



Nuclear Structure Contribution to the Lamb Shift of Muonic Deuterium in Pionless EFT

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Outline

- Introduction:
 - Nuclear radii and two-photon exchange (TPE) corrections
- Framework:
 - Pionless effective field theory (π EFT)
- In Practice:
 - Deuteron doubly virtual Compton scattering in π EFT
 - TPE corrections to the Lamb shift of μ D in π EFT
- Detour:
 - TPE corrections via sum rules supplemented by EFTs
- Summary&Outlook

Nuclear Radii from Muonic Atoms

- Lamb shift is sensitive to the finite size of the atomic nucleus

$$\Delta E_{2S}^{\text{f.s.}} = m_r^3 \alpha^4 r^2 / 12$$

- Extraction of the nuclear charge radius from atomic spectra is possible
- Muon is heavier \rightarrow closer to the nucleus \rightarrow much larger effect $\sim 10^7 \times$ electron
- Muonic atoms: better sensitivity to the radius, better precision
- Deuteron radius puzzle

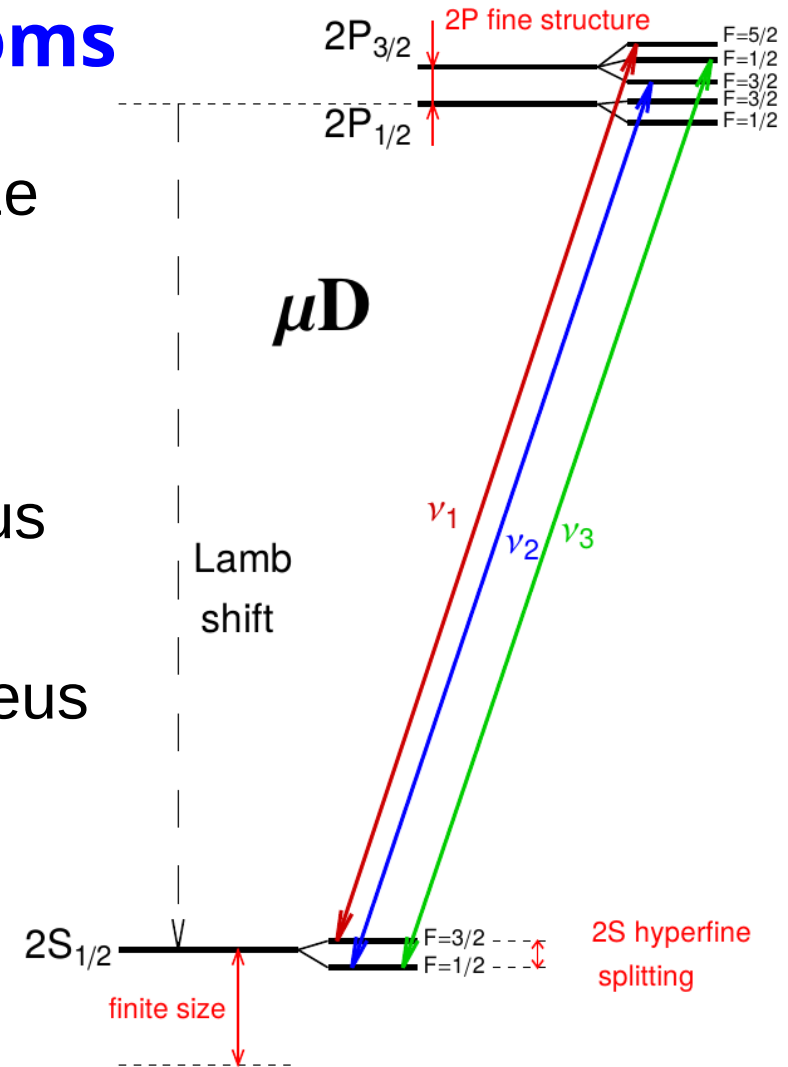
$$r_d(\mu\text{D}) = 2.12562(78) \text{ fm}$$

Pohl et al (2016)

$$r_d = 2.1424(21) \text{ fm}$$

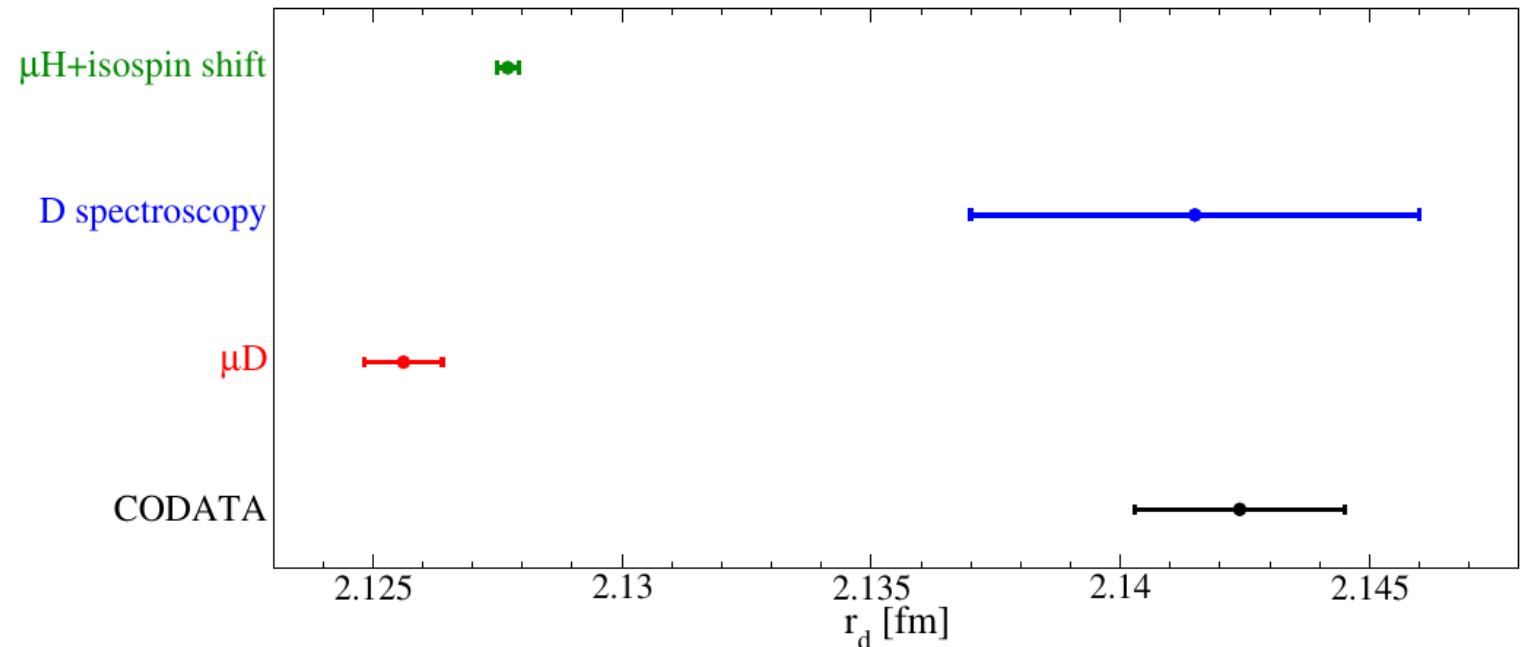
CODATA 2016

– 6..7 σ discrepancy



Deuteron Radius Puzzle

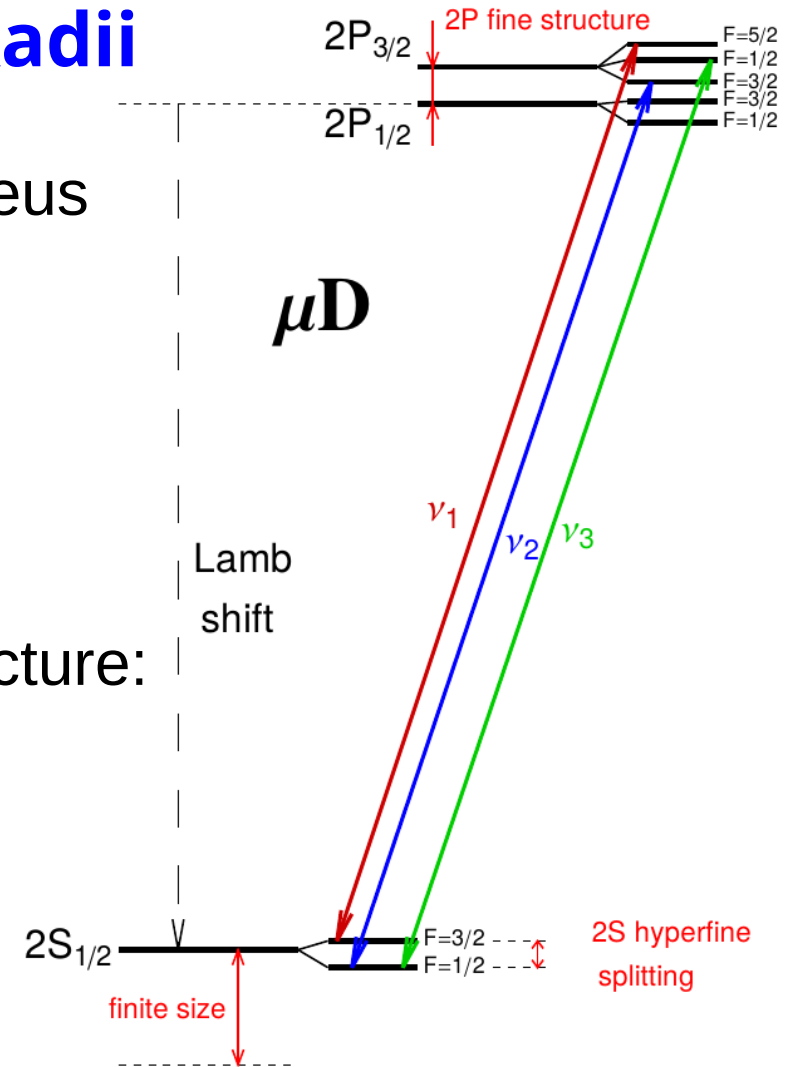
- Deuteron radius puzzle



- CODATA uses the isotopic shift to evaluate $r_d \rightarrow$ dependent on the proton radius puzzle
- One can extract r_d from (electronic) deuterium spectroscopy alone (with a bigger error) $\rightarrow 3.5\sigma$ discrepancy on its own

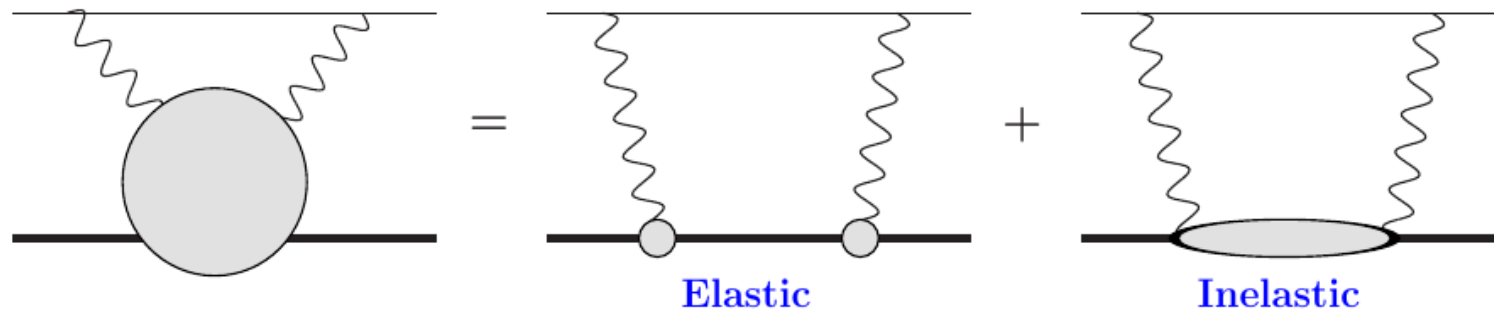
Nuclear Structure in Nuclear Radii

- Muon is heavier \rightarrow closer to the nucleus
- Better sensitivity to the radius!
- **Side effect:** better sensitivity to the detailed **structure of the nucleus**
- The leading effect of the nuclear structure:
 - two-photon exchange corrections
 - affect the extraction of nuclear radii
 - limit the theoretical accuracy



$$\Delta E_{LS}^{\text{theory}}(\mu\text{D}) = 228.7766(10) - 6.1103(3)r_d^2 + \Delta E_{\text{TPE}} \text{ meV}$$

Two-photon Exchange Corrections



- Elastic can be calculated in terms of e.m. deuteron FFs
- Ways to calculate inelastic
 - Hamiltonian approach (2nd order perturbation) Friar; Pachucki, Wienczek; Bacca, Barnea, Hernandez, Nevo Dinur, Ji
 - Sum rules (structure/response functions) Carlson, Gorshteyn, Vanderhaeghen
 - Nuclear VVCS amplitude
- We want to calculate the VVCS amplitude from pionless effective field theory – a systematically controlled approach
- I will also briefly discuss EFT input for sum rules

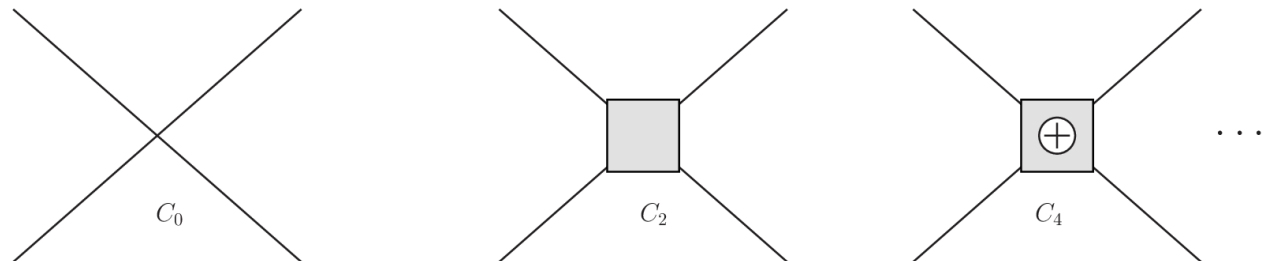
Pionless EFT

- Effective field theory of nucleon-nucleon interactions
- Works at very low energies/momenta $p \ll m_\pi$
- EFT of non-relativistic nucleons with contact interactions
 - Symmetries (isospin, Galilean, e.m. gauge, ...)
 - Lagrangians are ordered according to the counting

$$\mathcal{L} = N^\dagger \left[iD_0 + \frac{\vec{D}^2}{2M} \right] N + \hat{\mu} N^\dagger (\vec{\sigma} \cdot \vec{B}) N$$

Kaplan, Savage, Wise;
Chen, Rupak, Savage; ...

$$+ C_0 (N^\dagger P^i N_c) (N_c^\dagger P^i N) + \frac{1}{8} C_2 \left\{ (N^\dagger P^i N_c) \left(N_c^\dagger P^i \left[\overleftarrow{D}^2 + \overrightarrow{D}^2 - 2\overleftarrow{D} \cdot \overrightarrow{D} \right] N \right) + \text{H.c.} \right\}$$
$$+ \dots$$



Pionless EFT: Counting

- Nucleons are non-relativistic $\rightarrow E \simeq p^2/M$: energies are $O(p^2)$
- Loop integrals are $dE d^3p = O(p^5)$
- Nucleon propagators are $(E - p^2/2M)^{-1} = O(p^{-2})$
- Typical momenta $p \sim \gamma = \sqrt{ME_d} \simeq 45 \text{ MeV}$
- Expansion parameter $p/m_\pi \simeq \gamma/m_\pi \simeq 1/3$

$$\begin{aligned} \mathcal{L} = & N^\dagger \left[iD_0 + \frac{\vec{D}^2}{2M} \right] N + \hat{\mu} N^\dagger (\vec{\sigma} \cdot \vec{B}) N \\ & + C_0 (N^\dagger P^i N_c) (N_c^\dagger P^i N) + \frac{1}{8} C_2 \left\{ (N^\dagger P^i N_c) \left(N_c^\dagger P^i \left[\overleftarrow{D}^2 + \overrightarrow{D}^2 - 2\overleftarrow{D} \cdot \overrightarrow{D} \right] N \right) + \text{H.c.} \right\} \\ & + \dots \end{aligned}$$

- What is the counting for the coupling constants?
 - NN system has a low-lying bound/virtual state
 - Need to enhance constants: $C_{2n} = O(p^{-n-1})$

Pionless EFT: Resumming the LO

- $C_0 = O(p^{-1})$: leading-order loops need to be resummed

$$\mathcal{M}_{-1} = \underbrace{\begin{array}{c} \diagup \quad \diagdown \\ p^{-1} \end{array} + \begin{array}{c} \diagup \quad \text{loop} \quad \diagdown \\ p^{-1} \quad p^{-1} \end{array} + \begin{array}{c} \diagup \quad \text{loop} \quad \text{loop} \quad \diagdown \\ p^{-1} \quad p^{-1} \quad p^{-1} \end{array} + \dots}_{O(p^{-1})}$$

- Result: LO scattering amplitude reproduces the deuteron pole – the resummation is in fact a necessity!

$$\mathcal{M}_{-1} = \frac{4\pi}{M} \frac{1}{-\gamma - ik}$$

- Higher-order terms (effective range, ...) can be taken into account perturbatively, in the expansion in powers of p

Pionless EFT: Different Expansions Beyond LO

- Scattering amplitude in the deuteron channel

$$f(k) = \frac{1}{-\gamma - ik + \frac{1}{2}\rho_d(k^2 + \gamma^2) + w_2(k^2 + \gamma^2)^2 + \dots}$$

- Around the pole

$$f(k) = \frac{Z}{-\gamma - ik} + R(k) \quad \text{where } R(k) \text{ is regular at the pole and } Z = \frac{1}{1 - \gamma\rho_d}$$

- Reproduce the effective range at NLO (ρ -expansion)

$$f(k) = \frac{1}{(-\gamma - ik)[1 + \frac{1}{2}\rho_d(ik - \gamma) + \dots]} = \frac{1}{-\gamma - ik} \left[1 + \rho_d\gamma - \frac{1}{2}\rho_d(ik + \gamma) + \dots \right]$$

- The residue is reproduced perturbatively order-by-order

$$Z = 1 + (\gamma\rho_d) + (\gamma\rho_d)^2 + (\gamma\rho_d)^3 + \dots$$

Pionless EFT: Different Expansions Beyond LO

- Scattering amplitude in the deuteron channel

$$f(k) = \frac{1}{-\gamma - ik + \frac{1}{2}\rho_d(k^2 + \gamma^2) + w_2(k^2 + \gamma^2)^2 + \dots}$$

- Around the pole

$$f(k) = \frac{Z}{-\gamma - ik} + R(k) \quad \text{where } R(k) \text{ is regular at the pole and } Z = \frac{1}{1 - \gamma\rho_d}$$

- ... or reproduce the residue at NLO (Z-expansion)

$$f(k) = \frac{1}{(-\gamma - ik) \underbrace{\left[1 - \gamma\rho_d + \frac{1}{2}\rho_d(ik + \gamma) + \dots\right]}_{1/Z}} = \frac{Z}{-\gamma - ik} \left[1 - \frac{Z}{2}\rho_d(ik + \gamma) + \dots\right]$$

- Z expansion yields better convergence for low-energy observables (\rightarrow asymptotic normalisation of the deuteron wave function is correct already at NLO)

Phillips, Rupak, Savage

In Practice: Counting

- We want to perform a NNLO calculation of the Lamb shift

$$\Delta E_{2S} = \frac{\alpha}{2\pi^2 m_\mu} \phi_{2S}(0)^2 \int d^4 q \frac{f_L(-iq_0, Q^2) - 2(q_0^2/Q^2) f_T(-iq_0, Q^2)}{Q^2 \left(\frac{Q^4}{4m_\mu^2} + q_0^2 \right)}$$

$$T_{\text{VVCS}}^{\text{scalar}} = [\epsilon_0 \epsilon_0'^* f_L(\nu, Q^2) + \vec{\epsilon} \cdot \vec{\epsilon}'^* f_T(\nu, Q^2)] \vec{\epsilon}_d \cdot \vec{\epsilon}_d'^*$$

- Subtract the pole (elastic) pieces – calculated via emp. FFs
- The counting gives

$$f_L(\nu, Q^2) = p^{-2} + p^{-1} + \dots$$

$$f_T(\nu, Q^2) = p^0 + p^1 + \dots$$

- Together with the weighting in the integral, this means the transverse amplitude starts to contribute to LS only at N4LO

In Practice: Counting

- Relativistic corrections to one-nucleon vertex are counted at N4LO as well because of the $1/M^2$ suppression
- SD-mixing also starts only at N4LO
- The relevant Lagrangian up to NNLO [*]

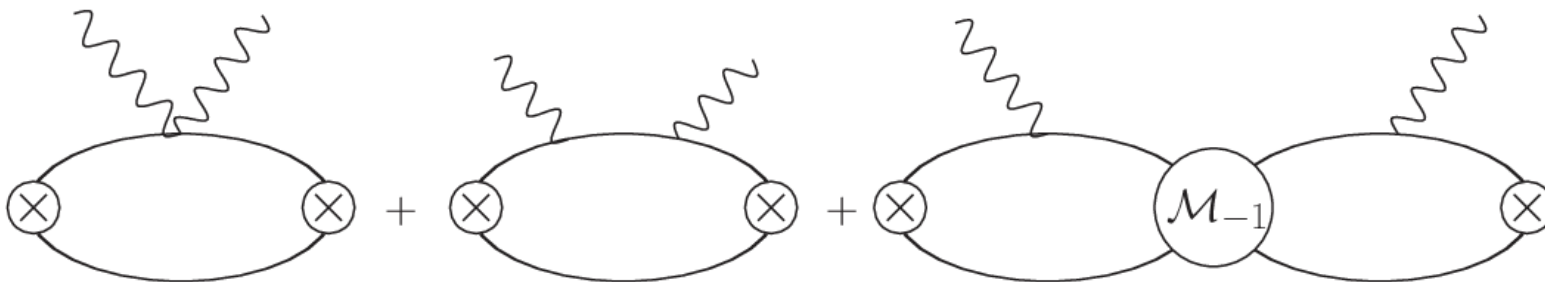
$$\begin{aligned}\mathcal{L} = & N^\dagger \left[iD_0 + \frac{\vec{D}^2}{2M} \right] N + \hat{\mu} N^\dagger (\vec{\sigma} \cdot \vec{B}) N \\ & + C_0 (N^\dagger P^i N_c) (N_c^\dagger P^i N) + \frac{1}{8} C_2 \left\{ (N^\dagger P^i N_c) \left(N_c^\dagger P^i \left[\overleftarrow{D}^2 + \vec{D}^2 - 2\overleftarrow{D} \cdot \vec{D} \right] N \right) + \text{H.c.} \right\} \\ & - \frac{C_4}{16} \left(N^\dagger \left[\overleftarrow{D}^2 + \vec{D}^2 - 2\overleftarrow{D} \cdot \vec{D} \right] P^i N_c \right) \left(N_c^\dagger P^i \left[\overleftarrow{D}^2 + \vec{D}^2 - 2\overleftarrow{D} \cdot \vec{D} \right] N \right) + \dots\end{aligned}$$

- In practice we calculate $f_T(\nu, Q^2)$ up to NLO completely
 - Add singlet NN interactions and two $NN\gamma$ contact terms
 - Cross check of the smallness of the $f_T(\nu, Q^2)$ contribution
 - Gauge invariance-conserving terms calculated up to NNLO

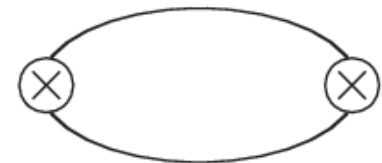
In Practice: Amplitude with Deuterons

- The reaction amplitude is given by the LSZ reduction

$$T = M \left[\frac{d\Sigma(E)}{dE} \Big|_{E=E_d} \right]^{-1}$$

$$M = \text{diagram 1} + \text{diagram 2} + \text{diagram 3} + \dots$$
The equation shows a series of Feynman diagrams for the reaction amplitude M. The first diagram is a loop with two vertices marked with an 'X' and a wavy line entering from the top. The second diagram is similar but with two wavy lines entering from the top. The third diagram consists of two loops connected by a central circle labeled M_{-1}, with wavy lines entering from the top of each loop. The series continues with an ellipsis.

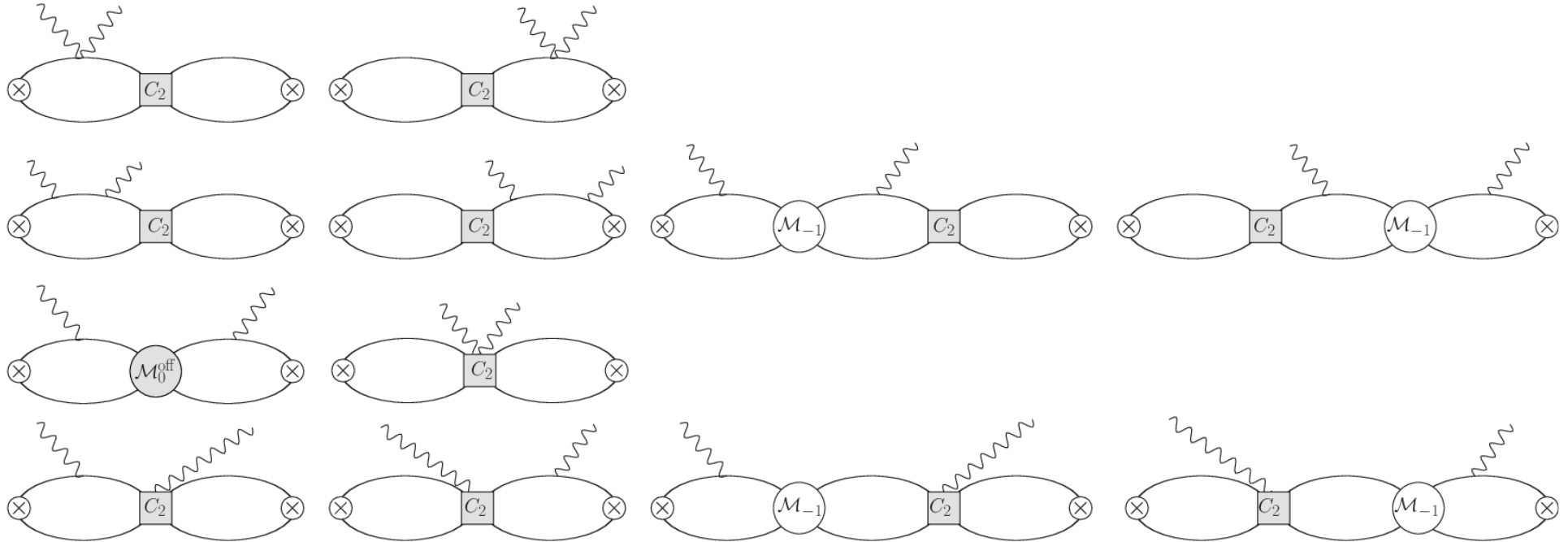
– irreducible VVCS graphs (here full LO in f_L ; crossed not shown)

$$\Sigma = \text{diagram 1} + \dots$$
The equation shows the first diagram of the series for the deuteron self-energy Sigma, which is a loop with two vertices marked with an 'X'. The series continues with an ellipsis.

– deuteron self-energy at LO

In Practice: NLO + NNLO

- NLO irreducible VVCS graphs (f_L ; crossed not shown)



- LO+NLO deuteron self-energy

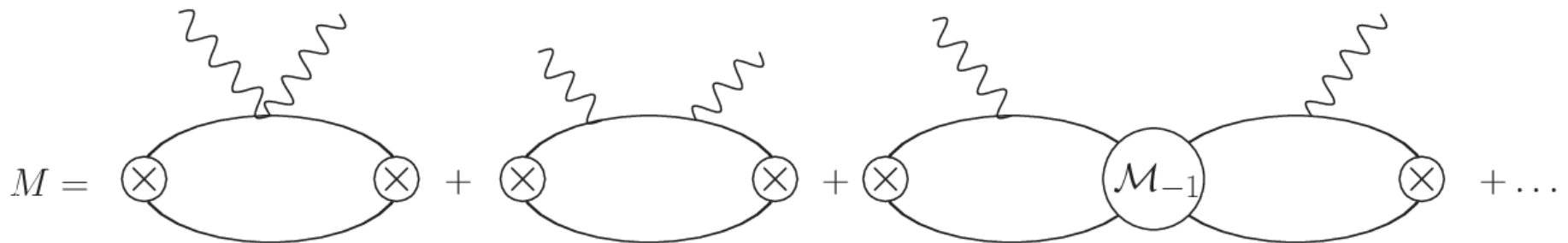
$$\Sigma = \text{[Diagram 1]} + \text{[Diagram 2]} + \dots$$

The first diagram is a single oval with two external wavy lines and vertices marked with an 'X'. The second diagram is two ovals connected by a central block labeled C_2 , with four external wavy lines and vertices marked with an 'X'.

- NNLO: ~30 irreducible VVCS graphs, 2 self-energy graphs

In Practice: Evaluation and Checks

- The diagrams are evaluated analytically (pionless EFT!)
- We use the PDS scheme to regularise the divergences
- The regularisation scale dependence has to vanish in the total amplitude
 - If it does not: look for an error (or a missing NN contact term)
- Current conservation/gauge invariance is checked explicitly



The diagram shows the expansion of the amplitude M as a sum of Feynman diagrams. The first term is a bubble diagram with two external wavy lines and two vertices marked with an 'X'. The second term is a bubble diagram with two external wavy lines and two vertices marked with an 'X'. The third term is a bubble diagram with two external wavy lines and two vertices marked with an 'X', with a central circle labeled \mathcal{M}_{-1} . The expansion continues with an ellipsis.

$$M = \text{bubble} + \text{bubble} + \text{bubble} \circ \mathcal{M}_{-1} + \dots$$

$$M = \epsilon_\mu \epsilon_\nu'^* M^{\mu\nu}$$

$$q_\mu M^{\mu\nu} = q_\nu M^{\mu\nu} = 0 \quad \text{individually at each order}$$

- all pieces that do not fulfil this have to cancel

In Practice: Contributions

- The expression for the residue is very simple:

$$\left. \frac{d\Sigma(E)}{dE} \right|_{E=E_d} = \frac{M^2}{8\pi\gamma} [1 - (Z - 1) + (Z - 1)^2 - (Z - 1)^3 + \dots]$$

$$\left[\left. \frac{d\Sigma(E)}{dE} \right|_{E=E_d} \right]^{-1} = \frac{8\pi\gamma}{M^2} [1 + (Z - 1) + 0 + 0 + \dots]$$

- The bulk of the contribution comes from the LO VVCS graphs + NLO residue correction
- NLO (+NNLO) VVCS graphs are small $\sim 1\%$
- Transverse and (some of) the relativistic terms are $\sim 1\%$

In Practice: the LS Integral

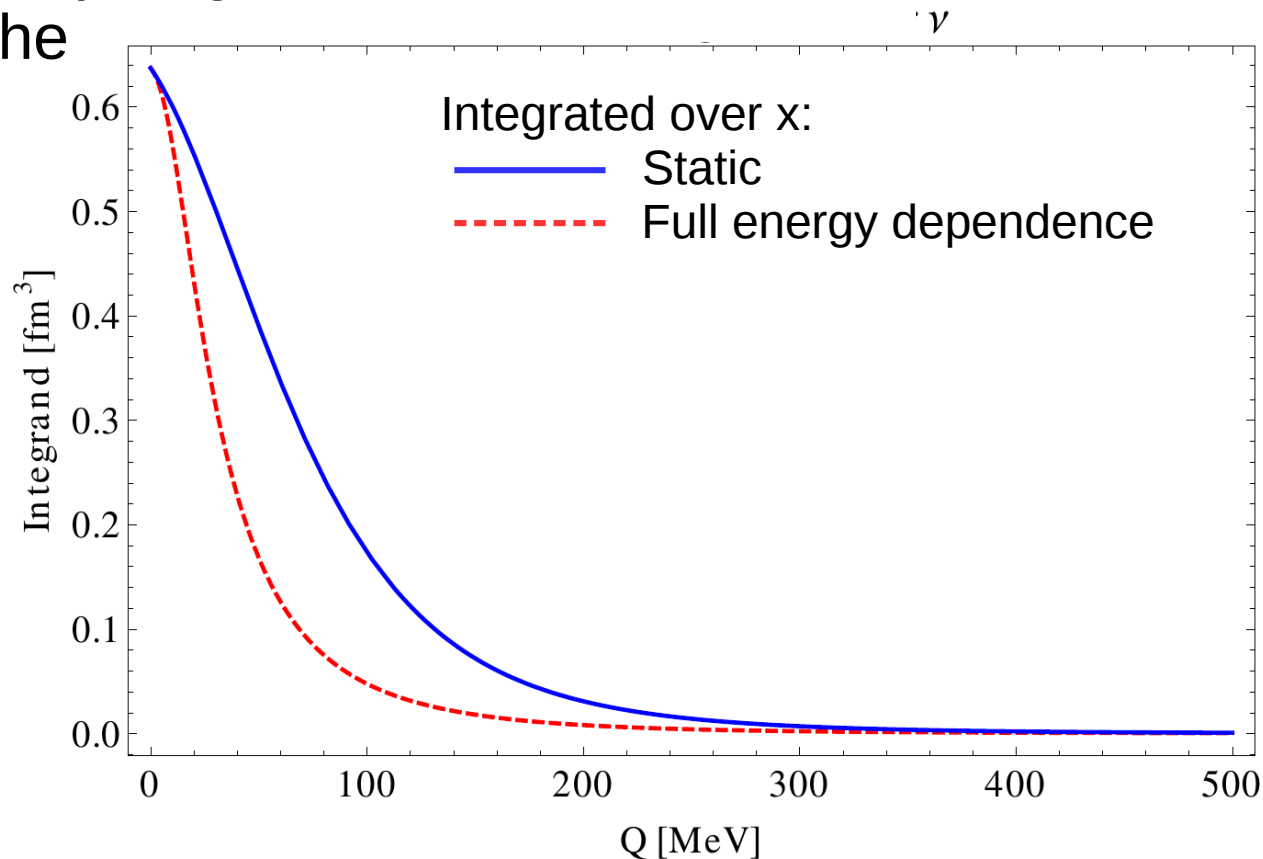
- The energy shift (non-pole f_L , Wick-rotated, hyperspherical)

$$\Delta E_{2S}^{\text{inel}} = \frac{2\alpha}{\pi m_\mu} [\phi_{2S}(0)]^2 \int_0^\infty \frac{dQ}{Q} \int_{-1}^1 dx \sqrt{1-x^2} \frac{f_L(-iQx, Q^2)}{\frac{Q^2}{4m_\mu^2} + x^2}$$

- Static limit $f_L(-iQx, Q^2) \rightarrow f_L(0, Q^2)$
works well for muonic hydrogen

but not here – keep the
energy dependence

- Better convergence
- High-momenta
contribution
is (hopefully)
under control



Results (Preliminary)

- Results for the μ D Lamb shift contribution

$$\Delta E_{2S}^{\text{inel}} = -1.58(0.04) \text{ meV}$$

- Other contributions:
 - Elastic: from deuteron FFs, well under control = $-0.415(2) \text{ meV}$
 - Coulomb: calculated in Hamiltonian approach = 0.262 meV

- Adding them up

$$\Delta E_{2S}^{\text{TPE}} = \Delta E_{2S}^{\text{inel}} + \Delta E_{2S}^{\text{elastic}} + \Delta E_{2S}^{\text{Coulomb}} = -1.73(0.05) \text{ meV}$$

- Compares well with other calculations
 - $1.675(0.04) \text{ meV}$ [Hernandez et al. 2019](#)
 - $1.680(0.016) \text{ meV}$ [Pachucki 2011](#)
 - $1.7096(0.02) \text{ meV}$ [Krauth et al. 2016 \(theory summary\)](#)
- Differences are most likely due to nucleon FFs (relativistic corrections), the present calculation is for pointlike nucleons

Sum Rules

- The sum rules connect the TPE correction to the target structure functions F_1 and F_2 [inclusive electron scattering]

$$\Delta E_{2S}^{\text{inel}} = -\frac{2\alpha^2}{M_d m_\mu} [\phi_{2S}(0)]^2 \int_0^\infty \frac{dQ^2}{Q^2} \int_{\nu_{\text{thr}}}^\infty \frac{d\nu}{\nu} \left[\tilde{\gamma}_1(\nu, Q^2) F_1(\nu, Q^2) + \frac{M_d \nu}{Q^2} \tilde{\gamma}_2(\nu, Q^2) F_2(\nu, Q^2) \right]$$

- There is also a subtraction contribution expressed as a weighted integral of $F_1(\nu, Q^2) - F_1(\nu, 0)$ Gorshteyn
- Use a parameterisation of the structure function, fit to the available data – a data-driven approach

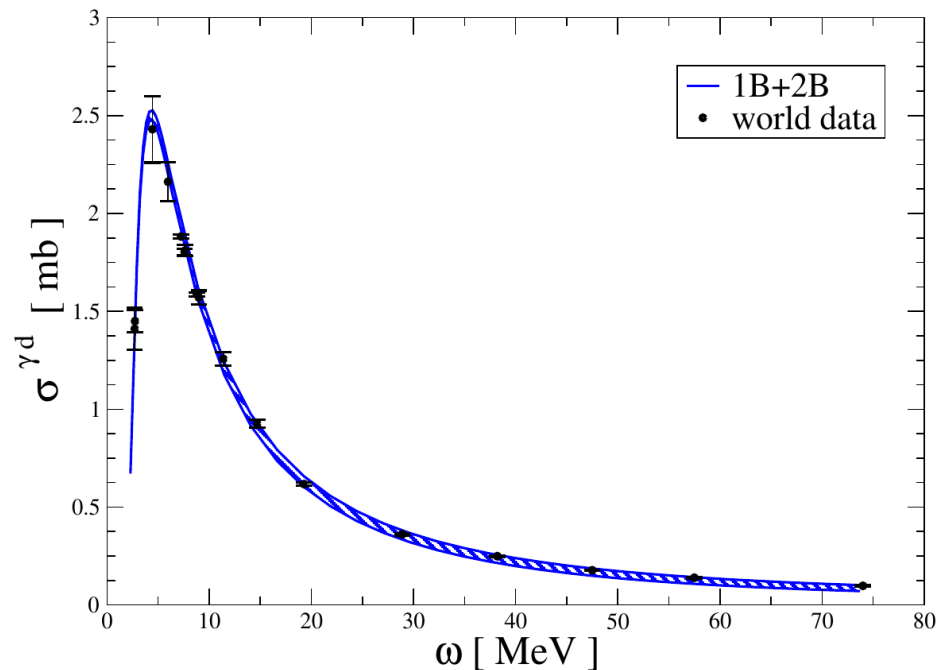
$$\Delta E_{2S}^{\text{inel}} = -1.749 \pm 0.740 \text{ meV} \quad \text{Carlson, Gorshteyn, Vanderhaeghen}$$

- Problem: too few quality low- Q^2 data \rightarrow big uncertainty
- Solution: use input from a theory (Chiral/pionless EFT)
- Verify theory using the available data, supplement the SR

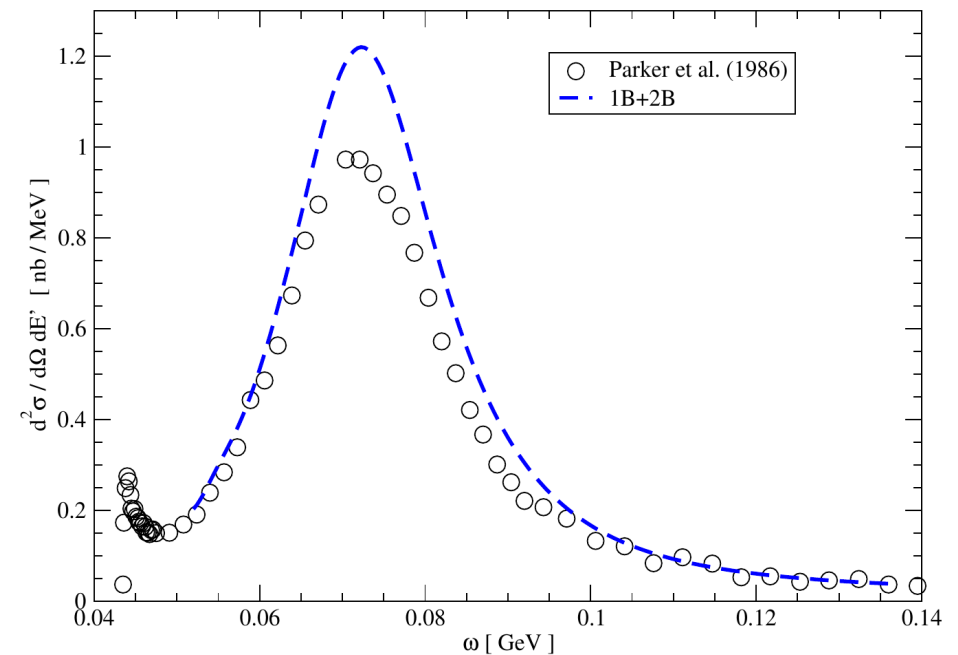
Sum Rules Aided by (Chiral) EFT

- NNLO Chiral EFT evaluation of the deuteron photo/electrodissociation cross section

Deuteron photodissociation



Electrodissociation $E=220$ MeV, $\theta=180^\circ$



- Data description is reasonably good so far
- Work in progress!
- Better data in low- Q^2 and low-energy region desirable!

Summary and Outlook

- Preliminary results for the TPE correction consistent with Hamiltonian approach and sum rules
- Full NNLO calculation, including error analysis (relativistic corrections, transverse contributions, ...) ready soon
- The most important relativistic corrections – due to the nucleon FFs – might consider to include them
- Coulomb [non-forward] distortions in EFT?
- The approach could be extended to the hyperfine splitting
- Sum rules with structure functions input from EFTs: verification of the theory and decreased error in the TPE correction evaluation