# Adams and Hoelbling fermions: numerical properties

## Philippe de Forcrand ETH Zürich & CERN & YITP

### with Aleksi Kurkela (McGill) & Marco Panero (Helsinki)

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#### ΞTH

Eidgenössische Technische Hochschule Zürich Swiss Federal Institute of Technology Zurich

## Motivation

- Light u, d quarks needed to simulate correct physics  $\rightarrow$  expensive
- Cost-saving: staggered fermions (1/4 d.o.f.)

$$S_{F} = \sum_{x} \bar{\chi}(x) \sum_{\mu} \eta_{\mu}(x) (U_{\mu}(x)\chi(x+\hat{\mu}) - U_{\mu}^{\dagger}(x-\hat{\mu})\chi(x-\hat{\mu})) + m_{q} \sum_{x} \bar{\chi}(x)\chi(x)$$

 $\eta_{\mu}=\pm 1;\,\{\gamma_{\mu},\gamma_{\nu}\}=2\delta_{\mu\nu}~\rightarrow~\prod_{4}\eta=-1$  around any plaquette

• Drawback:  $N_f = 4$  (degenerate when a = 0) "tastes"  $\rightarrow \sqrt{\det(D_{st})}$ "rooting is evil" Mike Creutz

- non-locality?
- 't Hooft vertex,  $U(1)_A$  breaking?
- staggered fermions "don't feel the topology"
- No quartet of low-lying eigenvalues  $\leftrightarrow$  no index theorem
- Adams to the rescue:  $N_f = 2$  staggered overlap fermions

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• Hoelbling: N<sub>f</sub> = 1
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## Construction

Two ingredients:

- 1: taste-dependent mass term (cf. Wilson term)
  - $\implies$  split tastes into several branches fine-tune for massless limit
- 2: plug into overlap (cf. Neuberger)
  - $\implies$  eliminate fine-tuning, etc.. large computational overhead

Look at 1 then 2

## Possible mass terms

# Golterman & Smit (1984)

1	<mark>0</mark> -link
$\gamma_{\mu}$	1-link
$\gamma_{\mu}\gamma_{\nu}$	<mark>2</mark> -link
γμγνγρ	<mark>3</mark> -link
Y1 Y2 Y3 Y4	4-link

• 1-link and 3-link do not satisfy  $\gamma_5$ -hermiticity  $\rightarrow$  determinant *complex* 

- can be fixed with 2 conjugate copies, ie. doubling  $N_f$
- 1-link: cancellations between kinetic & mass terms free spectrum on a circle!

0-link = KS; 2-link = Hoelbling; 4-link = Adams

Notation: 0-link  $\rightarrow$  *m*, 2-link  $\rightarrow$  *M*<sub>Hoelbling</sub>, 4-link  $\rightarrow$  *M*<sub>Adams</sub>  $\propto$   $\Gamma_{55}\Gamma_5$ 

• NB. Also can multiply above mass terms by  $\Gamma_{55} \equiv (-)^{\chi}$  ?

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### Free spectra and taste reduction I



## Free spectra and taste reduction II

• Select one branch by *x*-shift, ie. additive mass renormalization (cf. Wilson) Other branch(es) have masses O(1/a), ie. doublers



 $\Rightarrow$  can trade  $\Gamma_5 = \gamma_5 \otimes \mathbf{1} + \mathcal{O}(a)$  for  $\Gamma_{55} = \gamma_5 \otimes \gamma_5$  *exactly*  $\rightarrow$  index theorem

## Index from eigenvalue flow

- Index from flow of eigenvalues  $\lambda(m)$  of  $H(m) = \gamma_5(\not D + m)$ 
  - ullet ( $ot\!\!\!/ + m_0)$  has zero-mode  $|\Psi_0
    angle \, otherefore \,$  eigenvalue  $\lambda(m_0)=$  0 for  $H(m_0)$
  - Perturb *m* away from  $m_0 \rightarrow \text{eigenvalue displaced by } \langle \Psi_0 | \gamma_5(m-m_0) | \Psi_0 \rangle$  $\lambda(m) = \pm (m-m_0) \text{ (ie. crossing)}$  **IF**  $\langle \Psi_0 | \gamma_5 | \Psi_0 \rangle = \pm 1, \text{ ie. } | \Psi_0 \rangle \text{ chiral}$
- Alternative: can also consider flow for  $\hat{H}(\rho) = \Gamma_{55}(\not \! D_{st} + \rho M_{Adams})$

Adams' original proposal

•  $\mathcal{D}_{st}$  has near-chiral, near-zero modes •  $M_{Adams} = \Gamma_{55}\Gamma_5 \rightarrow \text{perturbation } \rho\Gamma_5 \approx \rho\gamma_5 \otimes \mathbf{1}$ 

"Classic" versus "original" ?

# Comparing two eigenvalue flows

### Adams' original



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## Index theorem

• Cold configuration: agreement with analytic result



### • Cooled Q = 1 instanton: $N_f \times Q$ crossings



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# Effects of gauge field fluctuations

### • $\beta = 6.0$ , Q = 1: eigenvalue gap closes, esp. Adams



# Effects of gauge field fluctuations

- The width of the spectrum fluctuates (shrinks)
  - $\rightarrow$  fine-tuning for massless quarks
- The gap in the spectrum fills up
  - $\rightarrow\,$  distinction between light modes and doublers blurred

# Eigenvalue spectra: Adams vs Hoelbling, 8<sup>4</sup>



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### Usefulness

Superior robustness of Hoelbling fermions: 2-link vs 4-link transporters

Can always smear links to suppress fluctuations: Thanks Stephan Dürr



### Pion mass

- Compute quark propagator G(x, y, z, t) from point source:
  - $16^3 \times 32$  lattices,  $\beta = 6.0$  quenched
  - Form  $\vec{p} = \vec{0}$  meson correlator (connected diagram only)  $C(t) = \sum_{xyz} G(x, y, z, t) \Gamma_{55} G(x, y, z, t)^{\dagger} \Gamma_{55} = \sum_{xyz} |G(x, y, z, t)|^2$
  - Look for effective mass plateau  $\rightarrow$  (lightest) pion mass ( $am_{\pi}$ )
- Monitor  $(am_{\pi})^2$  vs  $(am_q)$ : PCAC + additive mass renormalization

## **Pion correlator**



## Pion mass vs m<sub>a</sub>



### Pion correlator Hoelbling



## Overlap construction

• Just like Neuberger: 
$$D_{sov} = 1 + \gamma_5^{\prime\prime} \operatorname{sign}(H(-m_0))$$

with " $\gamma_5''=\Gamma_{55}=(-)^{x+y+z+t}$  (need " $\gamma_5''^2=1)$ 

- Potential advantages:
  - cheaper (4 times fewer d.o.f. per site)
  - more robust ?

And reduces  $N_f = 4$  to  $N_f = 2$  tastes without fine-tuning

### N.B. mo is really p in Adams' proposal [no mass shift]

# Free field: $U_{\mu}(x) = \mathbf{1} \ \forall \mathbf{x}, \mu$

Spectrum of kernel:  $\gamma_5 H_W(m_0 = -1)$  and  $\gamma_5 H_{Adams}(m_0 = -1)$ 



 $\gamma_5 \text{sign}(H) = \frac{D}{\sqrt{D^{\dagger}D}}$  projects eigenvalues of  $D = \gamma_5 H$  on unit circle

Adams: two low-*p* eigenmodes projected to -1, two projected to  $+1 \Rightarrow N_f = 2$ 



Adams comparable to Neuberger although kernel less local (4-link)



Adams comparable to Neuberger although kernel less local (4-link)

# Cost of applying operator

- Multiplication by D: about 2 times faster for Adams (no Dirac indices)
- Sign(*H*) [using CG, no deflation]:
  - about 8 times faster for Adams on easy cases
  - about 2-3 times faster on hard cases

Can optimize  $m_0$  and  $\rho$  in Adams' operator: not exploited yet

#### Also:

improved kinetic operator, link smearing (kinetic and/or mass), deflation, preconditioning, ...

## Cost of inversion: compare with Neuberger

Apples with apples:

- same gauge field (12<sup>4</sup>,  $\beta$  = 6.0)
- same basic algorithm (CG inner, CG outer)



Net CPU gain: factor 2-3 over Neuberger...

### Cost of inversion: compare with Neuberger

#### Adams versus Neuberger



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## Conclusions

- Index theorem, overlap etc.. work as advertised
- Compare taste-dependent mass terms *exhaustively* first then try overlap
- Here: 2-link (Hoelbling) more robust than 4-link (Adams)
- 2-link  $\rightarrow N_f = 2$  without fine-tuning ?
- Other possibilities not yet considered?
   eg. start with 8 tastes, split 1-3-3-1

## Sign problem?

# Thanks Mike, Tatsu, Taro

Take Q = 1 configuration  $\rightarrow 1+2+1$  real eigenvalues. Vary  $m_0 : \bigotimes$ 

