

# Field selection——a method to save the redundant cost

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# Motivation

Nowadays, lattice QCD simulation became more and more **expensive**

- Nuclear physics  
contraction costs grow exponentially in number of quarks
- Complex operator  
such as four point function, more volume average( $\propto V^3$ )
- Higher momentum  
due to sign-to-noisy problem, the statistics are increase exponentially
- Higher precision  
error reduce by factor  $1/\sqrt{N}$   
different ensembles to estimate the system error

Meanwhile, there are some **redundant cost** at our simulation process

- All mode average(AMA), truncated solver methods with a bias correction

$$C = \frac{1}{N_{LP}} \sum_{i=1}^{N_{LP}} C_i^{LP} + \frac{1}{N_{HP}} \sum_{i=1}^{N_{HP}} (C_i^{HP} - C_i^{LP})$$

T. Blum, T. Izubuchi, and E. Shintani, Phys. Rev. D88 (2013), no. 9 094503

$C^{LP}$  ( $C^{HP}$ ) are the low(high) precision result  $10^{-4}$  ( $10^{-8}$ ),  $N_{LP} \ll N_{HP}$

- **Field selection** ... ..

# Field selection

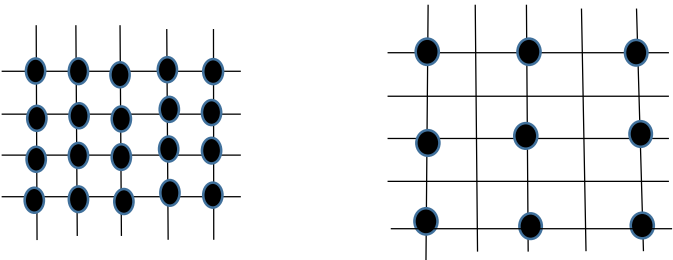
n point function

$$C^n = \sum_{x_1 \in V} \sum_{x_2 \in V} \dots \sum_{x_{n-1} \in V} O_1(t_1, x_1) O_2 \dots O_{n-1} O_n(t_n, x_n) \propto V^{n-1}$$

**Main idea:**  
 Highly correlation between different position.  
 Replace volume summation with subspace summation

## 1. sparse grid field selection

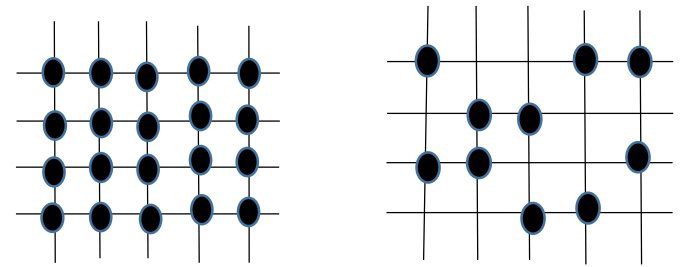
W. Detmold, D.J.Murphy, A.V.Pochinsky et al.  
 arXiv:1908.07050



$$\sum_{x \in V}$$

$$\sum_{x \in \Lambda_{grid}}$$

## 2. random field selection



$$\sum_{x \in V}$$

$$\sum_{x \in \Lambda_{random}}$$

# Two point function

- full correlation function

$$C_{full}^2 = \langle 0 | \sum_{x \in V} O(x, t) O^\dagger(x_0, t_0) | 0 \rangle$$

$$V = \{(n_1, n_2, n_3) | 0 < n_i < L\}$$

- random field selection

$$C_r^2 = \langle 0 | \sum_{x \in \Lambda_r} O(x, t) O^\dagger(x_0, t_0) | 0 \rangle$$

$$\Lambda_r = \{(n_1, n_2, n_3) | n_i \text{ randomly}\}$$

- grid field selection

$$C_g^2 = \langle 0 | \sum_{x \in \Lambda_g} O(x, t) O^\dagger(x_0, t_0) | 0 \rangle$$

$$\Lambda_g = \{(n_1, n_2, n_3) | n_i = 0 \text{ mod } N\} \quad (L/N)^3$$

$\Lambda_r$  is different at different t

For noncube number

source position must be in  $\Lambda_g$

$$\Lambda_g = \{(n_1, n_2, n_3) | n_i = 0 \text{ mod } N \ \&\& \ n_1 + n_2 + n_3 = 0 \text{ mod } 2N\} \quad (L/N)^3/2$$

$$\Lambda_g = \{(n_1, n_2, n_3) | n_i = 0 \text{ mod } N \ \&\& \ n_i + n_j = 0 \text{ mod } 2N\} \quad (L/N)^3/4$$

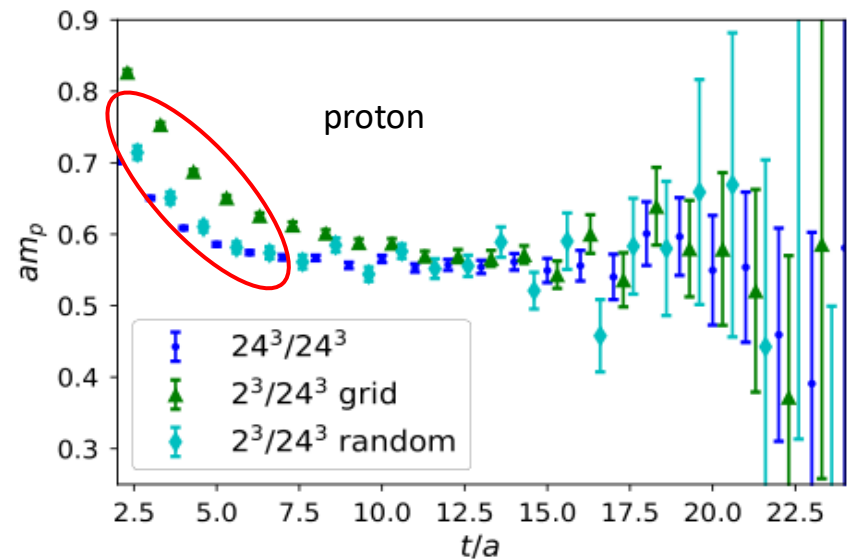
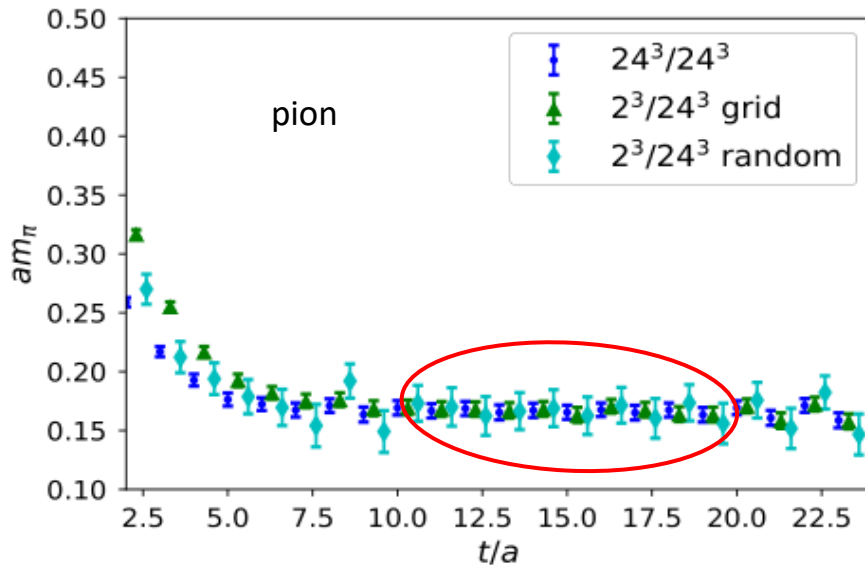
- numbers tested are  $N_{th} = \{24^3, 12^3, \frac{12^3}{2}, \frac{12^3}{4}, 6^3, \frac{6^3}{2}, 4^3, \frac{6^3}{4}, \frac{4^3}{2}, 3^3, \frac{4^3}{4}, 2^3, 4, 2, 1\}_{4/10}$

**Ensemble information** :  $24^3 \times 48$ ,  
 $N_f = 2 + 1 + 1$   $m_\pi \approx 350 \text{ MeV}$   
 clover-improved Wilson twist mass  
 ensemble, smearing propagators  
 are used

# Effective mass

The effective mass are obtained by equation

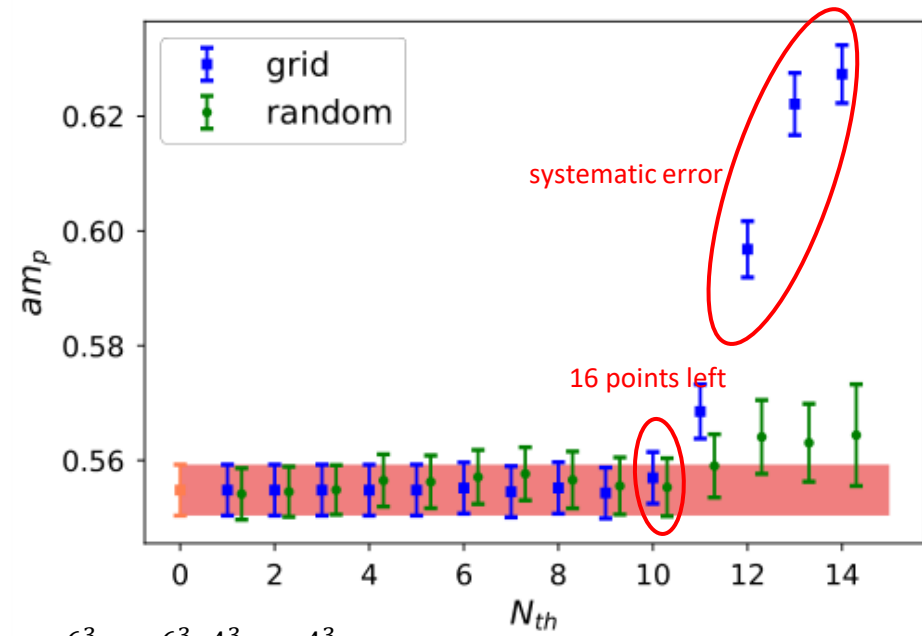
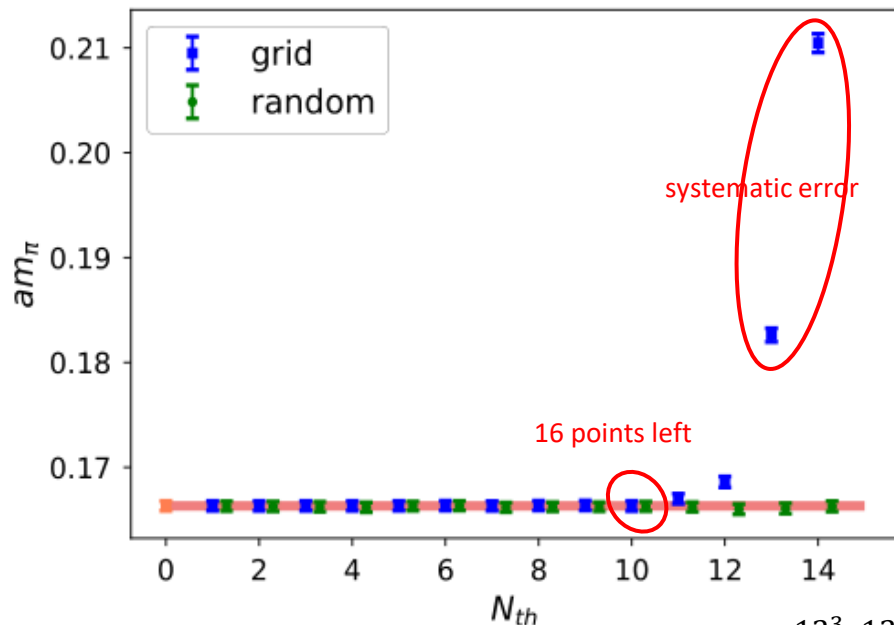
$$aE(t) = \begin{cases} \cosh^{-1} \left[ \frac{C(t-1) + C(t+1)}{2C(t)} \right], & \text{mesons} \\ \log \left[ \frac{C(t-1)}{C(t)} \right], & \text{baryons} \end{cases}$$



- at large  $t$ , central values from different method are **consistent**
- random field selection have **larger error** than grid
- at short  $t$ , grid field selection has **systematic error**

# Mass after correlated fit

- fitting window is decided by full correlation function
- fitting window is [8,22] for pion, [10,22] for proton



$$N_{th} = \{24^3, 12^3, \frac{12^3}{2}, \frac{12^3}{4}, 6^3, \frac{6^3}{2}, 4^3, \frac{6^3}{4}, \frac{4^3}{2}, 3^3, \frac{4^3}{4}, 2^3, 4, 2, 1\}$$

- 16 points can reproduce the full correlation function result
- Larger error in random field selection at effective mass plot disappear after fit
  - Effective mass plot only reflect local information
  - less correlated between different time slices
- Random field selection is used in remainder of talk

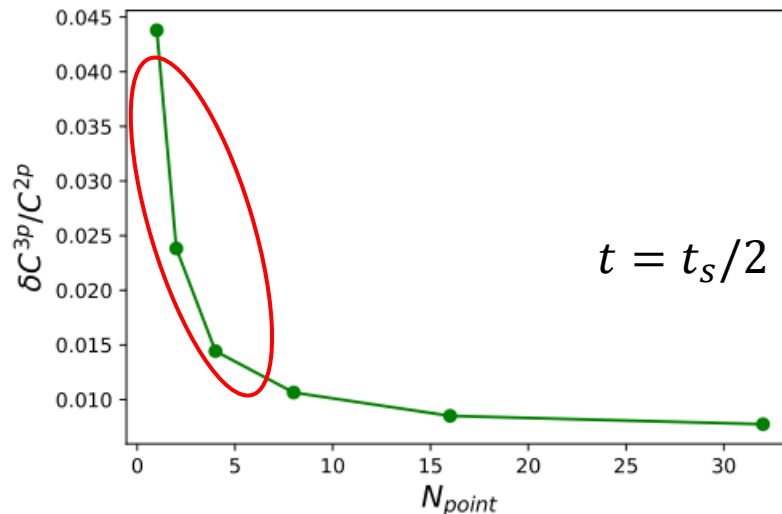
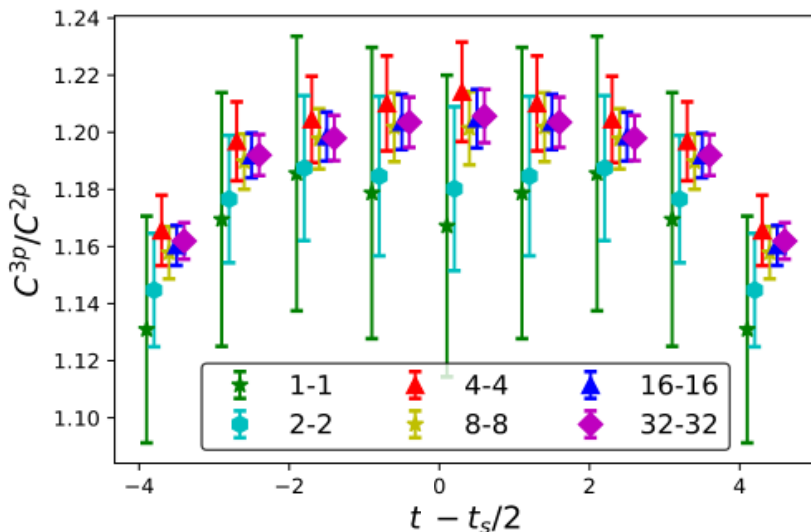
# Proton axial charge

- Construct correlated function like building block

$$C^{3pt} = \sum_{x \in V} \sum_{x_f \in \Lambda'} \sum_{x_i \in \Lambda} \langle O(t_f, x_f) A^\mu(t, x) O^\dagger(t_i, x_i) \rangle$$

$$C^{2pt} = \sum_{x_f \in \Lambda'} \sum_{x_i \in \Lambda} \langle O(t_f, x_f) O^\dagger(t_i, x_i) \rangle$$

- 32 propagators on each two time slice for 91 configuration
- test the number of point at source and sink



- error reduces fast at small number
- 32 points used at source and sink

# Fit method

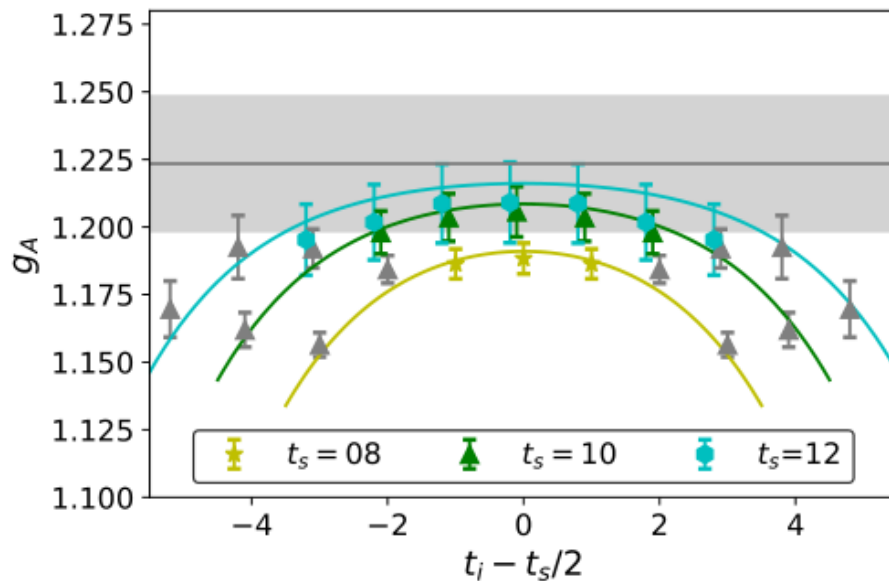
- consider first excited state
- two point function

$$C^{2pt} = |B_0|^2 e^{-E_0 t_s} + |B_1|^2 e^{-E_1 t_s}$$

- mass and amplitudes  $B_0, B_1, E_0, E_1$  are fitted from two point function
- Three point function

$$C^{3pt} = |B_0|^2 \langle 0|A|0 \rangle e^{-E_0 t_s} + |B_1|^2 \langle 1|A|1 \rangle e^{-E_1 t_s} \\ + B_0 B_1^* \langle 0|A|1 \rangle e^{-E_1 t_i} e^{-E_0(t_s - t_i)} + B_0^* B_1 \langle 1|A|0 \rangle e^{-E_0 t_i} e^{-E_1(t_s - t_i)}$$

- The three matrix elements are obtained from three point function by fitting multiple value of  $t$  and  $t_s$



$$g_A = 1.223(25)$$

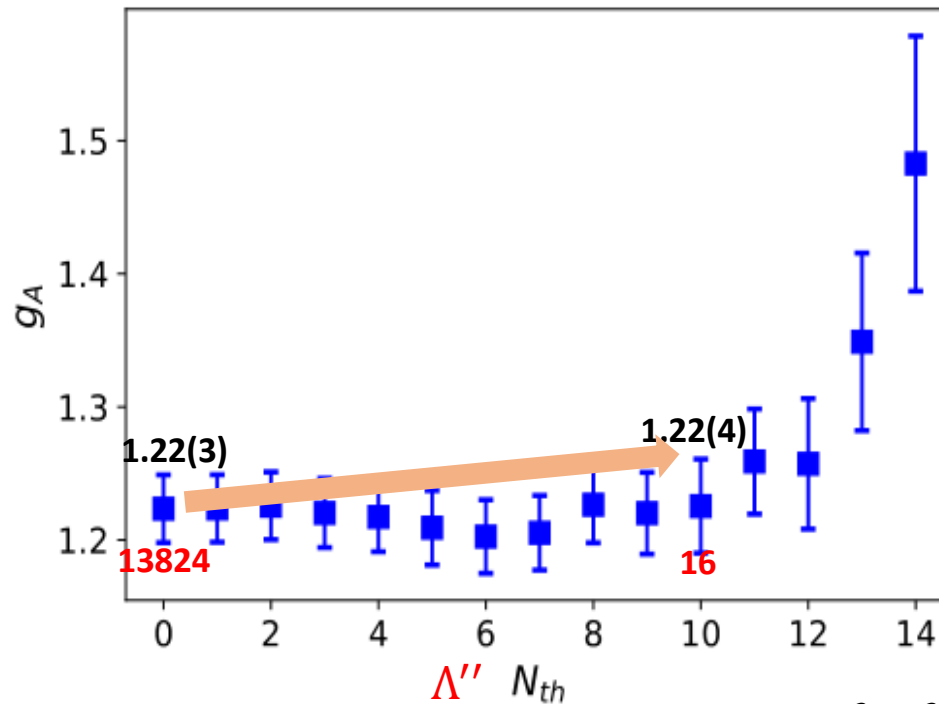


# proton axial charge

- Test the number of point in  $\Lambda''$

$$C^{3pt} = \sum_{x \in \Lambda''} \sum_{x_f \in \Lambda'} \sum_{x_i \in \Lambda} \langle O(t_f, x_f) A^\mu(t, x) O^\dagger(t_i, x_i) \rangle$$

- the number of point in  $\Lambda'$  and  $\Lambda$  are 32



- 16 points in  $\Lambda''$  reproduce the full lattice

$$N_{th} = \{24^3, 12^3, \frac{12^3}{2}, \frac{12^3}{4}, 6^3, \frac{6^3}{2}, 4^3, \frac{6^3}{4}, \frac{4^3}{2}, 3^3, \frac{4^3}{4}, 2^3, 4, 2, 1\}$$

# conclusions

- We have explored grid and random field selection method at two point function and three point function
- field selection almost not lose any precision by reducing the summation by a factor  $O(1000)$ 
  - The **contraction** are speeded up highly
  - The **storage of propagators** is reduced significantly
- With field selection , much **redundant cost** will be saved
- Paper to appear on arXiv soon...

**Thanks for listening**