Moments of structure functions for $N_f = 2$ near the physical point

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We report on our on-going study of the lower moments of iso-vector polarised and unpolarised structure functions, $g_A$ and $\langle x \rangle_{u-d}$, respectively, and the iso-vector scalar and tensor charge, for $N_f = 2$ non-perturbatively improved clover fermions. With pion masses which go down to about 150 MeV, we investigate finite volume effects and excited state contributions.

31st International Symposium on Lattice Field Theory - LATTICE 2013
July 29 - August 3, 2013
Mainz, Germany

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sink

Figure 1: The connected (left) and disconnected (right) contributions to nucleon three-point functions for a nucleon source at \( t = 0 \), sink at \( t_{\text{sink}} \) and operator insertion at \( t_{\text{ins}} \).

1. Introduction

In recent years it has become clear that \( g_A \) and \( \langle x \rangle_{u-d} \), benchmark quantities for lattice calculations of nucleon structure, are sensitive to a number of sources of systematic error - finite volume, non-physical pion mass, excited state contamination and finite lattice spacing. Continuing improvements in computing power, algorithms and analysis techniques mean systematics can now be investigated and in some cases removed.

In the following we present preliminary results for these quantities along with the iso-vector scalar and tensor charges, \( g_S \) and \( g_T \), respectively, for \( N_f = 2 \) ensembles including different lattice spacings, pion masses and volumes focusing on studying excited state effects. The iso-vector combination only requires the calculation of the connected quark diagram, see Figure 1. We have also determined the disconnected contributions to the scalar matrix element. This and the iso-vector generalised form factors for a range of operators are detailed in [1]. While many lattice simulations now include dynamical strange quarks, so far the strange quark has been found to play a minor role in nucleon structure and \( N_f = 2 \) simulations are still relevant.

2. Simulation details

The results were computed using ensembles generated by QCDSF and the Regensburg Group, with \( N_f = 2 \) degenerate flavours of dynamical sea quarks using the non-perturbatively improved clover action at two lattice spacings and a range of pion masses from \( m_\pi \sim 490 - 150 \) MeV, see Table 1. Two volumes are available for two values of \( m_\pi \), in particular at the near physical point.

The two-point and three-point functions were computed using Wuppertal smeared sources and sinks with APE smeared gauge links. For each ensemble the smearing was optimised to minimize the excited state contributions to the nucleon two-point function. The connected three-point functions were generated using the standard sequential propagator method which involves fixing the sink timeslice \( (t_{\text{sink}}) \). An alternative approach using stochastic estimates has been investigated, see [2]. The values of \( t_{\text{sink}} \) were chosen using ensemble II. For \( t_{\text{sink}} = 15a \sim 1 \) fm on this ensemble no significant excited state contributions to the quantities \( \langle x \rangle_{u-d} \) and \( g_A \) were found. This was checked by performing an excited state analysis with multiple \( t_{\text{sink}} \)s, described in the next section. \( t_{\text{sink}} = 15a \) was then used for all \( \beta = 5.29 \) ensembles and rescaled for \( \beta = 5.40 \). Multiple measure-
ments were performed on each configuration. Autocorrelations were investigated by binning the data with different bin sizes.

In order to reduce the discretisation effects to \( O(a^2) \), the operators, as well as the quark action, need to be (nonperturbatively) improved. The \( O(a) \) improved renormalised operators have the form [3]

\[
O^{\text{improv}} = Z_O \left[ (1 + b_O m_q) O + ac_O O' \right].
\]  

(2.1)

For \( g_A \), \( O = \bar{q} \gamma_{\mu} s q \), and the improvement term, \( ac_O O' = ac_A \partial_{\mu} \bar{q} \gamma_{5} q \), does not contribute for forward matrix elements. We use the \( Z_O \) factors determined non-perturbatively [4, 5] and the \( b_O \) factors from [6]. Setting \( c_O = 0 \), our values for \( g_A \) will have leading \( O(a^2) \) effects, while \( \langle x \rangle_{ ud } \), \( g_S \) and \( g_T \) have \( O(a) \).

3. Excited state fits

Excited state contributions to nucleon structure have been investigated by a number of groups recently, see, for example, [7, 8, 9, 10, 11]. Our analysis on ensemble II is similar to that performed in reference [11]. We fitted the two-point \((C_{2pt})\) and three-point \((C_{3pt})\) functions at multiple \( t_{\text{sink}} \)s simultaneously using the functional forms:

\[
C_{2pt}(t_{\text{sink}}) = \sum_{\beta} \langle \mathcal{N}(\vec{x}, t_{\text{sink}}) \mathcal{F}(\vec{0}, 0) \rangle = |Z_0|^2 e^{-m_0 t_{\text{sink}}} + |Z_1|^2 e^{-m_1 t_{\text{sink}}} + \ldots
\]  

(3.1)

\[
C_{3pt}(t_{\text{sink}}, t_{\text{ins}}) = \sum_{\beta} \langle \mathcal{N}(\vec{x}, t_{\text{sink}}) O(\vec{y}, t_{\text{ins}}) \mathcal{F}(\vec{0}, 0) \rangle
\]  

(3.2)

\[
= |Z_0|^2 \langle \mathcal{N}_0 O(\mathcal{N}_0) e^{-m_0 t_{\text{sink}}} + Z_1 \mathcal{Z}_0 \mathcal{N}_1 O(\mathcal{N}_0) e^{-m_1 t_{\text{sink}} - t_{\text{ins}}} + \ldots
\]  

(3.3)

\[
+ Z_0^2 \mathcal{Z}_1 \mathcal{N}_0 O(\mathcal{N}_1) e^{-m_0 t_{\text{sink}} - t_{\text{ins}}} + |Z_1|^2 \langle \mathcal{N}_1 O(\mathcal{N}_1) e^{-m_1 t_{\text{sink}} - t_{\text{ins}}} + \ldots
\]  

(3.4)

\[
= |Z_0|^2 e^{-m_0 t_{\text{sink}}} \left( B_0 + B_1 [e^{-\Delta m t_{\text{sink}} - t_{\text{ins}}} + e^{-\Delta m t_{\text{ins}}} + B_2 e^{-\Delta m t_{\text{sink}}} \right) + \ldots
\]  

(3.5)
where $\mathcal{N}$ destroys the nucleon, $O$ is the current insertion and $Z_i = \langle 0 | \mathcal{N} | N_i \rangle$. $|N_0\rangle$ and $|N_1\rangle$ represent the nucleon ground and first excited state, respectively. For the cases considered here, it is sufficient to consider zero initial and final momentum in order to extract forward matrix elements.

The matrix element of interest is given by $B_0 = \langle N_0 | O | N_0 \rangle$, while $B_1 \propto \langle N_1 | O | N_0 \rangle$ gives the transition matrix element from the ground to the first excited state and $B_2 \propto \langle N_1 | O | N_1 \rangle$ the first excited state matrix element. These fits can be compared to the traditional approach of fitting the ratio of three-point to two-point functions to a constant:

$$\frac{C_{3pt}(t_{sink}, t_{ins})}{C_{2pt}(t_{sink})} = B_0 + \ldots$$  \hfill (3.6)

We illustrate the results for the example of $\langle x \rangle_{u-d}$, which we found to have significant excited state contributions at small $t_{sink}$. Figure 2 displays the raw results for the six $t_{sink}$ values. Agreement is found for $t_{sink} \geq 11a$ for $t_{ins} - t_{sink}/2 \approx 0$. For $t_{sink} \geq 13a$, the plateau extends for several timeslices, however, as expected, the statistical errors also increase. Correlated fits to all three-point and two-point functions using Eqs. (3.2) and (3.5) (truncating after the first excited state) produced good $\chi^2/d.o.f \approx 1$ and were stable against changes in the fitting ranges to $C_{3pt}$ and $C_{2pt}$. An example of one of the combined fits is shown in Figure 2. For the smearing we have used, the ground and first excited state are the dominant contributions for $t_{sink}$ as small as $7a$. In Figure 3, a result of the combined fit is compared to the values obtained fitting the ratio $C_{3pt}/C_{2pt}$ to a constant for different $t_{sink}$; consistency is found between the two methods for $t_{sink} \geq 11a$. From this analysis we find our optimised smearing is sufficient to extract the ground state matrix element using a single $t_{sink} \geq 11a$ and we take the conservative choice of $15a$. We remark that with a less optimised smearing we found a $t_{sink} > 15a$ insufficient.

For the other ensembles, where we only have one $t_{sink}$, we check the size of the excited state contamination by performing fits using Eqs. 3.2 and 3.5 and the parameters proportional to the
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Figure 3: (Left) For ensemble II, $\langle x \rangle_{u-d}^{\text{MS}}$ extracted from constant fits to the ratio $C_{3pt}/C_{2pt}$ (Eq. (3.6)) for different $t_{\text{sink}}$ and extracted from the combined fit to $C_{3pt}$ and $C_{2pt}$, for all $t_{\text{sink}}$, including the first excited state, denoted by $C$. (Right) For ensemble IV, the ratio $C_{3pt}/C_{2pt}$ factor is shown (red triangles). The two lines indicate fits (i) constraining the excited state parameters using the results from ensemble II (blue line) and (ii) leaving $B_1$ free and $B_2 = 0$ (green line). For each of these fits the raw results for $C_{3pt}$ factor, divided by $|Z_0|^2 e^{-m_0 t_{\text{sink}}}$, is shown.

excited state matrix elements, $B_1$ and $B_2$, extracted using ensemble II (which we assume to be slowly varying with the quark mass and lattice spacing). Figure 3 shows such a constrained fit for ensemble IV. The value extracted for the ground state is consistent with a fit using the same functional form, but with $B_1$ as a free parameter and $B_2 = 0$. It is also consistent with the result of a constant fit to $C_{3pt}/C_{2pt}$.

The excited state fitting analysis was similarly successful for the other quantities of interest. We note that for $g_A$, significant excited state contributions were found for $t_{\text{sink}} = 7$ and 9, however, these contributions cancelled in the ratio $C_{3pt}/C_{2pt}$, giving a plateau for the ratio, even for the smallest $t_{\text{sink}}$. For $g_S$, the ground state was dominant for all $t_{\text{sink}}$ for the iso-vector combination in the ratio $C_{3pt}/C_{2pt}$. This appears to be due to a cancellation in the $u-d$ combination since for the iso-scalar matrix elements, not presented in this work, there were significant excited state contributions. The tensor charge analysis was similar to that for $\langle x \rangle_{u-d}$.

4. Results

In the following we present results obtained from a constant fit to the ratio $C_{3pt}/C_{2pt}$, for a single $t_{\text{sink}}$. Figure 4 shows our results for $g_A$, $\langle x \rangle_{u-d}$, $g_T$ and $g_S$ from all ensembles as a function of $m_\pi^2$. Recent work from other groups is indicated, where for $g_A$ and $\langle x \rangle_{u-d}$, due to the large number of previous determinations, we only compare with other $N_f = 2$ calculations. Recent $N_f = 2 + 1$ and $2 + 1 + 1$ results are reviewed in [12]. For the scalar and tensor charge most other results are for $N_f = 2 + 1$ and $2 + 1 + 1$.

\footnote{This parameter cannot be determined from fits to a single $t_{\text{sink}}}
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Figure 4: Results obtained from all ensembles as a function of $m_\pi^2$ (blue circles) compared to previous works. The physical point is indicated by a vertical line. For $g_A$ and $\langle x \rangle_{u-d}$ only $N_f = 2$ determinations are shown. For the former, the experimental result is shown as a black square, while for the latter, the expectations from PDF parameterisations of the NNPDF, ABM and MSTW groups are shown.

Considering $g_A$ first, one can see in Figure 4 that there is a significant dependence of our results on the volume and possibly $m_\pi$ and $a$. Consistency is found with the values from the Mainz Group [9] and also ETMC [13]. The discrepancy with reference [14], which uses some of the same configurations but in some cases different smearing and $t_{\text{sink}}$-values, is mostly likely due to excited state contamination. The near physical point in that study is computed on a subset of the statistics (same smearing and $t_{\text{sink}}$) for ensemble IV. For the higher statistics used in our work, $g_A$ is not consistent with the experimental value at this small volume ($Lm_\pi = 2.7$). The extent of finite volume effects at the near physical point will become clearer once we achieve full statistics on the larger volume, ensemble V ($Lm_\pi = 3.5$).

The effect of excited state contributions can also be seen in the results for $\langle x \rangle_{u-d}$. Our values lie significantly below earlier results of QCDSF [15, 16] and ETMC [17]. Such effects have been seen in previous works [7, 8, 9, 10, 11]. The two volumes at $\beta = 5.29$ and $m_\pi \sim 290$ MeV, and similarly at the near physical point, indicate that finite volume effects are not significant for this quantity. Similarly, the pion mass dependence seems to be mild. The remaining discrepancy with the predictions from PDF parameterisations may be due to lattice spacing effects.
The scalar coupling suffers from larger statistical errors than the other quantities considered here. Within the large error, there is no significant dependence on $m_\pi$, volume or lattice spacing. Our results are consistent with other recent determinations. A similar picture is found for the tensor charge, although the statistical errors are smaller in this case.

5. Outlook

Control of excited state contributions and simulation at near physical pion masses are first steps towards a precise determination of $g_A$ and $\langle x \rangle_{u-d}$. For $g_A$, a careful volume extrapolation is needed, while in both cases the continuum limit needs to be studied. We are extending our simulations with this aim.

6. Acknowledgements

This work is supported by the EU ITN STRONGnet and the DFG SFB/TRR 55. Computations were performed on SuperMUC of the Leibniz Computing Center, the Regensburg iDataCool cluster and the SFB/TR55 QPACE supercomputers. The Chroma software suite [18] was used extensively in this work along with the domain decomposition solver implementation of [19].

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