



The locality of the fourth root of staggered fermion determinant in the interacting case.

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The fourth root approximation in LQCD simulations with dynamical staggered fermions requires justification. We test its validity numerically in the interacting theory in a renormalization group framework.

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1. Introduction

Numerical simulations of dynamical fermions within the framework of the staggered formalism are both computationally cost effective and phenomenologically successful. Exploiting the advantageous properties of an improved staggered fermion formulation, various collaborations are performing high-precision lattice QCD calculations that are in excellent agreement with experimentally known measurements [1]. However, this success is clouded by the long-standing problematic issue of the validity of the fourth root approximation. The staggered fermion describes, in fact, four tastes on the lattice, so in order to study QCD with $N_f = 2$ or $N_f = 2 + 1$, a Boltzmann weighting is used that contains a fractional power of the fermion determinant. Because of tastebreaking at nonzero lattice spacing, taking the fractional power of the staggered determinant before restoring taste symmetry is conceptually nontrivial. The issue whether the fourth root prescription gives a lattice theory in the right universality class to reproduce QCD is yet unresolved and the phenomenological success of the staggered formulation has heated up the discussion. In the past years various numerical investigations [2] have addressed this concern. Recently a direct approach was adopted towards establishing whether the universality class is the right one: if a single-taste local fermion action can be found, whose determinant is equal to the fourth root of the staggered fermion determinant up to cutoff effects

$$\lim_{a \to 0} (\det D_{st})^{1/4} = \det D \det H \tag{1.1}$$

(where D is a local Dirac operator and H is local and contains only cutoff effects) then the fourth root prescription can be consistently fitted into the framework of a local field theory. In the free theory numerical [3] and analytical studies [4, 5] showed that such a local operator exists. In particular, in [5] Shamir applied renormalization group blocking to the free staggered operator D_0 in the spin \otimes taste representation. After n blocking steps Q_n , the fermionic degrees of freedom live on the coarse lattice with lattice spacing $a_c = 2^n a$ and the determinant of the staggered operator decomposes as det $D_{st}(a = 2^{-n}a_c) = \det D_n \det G_n^{-1}$. More details can be found in [5], where it was proved analytically that the blocked propagator $D_n^{-1} = \alpha^{-1} + Q_n D_0^{-1} Q_n^{\dagger}$ factorizes, in the limit $n \to \infty$, as $D_{\infty}^{-1} = D_{\text{Tg}} \otimes I$ and that $G_n^{-1} = D_0 + \alpha Q_n^{\dagger} Q_n$ is a local operator that contains only cutoff effects. This completes the proof that in the continuum limit, *i.e.* $a(=2^{-n}a_c) \to 0$, the decompositions of Eq. (1.1) holds in the free case.

2. Renormalization group transformation in the interacting case

In this section we briefly discuss the renormalization group (RG) program adopted to test the interacting case. The new complication is that a mapping from the one-component staggered fermion basis to the spin \otimes taste basis is not unique. We define the interacting theory in the one-component formalism. The first fermion RG blocking transformation is used to perform the transition to the spin \otimes taste representation:

$$(Q_{(1)}\psi)^{a\alpha}(2x) = \frac{1}{2} \sum_{r_{\mu}=0,1} (\gamma_1^{r_1} \gamma_2^{r_2} \gamma_3^{r_3} \gamma_4^{r_4})^{(a\alpha)} W_r(2x, 2x+r) \psi(2x+r), \tag{2.1}$$

where a and α are Dirac spin and taste indices respectively. We require the parallel transporters $W_r(2x,2x+r)$ to be a sum over all the shortest paths ¹ between the origin 2x and the other sixteen sites 2x + r of the hypercube. The sum over different paths is thought to reduce the breaking of hypercubic symmetry; however, the choice to parallel transport all the points within a given hypercube to the corner r = 0 unavoidably introduces some asymmetries. The subsequent RG steps are a covariant generalization of the arithmetic mean on a 2^4 hypercube used in [5]:

$$(Q_{(n)}\psi)(2x) = 2^{-4} \sum_{r_u=0,1} W_r(2x, 2x+r)\psi(2x+r) \qquad n > 1.$$
 (2.2)

The shortest paths that compose the parallel transporters $W_r(2x,2x+r)$ in Eq. (2.2) are constructed from links of the blocked lattice (with lattice spacing $2^n a$). These blocked links are built following a program suggested in [7]. In detail, the links that live on the fine lattice (with lattice spacing $2^{n-1} a$) are twice APE smeared and projected back to SU(3). The blocked links are then built multiplying two of these fine smeared links in line. Once the blocking kernels are defined we project the nondiagonal piece of the blocked propagator D_n^{-1} onto the spin \otimes taste Clifford space:

$$(Q_n D_0^{-1} Q_n^{\dagger})(x, y = 0) = \sum_{S, T} \Gamma_S \otimes \Gamma_T^{\dagger} M_{ST}(x, y = 0).$$
 (2.3)

Here x and y=0 2 live on the coarsest lattice and Q_n can be regarded as a 'big' blocking step $Q_n=Q_{(n)}Q_{(n-1)}\cdots Q_{(1)}$ that transforms the original lattice with lattice spacing a into the blocked lattice with lattice spacing $a_c=2^na$. The coefficients $M_{ST}(x,y=0)$ of the projection in Eq. (2.3) are evaluated numerically and averaged over two different lattice ensembles. We fixed Lorentz gauge before doing any RG blocking. Parameters of the simulations are shown in Table 1. Applying a different number of RG steps, n, on each of these two ensembles, the resulting blocked lattices have the same coarse-lattice spacing of $a_c=0.72$ fm. This allowed us to see how the coefficients M_{ST} scale with one additional RG blocking step.

unblocked lattice	$16^{3} \times 48$	$40^3 \times 96$
а	0.18 fm	0.09 fm
m_s a	0.125	0.05
number of RGT's	2	3
blocked lattice	$4^3 \times 12$	$5^3 \times 12$
a_c	0.72 fm	0.72 fm
number of cfgs	148	56

Table 1: Simulation parameters

As a consistency check on the blocking procedure described above, we evaluated the mass of the $(s\overline{s})$ meson. In fact, the spectrum on the blocked lattice is expected to be identical to the spectrum in the original theory. The meson mass can be evaluated directly from the coefficients $M_{ST}(x,y)$ of Eq. (2.3):

¹These are built using APE smeared links of the original lattice to be consistent with the RG program adopted for the subsequent RG steps.

²The fermion source is defined at the origin.

$$C_{s\overline{s}}^{T}(t) = \sum_{S,T',\vec{x}} \langle \operatorname{Tr}(\tau_{5}\Gamma_{T}\Gamma_{T'}\Gamma_{T}\tau_{5}\Gamma_{T'}^{\dagger}) |M_{ST'}(t,\vec{x})|^{2} \rangle, \tag{2.4}$$

where the Γ_S 's and the $\Gamma_{T,T'}$'s are the 16 Dirac matrices in the normal and adjoint $(\tau_{\mu} = \gamma_{\mu}^{\star})$ representation respectively. In Figure 1 we present the taste splittings for the meson masses evaluated on the two blocked lattices. At this level of statistics, the decrease in splitting from n=2 to n=3 is consistent with the expected $\mathcal{O}(a^2\alpha)$ or $\mathcal{O}(a^2\alpha^2)$, but the preliminary RG blocked splittings are systematically higher than values obtained in direct measurements at much lighter valence quark masses. Further study is needed.

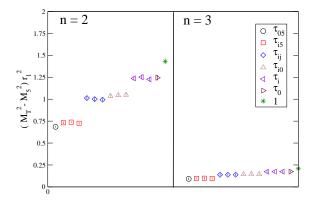


Figure 1: Pseudoscalar taste splittings after 2 and 3 RG blockings in units of r_1 , the distance at which the static quark potential satisfies $r^2 dV/dr = 1$.

3. Preliminary results

In this preliminary study we have analyzed for various displacements |x - y| the dominant terms in Eq. (2.3) that are also present in the free theory:

$$\sum_{\mu} (\gamma_{\mu} \otimes I) A_{\mu}^{(n)}(x,y) + (\gamma_{5} \otimes \tau_{5} \tau_{\mu}) B_{\mu}^{(n)}(x,y) + \sum_{\nu \neq \mu} (\gamma_{5} \otimes \tau_{5} \tau_{\nu}) B_{\nu}^{\prime (n)}(x,y) + (I \otimes I) C^{(n)}(x,y) + \cdots$$
(3.1)

Comparing Figures 2 and 3, it is remarkable to see that increasing by only one unit the number of RG blocking transformations the mass term $(I \otimes I)$ (\square) becomes larger than the taste breaking term $\gamma_5 \otimes \tau_5 \tau_i$ (\diamondsuit) for any direction i for displacement |x-y|=1. This is in agreement with what we expect from the free theory: in the limit $n \to \infty$ the only terms in M_{ST} that should survive are diagonal in taste space.

Other terms in Eq. (3.1) are seen in our simulations. Their contribution is 10 - 1000 times smaller than the leading kinetic term $A_{\mu}^{(n)}$ in Eq. (3.1) and this make it difficult to identify the terms that are statistically relevant. We need to increase the number of configurations to have a better understanding of the statistics. Besides this issue, it is crucial to know how these terms scale with n to determine whether the blocked propagator becomes diagonal in taste space in the interacting theory as it does in the free theory. However an element that can play an important role in this scaling is the breaking of the hypercubic symmetry, caused by the introduction of the parallel transporters, as it has been briefly discussed in the previous section. In fact, Figures 2

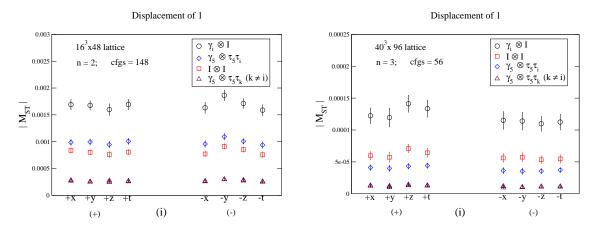


Figure 2: M_{ST} after n = 2 RG steps.

Figure 3: M_{ST} after n = 3 RG steps.

and 3 suggest that the blocking procedure is introducing some hypercubic asymmetries, since the coefficients $M_{ST}(x,y)$ are not perfectly degenerate under reflections about the principal axes. In order to investigate further these terms, present only in the interacting theory, it is necessary either to quantify the hypercubic symmetry breaking or to redefine the RG blocking transformations in order to guarantee hypercubic invariance (one suggestion is presented at this conference [6]).

4. Interacting theory vs free theory

In [5] the scaling properties of the taste breaking terms $B_{\mu}^{(n)}(p)$ were evaluated analytically. It was shown in momentum space that they scale like 2^{-n} , uniformly in p, so the blocked propagator becomes diagonal in taste space, when $n \to \infty$. Postponing the issue of the breaking of hypercubic symmetry to a later study, we consider here a linear combination of the amplitudes $|M_{ST}(x,y)|$ under reflections and rotations about the principal axes. In Figures 4 and 5 we show how the ratios between the taste violating terms and the leading kinetic term $A^{(n)}$ scale with n.

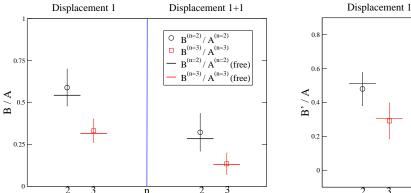


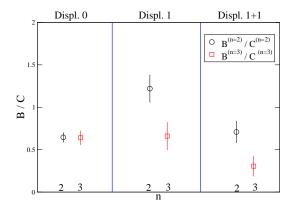
Figure 4: Taste breaking term $B^{(n)}$ over the kinetic term.

Displacement 1+1

Figure 5: Taste breaking term $B^{\prime(n)}$ over the kinetic term.

Our data show that they diminish when n is increased and that their magnitudes are also remarkably in agreement with the free theory represented by solid lines. In Figures 6 and 7 we show

the scaling properties of the taste breaking term divided by the mass term. The interacting theory agrees with the free theory for displacements |x-y|=1,2, however a discrepancy is seen at zero displacement where the interacting theory does not scale as expected. The zero displacement is not interesting when discussing locality thus this behavior, even if not well understood, does not spoil the good scaling properties we are seeing in the interacting theory.



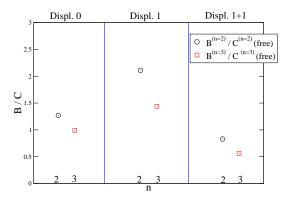


Figure 6: Taste breaking term divided by the mass term: interacting theory

Figure 7: Taste breaking term divided by the mass term: free theory.

The conclusion that we can draw from this preliminary study is that the blocked propagator in the interacting case is dominated by the same terms present in the free theory and that these terms scale as expected to make the decomposition of Eq. (1.1) possible. The statistical relevance and the scaling properties of other terms seen in the interacting theory is still under investigation.

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