer is bound by 1.62 mK ( $\hbar^2/ma^2 \approx 1.48$  mK).
- $E_3 \approx 131.84$  mK,  $E_3^* \approx 2.6502$  mK, giving a ratio of  $E_3^*/E_3 \approx 49.7$ .
- Recently this excited state was observed experimentally.

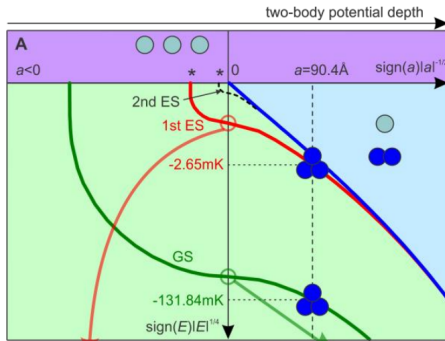


Theory: E. Hiyama and M. Kamimura, *Phys Rev A*. 85, 062505 (2012);

Experiment: M. Kunitski *et al.*, *Science* 348 551 (2015).

# $^4\text{He}$ Atoms

- The system of  $^4\text{He}$  atoms is known to be a natural candidate for universal physics, since  $a \approx 170.9a_0 \gg r_{vdW} \approx 9.5a_0$ .
- The  $^4\text{He}$  dimer is bound by 1.62 mK ( $\hbar^2/ma^2 \approx 1.48$  mK).
- $E_3 \approx 131.84$  mK,  $E_3^* \approx 2.6502$  mK, giving a ratio of  $E_3^*/E_3 \approx 49.7$ .
- Recently this excited state was observed experimentally.

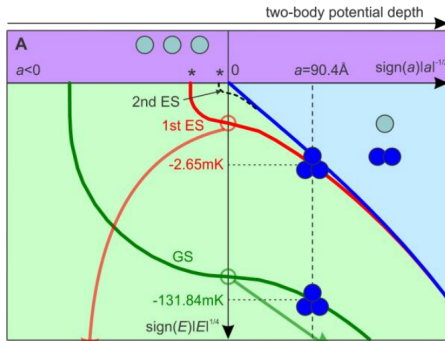


Theory: E. Hiyama and M. Kamimura, *Phys Rev A*. 85, 062505 (2012);

Experiment: M. Kunitski *et al.*, *Science* 348 551 (2015).

# $^4\text{He}$ Atoms

- The system of  $^4\text{He}$  atoms is known to be a natural candidate for universal physics, since  $a \approx 170.9a_0 \gg r_{vdW} \approx 9.5a_0$ .
- The  $^4\text{He}$  dimer is bound by 1.62 mK ( $\hbar^2/ma^2 \approx 1.48$  mK).
- $E_3 \approx 131.84$  mK,  $E_3^* \approx 2.6502$  mK, giving a ratio of  $E_3^*/E_3 \approx 49.7$ .
- Recently this excited state was observed experimentally.



Theory: E. Hiyama and M. Kamimura, *Phys Rev A*. **85**, 062505 (2012);

Experiment: M. Kunitski *et al.*, *Science* **348** 551 (2015).

# Effective Field Theory (EFT)

- Typically in physics we have an “underlying” theory, valid at a mass scale  $M_{hi}$ , but we want to study processes at momenta  $Q \approx M_{lo} \ll M_{hi}$ .
- For example, nuclear structure involves energies that are much smaller than the typical QCD mass scale,  $M_{QCD} \approx 1 \text{ GeV}$ .
- Effective Field Theory (EFT) is a framework to construct the interactions systematically. The high-energy degrees of freedom are integrated out, while the effective Lagrangian have the same symmetries as the underlying theory.
- The details of the underlying dynamics are contained in the interaction strengths.

# Effective Field Theory (EFT)

- Typically in physics we have an “underlying” theory, valid at a mass scale  $M_{hi}$ , but we want to study processes at momenta  $Q \approx M_{lo} \ll M_{hi}$ .
- For example, nuclear structure involves energies that are much smaller than the typical QCD mass scale,  $M_{QCD} \approx 1 \text{ GeV}$ .
- Effective Field Theory (EFT) is a framework to construct the interactions systematically. The high-energy degrees of freedom are integrated out, while the effective Lagrangian have the same symmetries as the underlying theory.
- The details of the underlying dynamics are contained in the interaction strengths.

# Effective Field Theory (EFT)

- Typically in physics we have an “underlying” theory, valid at a mass scale  $M_{hi}$ , but we want to study processes at momenta  $Q \approx M_{lo} \ll M_{hi}$ .
- For example, nuclear structure involves energies that are much smaller than the typical QCD mass scale,  $M_{QCD} \approx 1 \text{ GeV}$ .
- Effective Field Theory (EFT) is a framework to construct the interactions systematically. The high-energy degrees of freedom are integrated out, while the effective Lagrangian have the same symmetries as the underlying theory.
- The details of the underlying dynamics are contained in the interaction strengths.

# Effective Field Theory (EFT)

- Typically in physics we have an “underlying” theory, valid at a mass scale  $M_{hi}$ , but we want to study processes at momenta  $Q \approx M_{lo} \ll M_{hi}$ .
- For example, nuclear structure involves energies that are much smaller than the typical QCD mass scale,  $M_{QCD} \approx 1 \text{ GeV}$ .
- Effective Field Theory (EFT) is a framework to construct the interactions systematically. The high-energy degrees of freedom are integrated out, while the effective Lagrangian have the same symmetries as the underlying theory.
- The details of the underlying dynamics are contained in the interaction strengths.

# Pionless EFT

- The degrees of freedom in pionless EFT are the nucleons.
- We have to include all terms conserving our theory symmetries, order by order.
- For nucleons, the Leading Order (LO) is,

$$V_{LO} = a_1 + a_2 \sigma_i \cdot \sigma_j + a_3 \tau_i \cdot \tau_j + a_4 (\sigma_i \cdot \sigma_j) (\tau_i \cdot \tau_j)$$

where due to symmetry, only 2 are independent, corresponding to the two scattering lengths.

- The Next to Leading Order (NLO) is,

$$\begin{aligned} V_{NLO} &= b_1(k^2 + q^2) + b_2(k^2 + q^2) \sigma_i \cdot \sigma_j + b_3(k^2 + q^2) \tau_i \cdot \tau_j \\ &+ b_4(k^2 + q^2) (\sigma_i \cdot \sigma_j) (\tau_i \cdot \tau_j) \\ &q = p' - p, \quad k = p + p' \end{aligned}$$

here also only 2 parameters are independent, corresponding to the two effective ranges.

- $p$ -wave enters at  $N^3LO$ !



# Pionless EFT

- The degrees of freedom in pionless EFT are the nucleons.
- We have to include all terms conserving our theory symmetries, order by order.
- For nucleons, the Leading Order (LO) is,

$$V_{LO} = a_1 + a_2 \sigma_i \cdot \sigma_j + a_3 \tau_i \cdot \tau_j + a_4 (\sigma_i \cdot \sigma_j) (\tau_i \cdot \tau_j)$$

where due to symmetry, only 2 are independent, corresponding to the two scattering lengths.

- The Next to Leading Order (NLO) is,

$$\begin{aligned} V_{NLO} &= b_1(k^2 + q^2) + b_2(k^2 + q^2) \sigma_i \cdot \sigma_j + b_3(k^2 + q^2) \tau_i \cdot \tau_j \\ &+ b_4(k^2 + q^2) (\sigma_i \cdot \sigma_j) (\tau_i \cdot \tau_j) \\ &q = p' - p, \quad k = p + p' \end{aligned}$$

here also only 2 parameters are independent, corresponding to the two effective ranges.

- $p$ -wave enters at  $N^3LO$ !

# Pionless EFT

- The degrees of freedom in pionless EFT are the nucleons.
- We have to include all terms conserving our theory symmetries, order by order.
- For nucleons, the Leading Order (LO) is,

$$V_{LO} = a_1 + a_2 \sigma_i \cdot \sigma_j + a_3 \tau_i \cdot \tau_j + a_4 (\sigma_i \cdot \sigma_j) (\tau_i \cdot \tau_j)$$

where due to symmetry, only 2 are independent, corresponding to the two scattering lengths.

- The Next to Leading Order (NLO) is,

$$\begin{aligned} V_{NLO} &= b_1(k^2 + q^2) + b_2(k^2 + q^2) \sigma_i \cdot \sigma_j + b_3(k^2 + q^2) \tau_i \cdot \tau_j \\ &+ b_4(k^2 + q^2) (\sigma_i \cdot \sigma_j) (\tau_i \cdot \tau_j) \\ &q = p' - p, \quad k = p + p' \end{aligned}$$

here also only 2 parameters are independent, corresponding to the two effective ranges.

- $p$ -wave enters at  $N^3LO$ !

# Pionless EFT

- The degrees of freedom in pionless EFT are the nucleons.
- We have to include all terms conserving our theory symmetries, order by order.
- For nucleons, the Leading Order (LO) is,

$$V_{LO} = a_1 + a_2 \sigma_i \cdot \sigma_j + a_3 \tau_i \cdot \tau_j + a_4 (\sigma_i \cdot \sigma_j) (\tau_i \cdot \tau_j)$$

where due to symmetry, only 2 are independent, corresponding to the two scattering lengths.

- The Next to Leading Order (NLO) is,

$$\begin{aligned} V_{NLO} &= b_1(k^2 + q^2) + b_2(k^2 + q^2) \sigma_i \cdot \sigma_j + b_3(k^2 + q^2) \tau_i \cdot \tau_j \\ &+ b_4(k^2 + q^2) (\sigma_i \cdot \sigma_j) (\tau_i \cdot \tau_j) \\ &\quad \mathbf{q} = \mathbf{p}' - \mathbf{p}, \quad \mathbf{k} = \mathbf{p} + \mathbf{p}' \end{aligned}$$

here also only 2 parameters are independent, corresponding to the two effective ranges.

- $p$ -wave enters at  $N^3LO$ !

# Pionless EFT

- The degrees of freedom in pionless EFT are the nucleons.
- We have to include all terms conserving our theory symmetries, order by order.
- For nucleons, the Leading Order (LO) is,

$$V_{LO} = a_1 + a_2 \sigma_i \cdot \sigma_j + a_3 \tau_i \cdot \tau_j + a_4 (\sigma_i \cdot \sigma_j) (\tau_i \cdot \tau_j)$$

where due to symmetry, only 2 are independent, corresponding to the two scattering lengths.

- The Next to Leading Order (NLO) is,

$$\begin{aligned} V_{NLO} &= b_1(k^2 + q^2) + b_2(k^2 + q^2) \sigma_i \cdot \sigma_j + b_3(k^2 + q^2) \tau_i \cdot \tau_j \\ &+ b_4(k^2 + q^2) (\sigma_i \cdot \sigma_j) (\tau_i \cdot \tau_j) \\ &\quad \mathbf{q} = \mathbf{p}' - \mathbf{p}, \quad \mathbf{k} = \mathbf{p} + \mathbf{p}' \end{aligned}$$

here also only 2 parameters are independent, corresponding to the two effective ranges.

- $p$ -wave enters at  $N^3LO$ !

# Short-Range EFT for Bosonic system

- For spinless bosons, most of the terms are dropped, and we have at LO,

$$V_{LO} = a_1.$$

- At NLO,

$$V_{NLO} = b_1(k^2 + q^2).$$

- Terms proportional to  $k \cdot q$  are omitted due to time reversal symmetry.
- The LO term is iterated at NLO.

# Short-Range EFT for Bosonic system

- For spinless bosons, most of the terms are dropped, and we have at LO,

$$V_{LO} = a_1.$$

- At NLO,

$$V_{NLO} = b_1(k^2 + q^2).$$

- Terms proportional to  $k \cdot q$  are omitted due to time reversal symmetry.
- The LO term is iterated at NLO.

# Short-Range EFT for Bosonic system

- For spinless bosons, most of the terms are dropped, and we have at LO,

$$V_{LO} = a_1.$$

- At NLO,

$$V_{NLO} = b_1(k^2 + q^2).$$

- Terms proportional to  $\mathbf{k} \cdot \mathbf{q}$  are omitted due to time reversal symmetry.
- The LO term is iterated at NLO.

# Short-Range EFT for Bosonic system

- For spinless bosons, most of the terms are dropped, and we have at LO,

$$V_{LO} = a_1.$$

- At NLO,

$$V_{NLO} = b_1(k^2 + q^2).$$

- Terms proportional to  $\mathbf{k} \cdot \mathbf{q}$  are omitted due to time reversal symmetry.
- The LO term is iterated at NLO.



# Stochastic Variational Method I

- To solve the  $N$ -body Schrodinger equation, we use **correlated Gaussian basis**,

$$\psi(\eta_1, \eta_2 \dots \eta_{A-1}) = \sum_i c_i \mathcal{A} \exp(-\eta^T A_i \eta)$$

$\eta$  = Jacobi coordinates,  $A_i$  = matrix of  $(N-1) \times (N-1)$  numbers.

- Since  $a \gg \Lambda^{-1}$ , we need **large spread** of basis functions.
- Symmetrization gives factor of  $N!$  which limits the number of particles.
- Works for any angular momentum, for bosons and fermions. Spin and isospin can be introduced.

Y. Suzuki and K. Varga, "Stochastic Variational Approach to Quantum-Mechanical Few-Body Problems", Springer 1998.

# Stochastic Variational Method I

- To solve the  $N$ -body Schrodinger equation, we use **correlated Gaussian basis**,

$$\psi(\eta_1, \eta_2 \dots \eta_{A-1}) = \sum_i c_i \mathcal{A} \exp(-\eta^T A_i \eta)$$

$\eta$  = Jacobi coordinates,  $A_i$  = matrix of  $(N-1) \times (N-1)$  numbers.

- Since  $a \gg \Lambda^{-1}$ , we need **large spread** of basis functions.
- Symmetrization gives factor of  $N!$  which limits the number of particles.
- Works for any angular momentum, for bosons and fermions. Spin and isospin can be introduced.

Y. Suzuki and K. Varga, "Stochastic Variational Approach to Quantum-Mechanical Few-Body Problems", Springer 1998.

# Stochastic Variational Method I

- To solve the  $N$ -body Schrodinger equation, we use **correlated Gaussian basis**,

$$\psi(\eta_1, \eta_2 \dots \eta_{A-1}) = \sum_i c_i \mathcal{A} \exp(-\eta^T A_i \eta)$$

$\eta$  = Jacobi coordinates,  $A_i$  = matrix of  $(N-1) \times (N-1)$  numbers.

- Since  $a \gg \Lambda^{-1}$ , we need **large spread** of basis functions.
- Symmetrization gives factor of  **$N!$**  which limits the number of particles.
- Works for any angular momentum, for bosons and fermions. Spin and isospin can be introduced.

Y. Suzuki and K. Varga, "Stochastic Variational Approach to Quantum-Mechanical Few-Body Problems", Springer 1998.

# Stochastic Variational Method I

- To solve the  $N$ -body Schrodinger equation, we use **correlated Gaussian basis**,

$$\psi(\eta_1, \eta_2 \dots \eta_{A-1}) = \sum_i c_i \mathcal{A} \exp(-\eta^T A_i \eta)$$

$\eta$  = Jacobi coordinates,  $A_i$  = matrix of  $(N-1) \times (N-1)$  numbers.

- Since  $a \gg \Lambda^{-1}$ , we need **large spread** of basis functions.
- Symmetrization gives factor of  **$N!$**  which limits the number of particles.
- Works for any angular momentum, for bosons and fermions. Spin and isospin can be introduced.

Y. Suzuki and K. Varga, "Stochastic Variational Approach to Quantum-Mechanical Few-Body Problems", Springer 1998.

# Stochastic Variational Method II

- The matrix elements can be calculated **analytically** in most cases:

$$\langle A|A' \rangle = \left( \frac{(2\pi)^{N-1}}{\det B} \right)^{3/2}; \quad B = A + A'$$

$$\langle A|T_{int}|A' \rangle = 3\langle A|A' \rangle \text{Tr}[AB^{-1}A'\Pi]; \quad \Pi_{ij} = (2\mu_i)^{-1}\delta_{ij}$$

$$\langle A|V^{2b}|A' \rangle = \frac{\Lambda^2 C^{(0)}}{2m\sqrt{\pi}} \langle A|A' \rangle \sum_{i<j} \left( 1 + f_{ij}\Lambda^2/2 \right)^{-3/2}; \quad f_{ij} = \omega_{ij}^T B^{-1} \omega_{ij}, \quad r_{ij} = \omega_{ij}^T \eta$$

$$\langle A|V^{3b}|A' \rangle = \frac{\Lambda^2 D^{(0)}}{16m\pi^2} \langle A|A' \rangle \sum_{i<j<k} \sum_{cyc} \left( \det(I + \Lambda^2 F_{ijk}/2) \right)^{-3/2}; \quad F_{ijk} = \Omega_{ijk}^T B^{-1} \Omega_{ijk}$$

$$\Omega_{ijk} = (\omega_{ik} \ \omega_{jk})$$

# Stochastic Variational Method II

- The matrix elements can be calculated **analytically** in most cases:

$$\langle A|A' \rangle = \left( \frac{(2\pi)^{N-1}}{\det B} \right)^{3/2}; \quad B = A + A'$$

$$\langle A|T_{int}|A' \rangle = 3\langle A|A' \rangle \text{Tr}[AB^{-1}A'\Pi]; \quad \Pi_{ij} = (2\mu_i)^{-1}\delta_{ij}$$

$$\langle A|V^{2b}|A' \rangle = \frac{\Lambda^2 C^{(0)}}{2m\sqrt{\pi}} \langle A|A' \rangle \sum_{i<j} \left( 1 + f_{ij}\Lambda^2/2 \right)^{-3/2}; \quad f_{ij} = \omega_{ij}^T B^{-1} \omega_{ij}, \quad r_{ij} = \omega_{ij}^T \eta$$

$$\langle A|V^{3b}|A' \rangle = \frac{\Lambda^2 D^{(0)}}{16m\pi^2} \langle A|A' \rangle \sum_{i<j<k} \sum_{cyc} \left( \det(I + \Lambda^2 F_{ijk}/2) \right)^{-3/2}; \quad F_{ijk} = \Omega_{ijk}^T B^{-1} \Omega_{ijk}$$

$$\Omega_{ijk} = (\omega_{ik} \ \omega_{jk})$$

# Stochastic Variational Method II

- The matrix elements can be calculated **analytically** in most cases:

$$\langle A|A' \rangle = \left( \frac{(2\pi)^{N-1}}{\det B} \right)^{3/2}; \quad B = A + A'$$

$$\langle A|T_{int}|A' \rangle = 3\langle A|A' \rangle \text{Tr}[AB^{-1}A'\Pi]; \quad \Pi_{ij} = (2\mu_i)^{-1}\delta_{ij}$$

$$\langle A|V^{2b}|A' \rangle = \frac{\Lambda^2 C^{(0)}}{2m\sqrt{\pi}} \langle A|A' \rangle \sum_{i<j} \left( 1 + f_{ij}\Lambda^2/2 \right)^{-3/2}; \quad f_{ij} = \omega_{ij}^T B^{-1} \omega_{ij}, \quad r_{ij} = \omega_{ij}^T \eta$$

$$\langle A|V^{3b}|A' \rangle = \frac{\Lambda^2 D^{(0)}}{16m\pi^2} \langle A|A' \rangle \sum_{i<j<k} \sum_{cyc} \left( \det(I + \Lambda^2 F_{ijk}/2) \right)^{-3/2}; \quad F_{ijk} = \Omega_{ijk}^T B^{-1} \Omega_{ijk}$$

$$\Omega_{ijk} = (\omega_{ik} \ \omega_{jk})$$

# Stochastic Variational Method II

- The matrix elements can be calculated **analytically** in most cases:

$$\langle A|A' \rangle = \left( \frac{(2\pi)^{N-1}}{\det B} \right)^{3/2}; \quad B = A + A'$$

$$\langle A|T_{int}|A' \rangle = 3\langle A|A' \rangle \text{Tr}[AB^{-1}A'\Pi]; \quad \Pi_{ij} = (2\mu_i)^{-1}\delta_{ij}$$

$$\langle A|V^{2b}|A' \rangle = \frac{\Lambda^2 C^{(0)}}{2m\sqrt{\pi}} \langle A|A' \rangle \sum_{i<j} \left( 1 + f_{ij}\Lambda^2/2 \right)^{-3/2}; \quad f_{ij} = \omega_{ij}^T B^{-1} \omega_{ij}, \quad r_{ij} = \omega_{ij}^T \eta$$

$$\langle A|V^{3b}|A' \rangle = \frac{\Lambda^2 D^{(0)}}{16m\pi^2} \langle A|A' \rangle \sum_{i<j<k} \sum_{cyc} \left( \det(I + \Lambda^2 F_{ijk}/2) \right)^{-3/2}; \quad F_{ijk} = \Omega_{ijk}^T B^{-1} \Omega_{ijk}$$

$$\Omega_{ijk} = (\omega_{ik} \ \omega_{jk})$$



# Stochastic Variational Method II

- The matrix elements can be calculated **analytically** in most cases:

$$\langle A|A' \rangle = \left( \frac{(2\pi)^{N-1}}{\det B} \right)^{3/2}; \quad B = A + A'$$

$$\langle A|T_{int}|A' \rangle = 3\langle A|A' \rangle \text{Tr}[AB^{-1}A'\Pi]; \quad \Pi_{ij} = (2\mu_i)^{-1}\delta_{ij}$$

$$\langle A|V^{2b}|A' \rangle = \frac{\Lambda^2 C^{(0)}}{2m\sqrt{\pi}} \langle A|A' \rangle \sum_{i < j} \left( 1 + f_{ij}\Lambda^2/2 \right)^{-3/2}; \quad f_{ij} = \omega_{ij}^T B^{-1} \omega_{ij}, \quad \mathbf{r}_{ij} = \omega_{ij}^T \boldsymbol{\eta}$$

$$\langle A|V^{3b}|A' \rangle = \frac{\Lambda^2 D^{(0)}}{16m\pi^2} \langle A|A' \rangle \sum_{i < j < k} \sum_{cyc} \left( \det(I + \Lambda^2 F_{ijk}/2) \right)^{-3/2}; \quad F_{ijk} = \Omega_{ijk}^T B^{-1} \Omega_{ijk}$$

$$\Omega_{ijk} = (\omega_{ik} \ \omega_{jk})$$

# Stochastic Variational Method II

- The matrix elements can be calculated **analytically** in most cases:

$$\langle A|A' \rangle = \left( \frac{(2\pi)^{N-1}}{\det B} \right)^{3/2}; \quad B = A + A'$$

$$\langle A|T_{int}|A' \rangle = 3\langle A|A' \rangle \text{Tr}[AB^{-1}A'\Pi]; \quad \Pi_{ij} = (2\mu_i)^{-1}\delta_{ij}$$

$$\langle A|V^{2b}|A' \rangle = \frac{\Lambda^2 C^{(0)}}{2m\sqrt{\pi}} \langle A|A' \rangle \sum_{i<j} \left( 1 + f_{ij}\Lambda^2/2 \right)^{-3/2}; \quad f_{ij} = \omega_{ij}^T B^{-1} \omega_{ij}, \quad \mathbf{r}_{ij} = \omega_{ij}^T \boldsymbol{\eta}$$

$$\langle A|V^{3b}|A' \rangle = \frac{\Lambda^2 D^{(0)}}{16m\pi^2} \langle A|A' \rangle \sum_{i<j<k} \sum_{cyc} \left( \det(I + \Lambda^2 F_{ijk}/2) \right)^{-3/2}; \quad F_{ijk} = \Omega_{ijk}^T B^{-1} \Omega_{ijk}$$

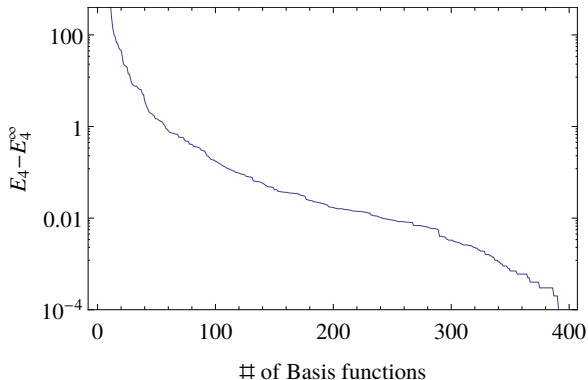
$$\Omega_{ijk} = (\omega_{ik} \ \omega_{jk})$$

# Stochastic Variational Method III

- To find the best  $A_i$ , we use the **Stochastic Variational Method (SVM)**:
- We add basis function one by one, or try to replace an exist basis function by a new one.
- We choose **randomly** the matrix  $A_i$  element by element, trying to minimize the energy.
- According to the variational principle, an **upper bound** for the ground (excited) state is achieved.

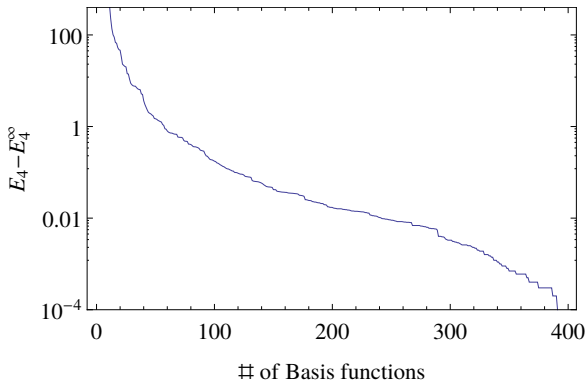
# Stochastic Variational Method III

- To find the best  $A_i$ , we use the **Stochastic Variational Method (SVM)**:
- We add basis function one by one, or try to replace an exist basis function by a new one.
- We choose **randomly** the matrix  $A_i$  element by element, trying to minimize the energy.
- According to the variational principle, an **upper bound** for the ground (excited) state is achieved.



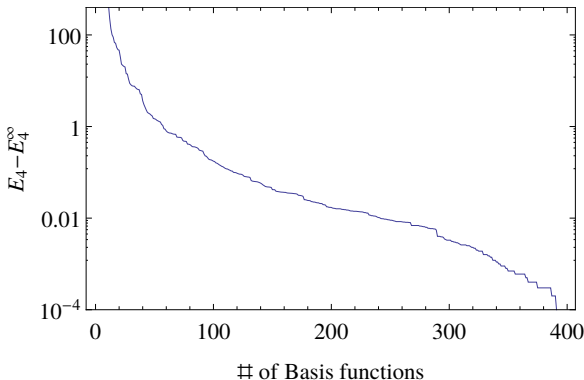
# Stochastic Variational Method III

- To find the best  $A_i$ , we use the **Stochastic Variational Method (SVM)**:
- We add basis function one by one, or try to replace an exist basis function by a new one.
- We choose **randomly** the matrix  $A_i$  element by element, trying to minimize the energy.
- According to the variational principle, an **upper bound** for the ground (excited) state is achieved.



# Stochastic Variational Method III

- To find the best  $A_i$ , we use the **Stochastic Variational Method (SVM)**:
- We add basis function one by one, or try to replace an exist basis function by a new one.
- We choose **randomly** the matrix  $A_i$  element by element, trying to minimize the energy.
- According to the variational principle, an **upper bound** for the ground (excited) state is achieved.



# Regularization I: non local potential

- At LO, we have only contact interaction,

$$V(r_{ij}) = g\delta(r_{ij}).$$

- This interaction needs **regularization** and **renormalization**.
- The bound state of two identical bosons (here  $\hbar = m = 1$ ),

$$-\nabla^2\psi(r) + g\delta(r)\psi(r) = -B_2\psi(r)$$

and in momentum space,

$$p^2\phi(p) + g \int \frac{d^3p'}{(2\pi)^3} \phi(p') = -B_2\phi(p)$$

- Therefore,

$$\frac{1}{g} = \int \frac{d^3p'}{(2\pi)^3} \frac{1}{p'^2 + B_2}$$

which **diverges!**

# Regularization I: non local potential

- At LO, we have only contact interaction,

$$V(r_{ij}) = g\delta(r_{ij}).$$

- This interaction needs **regularization** and **renormalization**.
- The bound state of two identical bosons (here  $\hbar = m = 1$ ),

$$-\nabla^2\psi(r) + g\delta(r)\psi(r) = -B_2\psi(r)$$

and in momentum space,

$$p^2\phi(p) + g \int \frac{d^3p'}{(2\pi)^3} \phi(p') = -B_2\phi(p)$$

- Therefore,

$$\frac{1}{g} = \int \frac{d^3p'}{(2\pi)^3} \frac{1}{p'^2 + B_2}$$

which **diverges!**



# Regularization I: non local potential

- At LO, we have only contact interaction,

$$V(r_{ij}) = g\delta(r_{ij}).$$

- This interaction needs **regularization** and **renormalization**.
- The bound state of two identical bosons (here  $\hbar = m = 1$ ),

$$-\nabla^2\psi(r) + g\delta(r)\psi(r) = -B_2\psi(r)$$

and in momentum space,

$$p^2\phi(p) + g \int \frac{d^3p'}{(2\pi)^3} \phi(p') = -B_2\phi(p)$$

- Therefore,

$$\frac{1}{g} = \int \frac{d^3p'}{(2\pi)^3} \frac{1}{p'^2 + B_2}$$

which **diverges!**

# Regularization I: non local potential

- At LO, we have only contact interaction,

$$V(r_{ij}) = g\delta(r_{ij}).$$

- This interaction needs **regularization** and **renormalization**.
- The bound state of two identical bosons (here  $\hbar = m = 1$ ),

$$-\nabla^2\psi(r) + g\delta(r)\psi(r) = -B_2\psi(r)$$

and in momentum space,

$$p^2\phi(p) + g \int \frac{d^3p'}{(2\pi)^3} \phi(p') = -B_2\phi(p)$$

- Therefore,

$$\frac{1}{g} = \int \frac{d^3p'}{(2\pi)^3} \frac{1}{p'^2 + B_2}$$

which **diverges!**

# Regularization I: non local potential

- To regularize, we can smear the interaction over a range of  $1/\Lambda$ ,

$$\delta_{\Lambda}(r) \xrightarrow{\Lambda \rightarrow \infty} \delta(r).$$

- Doing so for the incoming and outgoing momenta we have,

$$\frac{1}{g} = \int \frac{d^3p'}{(2\pi)^3} \frac{\exp(-2p'^2/\Lambda^2)}{p'^2 + B_2}$$

- Which can be expanded by powers of  $Q/\Lambda$ , ( $Q = \sqrt{B_2}$ )

$$g = \frac{8\sqrt{2}\pi^{3/2}}{\Lambda} \left( 1 + \sqrt{\pi} \frac{Q}{\Lambda} + \dots \right).$$

- With the price of **non-local** potential,

$$\langle r|V|r' \rangle = g\delta_{\Lambda}(r)\delta_{\Lambda}(r')$$

R.F. Mohr *et al.*, Ann. Phys. **321**, 225 (2006).

# Regularization I: non local potential

- To regularize, we can smear the interaction over a range of  $1/\Lambda$ ,

$$\delta_{\Lambda}(r) \xrightarrow{\Lambda \rightarrow \infty} \delta(r).$$

- Doing so for the incoming and outgoing momenta we have,

$$\frac{1}{g} = \int \frac{d^3p'}{(2\pi)^3} \frac{\exp(-2p'^2/\Lambda^2)}{p'^2 + B_2}$$

- Which can be expanded by powers of  $Q/\Lambda$ , ( $Q = \sqrt{B_2}$ )

$$g = \frac{8\sqrt{2}\pi^{3/2}}{\Lambda} \left( 1 + \sqrt{\pi} \frac{Q}{\Lambda} + \dots \right).$$

- With the price of **non-local** potential,

$$\langle r|V|r' \rangle = g\delta_{\Lambda}(r)\delta_{\Lambda}(r')$$

R.F. Mohr *et al.*, Ann. Phys. **321**, 225 (2006).

# Regularization I: non local potential

- To regularize, we can smear the interaction over a range of  $1/\Lambda$ ,

$$\delta_\Lambda(r) \xrightarrow{\Lambda \rightarrow \infty} \delta(r).$$

- Doing so for the incoming and outgoing momenta we have,

$$\frac{1}{g} = \int \frac{d^3p'}{(2\pi)^3} \frac{\exp(-2p'^2/\Lambda^2)}{p'^2 + B_2}$$

- Which can be expanded by powers of  $Q/\Lambda$ , ( $Q = \sqrt{B_2}$ )

$$g = \frac{8\sqrt{2}\pi^{3/2}}{\Lambda} \left( 1 + \sqrt{\pi} \frac{Q}{\Lambda} + \dots \right).$$

- With the price of non-local potential,

$$\langle r|V|r' \rangle = g\delta_\Lambda(r)\delta_\Lambda(r')$$

R.F. Mohr *et al.*, Ann. Phys. **321**, 225 (2006).

# Regularization I: non local potential

- To regularize, we can smear the interaction over a range of  $1/\Lambda$ ,

$$\delta_\Lambda(r) \xrightarrow{\Lambda \rightarrow \infty} \delta(r).$$

- Doing so for the incoming and outgoing momenta we have,

$$\frac{1}{g} = \int \frac{d^3p'}{(2\pi)^3} \frac{\exp(-2p'^2/\Lambda^2)}{p'^2 + B_2}$$

- Which can be expanded by powers of  $Q/\Lambda$ , ( $Q = \sqrt{B_2}$ )

$$g = \frac{8\sqrt{2}\pi^{3/2}}{\Lambda} \left( 1 + \sqrt{\pi} \frac{Q}{\Lambda} + \dots \right).$$

- With the price of **non-local** potential,

$$\langle r|V|r' \rangle = g\delta_\Lambda(r)\delta_\Lambda(r')$$

R.F. Mohr *et al.*, Ann. Phys. **321**, 225 (2006).

# Regularization II: local potential

- We can cut the momentum transfer  $q = p - p'$ , to get a **local** potential,

$$\langle r|V|r' \rangle = g\delta_\Lambda(r)\delta(r - r')$$

but now the two-body equation is to be solved numerically.

- The LEC is renormalized by fixing one observable, like the dimer binding energy or the scattering length, to its physical value.
- Using dimension less LEC,

$$V_{LO}(r) = \frac{4\pi\hbar^2}{m\Lambda} C^{(0)}(\Lambda)\delta_\Lambda(r), \quad C^{(0)}(\Lambda) = 2.38 \left( 1 + \frac{2.25}{a\Lambda} - \frac{4.68}{(a\Lambda)^2} + \dots \right)$$

# Regularization II: local potential

- We can cut the momentum transfer  $q = p - p'$ , to get a **local** potential,

$$\langle r|V|r'\rangle = g\delta_\Lambda(r)\delta(r - r')$$

but now the two-body equation is to be solved numerically.

- The LEC is renormalized by fixing one observable, like the dimer binding energy or the scattering length, to its physical value.
- Using dimension less LEC,

$$V_{LO}(r) = \frac{4\pi\hbar^2}{m\Lambda} C^{(0)}(\Lambda)\delta_\Lambda(r), \quad C^{(0)}(\Lambda) = 2.38 \left( 1 + \frac{2.25}{a\Lambda} - \frac{4.68}{(a\Lambda)^2} + \dots \right)$$



## Regularization II: local potential

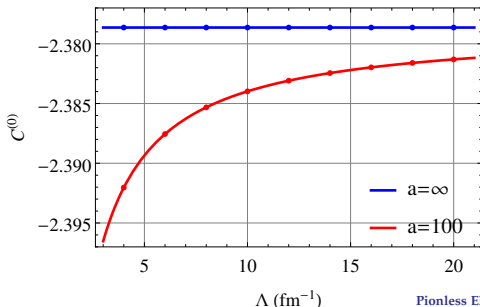
- We can cut the momentum transfer  $q = p - p'$ , to get a **local** potential,

$$\langle r|V|r' \rangle = g\delta_{\Lambda}(r)\delta(r - r')$$

but now the two-body equation is to be solved numerically.

- The LEC is renormalized by fixing one observable, like the dimer binding energy or the scattering length, to its physical value.
- Using dimension less LEC,

$$V_{LO}(r) = \frac{4\pi\hbar^2}{m\Lambda} C^{(0)}(\Lambda)\delta_{\Lambda}(r), \quad C^{(0)}(\Lambda) = 2.38 \left( 1 + \frac{2.25}{a\Lambda} - \frac{4.68}{(a\Lambda)^2} + \dots \right)$$



# Next to Leading Order

- At NLO, the LO term is iterated and 2-derivatives term is added:

$$V_{NLO}(r) = \frac{4\pi\hbar^2}{m\Lambda} \delta_\Lambda(\mathbf{r}) \left\{ C_0^{(1)}(\Lambda) + C_2^{(1)}(\Lambda) \left[ \overleftarrow{\nabla}^2 + \overrightarrow{\nabla}^2 \right] \right\}$$

- NLO term is to be taken in **perturbative way**.

- For energies,

$$\Delta E = \langle \psi_{LO} | V_{NLO} | \psi_{LO} \rangle$$

- For scattering amplitude - distorted wave Born approximation,

$$\Delta f_k = -\frac{m}{k^2} \int dr \psi_{LO}^2 V_{NLO}$$

$$f_k \approx \frac{1}{-a^{-1} - ik} \left( 1 - \frac{1}{-a^{-1} - ik} \left[ \frac{1}{2} r_{\text{eff}} k^2 + \frac{\delta a}{a^2} \right] \right)$$

# Next to Leading Order

- At NLO, the LO term is iterated and 2-derivatives term is added:

$$V_{NLO}(r) = \frac{4\pi\hbar^2}{m\Lambda} \delta_\Lambda(r) \left\{ C_0^{(1)}(\Lambda) + C_2^{(1)}(\Lambda) \left[ \overleftarrow{\nabla}^2 + \overrightarrow{\nabla}^2 \right] \right\}$$

- NLO term is to be taken in **perturbative way**.

- For energies,

$$\Delta E = \langle \psi_{LO} | V_{NLO} | \psi_{LO} \rangle$$

- For scattering amplitude - distorted wave Born approximation,

$$\Delta f_k = -\frac{m}{k^2} \int dr \psi_{LO}^2 V_{NLO}$$

$$f_k \approx \frac{1}{-a^{-1} - ik} \left( 1 - \frac{1}{-a^{-1} - ik} \left[ \frac{1}{2} r_{\text{eff}} k^2 + \frac{\delta a}{a^2} \right] \right)$$

# Next to Leading Order

- At NLO, the LO term is iterated and 2-derivatives term is added:

$$V_{NLO}(r) = \frac{4\pi\hbar^2}{m\Lambda} \delta_\Lambda(r) \left\{ C_0^{(1)}(\Lambda) + C_2^{(1)}(\Lambda) \left[ \overleftarrow{\nabla}^2 + \overrightarrow{\nabla}^2 \right] \right\}$$

- NLO term is to be taken in **perturbative way**.
- For energies,

$$\Delta E = \langle \psi_{LO} | V_{NLO} | \psi_{LO} \rangle$$

- For scattering amplitude - distorted wave Born approximation,

$$\Delta f_k = -\frac{m}{k^2} \int dr \psi_{LO}^2 V_{NLO}$$

- $$f_k \approx \frac{1}{-a^{-1} - ik} \left( 1 - \frac{1}{-a^{-1} - ik} \left[ \frac{1}{2} r_{\text{eff}} k^2 + \frac{\delta a}{a^2} \right] \right)$$

# Next to Leading Order

- At NLO, the LO term is iterated and 2-derivatives term is added:

$$V_{NLO}(r) = \frac{4\pi\hbar^2}{m\Lambda} \delta_\Lambda(\mathbf{r}) \left\{ C_0^{(1)}(\Lambda) + C_2^{(1)}(\Lambda) \left[ \overleftarrow{\nabla}^2 + \overrightarrow{\nabla}^2 \right] \right\}$$

- NLO term is to be taken in **perturbative way**.

- For energies,

$$\Delta E = \langle \psi_{LO} | V_{NLO} | \psi_{LO} \rangle$$

- For scattering amplitude - distorted wave Born approximation,

$$\Delta f_k = -\frac{m}{k^2} \int dr \psi_{LO}^2 V_{NLO}$$

$$f_k \approx \frac{1}{-a^{-1} - ik} \left( 1 - \frac{1}{-a^{-1} - ik} \left[ \frac{1}{2} r_{\text{eff}} k^2 + \frac{\delta a}{a^2} \right] \right)$$

# Next to Leading Order

- At NLO, the LO term is iterated and 2-derivatives term is added:

$$V_{NLO}(r) = \frac{4\pi\hbar^2}{m\Lambda} \delta_\Lambda(\mathbf{r}) \left\{ C_0^{(1)}(\Lambda) + C_2^{(1)}(\Lambda) \left[ \overleftarrow{\nabla}^2 + \overrightarrow{\nabla}^2 \right] \right\}$$

- NLO term is to be taken in **perturbative way**.

- For energies,

$$\Delta E = \langle \psi_{LO} | V_{NLO} | \psi_{LO} \rangle$$

- For scattering amplitude - distorted wave Born approximation,

$$\Delta f_k = -\frac{m}{k^2} \int dr \psi_{LO}^2 V_{NLO}$$

- 

$$f_k \approx \frac{1}{-a^{-1} - ik} \left( 1 - \frac{1}{-a^{-1} - ik} \left[ \frac{1}{2} r_{\text{eff}} k^2 + \frac{\delta a}{a^2} \right] \right)$$

# Two-boson system

- From **TTY** potential we have  $a = 189a_0$  and  $B_2 = 1.31$  mK.  
LO is fitted to  $a = 189a_0$ , NLO to  $r_{eff} = 14.2a_0$ .
- $r_{eff}/a \approx 8\%$ ;  $(r_{eff}/a)^2 \approx 0.5\%$

- Fitting to powers of  $Q_2/\Lambda$ , ( $B_A = AQ_A^2/2m$ ) we extract  
 $B_2^{LO} = 1.21$  mK,  $B_2^{NLO} = 1.30$  mK,  $B_2^{TTY} = 1.31$  mK

Tang, Toennies & Yiu, PRL 74, 1546 (1995)

# Two-boson system

- From **TTY** potential we have  $a = 189a_0$  and  $B_2 = 1.31$  mK.  
LO is fitted to  $a = 189a_0$ , NLO to  $r_{eff} = 14.2a_0$ .
- $r_{eff}/a \approx 8\%$ ;  $(r_{eff}/a)^2 \approx 0.5\%$

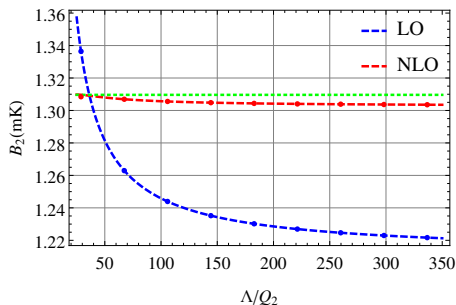
- Fitting to powers of  $Q_2/\Lambda$ , ( $B_A = A Q_A^2/2m$ ) we extract  
 $B_2^{LO} = 1.21$  mK,  $B_2^{NLO} = 1.30$  mK,  $B_2^{TTY} = 1.31$  mK

Tang, Toennies & Yiu, PRL 74, 1546 (1995)



# Two-boson system

- From **TTY** potential we have  $a = 189a_0$  and  $B_2 = 1.31$  mK.  
LO is fitted to  $a = 189a_0$ , NLO to  $r_{\text{eff}} = 14.2a_0$ .
- $r_{\text{eff}}/a \approx 8\%$ ;  $(r_{\text{eff}}/a)^2 \approx 0.5\%$



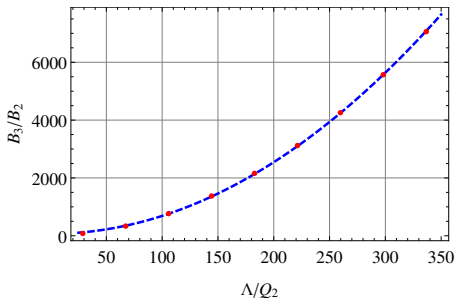
- Fitting to powers of  $Q_2/\Lambda$ , ( $B_A = A Q_A^2/2m$ ) we extract  
 $B_2^{LO} = 1.21$  mK,  $B_2^{NLO} = 1.30$  mK,  $B_2^{TTY} = 1.31$  mK

Tang, Toennies & Yiu, PRL 74, 1546 (1995)

# Three-boson system

Trying to calculate the trimer binding energy we get the **Thomas collapse**:

$$B_3 \propto \frac{\hbar\Lambda^2}{m}$$



- To stabilize the system, a 3-body counter term must be introduced at LO,

$$V_{LO}^{3N} = \frac{4\pi\hbar^2}{m\Lambda^4} D^{(0)} \sum_{i<j<k} \sum_{cyc} \delta_{\Lambda}(\mathbf{r}_{ij}) \delta_{\Lambda}(\mathbf{r}_{jk}),$$

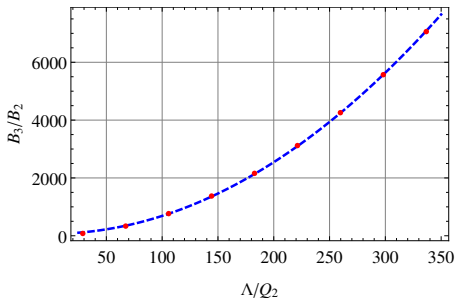
$\Lambda_*$  is a new momentum scale,  $D^{(0)} = f(a\Lambda, \Lambda/\Lambda_*)$

- $D^{(0)}$  is fixed by another observable.

# Three-boson system

Trying to calculate the trimer binding energy we get the **Thomas collapse**:

$$B_3 \propto \frac{\hbar\Lambda^2}{m}$$



- To stabilize the system, a 3-body counter term must be introduced at **LO**,

$$V_{LO}^{3N} = \frac{4\pi\hbar^2}{m\Lambda^4} D^{(0)} \sum_{i < j < k} \sum_{cyc} \delta_{\Lambda}(\mathbf{r}_{ij}) \delta_{\Lambda}(\mathbf{r}_{jk}),$$

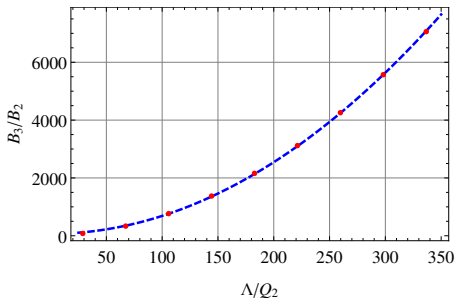
$\Lambda_*$  is a new momentum scale,  $D^{(0)} = f(a\Lambda, \Lambda/\Lambda_*)$

- $D^{(0)}$  is fixed by another observable.

# Three-boson system

Trying to calculate the trimer binding energy we get the **Thomas collapse**:

$$B_3 \propto \frac{\hbar\Lambda^2}{m}$$



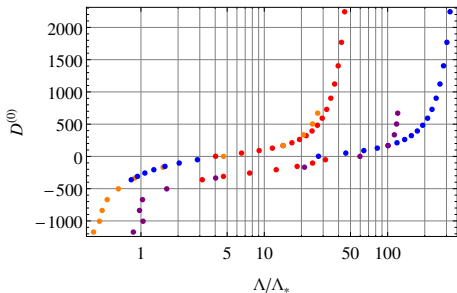
- To stabilize the system, a 3-body counter term must be introduced at **LO**,

$$V_{LO}^{3N} = \frac{4\pi\hbar^2}{m\Lambda^4} D^{(0)} \sum_{i < j < k} \sum_{cyc} \delta_{\Lambda}(\mathbf{r}_{ij}) \delta_{\Lambda}(\mathbf{r}_{jk}),$$

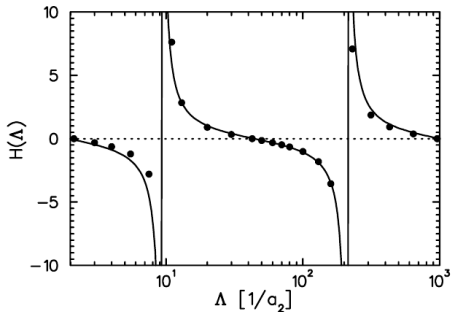
$\Lambda_*$  is a new momentum scale,  $D^{(0)} = f(a\Lambda, \Lambda/\Lambda_*)$

- $D^{(0)}$  is fixed by another observable.

# Three-boson system



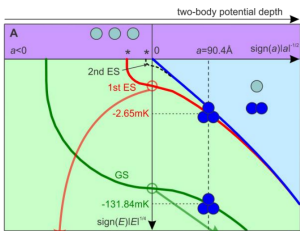
$\Lambda_3 = \Lambda_2$ , local, smooth cutoff



$\Lambda_3 \ll \Lambda_2$ , non-local, sharp cutoff  
P. F. Bedaque, H.W. Hammer, and U. van Kolck  
Phys. Rev. Lett. **82** 463 (1999).

$\Lambda_3 = \Lambda_2$ , non-local, smooth cutoff  
R.F. Mohr *et al.*, Ann. Phys. **321**, 225 (2006).

# Atom-dimer scattering



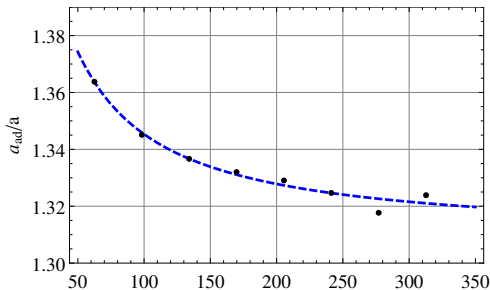
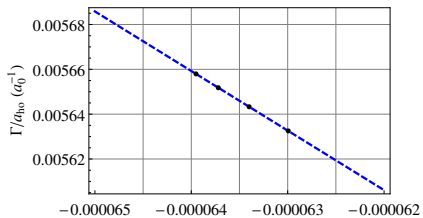
LO and NLO are fitted to  $B_3^* = 1.74B_2$ .

The atom-dimer scattering length is calculated in a trap, using

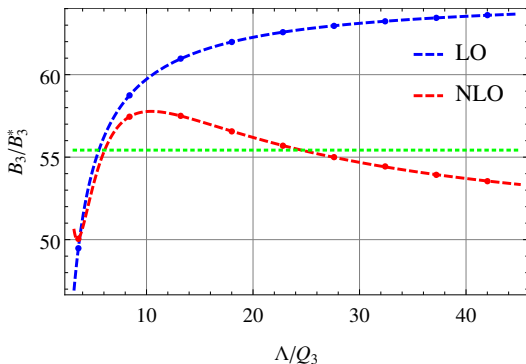
$$\frac{1}{\sqrt{2}} \frac{\Gamma[(1-\eta)/4]}{\Gamma[(3-\eta)/4]} = \frac{a_{ad}/a_{ho}}{1 - a_{ad}r_{ad}\eta/(4a_{ho}^2)}$$

$$a_{ho} = \sqrt{\hbar/(2\mu\omega)} \quad \mu = 2m/3 \quad \eta = 2(E_3 - E_2)/(\hbar\omega)$$

Suzuki *et al.*, PRA **80**, 033601 (2009); Stetcu *et al.*, Ann. Phys. **325**, 1644 (2010).



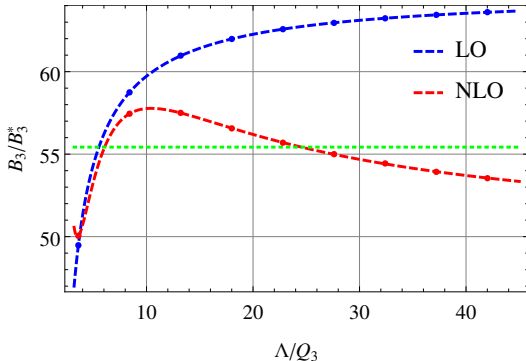
# Trimer ground state



- Fitting to powers of  $Q_3/\Lambda$ ,  $B_3(\Lambda) = B_{3\infty}(1 + \alpha \frac{Q_3}{\Lambda} + \beta \left(\frac{Q_3}{\Lambda}\right)^2 + \gamma \left(\frac{Q_3}{\Lambda}\right)^3)$

$B_{3\infty}/B_3^*$	$\alpha$	$\beta$	$\gamma$
64.83	-0.78	-	-
64.80	-0.78	-0.05	-
64.80	-0.80	0.36	-1.95

# Trimer ground state



- Fitting to powers of  $Q_3/\Lambda$ ,  $B_3(\Lambda) = B_{3\infty}(1 + \alpha \frac{Q_3}{\Lambda} + \beta (\frac{Q_3}{\Lambda})^2 + \gamma (\frac{Q_3}{\Lambda})^3)$  we extract

$$B_3^{LO}/B_3^* = 64.8, \quad B_3^{NLO}/B_3^* = 49.8, \quad B_3^{TTY}/B_3^* = 55.4$$

$$Q_3 r_{eff} \approx 0.6 \gg Q_2 r_{eff} \approx 0.08$$



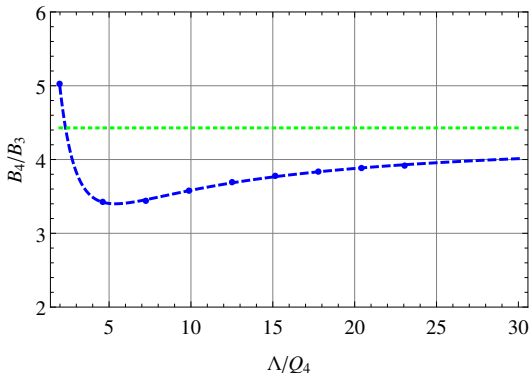
# Four-boson system

Are more terms needed to stabilize heavier systems?

$$B_4^{LO} / B_3 = 4.2(1), B_4^{TTY} / B_3 = 4.43$$

# Four-boson system

Are more terms needed to stabilize heavier systems?

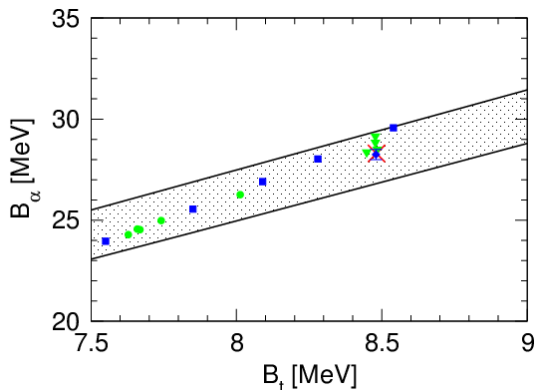


$$B_4^{LO}/B_3 = 4.2(1), B_4^{TTY}/B_3 = 4.43$$

# Tjon line

Another evidence is the **Tjon line**, the correlation between the binding energies of the triton and the  $\alpha$ -particle.

J.A. Tjon, Phys. Lett. B **56**, 217 (1975).

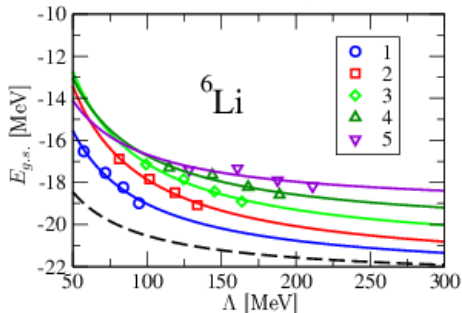


L. Platter, H.-W. Hammer, U.-G. Meissner, Phys. Lett. B **607**, 254 (2005).

## 5- and 6- boson system

Are more terms needed to stabilize heavier systems?

For nucleons,

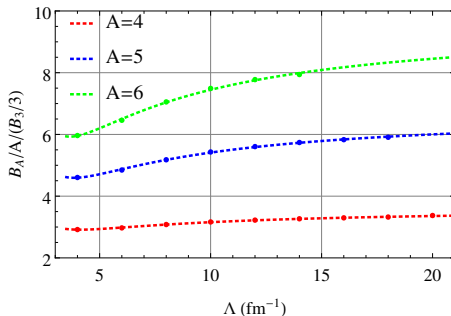


I. Stetcu, B.R. Barrett, and U. van Kolck, Phys. Lett. B **653**, 358 (2007).

A	x	x <sup>2</sup>	x <sup>3</sup>	Ref. [1]
4	9.63	10.22	9.94	9.88
5	17.51	19.08	18.40	18.05
6	24.37	27.46	24.98	26.28

# 5- and 6- boson system

Are more terms needed to stabilize heavier systems?

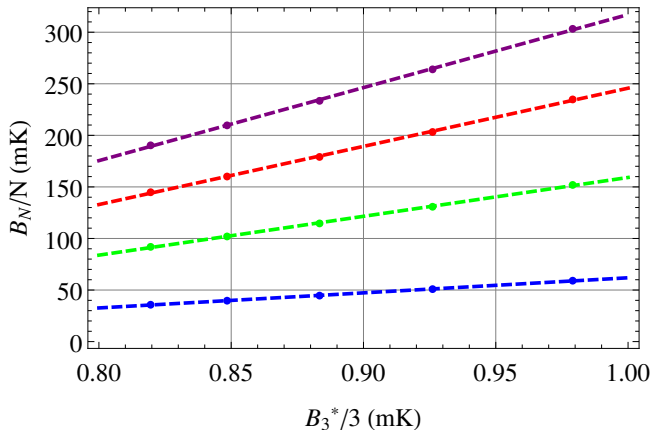


A	$x$	$x^2$	$x^3$	Ref. [1]
4	9.63	10.22	9.94	9.88
5	17.51	19.08	18.40	18.05
6	24.37	27.46	24.98	26.28

[1] J. von Stecher, J. Phys. B: At. Mol. Opt. Phys. **43**, 101002 (2010).

# Generalized Tjon-lines

Correlation between  $B_3^*$  to  $B_3$ ,  $B_4$ ,  $B_5$ , and  $B_6$ :



...Therefore, no 4, 5 or 6-body terms are needed at LO.

- At LO, the 2-body potential reads,

$$V_{LO} = a_1 + a_2 \sigma_i \cdot \sigma_j + a_3 \tau_i \cdot \tau_j + a_4 \sigma_i \cdot \sigma_j \tau_i \cdot \tau_j$$

- using the fermionic symmetry,

$$V_{LO} = C_S \hat{P}_S + C_T \hat{P}_T$$

where  $\hat{P}_\alpha$  is projection operator on channel  $\alpha$

- The 2-body LECs are fitted to the deuteron binding energy and the singlet  $^1S_0$   $np$  scattering length.
- The 3-body LEC is fitted to the triton binding energy
- No Coulomb interaction (should be NLO. See Koenig's talk next week).

- At LO, the 2-body potential reads,

$$V_{LO} = a_1 + a_2 \sigma_i \cdot \sigma_j + a_3 \tau_i \cdot \tau_j + a_4 \sigma_i \cdot \sigma_j \tau_i \cdot \tau_j$$

- using the fermionic symmetry,

$$V_{LO} = C_S \hat{P}_S + C_T \hat{P}_T$$

where  $\hat{P}_\alpha$  is projection operator on channel  $\alpha$

- The 2-body LECs are fitted to the deuteron binding energy and the singlet  $^1S_0$   $np$  scattering length.
- The 3-body LEC is fitted to the triton binding energy
- No Coulomb interaction (should be NLO. See Koenig's talk next week).



- At LO, the 2-body potential reads,

$$V_{LO} = a_1 + a_2 \sigma_i \cdot \sigma_j + a_3 \tau_i \cdot \tau_j + a_4 \sigma_i \cdot \sigma_j \tau_i \cdot \tau_j$$

- using the fermionic symmetry,

$$V_{LO} = C_S \hat{P}_S + C_T \hat{P}_T$$

where  $\hat{P}_\alpha$  is projection operator on channel  $\alpha$

- The 2-body LECs are fitted to the deuteron binding energy and the singlet  $^1S_0$   $np$  scattering length.
- The 3-body LEC is fitted to the triton binding energy
- No Coulomb interaction (should be NLO. See Koenig's talk next week).

- At LO, the 2-body potential reads,

$$V_{LO} = a_1 + a_2 \sigma_i \cdot \sigma_j + a_3 \tau_i \cdot \tau_j + a_4 \sigma_i \cdot \sigma_j \tau_i \cdot \tau_j$$

- using the fermionic symmetry,

$$V_{LO} = C_S \hat{P}_S + C_T \hat{P}_T$$

where  $\hat{P}_\alpha$  is projection operator on channel  $\alpha$

- The 2-body LECs are fitted to the deuteron binding energy and the singlet  $^1S_0$   $np$  scattering length.
- The 3-body LEC is fitted to the triton binding energy
- No Coulomb interaction (should be NLO. See Koenig's talk next week).

- At LO, the 2-body potential reads,

$$V_{LO} = a_1 + a_2 \sigma_i \cdot \sigma_j + a_3 \tau_i \cdot \tau_j + a_4 \sigma_i \cdot \sigma_j \tau_i \cdot \tau_j$$

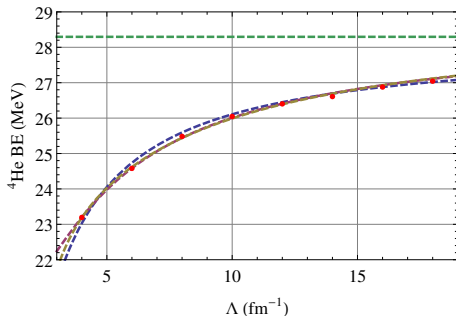
- using the fermionic symmetry,

$$V_{LO} = C_S \hat{P}_S + C_T \hat{P}_T$$

where  $\hat{P}_\alpha$  is projection operator on channel  $\alpha$

- The 2-body LECs are fitted to the deuteron binding energy and the singlet  $^1S_0$   $np$  scattering length.
- The 3-body LEC is fitted to the triton binding energy
- No Coulomb interaction (should be NLO. See Koenig's talk next week).

# $\alpha$ - ${}^4\text{He}$ nuclei



$$B_4(\Lambda) = B_\infty \left[ 1 + \alpha \frac{Q_4}{\Lambda} + \beta \left( \frac{Q_4}{\Lambda} \right)^2 + \gamma \left( \frac{Q_4}{\Lambda} \right)^3 + \dots \right]$$

$B_\infty$	$\alpha$	$\beta$	$\gamma$
28.16	-1.25	-	-
28.66	-1.77	3.17	-
28.88	-2.12	8.07	-19.21

# Conclusion

- A pionless EFT was constructed for few-body systems.
- The  ${}^4\text{He}$  atomic system was studied, and our EFT fits nicely the known results.
- The convergence of pionless EFT for  $A = 4, 5$  and  $6$  was studied.
- Generalized Tjon-lines were introduced, showing that at LO no 4,5 or 6-body term is needed.
- Similar results was shown for atomic nucleus.

# Conclusion

- A pionless EFT was constructed for few-body systems.
- The  $^4\text{He}$  atomic system was studied, and our EFT fits nicely the known results.
- The convergence of pionless EFT for  $A = 4, 5$  and  $6$  was studied.
- Generalized Tjon-lines were introduced, showing that at LO no 4,5 or 6-body term is needed.
- Similar results was shown for atomic nucleus.

# Conclusion

- A pionless EFT was constructed for few-body systems.
- The  $^4\text{He}$  atomic system was studied, and our EFT fits nicely the known results.
- The convergence of pionless EFT for  $A = 4, 5$  and  $6$  was studied.
- Generalized Tjon-lines were introduced, showing that at LO no 4,5 or 6-body term is needed.
- Similar results was shown for atomic nucleus.

# Conclusion

- A pionless EFT was constructed for few-body systems.
- The  ${}^4\text{He}$  atomic system was studied, and our EFT fits nicely the known results.
- The convergence of pionless EFT for  $A = 4, 5$  and  $6$  was studied.
- Generalized Tjon-lines were introduced, showing that at LO no 4,5 or 6-body term is needed.
- Similar results was shown for atomic nucleus.



# Conclusion

- A pionless EFT was constructed for few-body systems.
- The  $^4\text{He}$  atomic system was studied, and our EFT fits nicely the known results.
- The convergence of pionless EFT for  $A = 4, 5$  and  $6$  was studied.
- Generalized Tjon-lines were introduced, showing that at LO no 4,5 or 6-body term is needed.
- Similar results was shown for atomic nucleus.