













Quark transverse spin-momentum correlation of the nucleon from lattice QCD: the Boer-Mulders function



The Lattice Parton collaboration

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ABSTRACT: We present the first lattice QCD calculation of the quark transverse spin-momentum correlation, i.e., the naive time-reversal-odd Boer-Mulders function, of the nucleon, using large-momentum effective theory (LaMET). The calculation is carried out on an ensemble with lattice spacing $a = 0.098$ fm and pion mass 338 MeV, at various proton momenta up to 2.11 GeV. We have implemented perturbative matching up to the next-to-next-to-leading order together with a renormalization-group resummation improvement. The result exhibits the expected decay behavior with increasing transverse separation b_{\perp} . We also compare the results in the nucleon and pion.

KEYWORDS: Hadronic Spectroscopy, Structure and Interactions, Lattice QCD

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1 Introduction

Transverse-momentum-dependent parton distribution functions (TMDPDFs), together with generalized parton distributions (GPDs), play an essential role in mapping the three-dimensional structure of nucleons, known as the nucleon tomography program [1]. In contrast to GPDs which characterize the correlation between transverse position and longitudinal momentum of partons, TMDPDFs describe the correlation between transverse and longitudinal momenta of partons. At leading-twist accuracy, there are eight independent quark TMDPDFs which are characterized by their nucleon and quark polarizations [2]. Among them, the naive time-reversal-odd Sivers and Boer-Mulders functions, are of special interest. While the Sivers function characterizes the correlation of quark transverse momentum and nucleon transverse polarization, the Boer-Mulders function describes the correlation between quark transverse polarization and momentum.

Due to their T-odd nature, the factorization of processes involving Sivers and Boer-Mulders TMDPDFs requires the inclusion of other T-odd nonperturbative quantities such as T-odd fragmentation functions [3, 4]. This renders experimental measurements of these TMDPDFs very challenging. Nevertheless, phenomenological efforts have been made to extract these functions from azimuthal angle asymmetries in semi-inclusive deep inelastic scattering (SIDIS) and Drell-Yan processes [5–9]. As a complementary tool, first-principles nonperturbative approaches like lattice QCD can also offer valuable insights into these functions. To achieve this, it is essential to calibrate them using quantities for which both state-of-the-art experiments and lattice simulations can reliably estimate precision, such as the unpolarized quark TMDPDFs. This will allow us to trust other lattice results to a similar

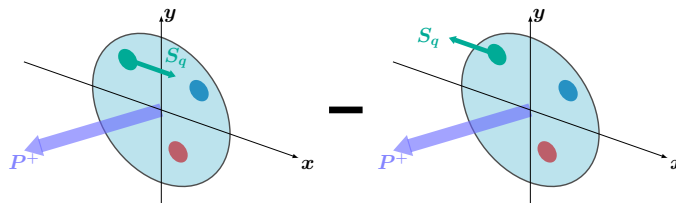


Figure 1. Illustration of the physical interpretation of the Boer-Mulders TMDPDF. S_q denotes the quark spin.

degree, even when they cannot yet be precisely tested by experiments. With this in mind, our long-term goal is to calculate all leading-twist TMDPDFs of hadrons using lattice QCD.

There have been exploratory lattice studies of the TMDPDFs, including the Sivers and Boer-Mulders functions, with the goal to access their moments by forming ratios of appropriate correlators [10–13]. More recently, the development of large-momentum effective theory (LaMET) [14–16] has made it possible to calculate the full TMDPDFs. Notably, a lattice calculation of the nucleon unpolarized quark TMDPDF [17] and the pion Boer-Mulders function [18] has been carried out within the LaMET framework. In contrast to the unpolarized quark TMDPDF presented in ref. [17], the pion Boer-Mulders function [18] shows clearly the expected decay with increasing transverse separations between the quark fields. This suggests that different TMDPDFs might receive rather different higher-twist contaminations, as analyzed in ref. [19].

In this work, we present a calculation of the isovector nucleon Boer-Mulders function $h_1^{\perp,u}(x, b_\perp) - h_1^{\perp,d}(x, b_\perp)$ using LaMET. An illustration of the function is shown in figure 1. The calculation is done on a lattice ensemble with lattice spacing $a = 0.098$ fm and pion mass 338 MeV. The proton momenta are up to 2.11 GeV. The lattice matrix elements are nonperturbatively renormalized in the short-distance ratio scheme [20] and extrapolated to large quasi-light-front (quasi-LF) distances following the same strategy as in refs. [17, 18]. After performing a Fourier transform to longitudinal momentum space, the quasi-Boer-Mulders function is matched to the standard Boer-Mulders function using a factorization formula presented in ref. [21]. The intrinsic soft function and Collins-Soper evolution kernel are taken from previous calculations on the same ensemble [22, 23], and the perturbative matching kernel is implemented up to the next-to-next-to-leading order (NNLO) using refs. [24, 25]. We also use renormalization-group resummation [26]. We then compare the Boer-Mulders functions in the nucleon and pion.

The rest of the paper is organized as follows: in section 2, we give a brief overview of the theoretical framework. In section 3, we present the details of our lattice calculation. In section 4, we discuss the numerical results for the Boer-Mulders TMDPDF in the nucleon and compare it with the same function in the pion. A summary is given in section 5.

2 Theoretical framework

To access the quark transverse spin-momentum correlation, or Boer-Mulders function, we start from the following subtracted quasi-TMDPDF matrix element [21, 27–29]

$$\tilde{f}(z, b_\perp, P^z, a) = \lim_{L \rightarrow \infty} \frac{\langle P | \hat{O}_\square(z, L, b_\perp) | P \rangle}{\sqrt{Z_E(2L + z, b_\perp, a)}}, \quad (2.1)$$

where $|P\rangle$ denotes an unpolarized nucleon state with momentum $P = (P^0, 0, 0, P^z)$, the numerator $\langle P|\hat{O}_\square(z, L, b_\perp)|P\rangle$ is often referred to as the unsubtracted quasi-TMDPDF matrix element with the quark correlator

$$\hat{O}_\square(z, L, b_\perp) = \bar{\psi}(b_\perp \hat{n}_\perp) i\sigma^{yt} \gamma_5 \mathcal{W}_\square(z, L, b_\perp) \psi(z \hat{n}_z), \quad (2.2)$$

where \mathcal{W}_\square represents a staple-shaped gauge link

$$\begin{aligned} \mathcal{W}_\square(b_\perp, L, z) &= U_z^\dagger((z+L)\hat{n}_z + b_\perp \hat{n}_\perp, b_\perp \hat{n}_\perp) \\ &\quad \times U_\perp((z+L)\hat{n}_z + b_\perp \hat{n}_\perp, (z+L)\hat{n}_z) U_z((z+L)\hat{n}_z, z \hat{n}_z), \\ U_i(\eta + s \hat{n}_i, \eta) &= \mathcal{P} \exp \left[-ig \int_0^s dt \hat{n}_i \cdot A(\eta^\mu + t \hat{n}_i^\mu) \right]. \end{aligned} \quad (2.3)$$

\hat{n}_z, \hat{n}_\perp are unit vectors along the spatial z and transverse directions, respectively. The staple-shaped gauge link \mathcal{W}_\square becomes future-pointing under a large boost $P^z \rightarrow \infty$, indicating that the computed TMDPDF is relevant for a SIDIS process and differs from that for Drell-Yan processes by a minus sign. Z_E is a flat rectangular Euclidean Wilson-loop along the n_z direction with length $2L + z$ and width b_\perp ,

$$\begin{aligned} Z_E(2L + z, b_\perp, a) &= \frac{1}{N_c} \text{Tr} \langle 0 | U_\perp^\dagger(-\vec{L} + \vec{b}; -b_\perp) \times U_z^\dagger(\vec{L} + \vec{z} + \vec{b}_\perp; -2L - z) \\ &\quad \times U_\perp(\vec{L} + \vec{z}; b_\perp) U_z(-\vec{L}; 2L + z) | 0 \rangle, \end{aligned} \quad (2.4)$$

in which $\vec{L} = L \hat{n}_z$, $\vec{z} = z \hat{n}_z$ and $\vec{b} = b \hat{n}_\perp$. The link length L is introduced to regulate the pinch-pole singularity associated with the longitudinal gauge links [16], and can be safely taken to infinity in \tilde{f} , as its dependence cancels in the ratio up to power corrections that are suppressed by L . In the subtracted quasi-TMDPDF, there are still residual logarithmic divergences originating from the endpoints of the quark correlator. They can be removed in the short-distance ratio scheme [20] where the renormalization factor is defined as a subtracted quasi-TMDPDF matrix element at small $z = z_0$ and $b_\perp = b_{\perp,0}$ within the perturbative region and for zero momentum,

$$Z_O(1/a, \mu) = \frac{\tilde{f}(z_0, b_{\perp,0}, 0, a)}{\tilde{h}_\Gamma^{\overline{\text{MS}}}(z_0, b_{\perp,0}, \mu)}. \quad (2.5)$$

Note that in the equation above we have included the conversion factor to $\overline{\text{MS}}$ scheme $\tilde{h}_\Gamma^{\overline{\text{MS}}}$, which is expected to be the same for the quark quasi-Boer-Mulders function and for the unpolarized quark quasi-TMDPDF [30] and takes the following form up to one-loop accuracy [20]

$$\tilde{h}_\Gamma^{\overline{\text{MS}}}(z, b_\perp, \mu) = 1 + \frac{\alpha_s(\mu) C_F}{2\pi} \left[\frac{1}{2} + \frac{3}{2} \ln \left(\frac{\mu^2 (b_\perp^2 + z^2) e^{\gamma_E}}{4} \right) - 2 \frac{z}{b_\perp} \arctan \frac{z}{b_\perp} \right], \quad (2.6)$$

with $\alpha_s = g^2/(4\pi)$, $C_F = 4/3$. The fully renormalized quasi-TMDPDF then reads

$$\tilde{f}_R(z, b_\perp, P^z, \mu) = Z_O^{-1}(1/a, \mu) \tilde{f}(z, b_\perp, P^z, a). \quad (2.7)$$

In the renormalization factor Z_O , the dependence on z_0 and $b_{\perp,0}$ is expected to cancel. However, this cancellation is never complete due to missing higher-order perturbative contributions. To reduce the dependence on z_0 or $b_{\perp,0}$, one can perform a renormalization group (RG) resummation [26] from some reference scale μ_0 to μ , where μ_0 is the physical scale that can be chosen as $\frac{2e^{-\gamma_E}}{\sqrt{b_{\perp}^2+z^2}}$. The evolved perturbative result, denoted as $\tilde{h}_{\Gamma}^{\overline{\text{MS}},\text{RGR}}(z, b_{\perp}, \mu)$, is then used in eq. (2.5) instead of $\tilde{h}_{\Gamma}^{\overline{\text{MS}}}(z, b_{\perp}, \mu)$:

$$\tilde{h}_{\Gamma}^{\overline{\text{MS}},\text{RGR}}(z, b_{\perp}, \mu) = \tilde{h}_{\Gamma}^{\overline{\text{MS}}}(z, b_{\perp}, \mu_0) \exp \left[\int_{\alpha_s(\mu_0)}^{\alpha_s(\mu)} d\alpha' \frac{\gamma_F(\alpha')}{\beta(\alpha')} \right]. \quad (2.8)$$

The renormalized quasi-TMDPDF in momentum space is given by the Fourier transform

$$\tilde{f}_R(x, b_{\perp}, P^z, \mu) = \int \frac{dz}{2\pi} e^{-iz(xP^z)} \tilde{f}_R(z, b_{\perp}, P^z, \mu). \quad (2.9)$$

It can be related to the physical Boer-Mulders TMDPDF by the following factorization formula [21]

$$\begin{aligned} \tilde{f}_R(x, b_{\perp}, \zeta_z, \mu) \sqrt{S_I(b_{\perp}, \mu)} &= H_{\Gamma} \left(\frac{\zeta_z}{\mu^2} \right) e^{\frac{1}{2} \ln(\zeta_z/\zeta) K(b_{\perp}, \mu)} f(x, b_{\perp}, \zeta, \mu) \\ &+ \mathcal{O} \left(\frac{\Lambda_{\text{QCD}}^2}{\zeta_z}, \frac{M^2}{P_z^2}, \frac{1}{b_{\perp}^2 \zeta_z} \right), \end{aligned} \quad (2.10)$$

where ζ is the rapidity scale, and $\zeta_z = (2xP_z)^2$. $S_I(b_{\perp}, \mu)$ denotes the intrinsic soft function which is associated with the emission of soft gluons [21], $K(b_{\perp}, \mu)$ is the Collins-Soper evolution kernel [31]. Both S_I and K can be calculated nonperturbatively in lattice QCD [22, 23, 31–41]. The $\mathcal{O} \left(\frac{\Lambda_{\text{QCD}}^2}{\zeta_z}, \frac{M^2}{P_z^2}, \frac{1}{b_{\perp}^2 \zeta_z} \right)$ term denotes power corrections. $H_{\Gamma} = e^h$ is the hard matching kernel which has been calculated up to the next-to-next-to-leading order [16, 24, 25, 28, 29].

One can also perform an RG resummation to deal with the potentially large double logarithms $\sim \alpha_s^n \ln^{2n}(\zeta_z/\mu^2)$ in the hard kernel H_{Γ} and it improves the reliability of the perturbative matching. The RG resummation starts from a fixed order perturbation series at the physical scale ζ_z and evolves it to the renormalization scale μ

$$H \left(\alpha_s(\mu), \frac{\zeta_z}{\mu^2} \right) = H \left(\alpha_s(\sqrt{\zeta_z}) \right) \exp \left\{ \int_{\sqrt{\zeta_z}}^{\mu} \frac{d\mu'}{\mu'} \left[\Gamma_{\text{cusp}}(\alpha_s(\mu')) \ln \frac{\zeta_z}{\mu'^2} + \gamma_C(\alpha_s(\mu')) \right] \right\}, \quad (2.11)$$

where Γ_{cusp} and γ_C denote the cusp anomalous dimension and the single log anomalous dimension, respectively. Their explicit expressions can be found in refs. [17, 18]. The scale $\sqrt{\zeta_z}$ is varied between $0.8\sqrt{\zeta_z}$ and $1.2\sqrt{\zeta_z}$ to estimate systematic uncertainties related to the scale choice.

3 Lattice calculation

3.1 Lattice setup

In this work, we use the lattice ensemble X650 generated by the CLS collaboration [42] using 2+1 flavor dynamical clover fermions and tree-level Symanzik improved gauge action.

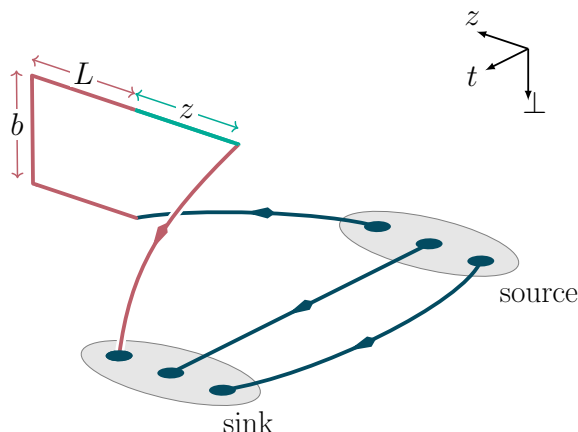


Figure 2. Illustration of the nucleon three-point function. The time direction is from source to sink.

Ensemble	$a(\text{fm})$	$L^3 \times T$	$m_\pi(\text{MeV})$	$m_\pi L$	$N_{\text{conf.}}$
X650	0.098	$48^3 \times 48$	338	8.1	1250

Table 1. The lattice ensemble used in this calculation [45].

Ensemble	t_{sep}/a	z_{max}/a	$b_{\perp, \text{max}}/a$	L/a
X650	{6, 7, 8, 9, 10}	18	3	{6, 8, 10}

Table 2. Source-sink separations t_{sep}/a , maximum longitudinal and transverse separations z_{max}/a and $b_{\perp, \text{max}}/a$ of the quark fields in the quasi-TMDPDF defined in eq. (2.1) and staple link length L/a used in the analysis.

The lattice spacing is $a = 0.098 \text{ fm}$ and pion mass is $m_\pi = 338 \text{ MeV}$. Two different types of smearing are adopted to improve the signal-to-noise ratio: the momentum smearing source technique [43] to improve the signal for a fast moving nucleon and HYP smearing to improve the signal for non-local operators with large separations [44].

We use the sequential source method with fixed sink to facilitate the calculation of the nucleon three-point correlator, as illustrated in figure 2, where we also show the staple-shaped gauge link defined in eq. (2.3) with length parameters z , b_\perp , L . To increase the number of measurements, we put two sources in the temporal direction and 2, 2, 1 sources in the x , y , z directions, respectively. Details of the lattice setup and parameters are collected in table 1. The bare matrix elements are calculated with the nucleon carrying different spatial momenta: $P^z = \{1.32, 1.58, 1.84, 2.11\} \text{ GeV}$. The bare matrix element of a nucleon at rest is also needed to extract the renormalization factor, as shown in eq. (2.5). In table 2, we give details on the temporal source-sink separations, the maximum of the longitudinal separation z_{max} , transverse separation b_{max} , as well as the staple length L .

3.2 Dispersion relation

To estimate the discretization effects of the X650 ensemble using the nucleon dispersion relation and extract the quasi-TMDPDF matrix element in eq. (2.1), we calculate the

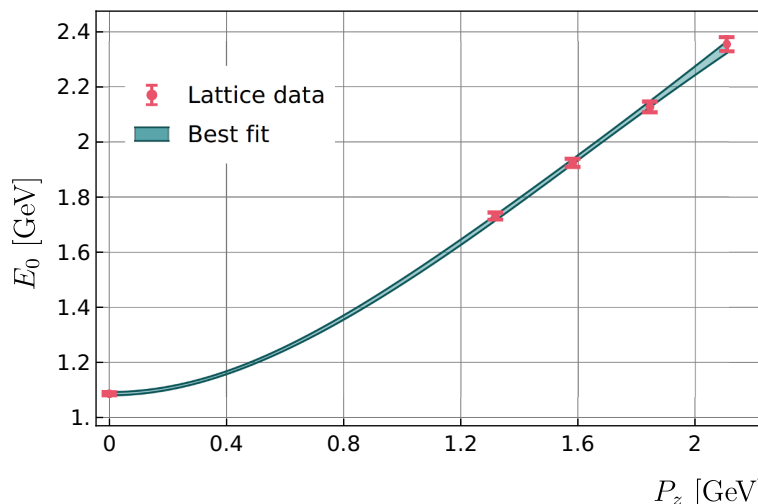


Figure 3. The dispersion relation of the nucleon on the X650 ensemble.

two-point function $C_2(P^z, t_{\text{sep}})$ defined as

$$C_2(P^z, t_{\text{sep}}) = \sum_{\vec{x}} e^{-i\vec{P}\cdot\vec{x}} T_u \langle \chi(t_{\text{sep}}, \vec{x}) \bar{\chi}(0, \vec{0}) \rangle, \quad (3.1)$$

where $T_u = (1 + \gamma^t)/2$ projects out the unpolarized nucleon state. $\chi = \epsilon^{abc} u_a (u_b^T C \gamma_5 d_c)$ is the nucleon interpolation operator. By considering the ground state and first excited state, the two-point function can be parameterized as

$$C_2(P^z, t_{\text{sep}}) = c_0 e^{-E_0 t_{\text{sep}}} (1 + c_1 e^{-\Delta E t_{\text{sep}}}), \quad (3.2)$$

where E_0 denotes the ground state energy of the nucleon, ΔE is the energy shift between the first excited state and the ground state. The dispersion relation of the ground state on the lattice reads

$$E_0(P^z) = \sqrt{M^2 + c_1 P_z^2 + c_2 P_z^4 a^2}, \quad (3.3)$$

where c_1 and c_2 take into account the discretization effect and reduce to 1 and 0 in the continuum limit $a \rightarrow 0$, respectively.

By fitting two-point functions at various momenta using eq. (3.2), we obtain the nucleon’s ground state energy E_0 . E_0 for various P_z and the fitted curve are plotted in figure 3 with $c_1 = 1.073(40)$, $c_2 = -0.094(50)$, which are consistent with the continuum limit within 2σ . This indicates that the discretization effect is under control. For the lack of resources we leave more detailed investigations of discretization effects and pion mass dependence using multiple lattice ensembles for future work. Such an investigation is essential for a precise lattice determination of the nucleon Boer-Mulders function. The extracted ground state nucleon mass M is 1.087(6) GeV.

3.3 Wilson loop from lattice calculations

The Wilson loop $Z_E(r = 2L + z, b_\perp)$ is obtained by calculating the expectation value of a closed rectangular gauge link with side lengths r and b_\perp . At large r and b_\perp , the signal-to-noise

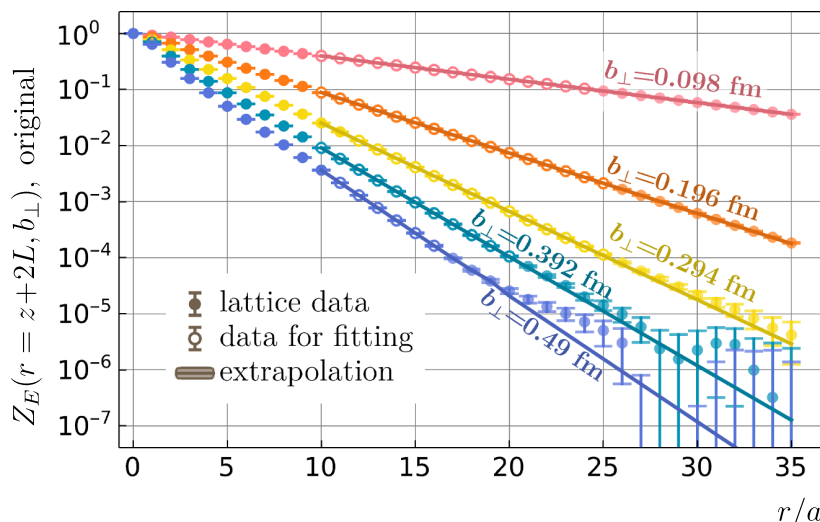


Figure 4. The Wilson loop and its extrapolation with respect to r and b_{\perp} . The points and error bars show the original calculations of the Wilson loop, with worse signals at larger (r, b_{\perp}) . The empty points indicate the fit ranges for extrapolation. The bands show extrapolated results by fitting the effective energy of Z_E . We extrapolate the Wilson loop up to $r = 60a$.

ratio deteriorates, as shown in figure 4. The lattice values of Z_E even turn negative at certain points of large r and b_{\perp} . However, we know that the Wilson loop has to be positive. To avoid this, we use the same strategy as in ref. [20] to fit the effective energies of Z_E which give the QCD static potential in the form $Z_E(2L + z, b_{\perp}) = c(b_{\perp})e^{-V(b_{\perp})(2L+z)}$ with $V(b_{\perp})$ being the static potential. We then extrapolate the result to large values of r . The extrapolated Wilson loop is shown as bands in figure 4. In our calculation, we use the extrapolated Wilson loop to obtain the subtracted quasi-TMDPDF matrix element.

3.4 Unsubtracted quasi-TMDPDF matrix elements from lattice calculations

To extract the unsubtracted quasi-TMDPDF matrix elements, we calculate the three-point function $C_3(P^z, t_{\text{sep}}, t)$

$$C_3(P^z, t_{\text{sep}}, t) = \sum_{\vec{x}} e^{-i\vec{P}\cdot\vec{x}} T_u \left\langle \chi(t_{\text{sep}}, \vec{x}) \sum_{\vec{y}} \hat{O}(t, \vec{y}) \bar{\chi}(0, \vec{0}) \right\rangle, \quad (3.4)$$

where $\hat{O}(t, \vec{y})$ denotes the bilinear operator $\hat{O}_{\square}(z, L, b_{\perp})$ in eq. (2.1) that is inserted at the discrete time slice t . We parameterize the three-point function using a two-state fit

$$C_3(P^z, t_{\text{sep}}, t) = c_0 e^{-E_0 t_{\text{sep}}} \times [\tilde{h}_0 + c_2(e^{-\Delta E t} + e^{-\Delta E(t_{\text{sep}}-t)}) + c_3 e^{-\Delta E t_{\text{sep}}}], \quad (3.5)$$

and form the ratio of the three- and two-point function

$$\frac{C_3}{C_2} = \frac{\tilde{h}_0 + c_2(e^{-\Delta E t} + e^{-\Delta E(t_{\text{sep}}-t)}) + c_3 e^{-\Delta E t_{\text{sep}}}}{1 + c_1 e^{-\Delta E t_{\text{sep}}}}, \quad (3.6)$$

where $\tilde{h}_0 = \langle P | \hat{O}_{\square}(z, L, b_{\perp}) | P \rangle / 2E_0(P^z)$ yields the desired matrix element, $E_0(P^z)$ is the ground state energy. For simplicity, we denote the r.h.s. of eq. (3.6) as $R(t, t_{\text{sep}})$ hereafter.

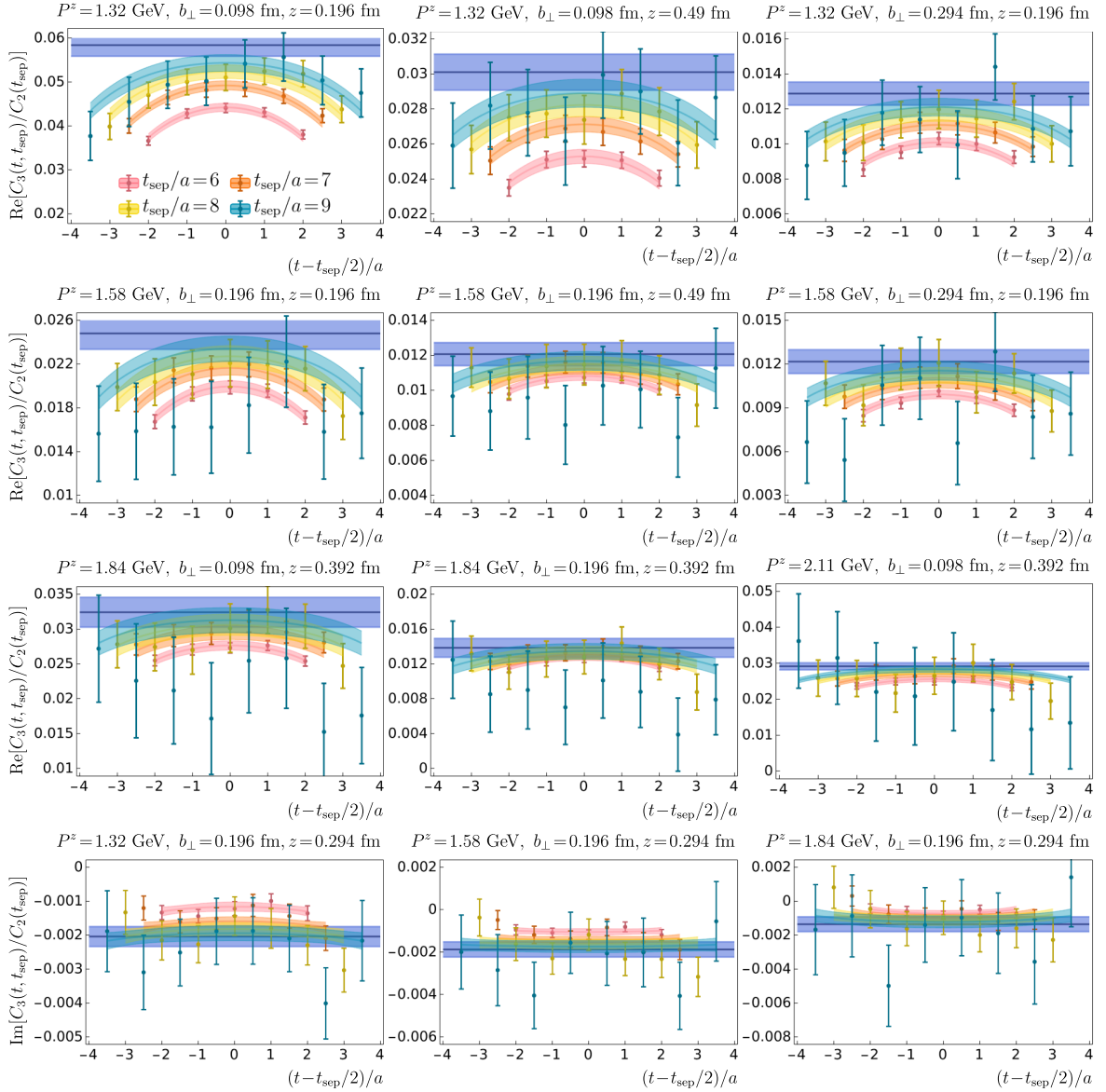


Figure 5. The real and imaginary parts of the ratio $R(t, t_{\text{sep}})$ versus $t - t_{\text{sep}}/2$ for various momenta and (z, b_{\perp}) , with $L = 8a$. The straight grey band shows the central value and error of \tilde{h}_0 by fitting to the lattice data. The curved bands are $R(t, t_{\text{sep}})$ values reconstructed from the fitted parameters. The values of t_{sep} involved in the fitting are $t_{\text{sep}} = \{6, 7, 8, 9\}a = \{0.588, 0.686, 0.784, 0.882\}$ fm.

We use bootstrap resampling in our analysis to treat correlations in the dataset. A joint fit of the two-point function eq. (3.2) and the ratio eq. (3.6) is done after resampling data for all combinations of (P^z, z, L, b_{\perp}) . The points at $t = 0$ and $t = t_{\text{sep}}$ are excluded in order to reduce contamination from excited states. Considering the bad signal-to-noise ratio for larger t_{sep} , we use $t_{\text{sep}}/a = \{6, 7, 8, 9\}$ for nonzero momenta.

As an illustrative example, we plot in figure 5 the lattice data and fit results of the real and imaginary part of the ratio $R(t, t_{\text{sep}})$ with $L = 8a$. The scattered points and error bars are lattice data corresponding to (t_{sep}, t) that are used in the fitting. The grey bands

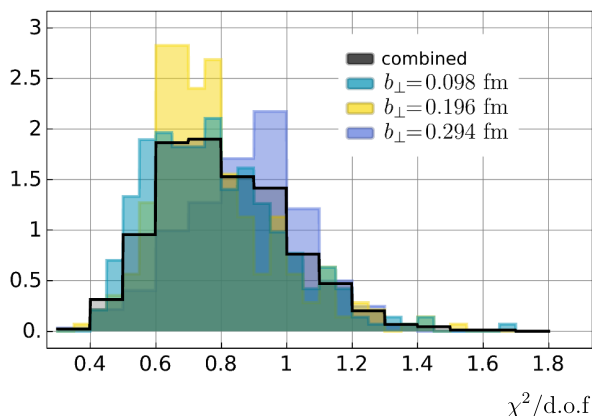


Figure 6. Probability distributions of $\chi^2/\text{d.o.f.}$. The distributions are normalized to 1.

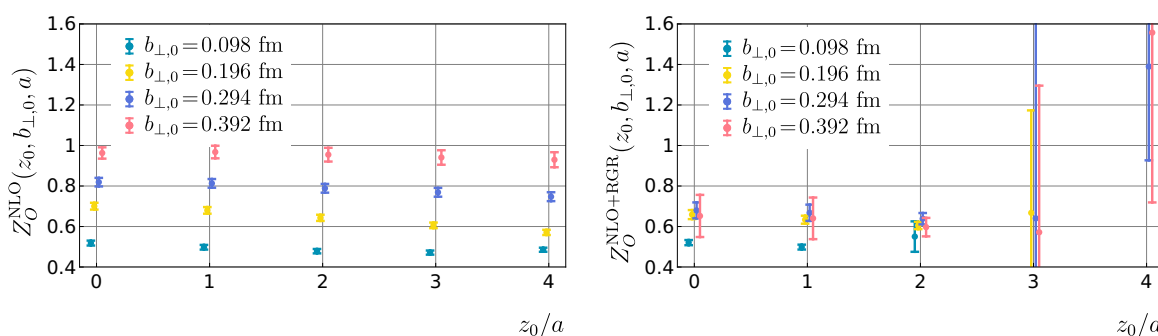


Figure 7. The renormalization factor Z_O , with the NLO result as well as the NLO+RGR result at $P^z = 0$ and various $b_\perp = b_{\perp,0}$ and $z = z_0$.

show the central values and errors of \tilde{h}_0 . The curved bands are the reconstructed $R(t, t_{\text{sep}})$ from the fitted parameters.

To assess the quality of the combined fits, we plot histograms of $\chi^2/\text{d.o.f.}$ in figure 6. The distributions have been normalized to 1.

3.5 Renormalization and L dependence of the subtracted quasi-TMDPDF matrix elements

To obtain the fully renormalized quasi-TMDPDF matrix element, we still need the logarithmic renormalization factor Z_O . It is computed from the zero-momentum matrix element of the same quark quasi-TMDPDF operator with $z = z_0$, $b_\perp = b_{\perp,0}$ being chosen within the perturbative region, and combined with the perturbative conversion factor. In figure 7, we show Z_O obtained from the NLO perturbative result of $\tilde{h}_\Gamma^{\overline{\text{MS}}}$ in the left panel and from the NLO+RGR result in the right panel. We can see a window with $z_0 \leq 3a$ and $b_{\perp,0}/a = \{2, 3\}$ where the dependence of Z_O on z_0 and $b_{\perp,0}$ is significantly reduced. Therefore, we use the evolved perturbative result $\tilde{h}_\Gamma^{\overline{\text{MS}},\text{RGR}}$ in eq. (2.8) to calculate Z_O , where we average over z_0 and $b_{\perp,0}$ in the window region. The result is $0.660(16)(24)$, where the number in the first parenthesis is the statistical error, while the number in the second parenthesis is the systematical error estimated from varying the RG evolution scale from $0.8\mu_0$ to $1.2\mu_0$, with $\mu_0 = 2e^{-\gamma_E}/\sqrt{b_{\perp,0}^2 + z_0^2}$.

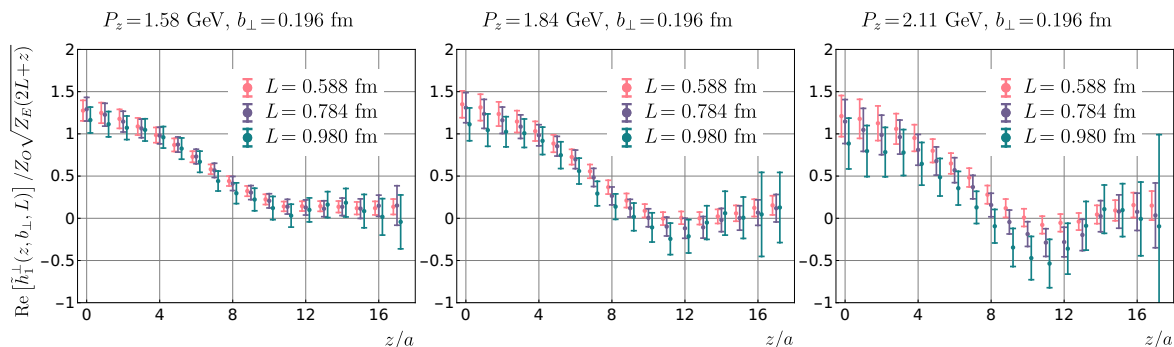


Figure 8. The renormalized quasi-TMDPDF $\tilde{h}_1^\perp(z, b_\perp, P^z, L)$, taking $b_\perp = 0.196$ fm as an example. The convergence with increasing L indicated the existence of an infinite L limit.

After dividing the unsubtracted quasi-TMDPDF matrix element by $\sqrt{Z_E(2L+z)}$ and Z_O , we find that the fully renormalized result reaches a reasonable plateau in the interval $6 \leq L/a \leq 10$. As an example, we show in figure 8 the renormalized quasi-TMDPDF matrix elements for various (P^z, b_\perp, L) versus z . Convergence can be observed for various L . The systematic error from varying μ_0 has been considered in the renormalization. In the following, we take the result at $L = 8a = 0.784$ fm as an approximation of the result in the infinite L limit, and take the difference between results at $L = 8a = 0.784$ fm and $L = 10a = 0.980$ fm as an estimate of the systematic uncertainty due to finite L . The resulting systematic uncertainty is shown in figure 12.

3.6 Large λ extrapolation of quasi-TMDPDF in coordinate space

As we can see from figure 8, the error of the quasi-TMDPDF increases rapidly at large $\lambda = zP^z$. In order to facilitate the subsequent Fourier transform, we perform an extrapolation to large λ using the same ansatz as that used in ref. [17]

$$\tilde{h}_{\text{extra}}(\lambda) = \left[\frac{m_1}{(-i\lambda)^{n_1}} + e^{i\lambda} \frac{m_2}{(i\lambda)^{n_2}} \right] e^{-\lambda/\lambda_0}, \quad (3.7)$$

where all parameters $m_{1,2}$, $n_{1,2}$ and λ_0 depend on b_\perp . The algebraic terms account for a power law behavior in the endpoint. Convergence can be observed for various regions of x , while the exponential decay is based on the expectation that the correlation length of the function (denoted as λ_0) is finite at finite momentum. We perform an independent extrapolation for each b_\perp . In figure 9, we show some examples of extrapolation for various $P^z = \{1.58, 1.84, 2.11\}$ GeV and $b_\perp = 0.196$ fm. For $P^z = 1.58$ GeV the imaginary part is also plotted. In these figures, both the original data points of \tilde{h}_1^\perp and the extrapolation bands from the fitting eq. (3.7) are shown. We indicate in each plot the region chosen to fit the extrapolation form. The extrapolation results agree with lattice data in the intermediate λ region and give smooth curves with much reduced errors for large λ and we replace the lattice data with $\tilde{h}_{\text{extra}}(\lambda)$ starting from the beginning of the fitted region.

3.7 Matching to physical Boer-Mulders TMDPDF

With the extrapolation at large λ , we can Fourier transform the Boer-Mulders quasi-TMDPDF to momentum space and apply the matching formula in eq. (2.10) to obtain the physical

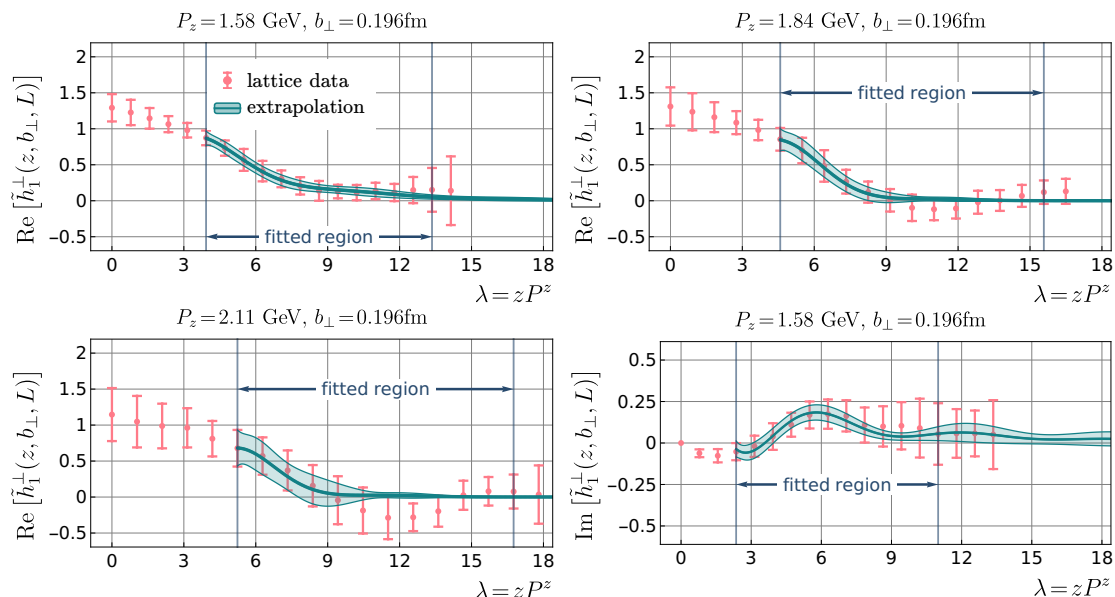


Figure 9. Illustration of extrapolation for various (P^z, b_\perp) . The scattered points and error bars are central values and errors of the renormalized quasi-TMDPDF \tilde{h}_1^+ . The bands denote the extrapolations based on eq. (3.7). The regions used for the fitting are also indicated.

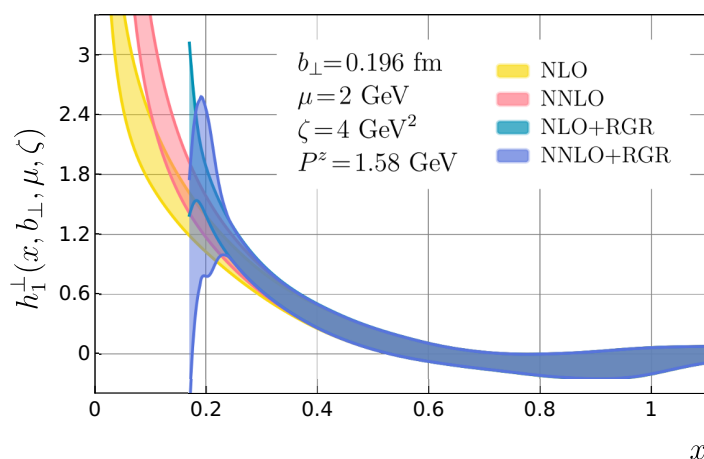


Figure 10. Illustration of the Boer-Mulders function by implementing the NLO, NLO+RGR, NNLO and NNLO+RGR correction, taking $b_\perp = 0.196$ fm, $P^z = 1.58$ GeV as an example.

Boer-Mulders TMDPDF. The intrinsic soft function and Collins-Soper kernel have been calculated for X650 in [22]. Here we adopt these results in our matching procedure.

In figure 10, we show the effects of perturbative matching in NLO, NLO+RGR, NNLO and NNLO+RGR, taking the $b_\perp = 0.196$ fm, $P^z = 1.58$ GeV case as an example. For the results with RGR improvement, we cut the curves at small x where $2xP^z \lesssim \Lambda_{\text{QCD}}$. As can be seen from the figure, at $x \gtrsim 0.2$ the results using the NLO, NLO+RGR, NNLO, NNLO+RGR matching are consistent with each other.

In figure 11, we show the momentum dependence of physical Boer-Mulders functions at different b_\perp values, where we also show the negative x region which corresponds to the

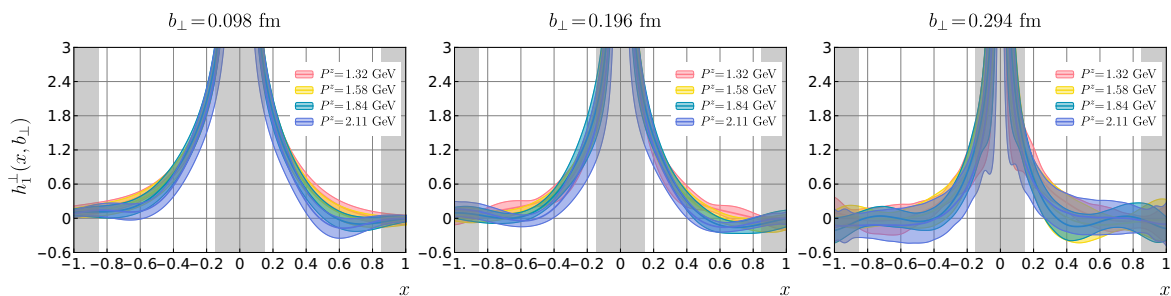


Figure 11. Comparison of physical Boer-Mulders TMDPDFs at different P^z and b_\perp , where the NNLO matching has been implemented. Both statistical and systematic errors have been included. Shaded regions are estimated unreliable regions where power corrections also become important.

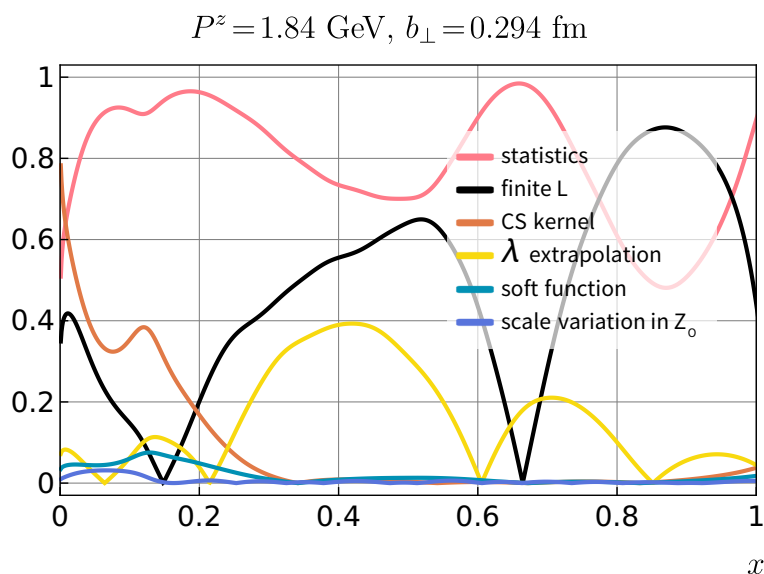


Figure 12. Different sources of uncertainties, taking $P^z = 1.84$ GeV and $b_\perp = 0.294$ fm as an example.

antiquark distribution. Note that a slight asymmetry can be observed under $x \rightarrow -x$, in contrast with the pion Boer-Mulders function which is symmetric under $x \rightarrow -x$ due to isospin symmetry [18]. Given that the nucleon momenta in this calculation are not significantly larger than the nucleon mass, we do not expect a reliable extrapolation to infinite momentum. Therefore, we choose to present the results at different momenta rather than performing an infinite momentum extrapolation. From figure 11, we can see that the bands show a reasonable convergence behavior with increasing momentum. We only show the results up to $b_\perp = 0.294$ fm, as it becomes difficult to control the uncertainties beyond that. The shaded regions are unreliable regions estimated from the breakdown of RGR. In these regions, power corrections in the factorization formula also become important and have to be taken into account.

The error bands in figure 11 include both statistical and systematic uncertainties, where the latter are estimated from uncertainties of the intrinsic soft function, Collins-Soper kernel, extrapolation by shifting the fitting region by a , and scale variation in RGR. In figure 12,

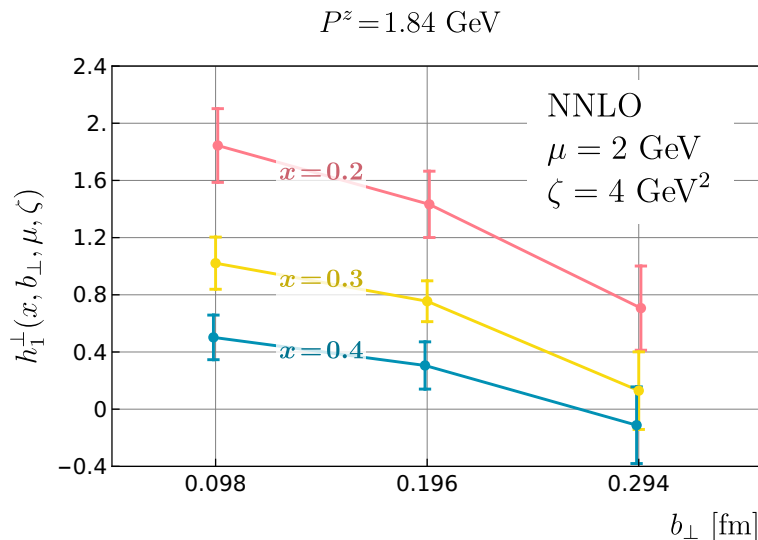


Figure 13. Dependence of the nucleon Boer-Mulders function on b_{\perp} , taking $P^z = 1.84$ GeV, $x = 0.2$, 0.3 and 0.4 as example.

we plot the ratio of each uncertainty and the combined uncertainty, taking the data at $P^z = 1.84$ GeV and $b_{\perp} = 0.294$ fm as an example.

In figure 13, we show the dependence of the nucleon Boer-Mulders function on b_{\perp} . A decay trend can be observed with increasing b_{\perp} . This is in contrast with the unpolarized quark TMDPDF result in ref. [17], where the decay with b_{\perp} is not obvious. This suggests that the Boer-Mulders function might receive smaller higher-twist contributions from the unpolarized quark TMDPDF, and keeping these contributions under control will be crucial in extracting a reliable result for the TMDPDF.

4 Comparison with pion Boer-Mulders function

In ref. [18], we have calculated the Boer-Mulders TMDPDF of the pion on the same ensemble X650. In figure 14, we show a comparison plot between the results for the nucleon and pion. Since no infinite momentum extrapolation has been performed in the present work due to poor signal-to-noise ratios at large nucleon momenta, our comparison is based on lattice data at the same P^z and b_{\perp} . We find that the nucleon and pion Boer-Mulders functions share the same sign, as discussed in ref. [46]. Moreover, the function decreases more rapidly with the momentum fraction x for the nucleon than for the pion. In addition, the Boer-Mulders function exhibits an asymmetry between quark and antiquark combinations for the nucleon, but it is symmetric for the pion due to isospin symmetry. Drawing further quantitative conclusions would require a detailed study of the continuum and infinite momentum limits of the nucleon Boer-Mulders function, which we will address in future publications.

5 Summary and outlook

In summary, we have presented an exploratory study of the nucleon Boer-Mulders TMDPDF using LaMET. The calculation was done on ensemble X650 with a single lattice spacing

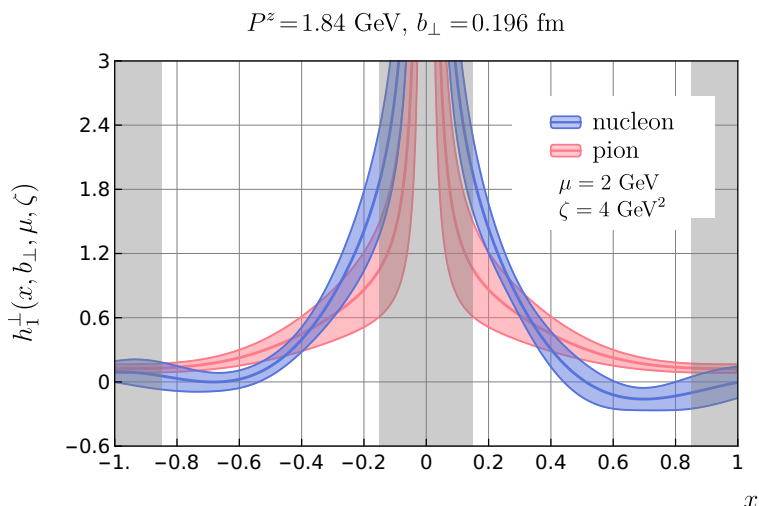


Figure 14. Comparison of the Boer-Mulders functions for the nucleon and pion, taking $P^z = 1.84$ GeV, $b_\perp = 0.196$ fm as an example. Shaded regions are estimated unreliable regions where power corrections also become important.

and unphysical pion mass. Our results provide preliminary insights into the Boer-Mulders function. For example, they show a decay behavior with increasing transverse separation b_\perp , and have the same sign for the nucleon and pion, with the former decaying more rapidly with the momentum fraction x than the latter. More definitive conclusions require further investigation. However, following our work, it should be relatively straightforward to identify the detailed properties of the Boer-Mulders function.

In the future, we plan to improve our calculation in the following aspects. We aim to explore the continuum, infinite momentum as well as the physical limits by collecting data on other ensembles with different lattice spacings and pion masses. To achieve this, we also need to calculate the intrinsic soft function and the Collins-Soper kernel on such ensembles. Once this is done, we will be ready to compare our lattice results with phenomenological results obtained from fitting to experimental data.

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Code Availability Statement. This article has no associated code or the code will not be deposited.

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