

# **Subtractive renormalization of the NN interaction and deuteron electro-disintegration**

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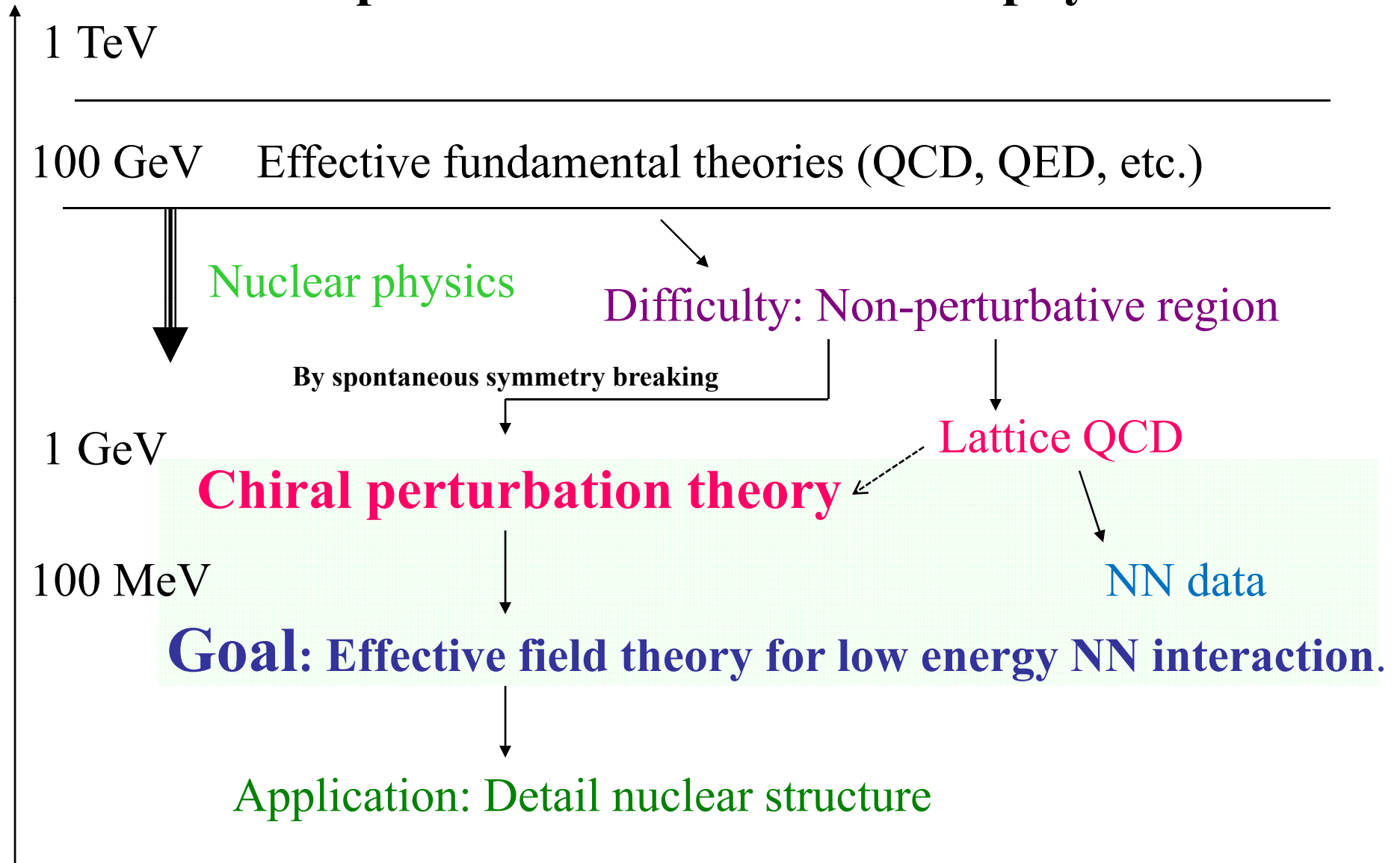
**Ohio University**

**C. J. Yang, C. Elster, and D. R. Phillips**

**PRC. 77 014002(2008), PRC.80 034002, PRC.80 044002**

**Summary:PoS(CD09)064 (arXiv:0909.5414)**

# A simplified overview of nuclear physics



# Approach to NN interaction

## Chiral EFT on NN sector:

- Lagrangian based on CHPT  
( $\pi N, \pi\pi$ , perturbative).
- Main category:

**perturbative,**

(no iteration of all diagrams)

Kaplan, Savage, Wise, Birse,  
Beane, etc.

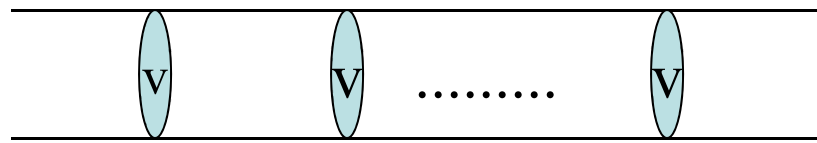
non-perturbative.

(Lippmann-Schwinger eq.)

Ordonez, Ray, van Klock,  
Epelbaum, Entem, Machleidt,  
Valderrama, Arriola, etc.


# NN interaction and renormalization

Lippmann-Schwinger eq. (LSE)



$$T(k, k; E) = \frac{e^{i\delta} \sin \delta}{-2\mu k}$$

$$T_l(k', k; E) = V_l(k', k; E) + \frac{2}{\pi} \int_0^\Lambda dp p^2 \frac{V_l(k', p; E) T_l(p, k; E)}{E - E_p + i\epsilon}$$

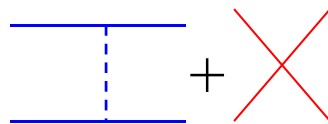
 Includes **long range part** (pion exchange) + **contact term**

Expanded in power of momentum.

In s-waves, LO:  $C_{S(T)}$ .

Represent short range (high energy) physics.

Cutoff in LSE



$C_{S(T)}(\Lambda)$  are determined by fitting to physical observables.


**A successful renormalization:** For different  $\Lambda$  (or at least in some interval), we should be able to get the same physical result.

# The goal of this work

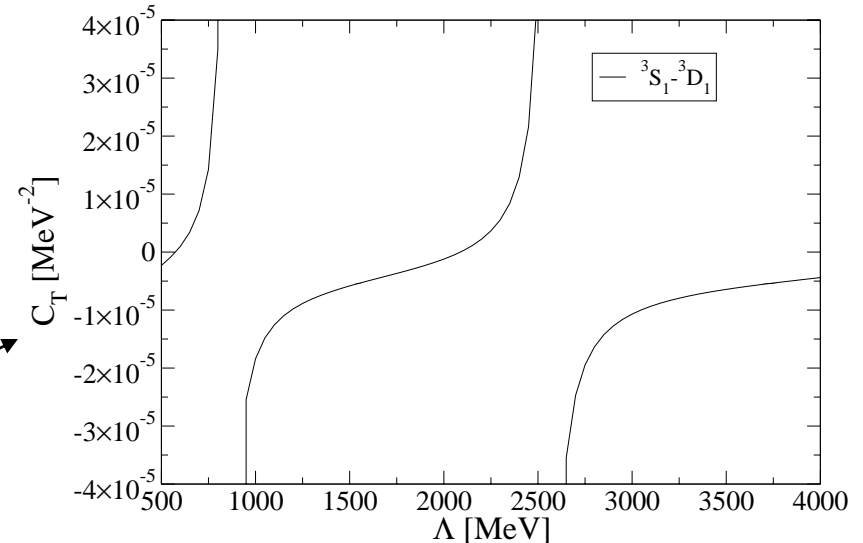
- We want to evaluate whether (or under what conditions)  $\chi$ ET, as an EFT (like QED or QCD), is really improved order by order, after renormalization.
  - \* More and more counter terms (with free parameters fitted to data), which somehow make it difficult to see whether the improvement is coming from the inclusion of higher order or the fit.
  - \* Many EFTs, each needs to be checked (RG) order by order.  
Pionful  $\rightarrow$  ~~analytic~~  $\rightarrow$  need quick and good way to pin down LEC's.
    1. What kind(order) of contact term should one adopt?
    2. Cutoff-indep. in phase shift  $\leftrightarrow$  renormalize successfully.
    3. What is the highest  $\Lambda$  can one use?
- **We invent a subtractive renormalization scheme to achieve our goal. Advantage: go to high  $\Lambda$ , eliminate correlation b/w LECs.**

# Determining LEC's in the contact terms

## 1. Fitting (Epelbaum, Machleidt, Valderrama, etc.)

- Previously, renormalize  $V$  by adjusting the unknown constants to fit data, e.g. fit scattering length, effective range or phase shifts.
- Problem: 1. fine-tuning. 2. 

Problem of fitting



## 2. Subtraction method:

**Leading order\*:** \*C. J. Yang, C. Elster, and D. R. Phillips, PRC. 77 014002(2008)

- By subtracting two LSE to cancel the divergent part ( $C_{S(T)}$ ).
- Direct input of the scattering length  $a_{S,T}$ .
- This was done by relating the on-shell  $t(p_0, p_0; E)$  for an arbitrary  $E$  to the fully-off-shell  $t(p', p; 0)$ .

# Method: Three steps

$$T(p, 0; 0) = V(p, 0; 0) + C + \frac{2}{\pi} \int_0^\Lambda dp' p' \left[ \frac{V(p, p'; 0) + C}{-p'^2/2\mu} \right] T(p', 0; 0) \quad (1) \quad T(0, 0; 0) = \frac{f_0}{-2\mu k} = \frac{a}{2\mu}$$

$$T(0, 0; 0) = V(0, 0; 0) + C + \frac{2}{\pi} \int_0^\Lambda dp' p' \left[ \frac{V(0, p'; 0) + C}{-p'^2/2\mu} \right] T(p', 0; 0) \quad (2)$$

$$(1) - (2): T(p, 0; 0) - \frac{a}{2\mu} = V(p, 0; 0) + \frac{2}{\pi} \int_0^\Lambda dp' p'^2 \left[ \frac{V(p, p'; 0) - V(0, p'; 0)}{-p'^2/2\mu} \right] T(p', 0; 0)$$

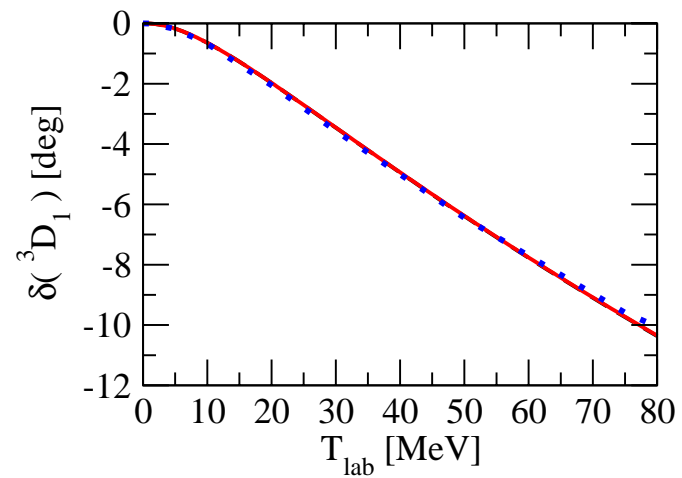
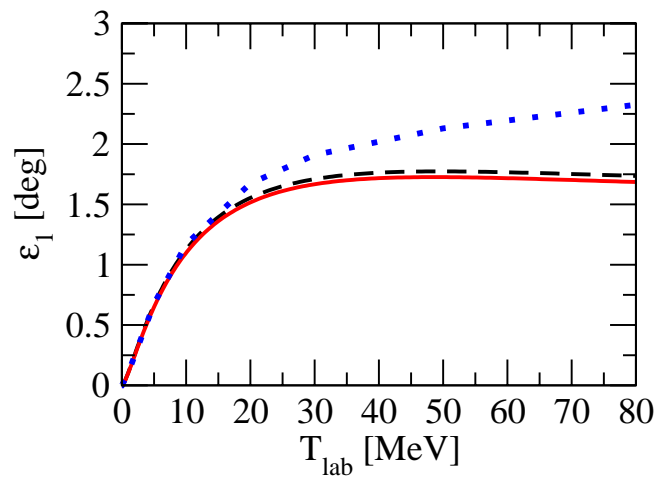
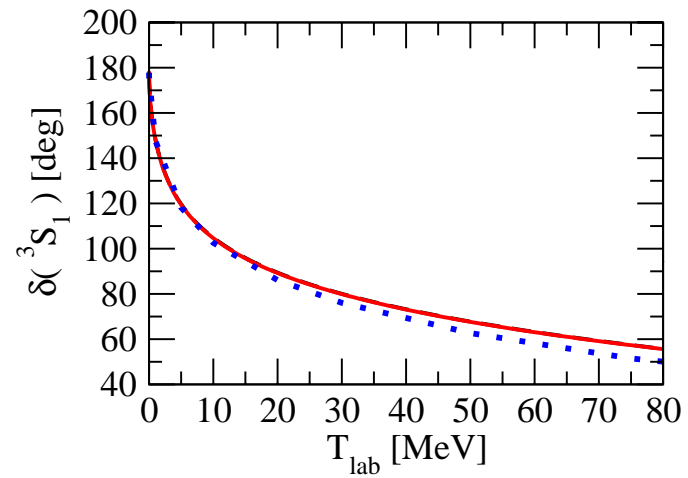
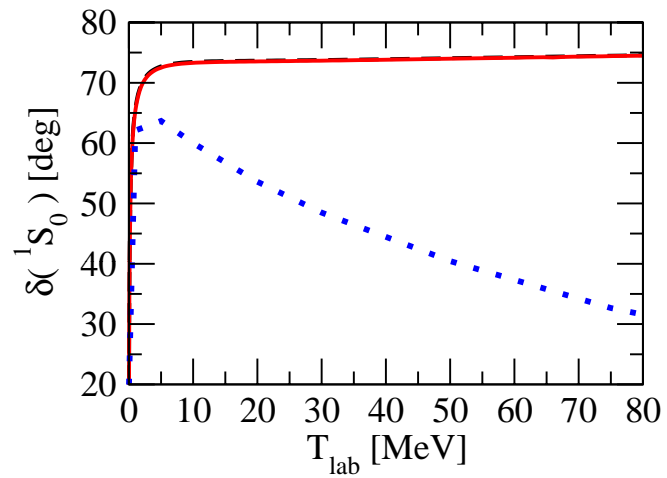
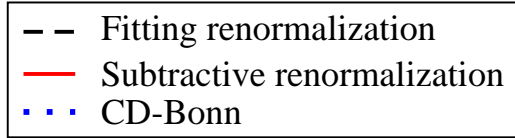
$$T(p, p^*; 0) - \frac{a}{2\mu} = V(p, p^*; 0) + C + \frac{2}{\pi} \int_0^\Lambda dp' p'^2 \left[ \frac{V(p, p'; 0) + C}{-p'^2/2\mu} \right] T(p', p^*; 0) \quad (3)$$

$$T(0, p^*; 0) - \frac{a}{2\mu} = V(0, p^*; 0) + C + \frac{2}{\pi} \int_0^\Lambda dp' p'^2 \left[ \frac{V(0, p'; 0) + C}{-p'^2/2\mu} \right] T(p', p^*; 0) \quad (4)$$

$$(3) - (4): T(p, p^*; 0) - T(0, p^*; 0) = V(p, p^*; 0) - V(0, p^*; 0) + \frac{2}{\pi} \int_0^\Lambda dp' p'^2 \left[ \frac{V(p, p'; 0) - V(0, p'; 0)}{-p'^2/2\mu} \right] T(p', p^*; 0)$$

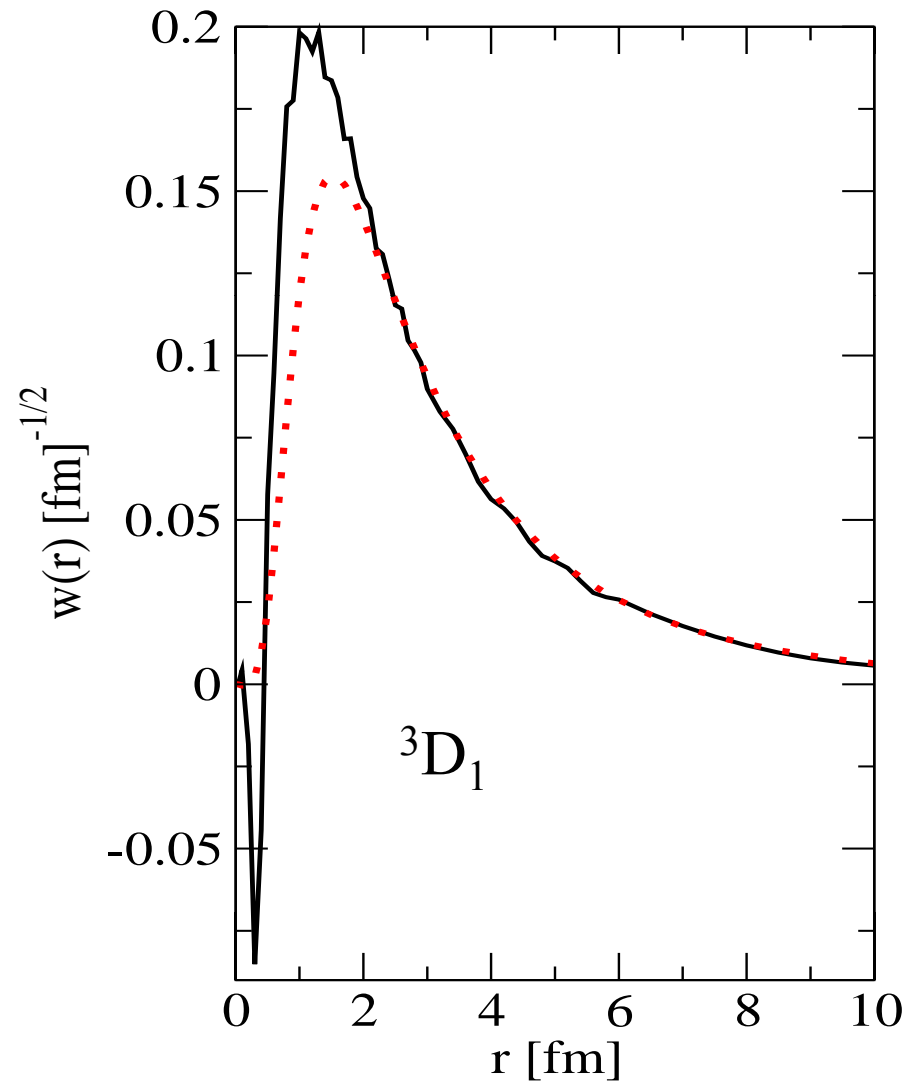
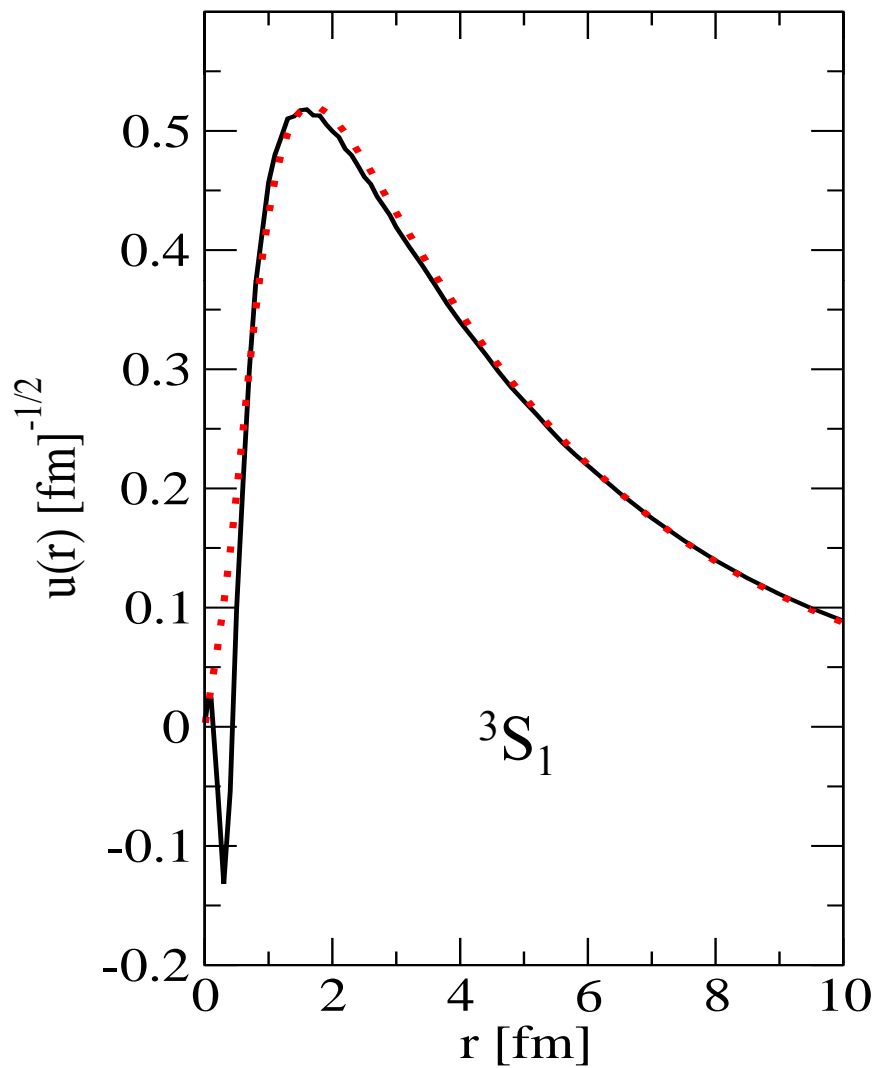
*Final step:*  $[1 + T(0)(G(0) - G(E))]T(E) = T(0)$

# LO Results





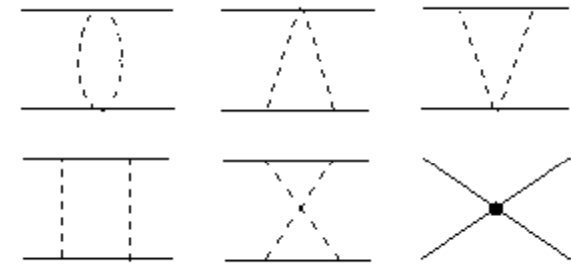
# Wave function



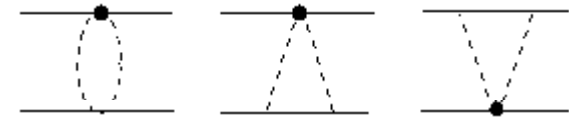
**So subtraction works for the leading  
order.**

**Next: Go to NNLO**

- In NLO( $Q^2$ ) and NNLO( $Q^3$ ) we have TPE, which diverges as  $Q^{2(3)}$ .  $\Rightarrow$  Include the  $O(Q^2)$  contact term to renormalize it.
- $V = \text{OPE} + \text{TPE} + \lambda + \gamma^* [O(Q^2)]$ .
- Further develop our subtraction technique to solve it.



NLO( $Q^2$ )



NNLO( $Q^3$ )

Use dimensional regularization (DR) or spectral function regularization (SFR) to regularize the diverge loop integral.

**Contact terms:** from LO to NLO/NNLO

In p-waves: none  $\rightarrow$   $C_{pp'}$ .

s-waves:  $C_{S(T)} \rightarrow \lambda_{II'} + \gamma_{II'} [O(Q^2)]$ .

## Subtraction: (up to) NNLO for p-waves

Need to cancel out :  $C^{SJ} p' k$ . Idea: get  $t(p', p; E) \leftarrow t(p', k; 0) \leftarrow \alpha_{11}$

Note:  $t_{l,1}(p', p; E) \sim (\dots) p'^1 p^1$

Step 1 of subtraction:

**Key:** 1. Divide by  $p' k$  and then take limit. 2.  $\lim_{k \rightarrow 0} \left[ \frac{t(k, k; 0)}{k k} \right] = \frac{\alpha_{11}}{M}$

$$\lim_{k \rightarrow 0} \left[ \frac{t(p', k; 0)}{p' k} \right] = \lim_{k \rightarrow 0} \left[ \frac{v^{LR}(p', k)}{p' k} + C_{11} \right] + \frac{2}{\pi} M \lim_{k \rightarrow 0} \left[ \frac{1}{p' k} \int_0^\Lambda dp'' p''^2 \frac{(v^{LR}(p', p'') + C_{11} p' p'') t(p'', k; 0)}{-p''^2} \right] \quad (1)$$

$$\lim_{k \rightarrow 0} \left[ \frac{t(k, k; 0)}{k k} \right] = \lim_{k \rightarrow 0} \left[ \frac{v^{LR}(k, k)}{k k} + C_{11} \right] + \frac{2}{\pi} M \lim_{k \rightarrow 0} \left[ \frac{1}{k k} \int_0^\Lambda dp'' p''^2 \frac{(v^{LR}(k, p'') + C_{11} k p'') t(p'', k; 0)}{-p''^2} \right] \quad (2)$$

$$p' [(1) - (2)] \Rightarrow \lim_{k \rightarrow 0} \left[ \frac{t(p', k; 0)}{k} \right] - \frac{\alpha_{11}}{M} p' = \lim_{k \rightarrow 0} \left[ \frac{v^{LR}(p', k)}{k} \right] - p' \lim_{k \rightarrow 0} \left[ \frac{v^{LR}(k, k)}{k k} \right]$$

$$+ \frac{2}{\pi} M \lim_{k \rightarrow 0} \left[ \int_0^\Lambda \frac{dp'' p''^2 (v^{LR}(p', p'') - \lim_{k \rightarrow 0} \left[ \frac{v^{LR}(k, p'')}{k} \right] p') \lim_{k \rightarrow 0} \left[ \frac{t(p'', k; 0)}{k} \right]}{-p''^2} \right] \quad \text{get } \lim_{k \rightarrow 0} \left[ \frac{t(p', k; 0)}{k} \right].$$

*Step 2 of subtraction:*

Use result of step 1, i.e.,  $\lim_{k \rightarrow 0} [t(p', k; 0)/k]$  as input.

Do similar subtraction again.  $\Rightarrow$  get  $\frac{t(p'', p'; 0)}{p'}$ , or,  $t(p'', p'; 0) \equiv t(0)$ .

Final step:

Use  $t(E) = t(0) + t(0)[g(E) - g(0)]t(E)$ .

$\Rightarrow$  get on-shell t-matrix  $t(E)$ .

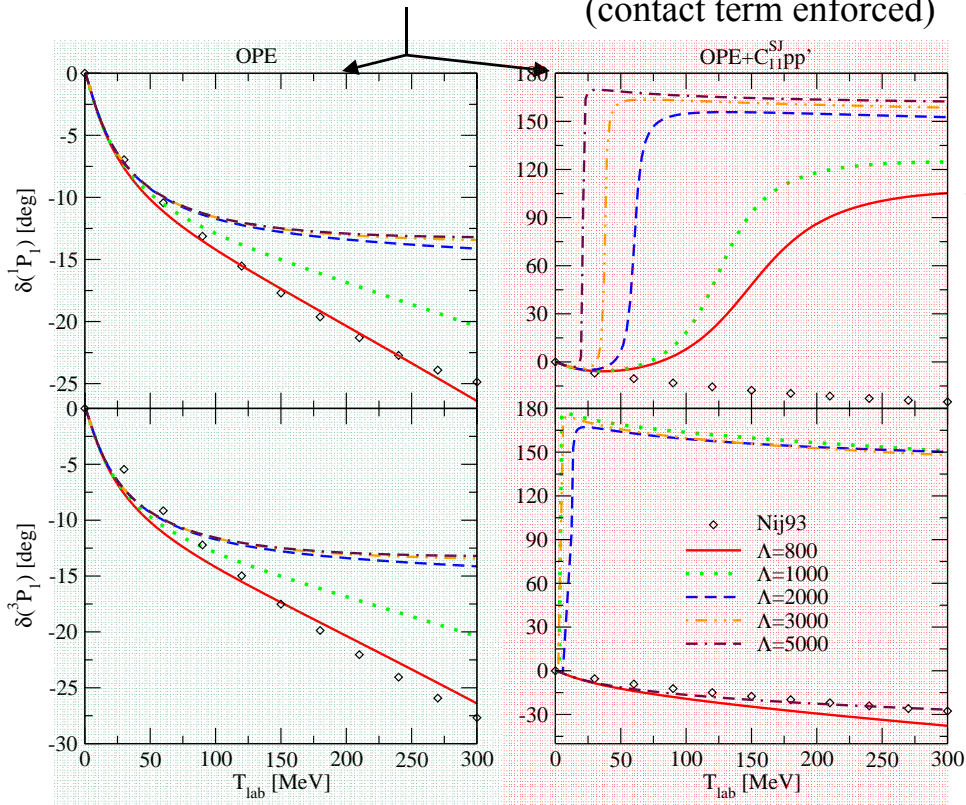
# Results at Leading order(LO):

See also: Nogga, Timmermans and van Kolck (2005), Valderrama (2006).

## unrenormalized v.s. renormalized (Fail of Weinberg counting)

singular & repulsive

(contact term enforced)

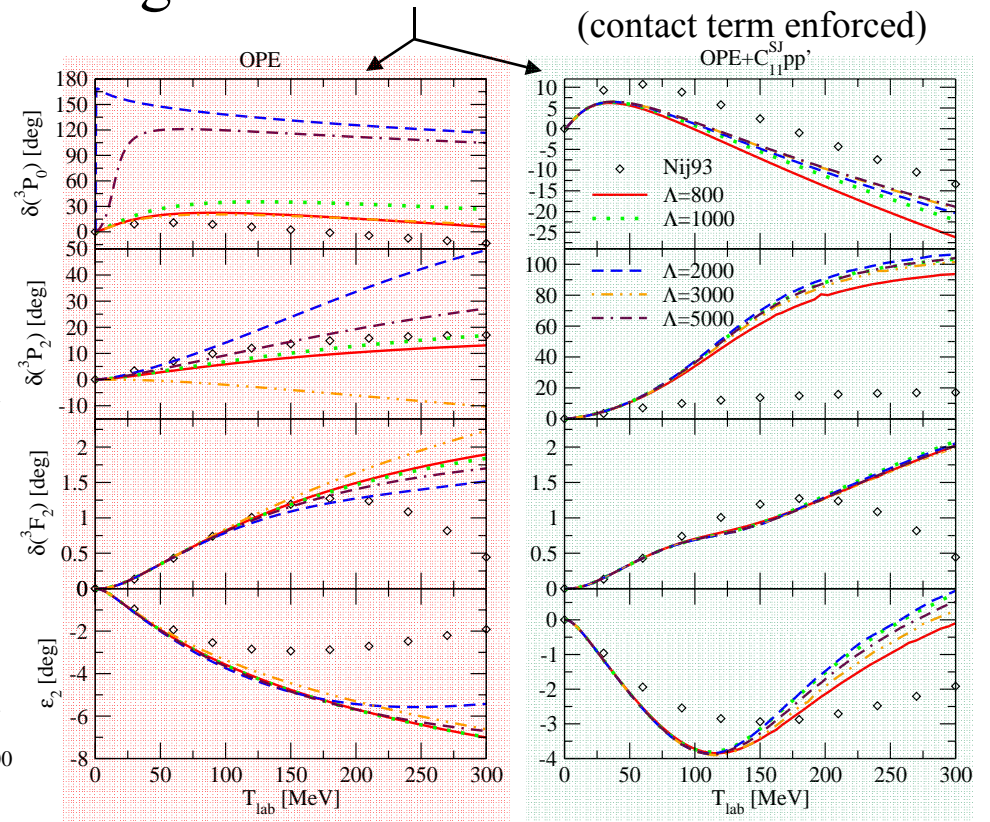


convergent

Divergent w.r.t.  $\Lambda$

singular & attractive

(contact term enforced)



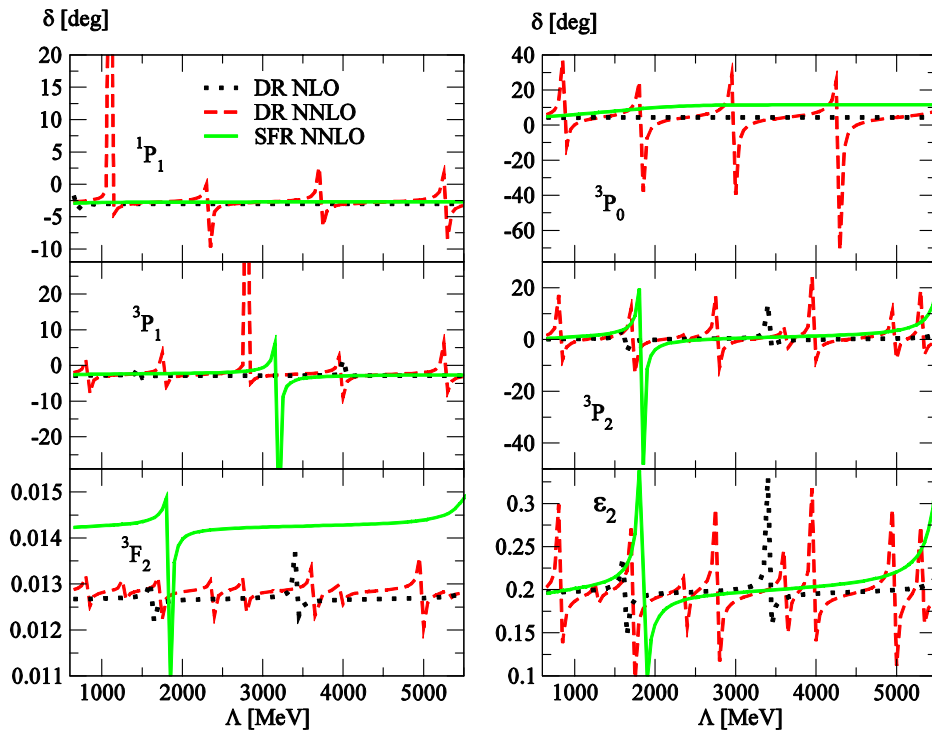
Divergent w.r.t.  $\Lambda$

convergent

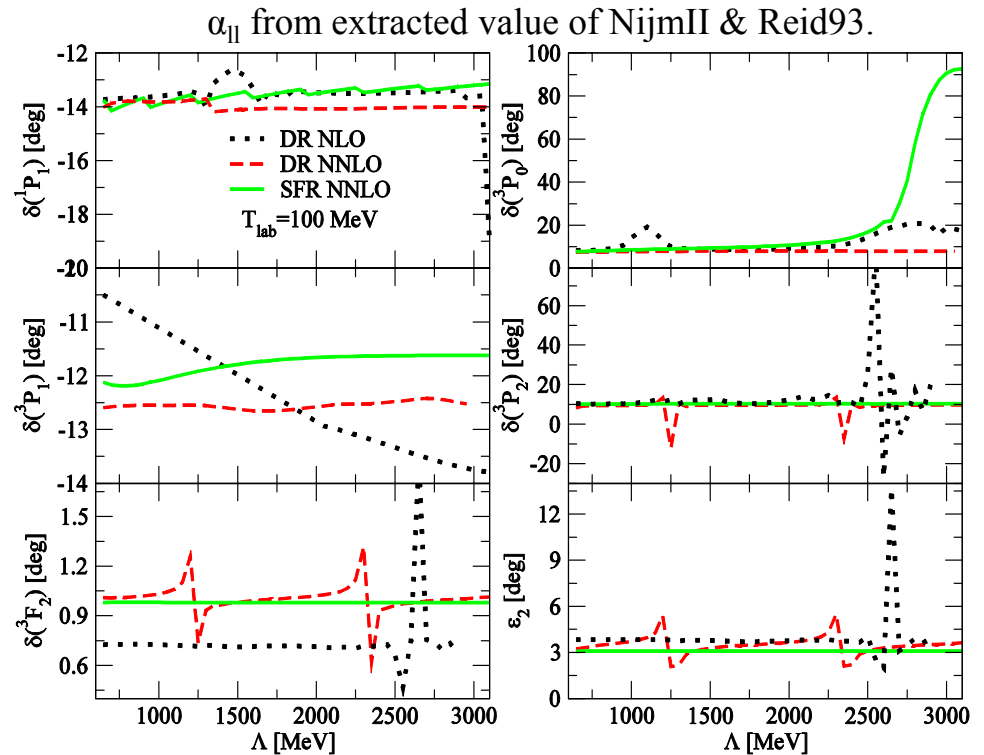
$\alpha_{11}$  from extracted value of NijmII & Reid93.

# $r \rightarrow 0$ connection for the higher order

Un-renormalized



Renormalized



	$^1P_1$	$^3P_0$	$^3P_1$	$^3P_2$
OPE	U	R	U	R
NLO (DR)	U	U	R	R
NNLO (DR)	R	R	R	*
NLO (DR) + NNLO (SFR)	U	U	R	R
NNLO (SFR)	U	U	R	R

*Need counter term to stabilize?*

*Attractive: Yes (R on the left).*

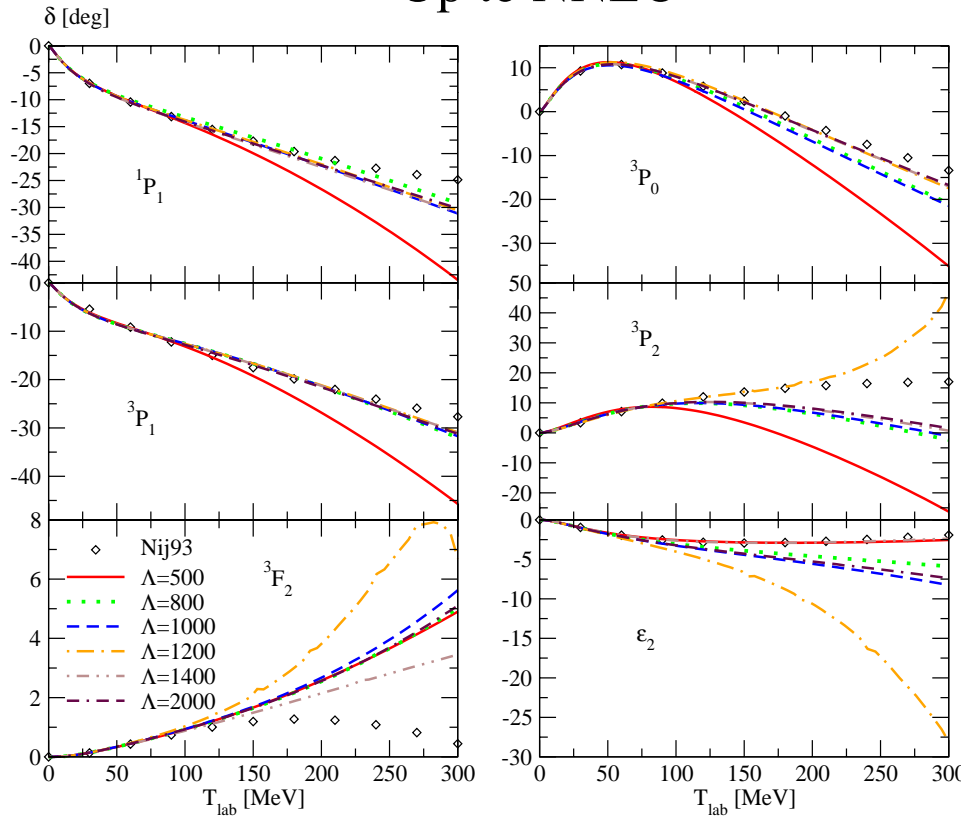
*Repulsive: No (U on the left).*

(See: Valderrama, et al (2006) for same thing in S-waves.)

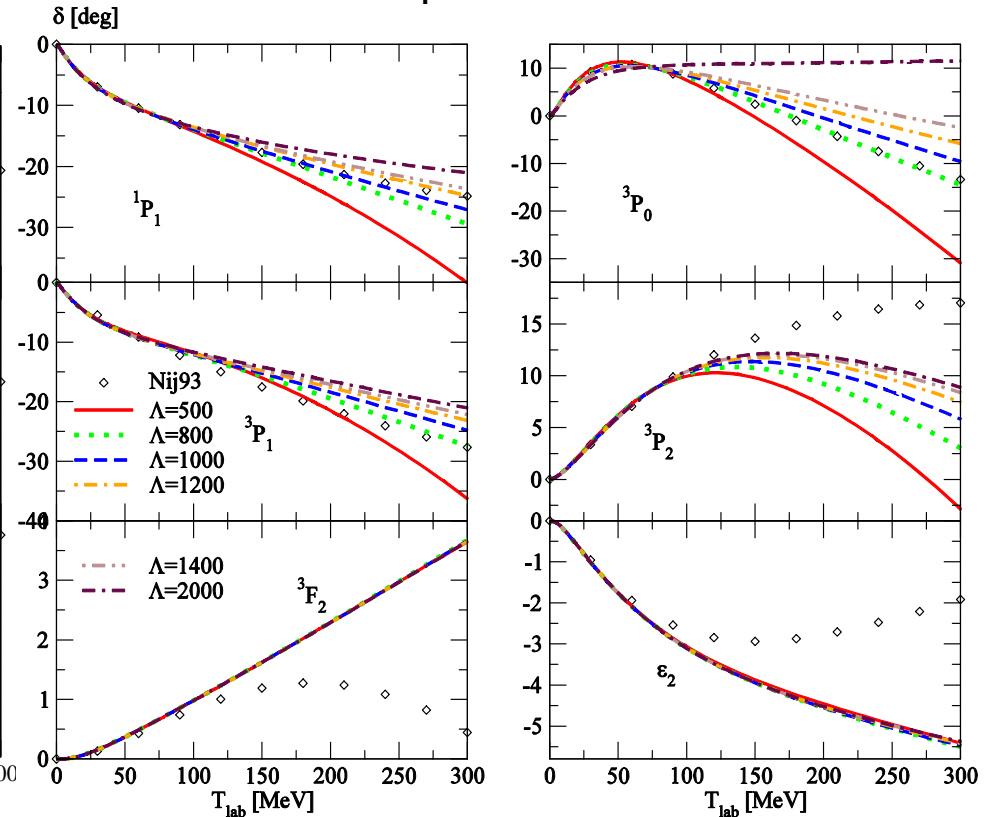
# Best fit: by adjusting $\alpha^{\text{SJ}}$

For DR TPE, need to vary  $\alpha^{\text{SJ}}$  away from Nijm value up to 30% in some channels.

Dimensional regularization (DR)  
Up to NNLO



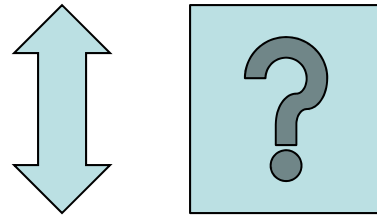
Spectral function regularization (SFR)  
Up to NNLO





**Even if we can fit to the data, does this necessarily mean that the renormalization is done successfully?**

**Cutoff-indep. in phase shift**

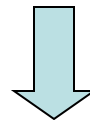
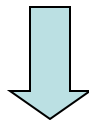


**Successful Renormalization.**

# $\alpha_{\text{best}}$ v.s. $\Lambda$ (DR TPE)

Dimensional regularization (DR) up to NNLO

$\Lambda$ [MeV]	$\alpha_{11}^{01}(^1P_1)$	$\alpha_{11}^{10}(^3P_0)$	$\alpha_{11}^{11}(^3P_1)$	$\alpha_{11}^{12}(^3P_2)$
500	3.6405	-4.3385	3.1326	-0.59589
800	3.6455	-4.8930	2.9524	-0.47614
1000	3.6593	-4.9346	2.9552	-0.47028
1200	<u>3.6593</u>	<u>-4.5049</u>	<u>3.3613</u>	<u>-0.42356</u>
1400	3.6732	-3.4098	4.1098	-0.45942
1600	3.6732	-2.5089	4.4564	-0.46142
1800	3.6732	-2.3425	4.0405	-0.46142
2000	3.6732	-2.9386	3.4819	-0.45642



$\Lambda_c \sim 1\text{GeV}$

(above that, renormalization point dependence show up.)

# Summary of p-waves results

## Whether need contact term to reach $\Lambda$ independence?

=> exactly depends on the singularity structure of  $V(r \rightarrow 0)$ .

**Attractive: Yes.**

**Repulsive: No.**

## Successfully renormalized?

1. There is a critical cutoff  $\Lambda_C \sim 1$  GeV for DR TPE up to NNLO, after that the contact term in LSE dominate the result.
2. Replacing the whole DR NNLO TPE by SFR brings  $\Lambda_C$  up to 2.5 GeV.
3.  $\Lambda_C$  is in the same order of  $\Lambda_\chi$  ( $\sim 1$  GeV).

## Applications:

**This scheme works for any contact term has the form  $C p^l p'^l$ ,  
=> can be applied in D-waves and F-waves.**

# NNLO in S-waves

- Adopt more subtractions to perform the renormalization.
- In S-wave, contact term (up to  $Q^2$ ) has the following structure:

$\lambda + C_2(p^2 + p'^2)$ .....for  $^1S_0$ . (two unknowns)  $\Rightarrow$  solve by 1 sub + 1 fit.

$\begin{pmatrix} \lambda + C_2(p^2 + p'^2) & \lambda_1 p'^2 \\ \lambda_1 p^2 & 0 \end{pmatrix}$ ....for  $^3S_1 - ^3D_1$ . (three unknowns)  $\Rightarrow$  solve by 2 sub + 1 fit.

-

# Method

- Combine previous ideas:
  1. Use first subtraction to eliminate  $\lambda$ .
  2. Use previous formula to relate  $t(E^*)$  to  $t(E)$ .
  3. For  ${}^3S_1$ - ${}^3D_1$ , apply the idea of dividing  $p^l p'^l$  in LSE to eliminate  $\lambda_t p^2$ .
  4. Input:  $a_0$ ,  $\delta(E^*)$  for the singlet; with additional  $\alpha_{20}$  for triplet.

# 1 Subtraction + 1 fitting

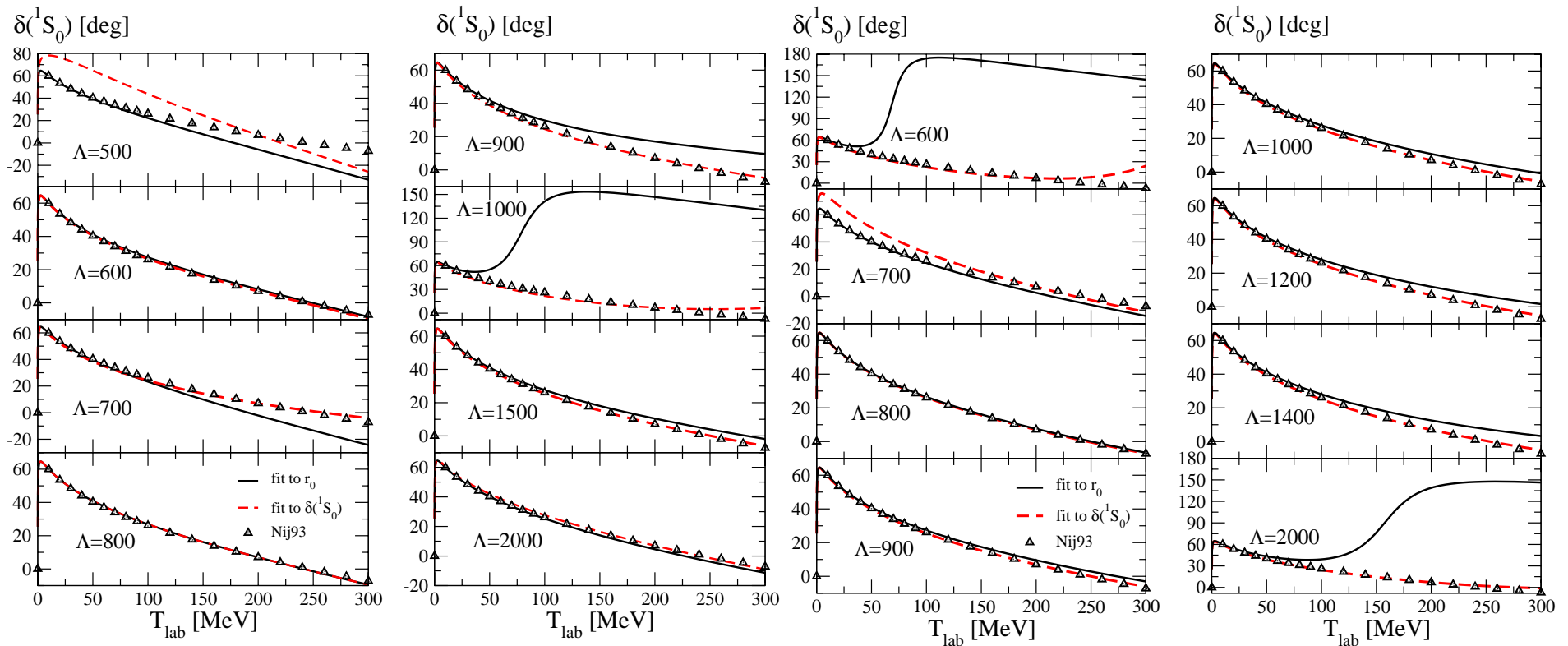
- For contact term at NLO/NNLO, we need to replace one input by one fitting (due to the fact that  $C_2(p^2+p'^2)$  term has *no* on-shell observable corresponds to it.)
- So, we need 1 fitting (fitted either to  $r_0$  or  $\delta(E^*)$  ).

# Results: Singlet S-wave

At higher  $\Lambda$ , has fitting point dependence.

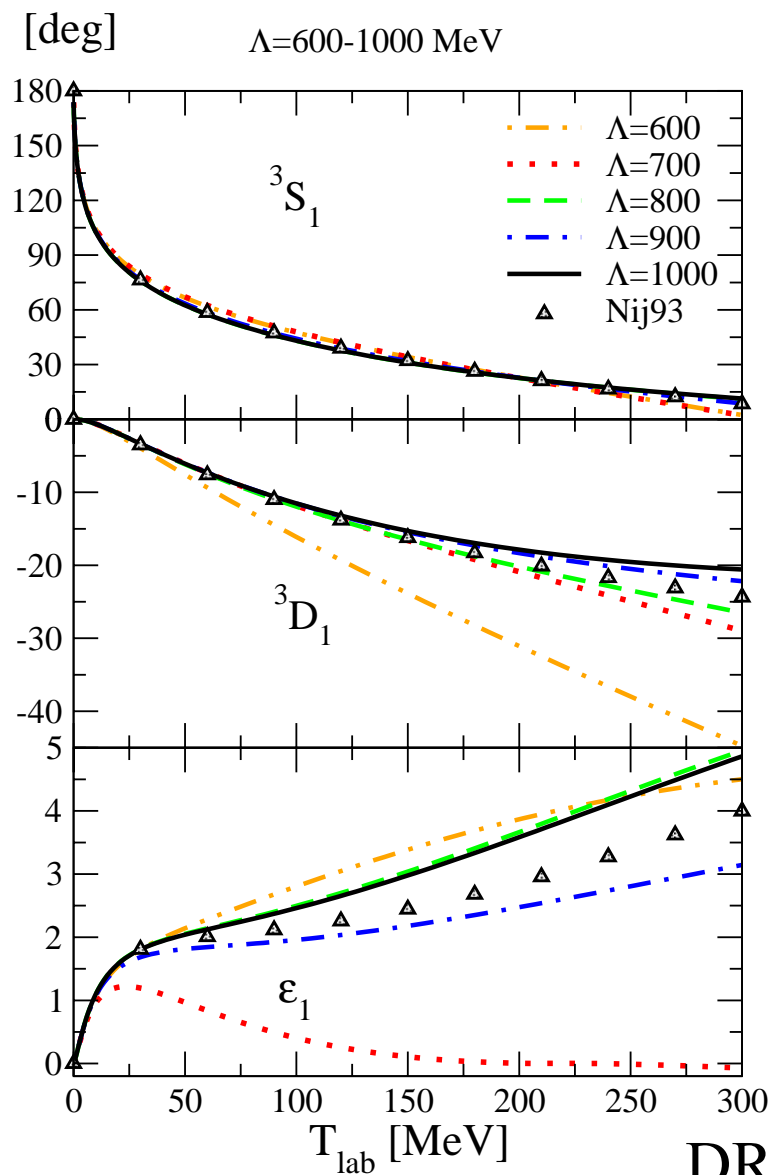
DR NNLO(fails at 1GeV)

Full SFR NNLO(fails at 2 GeV)

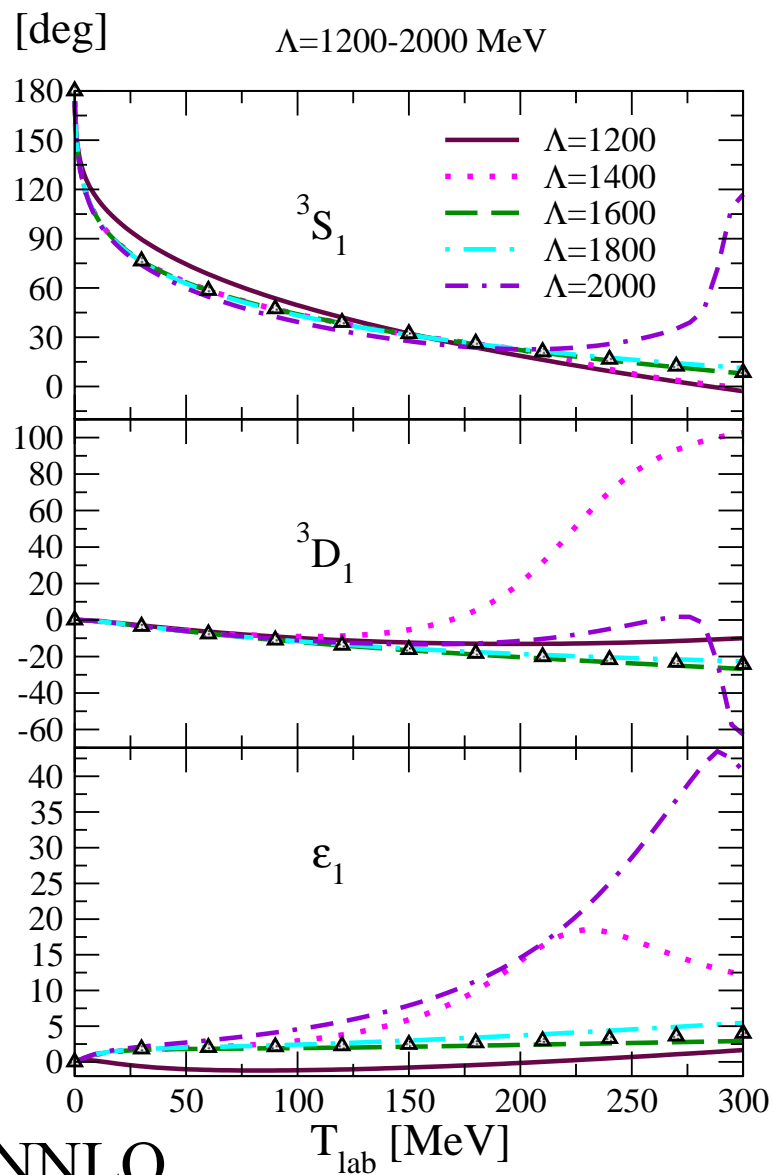


$\Lambda$  cannot be too low ( $\leq 500$  MeV) also.

# Results: Triplet S-wave (2 sub+1 fitting)

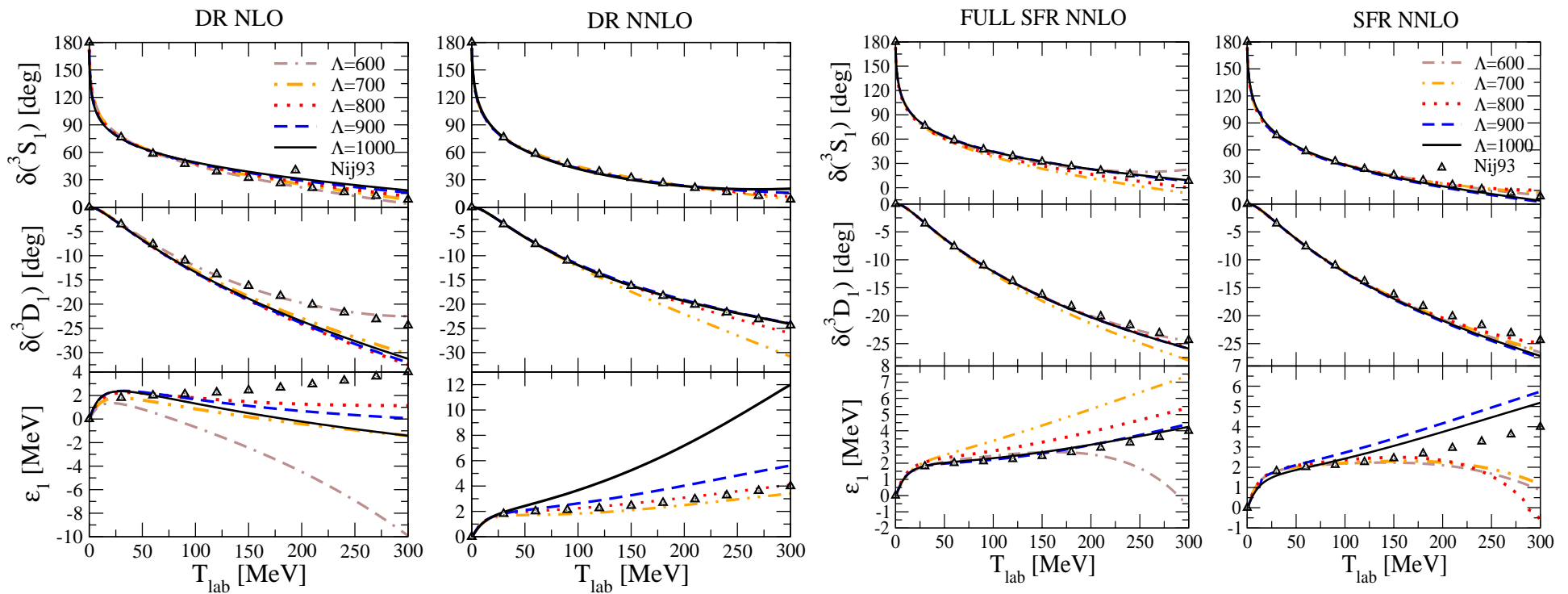


DR NNLO





# Allow all 3 constants to fit freely



**3 fitting > 2 sub + 1 fitting .**

$\Rightarrow$  The two low energy inputs prohibits the best fit.

## **Additional: Energy dependent contact term**

- Replace  $(p^2+p'^2)$  by  $E$ .
- Why?

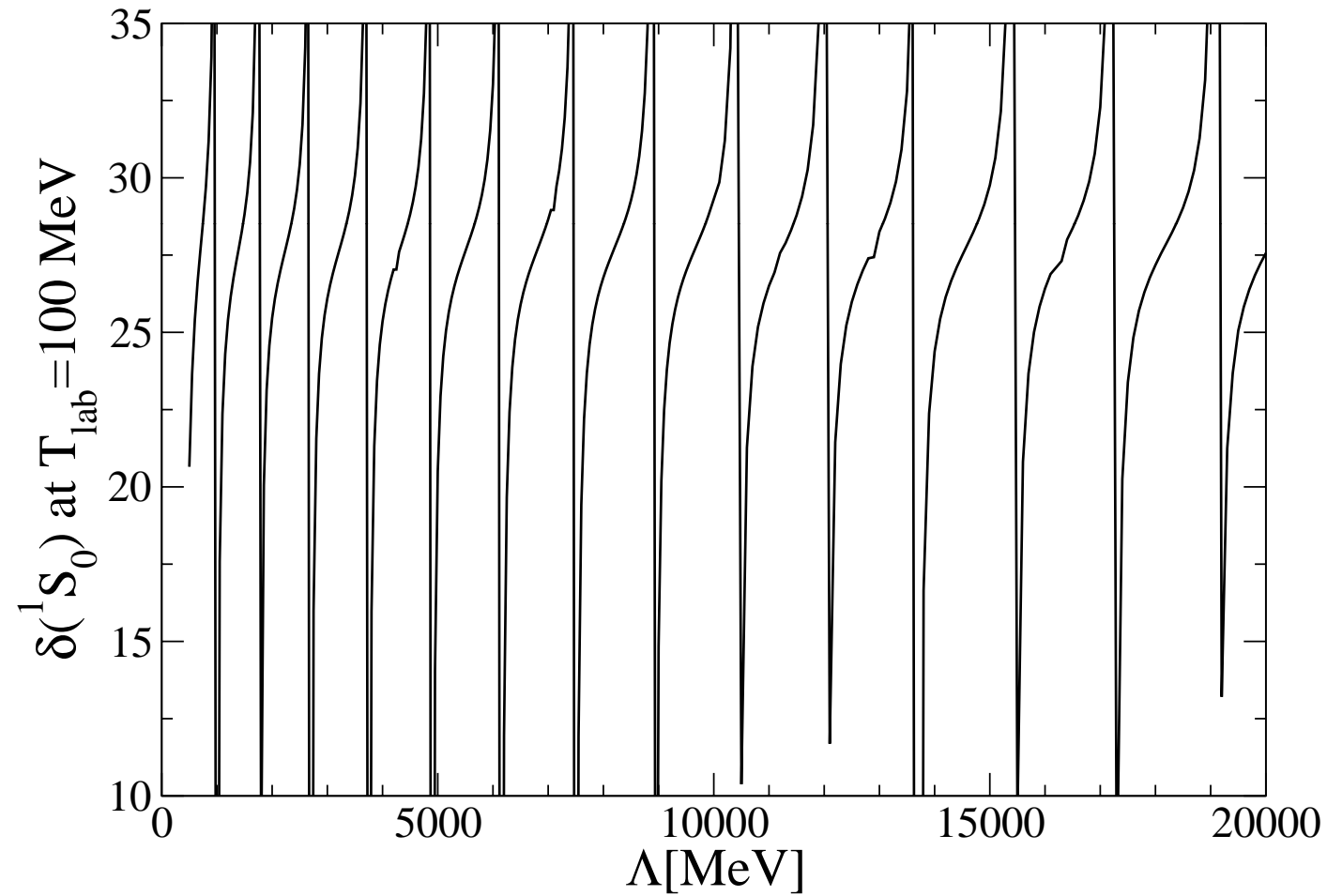
=> See whether the E-dep. term fit data better.

$\lambda + \gamma E$ .....for  $^1S_0$ . (two unknowns) => solve by 2 sub.

$\begin{pmatrix} \lambda + \gamma E & \lambda_t p'^2 \\ \lambda_t p^2 & 0 \end{pmatrix}$ .....for  $^3S_1 - ^3D_1$ . (three unknowns) => solve by 3 sub.

Result: *When it fits, it fits better (than p-dep). But,*

# Oscillation with respect to $\Lambda$



# S-waves Summary

- **P-dep. contact term:**

For  $^1S_0$  the indep. of renormalization point breaks once  $\Lambda > 1 \sim 1.2$  GeV for DR NNLO, and  $\Lambda \sim 2000$  for SFR NNLO.

For  $^3S_1$ - $^3D_1$ , fit breaks down at about  $\Lambda \sim 1.2$  GeV in general.

- **E-dep. contact term**, there is oscillatory behavior.

1. Singlet channel: The first diverged phase shift appears at  $\Lambda \sim 1000$  MeV for DR NNLO

$\Lambda \sim 2000$  MeV for Full SFR NNLO.

2. Triplet channel: The first diverged phase shift appears at  $\Lambda \sim 1200$  MeV for DR NNLO and

$\Lambda \sim 2300$  MeV for Full SFR NNLO.

# Conclusion on NN scattering

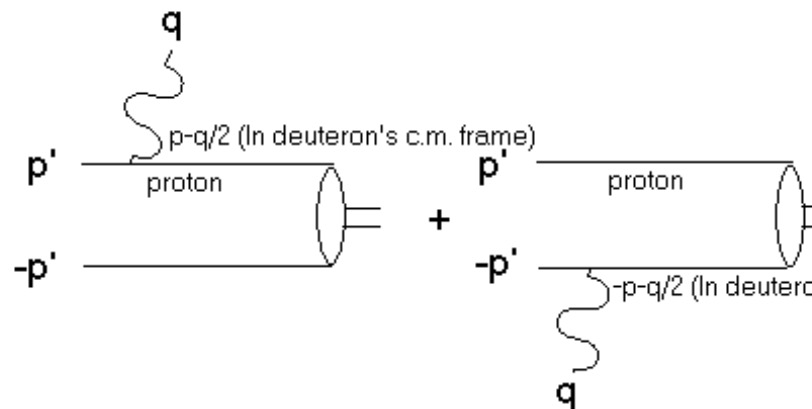
- Subtraction method provides:
  1. An easy way to go to high cutoff in LSE.
  2. A clean information of the dependence of results on the low energy observable ( $\alpha^{SJ}$ ).
- We found:
  1. **There is a critical cutoff  $\Lambda_C \sim 1$  GeV for DR TPE up to NNLO, after that the contact term in LSE dominate the result.**
  2. **Replacing the whole DR NNLO TPE by SFR brings  $\Lambda_C$  up to 2 GeV.**

# Deuteron electro-disintegration

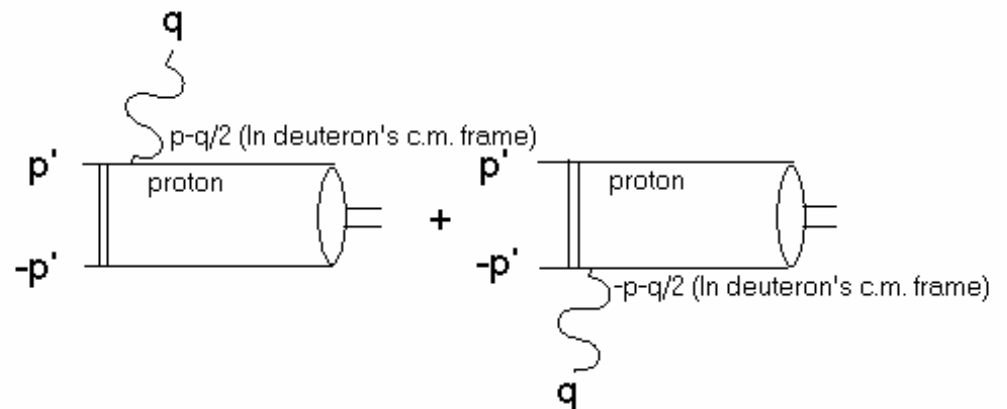
Motivation:

- Further test the wfs and t-matrices from  $\chi$ pt potential.
- Look for cutoff-dependence in other physical process.
- Compare results to traditional models.

Impulsive approximation



Final state interaction



# Longitudinal response function

$$\frac{d\sigma}{d\Omega_s} = c\{\rho_L f_L + \rho_T f_T + \rho_{LT} f_{LT} + \rho_{TT} f_{TT} + h(\rho'_{LT} f'_{LT} + \rho'_T f'_T)\}.$$

Where  $h$  is the electron polarization degree,  $\rho(f)_{L,T,LT,TT}$  describe the lepton(hadron) tensor.

$c = \frac{\alpha}{6\pi^2} \frac{k_2^2}{k_1^2 q_0^4}$ , where  $k_{1,2}^L$  is the abs. value of the electron momentum in the Lab frame.

$$f_L = -\pi^2 2\alpha p_{np} M_N |M_{T_f S' M_s'}(q, 0; \omega)|^2$$

$$M_{T_f S' M_s'}(q, 0; \omega) = \langle \mathbf{p}' S' M_s' T_f | J(\mathbf{q}; 0) | M \rangle + \langle \mathbf{p}' S' M_s' T_f | t(E') G_0(E') J(\mathbf{q}; 0) | M \rangle$$

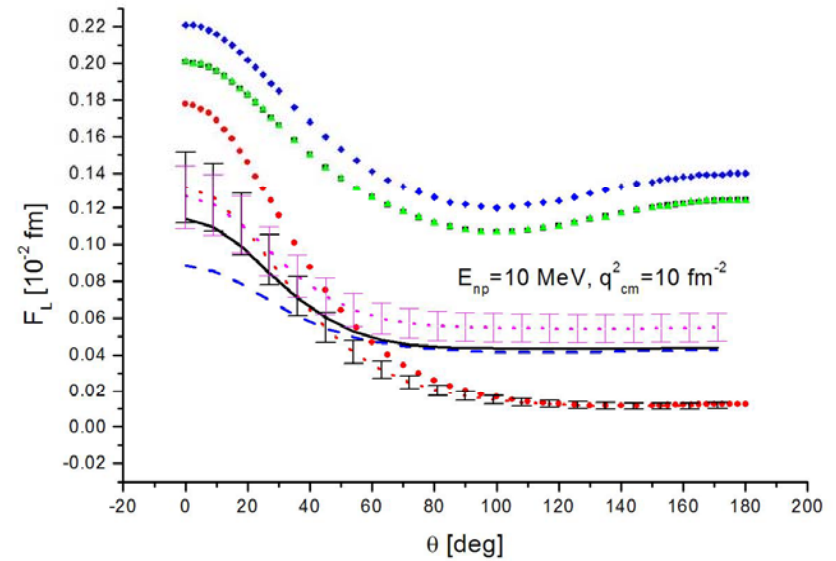
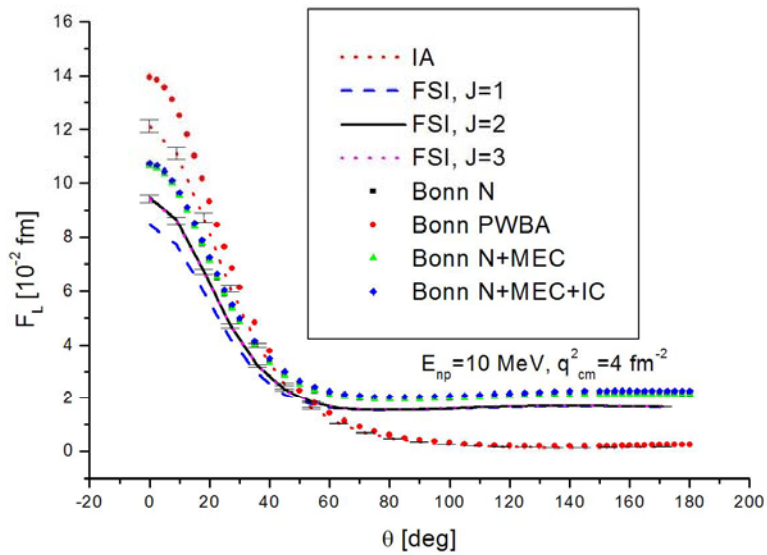
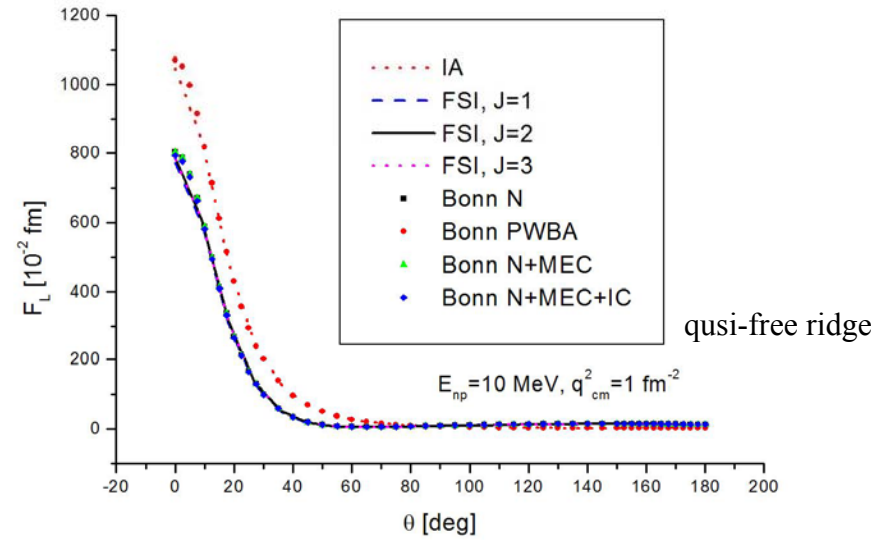
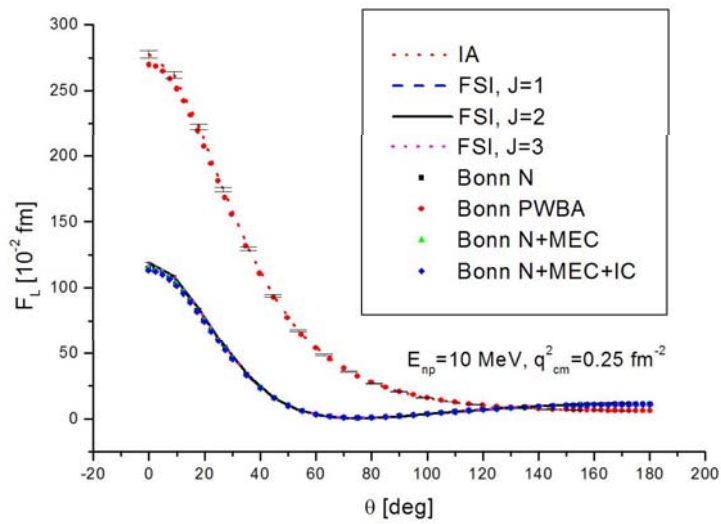
Impulsive approximation (IA)

$$\sim \Psi_d(|\mathbf{p}' - \mathbf{q}/2|)$$

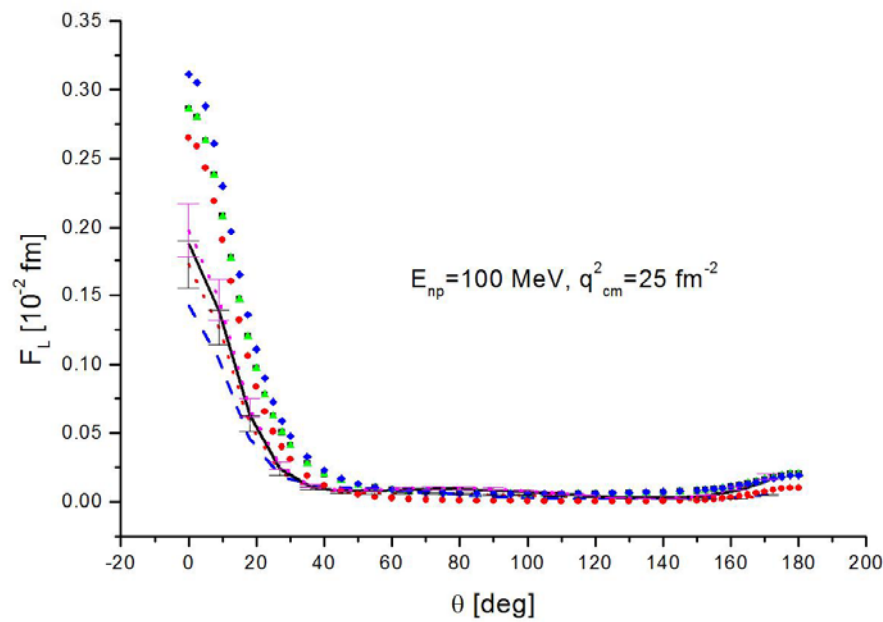
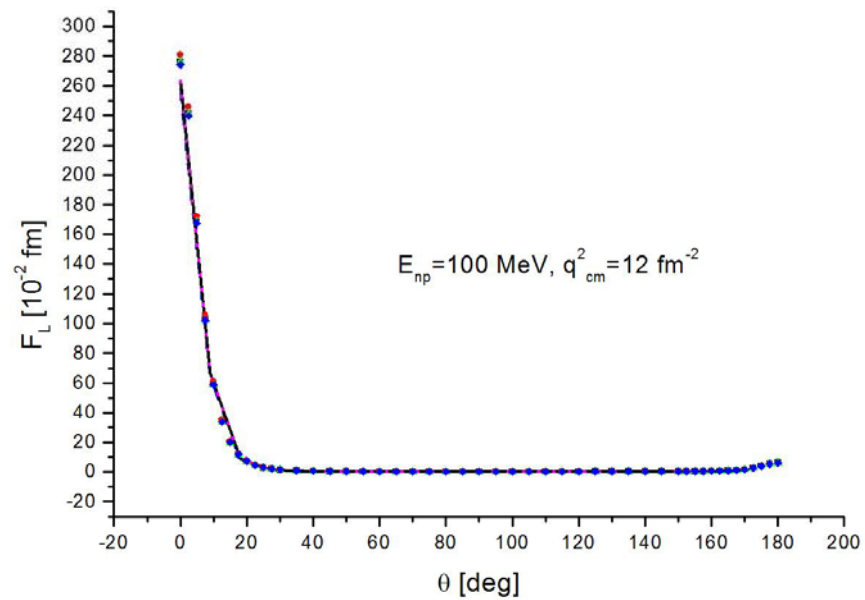
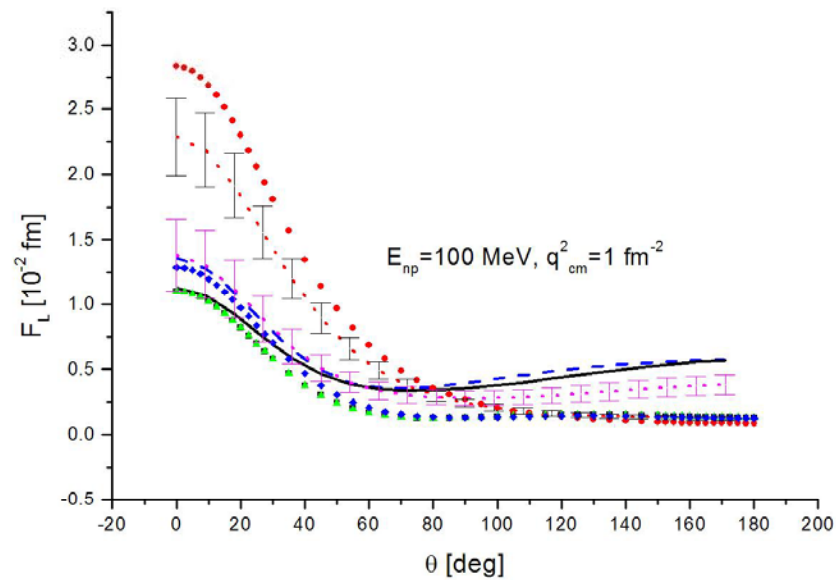
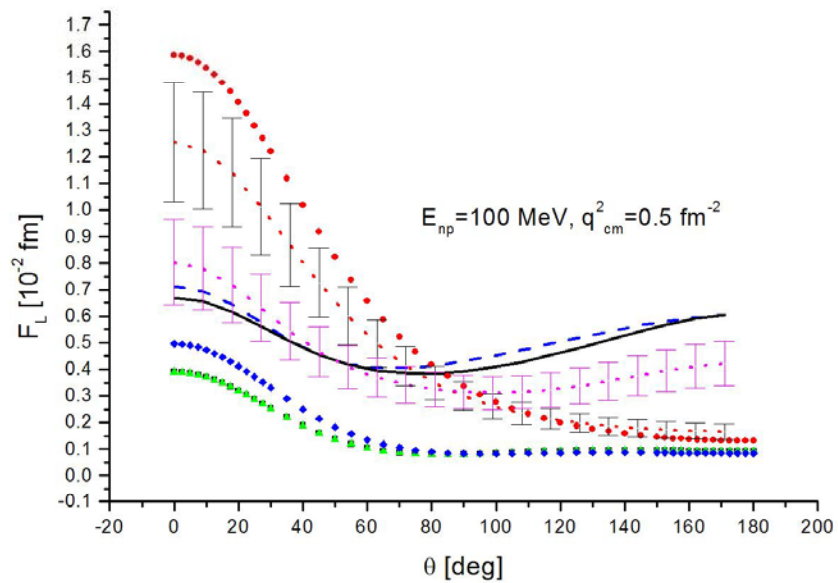
Final state interaction (FSI)

$$\sim \int d^3\mathbf{p} t(\mathbf{p}', \mathbf{p}; E') (E' - \mathbf{p}^2/M + i\epsilon)^{-1} \Psi_d(|\mathbf{p} - \mathbf{q}/2|)$$

# Results







# Conclusion

$E_{np}$ (MeV) \ $q_{cm}^2$ (fm <sup>-2</sup> )	10	30	60	100	120	160
0.25(q~100MeV)	✓ ✓ ✓	✓ ⇕ ✓	⇕ ⇕ ✓	✗ ✗ ✓	✗ ✗ ✓	✗ ✗ ✓
0.5(q~100MeV)	✓ ✓ ✓	✓ ✓ ✓	⇕ ✓ ✓	✗ ✗ ✓	✗ ✗ ✓	✗ ✗ ✓
1(q~100MeV)	✓ ✓ ✗	✓ ✓ ⇕	⇕ ✓ ✓	✗ ✗ ✓	✗ ✗ ✓	✗ ✗ ✓
4(q~100MeV)	⇕ ⇕ ⇕	✓ ✓ ✗	✓ ✓ ✗	⇕ ⇕ ⇕	⇕ ⇕ ⇕	⇕ ⇕ ✓
10(q~100MeV)	✗ ✗ ✓	⇕ ✗ ⇕	✓ ✓ ✗	✓ ✓ ✗	✓ ✓ ✗	✓ ✓ ✗
16(q~100MeV)	✗ ✗ ✓	⇕ ✗ ✓	⇕ ⇕ ⇕	✓ ✓ ✗	✓ ✓ ✗	✓ ✓ ✗
25(q~100MeV)	✗ ✗ ✓	✗ ✗ ✓	✗ ✗ ✓	✗ ✗ ⇕	⇕ ✗ ⇕	⇕ ⇕ ✗

$\Lambda$ -ind.?

Match Bonn result?

FSI important?

# Future Works and application

- Apply the results for the deuteron electro-disintegration calculation.
- Use the same technique to evaluate the NN problem in perturbative way (with Bingwei Long).
- Evaluate the LEC's for the N<sup>3</sup>LO:  
=> eliminate the correlations between them cause by performing a best fit in the phase shift.

## Application:

Our method enables us to obtain an EFT NN potential at “any” cutoff; it is faster (don't need to search for the L.E.C.'s) and with a more reasonable renormalization procedure.

⇒ This  $V_{\text{EFT}}$  can be used in further few body and nuclear structure calculation (evaluate  $\Lambda$ -dep., etc.).

**THANK YOU!**

# Singular potential

Consider time-independent Schrodinger eq. with  $V = \frac{-\lambda}{r^n}$  as  $r \sim 0$ :

$KE \sim \frac{\hbar^2}{mr^2}$ ,  $PE \sim \frac{-\lambda}{r^n}$ , Then for  $n > 2$ , total E can take arbitrarily large negative values for very small r.

$\Rightarrow$  Simplify the equation: (ignore m,  $\hbar$ , set  $x = r/r_0$ )

S wave, radial part with  $E=0$ :  $u'' + \frac{\lambda}{r^n} u = 0$

For  $n=1$ , which is Coulomb, O.K.

For  $n=2$ , general solution is:  $u = Ax^{1/2 + \sqrt{1/4 - \lambda}} + Bx^{1/2 - \sqrt{1/4 - \lambda}}$ ,

so for  $\lambda < 1/4$ , by choosing A, B, we can make stationary state.

for  $\lambda > 1/4$ , **singular potential.**

For  $n > 3$ , **singular potential.**

$$T(0) = V + VG(0)T(0) \quad (5) \quad \text{*Notation: } G(E) \equiv [E - H_0 + i\varepsilon]^{-1}$$

$$T(E) \equiv T(p, p^*, E)$$

$$T = V + VGV + VGVGV + \dots = V + VG(\underbrace{V + VGV + \dots}_T) = V + VGT.$$

$$\text{or, } = V + (\underbrace{V + VGV + \dots}_T)GV = V + TGV$$

(5) becomes  $T(0) = V + T(0)G(0)V$  ←

$$T(E) = V + VG(E)T(E) = V(1 + G(E)T(E)),$$

$$\Rightarrow T(E)(1 + G(E)T(E))^{-1} = V$$

$$T(0) = (1 + T(0)G(0))T(E)(1 + G(0)T(E))^{-1}$$

$$T(0)(1 + G(E)T(E)) = (1 + T(0)G(0))T(E)$$

$$T(0) + T(0)G(E)T(E) = T(E) + T(0)G(0)T(E)$$

$$T(E) = T(0) + T(0)[G(E) - G(0)]T(E)$$

$$[1 + T(0)(G(0) - G(E))]T(E) = T(0)$$

=> Get T(E) from T(0)

