

Effective Field Theory for Few-Boson Systems

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EMMI Rapid Reaction Task Force:

The systematic treatment of the Coulomb interaction in few-body systems

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Motivation

- Pionless EFT is very effective for low-energy physics in the few-nucleon case.
- A 3-body interaction is needed at LO, but no 4-body interaction is needed.
- This holds also for the few-boson system.
- Are higher-body forces needed at LO to describe systems with more bodies?
- What is the regime of validity of the EFT as the number of particles increase?

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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_2 .

$$\lim_{k \rightarrow 0} k \cot \delta(k) = -\frac{1}{a_2} + \frac{1}{2}r_2k^2$$

From Sakurai's book

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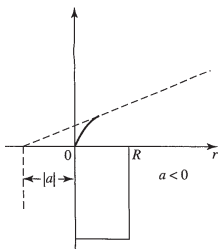
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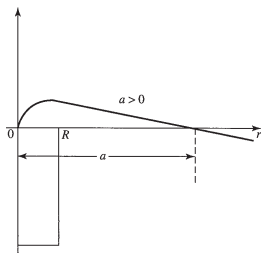
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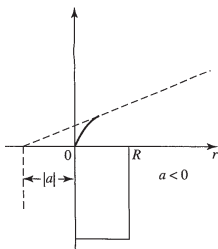


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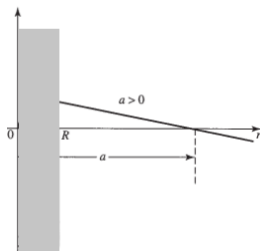
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Universality

- Generally, $a_2 \approx r_2 \approx R$.
Universal systems are fine-tuned to get $a_2 \gg r_2, R$.
- Corrections to universal theory are of order of r_2/a_2 and R/a_2 .
- For $a_2 > 0$, we have universal dimer with energy $E = -\hbar^2/ma_2^2$.
- ^4He Atoms: $a_2 \approx 170.9a_0$, (a_0 = the Bohr radius), is much larger than its van der Waals radius, $r_{vdW} \approx 9.5a_0$.
- 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,

$$a_2(B) = a_{bg} \left(1 + \frac{\Delta}{B - B_0} \right)$$

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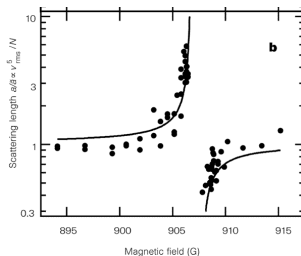
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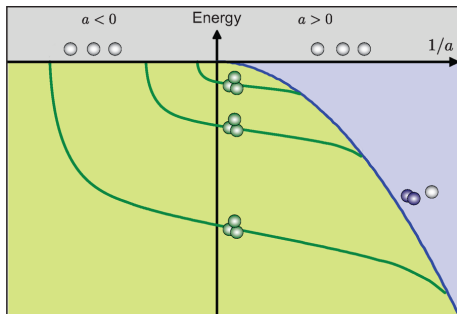
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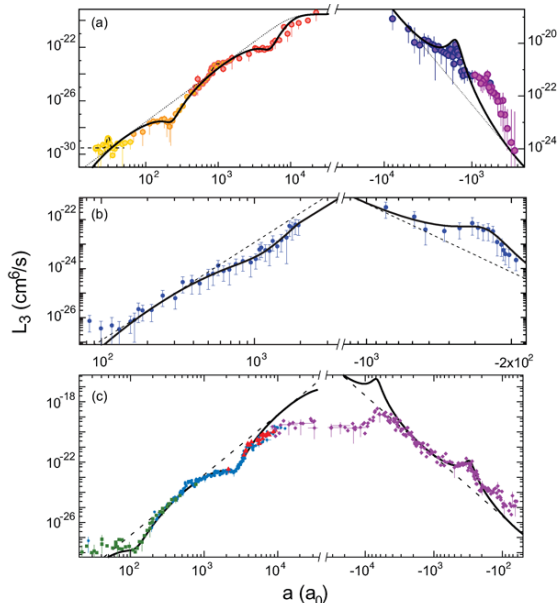
Efimov Physics

- The unitary limit: $E_2 = 0$, $a_2 \rightarrow \infty$.
- 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}K
M. Zaccanti *et al.*,
Nature Phys. **5**, 586 (2009).

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

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

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

Effective Field Theory (EFT)

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- For example, nuclear structure involves energies that are much smaller than the typical QCD mass scale, $M_{QCD} \approx 1$ 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.

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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 !

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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; the NLO term is treated as perturbation.

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Stochastic Variational Method I

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

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

η = 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 bosons and fermions. Arbitrary angular momentum and parity, as well as spin and isospin, can be introduced.

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Stochastic Variational Method II

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

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

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

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

Stochastic Variational Method II

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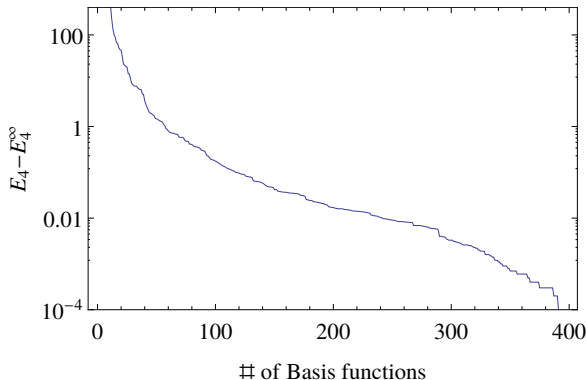
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- To find the best A_i , we use the **Stochastic Variational Method (SVM)**:
- We add basis function one by one, or try to replace an exist basis function by a new one.
- We choose **randomly** the matrix A_i element by element, trying to minimize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

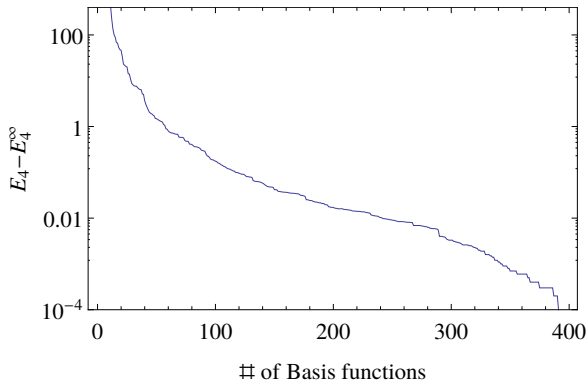
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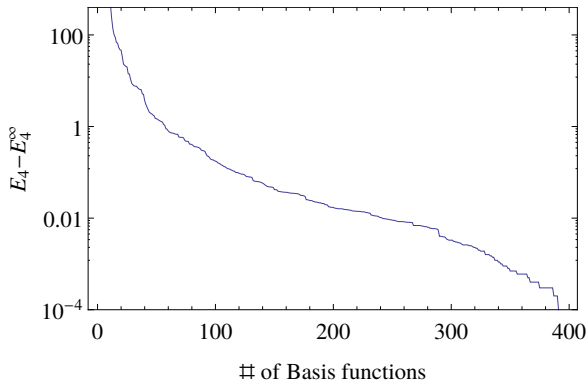
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Regularization I: non local potential

- At LO, we have only contact interaction,

$$V(r_{ij}) = \tilde{C}^{(0)} \delta(r_{ij}).$$

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

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and in momentum space,

$$\frac{p^2}{m} \phi(p) + \tilde{C}^{(0)} \int \frac{d^3 p'}{(2\pi)^3} \phi(p') = -B_2 \phi(p)$$

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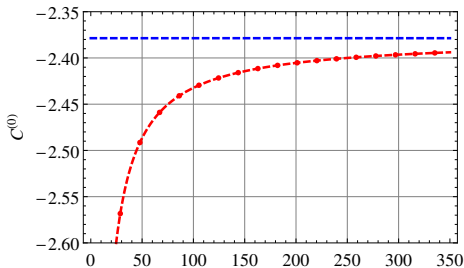
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- At NLO, the LO term is iterated and 2-derivatives term is added:

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

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

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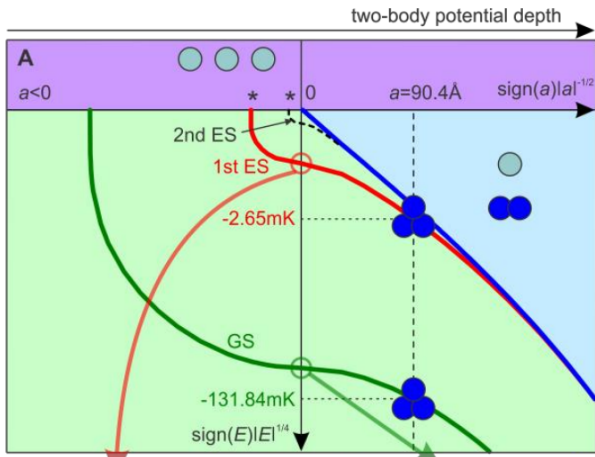
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^4He Atoms

- To be definite, we focus on ^4He atomic systems.



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

^4He Atoms

Length scales (in Å) for the ^4He atoms:

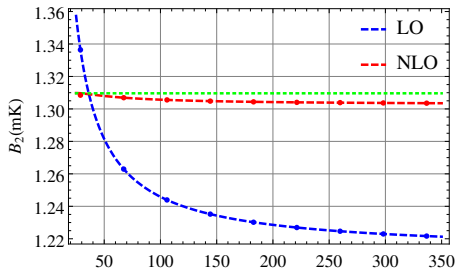
	LM2M2	TTY	PCKLJS
a_2	100.23	100.01	90.42(92)
r_2	7.326	7.329	7.27
r_{vdW}	5.378	5.378	5.378

Binding energies (in mK) of ^4He clusters:

	LM2M2	TTY	PCKLJS	experiment
B_2	1.3094	1.3096	1.6154	$1.3_{-0.19}^{+0.25}$; 1.76(15)
B_3^*	2.2779	2.2761	2.6502	
$B_3^* - B_2$	0.9685	0.9665	1.0348	0.98(2)
B_3	126.50	126.16	131.84	
B_4^*	127.42		132.70	
B_4	559.22		573.90	

Two-boson system

- Breakdown scale $\sim r_{\text{vdw}} \sim r_2/2 \sim 7a_0$.



$$B_2 = B_{LO} \left[1 + \mathcal{O} \left(\frac{Q_2}{M}, \frac{Q_2}{\Lambda} \right) \right]; B_2(\Lambda) = B_{2\infty} \left[1 + \alpha \frac{Q_2}{\Lambda} + \beta \left(\frac{Q_2}{\Lambda} \right)^2 + \gamma \left(\frac{Q_2}{\Lambda} \right)^3 \right]$$

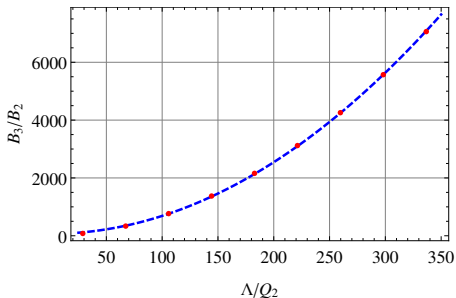
$B_{2\infty}(mK)$	α	β	γ
1.21	2.86	—	—
1.21	2.76	5.84	—
1.21	2.76	5.52	12.80
1.31			

$$\alpha Q_2 r_2 / 2 \approx 10\%$$

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,

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

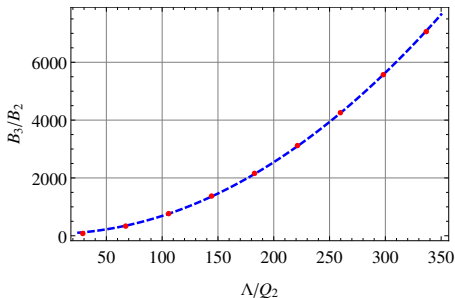
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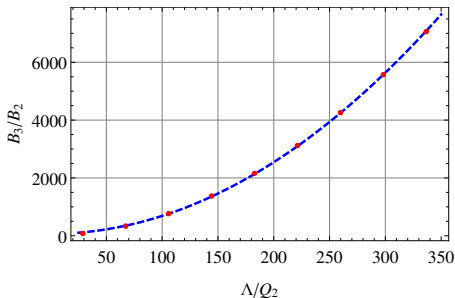
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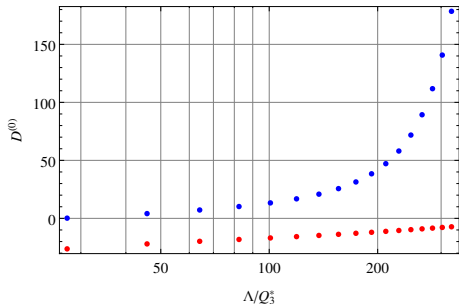
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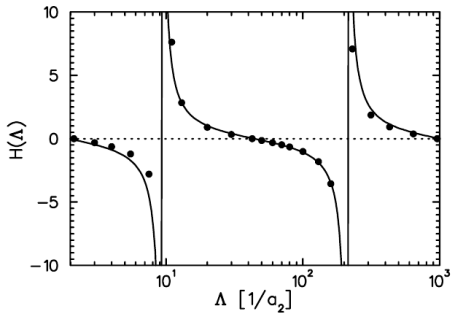
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$\Lambda_3 = \Lambda_2$, local, smooth cutoff



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

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Atom-dimer scattering

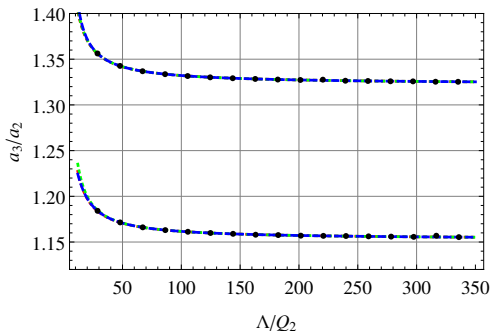
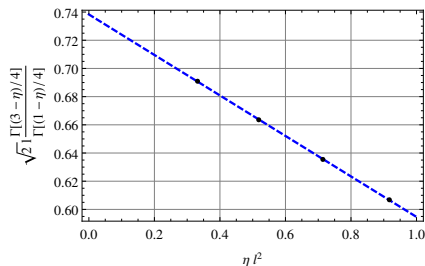
3-body LEC is fitted to B_3^* .

a_3 , the atom-dimer scattering length is calculated in a trap, using

$$\sqrt{2}l \frac{\Gamma[(3-\eta)/4]}{\Gamma[(1-\eta)/4]} \simeq \frac{a_2}{a_3} \left(1 - \frac{a_3 r_3}{4a_2^2} \eta l^2 \right),$$

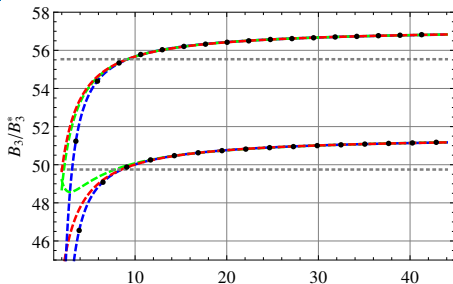
$$\eta = 2(E_3 - E_2)/\omega, l = a_2/a_{ho}, a_{ho} = 1/\sqrt{2\mu\omega}, \mu \simeq 2m/3$$

Busch *et al.*, Found. Phys. **28** 549 (1998); Stetcu *et al.*, Ann. Phys. **325**, 1644 (2010).



$$a_2 = 1.151(2) \text{ (LM2M2)}$$

Trimer ground state



$$B_3(\Lambda) = B_{3\infty} \left[1 + \alpha \frac{Q_3}{\Lambda} + \beta \left(\frac{Q_3}{\Lambda} \right)^2 + \gamma \left(\frac{Q_3}{\Lambda} \right)^3 \right]$$

LM2M2

PCKLJS

$B_3(\infty)/B_3^*$	α	β	γ
57.22	-0.26	—	—
57.23	-0.25	0.09	—
57.21	-0.26	0.20	-2.04
55.53			

$B_3(\infty)/B_3^*$	α	β	γ
51.51	-0.29	—	—
51.56	-0.34	0.49	—
51.52	-0.31	0.47	-2.80
49.75			

$$\alpha Q_3 r_2 / 2 \approx 10\%$$

Four-boson system

Are more terms needed to stabilize heavier systems?

LM2M2

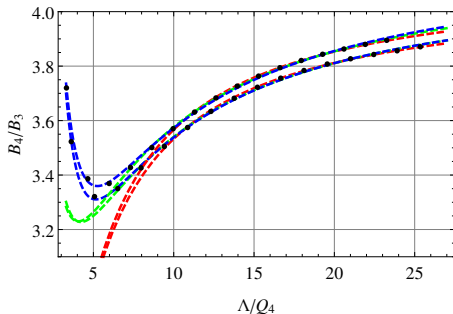
$B_4/B_3(\infty)$	α	β	γ
4.14	-1.38	—	—
4.21	-1.90	3.86	—
4.24	-2.02	4.07	4.31
4.42			

PCKLJS

$B_4/B_3(\infty)$	α	β	γ
4.09	-1.36	—	—
4.17	-1.90	4.02	—
4.16	-1.80	2.32	8.00
4.35			

Four-boson system

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LM2M2

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PCKLJS

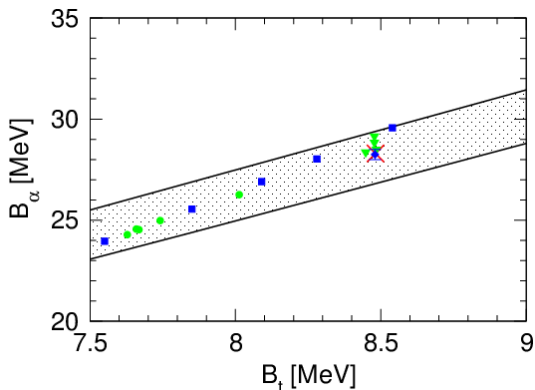
$B_4/B_3(\infty)$	α	β	γ
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4.17	-1.90	4.02	—
4.16	-1.80	2.32	8.00
4.35			

$$Q_4 r_2 / 2 \approx 55\%$$

Tjon line

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

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

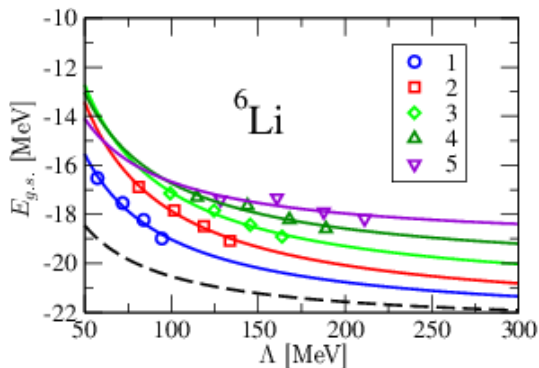


L. Platter, H.-W. Hammer, U.-G. Meissner, Phys. Lett. B **607**, 254 (2005).

Heavier system

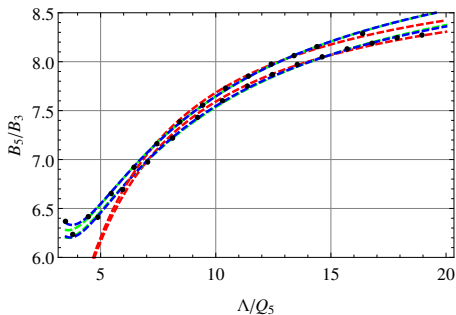
Are more terms needed to stabilize heavier systems?

For nucleons,



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

Five-boson system



LM2M2

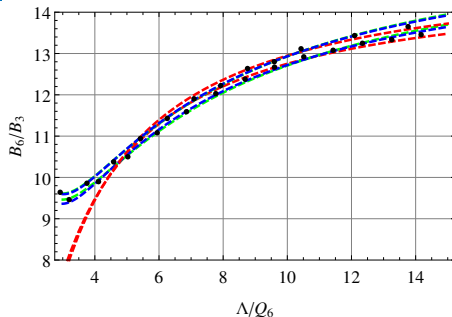
$B_5/B_3(\infty)$	α	β	γ
9.16	-1.61	—	—
9.61	-2.49	4.48	—
9.60	-2.46	4.06	1.38
10.33			

PCKLJS

$B_5/B_3(\infty)$	α	β	γ
9.02	-1.58	—	—
9.41	-2.43	4.31	—
9.32	-2.20	2.63	3.47

$$Q_5 r_2 / 2 \approx 75\%$$

Six-boson system



LM2M2

$B_6/B_3(\infty)$	α	β	γ
15.3	-1.53	—	—
16.4	-2.48	3.73	—
16.3	-2.35	2.99	1.15
18.41			

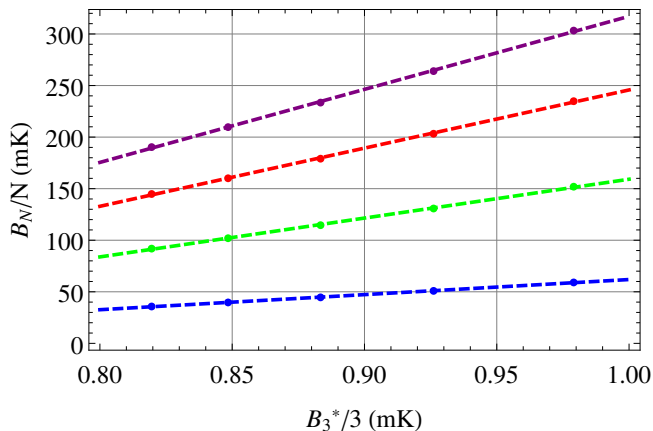
PCKLJS

$B_6/B_3(\infty)$	α	β	γ
14.9	-1.46	—	—
16.1	-2.47	3.69	—
15.7	-2.14	1.92	2.52

$$Q_6 r_2 / 2 \approx 85\%$$

Generalized Tjon-lines

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



Conclusion

- A pionless EFT was constructed for few-body systems.
- The ^4He 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.

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