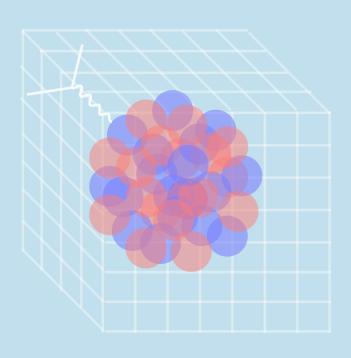
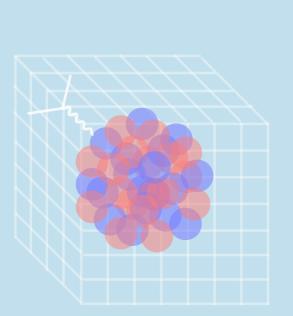
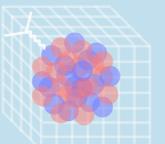
**ZOHREH DAVOUDI** UNIVERSITY OF MARYLAND AND RIKEN FELLOW

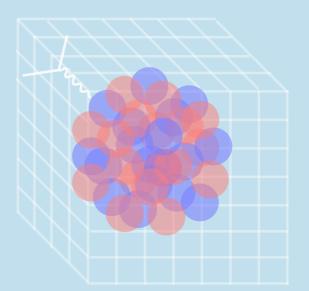
## LATTICE QCD AND NUCLEON(US) STRUCTURE

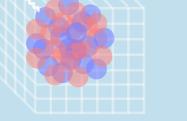
THE 2019 CFNS SUMMER SCHOOL ON THE PHYSICS OF THE ELECTRON ION COLLIDER

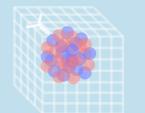


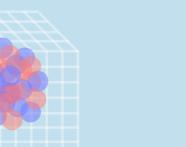


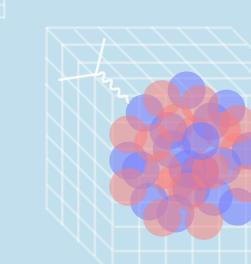










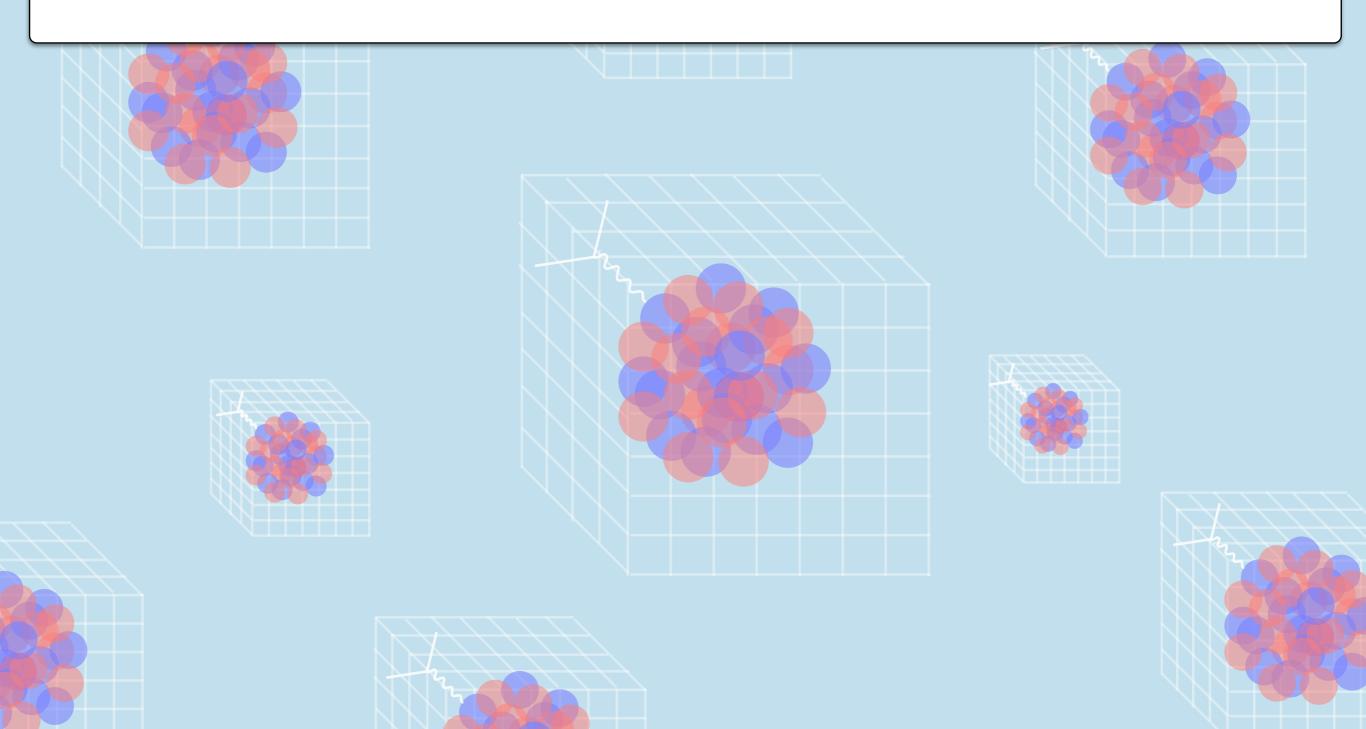


#### LECTUTE I: LATTICE QCD FORMALISM AND METHODOLOGY

#### LECTUTE II: NUCLEON STRUCTURE FROM LATTICE QCD

### LECTUTE III: TOWARDS NUCLEAR STRUCTURE FROM LATTICE QCD

#### LECTUTE I: LATTICE QCD FORMALISM AND METHODOLOGY

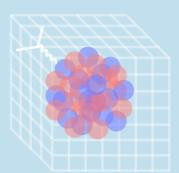


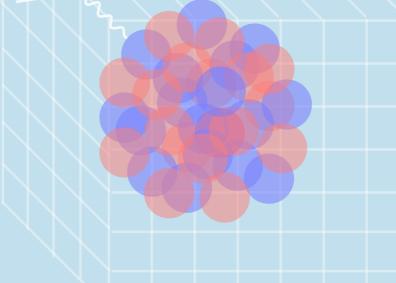
Quantum chromodynamics (QCD) in continuum:

# QCD is a SU(3) Yang-Mills theory augmented with several flavors of massive quarks: Quark kinetic and mass term Quark/gluon interactions $\mathcal{L}_{QCD} = \sum_{f=1}^{N_f} \left[ \bar{q}_f (i\gamma^\mu \partial_\mu - m_f) q_f - g A^i_\mu \bar{q}_f \gamma^\mu T^i q_f \right]$

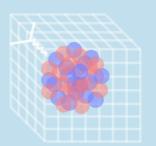
$$-\frac{1}{4}F^{i}_{\mu\nu}F^{i\mu\nu} + \frac{g}{2}f_{ijk}F^{i}_{\mu\nu}A^{i\mu}A^{j\nu} - \frac{g^2}{4}f_{ijk}f_{klm}A^{j}_{\mu}A^{k}_{\nu}A^{l\mu}A^{m\nu}$$

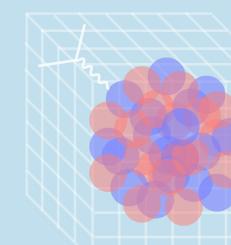
Gluons kinetic and interaction terms



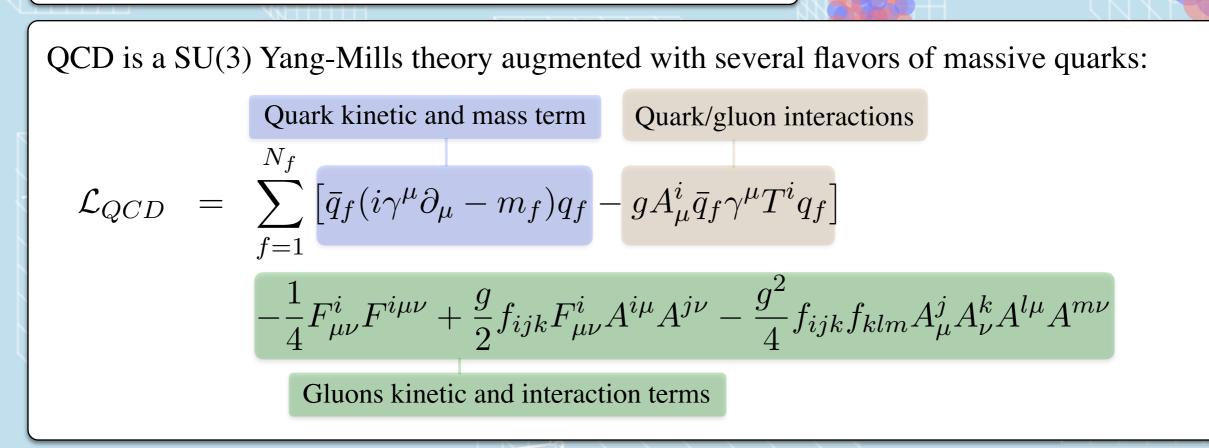








Quantum chromodynamics (QCD) in continuum:



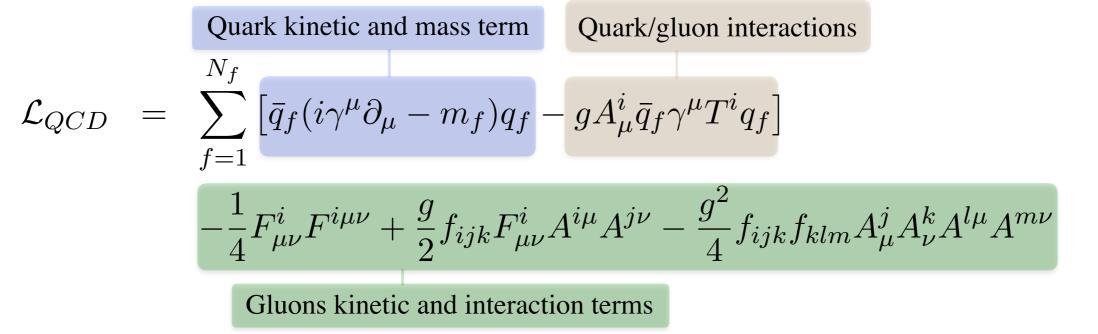
Observe that:

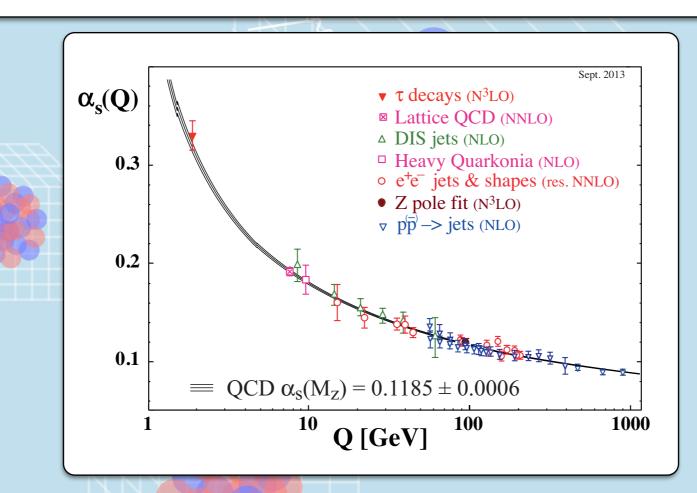
- i) There are only  $1 + N_f$  input parameters plus QCD coupling. Fix them by a few quantities and all strongly-interacting aspects of nuclear physics is predicted (in principle)!
- ii) QCD is asymptotically free such that:  $\alpha_s(\mu') = \frac{1}{2b_0 \log \frac{\mu'}{\Lambda_{OCD}}}$

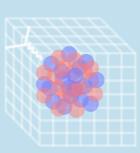
Positive constant for  $N_f \leq 16$ 

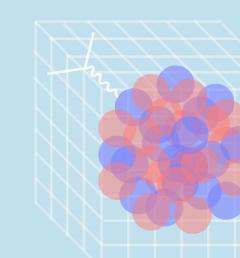
Quantum chromodynamics (QCD) in continuum:

QCD is a SU(3) Yang-Mills theory augmented with several flavors of massive quarks:









Let's enumerate the steps toward numerically simulating this theory nonperturbatively...

**Step I**: Discretize the QCD action in both space and time. Wick rotate to imaginary times. Consider a finite hypercubic lattice.

Step II: Generate a large sample of thermalized decorrelated vacuum configurations.

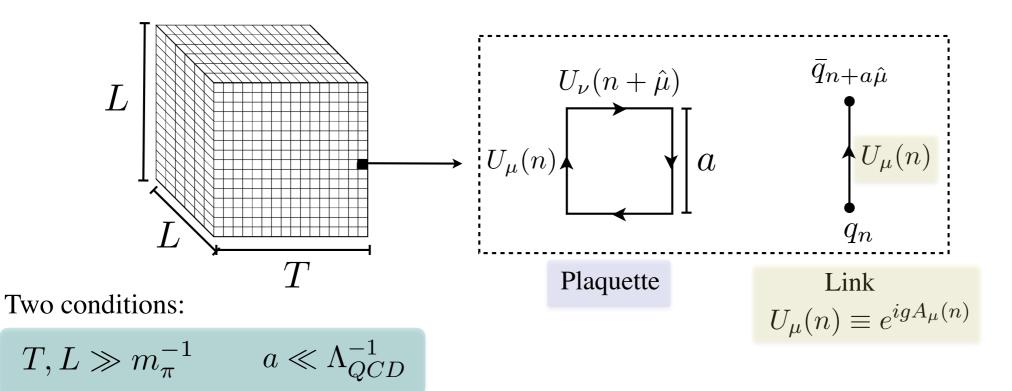
**Step III**: Form the correlation functions by contracting the quarks. Need to specify the interpolating operators for the state under study.

Step IV: Extract energies and matrix elements from correlation functions.

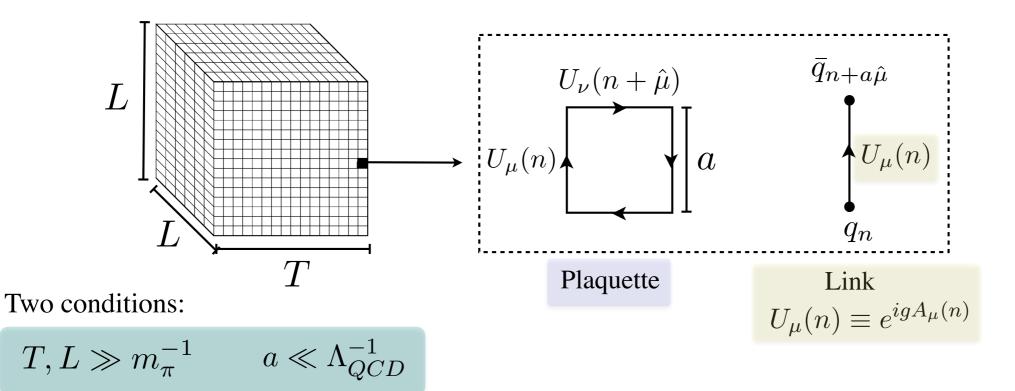
**Step V**: Make the connection to physical observables, such as scattering amplitudes, decay rates, etc.

See e.g., ZD, arXiv:1409.1966 [hep-lat]

**Step I**: Discretize the QCD action in both space and time. Wick rotate to imaginary times. Consider a finite hypercubic lattice.



**Step I**: Discretize the QCD action in both space and time. Wick rotate to imaginary times. Consider a finite hypercubic lattice.



An example of a discretized action by K. Wilson:

$$= 2/g^{2}$$

$$S_{\text{Wilson}}^{(E)} = \frac{\beta}{N_{c}} \sum_{n} \sum_{\mu < \nu} \Re \text{Tr}[\mathbb{1} - P_{\mu\nu;n}] \qquad \text{Wilson parameter. Gives the naive action if set to zero and has doublers problem.}$$

$$- \sum_{n} \bar{q}_{n}[\overline{m}^{(0)} + 4]q_{n} + \sum_{n} \sum_{\mu} \left[ \bar{q}_{n} \frac{r - \gamma_{\mu}}{2} U_{\mu}(n) q_{n+\hat{\mu}} + \bar{q}_{n} \frac{r + \gamma_{\mu}}{2} U_{\mu}^{\dagger}(n-\hat{\mu}) q_{n-\hat{\mu}} \right]$$

For discussions of actions consistent with chiral symmetry of continuum see: Kaplan, arXiv:0912.2560 [hep-lat].

 $\langle \hat{\mathcal{O}} \rangle = \frac{1}{\mathcal{Z}} \int \mathcal{D}U_{\mu} \mathcal{D}q \mathcal{D}\bar{q} \ e^{-S_{\text{lattice}}^{(G)}[U] - S_{\text{lattice}}^{(F)}[U,q,\bar{q}]} \ \hat{\mathcal{O}}[U,q,\bar{q}]$ 

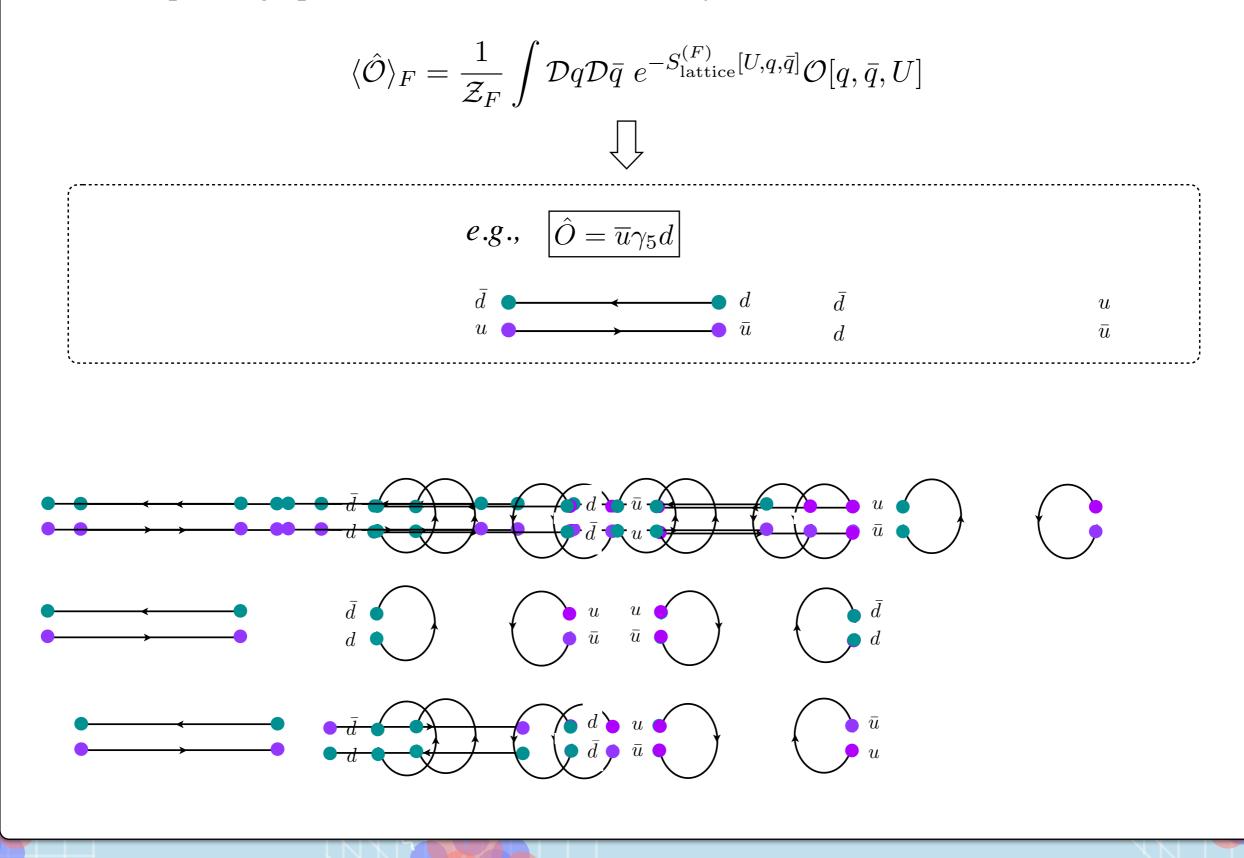
Quark part of expectation values

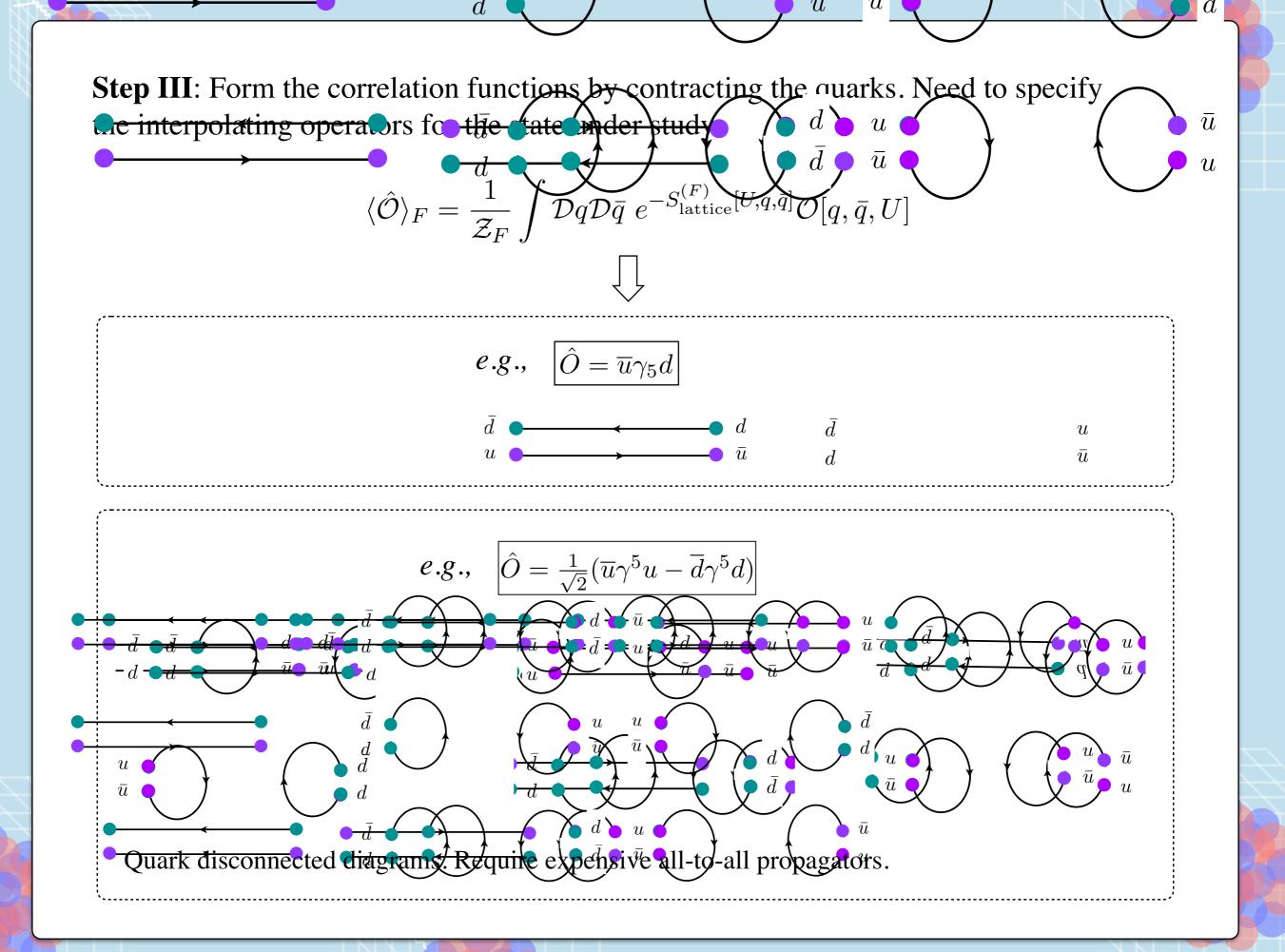
$$\begin{split} &\langle \hat{\mathcal{O}} \rangle = \frac{1}{\mathcal{Z}} \int \mathcal{D}U_{\mu} \mathcal{D}q \mathcal{D}\bar{q} \ e^{-S_{\text{lattice}}^{(G)}[U] - S_{\text{lattice}}^{(F)}[U,q,\bar{q}]} \ \hat{\mathcal{O}}[U,q,\bar{q}] \\ & \text{Quark part of expectation values} \\ \\ & \text{Define: } \langle \hat{\mathcal{O}} \rangle_{F} = \frac{1}{\mathcal{Z}_{F}} \int \mathcal{D}q \mathcal{D}\bar{q} \ e^{-S_{\text{lattice}}^{(F)}[U,q,\bar{q}]} \mathcal{O}[q,\bar{q},U] \\ & \mathcal{Z}_{F} = \int \mathcal{D}q \mathcal{D}\bar{q} \ e^{-S_{\text{lattice}}^{(F)}[U,q,\bar{q}]} = \prod_{f} \det D_{f} \ \text{Dirac matrix} \end{split}$$

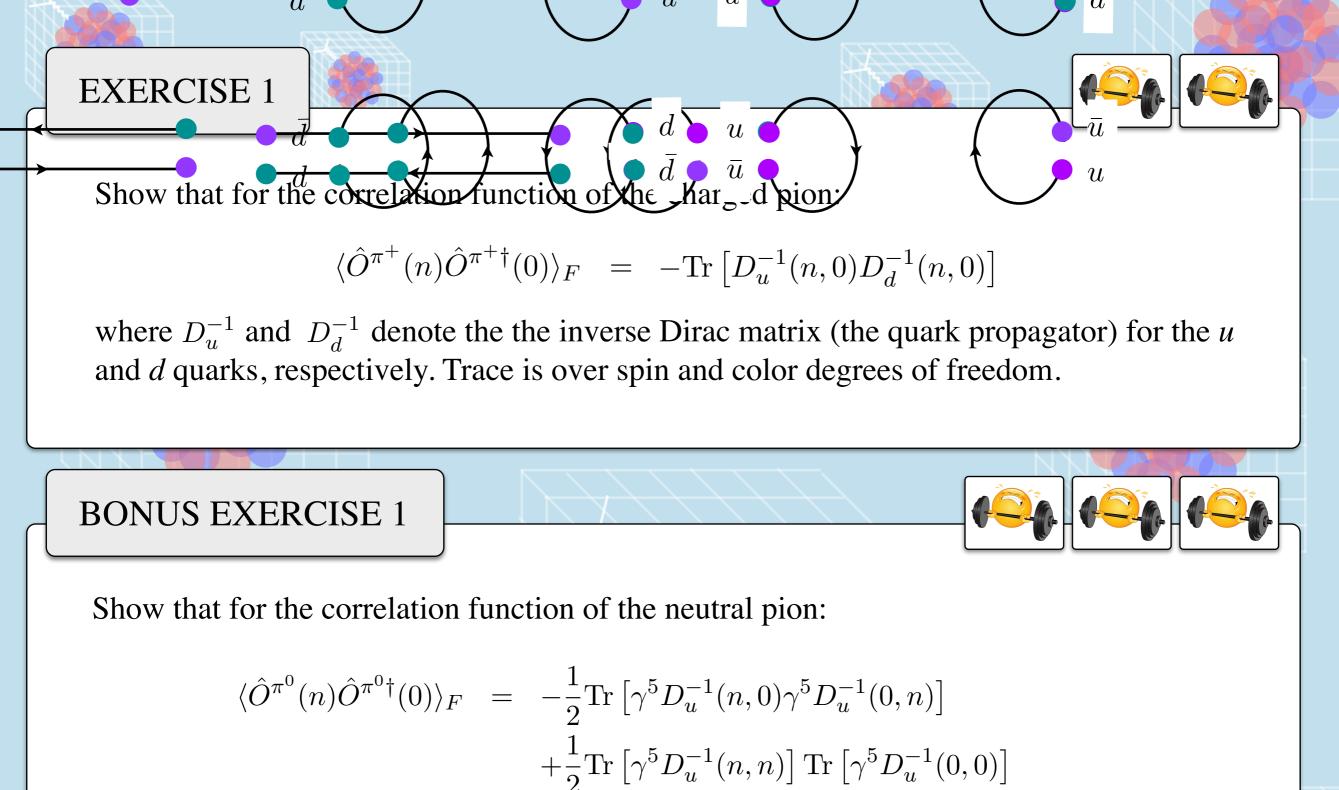
**Step III**: Form the correlation functions by contracting the quarks. Need to specify the interpolating operators for the state under study.

$$\langle \hat{\mathcal{O}} \rangle_F = \frac{1}{\mathcal{Z}_F} \int \mathcal{D}q \mathcal{D}\bar{q} \ e^{-S_{\text{lattice}}^{(F)}[U,q,\bar{q}]} \mathcal{O}[q,\bar{q},U]$$

**Step III**: Form the correlation functions by contracting the quarks. Need to specify the interpolating operators for the state under study.







$$-\frac{1}{2}\operatorname{Tr}\left[\gamma^{5}D_{u}^{-1}(n,n)\right]\operatorname{Tr}\left[\gamma^{5}D_{d}^{-1}(0,0)\right] + \left\{u \leftrightarrow d\right\}$$

First on why steps II and III are expensive...



Example: Consider a lattice with: L/a = 48, T/a = 256

First on why steps II and III are expensive...



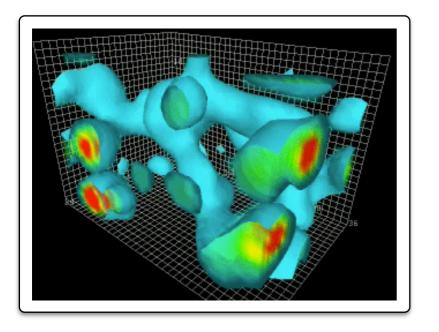
Example: Consider a lattice with: L/a = 48, T/a = 256

Sampling SU(3) matrices. Already for one sample requires storing

 $8 \times 48^3 \times 256 = 226, 492, 416$ 

c-numbers in the computer!

Requires tens of thousands of uncorrelated samples. Molecular-dynamics-inspired hybrid Monte Carlo sampling algorithms often used.



First on why steps II and III are expensive...

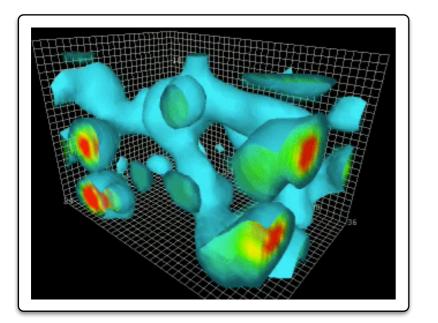
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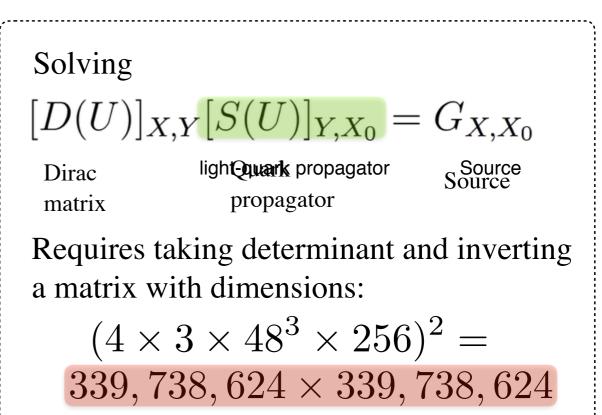
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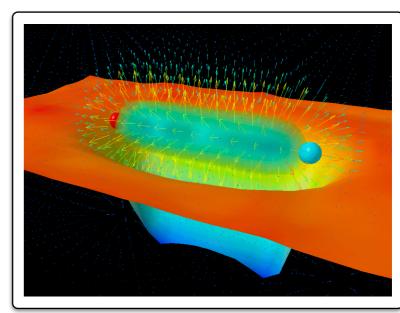
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Step IV: Extract energies and matrix elements from correlation functions

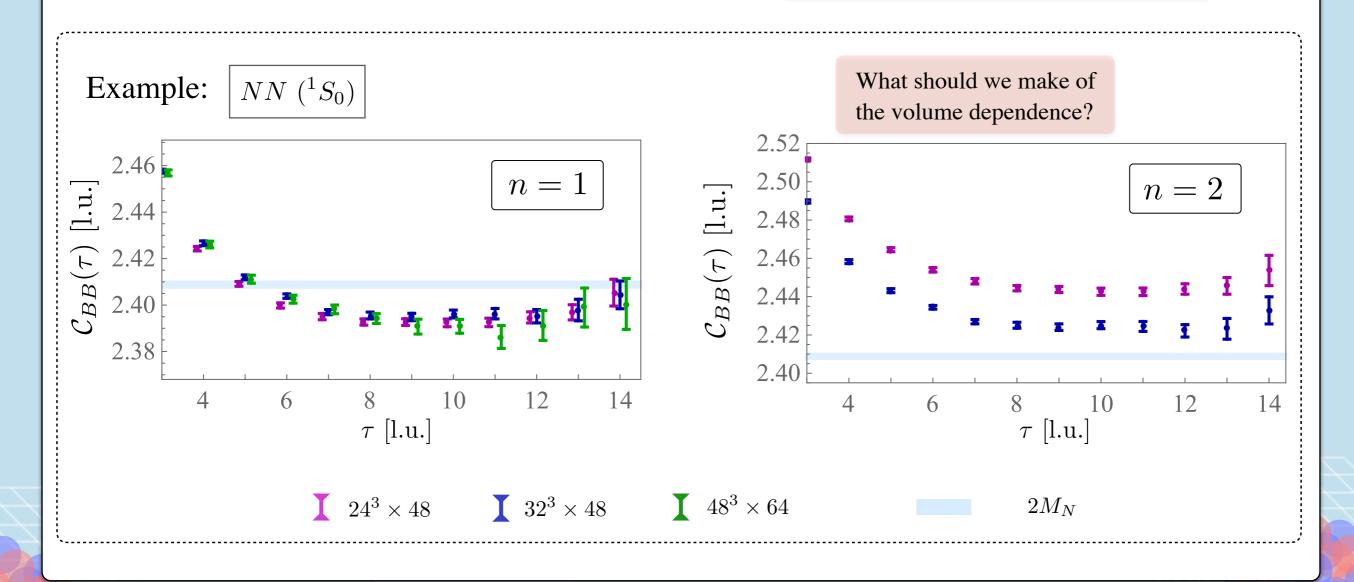
$$C_{\hat{\mathcal{O}},\hat{\mathcal{O}}'}(\tau;\mathbf{d}) = \sum_{\mathbf{x}} e^{2\pi i \mathbf{d} \cdot \mathbf{x}/L} \langle 0 | \hat{\mathcal{O}}'(\mathbf{x},\tau) \hat{\mathcal{O}}^{\dagger}(\mathbf{0},0) | 0 \rangle = \mathcal{Z}_{0}' \mathcal{Z}_{0}^{\dagger} e^{-E^{(0)}\tau} + \mathcal{Z}_{1}' \mathcal{Z}_{1}^{\dagger} e^{-E^{(1)}\tau} + \dots$$

Ground state and a tower of excited states are, in principle, accessible!

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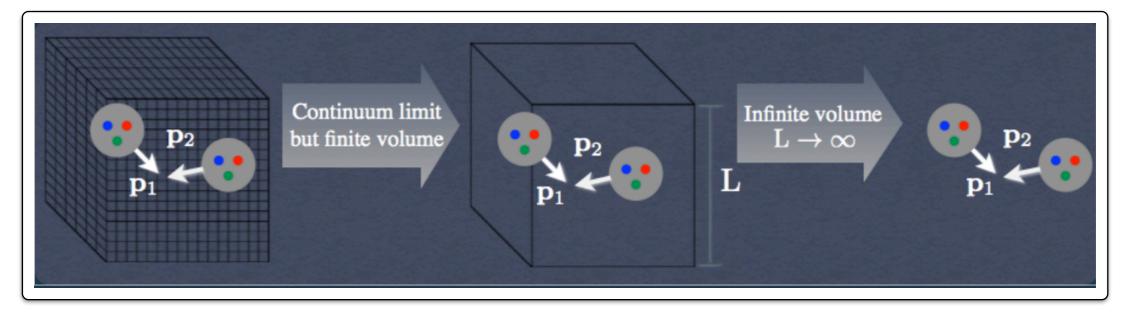
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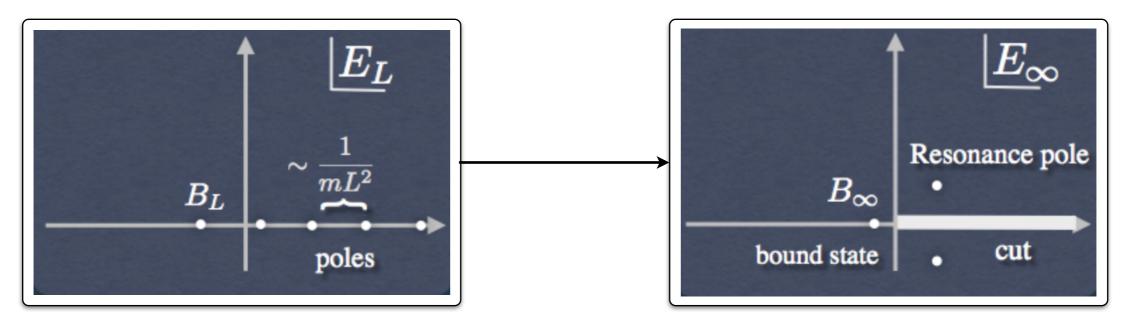


Beane et al (NPLQCD), arXiv:1705.09239, Wagman et al (NPLQCD), arXiv:1706.06550.

**Step V**: Make the connection to physical observables, such as scattering amplitudes, decay rates, etc. Still not fully developed and presents challenge in multi-hadron systems.

#### Example: two-hadron scattering

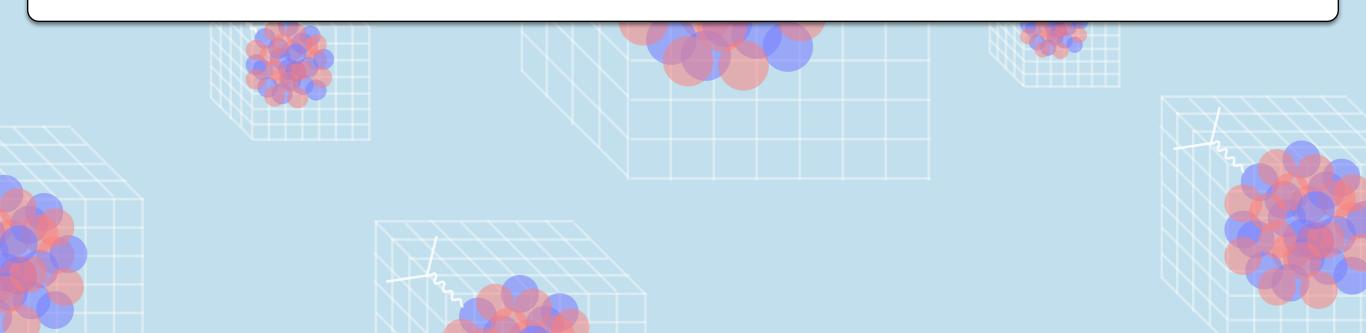




Let's discuss in greater depth step V:

**Step V**: make the connection to physical observables, such as scattering amplitudes, decay rates, etc.

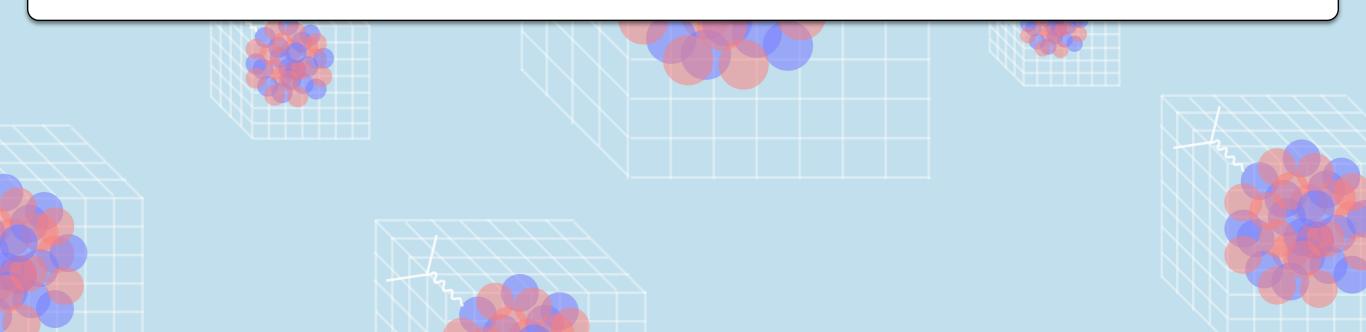
- i) Finite-volume effects in the single-hadron sector
- ii) Finite-volume formalism for two-hadron elastic scattering
- iii) Finite-volume formalism for coupled-channel two-hadron elastic scattering and resonances
- iv) Finite-volume formalism for transition amplitudes and resonance form factors
- v) Finite-volume formalism for three-hadron scattering and resonances
- vi) Finite-volume effects in lattice QED+QCD studies of hadrons



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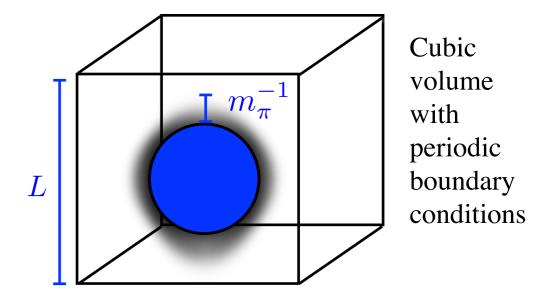
- v) Finite-volume formalism for three-hadron scattering and resonances
- vi) Finite-volume effects in lattice QED+QCD studies of hadrons



QCD finite-volume corrections to single-hadron observables are exponentially suppressed.

Example: The mass of the nucleon

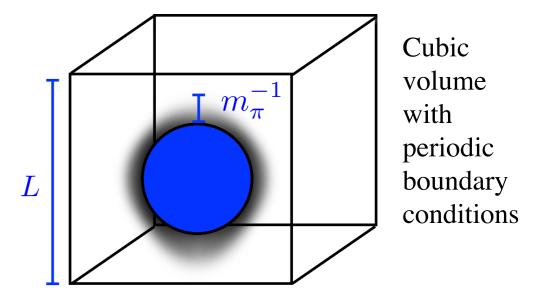
Finite-volume effects are IR in nature. Can use an effective field theory description to characterize them.

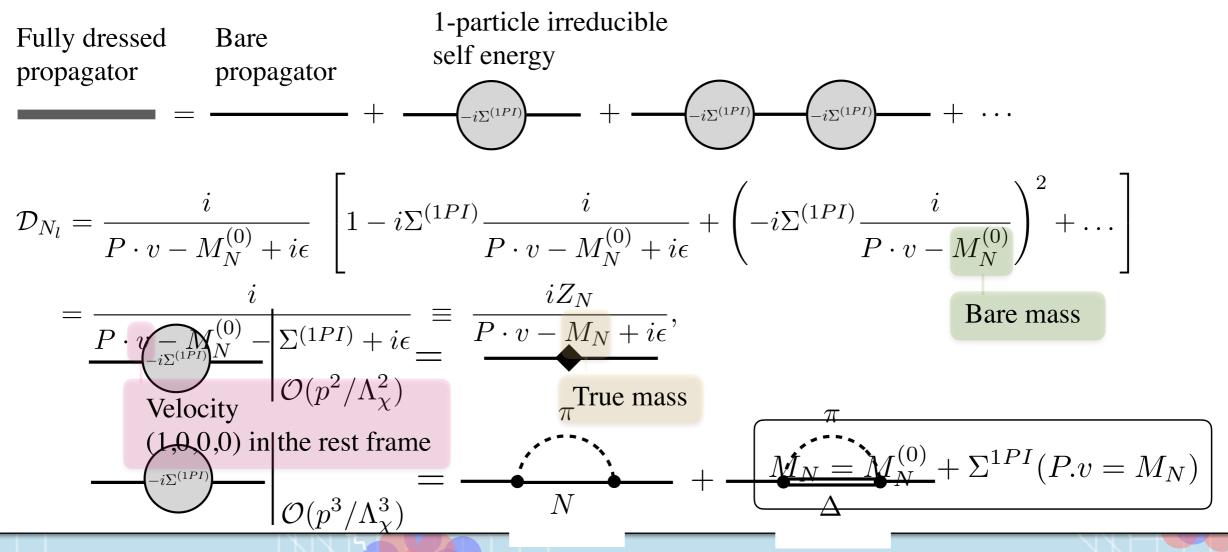


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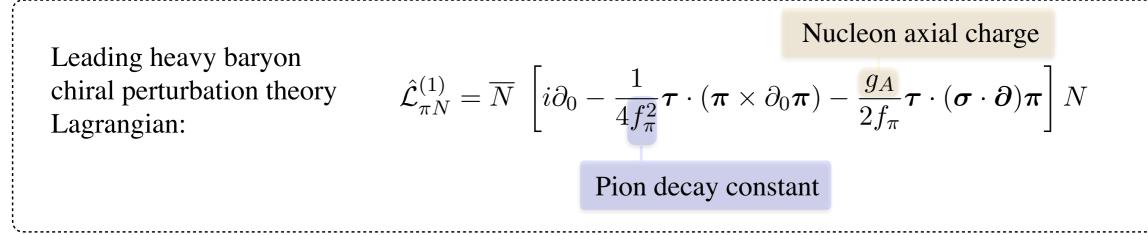
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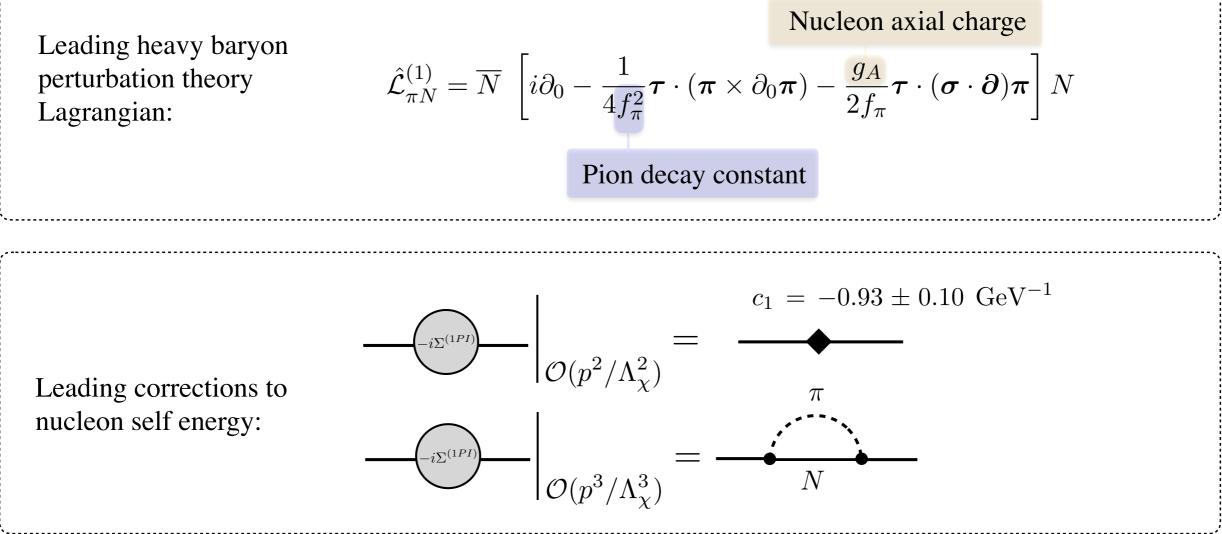


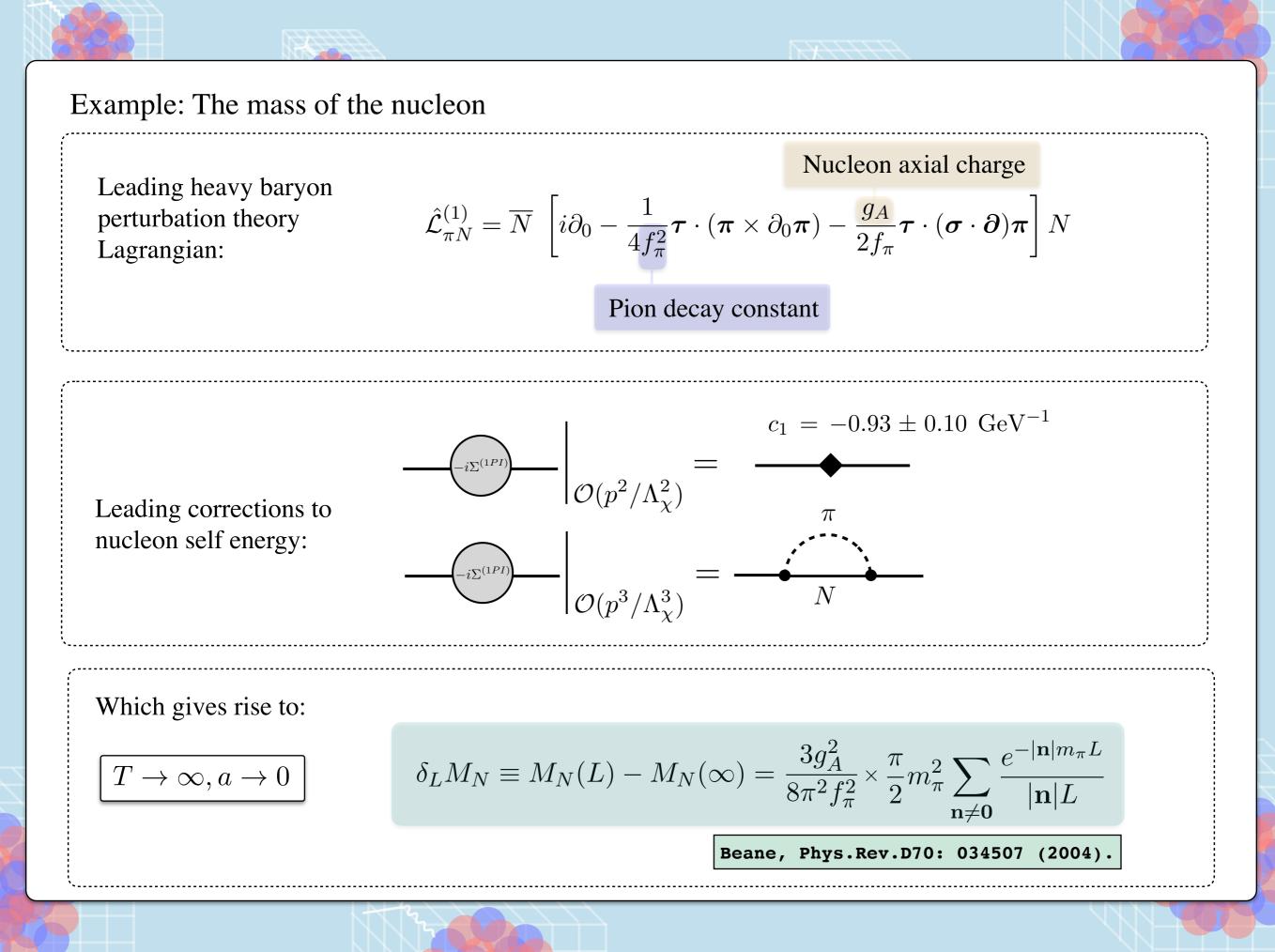


#### Example: The mass of the nucleon



## Example: The mass of the nucleon





#### EXERCISE 2

Plot the first few terms in the expression for the corrections to the nucleon mass as a function of the spatial extent of the volume. How large the volume must be such that correction to the mass of the nucleon are sub-percent?

**BONUS EXERCISE 2** 

Drive the expression for the volume corrections to the mass of the nucleon at leading order in heavy-baryon chiral perturbation theory. The first step is to realize that the volume corrections arise from the loop integral where the integration over a continuous momentum is replaced by a summation over discretized momenta in a periodic cubic volume:

$$\int \frac{dk_0}{2\pi} \frac{d^3k}{(2\pi)^3} \to \int \frac{dk_0}{2\pi} \frac{1}{L^3} \sum_{\mathbf{k}=2\pi\mathbf{n}/L} \text{with } \mathbf{n} \in \mathbb{Z}$$

The second step is to make use of Poisson resumption formula:

$$\frac{1}{L^3} \sum_{k} f(\mathbf{k}) = \int \frac{d^3k}{(2\pi)^3} f(\mathbf{k}) + \sum_{m \neq 0} \int \frac{d^3k}{(2\pi)^3} f(\mathbf{k}) e^{i\mathbf{k}.\mathbf{m}L}$$

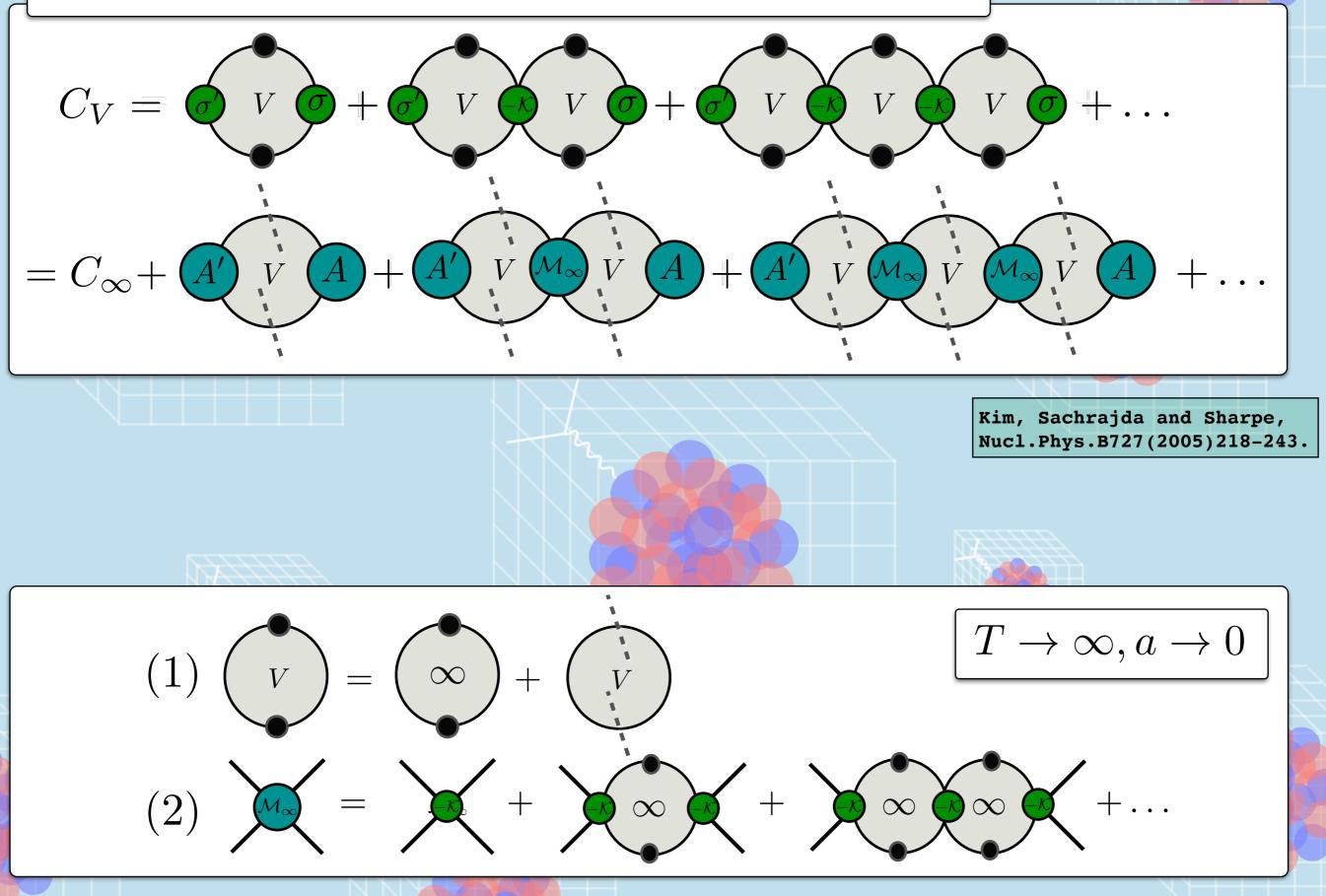
where m is another integer three-vector.

**Step V**: make the connection to physical observables, such as scattering amplitudes, decay rates, etc.

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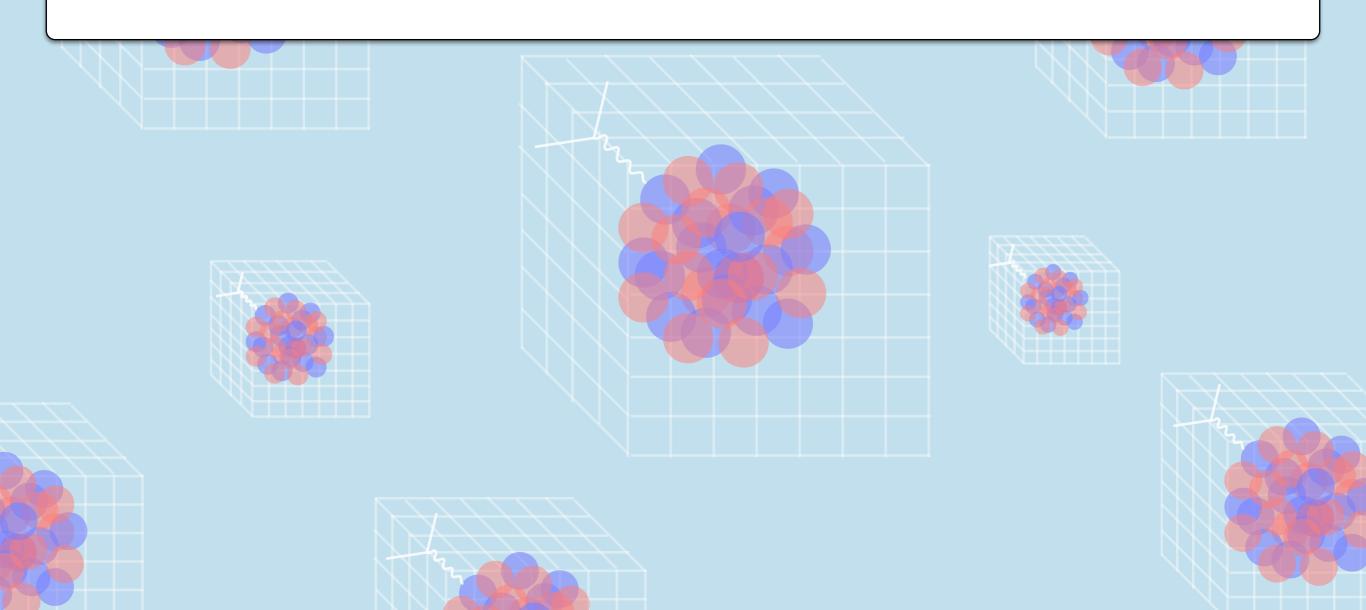


Let's derive the Luescher's formula first. A QFT derivation goes as follows:

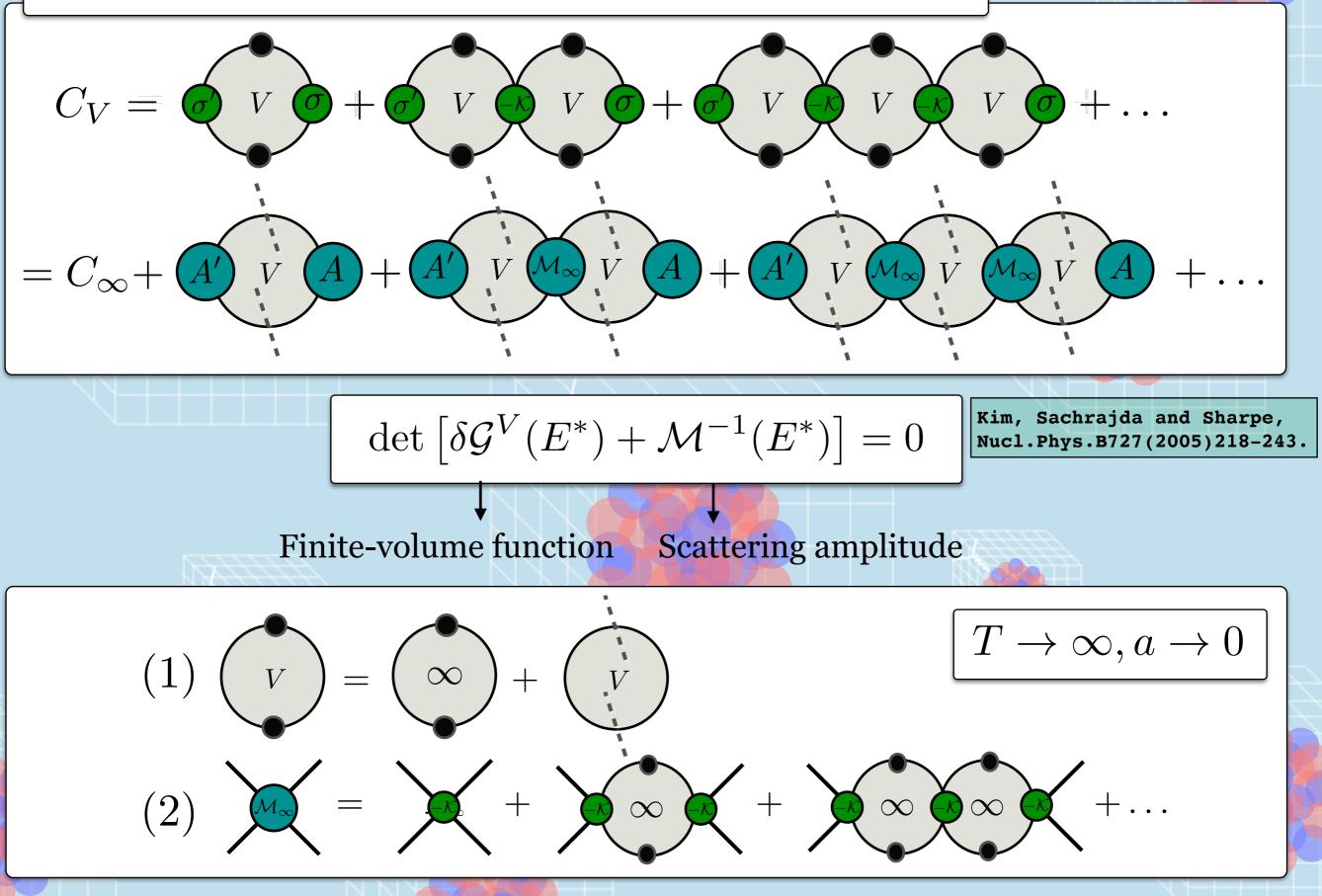


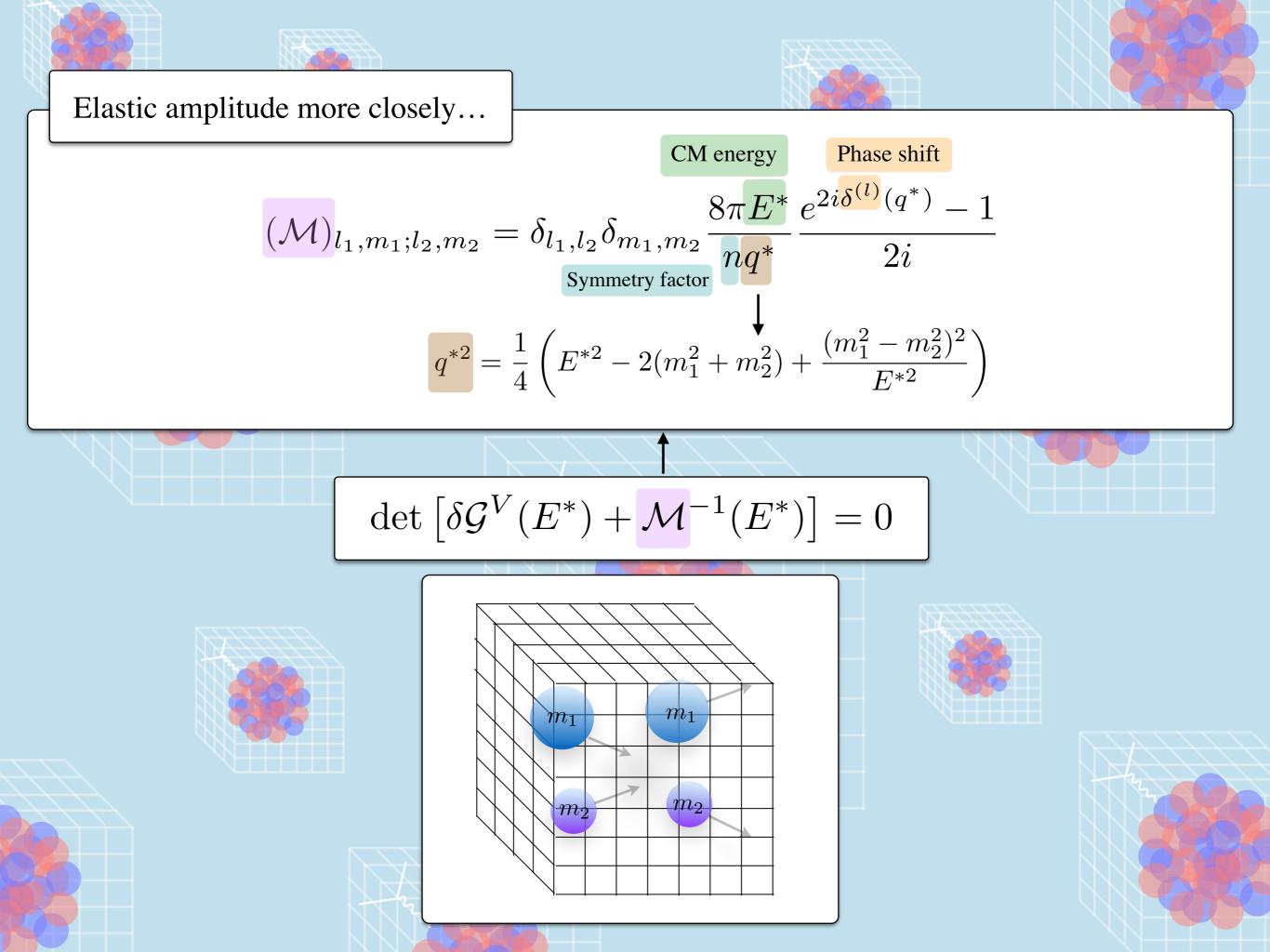


By rearranging the diagrams in  $C_V$  (the first line in the upper panel) using the relations in the lower panel, verify the expansion in the second line in the upper panel. What is the relation between  $\sigma(\sigma')$  and A(A')?



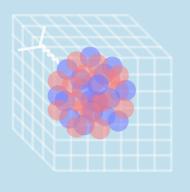
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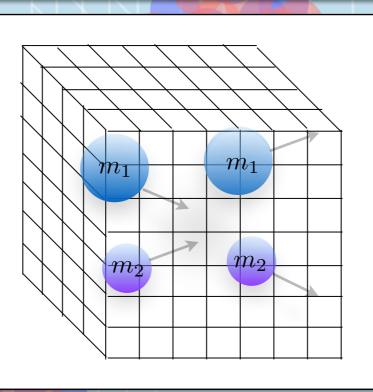


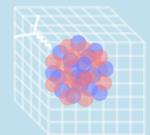


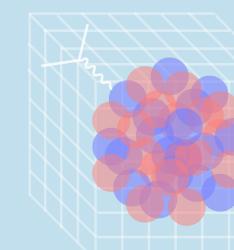
Finite-volume function more closely...

$$\det \left[ \delta \mathcal{G}^V(E^*) + \mathcal{M}^{-1}(E^*) \right] = 0$$

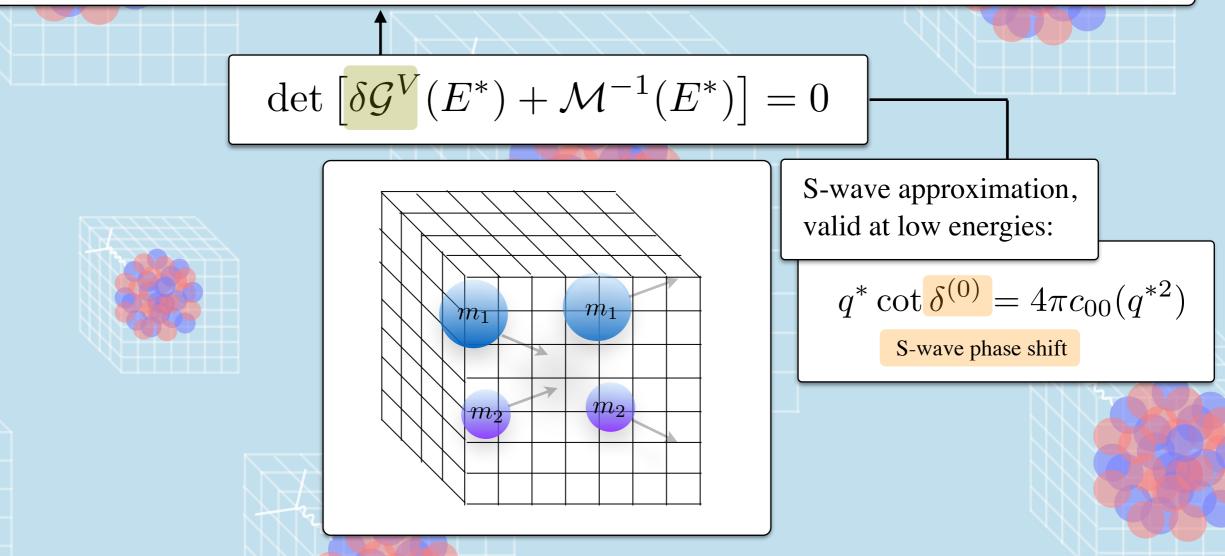


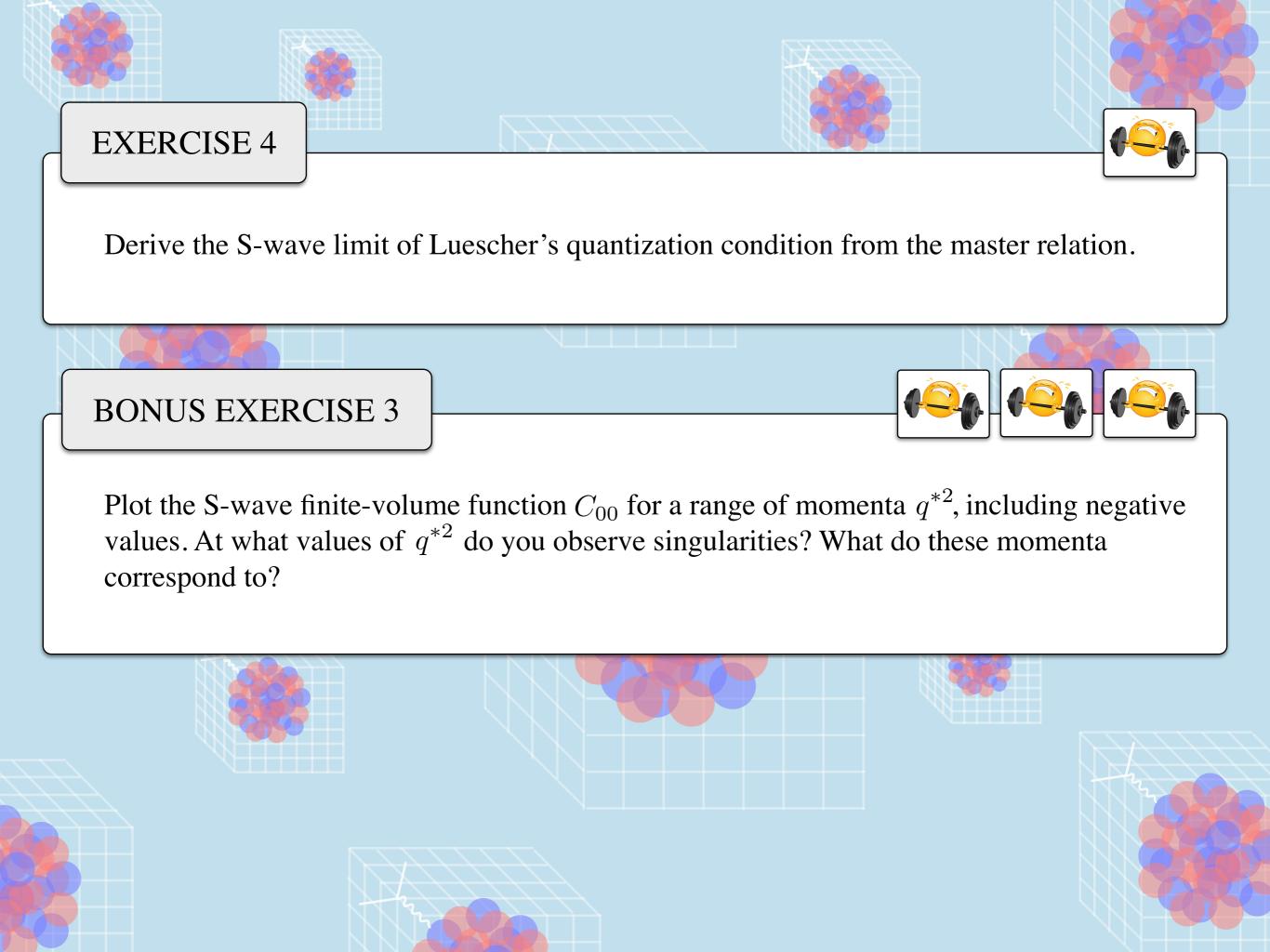






Finite-volume function more closely...





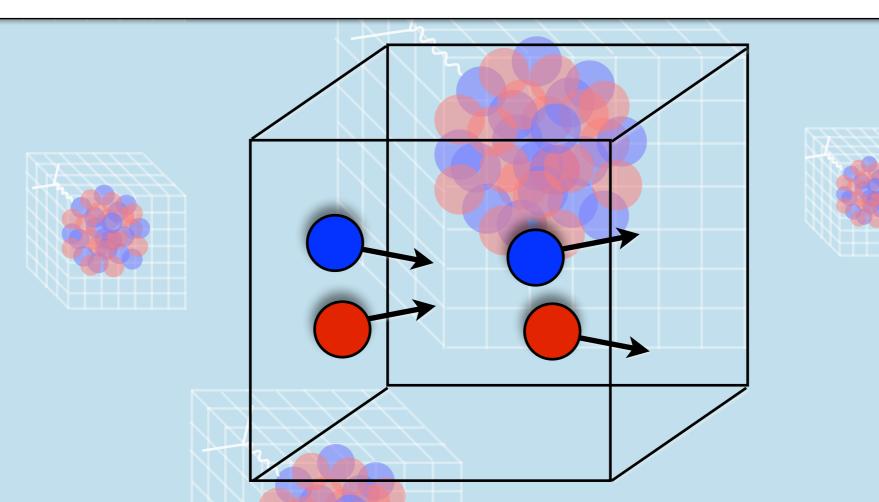
Now let's see an application of Luescher's method to obtain elastic scattering amplitudes of two hadrons from lattice QCD: Wagman et al.(NPLQCD), Phys.Rev.D 96,114510(2017).

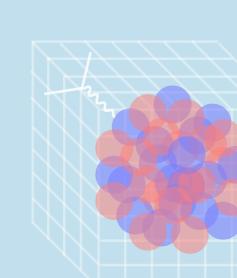
Two-baryon states with SU(3) symmetry

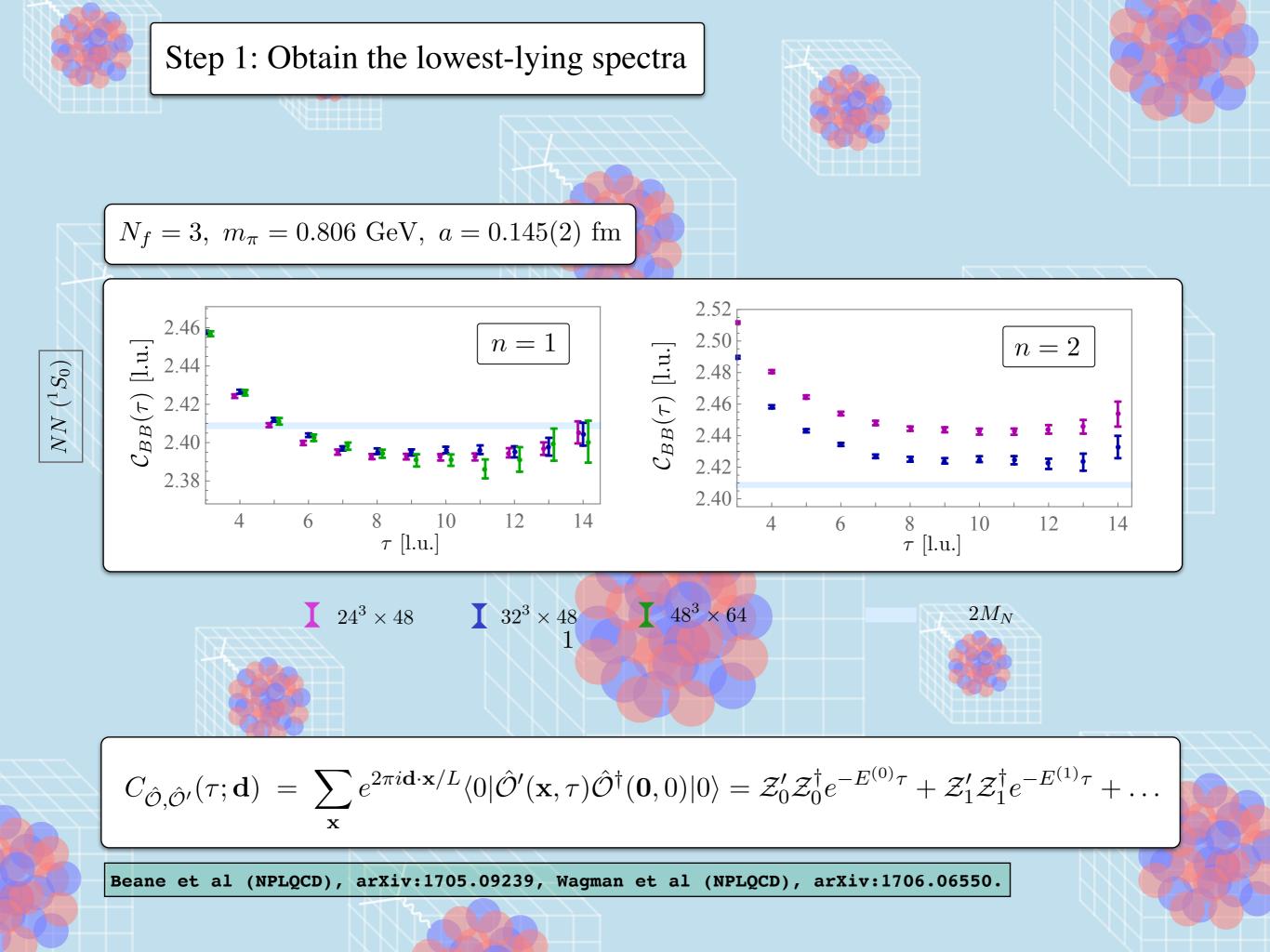
$$\{n, p, \Sigma^+, \Sigma^0, \Sigma^-, \Xi^0, \Xi^+, \Lambda\}$$

SU(3) decomposition of states:  $8 \otimes 8 = 27 \oplus 10 \oplus \overline{10} \oplus 8_S \oplus 8_A \oplus 1$ 

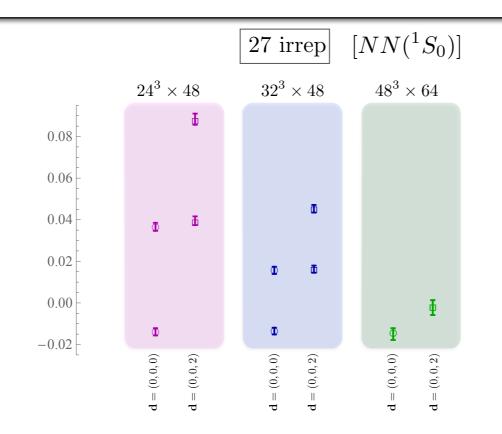
Let's see what these states are...

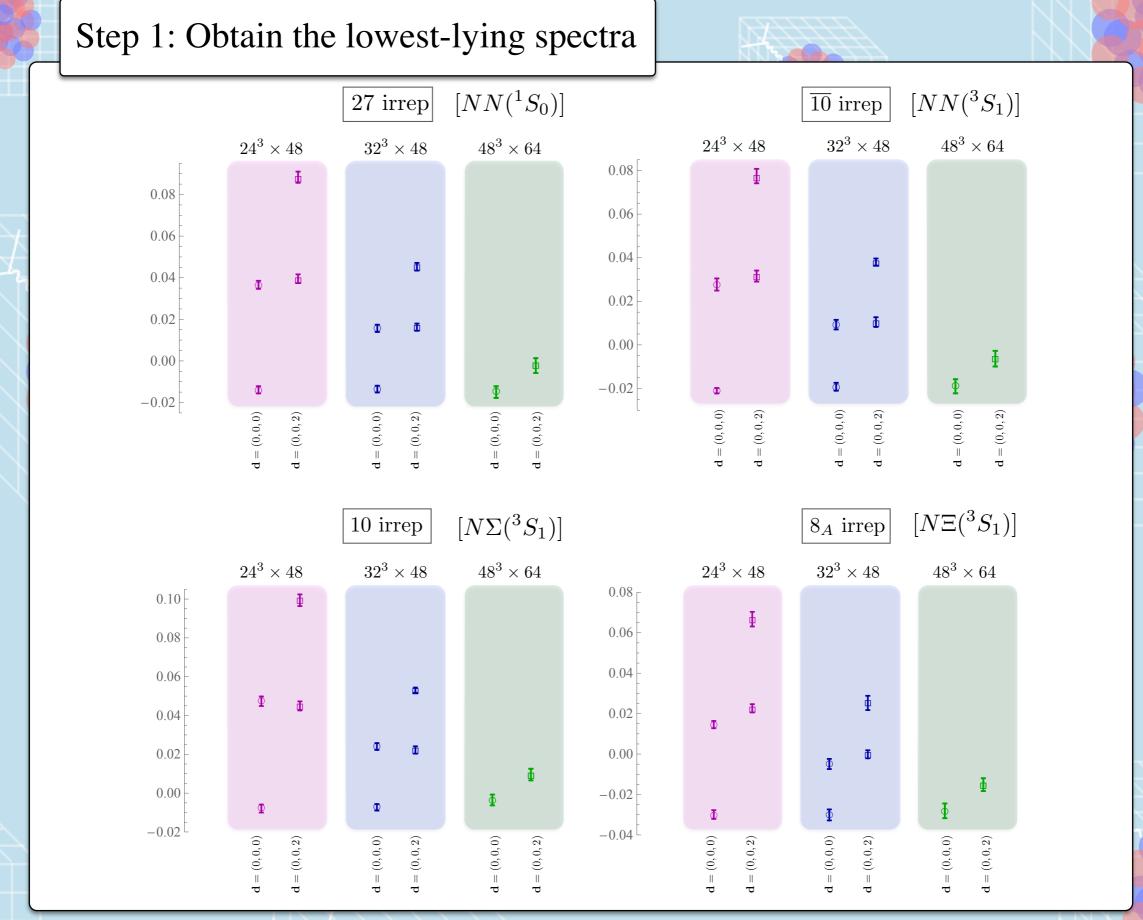






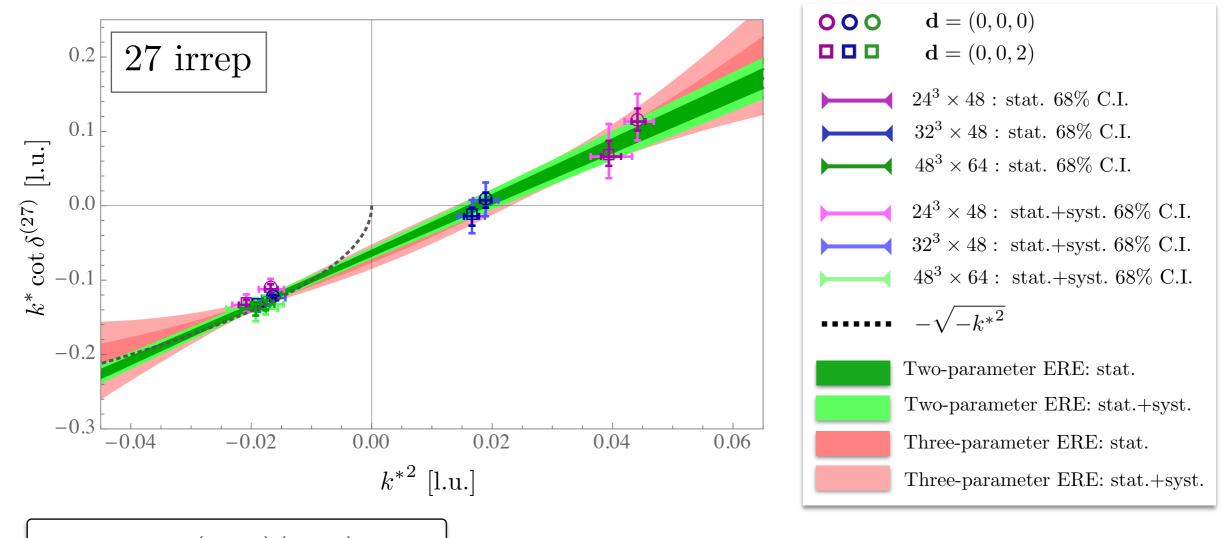
### Step 1: Obtain the lowest-lying spectra





Step 2: Feed the energies to the Luescher's equation and obtain the S-wave scattering phase shifts.

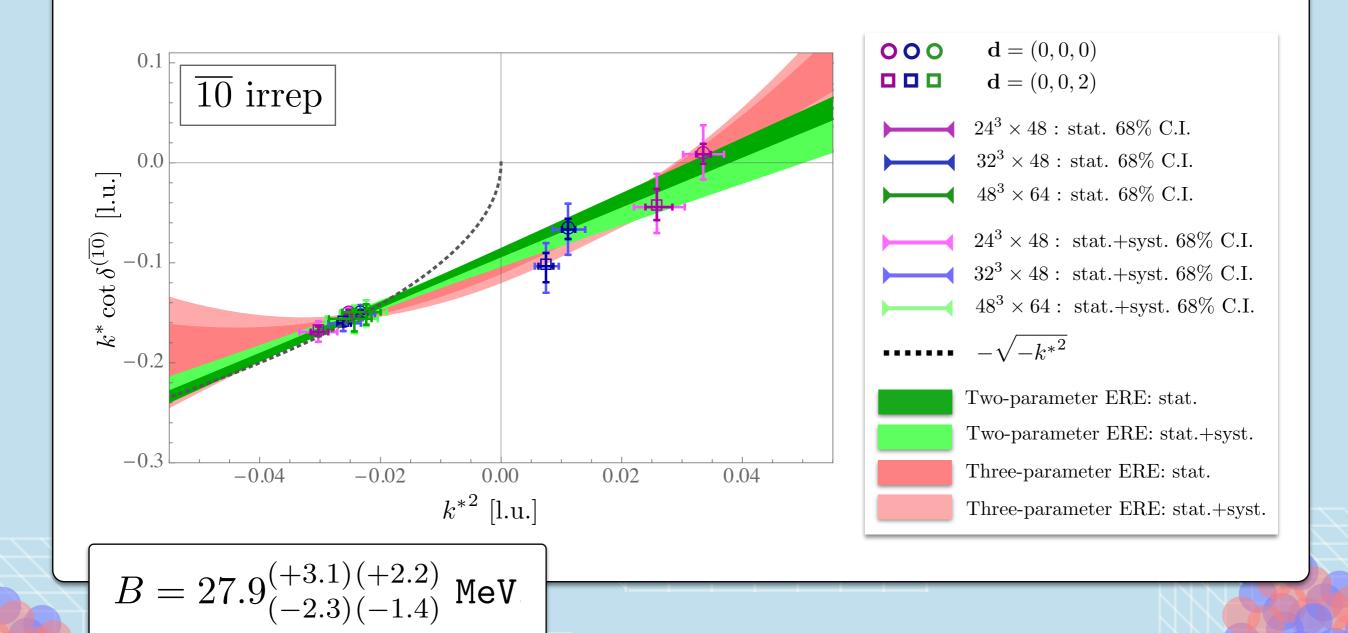
$$N_f = 3, \ m_\pi = 0.806 \text{ GeV}, \ a = 0.145(2) \text{ fm}$$



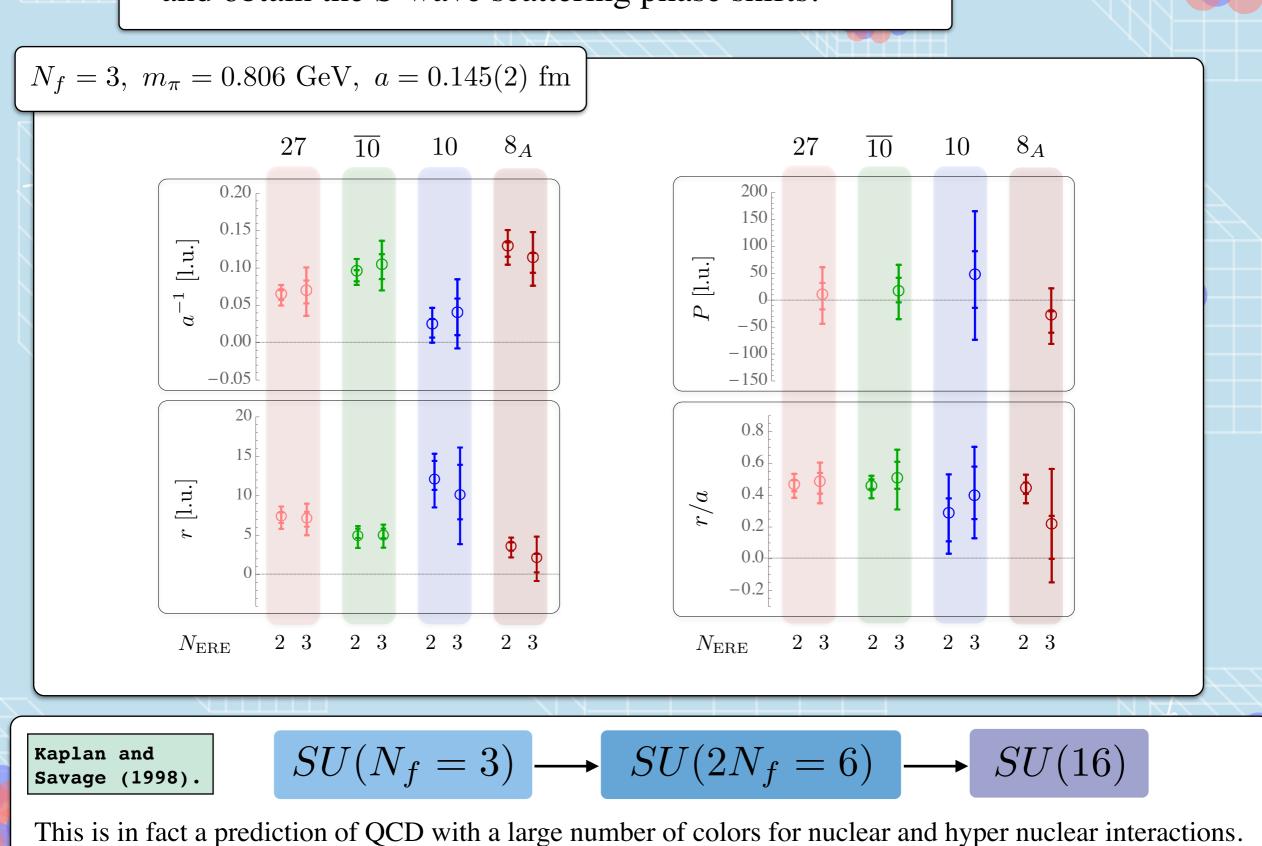
$$B = 20.6^{(+1.8)(+2.8)}_{(-2.4)(-1.6)} \; {
m MeV}$$

Step 2: Feed the energies to the Luescher's equation and obtain the S-wave scattering phase shifts.

$$N_f = 3, \ m_\pi = 0.806 \text{ GeV}, \ a = 0.145(2) \text{ fm}$$



Step 3: Feed the energies to the Luescher's equation and obtain the S-wave scattering phase shifts.



Let's discuss in greater depth step V:

**Step V**: make the connection to physical observables, such as scattering amplitudes, decay rates, etc.

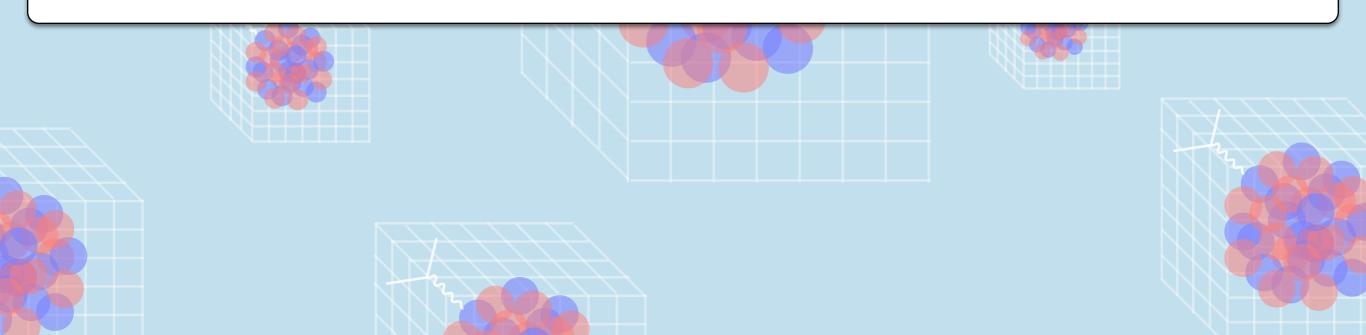
i) Finite-volume effects in the single-hadron sector

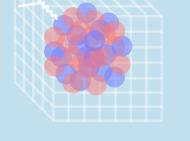
ii) Finite-volume formalism for two-hadron elastic scattering

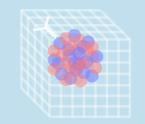
iii) Finite-volume formalism for coupled-channel two-hadron elastic scattering and resonances

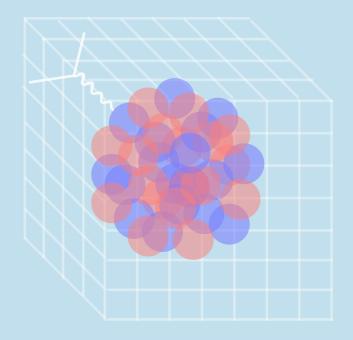
iv) Finite-volume formalism for transition amplitudes and resonance form factors

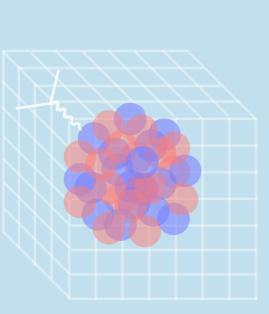
- v) Finite-volume formalism for three-hadron scattering and resonances
- vi) Finite-volume effects in lattice QED+QCD studies of hadrons



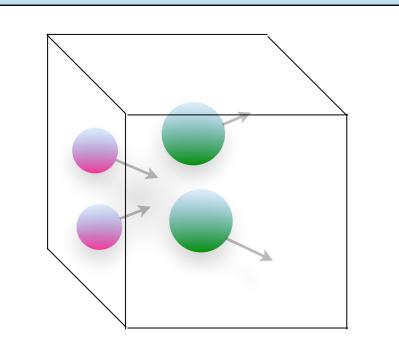


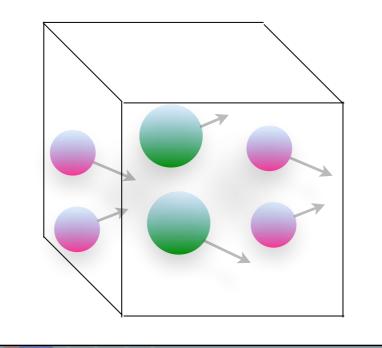


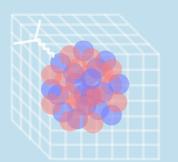


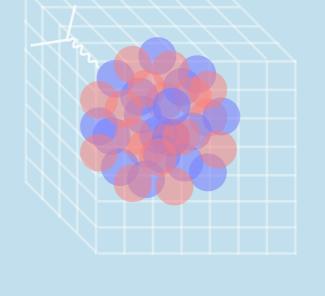


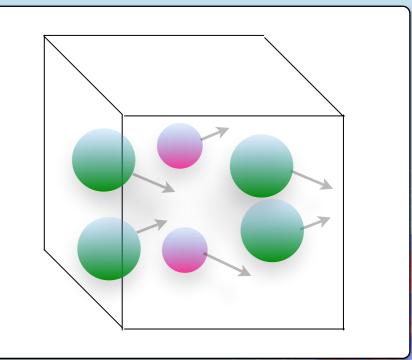




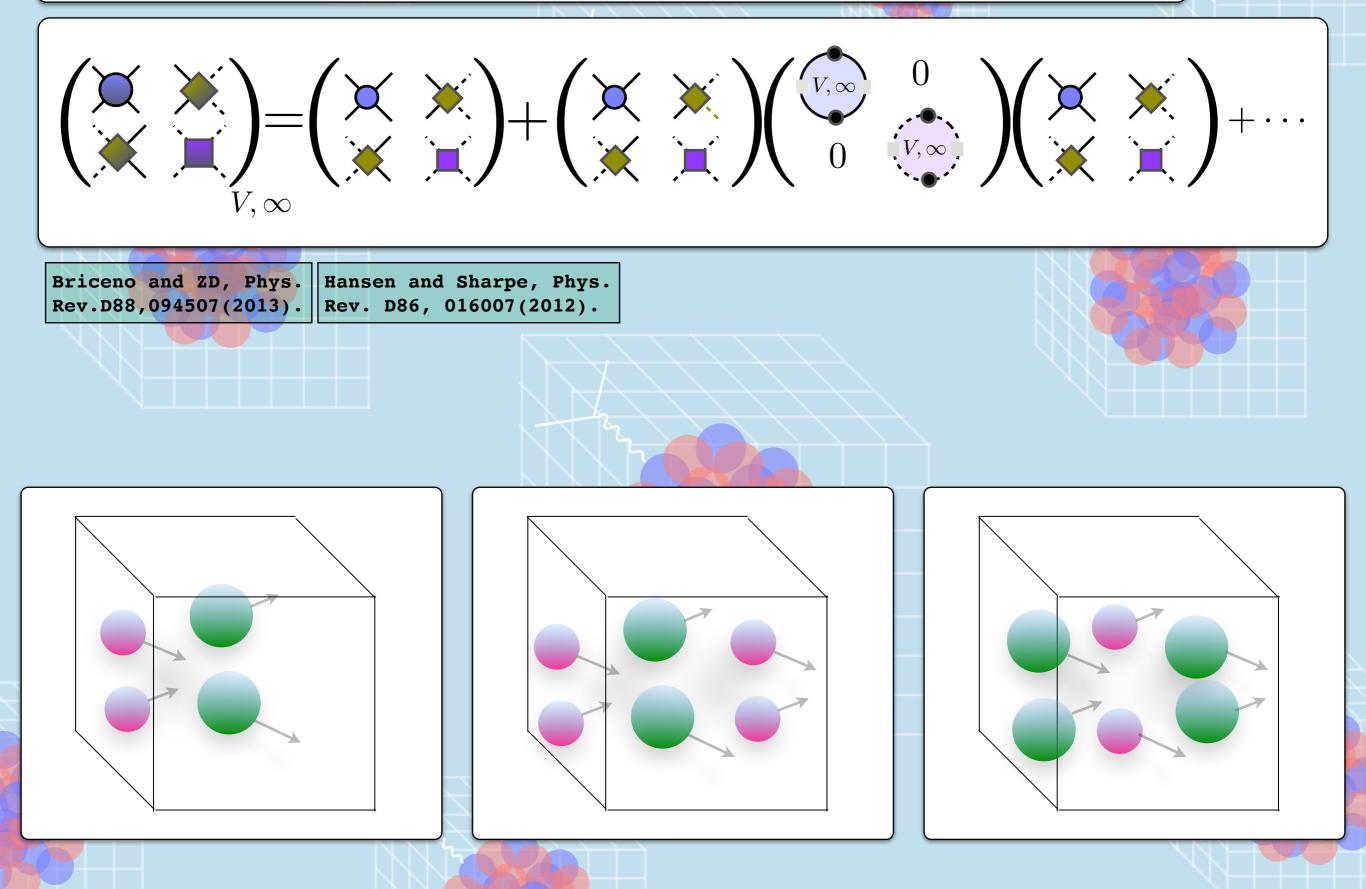




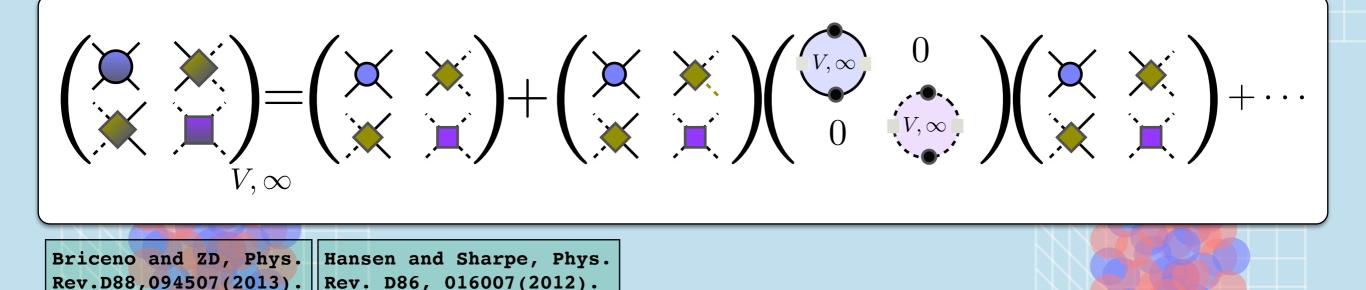




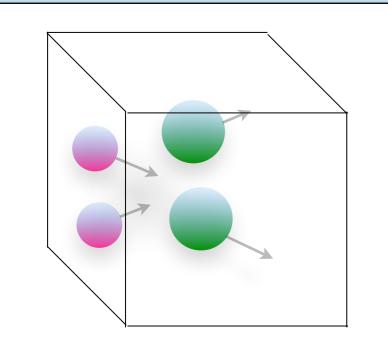
Coupled-channel generalization of luescher's formula is straightforward. Requires upgrading amplitudes and finite-volume functions to matrices in the channel space:

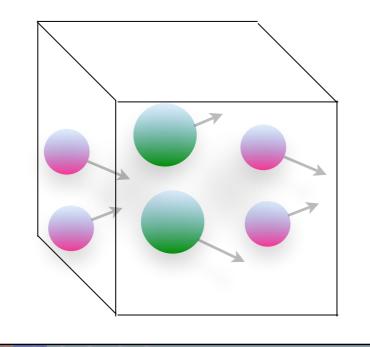


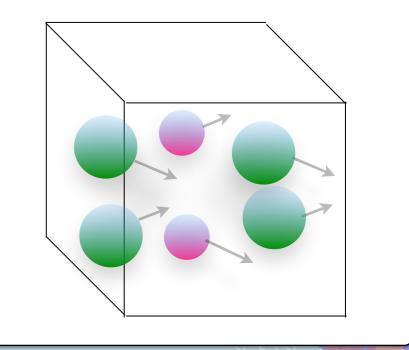
Coupled-channel generalization of luescher's formula is straightforward. Requires upgrading amplitudes and finite-volume functions to matrices in the channel space:



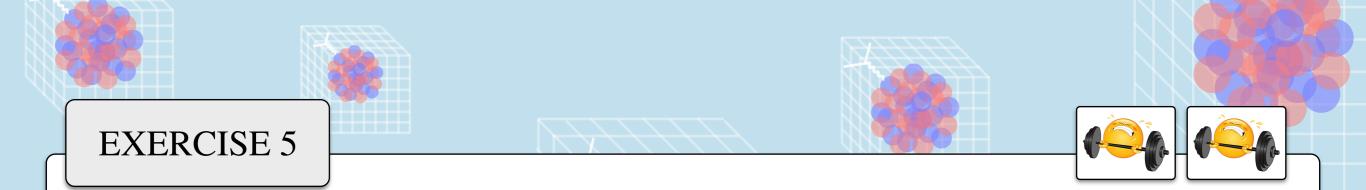
 $\operatorname{Det}\left[\delta\mathcal{G}^{V}(E^{*}) + \mathcal{M}^{-1}(E^{*})\right] = 0$ 







Coupled-channel generalization of luescher's formula is straightforward. Requires upgrading amplitudes and finite-volume functions to matrices in the channel space:



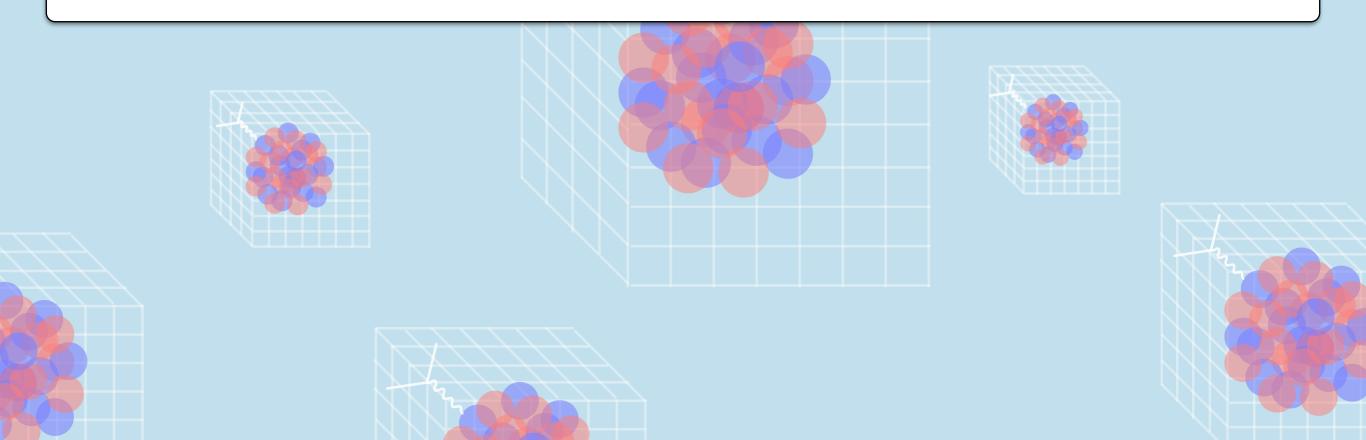
Derive the manifestly real form of a coupled two-channel scattering in the S-wave limit:

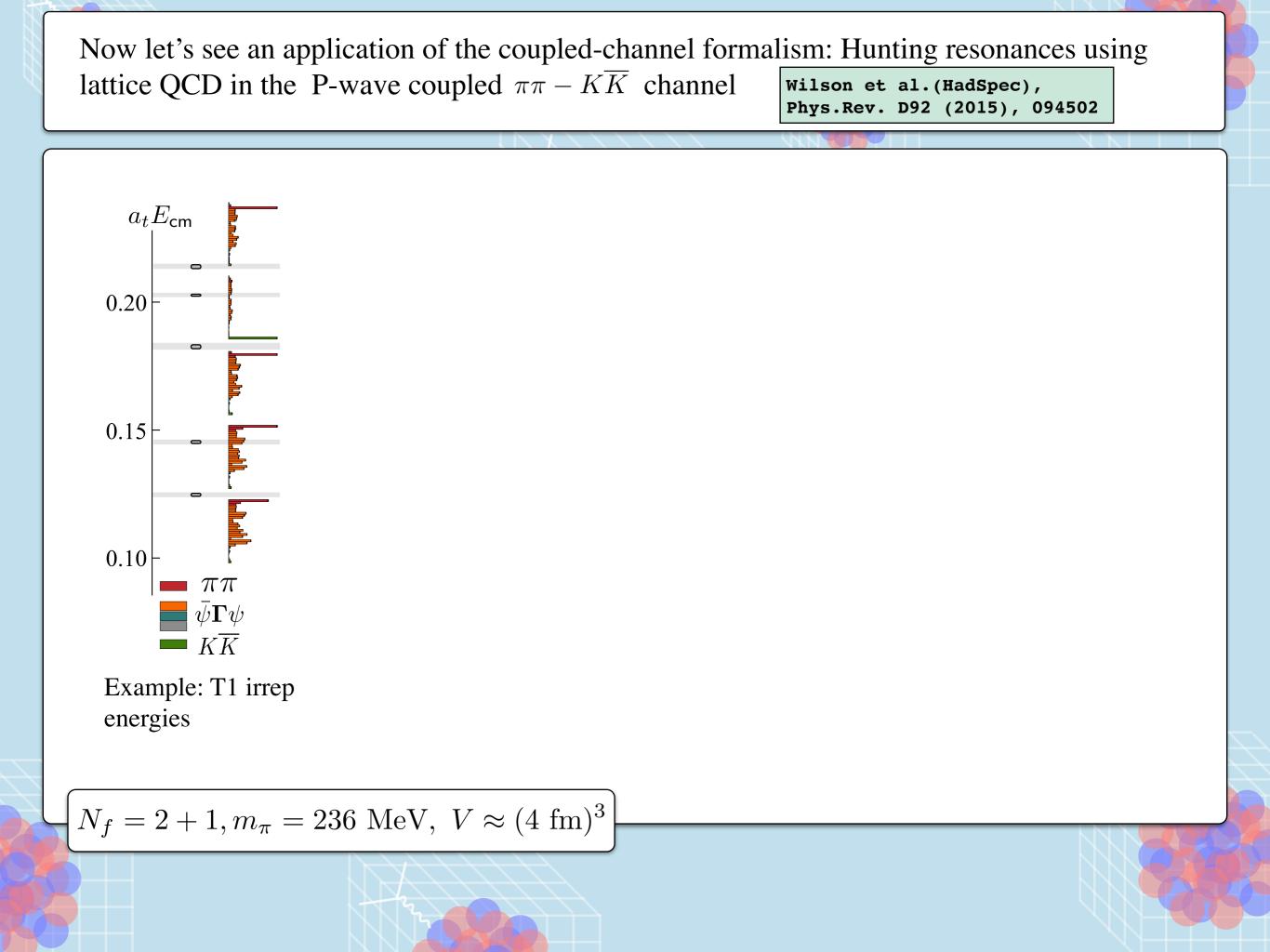
$$\cos 2\bar{\epsilon}\cos\left(\phi_1^P + \delta_1 - \phi_2^P - \delta_2\right) = \cos\left(\phi_1^P + \delta_1 + \phi_2^P + \delta_2\right)$$

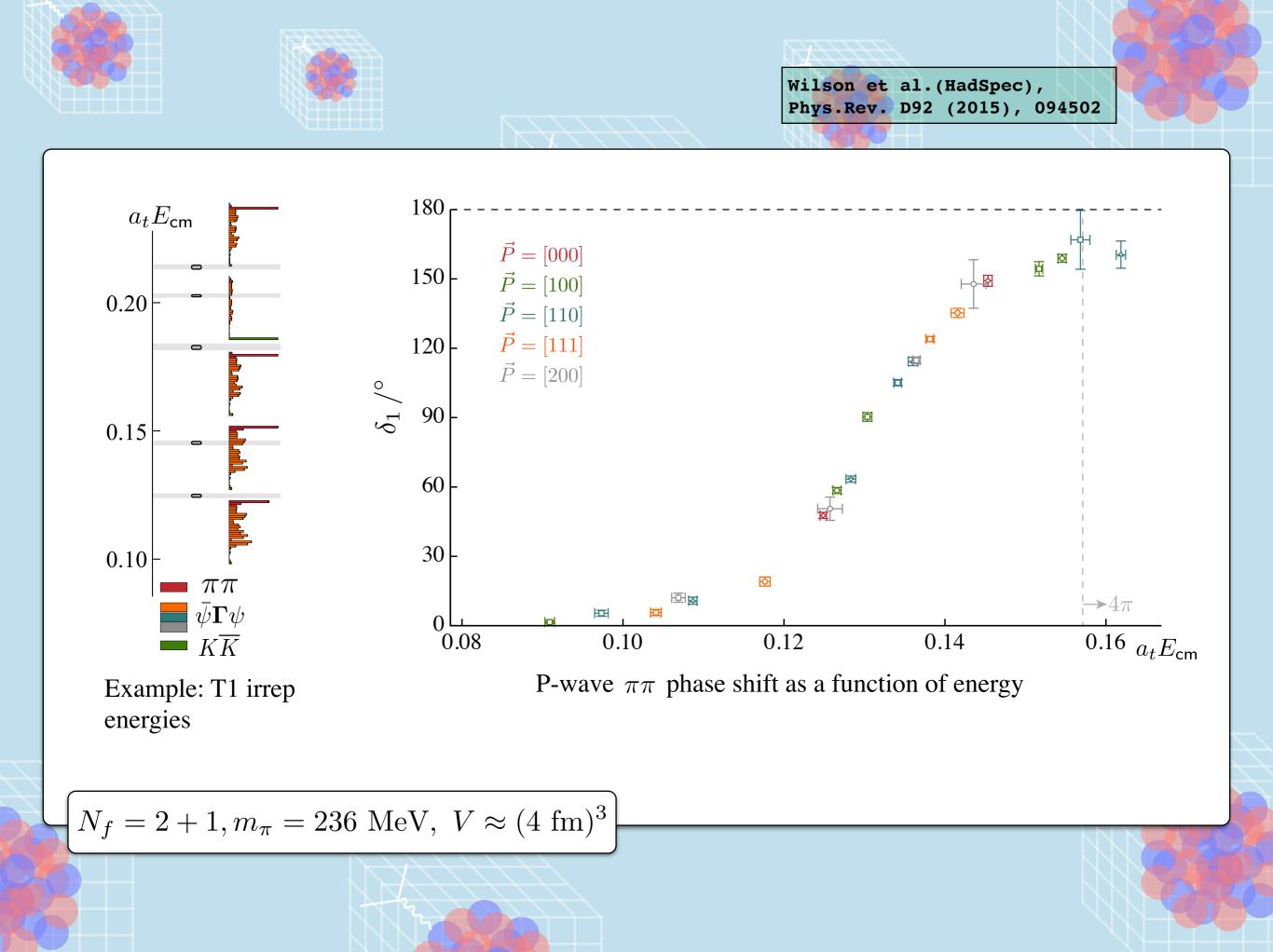
Here 1 and 2 indices refer to the two channels and superscript (0) is removed from the S-wave phase shifts for brevity. The finite-volume phase function  $\phi_i^P$  is defined as:

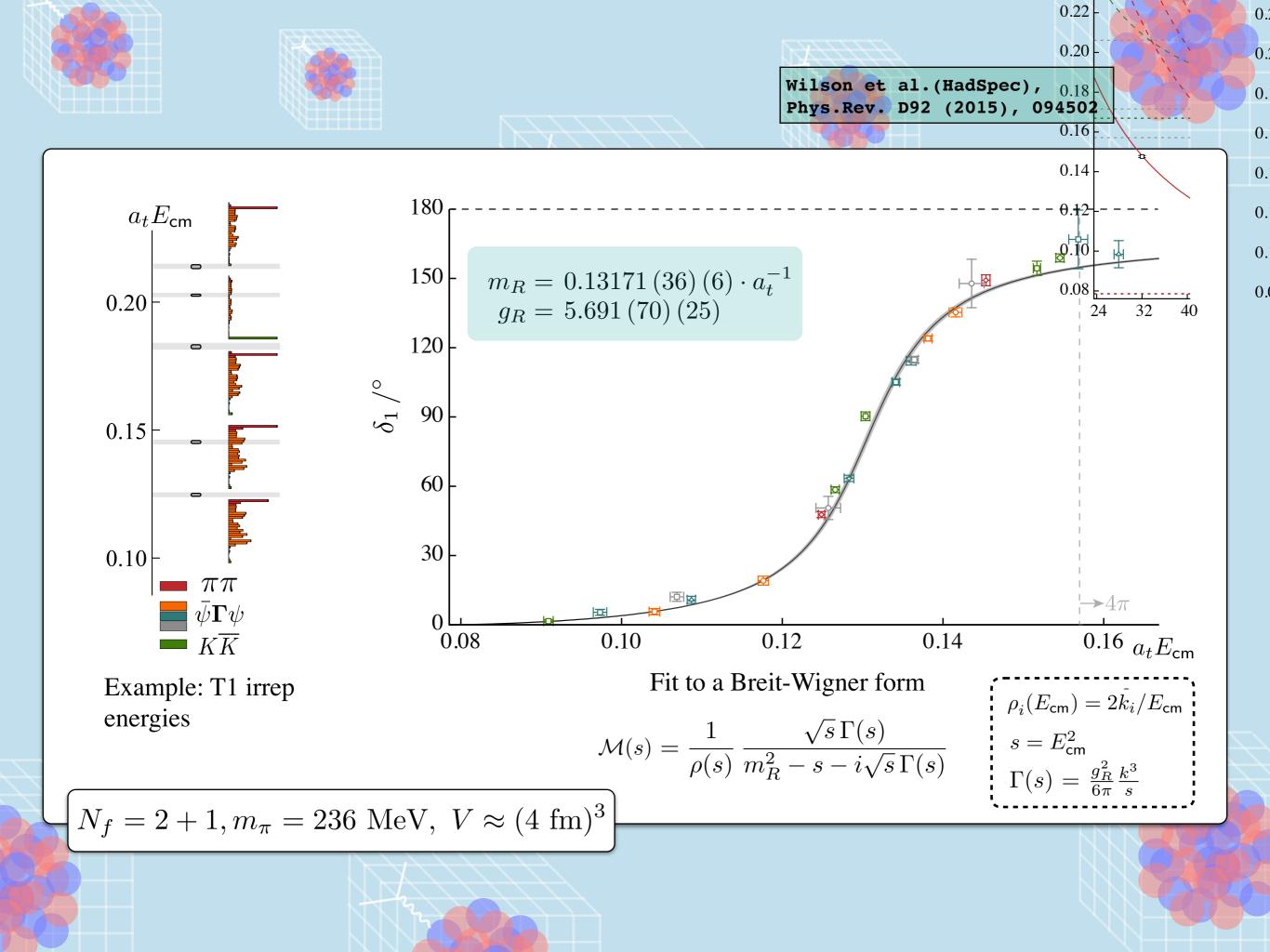
$$q_i^* \cot(\phi_i^P) \equiv -4\pi c_{00}^P(q_i^{*2})$$

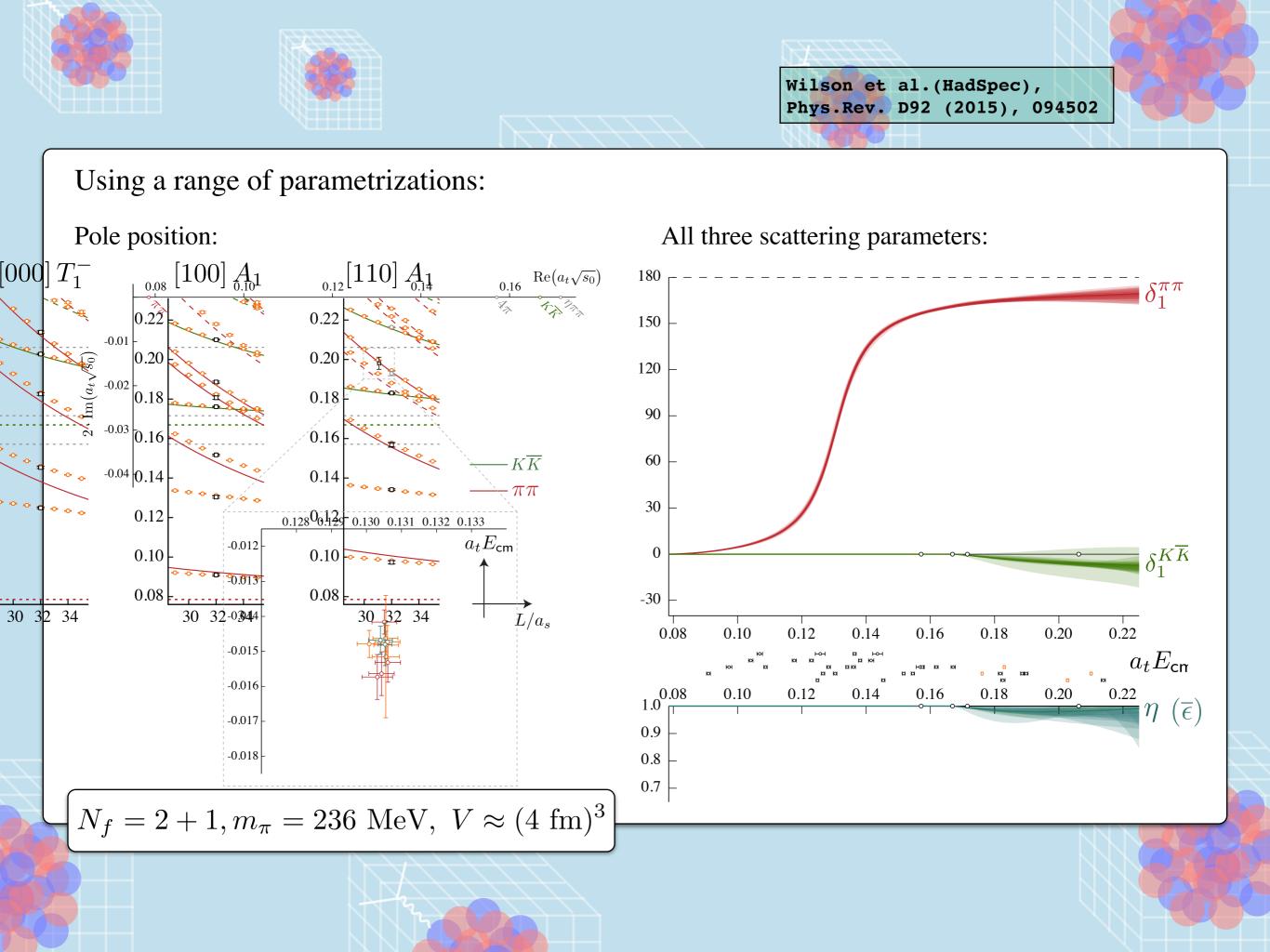
for i=1,2. This is a generic result: Luescher's "quantization condition" is a real condition.

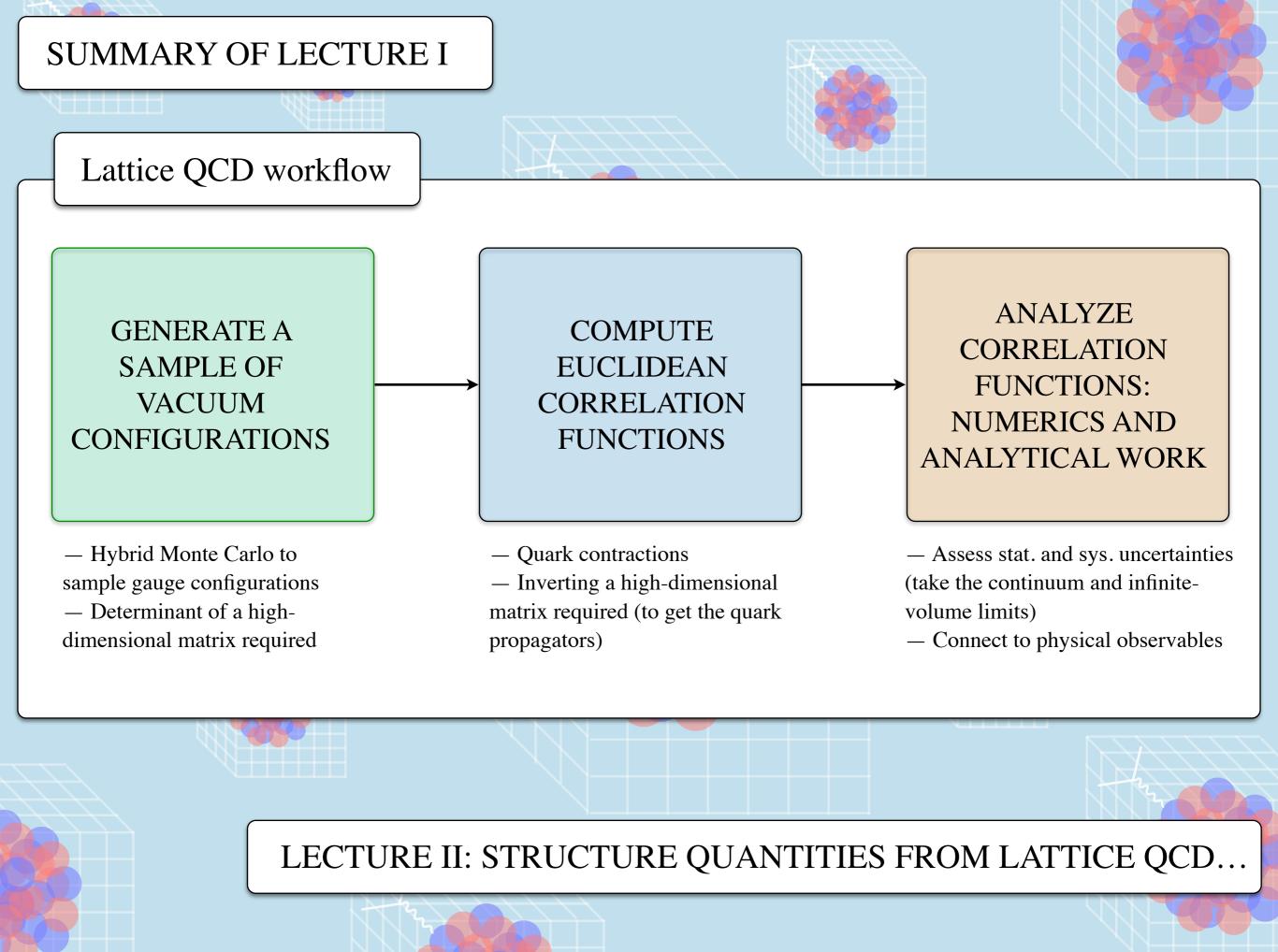


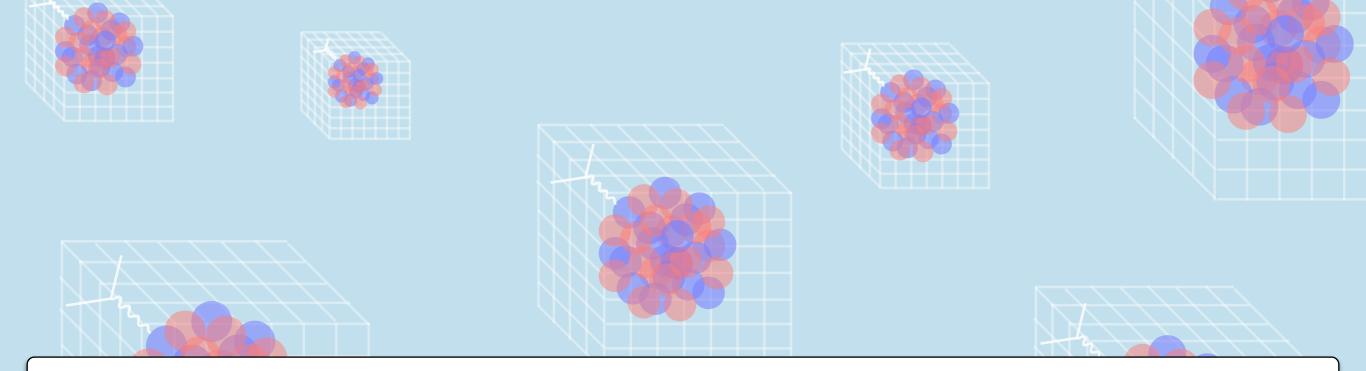




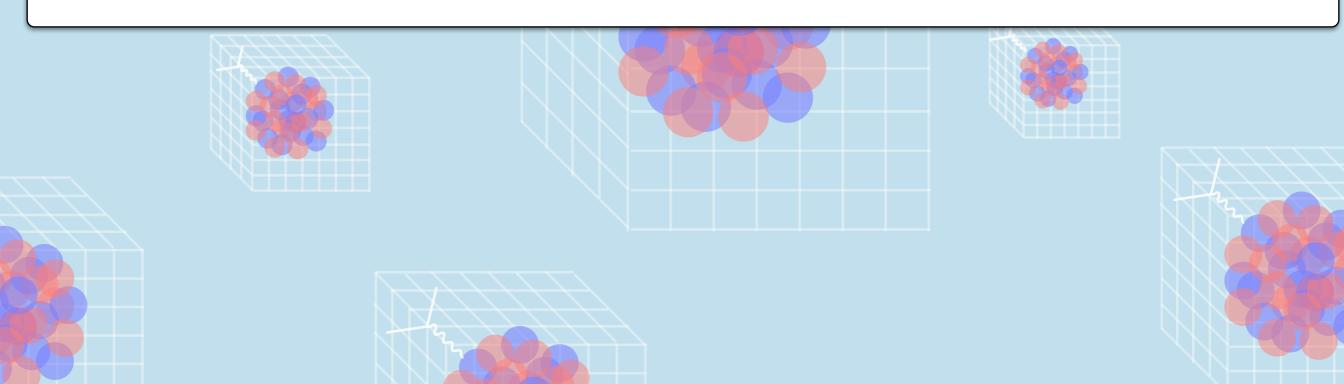








#### LECTUTE II: NUCLEON STRUCTURE FROM LATTICE QCD



Let's enumerate some of the methods that give access to structure quantities in general:

#### Three(four)-point functions

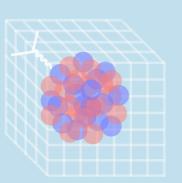
For e.g., form factors, moments of structure functions, Compton amplitude, transition amplitudes

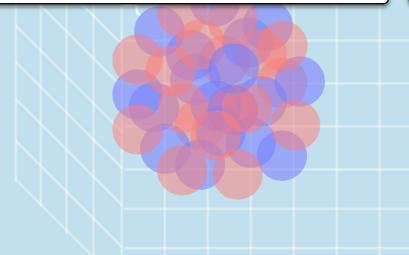
## Background-field methods

For e.g., EM moments and polarizabilities, charge radius, form factors and transition amplitudes.

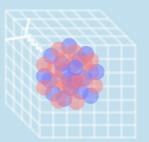
# Feynman-Hellmann inspired methods

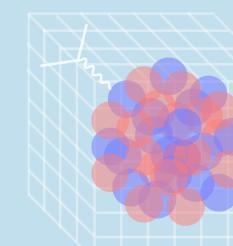
Similar to background fields. For e.g., axial charge, form factors, EM moments, transition amplitudes











Let's enumerate some of the methods that give access to structure quantities in general:

#### Three(four)-point functions

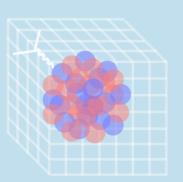
For e.g., form factors, moments of structure functions, Compton amplitude, transition amplitudes

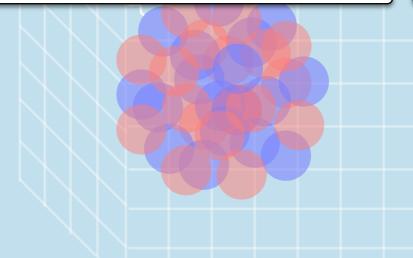
# Background-field methods

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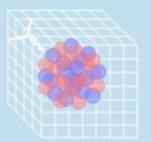
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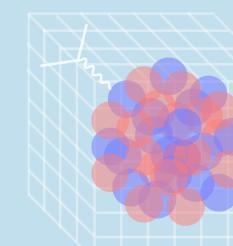
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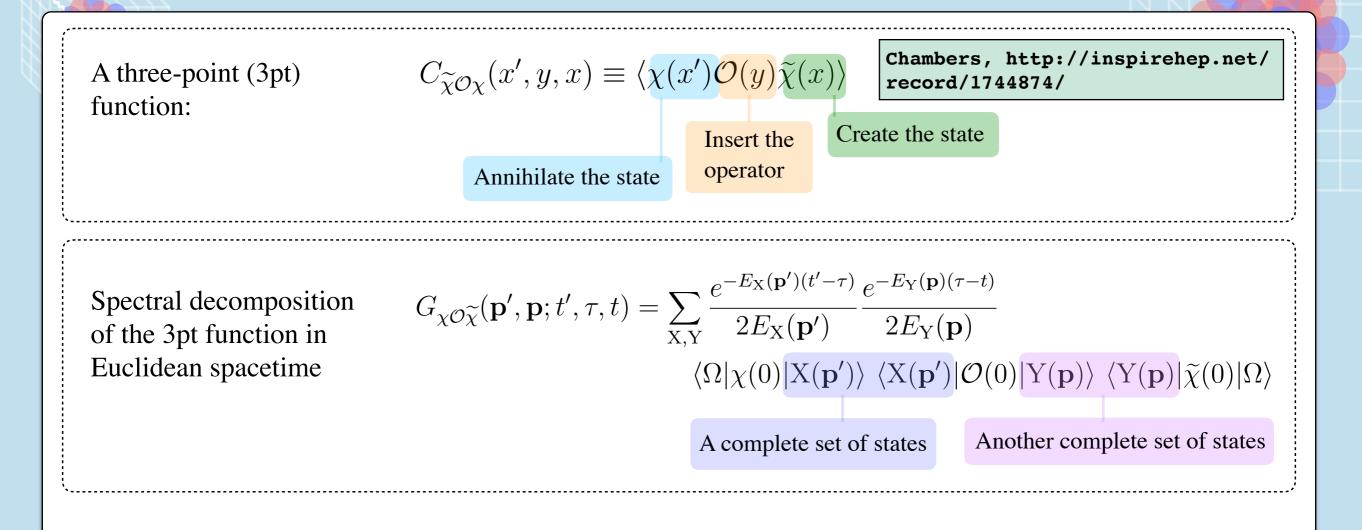


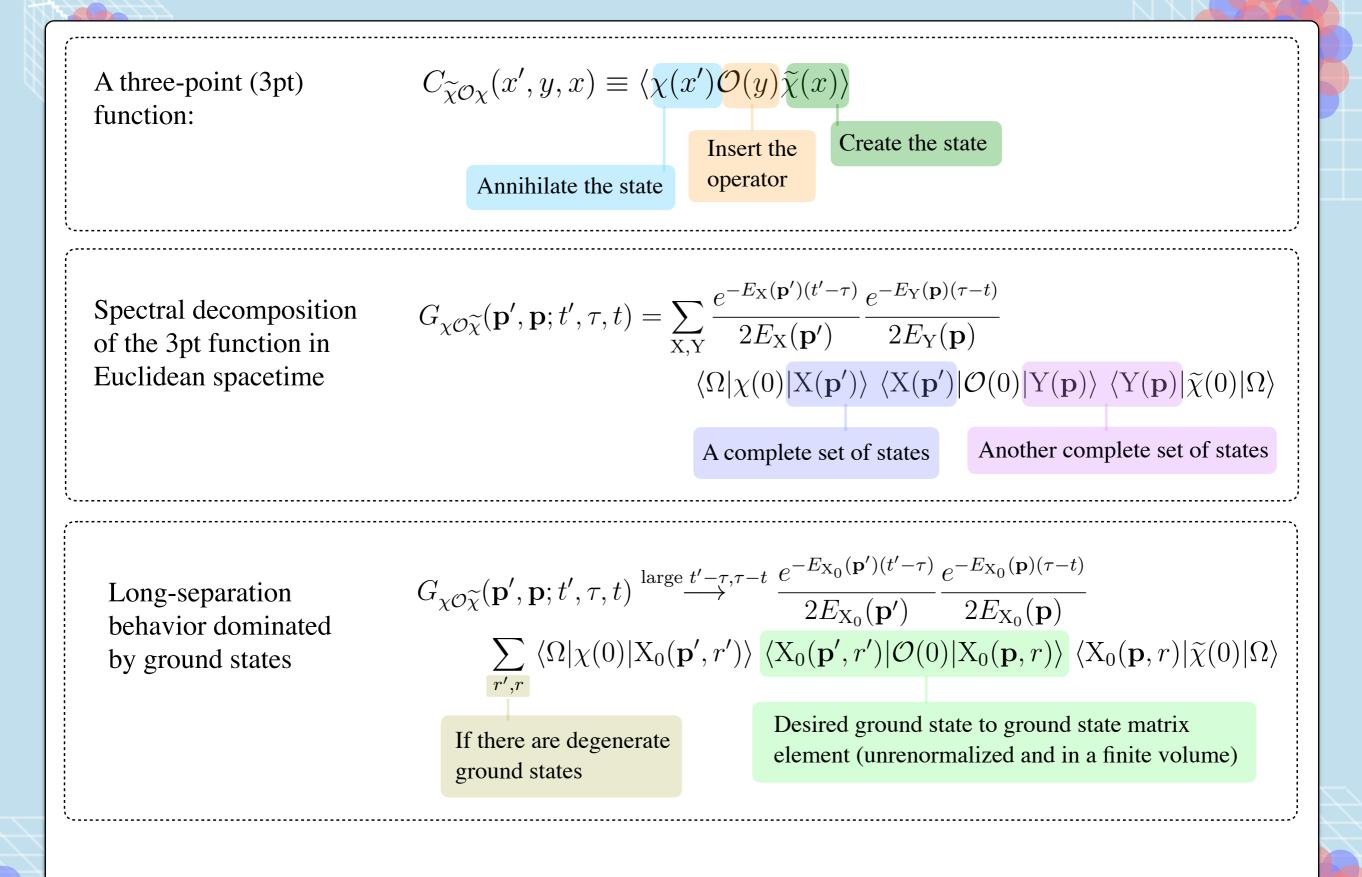


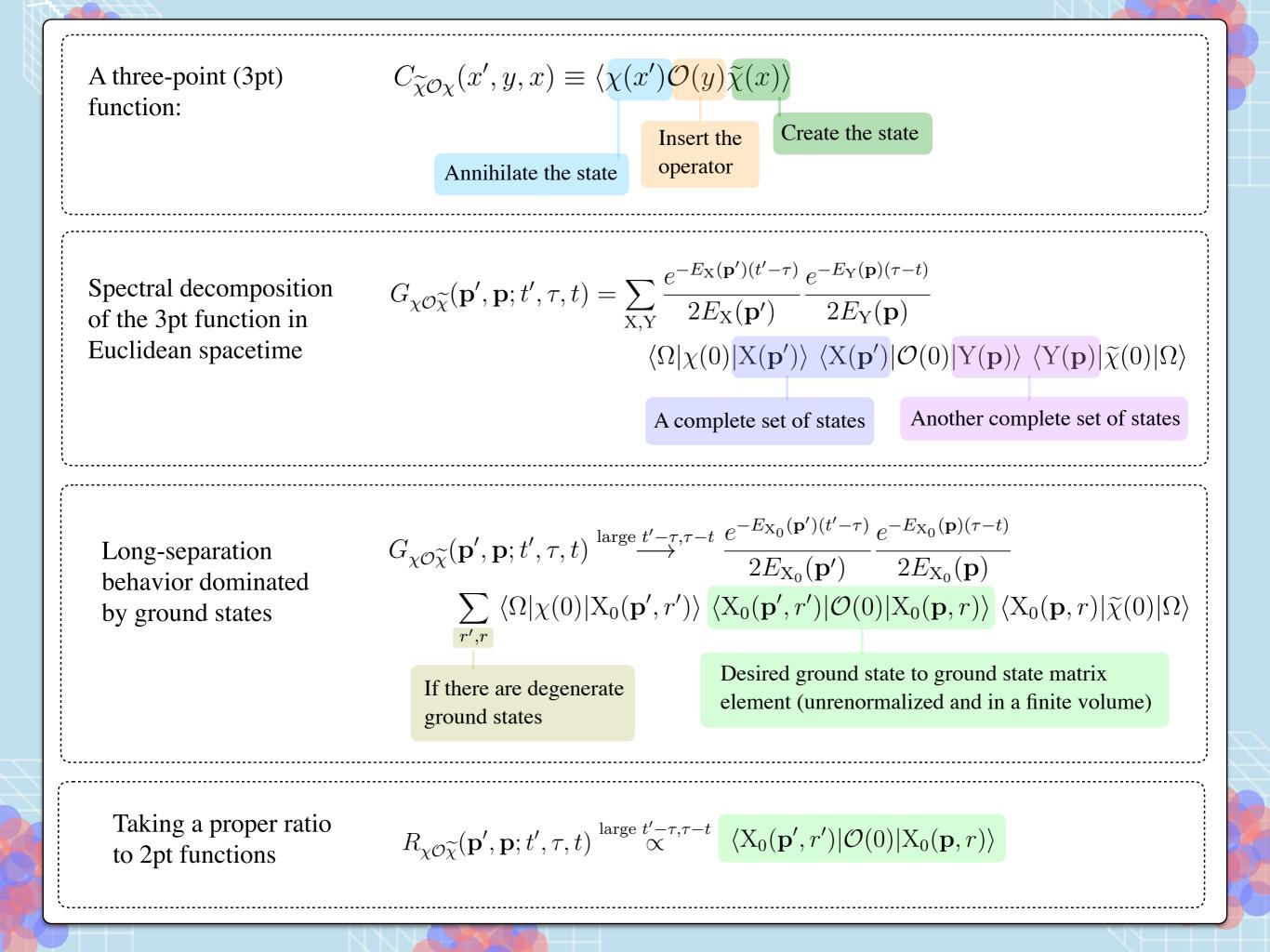




A three-point (3pt) $C_{\hat{\lambda}}$ function:	$\chi_{\mathcal{O}\chi}(x',y,x) \equiv \langle \chi(x) \rangle$		Chambers, http record/174487 Create the state	p://inspirehep.net/ 4/
	Annihilate the state	operator		
<				







#### **EXERCISE 6**



If the computational resources do not allow large source, operator and sink time separations to be achieved, one should worry about the effect of excited states. One way to have more confidence over the extracted ground state to ground state matrix element is to perform a multi-exponential fits to the ratio of 3pt to 2pt functions as a function of both the source-sink and the source-operator separations. Assume that both the ground state and the first excited states contribute significantly to such a ratio. Write down a generic form for such a multi-exponential function.



In the above exercise, sum over the time insertions of the operator and write down a new form for the ratio of 3pt to 2pt functions, which now is only a function of the source-sink time separation. This is referred to as the summation method in literature.

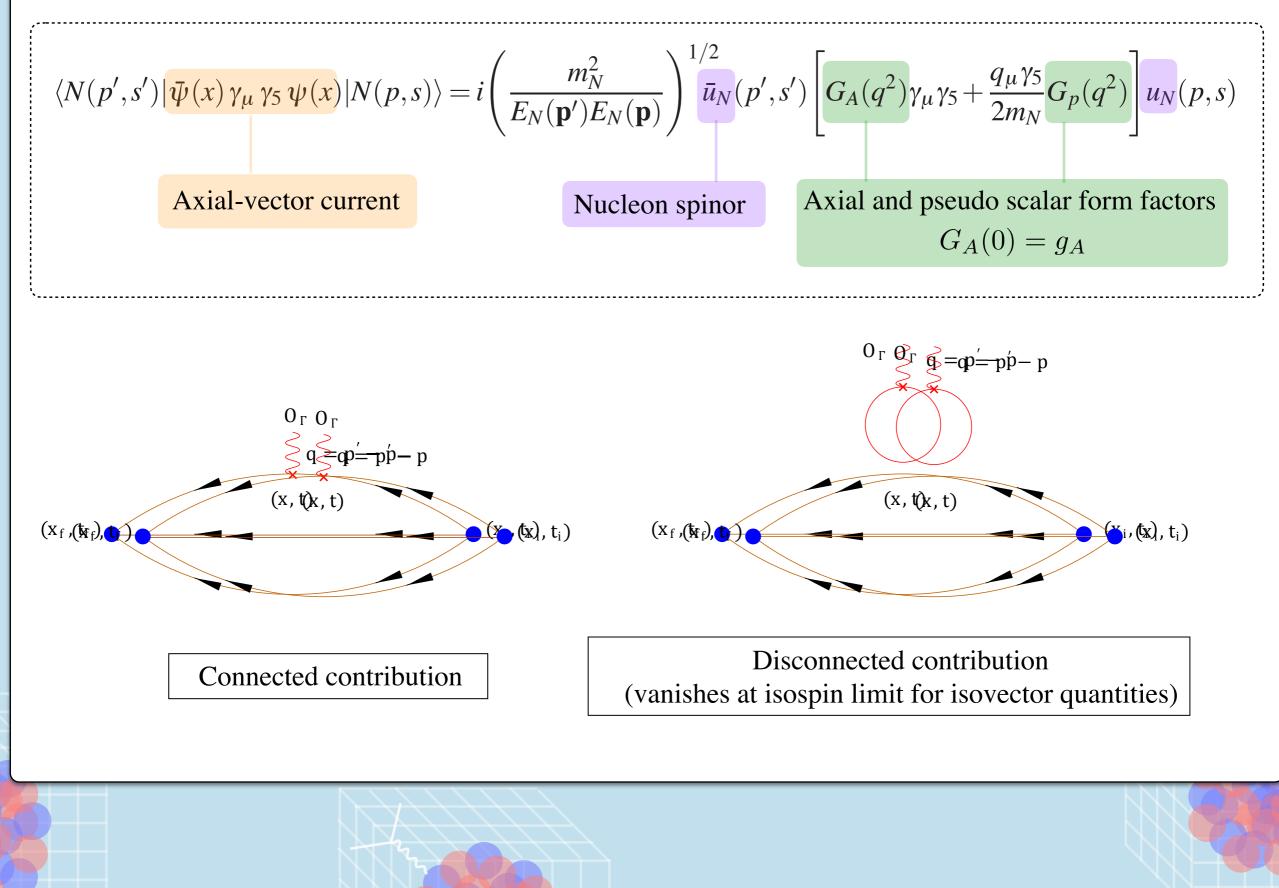
Example: The application of 3pt function method to obtain the axial charge/form factors of the nucleon

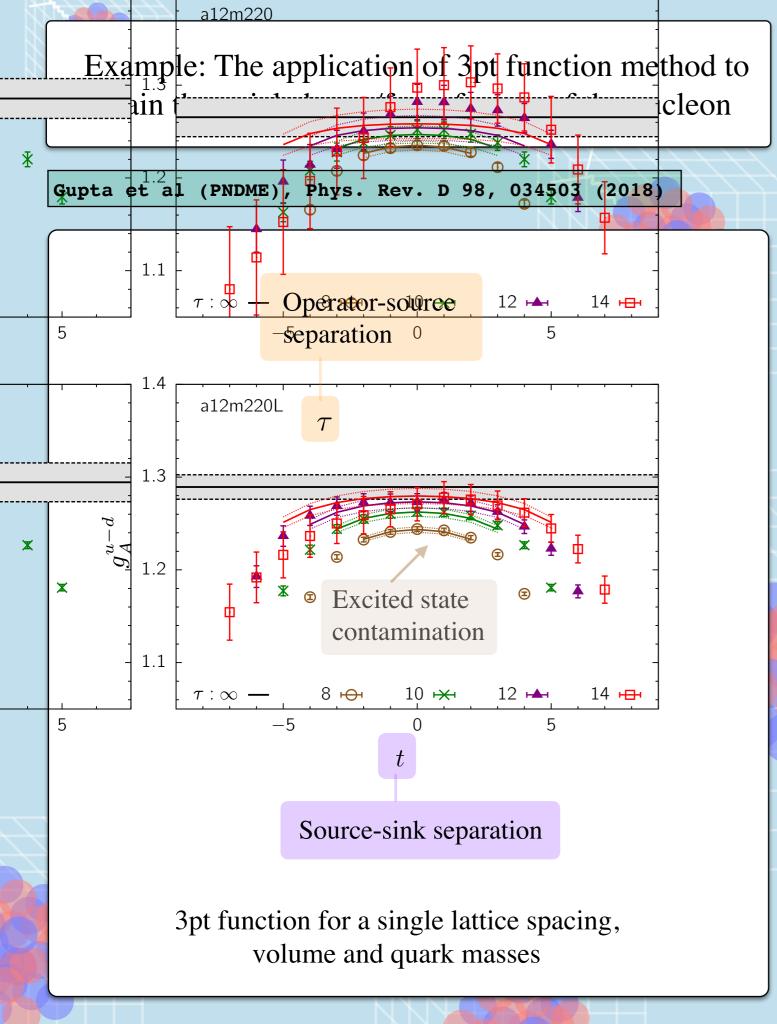
Constantinou, arXiv:1411.0078 [hep-lat].

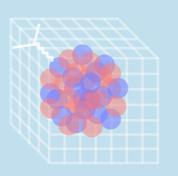
$$\langle N(p',s') | \bar{\psi}(x) \gamma_{\mu} \gamma_{5} \psi(x) | N(p,s) \rangle = i \left( \frac{m_{N}^{2}}{E_{N}(\mathbf{p}')E_{N}(\mathbf{p})} \right)^{1/2} \bar{u}_{N}(p',s') \left[ G_{A}(q^{2}) \gamma_{\mu} \gamma_{5} + \frac{q_{\mu} \gamma_{5}}{2m_{N}} G_{p}(q^{2}) \right] u_{N}(p,s)$$
Axial-vector current
Nucleon spinor
Axial and pseudo scalar form factors
$$G_{A}(0) = g_{A}$$

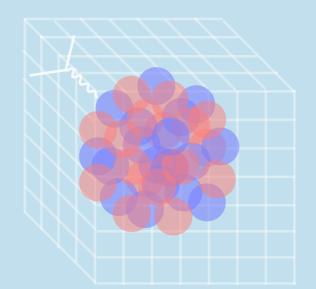
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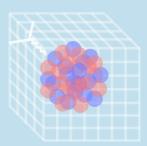
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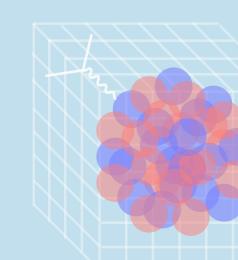


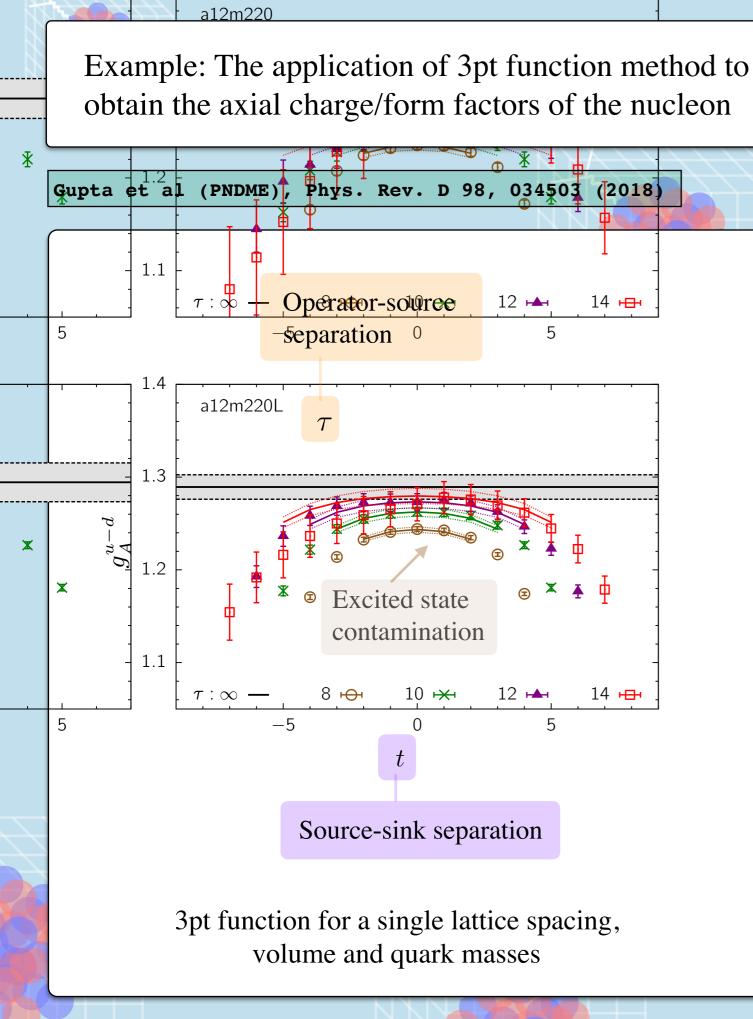


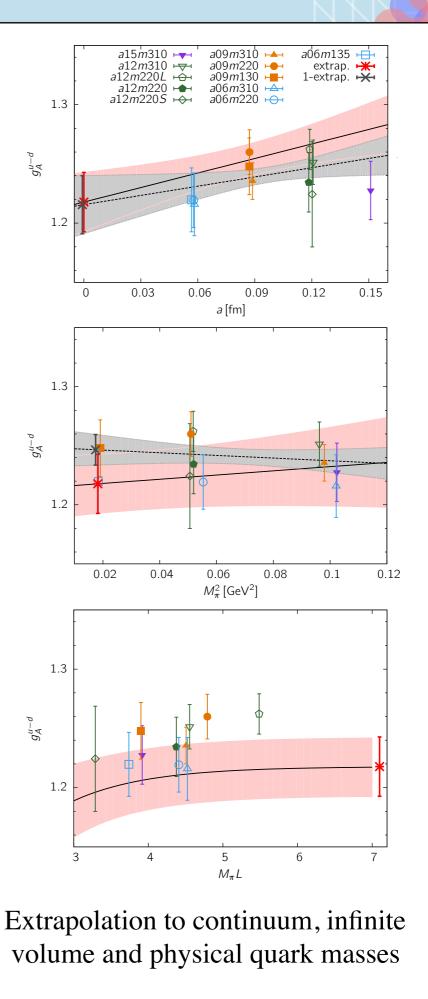




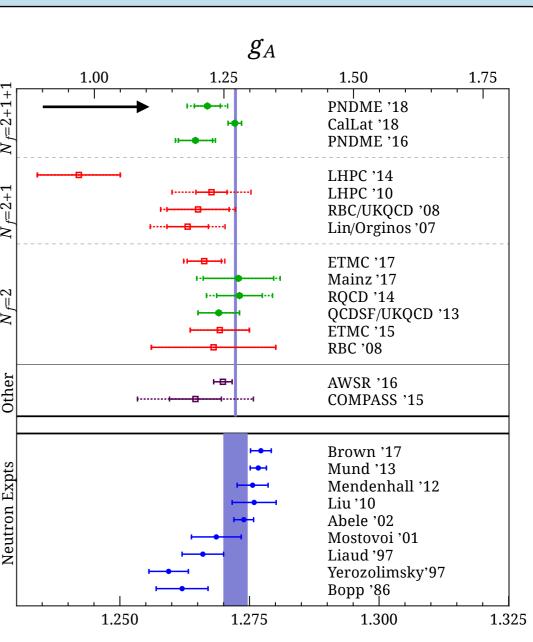




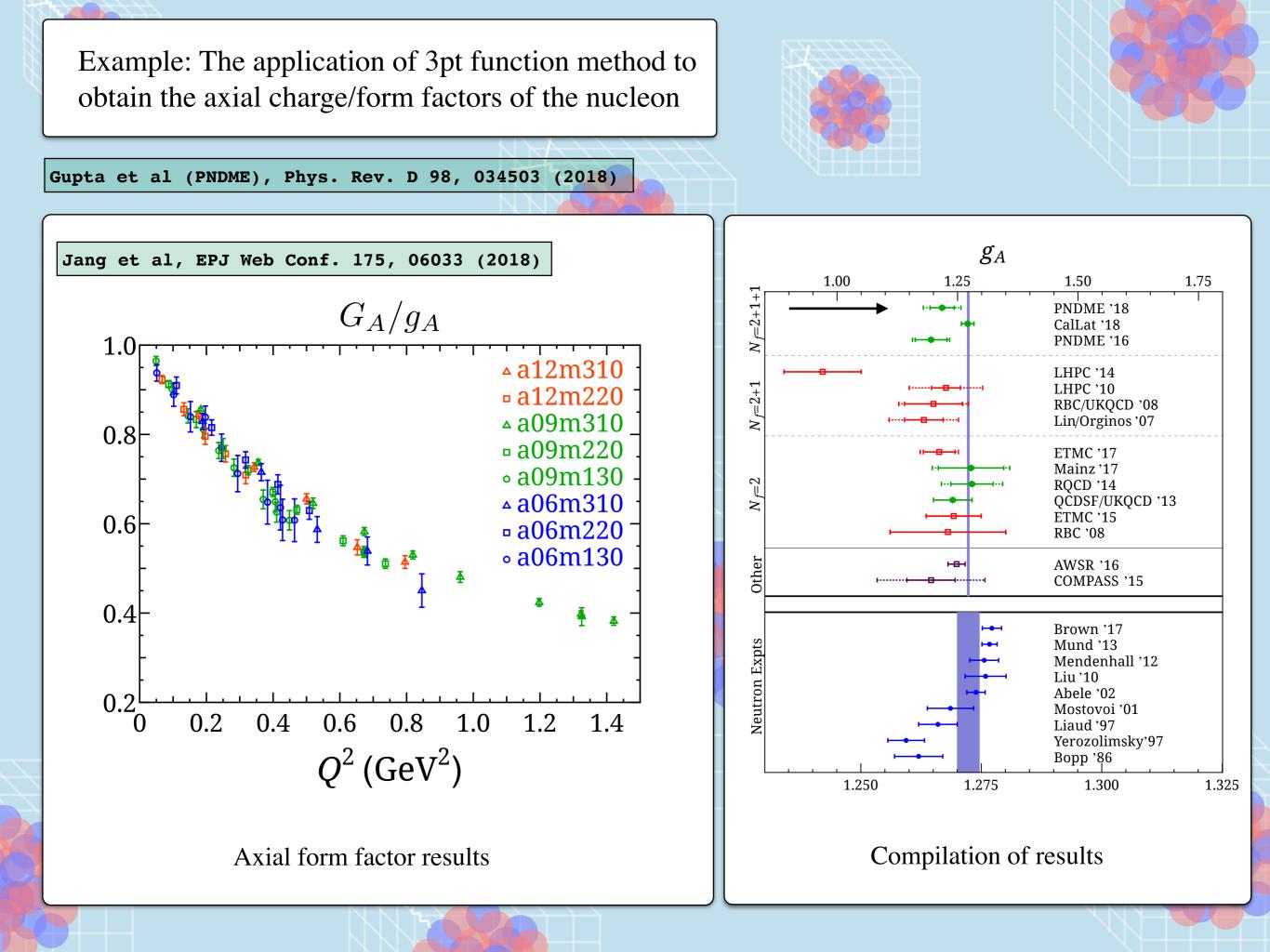




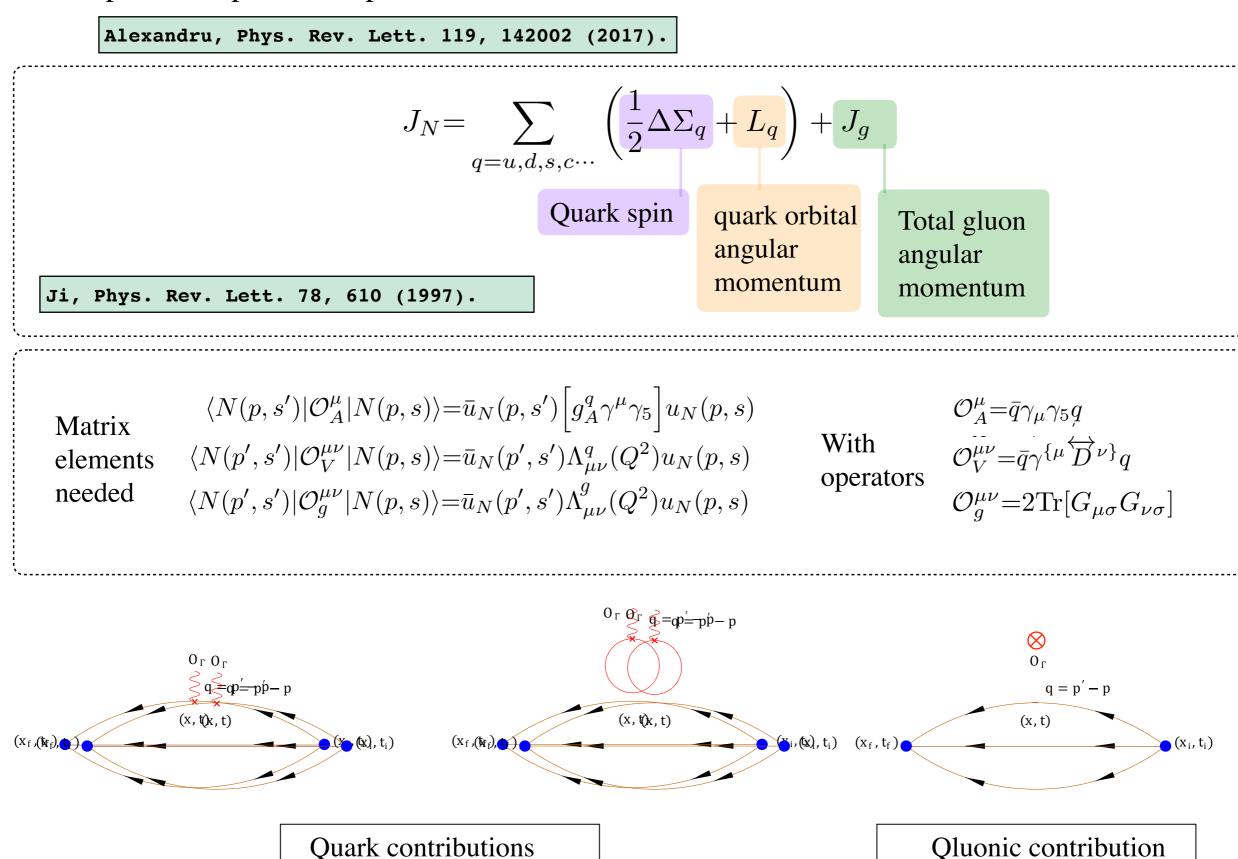
Example: The application of 3pt function method to obtain the axial charge/form factors of the nucleon Gupta et al (PNDME), Phys. Rev. D 98, 034503 (2018) 1.00  $N_{f}=2+1+1$ f = 2 + 1Z  $N_{f=2}$ Other Neutron Expts 1.250



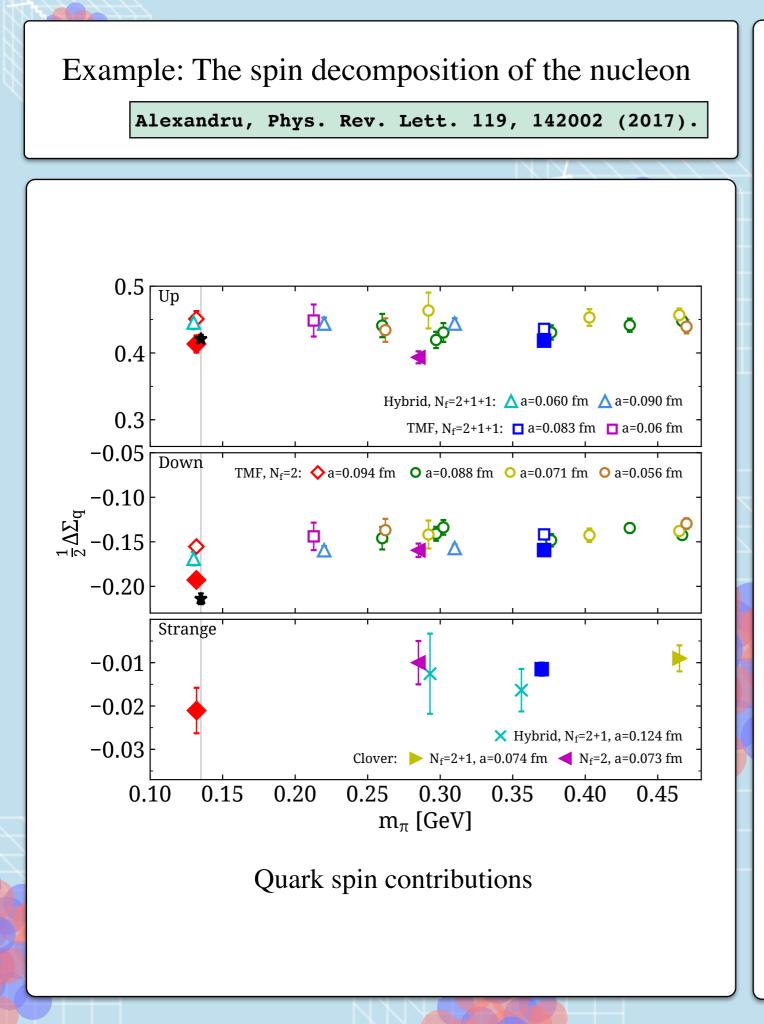
Compilation of results

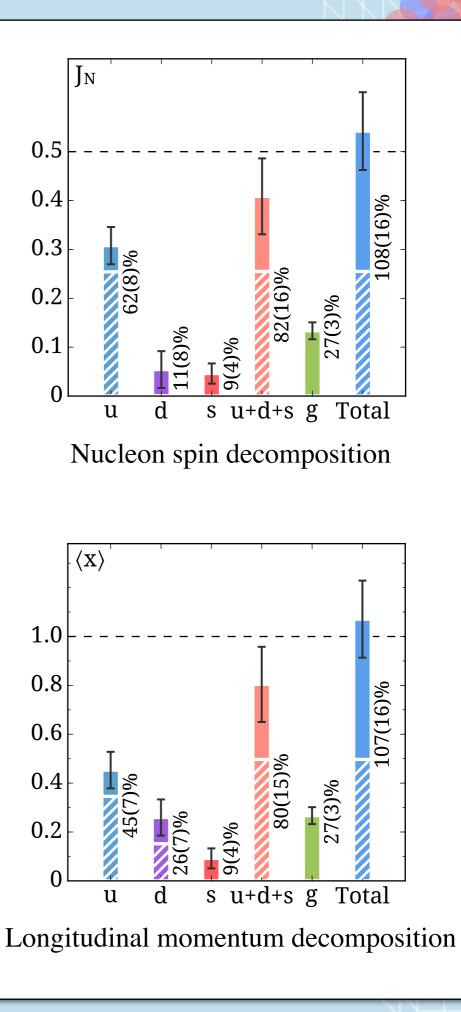


#### Example: The spin decomposition of the nucleon



 $(X_f)$ 





Let's enumerate a some of the methods that give access to structure quantities in general:

#### Three(four)-point functions

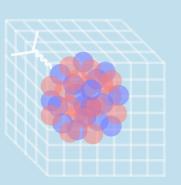
For e.g., form factors, moments of structure functions, Compton amplitude, transition amplitudes

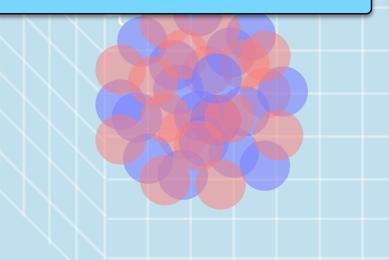
# Background-field methods

For e.g., EM moments and polarizabilities, charge radius, form factors and transition amplitudes.

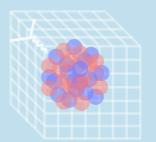
# Feynman-Hellmann inspired methods

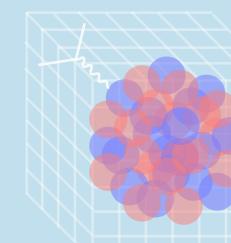
Similar to background fields. For e.g., axial charge, form factors, EM moments, transition amplitudes







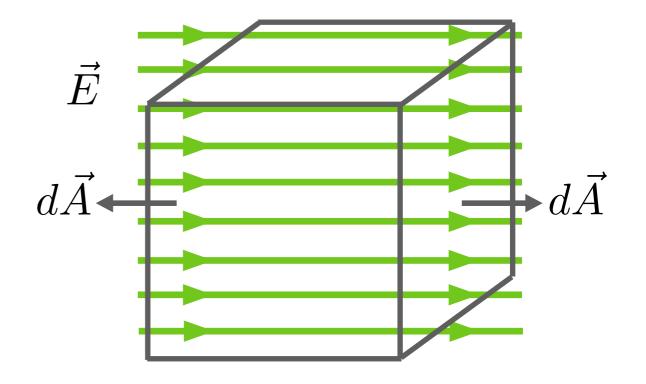


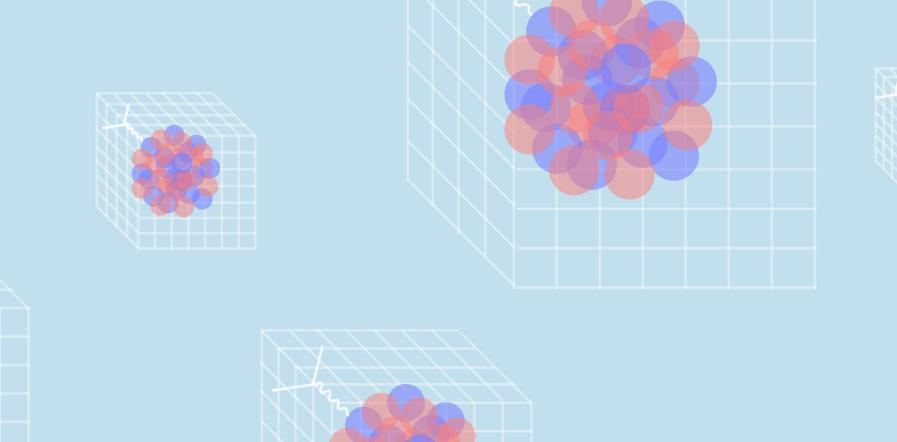


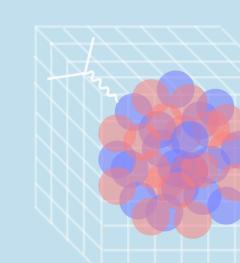
Background fields are non-dynamical, i.e., there will be no pair creation and annihilation in vacuum with a classical EM background field. This mean the photon zero mode is no problem: it is absent in the calculation!

$$U^{(\text{QCD})} \to U^{(\text{QCD})} \times U^{(\text{QED})}$$

Modify the links when forming the quark propagators.







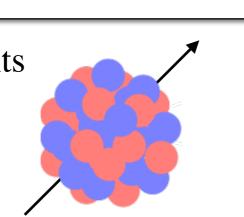
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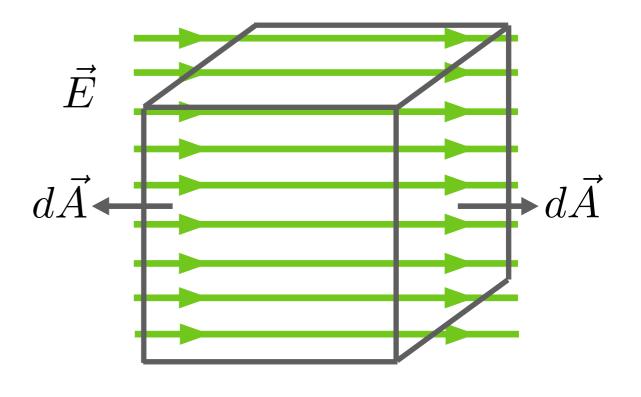
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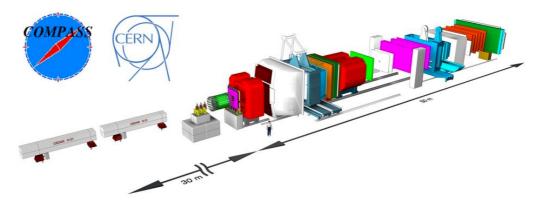
Traditionally they are used for constraining the response of hadrons/nuclei to external probes:

Magnetic moments



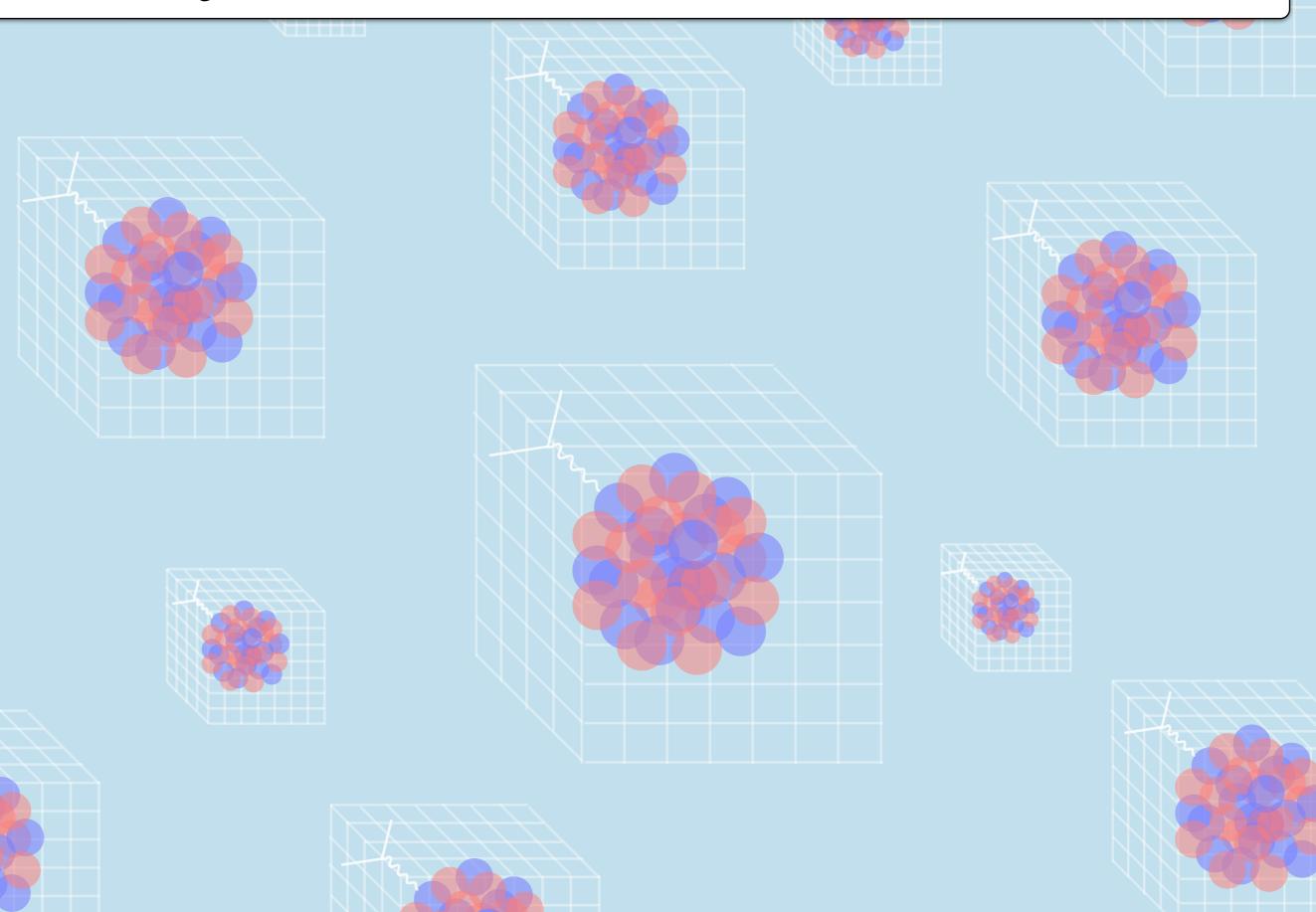


Electric and magnetic polarizabilities



See e.g., BEANE et al (NPLQCD), Phys.Rev.Lett. 113 (2014) 25, 252001 and Phys.Rev. D92 (2015) 11, 114502. for nuclear-physics calculations.

What does the requirement of periodicity impose on background fields? Let's consider a uniform background field.



What does the requirement of periodicity impose on background fields? Let's consider a uniform background field.

$$A_{\mu} = \left(-E \times (\mathbf{x}_{3} - \begin{bmatrix} \mathbf{x}_{3} \\ L \end{bmatrix} L), \mathbf{0} \right) \rightarrow \mathbf{E} = E\mathbf{x}_{3}$$

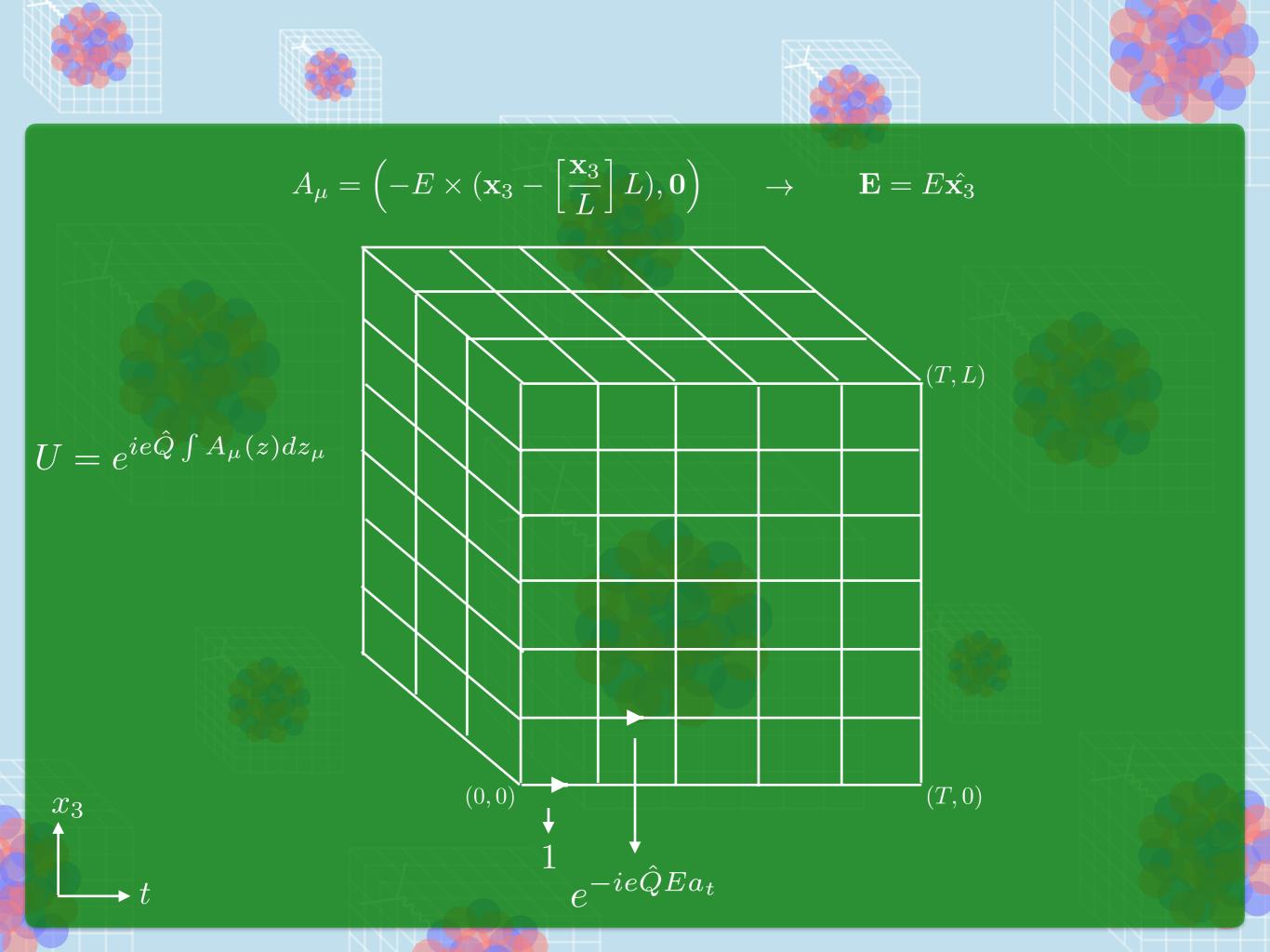
$$U = e^{ieQ \int A_{\mu}(z)dz_{\mu}}$$

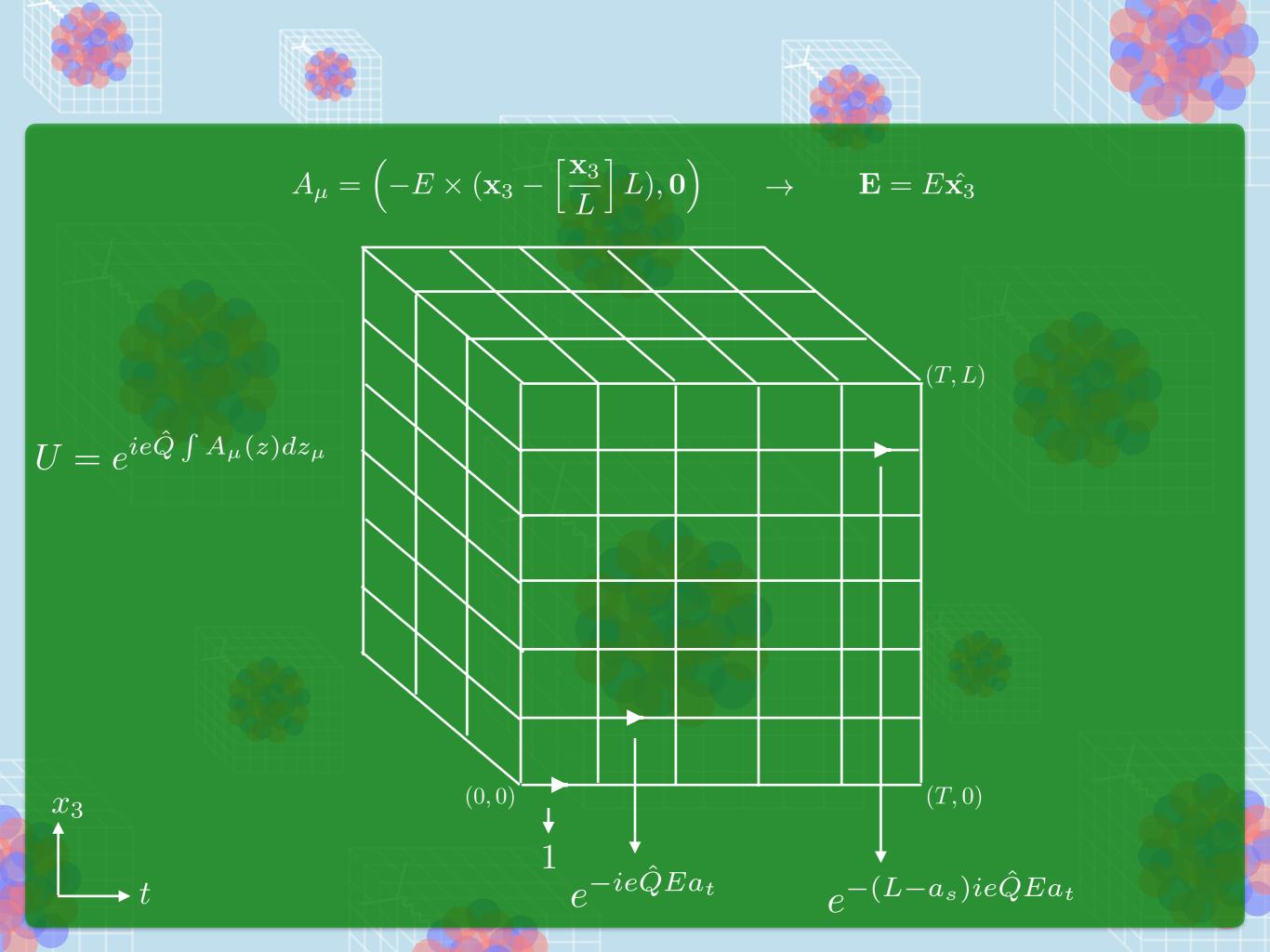
$$(0,0)$$

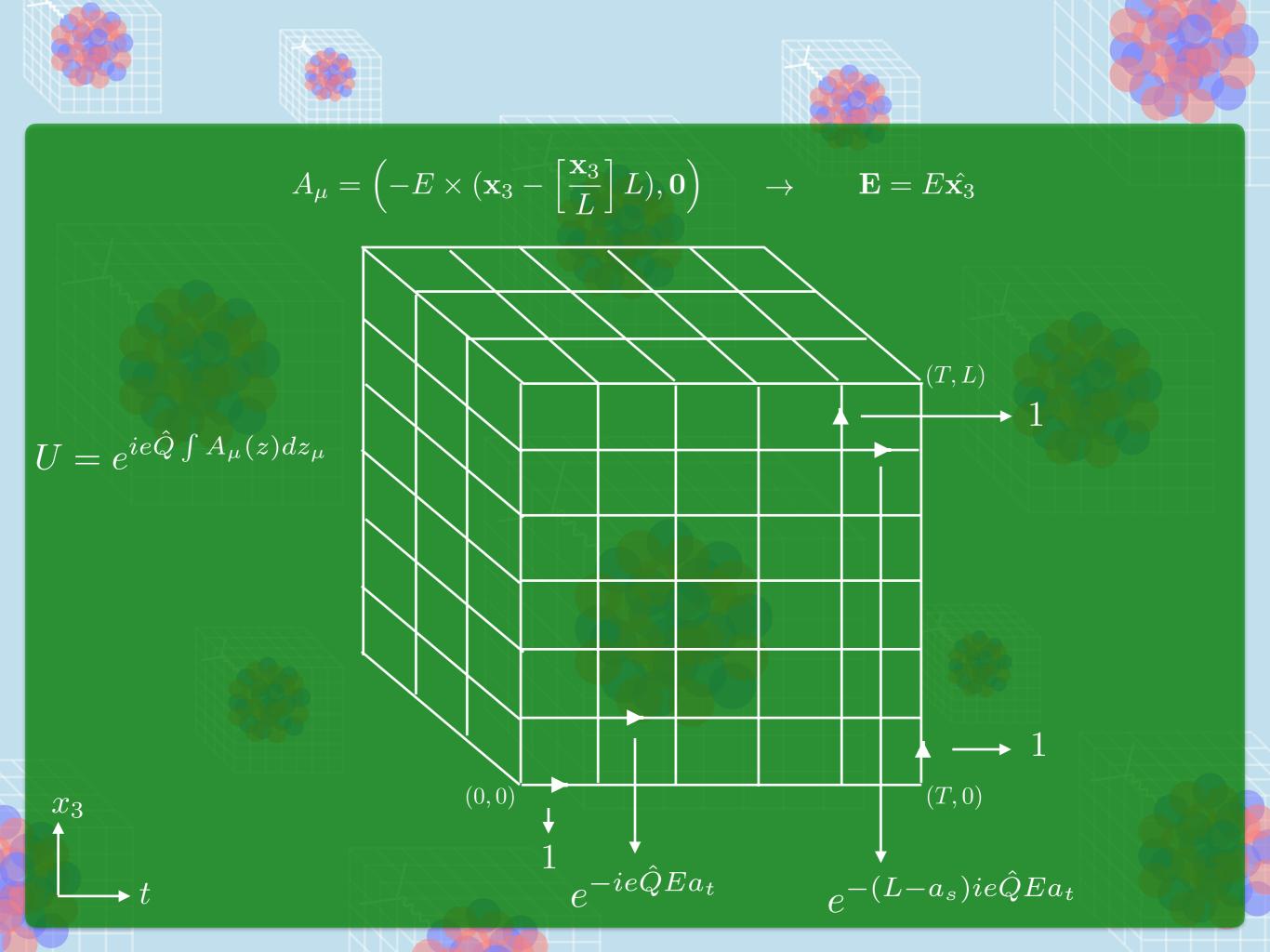
$$(0,0)$$

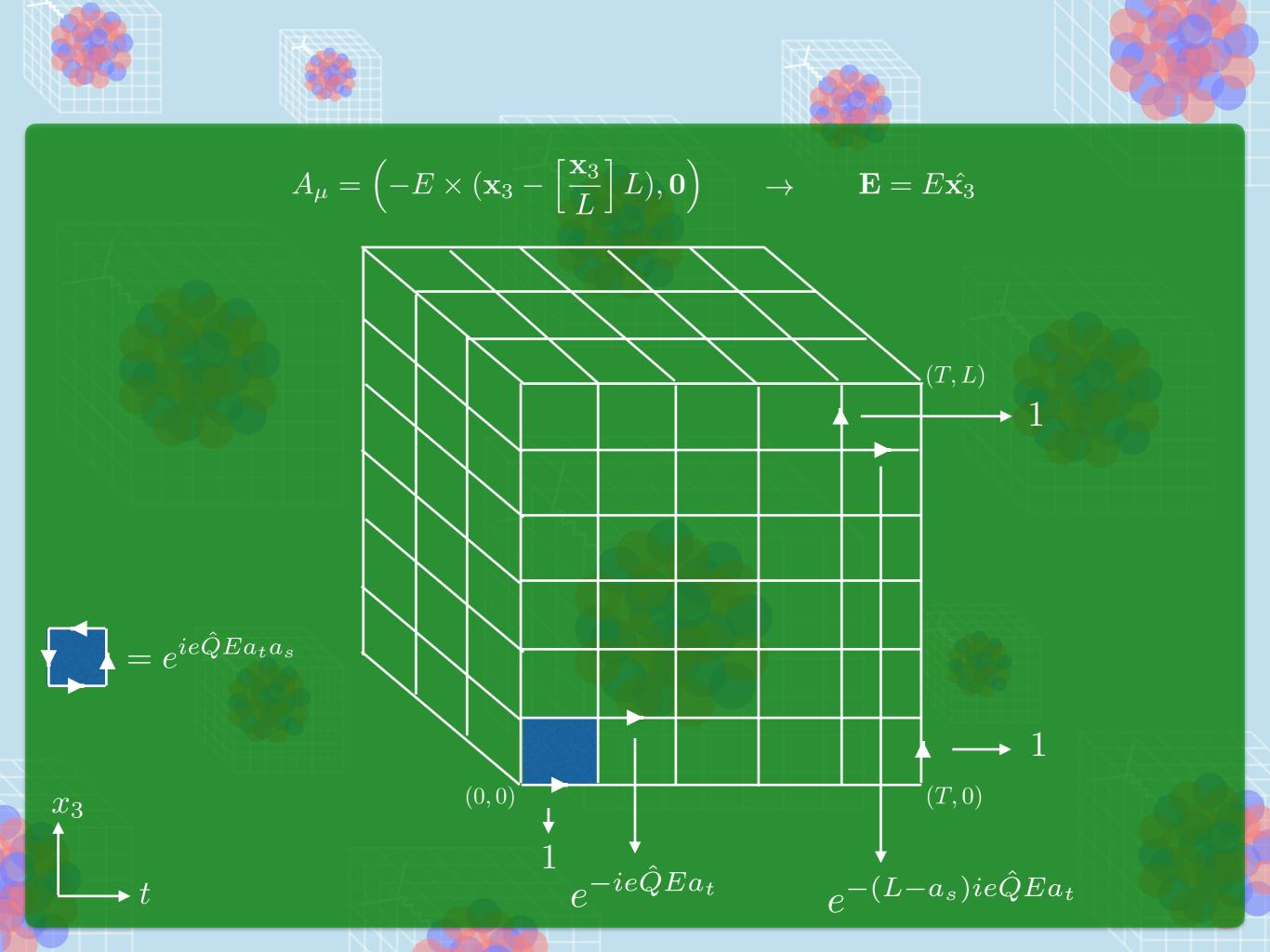
$$(T,0)$$

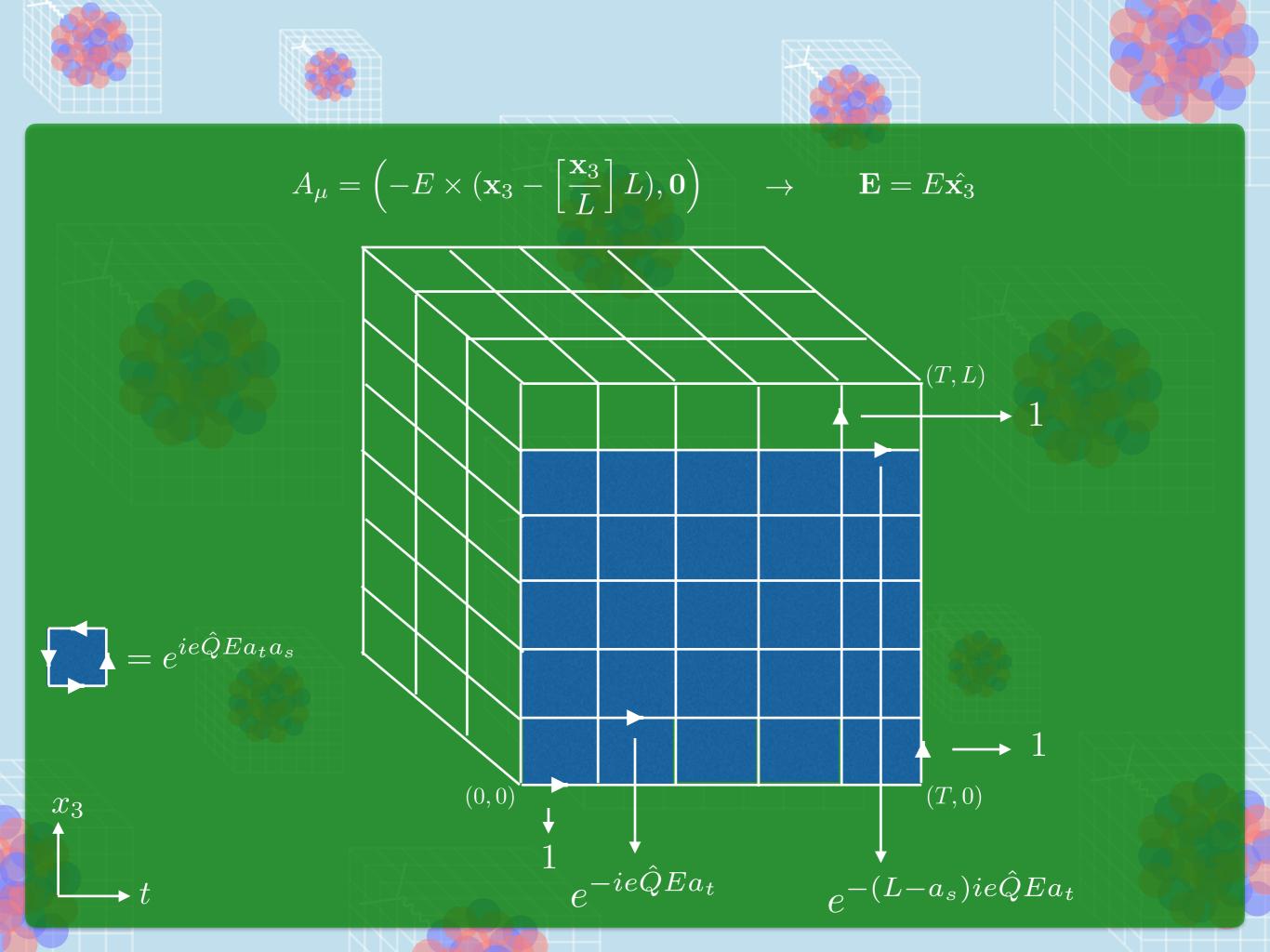
$$(T,0)$$



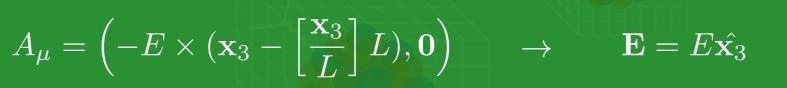




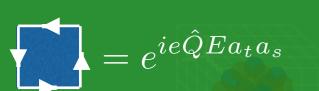


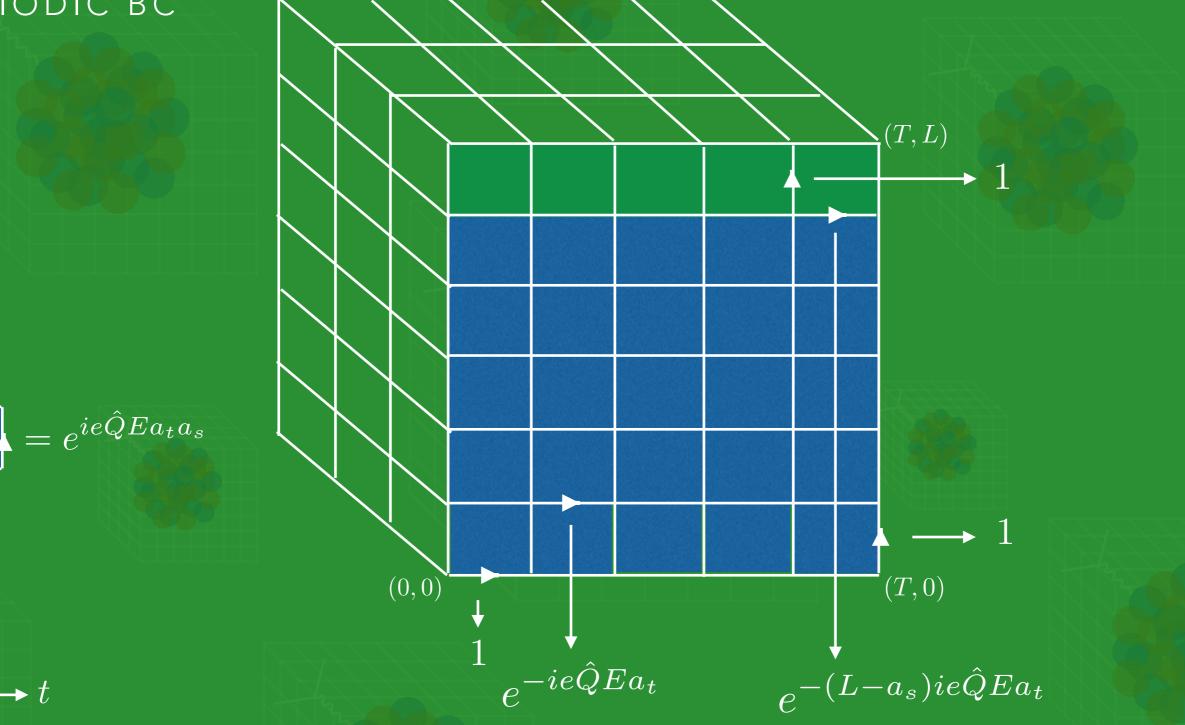




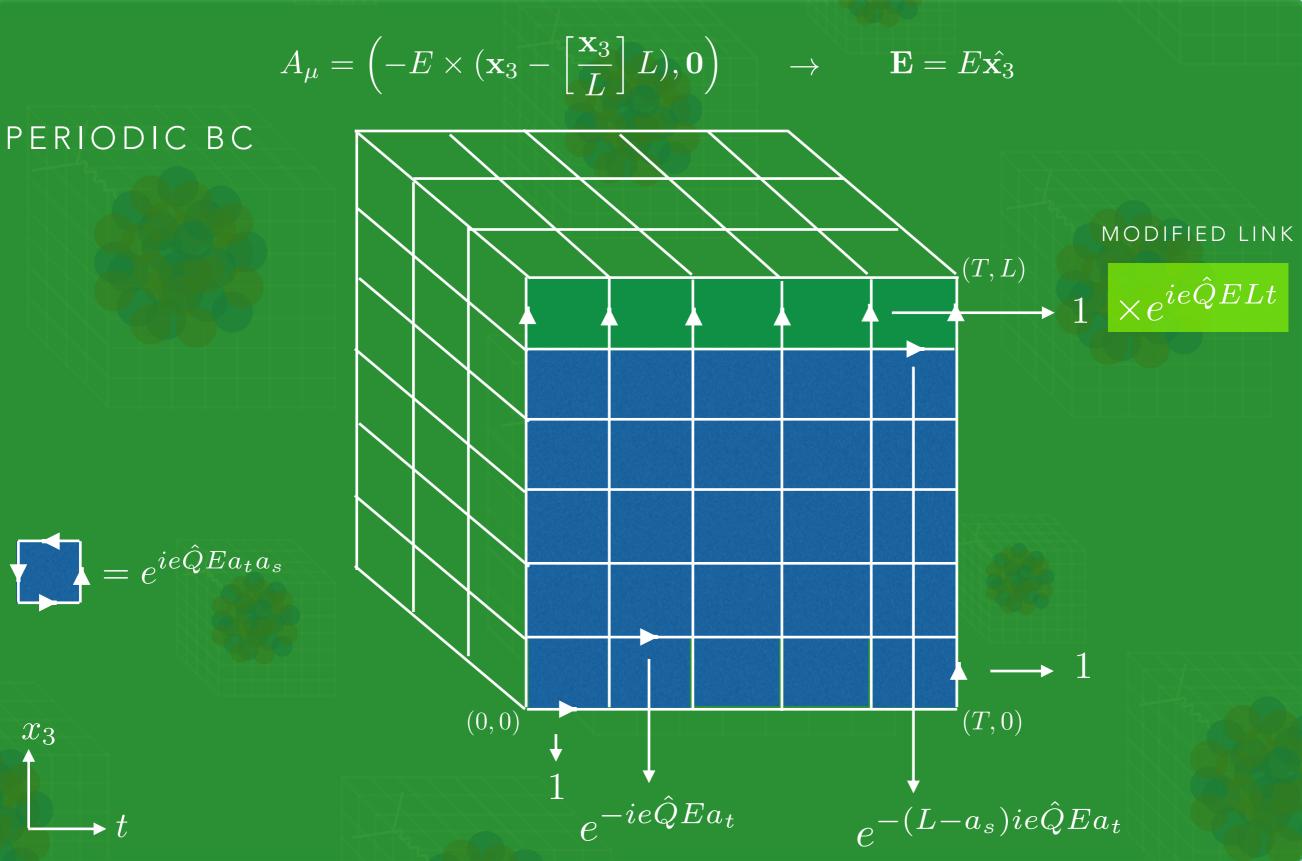


### PERIODIC BC

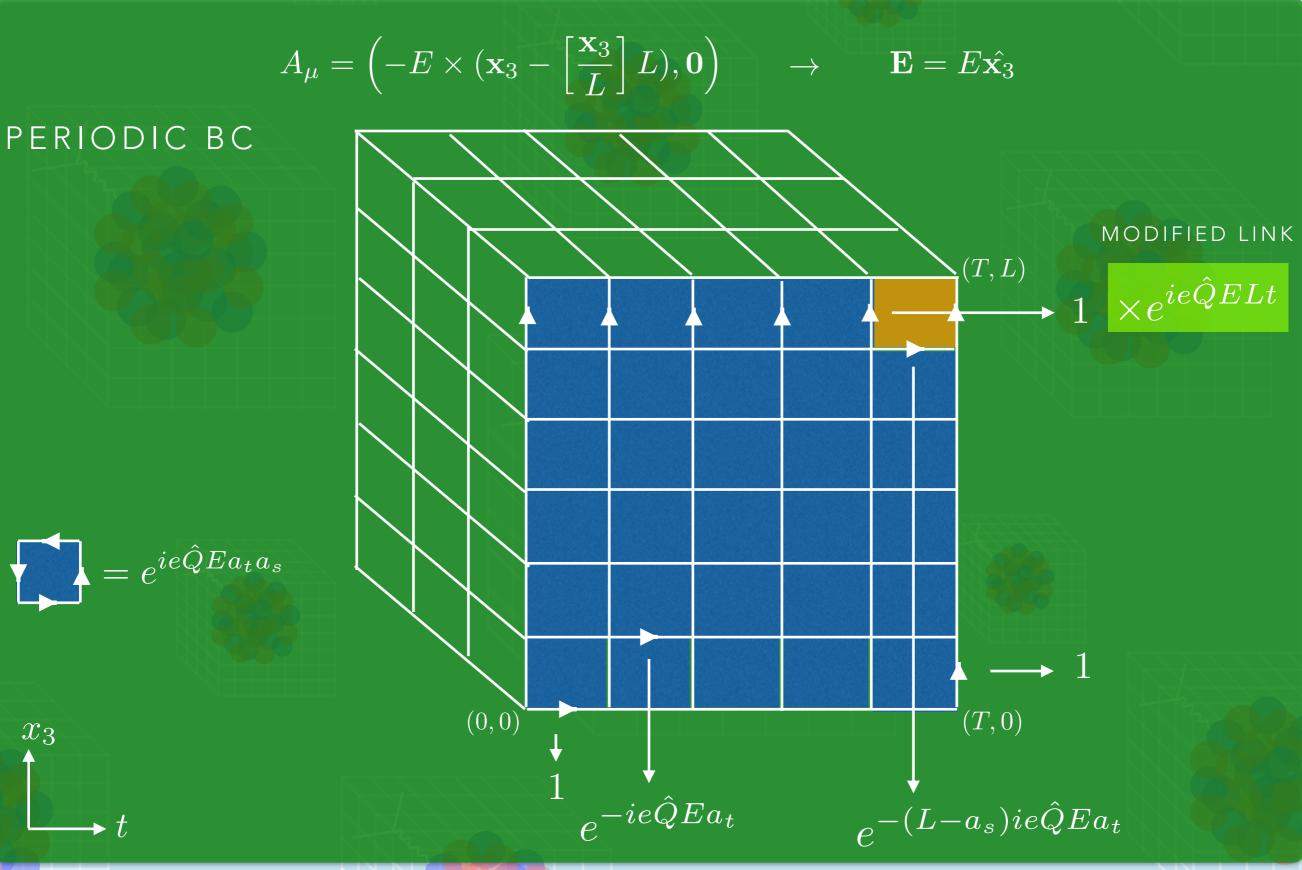


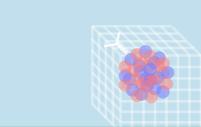


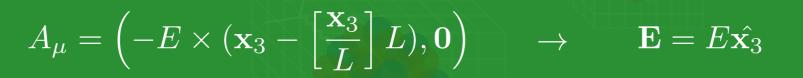












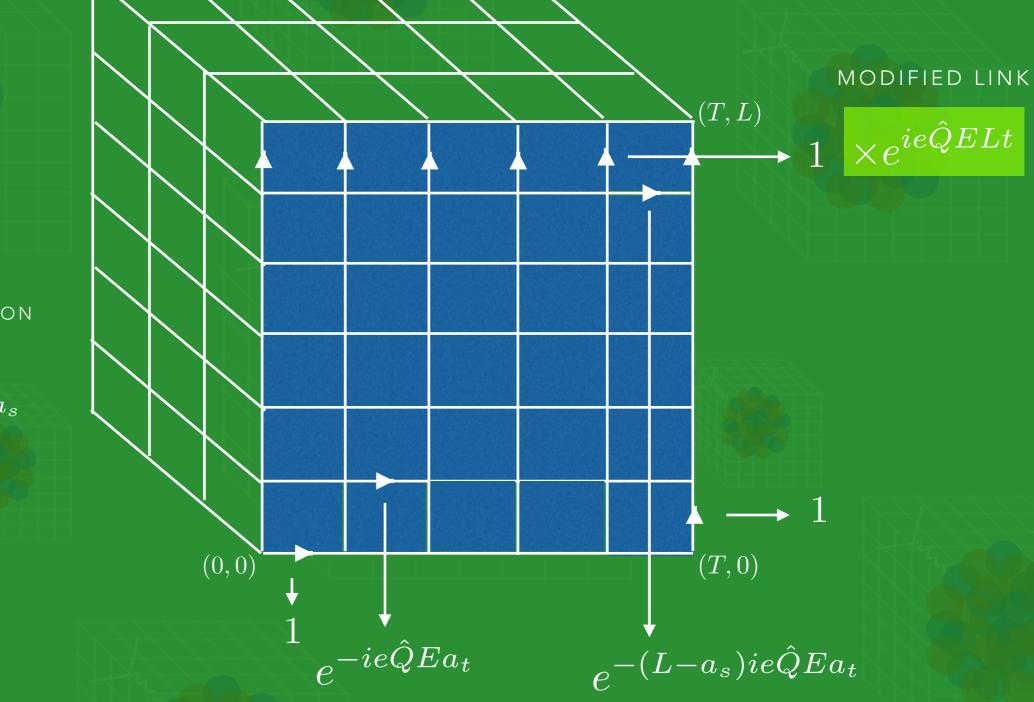
#### PERIODIC BC

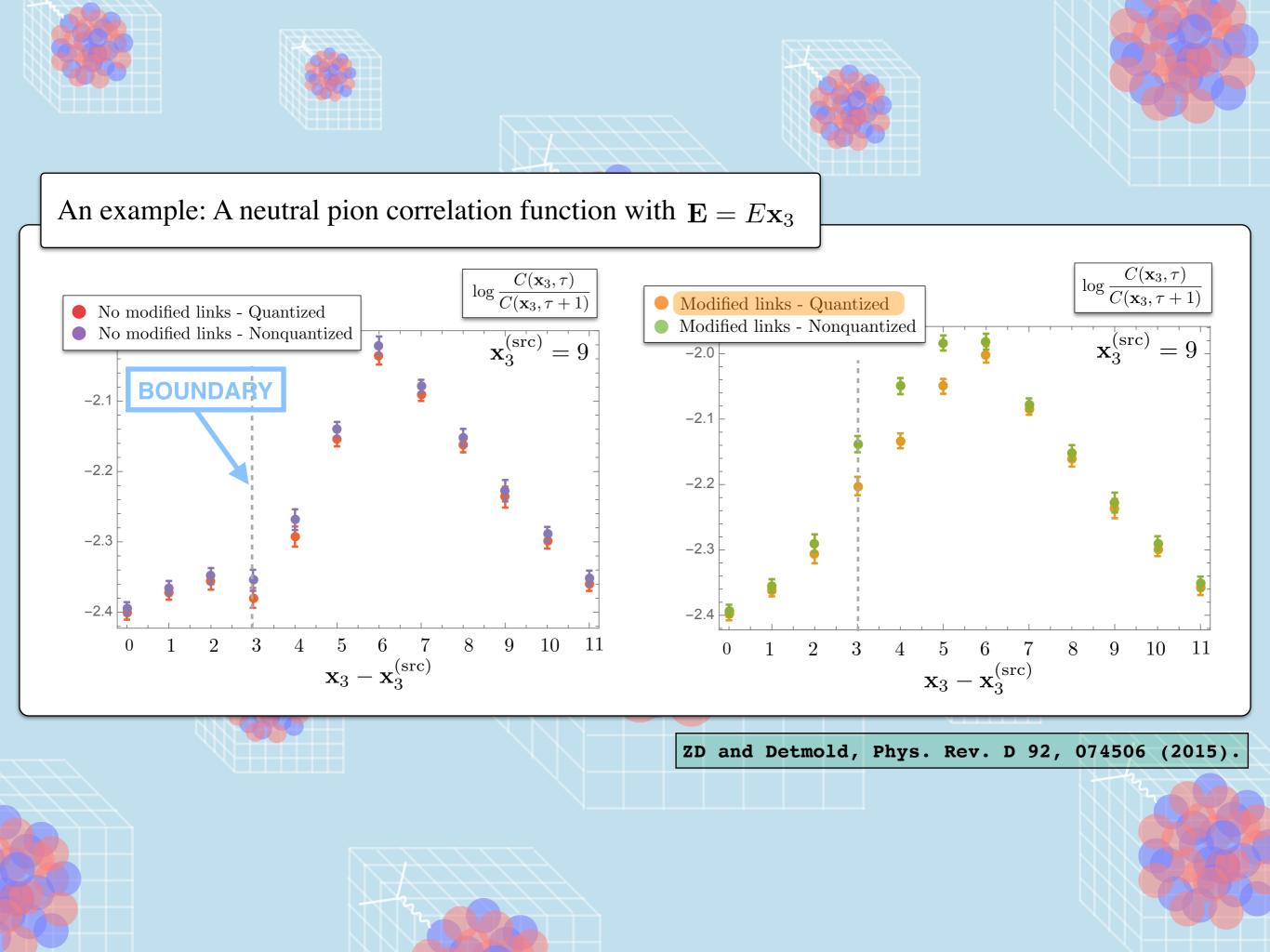
 $e^{ie\hat{Q}ELT} = 1$ 

 $E = \frac{2\pi n}{e\hat{Q}TL}$ 

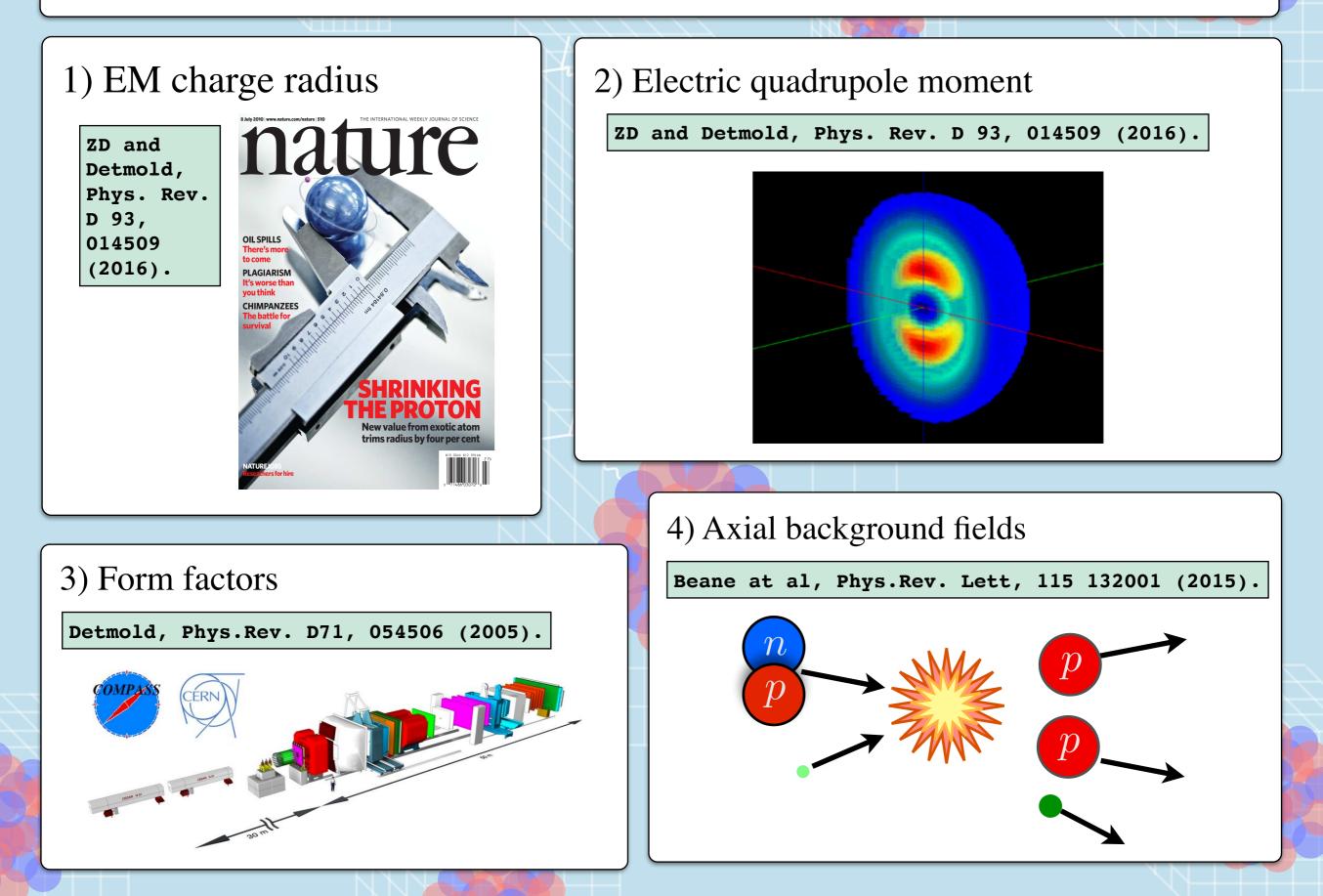
### 'T HOOFT QUANTIZATION CONDITION

$$\mathbf{T} = e^{ie\hat{Q}Ea_ta_s}$$





Various other structure properties of hadrons and nuclei, as well as their transitions, can be studied using more complex background fields:

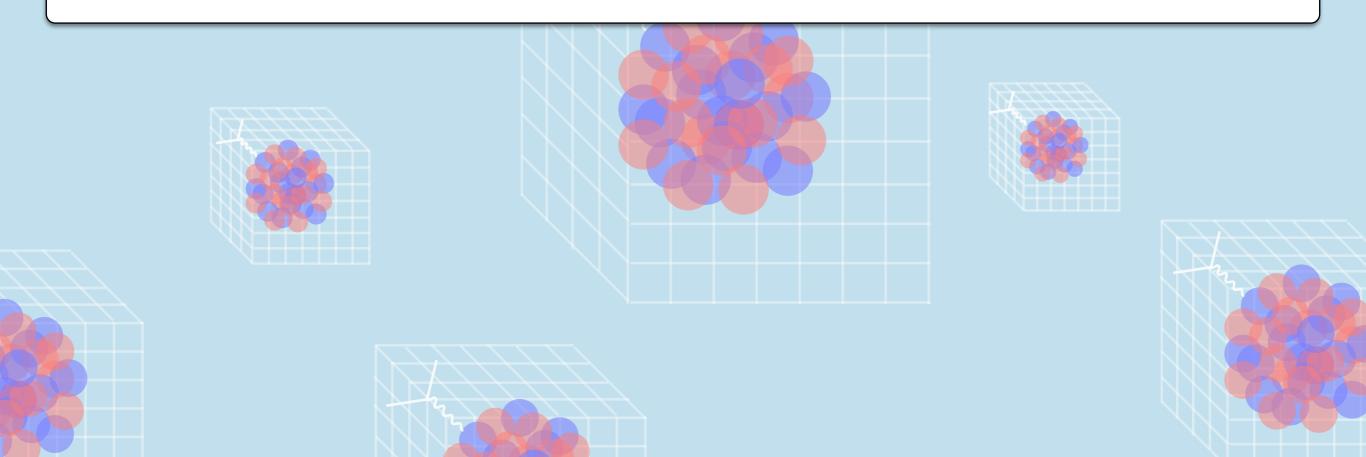


Consider a non-uniform background electric field that is produced by a background gauge potential

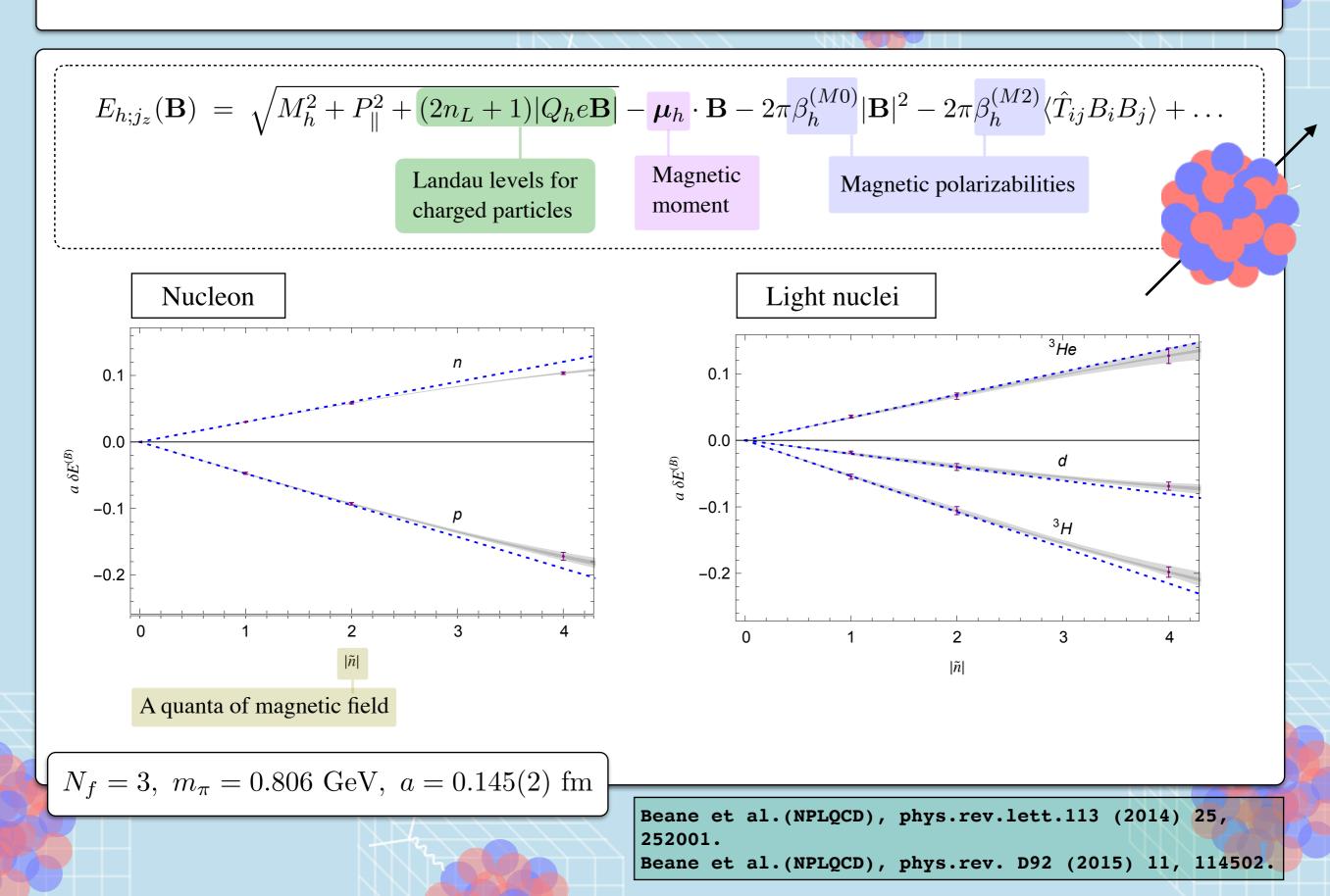
$$A_{\mu} = \left(-\frac{E_0}{2}(\mathbf{x}_3 - R - \left[\frac{\mathbf{x}_3}{L}\right]L)^2, \mathbf{0}\right)$$

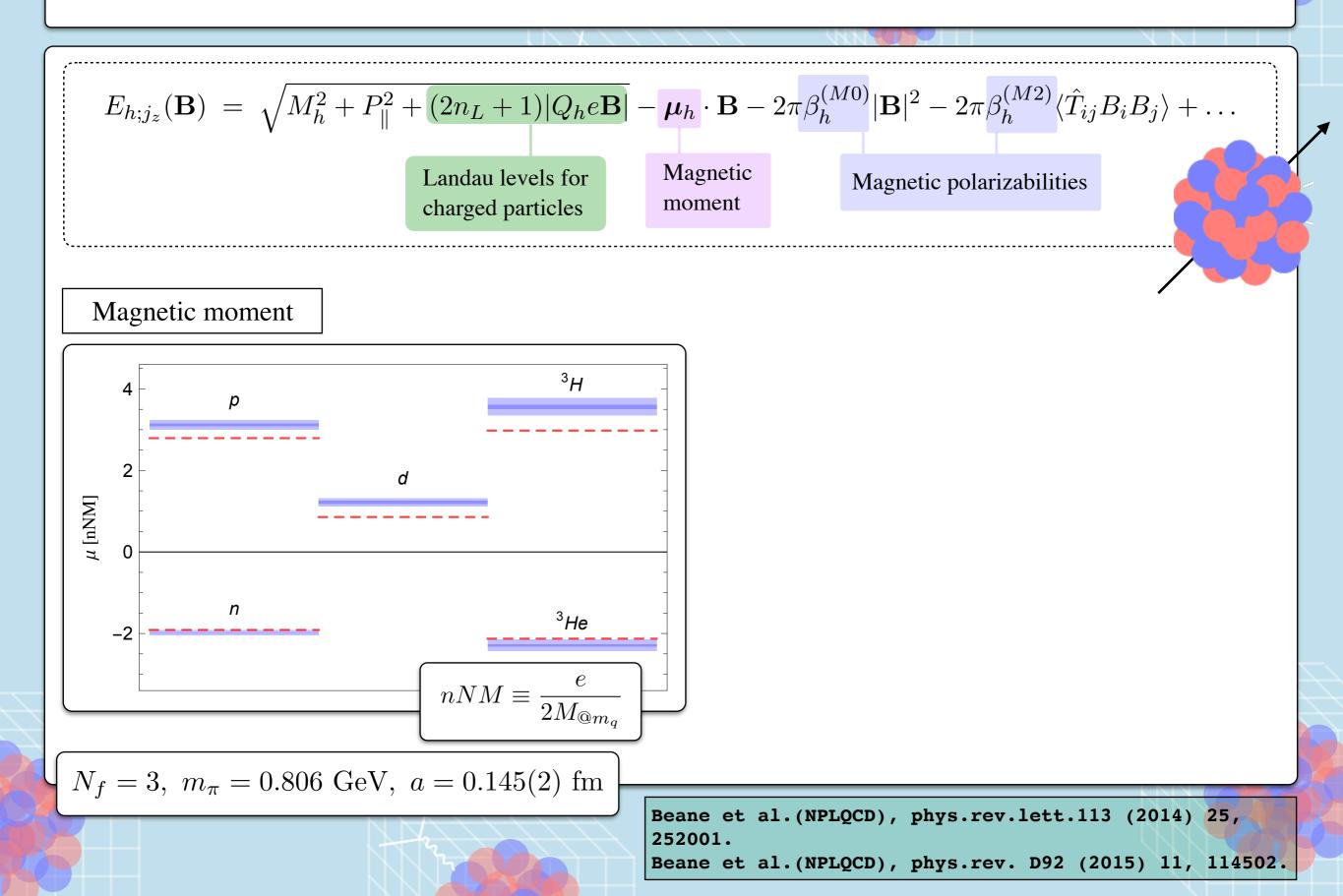
EXERCISE 7

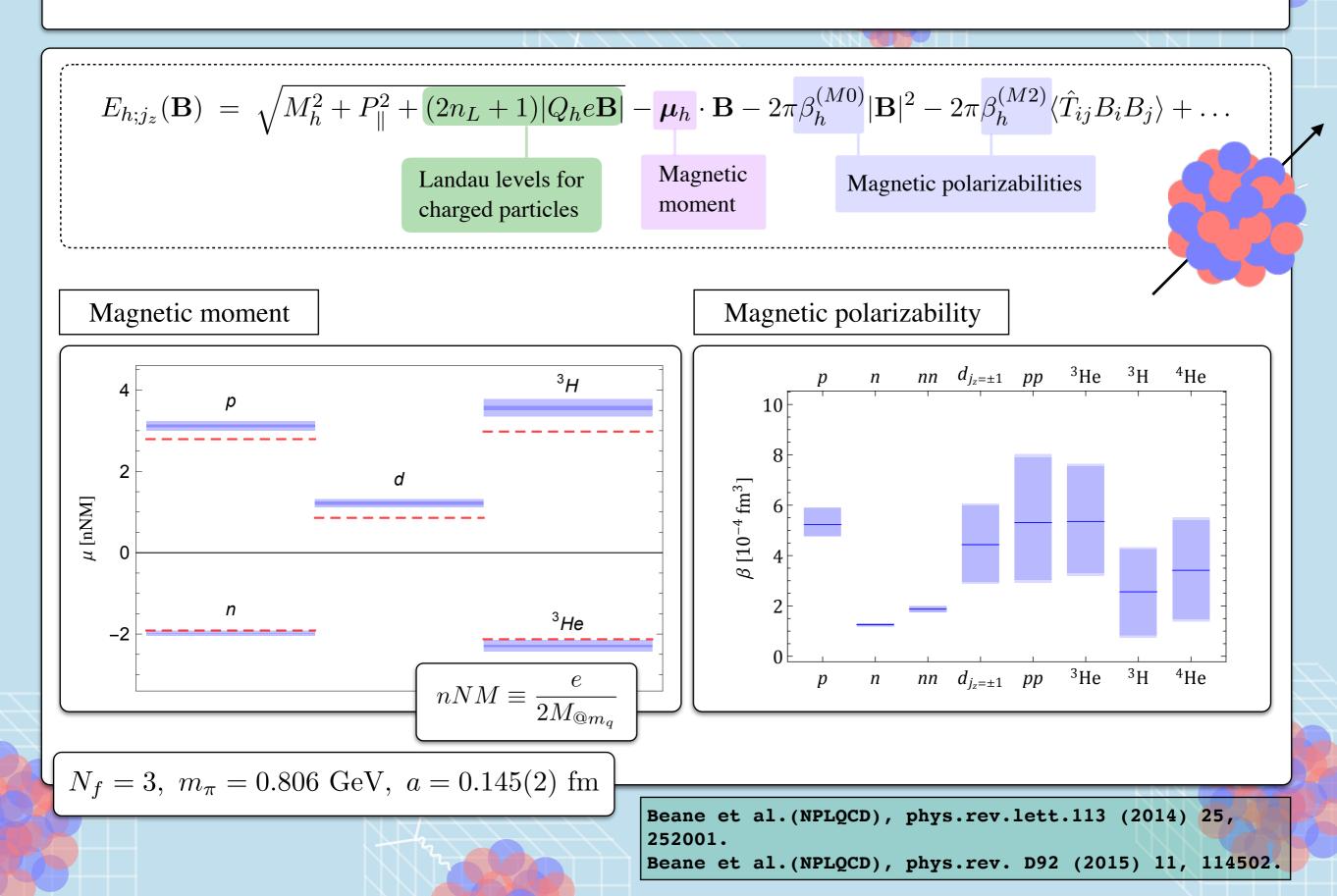
where  $0 \le R < L$ . Derive the prescription for the modified links as well as a quantization condition on the slope of the electric field strength,  $E_0$ , along direction  $\hat{\mathbf{x}}_3$ .



$$\begin{aligned} F_{h;j_{z}}(\mathbf{B}) &= \sqrt{M_{h}^{2} + P_{\parallel}^{2} + (2n_{L} + 1)|Q_{h}e\mathbf{B}|} - \mu_{h} \cdot \mathbf{B} - 2\pi\beta_{h}^{(M0)}|\mathbf{B}|^{2} - 2\pi\beta_{h}^{(M2)}\langle\hat{T}_{ij}B_{i}B_{j}\rangle + \dots \\ \text{Landau levels for charged particles} & Magnetic moment & Magnetic polarizabilities & Mag$$







Let's enumerate a some of the methods that give access to structure quantities in general:

#### Three(four)-point functions

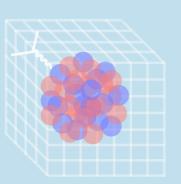
For e.g., form factors, moments of structure functions, Compton amplitude, transition amplitudes

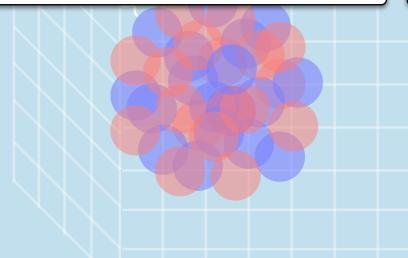
## Background-field methods

For e.g., EM moments and polarizabilities, charge radius, form factors and transition amplitudes.

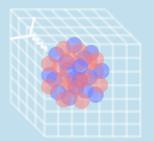
# Feynman-Hellmann inspired methods

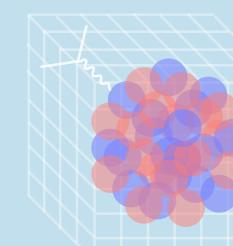
Similar to background fields. For e.g., axial charge, form factors, EM moments, transition amplitudes

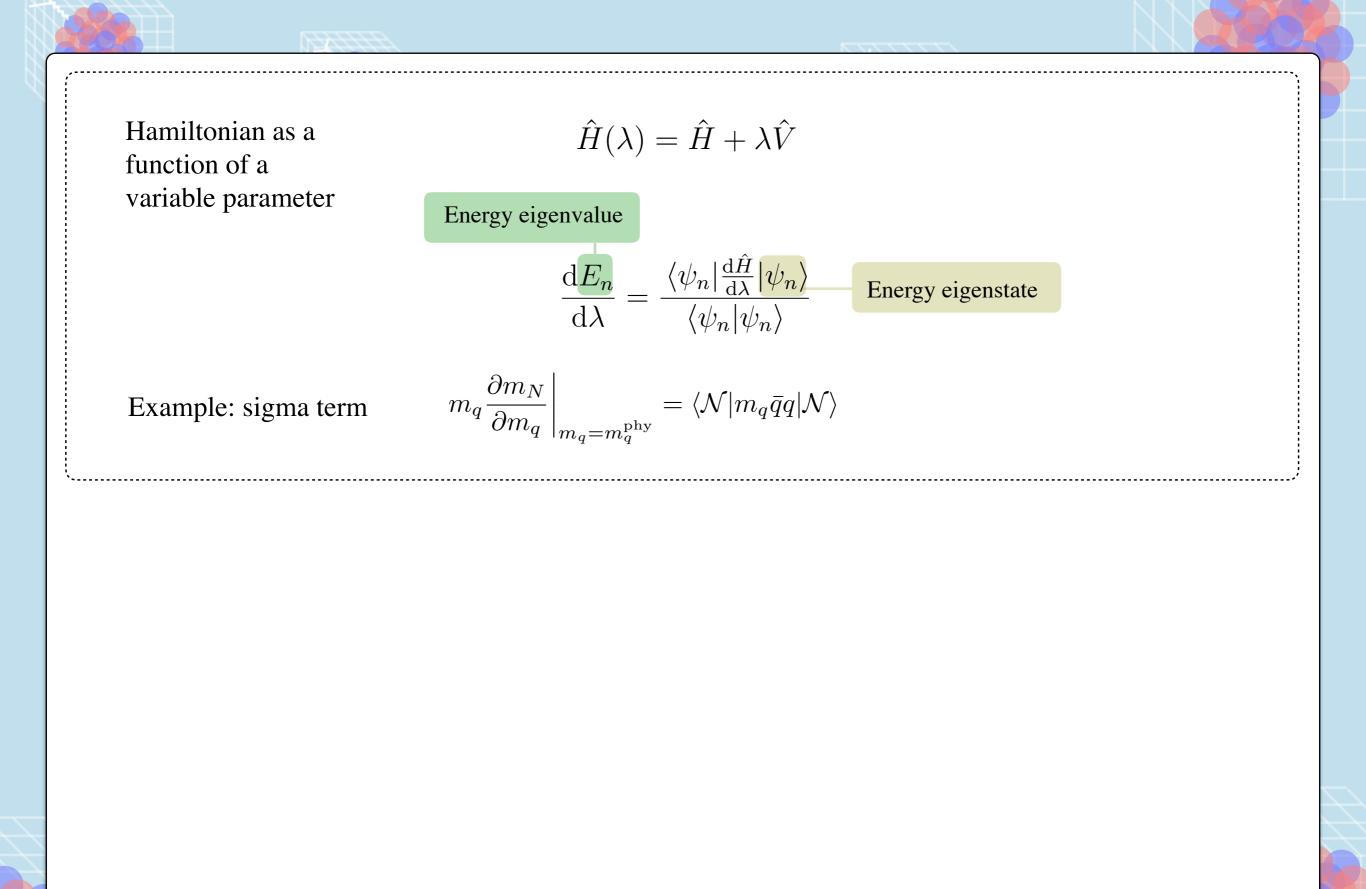


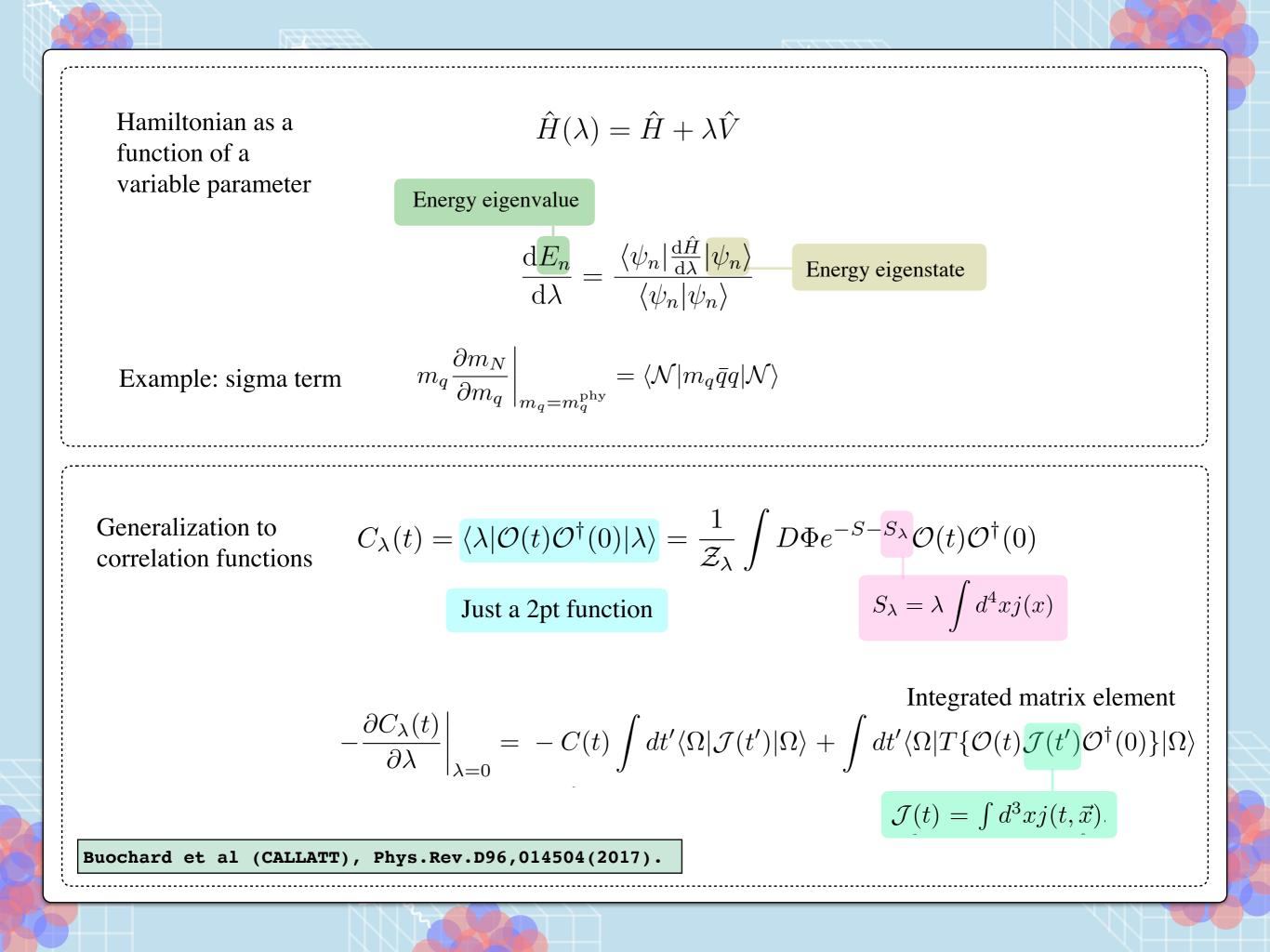












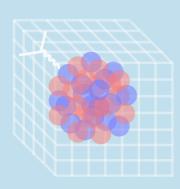
Example: axial charge of the nucleon and triton!

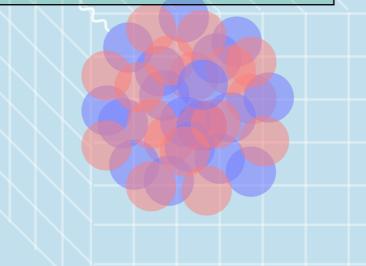
Since the operator here is a quark bilinear, a clever to implement this is by modifying the quark propagator.

 $S_{\lambda_{q};\Gamma}^{(q)}(x,y) = S^{(q)}(x,y) + \lambda_{q} \int dz \ S^{(q)}(x,z)\Gamma S^{(q)}(z,y)$ 

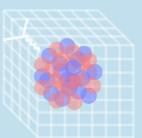
Savage et al (NPLQCD), Phys.Rev.Lett.119,062002(2017).

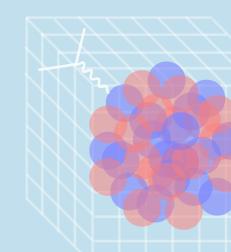
Buochard et al (CALLATT), Phys.Rev.D96,014504(2017).









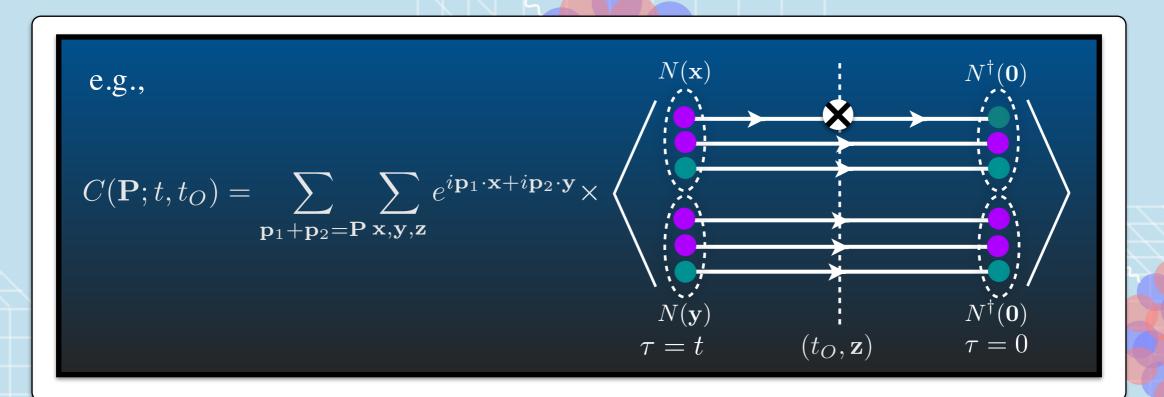


Example: axial charge of the nucleon and triton!

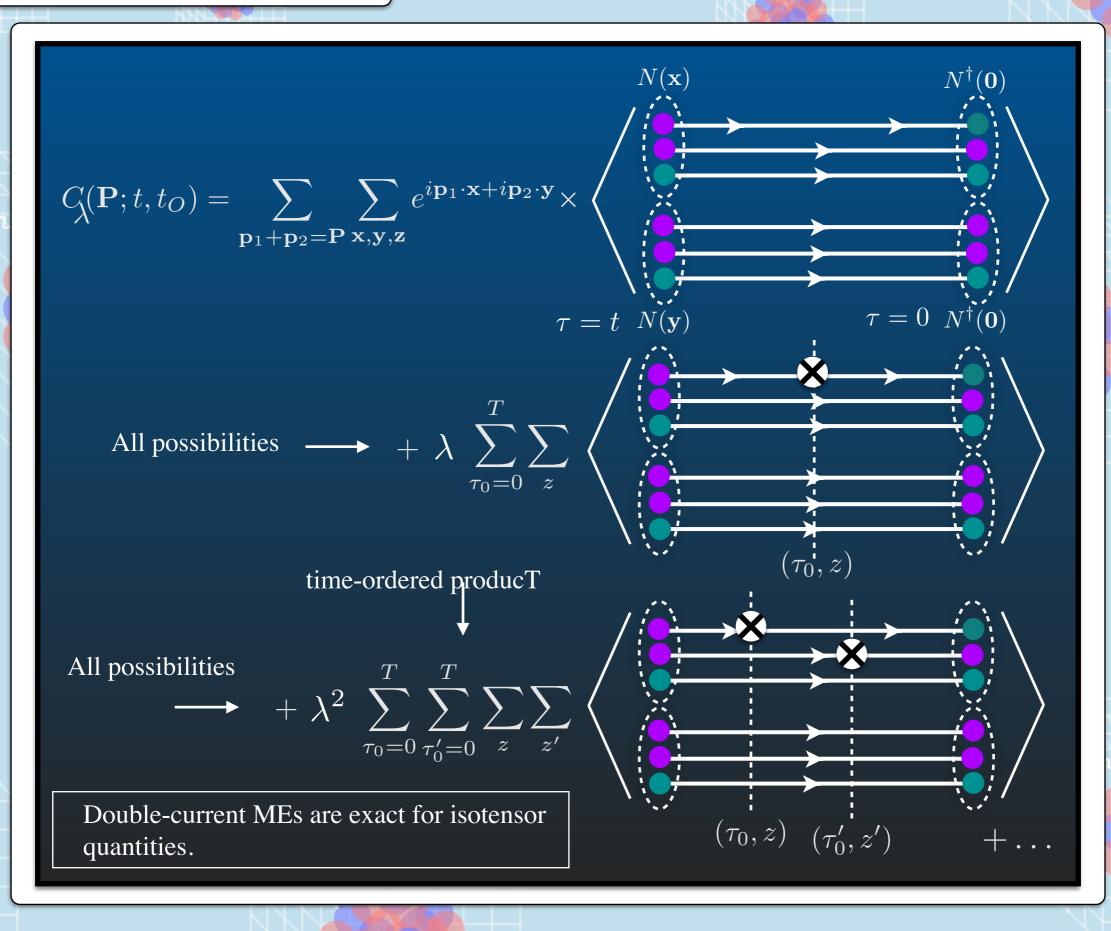
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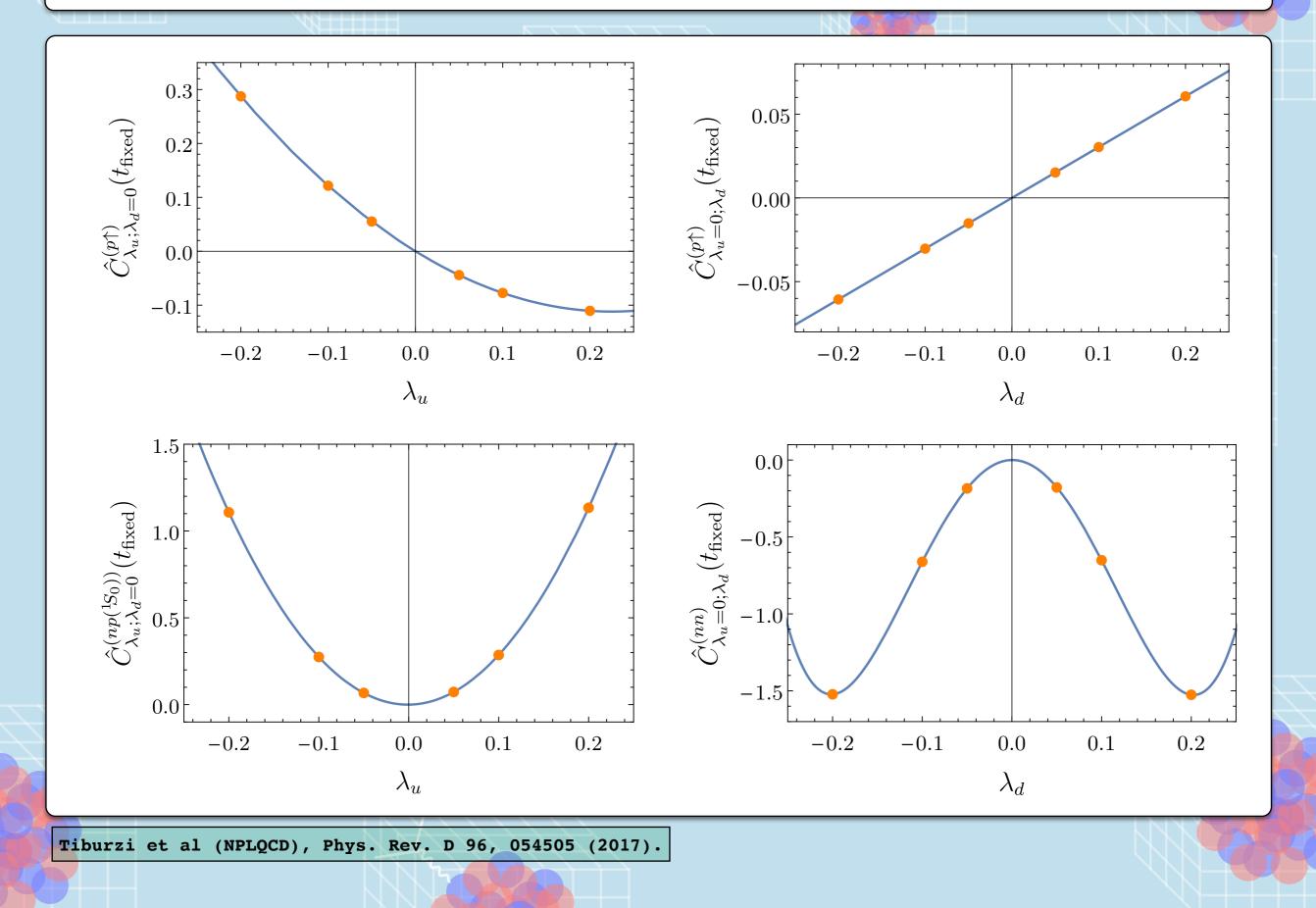
Buochard et al (CALLATT), Phys.Rev.D96,014504(2017).







#### Matrix elements from a compound propagator/background field



2pt functions we calculate is:  $\begin{aligned}
\text{Has information about the} \\
\text{matrix element we want}
\end{aligned}$   $\begin{aligned}
C_{\lambda_{u};\lambda_{d}=0}^{(p\uparrow)}(t) &= \sum_{\boldsymbol{x}} \left( \left\langle 0|\chi_{p\uparrow}(\boldsymbol{x},t)\chi_{p\uparrow}^{\dagger}(0)|0\rangle + \lambda_{u}\sum_{\boldsymbol{y}}\sum_{t_{1}=0}^{t} \left\langle 0|\chi_{p\uparrow}(\boldsymbol{x},t)J_{3}^{(u)}(\boldsymbol{y},t_{1})\chi_{p\uparrow}^{\dagger}(0)|0\rangle \right. \right) + d_{2}\lambda_{u}^{2}, \\
C_{\lambda_{u}=0;\lambda_{d}}^{(p\uparrow)}(t) &= \sum_{\boldsymbol{x}} \left( \left\langle 0|\chi_{p\uparrow}(\boldsymbol{x},t)\chi_{p\uparrow}^{\dagger}(0)|0\rangle + \lambda_{d}\sum_{\boldsymbol{y}}\sum_{t_{1}=0}^{t} \left\langle 0|\chi_{p\uparrow}(\boldsymbol{x},t)J_{3}^{(d)}(\boldsymbol{y},t_{1})\chi_{p\uparrow}^{\dagger}(0)|0\rangle \right. \right),
\end{aligned}$ 

Tiburzi et al (NPLQCD), Phys. Rev. D 96, 054505 (2017).

2pt functions we calculate is:  

$$C_{\lambda_{u};\lambda_{d}=0}^{(p\uparrow)}(t) = \sum_{x} \left( \left\langle 0 | \chi_{p\uparrow}(x,t) \chi_{p\uparrow}^{\dagger}(0) | 0 \right\rangle + \lambda_{u} \sum_{y} \sum_{t_{1}=0}^{t} \left\langle 0 | \chi_{p\uparrow}(x,t) J_{3}^{(u)}(y,t_{1}) \chi_{p\uparrow}^{\dagger}(0) | 0 \right\rangle \right) + d_{2}\lambda_{u}^{2},$$

$$C_{\lambda_{u}=0;\lambda_{d}}^{(p\uparrow)}(t) = \sum_{x} \left( \left\langle 0 | \chi_{p\uparrow}(x,t) \chi_{p\uparrow}^{\dagger}(0) | 0 \right\rangle + \lambda_{d} \sum_{y} \sum_{t_{1}=0}^{t} \left\langle 0 | \chi_{p\uparrow}(x,t) J_{3}^{(u)}(y,t_{1}) \chi_{p\uparrow}^{\dagger}(0) | 0 \right\rangle \right),$$
Taking the linear term:  

$$C_{\lambda_{u};\lambda_{d}=0}^{(p\uparrow)}(t) \Big|_{\mathcal{O}(\lambda_{u})} = \sum_{x,y} \sum_{t_{1}=0}^{t} \left\langle 0 | \chi_{p\uparrow}(x,t) J_{3}^{(u)}(y,t_{1}) \chi_{p\uparrow}^{\dagger}(0) | 0 \right\rangle$$

$$= \sum_{n,m} \sum_{x,y} \sum_{t_{1}=0}^{t} \left\langle 0 | \chi_{p\uparrow}(x,t) | n \rangle \langle n | J_{3}^{(u)}(y,t_{1}) | m \rangle \langle m | \chi_{p\uparrow}^{\dagger}(0) | 0 \rangle$$
Insert two complete set of states

2pt functions we calculate is:  

$$C_{\lambda_{u};\lambda_{\sigma}=0}^{(p\uparrow)}(t) = \sum_{x} \left( \langle 0|\chi_{p\uparrow}(x,t)\chi_{p\uparrow}^{\dagger}(0)|0\rangle + \lambda_{u}\sum_{y}\sum_{l_{1}=0}^{t} \langle 0|\chi_{p\uparrow}(x,t)J_{3}^{(u)}(y,t_{1})\chi_{p\uparrow}^{\dagger}(0)|0\rangle \right) + d_{2}\lambda_{u}^{2},$$

$$C_{\lambda_{u}=0;\lambda_{d}}^{(p\uparrow)}(t) = \sum_{x} \left( \langle 0|\chi_{p\uparrow}(x,t)\chi_{p\uparrow}^{\dagger}(0)|0\rangle + \lambda_{d}\sum_{y}\sum_{l_{1}=0}^{t} \langle 0|\chi_{p\uparrow}(x,t)J_{3}^{(u)}(y,t_{1})\chi_{p\uparrow}^{\dagger}(0)|0\rangle \right),$$
Taking the linear term:  

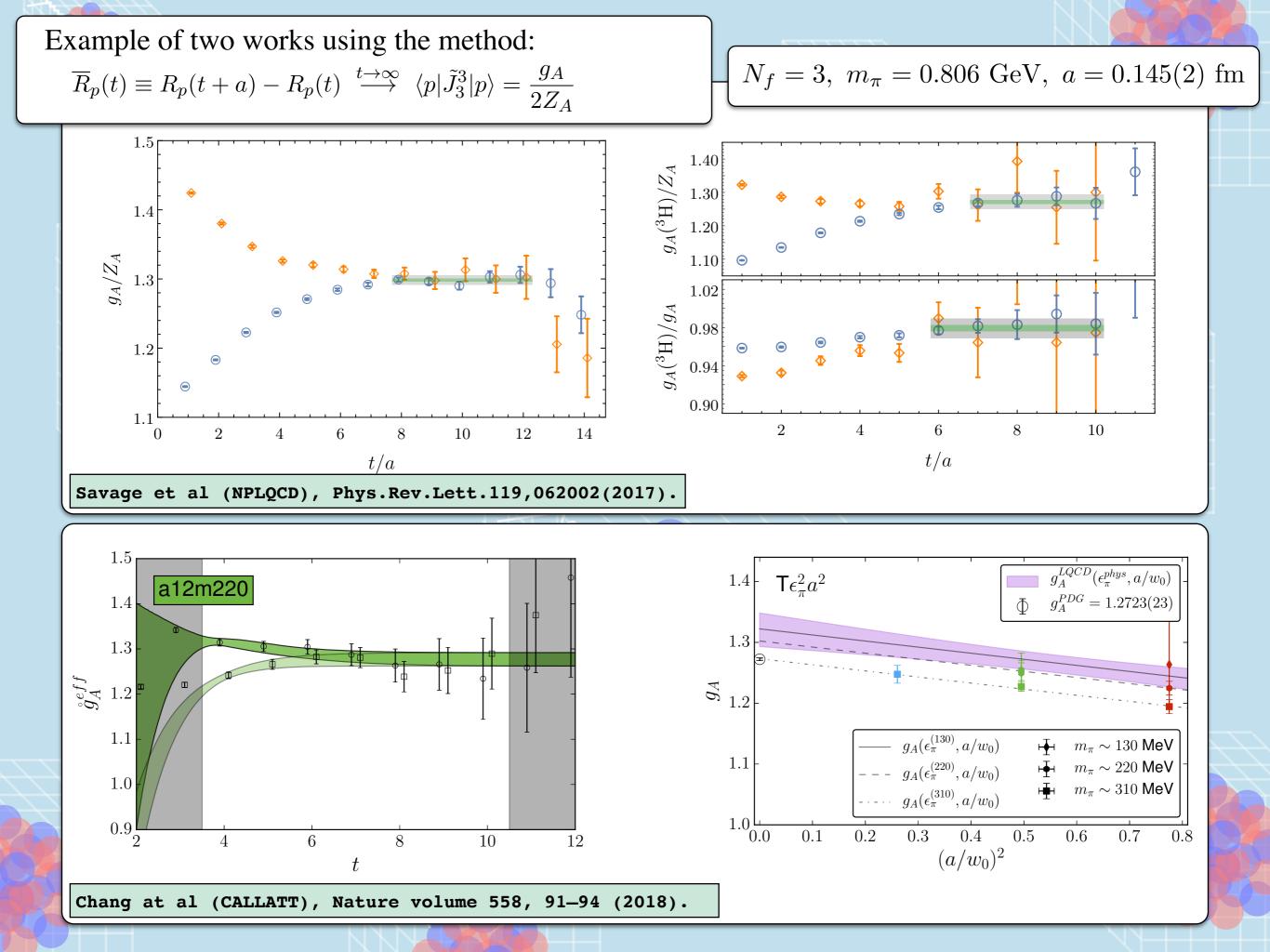
$$C_{\lambda_{u};\lambda_{d}=0}^{(p\uparrow)}(t)\Big|_{\mathcal{O}(\lambda_{u})} = \sum_{x,y}\sum_{l_{1}=0}^{t} \langle 0|\chi_{p\uparrow}(x,t)J_{3}^{(u)}(y,t_{1})\chi_{p\uparrow}^{\dagger}(0)|0\rangle$$

$$= \sum_{n,m}\sum_{x,y}\sum_{l_{1}=0}^{t} \langle 0|\chi_{p\uparrow}(x,t)M_{3}^{(u)}(y,t_{1})\chi_{p\uparrow}^{\dagger}(0)|0\rangle$$
Insert two complete set of states  
Isolate the ground state matrix element:  

$$C_{\lambda_{u};\lambda_{d}=0}^{(p\uparrow)}(t)\Big|_{\mathcal{O}(\lambda_{u})} = \sum_{l_{1}=0}^{t}\sum_{n,m}z_{n}z_{m}^{\dagger}e^{-E_{n}(t-t_{1})}e^{-E_{m}t_{1}}\langle n|J_{3}^{(u)}|m\rangle$$
Energy gap to excited states  

$$\frac{t\to\infty}{|z_{0}|^{2}}e^{-E_{0}t}\left[c+t(p\uparrow)J_{3}^{(u)}|p\uparrow\right) + \mathcal{O}(e^{-\delta t})\right]$$
The desired matrix element:

Tiburzi et al (NPLQCD), Phys. Rev. D 96, 054505 (2017).



Let's enumerate a some of the methods that give access to structure quantities in general:

#### Three(four)-point functions

For e.g., form factors, moments of structure functions, Compton amplitude, transition amplitudes

### Background-field methods

For e.g., EM moments and polarizabilities, charge radius, form factors and transition amplitudes.

### Feynman-Hellmann inspired methods

Similar to background fields. For e.g., axial charge, form factors, EM moments, transition amplitudes

We did not discuss many other interesting directions in the field, e.g.,

Moments of structure functions

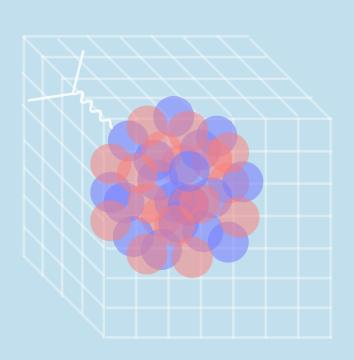
Hadron tensor through inverse transform methods

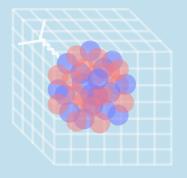
Quasi-PDFs and pseudo-PDFs

GPDs, TMDs, gluonic observables, etc.

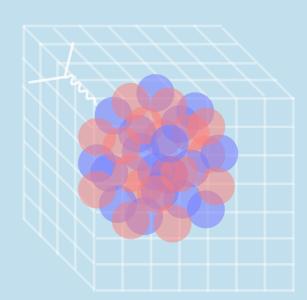
LECTURE III: NUCLEAR STRUCTURE, CHALLENGES AND PROGRESS

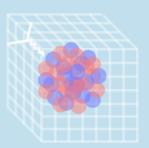
What about nuclear observables? Let's see the application of these methods to two examples:

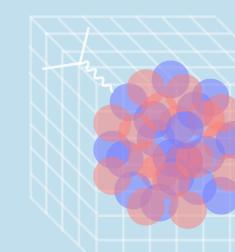


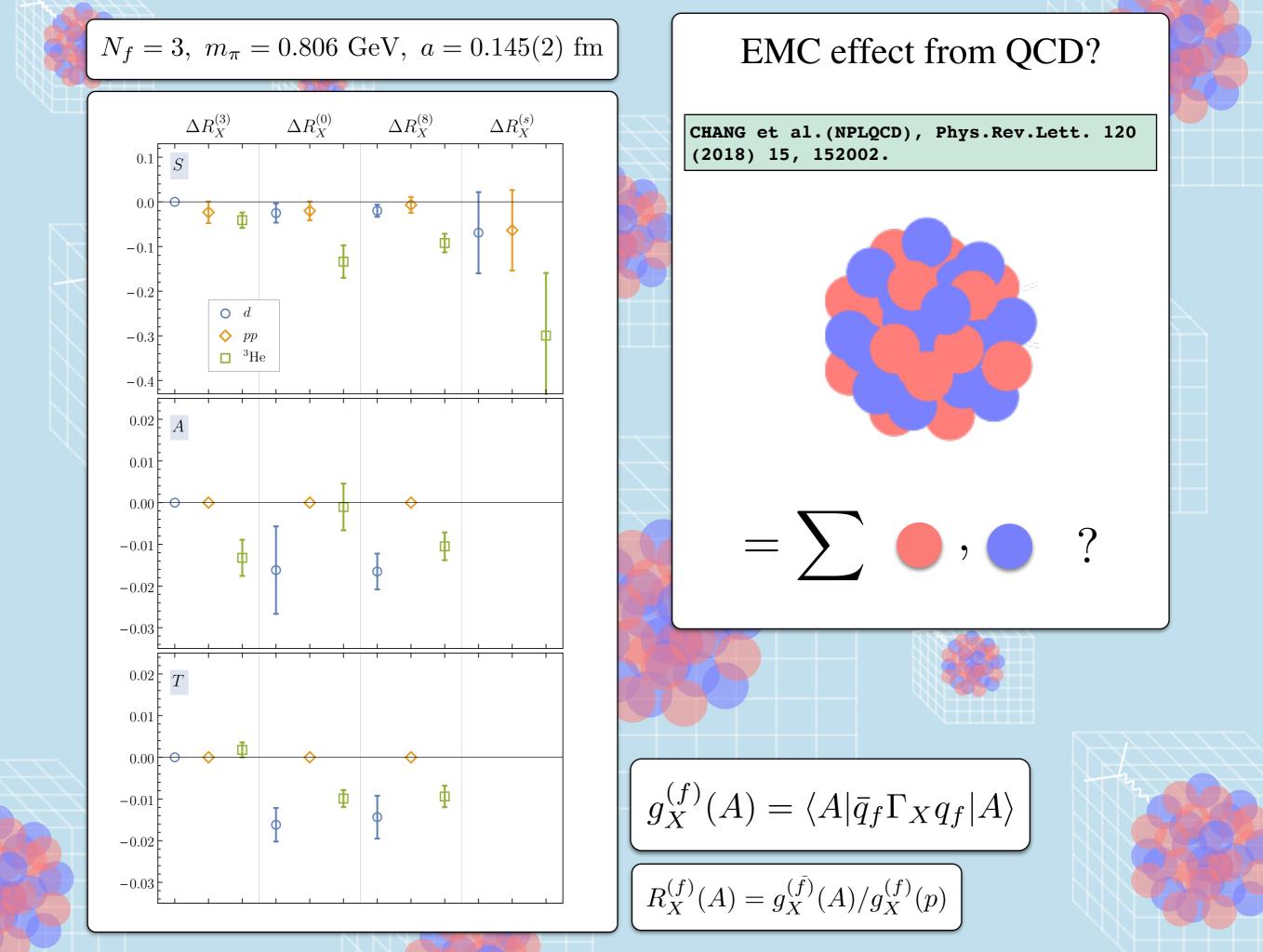


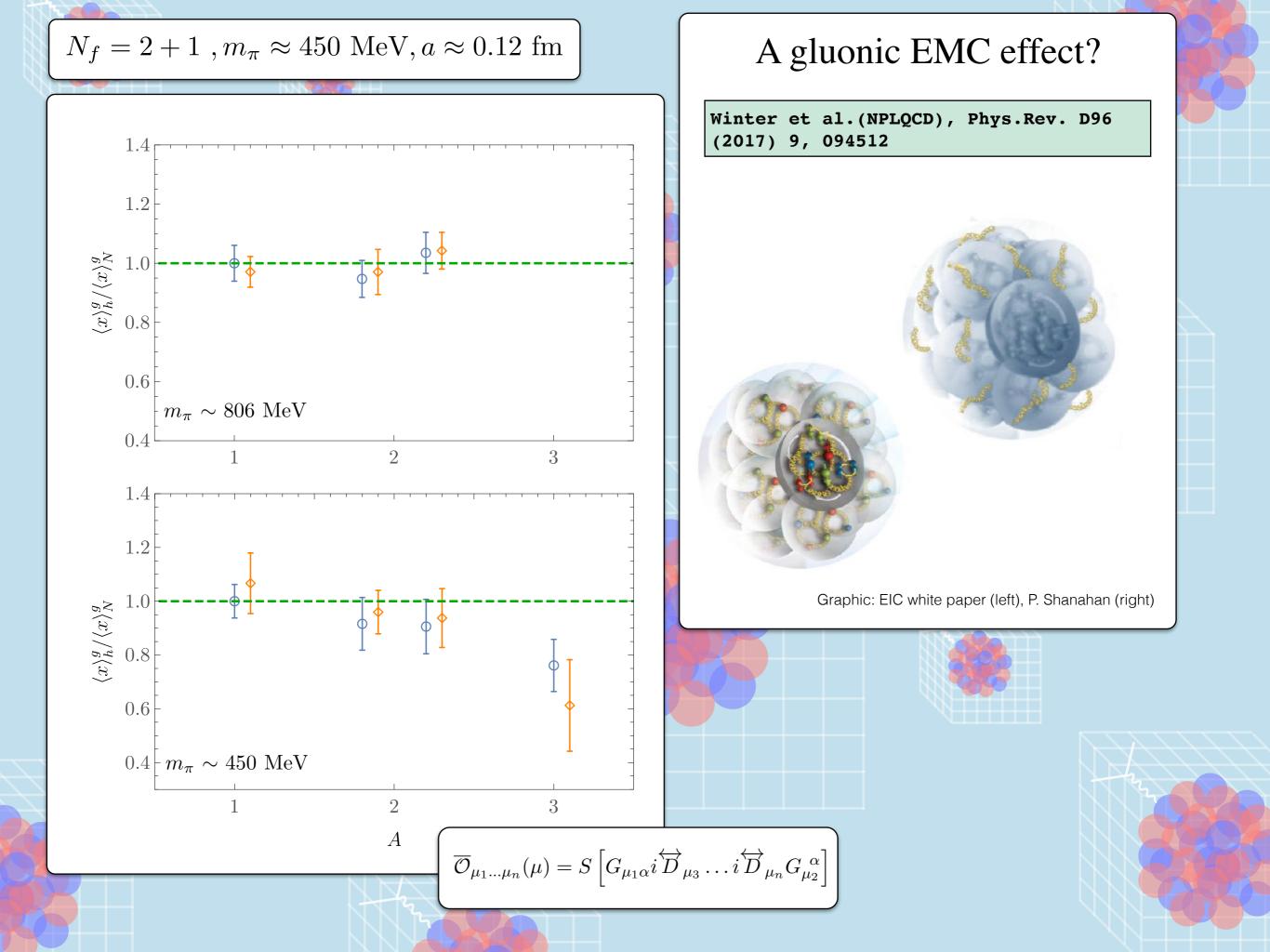


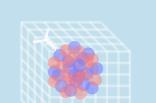














Three features make lattice QCD calculations of nuclei hard:

i) The complexity of systems grows rapidly with the number of quarks.

Detmold and Orginos, Phys. Rev. D 87, 114512 (2013). See also: Detmold and Savage, Phys.Rev.D82 014511 (2010). Doi and Endres, Comput. Phys. Commun. 184 (2013) 117.

ii) Excitation energies of nuclei are much smaller than the QCD scale.

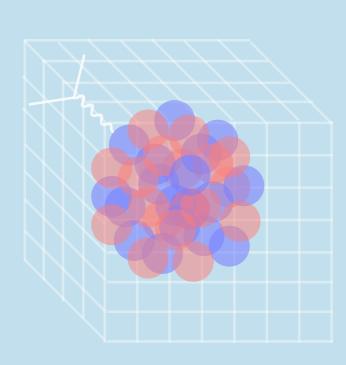
Beane at al (NPLQCD), Phys.Rev.D79 114502 (2009). Beane, Detmold, Orginos, Savage, Prog. Part. Nucl. Phys. 66 (2011). Junnakar and Walker-Loud, Phys.Rev. D87 (2013) 114510. Briceno, Dudek and Young, Rev. Mod. Phys. 90 025001.

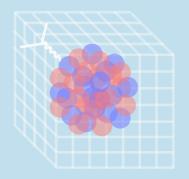
iii) There is a severe signal-to-noise degradation.

Paris (1984) and Lepage (1989).

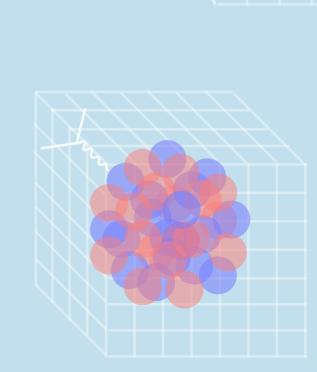
Wagman and Savage, Phys. Rev. D 96, 114508 (2017). Wagman and Savage, arXiv:1704.07356 [hep-lat].

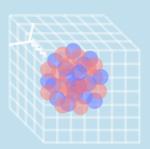
## i) The complexity of systems grows rapidly with the number of quarks.

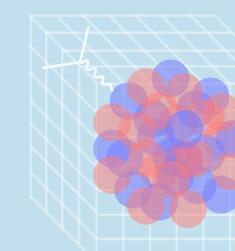


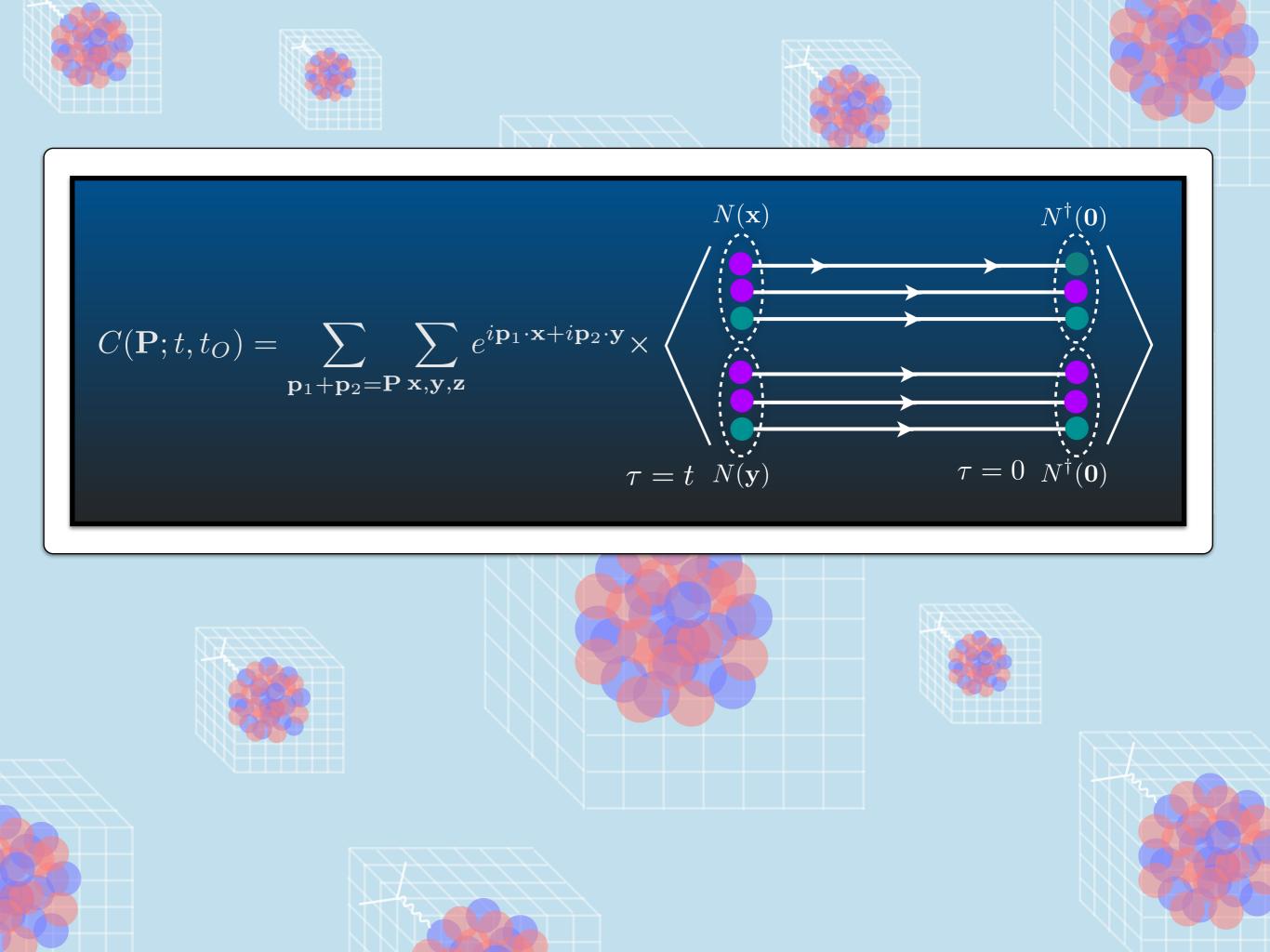


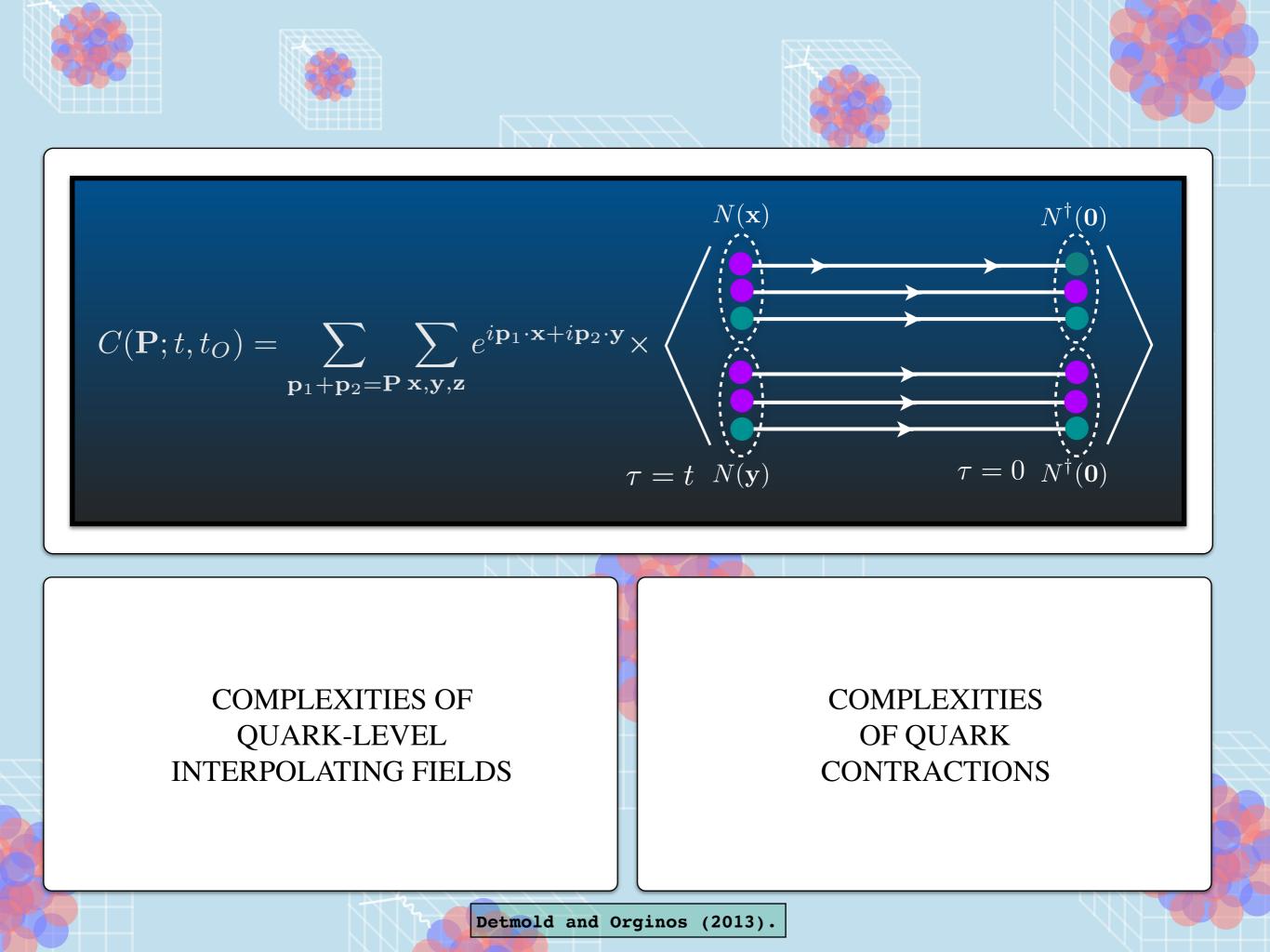










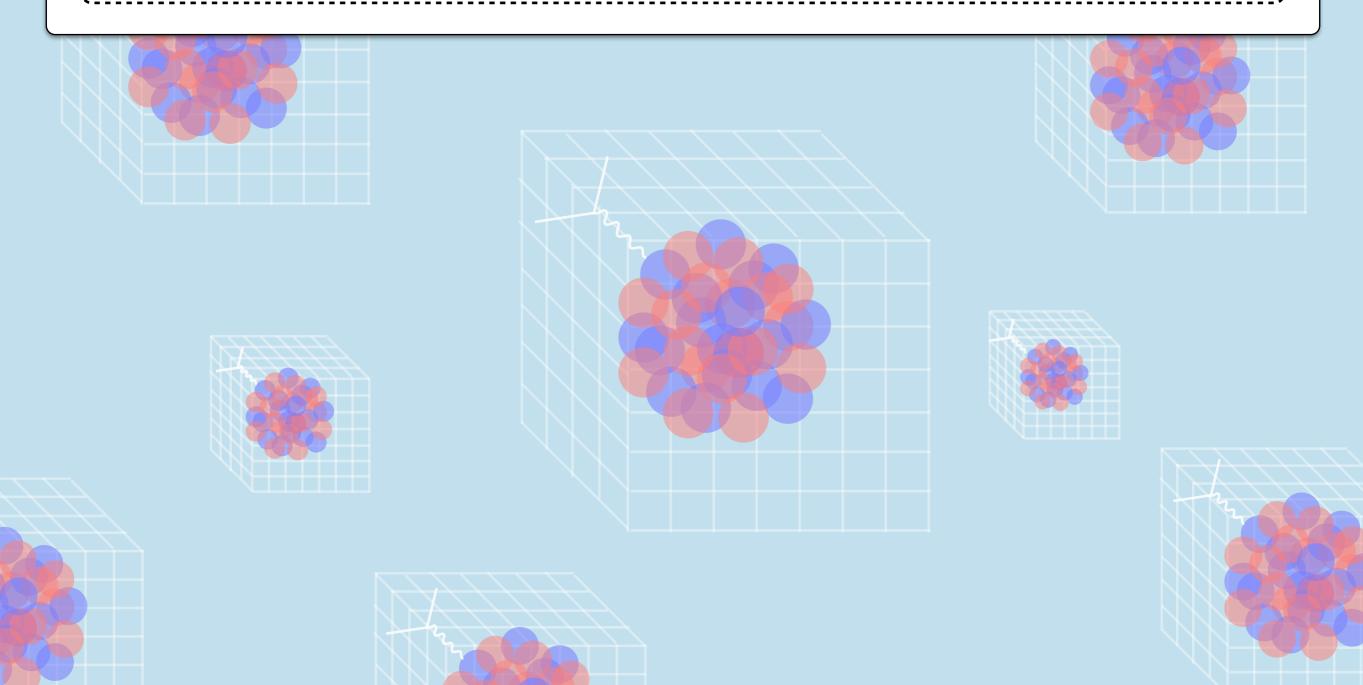


Naively the number of quark contractions for a nucleus goes as:

 $(2N_p + N_n)! (N_p + 2N_n)!$ 

How bad is this? Example: Consider radium-226 isotope. the number of contractions required is ~  $10^{1425}$ 





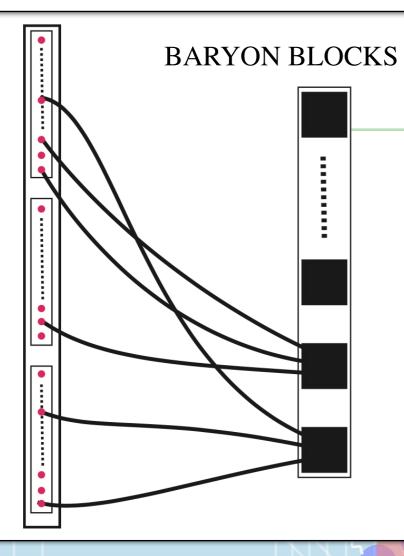
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An example of a more efficient algorithm:



$$\mathcal{B}_{b}^{a_{1},a_{2},a_{3}}(\mathbf{p},t;x_{0}) = \sum_{\mathbf{x}} e^{i\mathbf{p}\cdot\mathbf{x}} \sum_{k=1}^{N_{B(b)}} \tilde{w}_{b}^{(c_{1},c_{2},c_{3}),k}$$
$$\sum_{\mathbf{i}} \epsilon^{i_{1},i_{2},i_{3}} S(c_{i_{1}},x;a_{1},x_{0}) S(c_{i_{2}},x;a_{2},x_{0}) S(c_{i_{3}},x;a_{3},x_{0})$$

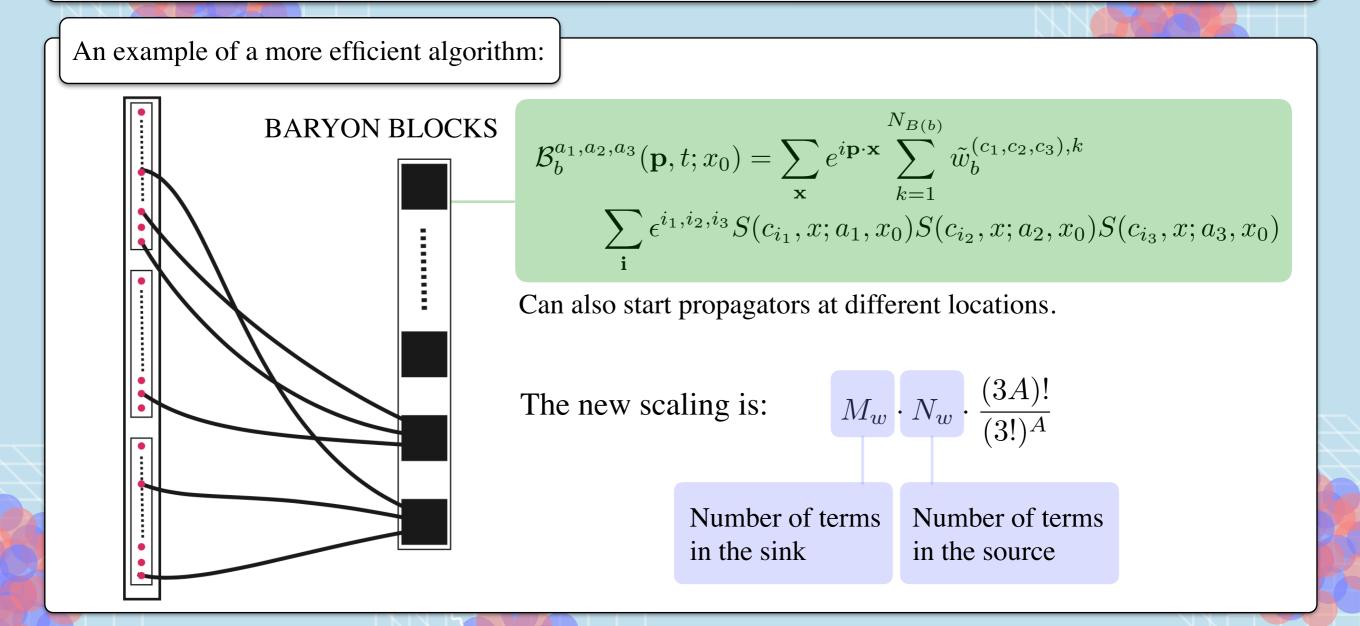
Can also start propagators at different locations.

Naively the number of quark contractions for a nucleus goes as:

 $(2N_p + N_n)! (N_p + 2N_n)!$ 

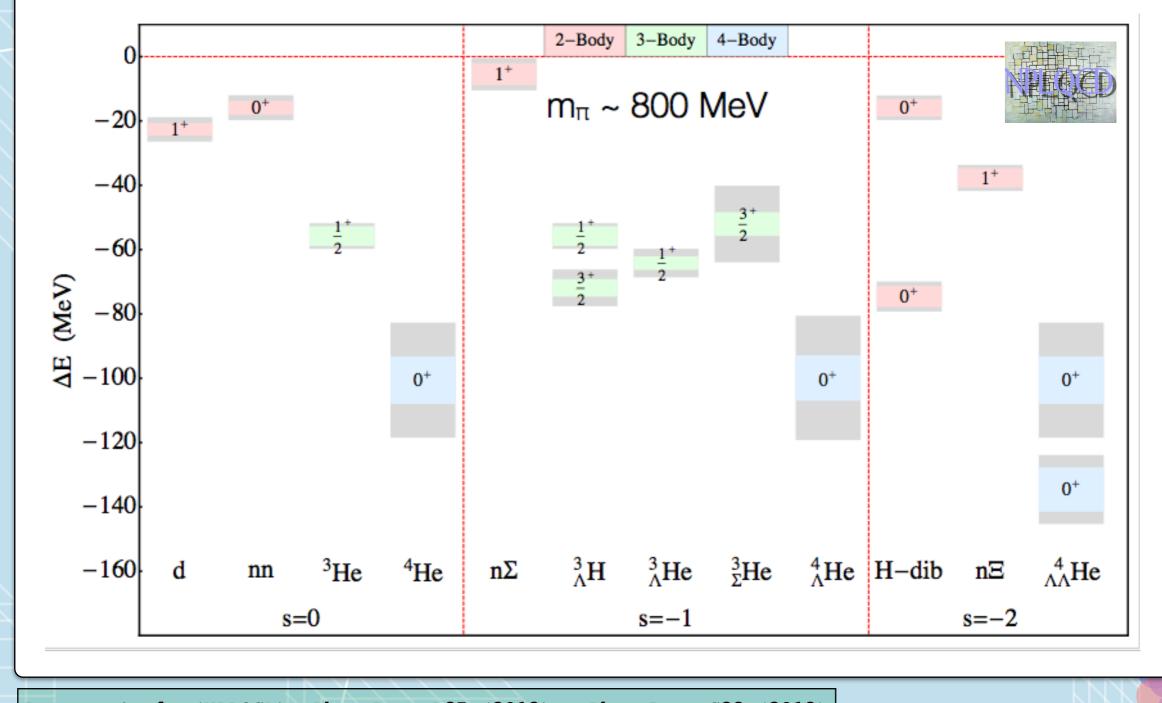
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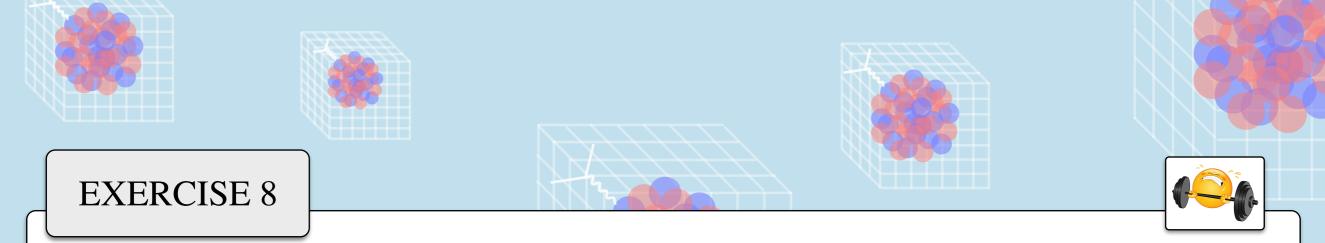


## Nuclei obtained from such an approach (at a heavier quark masses)

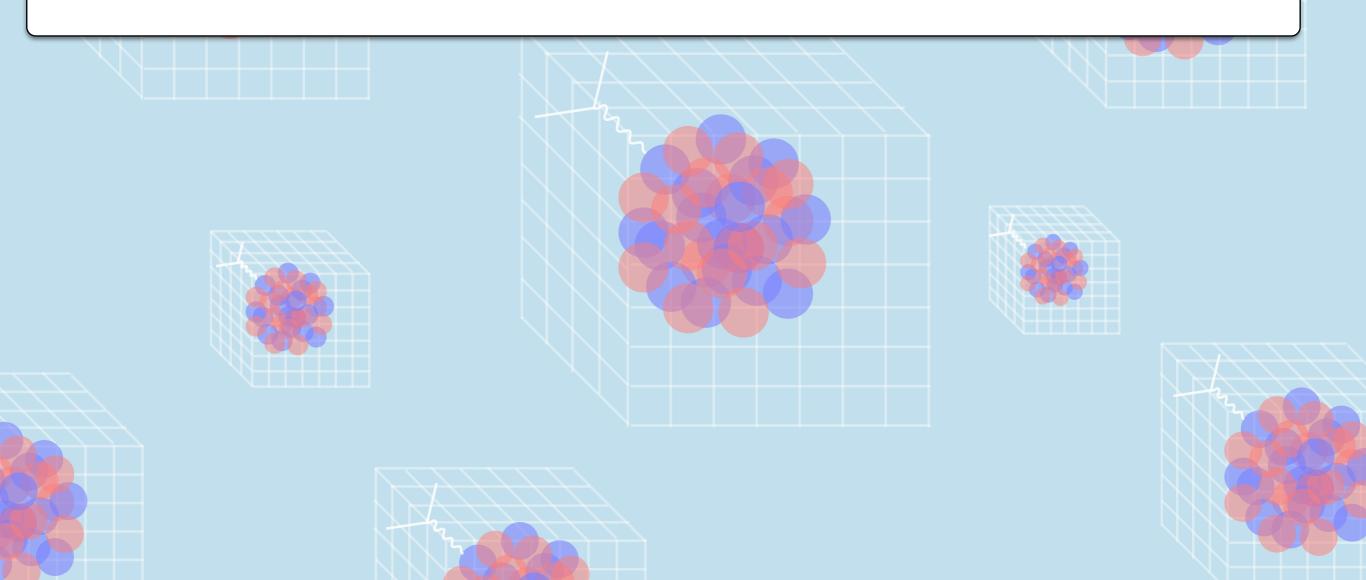
 $N_f = 3, \ m_\pi = 0.806 \text{ GeV}, \ a = 0.145(2) \text{ fm}$ 



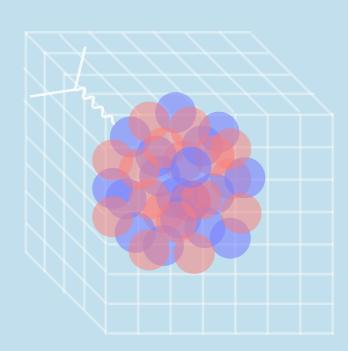
Beane, et al. (NPLQCD), Phys.Rev. D87 (2013) , Phys.Rev. C88 (2013)

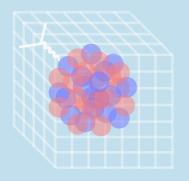


According to the naive counting, how many contractions are required for a nucleus at the source and sink with atomic numbers A = 4, 8, 12, 16? How many contractions are there with the use of the efficient algorithm described? There are even more optimal algorithms that lead to a polynomial scaling with the number of the quarks.

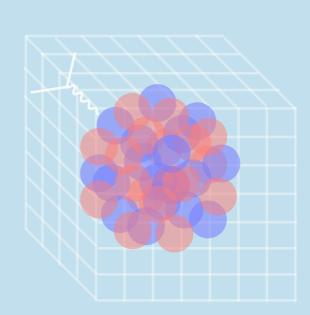


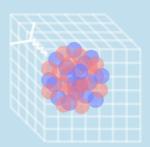
ii) Excitation energies of nuclei are much smaller than the QCD scale.

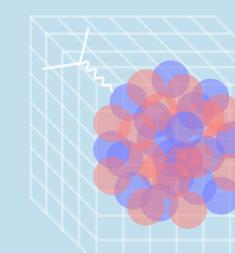


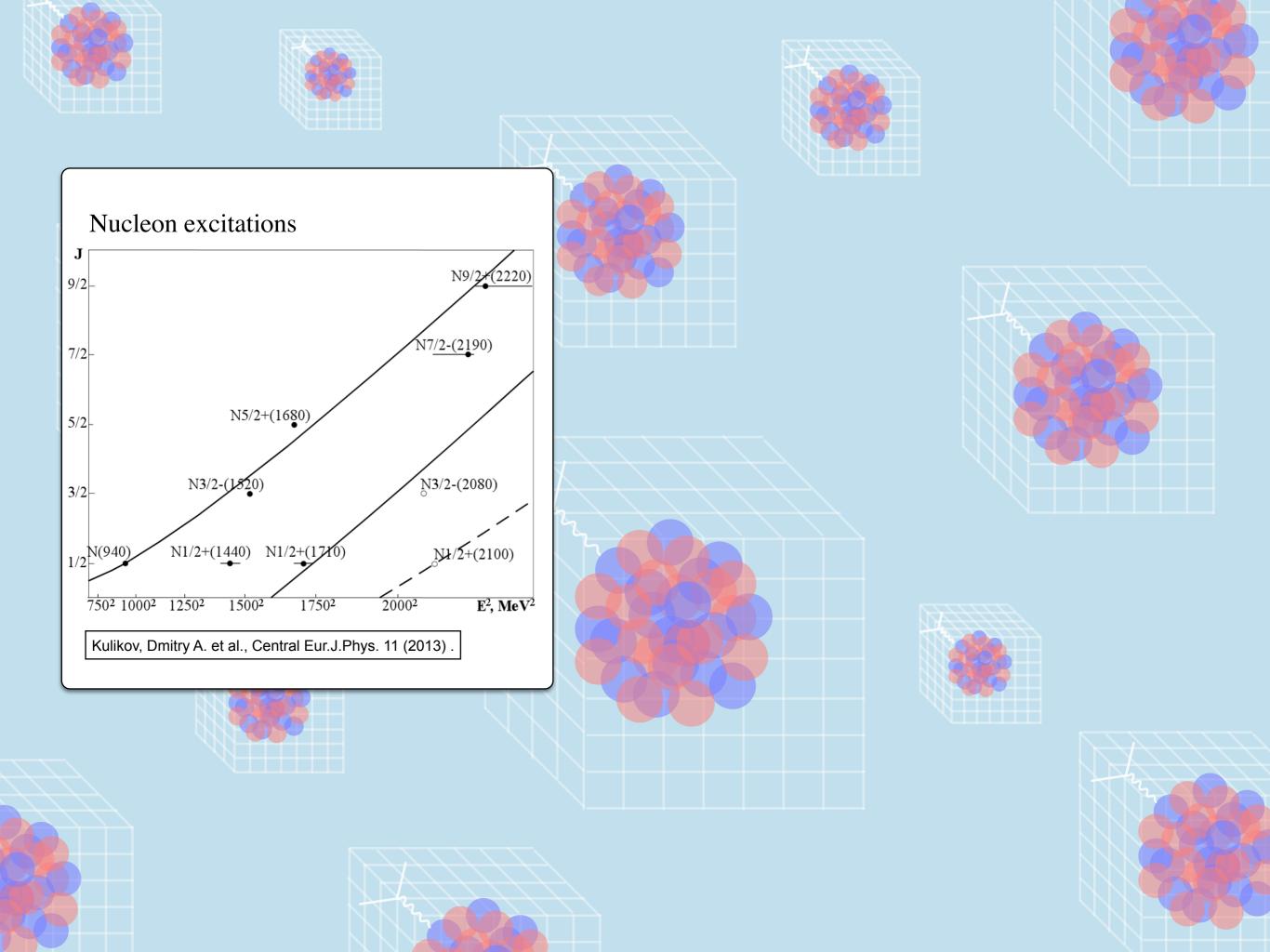


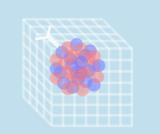


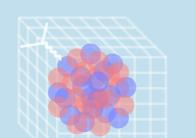


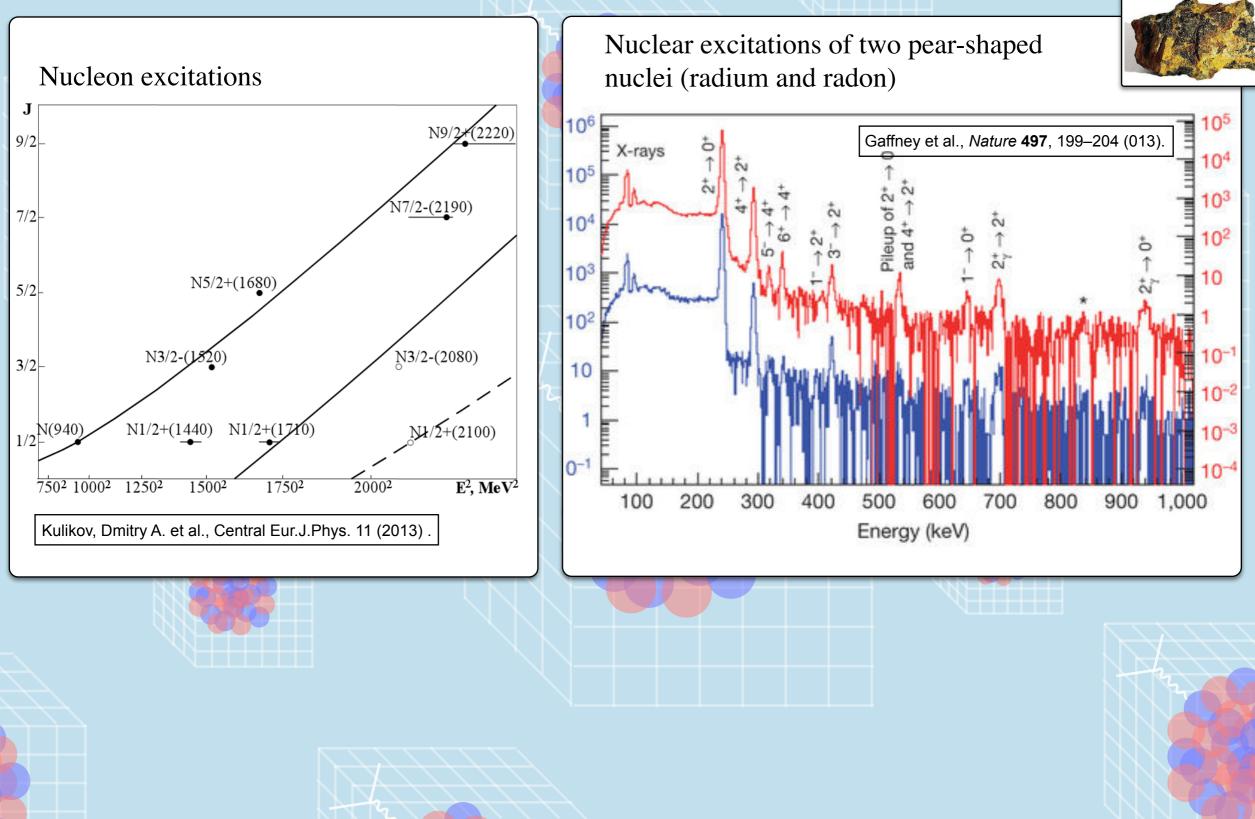


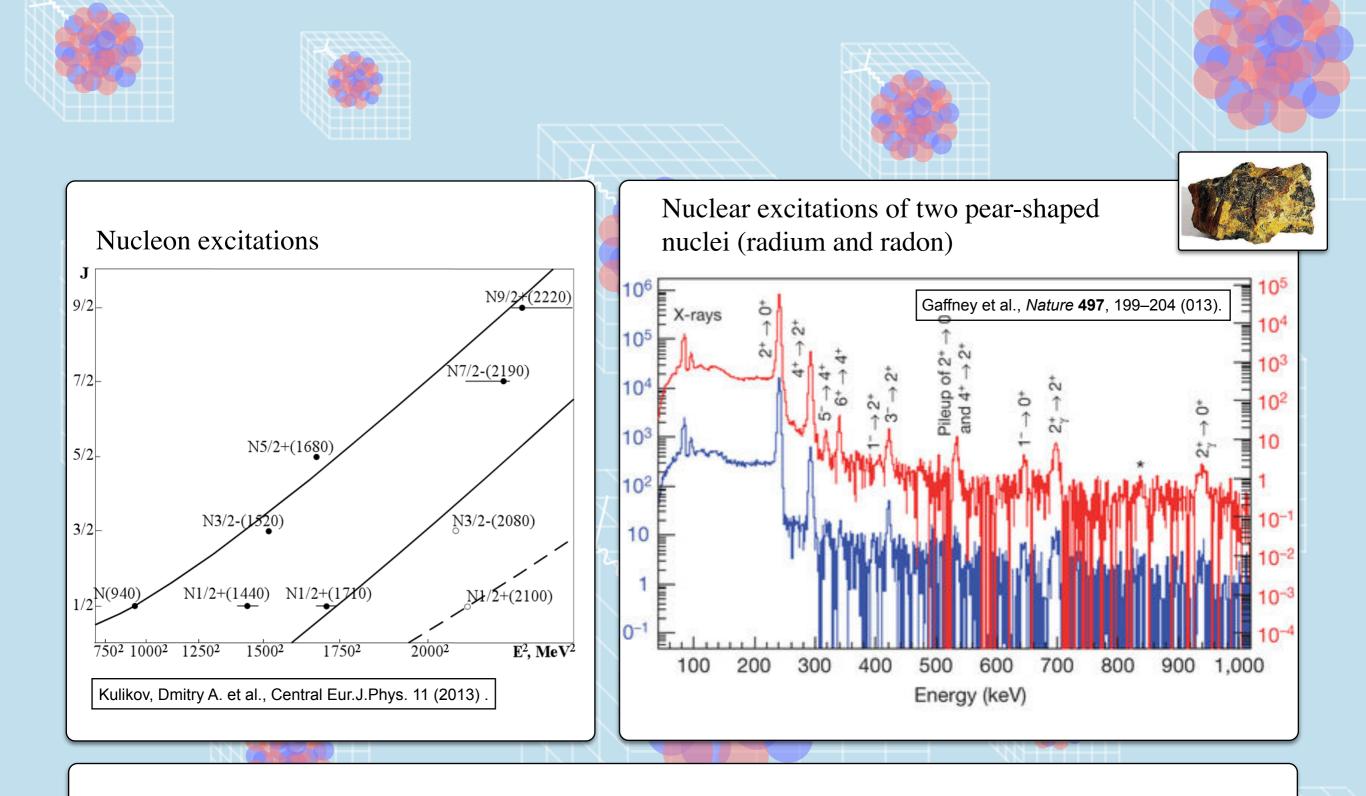




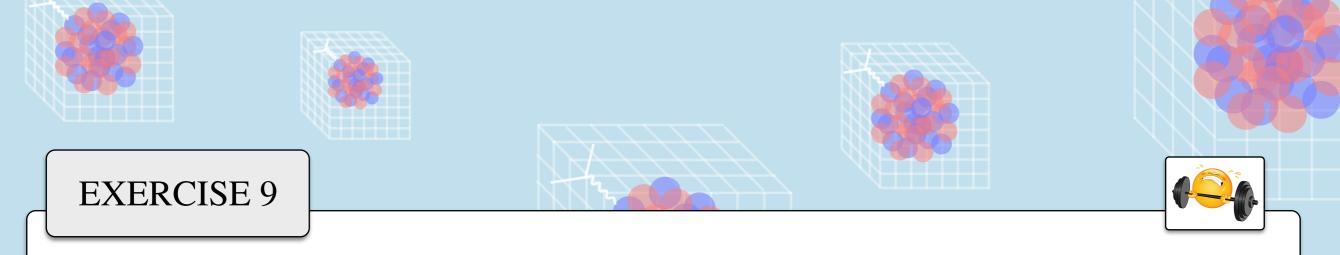




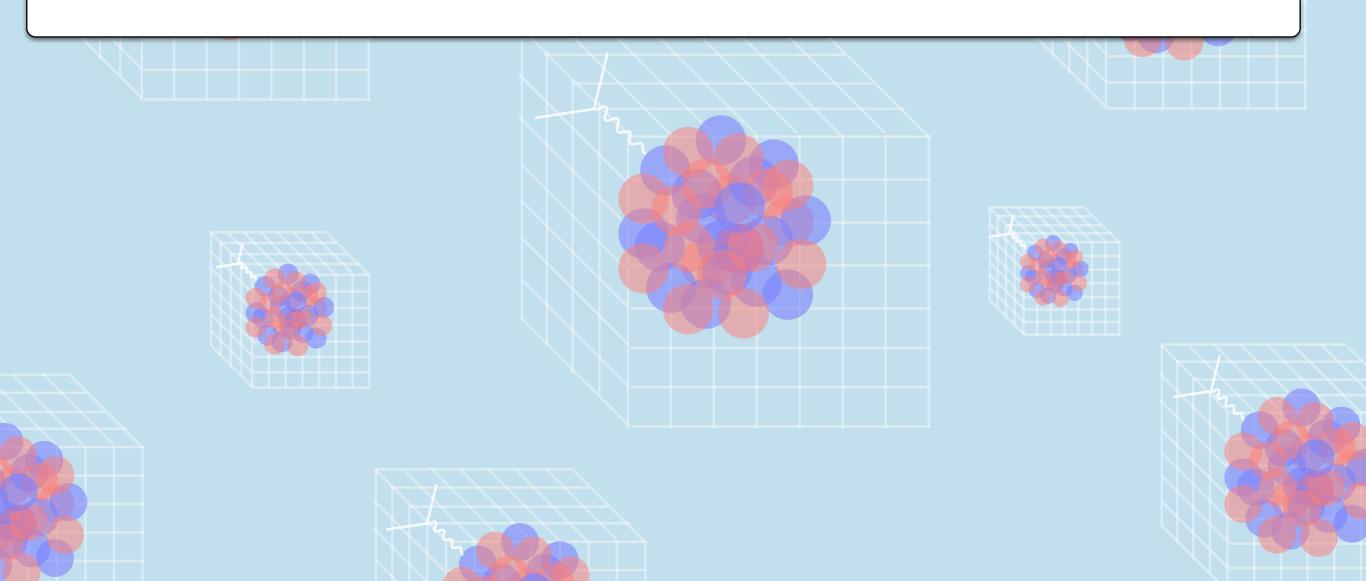


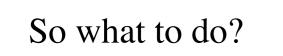


Getting radium directly from QCD will remain challenging for a long time! One should first compute A = 2, 3, 4 systems well. This is till not that easy:  $B_d = 2$  MeV!



With a given amount of computational resources, you have achieved a 1% statistical uncertainty on the extracted mass of the nucleon from your lattice QCD calculation. By what factor should you increase your computing resources (your statistics) to also achieve a 1% statistical uncertainty on the binding energy of the deuteron?



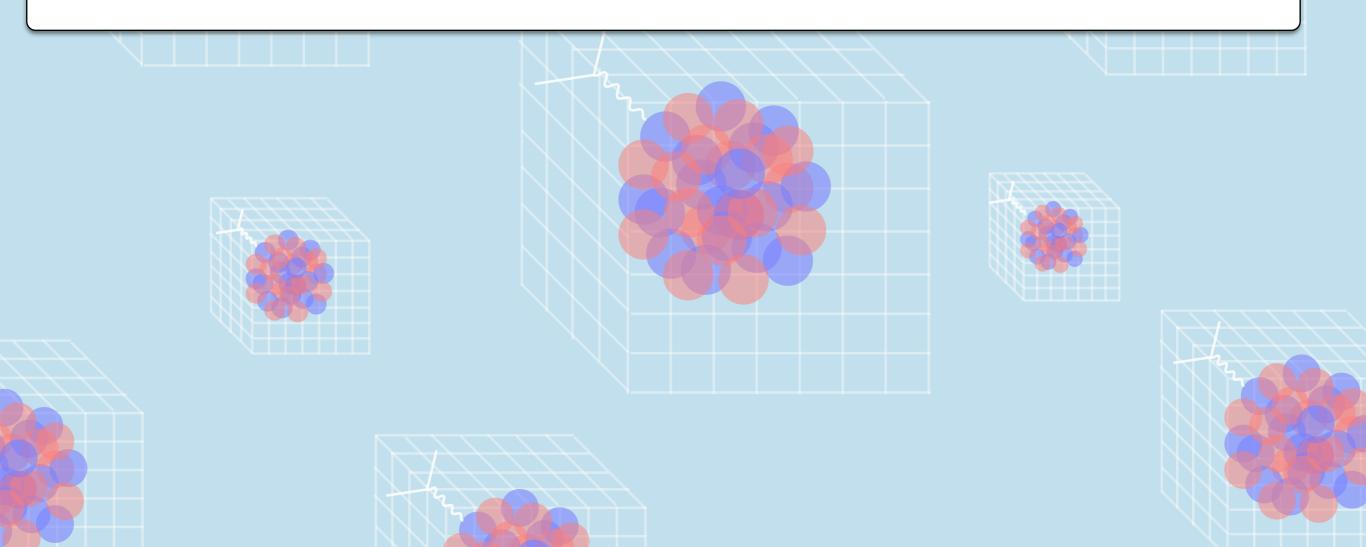


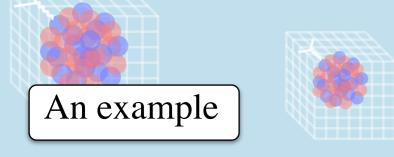
- With the most naive operators with similar overlaps to all states, unreasonably large times are needed to resolve nuclear energy gaps.
- The key to success of this program is in the use of good interpolating operators for nuclei. Since nucleons retain their identity in nuclei, forming baryon blocks at the sink turns out to be very advantageous.
- Ideally need to use a large set of operators for a variational analysis, but this has remained too costly in nuclear calculations. Applications in mesonic sector: Briceno, Dudek and Young, Rev. Mod. Phys. 90 025001.
  - Methods such as matrix Prony that eliminate the excited states in linear combinations of interpolators or correlations functions have shown to be useful.

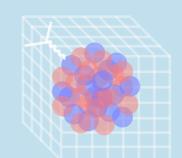
A good review: Beane, Detmold, Orginos, Savage, Prog. Part. Nucl. Phys. 66 (2011).



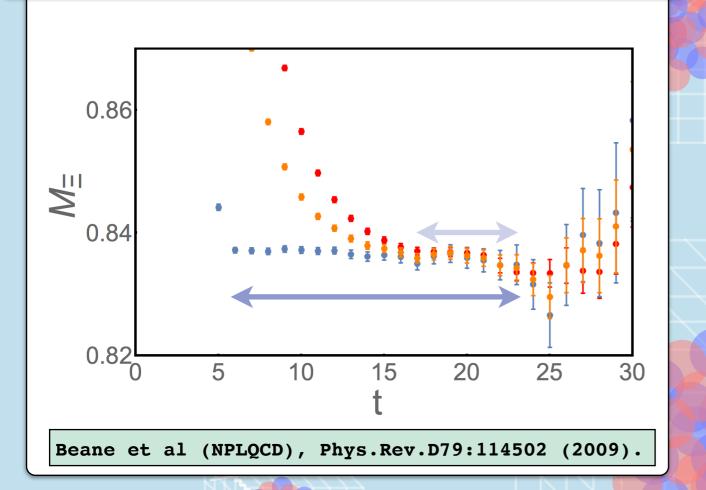
Consider a simple two-state model in the spectral decomposition of an Euclidean two-point function. Demonstrate that the time scale to reach the ground state of the model with a finite statistical precision can depend highly on the corresponding overlap factor for the state. It is sufficient to show this numerically and for a set of chosen energies and overlap factors.

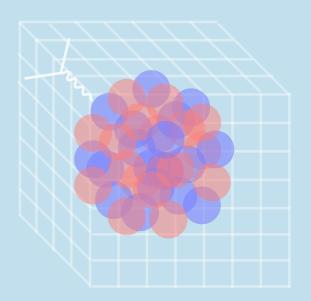


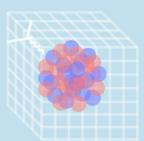


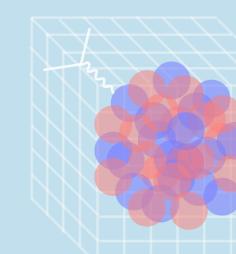


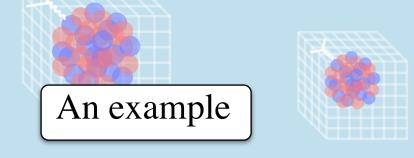
Linear combos. at the level of correlation functions

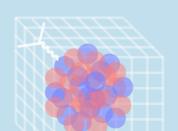




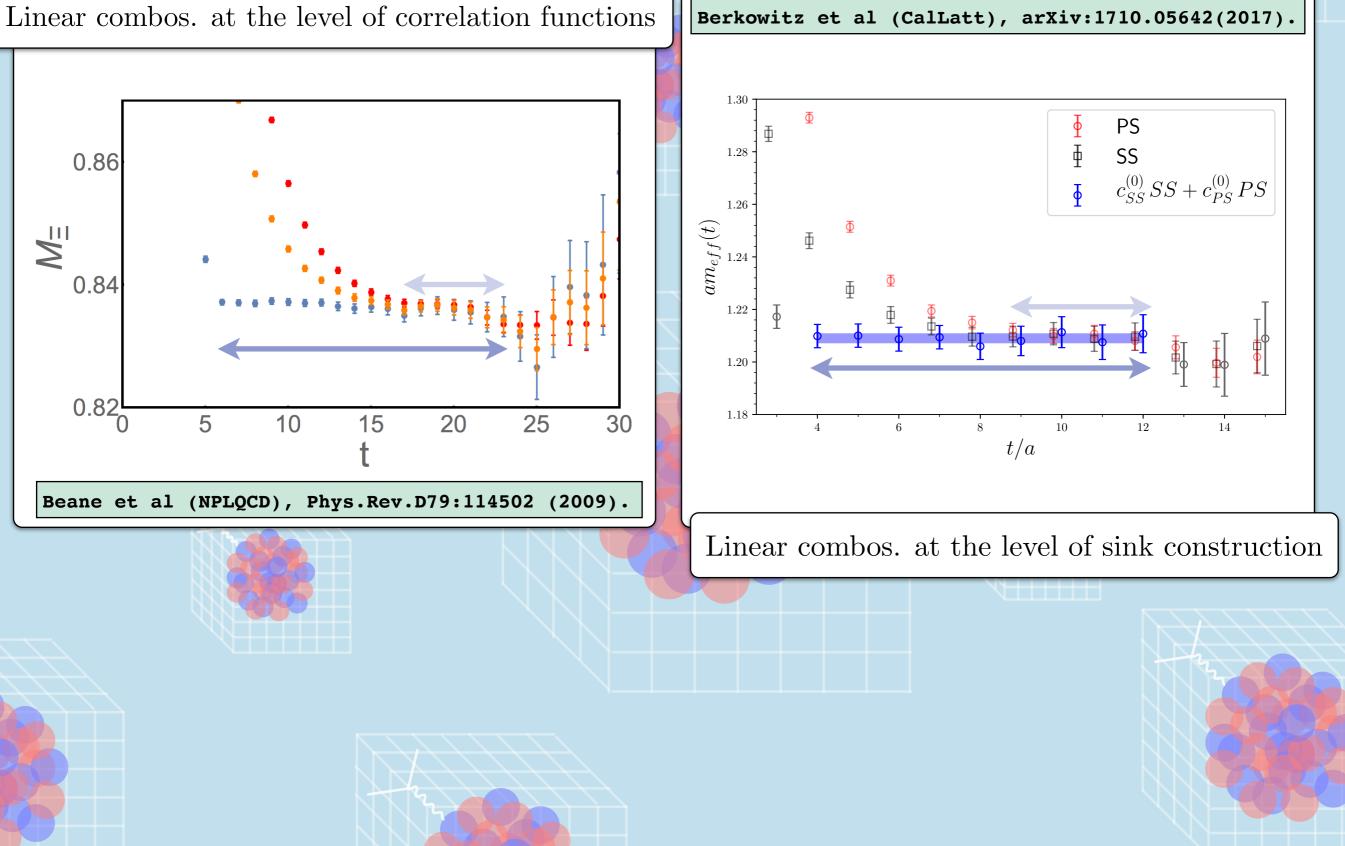




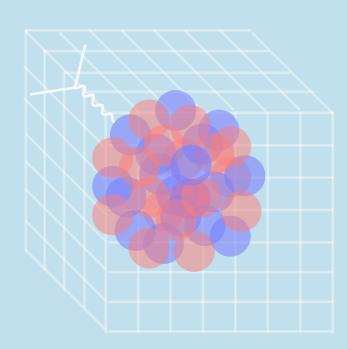


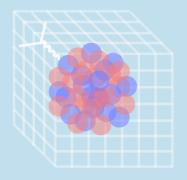


Berkowitz et al (CalLatt), arXiv:1710.05642(2017).

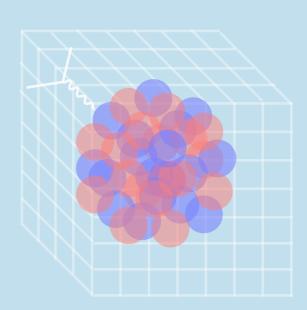


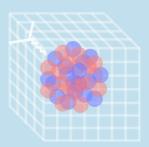
## iii) There is a severe signal-to-noise degradation.

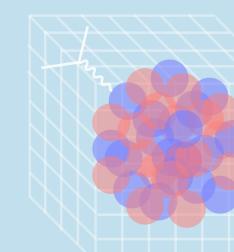


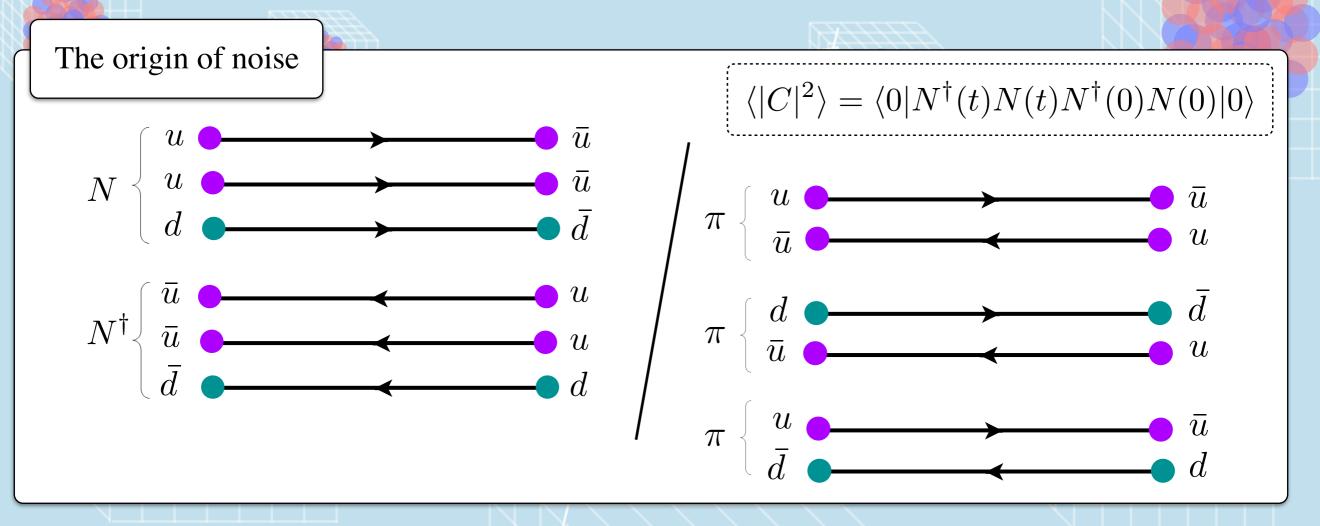


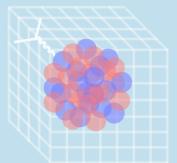


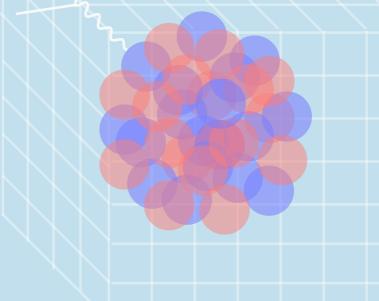




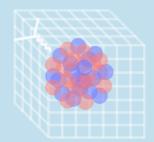


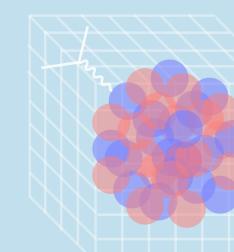


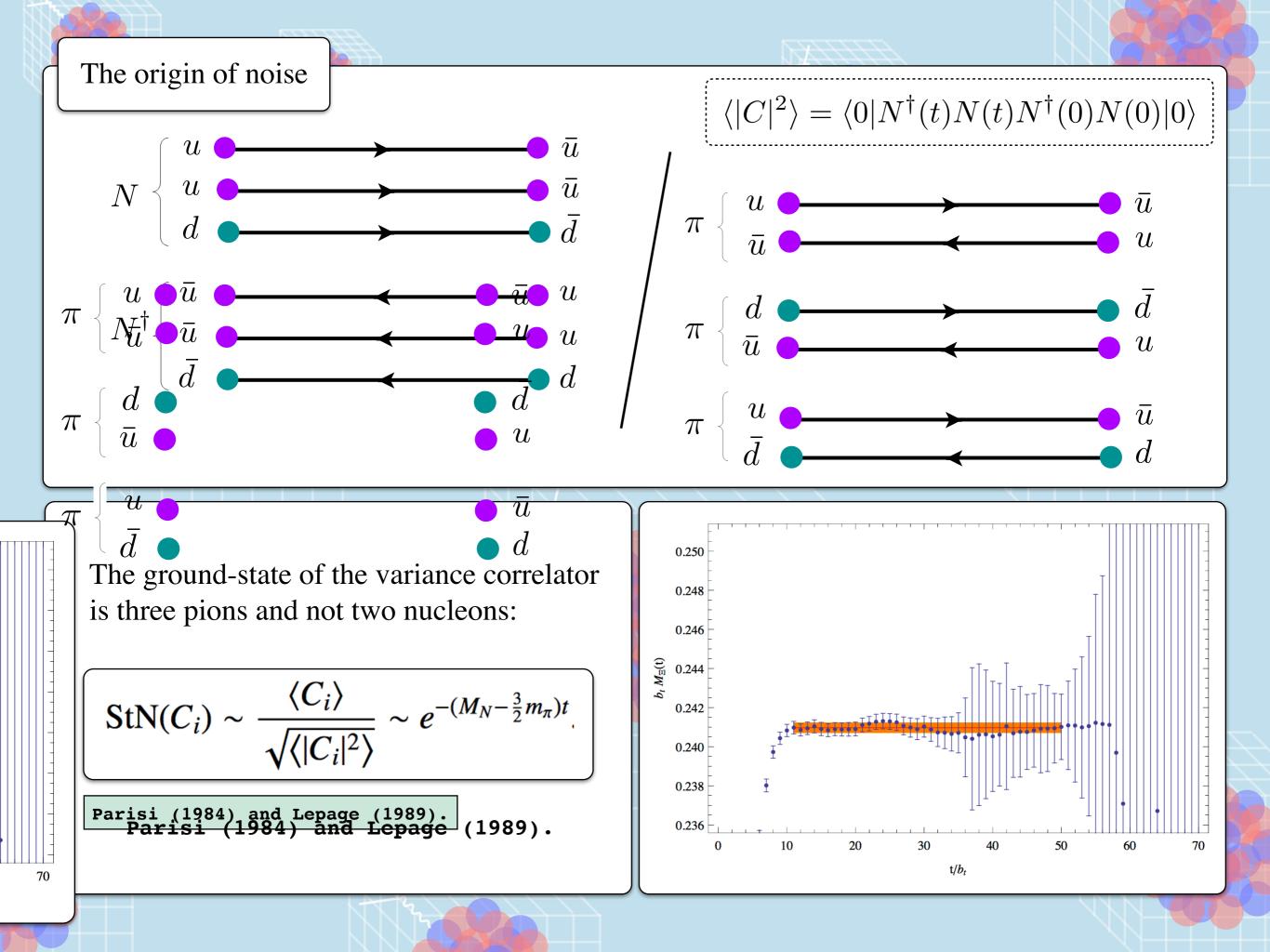


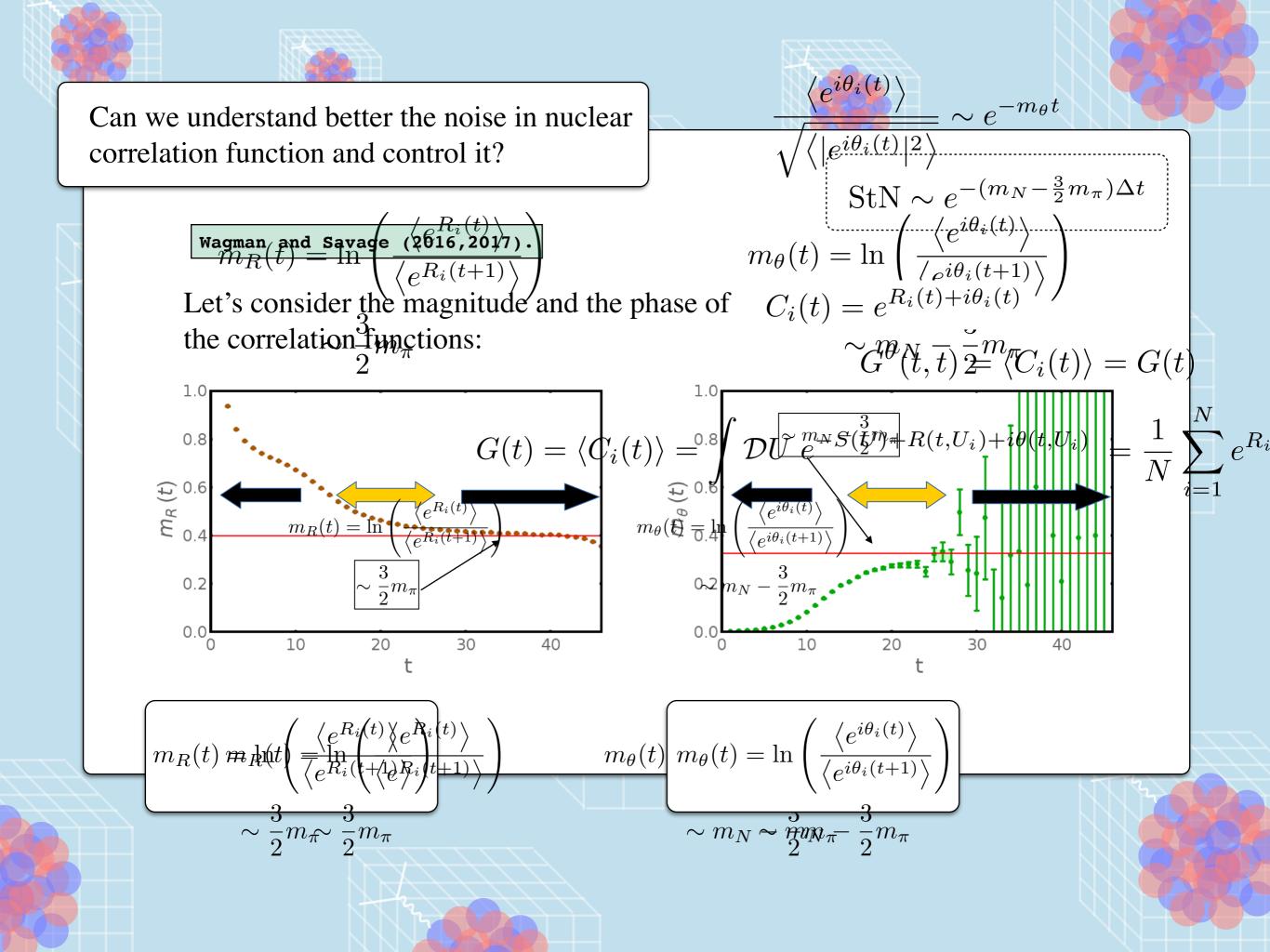


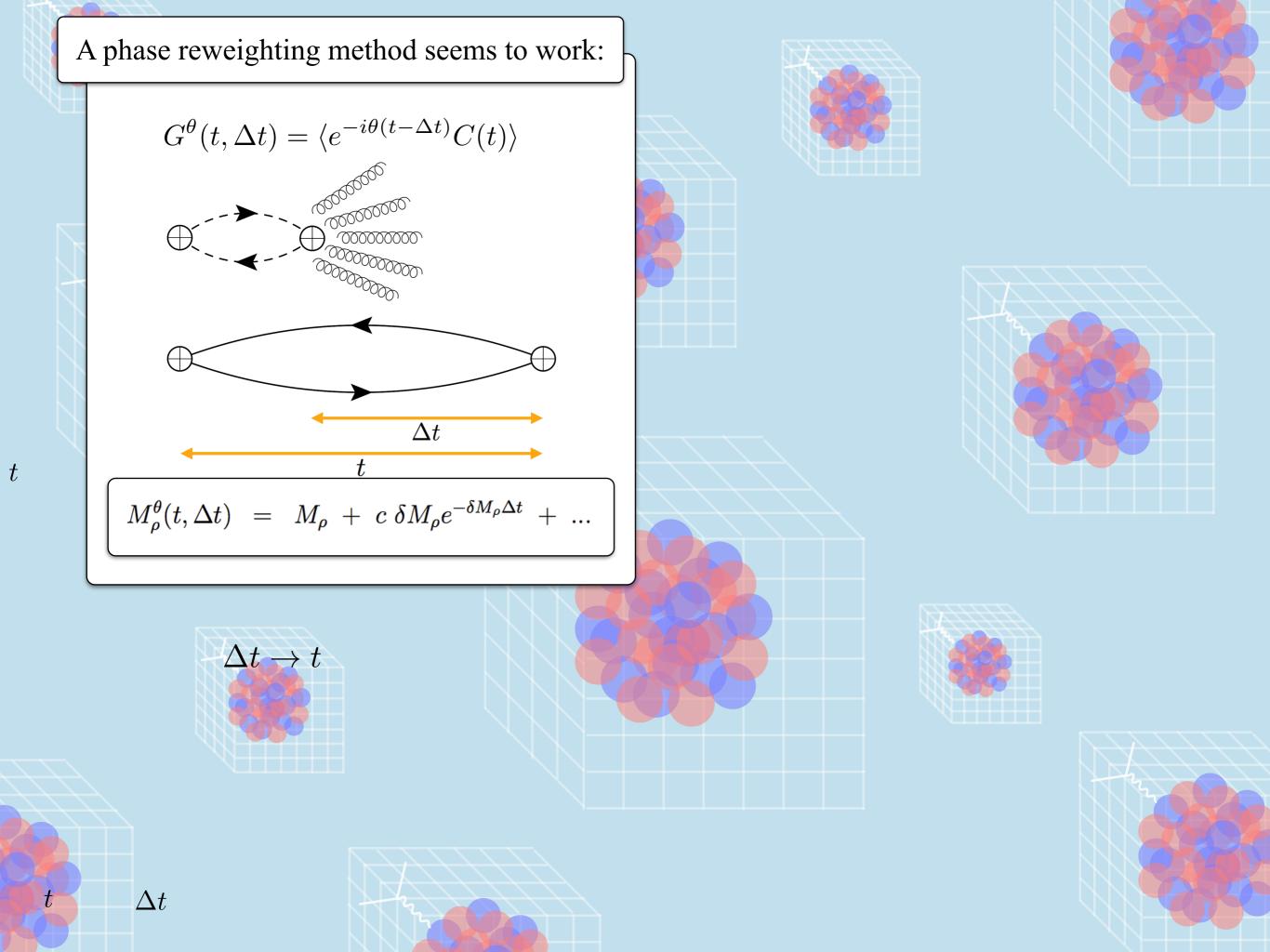


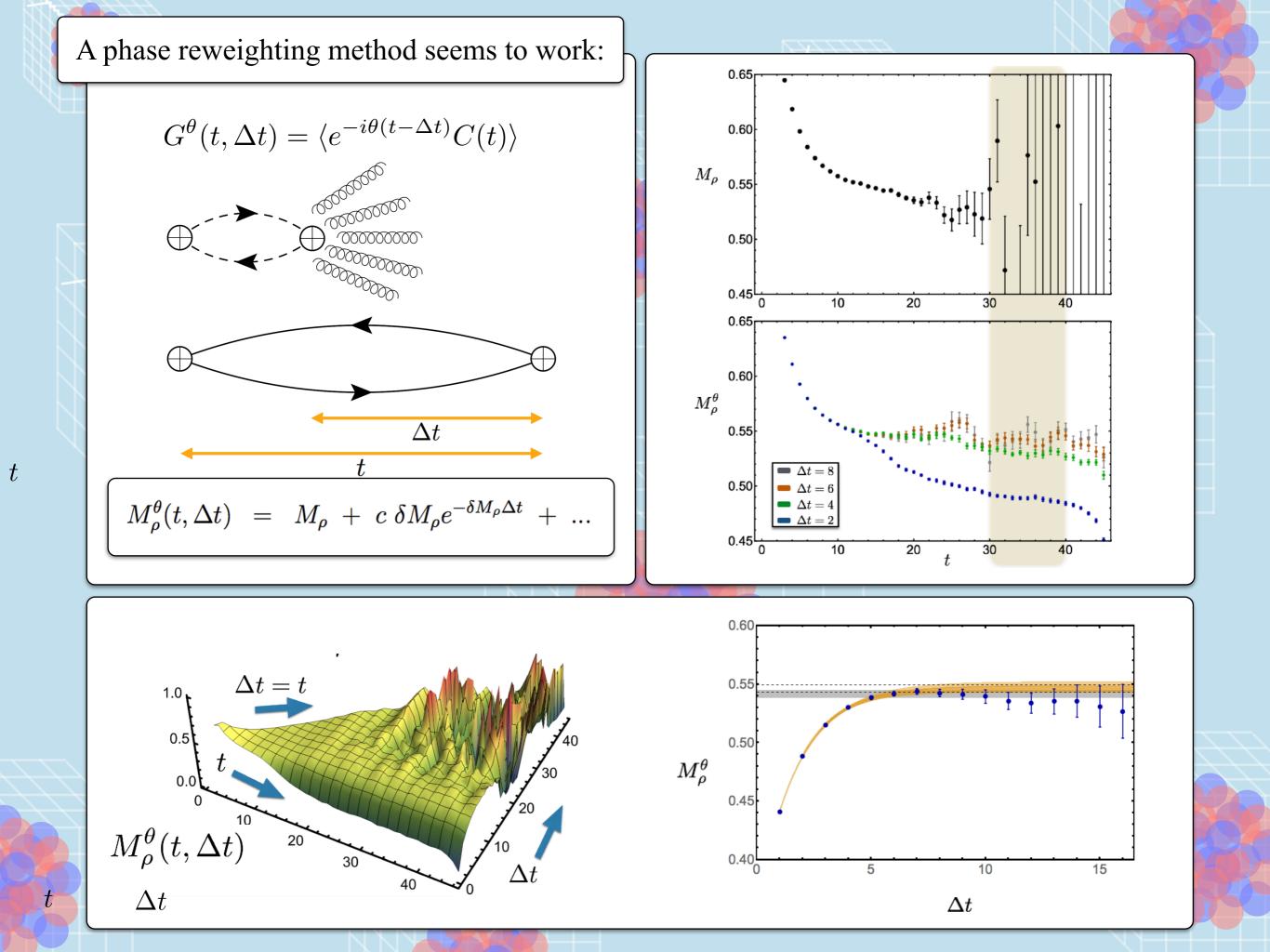












DESPITE CHALLENGES, GREAT PROGRESS HAS BEEN MADE. LQCD IS ON TRACK TO DELIVER RESULTS ON IMPORTANT QUANTITIES FOR THE EIC PHYSICS.

