



Nucleon isovector momentum fraction, helicity and transversity moment using Lattice QCD

Santanu Mondal,^{*a,c*,*} Tanmoy Bhattacharya,^{*a*} Rajan Gupta,^{*a*} Bálint Joó,^{*b*} Huey-Wen Lin,^{*c,d*} Sungwoo Park,^{*a*} Frank Winter^{*e*} and Boram Yoon^{*f*}

- ^aLos Alamos National Laboratory, Theoretical Division T-2, Los Alamos, NM 87545
- ^bOak Ridge Leadership Computing Facility, Oak Ridge National Laboratory, Oak Ridge, TN 37831, USA
- ^cDepartment of Physics and Astronomy, Michigan State University, MI, 48824, USA
- ^dDepartment of Computational Mathematics, Science and Engineering, Michigan State University, MI, 48824, U.S.A
- ^e Jefferson Lab, 12000 Jefferson Avenue, Newport News, Virginia 23606, USA
- ^f Los Alamos National Laboratory, Computer Computational and Statistical Sciences, CCS-7, Los Alamos, NM 87545

E-mail: mondals2@msu.edu

We present our recent high precision calculations [1, 2] of the first moment of nucleon isovector polarized, unpolarized and transversity distributions, i.e., momentum fraction, helicity and transversity moment, respectively. We use the standard method for the calculation of these moments (via matrix elements of twist two operators), and carry out a detailed analysis of the sources of systematic uncertainty, in particular of excited state contributions. Our calculations have been performed using two different lattice setups (Clover-on-HISQ and Clover-on-Clover), each with several ensembles. They give consistent results that are in agreement with global fit analyses.

The 38th International Symposium on Lattice Field Theory, LATTICE2021 26th-30th July, 2021 Zoom/Gather@Massachusetts Institute of Technology

*Speaker

© Copyright owned by the author(s) under the terms of the Creative Commons Attribution-NonCommercial-NoDerivatives 4.0 International License (CC BY-NC-ND 4.0).

Ensemble	a	M_{π}	$L^3 \times T$	$M_{\pi}L$	τ/a	N _{conf}
ID	(fm)	(MeV)				-
a15m310	0.1510(20)	320.6(4.3)	$16^3 \times 48$	3.93	{5, 6, 7, 8, 9}	1917
a12m310	0.1207(11)	310.2(2.8)	$24^3 \times 64$	4.55	{8, 10, 12, 14}	1013
a12m220	0.1184(09)	227.9(1.9)	$32^3 \times 64$	4.38	{8, 10, 12, 14}	1156
a12m220L	0.1189(09)	227.6(1.7)	$40^3 \times 64$	5.49	{8, 10, 12, 14}	1000
a09m310	0.0888(08)	313.0(2.8)	$32^3 \times 96$	4.51	{10, 12, 14, 16}	2263
a09m220	0.0872(07)	225.9(1.8)	$48^3 \times 96$	4.79	{10, 12, 14, 16}	960
a09m130	0.0871(06)	138.1(1.0)	$64^3 \times 96$	3.90	{10, 12, 14, 16}	1041
a06m310W	0.0582(04)	319.6(2.2)	$48^3 \times 144$	4.52	{18, 20, 22, 24}	500
a06m135	0.0570(01)	135.6(1.4)	$96^3 \times 192$	3.7	{16, 18, 20, 22}	751

Table 1: Lattice parameters of the 2 + 1 + 1-flavor HISQ ensembles generated by the MILC collaboration [3] and analyzed in this study. We give the lattice spacing *a*, pion mass M_{π} , lattice size $L^3 \times T$, the values of source-sink separation τ simulated, and the number of configurations analyzed.

1. Introduction

In the realm of QCD, among the key quantities to quantitatively characterize the rich and complex structure of hadrons are a number of universal, non-perturbative distribution functions. These are, parton distribution functions (PDFs), transverse momentum dependent PDFs (TMDs), generalized parton distributions (GPDs) and distribution amplitudes (DAs). For many years, there have been steady efforts to obtain these distributions both from experiments and theory. On the theoretical side, information on the moments of the distribution functions can be obtained via first principle Lattice QCD calculations. Subsequent to the proposal by Ji in 2013 [4], there have also been significant progress towards accessing the distributions themselves on the lattice [5].

The distributions are not measured directly in experiments, and phenomenological analyses including different theoretical inputs are needed to extract them from experimental data. In cases where both lattice results and phenomenological analyses of experimental data (global fits) exist, one can compare them to validate the control over systematics in the lattice calculations, and on the other hand provide a check on the phenomenological process used to extract these observables from experimental data [6, 7]. In other cases, lattice results are predictions.

Even for the best studied quantity on the lattice, the isovector momentum fraction $\langle x \rangle_{u-d}$, the data had large statistical and systematic uncertainties prior to 2018 [6]. Here we present high-precision lattice calculations for the isovector momentum fraction, helicity moment $\langle x \rangle_{\Delta u-\Delta d}$ and transversity moment $\langle x \rangle_{\delta u-\delta d}$ using two different lattice setups and controlling the sources of systematic uncertainties. Our study shows, that the lattice data for these three moments are now of quality comparable to that for nucleon charges (zeroth moments).

2. Lattice set up

We present our calculations of the three moments using two different lattice setups: (i) the Clover-on-HISQ calculation (PNDME 20) published in [1] was performed using nine HISQ ensembles generated by the MILC collaboration [3], whose parameters are summarized in Table 1. They

Ensemble	а	M_{π}	$L^3 \times T$	$M_{\pi}L$	τ/a	N _{conf}
ID	(fm)	(MeV)				
a127m285	0.127(2)	285(3)	$32^3 \times 96$	5.85	{8, 10, 12, 14}	2001
a094m270	0.094(1)	270(3)	$32^3 \times 64$	4.11	{10, 12, 14, 16, 18}	1464
a094m270L	0.094(1)	269(3)	$48^{3} \times 128$	6.16	{10, 12, 14, 16}	4501
a091m170	0.091(1)	169(2)	$48^3 \times 96$	3.74	{8, 10, 12, 14, 16}	4015
a091m170L	0.091(1)	169(2)	$64^3 \times 128$	5.08	{8, 10, 12, 14, 16}	1533
a073m270	0.0728(8)	272(3)	$48^3 \times 128$	4.8	{11, 13, 15, 17, 19}	4477
a071m170	0.0707(8)	167(2)	$72^3 \times 192$	4.26	{15, 17, 19, 21}	2100
a071m130	0.0707(8)	127(1)	$96^3 \times 192$	4.36	{13, 15, 17, 19, 21}	440
a056m280	0.056(1)	280(5)	$64^3 \times 192$	5.09	{18, 21, 24, 27, 30}	1723

Table 2: Lattice parameters of the 2 + 1-flavor clover ensembles generated by the JLab/W&M/LANL/MIT collaboration and analyzed in this study. We give the lattice spacing *a*, pion mass M_{π} , lattice size $L^3 \times T$, the values of source-sink separation τ simulated, and the number of configurations analyzed.

cover a range of lattice spacings $(0.057 \le a \le 0.15)$ fm, pion masses $(135 \le M_{\pi} \le 310)$ MeV and lattice sizes $(3.7 \le M_{\pi}L \le 5.5)$. For more details of the lattice methodology, the strategies for the calculations and the analyses see [1] and references therein.

(ii) The Clover-on-Clover calculation, published in [2] (NME 20), used seven Clover ensembles generated by the JLab/W&M/LANL/MIT collaboration [8]. Here (NME 21) we include two new ensembles, one at physical M_{π} and the other at smaller lattice spacing a = 0.056 fm which gives us better control over chiral and continuum extrapolations, respectively. Note that the data for the two new ensembles are preliminary and a071m130 is statistics limited. The parameters of these Clover ensembles are summarized in Table 2. They also cover a range of lattice spacings ($0.056 \le a \le 0.127$) fm, pion masses ($127 \le M_{\pi} \le 285$) MeV and lattice sizes ($3.74 \le M_{\pi}L \le 5.85$).

We construct the correlation functions needed to calculate the matrix elements using Wilsonclover fermions for both lattice setups. The Clover-on-HISQ formulation is non-unitary and can suffer from the problem of exceptional configurations at small, but a priori unknown, quark masses. However, we have not found evidence for such exceptional configurations on any of the nine ensembles analyzed in this work.

3. Lattice correlators and moments

The light quark operators $(q \in \{u, d\})$ used to calculate the moments are:

$$\langle x \rangle_{u-d} : O_{V^3}^{44} = \overline{q}(\gamma^4 \overleftrightarrow{D}^4 - \frac{1}{3}\gamma \cdot \overleftrightarrow{\mathbf{D}})\tau^3 q$$
 (1)

$$\langle x \rangle_{\Delta u - \Delta d} \quad : \quad O_{A^3}^{34} = \overline{q} \gamma^{\{3} \overleftrightarrow{D}^4\} \gamma^5 \tau^3 q \tag{2}$$

$$\langle x \rangle_{\delta u - \delta d} \quad : \quad O_{T^3}^{124} = \overline{q} \sigma^{[1\{2]} \overleftrightarrow{D}^{4\}} \tau^3 q \,. \tag{3}$$

From their matrix elements within the ground state of the nucleon, the moments are given by:

$$\langle 0|O_{V^3}^{44}|0\rangle = -M_N \langle x \rangle_{u-d}, \qquad (4)$$

$$\langle 0|O_{A^3}^{34}|0\rangle = -\frac{iM_N}{2} \langle x \rangle_{\Delta u - \Delta d}, \qquad (5)$$

$$\langle 0|O_{T^3}^{124}|0\rangle = -\frac{iM_N}{2} \langle x \rangle_{\delta u - \delta d}, \qquad (6)$$

where M_N is the nucleon mass. The nucleon interpolating operator N used is

$$\mathcal{N} = \epsilon^{abc} \left[q_1^{aT}(x) C \gamma^5 \frac{(1 \pm \gamma_4)}{2} q_2^b(x) \right] q_1^c(x) , \qquad (7)$$

where $\{a, b, c\}$ are color indices, $q_1, q_2 \in \{u, d\}$ and $C = \gamma_0 \gamma_2$ is the charge conjugation matrix. The nonrelativistic projection $(1 \pm \gamma_4)/2$ is inserted to improve the signal, with the plus and minus signs applied to the forward and backward propagation in Euclidean time, respectively. At zero momentum, this operator couples only to the spin $\frac{1}{2}$ states. The zero momentum two-point and three-point nucleon correlation functions are defined as

$$\boldsymbol{C}_{\alpha\beta}^{2pt}(\tau) = \sum_{\boldsymbol{x}} \langle 0 | \mathcal{N}_{\alpha}(\tau, \boldsymbol{x}) \overline{\mathcal{N}}_{\beta}(0, \boldsymbol{0}) | 0 \rangle$$
(8)

$$\boldsymbol{C}_{O,\alpha\beta}^{3pt}(\tau,t) = \sum_{\boldsymbol{x}',\boldsymbol{x}} \langle 0|\mathcal{N}_{\alpha}(\tau,\boldsymbol{x})O(t,\boldsymbol{x}')\overline{\mathcal{N}}_{\beta}(0,\boldsymbol{0})|0\rangle$$
(9)

where α , β are spin indices. The source is placed at time slice 0, the sink is at τ and the onederivative operators inserted at time slice *t*. Data have been accumulated for the values of τ specified in Tables 1 and 2, and, in each case, for all intermediate times $0 \le t \le \tau$.

4. Controlling the excited state contamination

A major challenge to precision results is removing the contribution of excited states in the three-point functions. These occur because the lattice nucleon interpolating operator, couples to the nucleon, all its excitations and to multi particle states with the same quantum numbers. The strategy to remove these artifacts are described in Refs. [1, 2]: reduce ESC by using smeared sources in the generation of quark propagators and then fit the data at multiple source-sink separations τ using the spectral decomposition of the correlation functions keeping as many excited states as possible without over-parameterizing the fits. The spectral decomposition of the zero-momentum two-point function, C_{2pt} , truncated at four states, is given by

$$C_{2\text{pt}}(\tau) = \sum_{i=0}^{3} |\mathcal{A}_i|^2 e^{-M_i \tau} \,. \tag{10}$$

We fit the data over the largest time range, $\{\tau_{min}-\tau_{max}\}$, allowed by statistics, i.e., by the stability of the covariance matrix, to extract the masses M_i and the amplitudes \mathcal{A}_i for the creation/annihilation of the four states by the interpolating operator. We perform two types of four-state fits. In the fit denoted {4}, we use the empirical Bayesian technique described in the Ref. [9] to stabilize the three

excited-state parameters. In the second fit, denoted $\{4^{N\pi}\}$, we use a normally distributed prior for M_1 , centered at the lower of the non-interacting energy of $N(-1)\pi(1)$ or the $N(0)\pi(0)\pi(0)$ state, and with a width of 0.04–0.05 in lattice units.

In the fits to the two-point functions, the {4} and $\{4^{N\pi}\}$ strategies cannot be distinguished on the basis of the χ^2 /dof. In fact, the full range of M_1 values between the two estimates, from {4} and $\{4^{N\pi}\}$, are viable on the basis of χ^2 /dof alone. The same is true of the values for M_2 , indicating a large flat region in parameter space. Because of this large region of possible values for the excited-state masses, M_i , we carry out the full analysis with three strategies that use different estimates of M_1 and investigate the sensitivity of the results on them.

The analysis of the three-point functions, $C_O^{3\text{pt}}$, is performed retaining up to three states $|i\rangle$ in the spectral decomposition:

$$C_O^{\text{3pt}}(\tau;t) = \sum_{i,j=0}^2 |\mathcal{A}_i| |\mathcal{A}_j| \langle i|O|j \rangle e^{-M_i t - M_j(\tau - t)} .$$

$$\tag{11}$$

To remove the ESC and extract the desired ground-state matrix element, $\langle 0|O|0\rangle$, we make a simultaneous fit in *t* and τ . In choosing the set of points, $\{t, \tau\}$, to include in the final fit, we attempt to balance statistical and systematic errors. First, we neglect t_{skip} points next to the source and sink in the fits as these have the largest ESC. Next, noting that the data at smaller τ have exponentially smaller errors but larger ESC, we pick the largest three values of τ for all seven ensembles. Since errors in the data grow with τ , we partially compensate for the larger weight given to smaller τ data by choosing t_{skip} to be the same for all τ , i.e., by including increasingly more *t* points with larger τ , the weight of the larger τ data points is increased. Most of our analysis uses a 3*-fit, which is a three-state fit with the term containing $\langle 2|O|2 \rangle$ set to zero, as it is undetermined and its inclusion results in an overparameterization based on the Akaike information criteria [10].

To investigate the sensitivity of $\langle 0|O|0 \rangle$ to possible values of M_i we carry out the full analysis with three strategies using the mnemonic $\{m, n\}$ to denote an *m*-state fit to the two-point function and an *n*-state fit to the three-point function. Figure 1 shows an example of difference in estimates from the three fit strategies for $\langle x \rangle_{u-d}$ from a073m270.

- {4, 3*}: The spectrum is taken from a {4} state fit to the two-point function using Eq. (10) and then a {3*} fit is made to the three-point function using Eq. (11). Both fits are made within a single jackknife loop. This is the standard strategy, which assumes that the same set of states are dominant in the two- and three-point functions.
- {4^{Nπ}, 3*}: The excited state spectrum is taken from a four-state fit to the two-point function but with a narrow prior for the first excited state mass taken to be the energy of a non-interacting N(p = 1)π(p = -1) state (or N(0)π(0)π(0) that has roughly the same energy). This spectrum is then used in a {3*} fit to the three-point function. This variant of the {4, 3*} strategy assumes that the lowest of the theoretically allowed tower of Nπ (or Nππ) states contributes.
- {4, 2^{free} }: The only parameters taken from the {4} state fit are the ground state amplitude \mathcal{R}_0 and mass M_0 , whose determination is robust. In the two-state fit to the three-point function,

the mass of the first excited state, M_1 , is left as a free parameter, ie, the most important determinant of ESC, M_1 , is obtained from the fit to the three-point function. The relative limitation of the {4, 2^{free}} strategy is that, with the current data, we can only make two-state fits to the three-point functions, i.e., include only one excited state.



Figure 1: Data of the ratio $\frac{C_0^{3pt}(\tau,t)}{C^{2pt}(\tau)}$ scaled with the kinematic factors to give $\langle x \rangle_{u-d}$ for the ensemble a073m270. The three panels show fits to the data with the largest three values of τ using three strategies: {4, 3*} (left panel), {4^N π , 3*} (middle panel) and {4, 2^{free}} (right panel). The fits are performed using data for the largest three values of τ . For each τ , the line in the same color as the data points is the result of the fit used to obtain the ground state matrix element. The blue band in each plot indicate the value of the moment obtained via the ground state matrix element from the fit.

5. Chiral, continuum and finite volume (CCFV) extrapolations

To obtain the final, physical results at $M_{\pi} = 135$ MeV, $M_{\pi}L \rightarrow \infty$ and a = 0, we make a simultaneous CCFV fit of data renormalized in the \overline{MS} scheme at $\mu = 2$ GeV keeping only the leading correction term in each variable:

$$\langle x \rangle (M_{\pi}; a; L) = c_1 + c_2 a + c_3 M_{\pi}^2 + c_4 \frac{M_{\pi}^2 e^{-M_{\pi}L}}{\sqrt{M_{\pi}L}} \,. \tag{12}$$

Note that, in both lattice setups the discretization errors start with a term linear in a. The results of the CCFV fits in case of PNDME 20 show that the finite volume correction term, c_4 , is not



Figure 2: Clover-on-HISQ data for $\langle x \rangle_{u-d}$ obtained via the fit strategy {4, 3*}, renormalized in the \overline{MS} scheme at $\mu = 2$ GeV, for all nine ensembles (PNDME 20). The blue band shows the CC fit result. In the left panel the fit is evaluated at $M_{\pi} = 135$ MeV and plotted versus *a*, while in the right panel it shows the result versus M_{π}^2 evaluated at a = 0.



Figure 3: Clover-on-Clover data for $\langle x \rangle_{u-d}$ obtained via the fit strategy {4, 3*}, renormalized in the *MS* scheme at $\mu = 2$ GeV, for seven ensembles ensembles (first row, NME 20) and for nine ensembles (second row, NME 21). The pink band shows the CCFV fit result. (Left) result evaluated at $M_{\pi} = 135$ MeV and $M_{\pi}L = \infty$ and plotted versus *a*. (Middle) result plotted versus M_{π}^2 and evaluated at a = 0 and $M_{\pi}L = \infty$. (Right) result plotted versus $M_{\pi}L$ and evaluated at a = 0 and $M_{\pi}L = \infty$.

constrained. Therefore, for PNDME 20, we use CC fits (i.e., with $c_4 = 0$ in Eq. 12) to obtain the final results. As an example, in Fig. 2 we present the PNDME 20 CC fit to $\langle x \rangle_{u-d}$ obtained via $\{4, 3^*\}$ fit strategy.

The NME 20 and 21 data are sensitive to the finite volume corrections. Therefore, we use CCFV fits to obtain the final results as shown in Fig. 3 for $\langle x \rangle_{u-d}$ obtained via {4, 3*} fit strategy.

In both lattice formulations and for all three moments, we find only a small positive slope with respect to both the lattice spacing and M_{π} . The main difference between the PNDME 20 and the NME 20/NME 21 results is the ~ 10% decrease due to the finite volume correction in the latter.

6. Results, comparison with the world data and conclusions

Moment	PNDME 20	NME 20	NME 21
$\langle x \rangle_{u-d}$	0.173(14)(07)	0.155(17)(20)	0.156(12)(20)
$\langle x \rangle_{\Delta u - \Delta d}$	0.213(15)(22)	0.183(14)(20)	0.185(12)(20)
$\langle x \rangle_{\delta u - \delta d}$	0.208(19)(24)	0.220(18)(20)	0.209(15)(20)

Fal	ble	3:	H	Final	resu	lts	for	the	moments	5.
-----	-----	----	---	-------	------	-----	-----	-----	---------	----

In Table 3 we compare PNDME 20, NME 20 and NME 21 results. We choose the results from CCFV/CC fits of the moments obtained via $\{4, 3^*\}$ fit strategy of the correlators. The first number inside brackets is overall statistical error. We will also take half the spread in results between



Figure 4: A comparison of results from lattice QCD calculations with dynamical fermions and global fits (below the black line). The left panel compares results for the momentum fraction, the middle for the helicity moment, and the right for the transversity moment. Our NME 21 result (preliminary) is also shown as the blue band to facilitate comparison.

 $\{4^{N\pi}, 3^*\}$ and the $\{4, 2^{\text{free}}\}$ fit strategies as a second uncertainty (the second number inside the brackets) to account for possible unresolved bias from incomplete control over ESC.

Our NME results are consistent with the PNDME 20 results. This is a valuable check of the PNDME 20 calculation that uses the non-unitary clover-on-HISQ lattice formulation. For $\langle x \rangle_{u-d}$ and $\langle x \rangle_{\Delta u-\Delta d}$ the NME results are $\approx 1\sigma$ smaller than the PNDME. A large part of the difference is due to the finite-volume correction in the NME results. There is reduction of statistical errors for all three quantities on going from NME 20 to NME 21. This is due to adding data from two new ensembles a071m130 and a056m280 in NME 21 which gives larger ranges in both the lattice-spacing and the pion mass in the CCFV fits.

An updated comparison of our results with other lattice calculations and phenomenological global fit estimates is given in Fig. 4. They are in good agreement with other recent lattice results by ETMC [11, 12], Mainz [13] and χ QCD [14] collaborations. Our estimate for the momentum fraction is in good agreement with most global fit estimates but has much larger error. The three estimates for the helicity moment from global fits have a large spread, and our estimate is consistent with the smaller error estimates. Lattice estimates for the transversity moment are a prediction.

7. Acknowledgements

We thank the MILC Collaboration for sharing the HISQ ensembles. The calculations used the Chroma software suite [15]. We gratefully acknowledge computing resources provided by NERSC, OLCF at Oak Ridge, USQCD and LANL Institutional Computing. Support for this work was provided the U.S. DOE Office of Science, HEP and NP, the NSF, and by the LANL LDRD program. The work of SM and HL are partially supported by the US National Science Foundation under grant PHY 1653405 "CAREER: Constraining Parton Distribution Functions for New-Physics Searches" and by the Research Corporation for Science Advancement through the Cottrell Scholar Award.

References

- S. Mondal, R. Gupta, S. Park, B. Yoon, T. Bhattacharya and H. W. Lin, Phys. Rev. D 102, no.5, 054512 (2020) doi:10.1103/PhysRevD.102.054512 [arXiv:2005.13779 [hep-lat]].
- [2] S. Mondal *et al.* [Nucleon Matrix Elements (NME)], JHEP **21**, 004 (2020) doi:10.1007/JHEP04(2021)044 [arXiv:2011.12787 [hep-lat]].
- [3] A. Bazavov et al., [MILC collaboration], Phys. Rev. **D87**, 054505 (2013) doi:10.1103/PhysRevD.87.054505 [arxiv:1212.4768 [hep-lat]].
- [4] X. Ji, Phys. Rev. Lett. 110, 262002 (2013) doi:10.1103/PhysRevLett.110.262002
 [arXiv:1305.1539 [hep-ph]].
- [5] K. Cichy, [arXiv:2110.07440 [hep-lat]].
- [6] H. W. Lin, E. R. Nocera, F. Olness, K. Orginos, J. Rojo, A. Accardi, C. Alexandrou, A. Bacchetta, G. Bozzi and J. W. Chen, *et al.* Prog. Part. Nucl. Phys. **100**, 107-160 (2018) doi:10.1016/j.ppnp.2018.01.007 [arXiv:1711.07916 [hep-ph]].
- [7] M. Constantinou, A. Courtoy, M. A. Ebert, M. Engelhardt, T. Giani, T. Hobbs, T. J. Hou, A. Kusina, K. Kutak and J. Liang, *et al.* Prog. Part. Nucl. Phys. **121**, 103908 (2021) doi:10.1016/j.ppnp.2021.103908 [arXiv:2006.08636 [hep-ph]].
- [8] R. Edwards, R. Gupta, B. Joo, K. Orginos, D. Richards, F. Winter, and B. Yoon, "U.S. 2+1 flavor clover lattice generation program," (2016), unpublished.
- [9] B. Yoon, Y. C. Jang, R. Gupta, T. Bhattacharya, J. Green, B. Joó, H. W. Lin, K. Orginos, D. Richards and S. Syritsyn, *et al.* Phys. Rev. D **95**, no.7, 074508 (2017) doi:10.1103/PhysRevD.95.074508 [arXiv:1611.07452 [hep-lat]].
- [10] H. Akaike, IEEE Transactions on Automatic Control, vol. 19, no. 6, pp. 716-723, December 1974, doi: 10.1109/TAC.1974.1100705.
- [11] C. Alexandrou, S. Bacchio, M. Constantinou, J. Finkenrath, K. Hadjiyiannakou, K. Jansen, G. Koutsou, H. Panagopoulos and G. Spanoudes, Phys. Rev. D 101, no.9, 094513 (2020) doi:10.1103/PhysRevD.101.094513 [arXiv:2003.08486 [hep-lat]].
- [12] C. Alexandrou, S. Bacchio, M. Constantinou, P. Dimopoulos, J. Finkenrath, R. Frezzotti, K. Hadjiyiannakou, K. Jansen, B. Kostrzewa and G. Koutsou, *et al.* Phys. Rev. D 101, no.3, 034519 (2020) doi:10.1103/PhysRevD.101.034519 [arXiv:1908.10706 [hep-lat]].
- [13] T. Harris, G. von Hippel, P. Junnarkar, H. B. Meyer, K. Ottnad, J. Wilhelm, H. Wittig and L. Wrang, Phys. Rev. D 100, no.3, 034513 (2019) doi:10.1103/PhysRevD.100.034513 [arXiv:1905.01291 [hep-lat]].
- [14] Y. B. Yang, J. Liang, Y. J. Bi, Y. Chen, T. Draper, K. F. Liu and Z. Liu, Phys. Rev. Lett. 121, no.21, 212001 (2018) doi:10.1103/PhysRevLett.121.212001 [arXiv:1808.08677 [hep-lat]].
- [15] R. G. Edwards *et al.* [SciDAC, LHPC and UKQCD], Nucl. Phys. B Proc. Suppl. **140**, 832 (2005) doi:10.1016/j.nuclphysbps.2004.11.254 [arXiv:hep-lat/0409003 [hep-lat]].