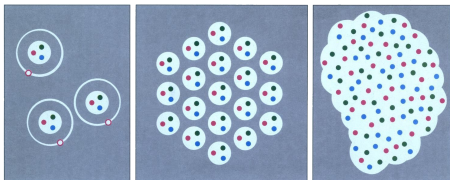


2 Flavor Formulation for Strong Coupling LQCD

Wolfgang Unger, ETH Zürich
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SEWM 2012, Swansea

10.07.2012



ETH

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Swiss Federal Institute of Technology Zurich

1 Motivation for Strong Coupling LQCD in Continuous Time

- Continuous Time Limit and $a/a_t = f(\gamma)$
- Continuous Time Partition Function $Z(\beta)$

2 Hamiltonian Formulation

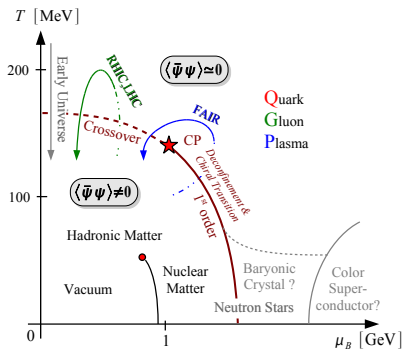
- Spin Representation
- Stochastic Series Expansion

3 Application: 2 flavor SC-LQCD

- Generalization of Spin Representation
- Preliminary Results

Why Strong Coupling Lattice QCD?

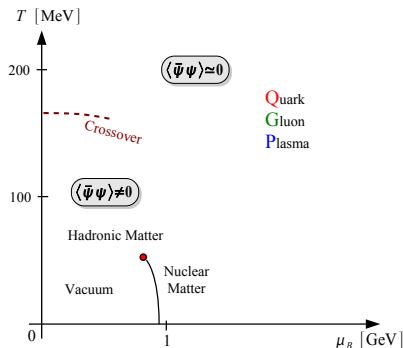
The QCD (μ , T) **phase diagram**:



- rich phase structure conjectured
- chiral and deconfinement transition
- QGP at high temperatures
- exotic matter at high density

Why Strong Coupling Lattice QCD?

The QCD (μ , T) **phase diagram**:



- because of the sign problem: very little is known
- so far: agreement on crossover temperature T_c at zero density

QCD has a severe **sign problem** for non-zero chemical potential $\mu = \frac{1}{3}\mu_B$:

- fermions anti-commute:

$$\gamma_5(i\not{p} + m + \mu\gamma_0)\gamma_5 = (i\not{p} + m + \mu\gamma_0)^\dagger$$

- the **fermion determinant** $\det M(\mu)$ becomes **complex!**

$$e^{-S_f} = \det M(\mu) = \overline{\det M(-\bar{\mu})}$$

- little hope that it can be circumvented:
 - Taylor expansion,
 - imaginary μ with analytic continuation,
 - reweighting method
 are all **limited to small** $\mu/T \lesssim 1$

Why Strong Coupling Lattice QCD?

Look at Lattice QCD in a regime where the **sign problem** can be made mild:

$$\text{Strong Coupling Limit: } \beta = \frac{2N_c}{g^2} \rightarrow 0$$

- allows to integrate out the gauge fields completely, as **link integration factorizes**
 \Rightarrow no fermion determinant
- drawback: strong coupling limit is converse to asymptotic freedom, lattice is maximally coarse

Strong coupling LQCD shares important features with QCD:

- exhibits **confinement**, i.e. only color singlet degrees of freedom survive:
 - **mesons** (represented by monomers and dimers)
 - **baryons** (represented by oriented self-avoiding loops)
- and **spontaneous chiral symmetry breaking/restoration**: (restored at T_c)
 \Rightarrow SC-LQCD is a great laboratory to study the full (μ, T) **phase diagram**

SC-LQCD is a useful toy model for **nuclear matter**

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SC-LQCD is a useful toy model for **nuclear matter**

SC-LQCD is a **1-parameter deformation of QCD**

SC-LQCD at finite temperature

How to vary the temperature?

- $aT = 1/N_\tau$ is discrete with N_τ even
- $aT_c \simeq 1.5$, i.e. $N_\tau^c < 2 \Rightarrow$ we cannot address the phase transition!

Solution: introduce an **anisotropy** γ in the Dirac couplings:

$$\mathcal{Z}(m_q, \mu, \gamma, N_\tau) = \sum_{\{k, n, \ell\}} \prod_{b=(x, \mu)} \frac{(3 - k_b)!}{3! k_b!} \gamma^{2k_b \delta_{\mu 0}} \prod_x \frac{3!}{n_x!} (2am_q)^{n_x} \prod_\ell w(\ell, \mu)$$

$k_b \in \{0, \dots, N_c\}, n_x \in \{0, \dots, N_c\}$

Should we expect $a/a_\tau = \gamma$, as suggested at weak coupling?

- **No:** meanfield predicts $a/a_\tau = \gamma^2$, since $\gamma_c^2 = N_\tau \frac{(d-1)(N_c+1)(N_c+2)}{6(N_c+3)}$

\Rightarrow sensible, N_τ -independent definition of the temperature:

$$aT \simeq \frac{\gamma^2}{N_\tau}$$

- Moreover, SC-LQCD partition function is a function of γ^2

However: **precise correspondence between a/a_τ and γ^2 not known**

SC-LQCD at finite Temperature and Continuous Time:

Strategy for **unambiguous** answer: the **continuous Euclidean time limit** (CT-limit):

$$N_\tau \rightarrow \infty, \quad \gamma \rightarrow \infty, \quad \gamma^2/N_\tau \equiv aT \text{ fixed}$$

- same as in analytic studies: $a_\tau = 0$, $aT = \beta^{-1} \in \mathbb{R}$

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Several **advantages** of continuous Euclidean time approach:

- ambiguities arising from the functional dependence of observables on the anisotropy parameter will be circumvented, **only one parameter** setting the temperature
- no need to perform the continuum extrapolation $N_\tau \rightarrow \infty$
- allows to estimate critical temperatures more precisely, with a faster algorithm (about 10 times faster than $N_t = 16$ at T_c)
- baryons become static in the CT-limit, the **sign problem is completely absent!**

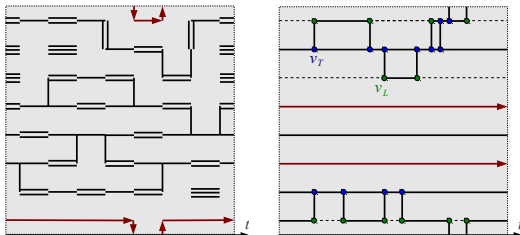
Continuous Time Partition Function, 1 flavor

Partition function in inverse temperature $\beta = 1/aT$ and in the **chiral limit**:

$$\mathcal{Z}(\beta, \mu) = \sum_{\kappa \in 2\mathbb{N}} \frac{(\beta/2)^\kappa}{\kappa!} \sum_{C \in \Gamma_\kappa} v_L^{n_L(C)} v_T^{n_T(C)} e^{3\beta\mu B(C)}, \quad v_L = 1, v_T = 2/\sqrt{3}$$

Typical (2-dimensional) configurations in discrete and continuous time at the same temperature:

- multiple spatial dimer become resolved into **single spatial dimers** as $a_t \rightarrow 0$
- baryons** become **static** in continuous time!



- $\kappa = \frac{1}{2} \sum_x n_L(x) + n_T(x)$ is the **number of spatial dimers**, B is baryon number
- weight of configuration given by number of spatial dimers and **vertices** v_L, v_T **regardless of time coordinates**

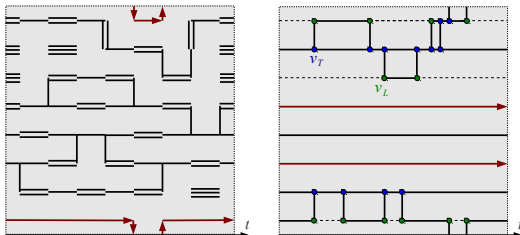
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- sum over all spatial dimer time coordinates $\sim N_\tau/2 \Rightarrow$ **expansion in** $\beta = N_\tau/\gamma^2$
- Γ_κ is the set of equivalence classes of configurations with κ spatial dimers, time coordinates of spatial dimers irrelevant

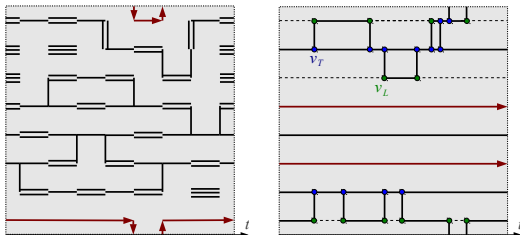
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- each term Γ_κ is represented by a **world line configuration**
- allows to apply QMC techniques: **continuous time worm** (Beard & Wiese), loop cluster algorithm (Evertz et al.), **stochastic series expansion** (Sandvik)

Mapping of 1-flavor $SU(N_c)$ to a spin system

Continuous time methods can be applied to any gauge group $SU(N_c)$:

- baryons become static for $N_c \geq 3$
- mesonic discrete time chains **classified by parity**:

	Discrete Time Chains	Parity Composition	Spin Composition	Example Configuration
U(1)				
U(2)				
U(3)				

⇒ mesonic CT line types **classified by "spin"**: $S = -N_c/2 \dots N_c/2$
 (remnant of staggered even/odd ordering), $\Delta S = \pm 1$ (absorption/emission)

- generalizes to arbitrary $U(N_c)$

Stochastic Series Expansion

Idea: rewrite partition function, based on decomposition in diagonal and non-diagonal elements $\mathcal{H} = \mathcal{H}_1 + \mathcal{H}_2$, truncation L :

$$\mathcal{Z}(\beta) = \text{Tr} \left\{ e^{-\beta \mathcal{H}} \right\} = \sum_{\chi} \sum_{S_L} \frac{\beta^{\kappa} (L - \kappa)!}{L!} \left\langle \chi \left| \prod_{i=1}^L \mathcal{H}_{a_i, b_i} \right| \chi \right\rangle, \quad \begin{aligned} \mathcal{H}_{1,b} &= \varepsilon \mathbb{1}, \quad \varepsilon \geq 0 \\ \mathcal{H}_{2,b} &= \frac{1}{2} S_x^+ S_y^- \end{aligned}$$

with S_L a **time-ordered sequence** of operator-indices: $S_L = [a_1, b_1], [a_2, b_2], \dots, [a_L, b_L]$

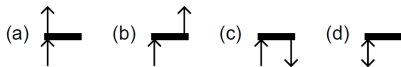
- $a_i = 0$: identity, $a_i = 1, 2$: diagonal/non-diagonal matrix element
- $b_i = \langle x, y \rangle \in Vd$ denotes a bond

Two kinds of **updates**:

1 **changing order in β** , $\kappa \mapsto \kappa \pm 1$:

$$P([1, b]_p \mapsto [0, 0]_p) = \frac{L - \kappa + 1}{Vd \beta \langle \chi | \mathcal{H}_{1,b} | \chi \rangle}, \quad P([0, 0]_p \mapsto [1, b]_p) = \frac{Vd \beta \langle \chi | \mathcal{H}_{1,b} | \chi \rangle}{L - \kappa}$$

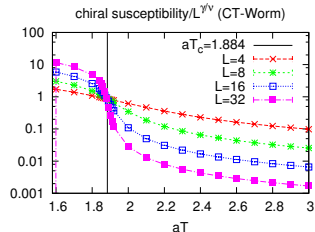
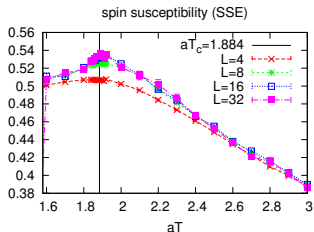
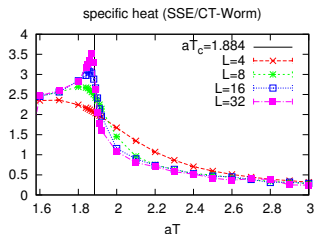
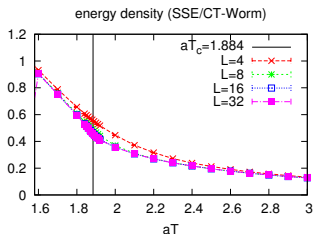
2 **operator loop**: visit bonds b_i successively from an input leg, determine output leg with heatbath probability $\langle \chi_x \chi_y | \mathcal{H}_{a_i b_i} | \chi'_x \chi'_y \rangle$



L is set larger than $\kappa_{\max} \Rightarrow$ **SSE is approximation free** (like CT-Worm)

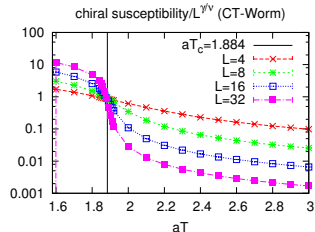
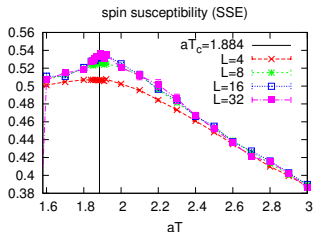
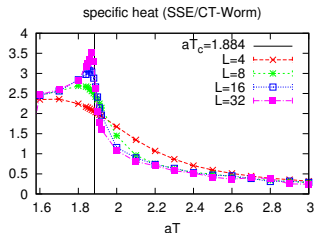
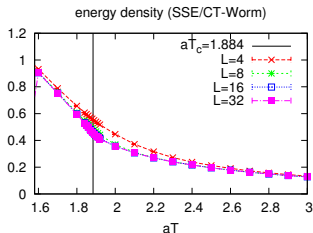
Observables for the Chiral Transition

- CT-Worm: energy density/specific heat, chiral susceptibility
- SSE: energy density/specific heat, “spin” susceptibility



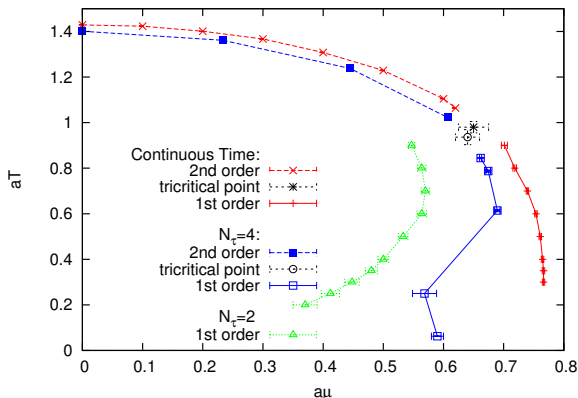
Observables for the Chiral Transition

- CT-Worm: energy density/specific heat, chiral susceptibility
- SSE: energy density/specific heat, **“spin” susceptibility feels chiral transition**



1-flavor SC-QCD Phase Diagram

Comparison of phase diagram in continuous time with $N_\tau = 4$ data (M. Fromm, 2010) via identification $a\mu = \gamma^2 a_\tau \mu$, studied via Worm algorithm [hep-lat/1111.1434]:

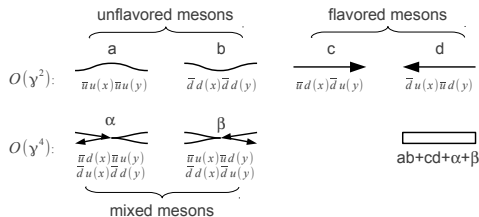


- behavior at low μ agrees well, location of TCP agrees within errors
- no re-entrance seen at small aT (also confirmed by Ohnishi *et al.*, LAT2012)

Application: Generalization of SC-LQCD to 2 chiral flavors!

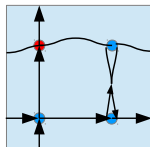
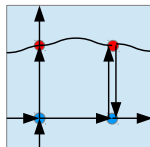
Aim: obtain phase diagram for 2-flavor SC-LQCD, where **pion exchange** may play a crucial role for nuclear transition, but:

- at present, no 2-flavor formulation for staggered SC-LQCD suitable for MC
- already the mesonic sector has a severe **sign problem** (worse than for finite μ HMC)
- 2 new types of mixed dimers give negative sign in mesonic loops already for U(2):



positive weight: 1/4

negative weight: -1/2



Observation in **continuous time** formulation:

- **static lines** for 2 staggered flavors have all **positive weight!**
- again: only single spatial dimers (no α , β spatial dimers)
- Hamiltonian formulation feasible

Continuous Time Transition Rules for $N_f = 2$

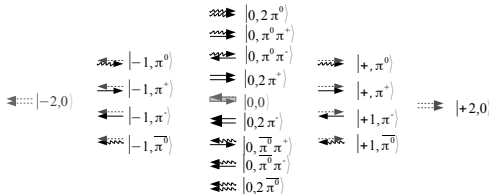
Flavored static lines:

- new classification in terms of

quantum numbers

$|S^z, Q_{\pi^0}, Q_{\pi^+}\rangle$

- $N_c = 2$: in total 19 types of lines, 18 have weight $1/4$ per a_t , "vacuum state" $|0, 0, 0\rangle$ has weight $1/\sqrt{8}$



- state multiplicities = $N_c + 1 + \frac{2}{3} N_c (N_c + 1) (N_c + 2)$, 44 states for $N_c = 3$
- "spin" S^z counts number of emission/absorption events (remnant of even/odd decomposition of lattice for staggered fermions) $S^z = -\frac{1}{2} N_c N_f, \dots, +\frac{1}{2} N_c N_f$
- two "charges" $Q_i = -N_c, \dots, +N_c$ denote the flavor content
- spin/charge conservation:** transitions at spatial dimers, raising charges at one site, lowering at a neighboring site:

$$|\Delta S^z| = 1, \quad |\Delta Q_{\pi^0}| + |\Delta Q_{\pi^+}| = 1$$

Hamiltonian for $N_f = 2$

$$\mathcal{H} = \frac{1}{2} \sum_{\langle x,y \rangle} \left(J_{\pi^0(x)}^+ J_{\pi^0(y)}^- + J_{\bar{\pi}^0(x)}^+ J_{\bar{\pi}^0(y)}^- + J_{\pi^+(x)}^+ J_{\pi^+(y)}^- + J_{\pi^+(x)}^+ J_{\pi^-(y)}^- \right)$$

Absorption ($J_{\pi_i}^+$, lower left triangle) and Emission ($J_{\pi_i}^-$, upper triangle), state vector:

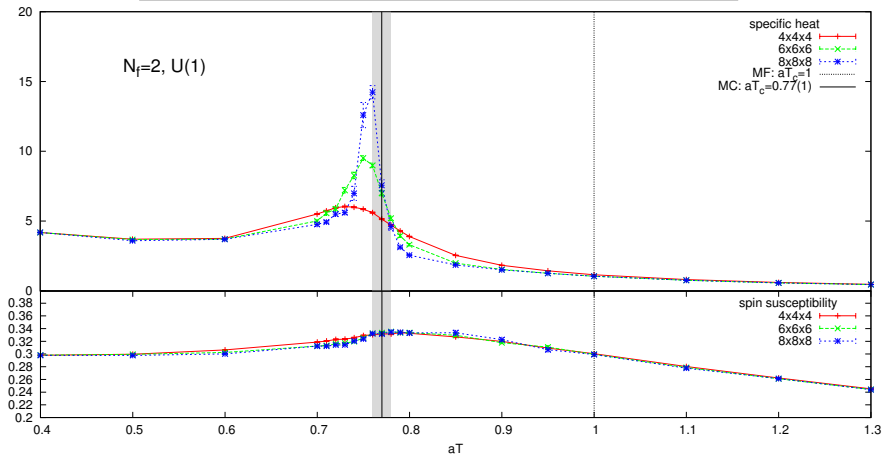
$$J_{\pi_i}^{+/-} = \begin{pmatrix} \begin{array}{c|c|c|c} \pi^0 & \pi^0 & \pi^+ & \pi^+ \\ \pi^0 & & & \\ \pi^+ & & & \\ \pi^- & & & \end{array} & \begin{array}{c|c|c|c} \pi^0 & & & \\ \bar{\pi}^0 & \pi^+ & \bar{\pi}^- & \\ \pi^+ & \bar{\pi}^+ & \bar{\pi}^0 & \\ \pi^- & \bar{\pi}^- & \bar{\pi}^0 & \bar{\pi}^0 \end{array} \\ \hline \begin{array}{c|c|c|c} \pi^0 & & & \\ \bar{\pi}^0 & & & \\ \pi^+ & & & \\ \pi^- & & & \end{array} & \begin{array}{c|c|c|c} \pi^0 & & & \\ \bar{\pi}^0 & \pi^+ & \bar{\pi}^- & \\ \pi^+ & \bar{\pi}^+ & \bar{\pi}^0 & \\ \pi^- & \bar{\pi}^- & \bar{\pi}^0 & \bar{\pi}^0 \end{array} \\ \hline \begin{array}{c|c|c|c} \pi^0 & & & \\ \bar{\pi}^0 & & & \\ \pi^+ & & & \\ \pi^- & & & \end{array} & \begin{array}{c|c|c|c} \pi^0 & \bar{\pi}^- & \bar{\pi}^+ & \\ \pi^0 & & & \\ \pi^- & \pi^+ & \bar{\pi}^0 & \\ \pi^+ & \pi^- & \bar{\pi}^0 & \bar{\pi}^0 \end{array} \\ \hline \begin{array}{c|c|c|c} \pi^0 & & & \\ \bar{\pi}^0 & & & \\ \pi^+ & & & \\ \pi^- & & & \end{array} & \begin{array}{c|c|c|c} \pi^0 & & & \\ \bar{\pi}^0 & & & \\ \pi^+ & & & \\ \pi^- & & & \end{array} \end{pmatrix}, \quad \chi = \begin{pmatrix} -2, 0 \\ -1, \pi^0 \\ -1, \bar{\pi}^0 \\ -1, \pi^+ \\ -1, \pi^- \\ \hline 0, 2\pi^0 \\ 0, 2\bar{\pi}^0 \\ 0, 2\pi^+ \\ 0, 2\pi^- \\ \hline 0, 0 \\ \hline 0, \pi^0 \pi^+ \\ 0, \pi^0 \pi^- \\ 0, \bar{\pi}^0 \pi^+ \\ 0, \bar{\pi}^0 \pi^- \\ \hline +1, \pi^0 \\ +1, \bar{\pi}^0 \\ +1, \pi^+ \\ +1, \pi^- \\ +2, 0 \end{pmatrix}$$

- vertex weights are $v_{\pi_i} = 1$ for vertices not mixing the two charges Q_i and $v_{\hat{\pi}_i} = \frac{1}{\sqrt{2}}$ for vertices mixing the charges

Preliminary Result for 2-flavor SC-LQCD

Comparison of aT_c from MC data with mean field:

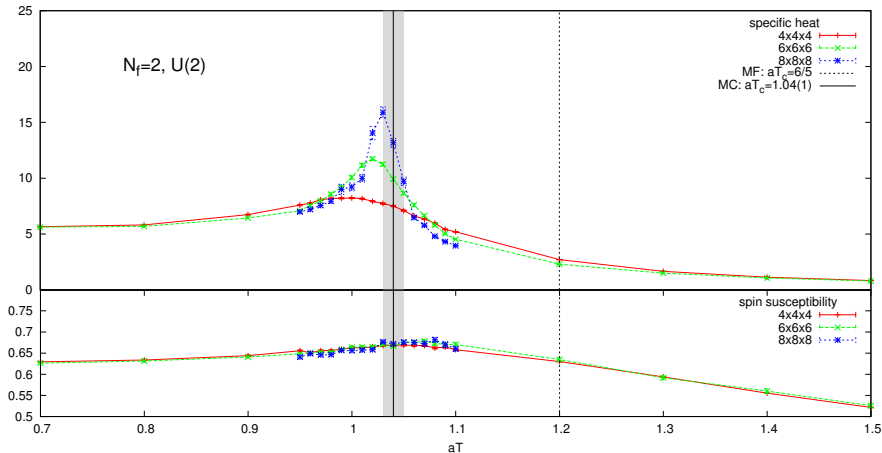
N_f	$N_c = 1$	$N_c = 2$	$N_c = 3$
1	3/2 [1.102(1)]	4/2 [1.467(1)]	5/2 [1.884(1)]
2	5/5 [0.77(1)]	6/5 [1.04(1)]	7/5



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Summary

Achievements:

- CT partition function: new formulation as a **quantum spin system!**
- “spin” formulation and Hamiltonian follow from conservation laws for even/odd chains of time-like dimers and flavors - this **generalizes to arbitrary N_c, N_f**
- new observable: spin susceptibility, sensitive to chiral transition
- quantum Monte Carlo applicable: e.g. continuous time worm or **stochastic series expansion** (most convenient)
- now also applied to U(2) with two flavors (incorporates pion exchange)
- extension to SU(3) with finite baryon chemical potential straightforward (Hamiltonian worked out, but no simulations yet)

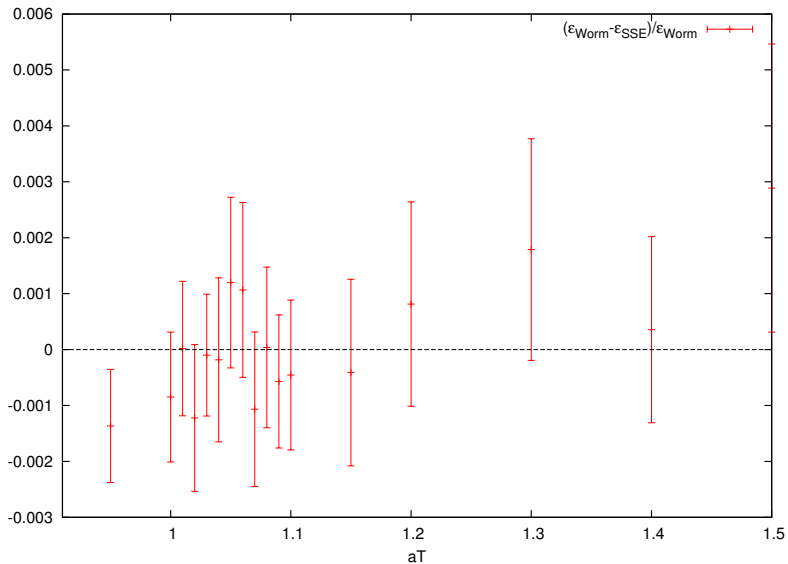
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Soon: obtain full 2-flavor SC-LQCD phase diagram in (μ, T) -plane!

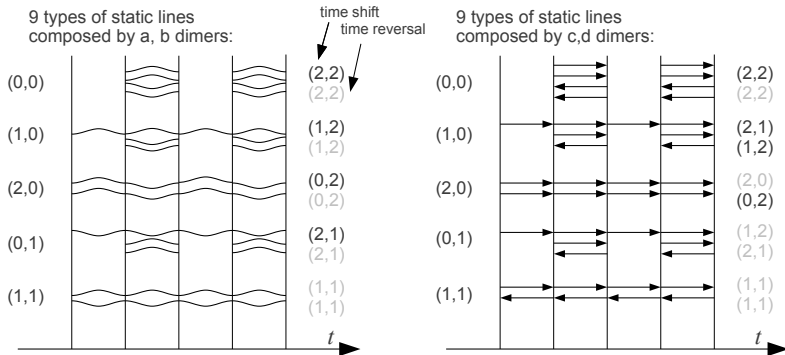
Comparison of SSE and CT-Worm



Static Line Rules

Combine temporal dimers of alternating orders in γ^2 (here for $N_c = 2$):

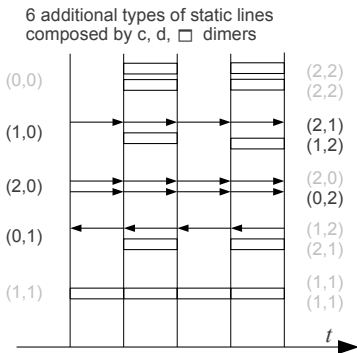
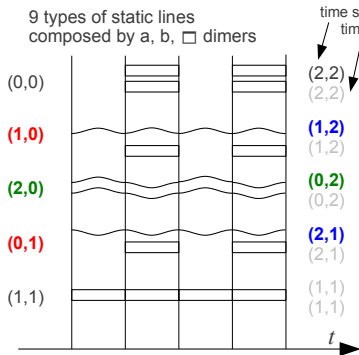
- first: consider (a, b) and (c, d) dimers separately
- then: resum them to obtain flux representation



Static Line Rules

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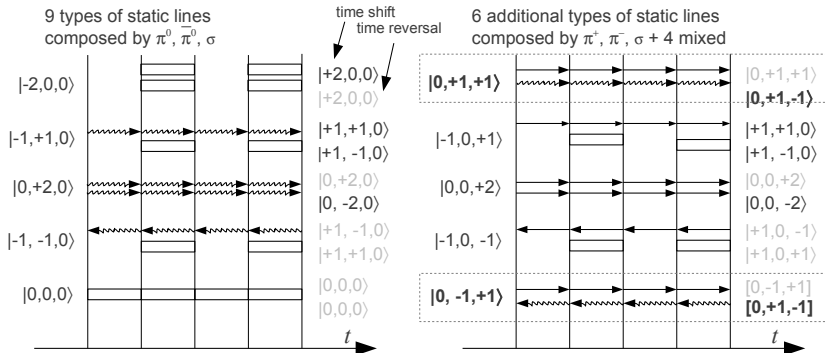
- resummation of $a - b - a$ and $b - a - b$ chains
- resummation of $ab + cd + \alpha + \beta$ into \square dimers



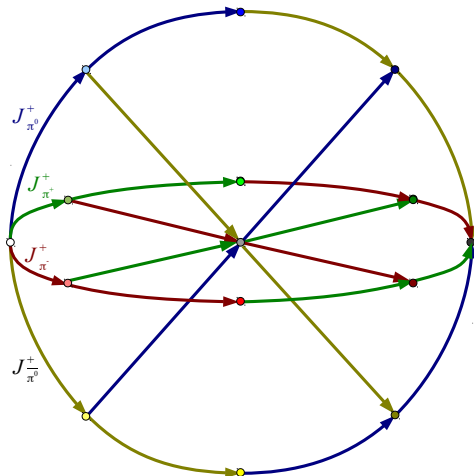
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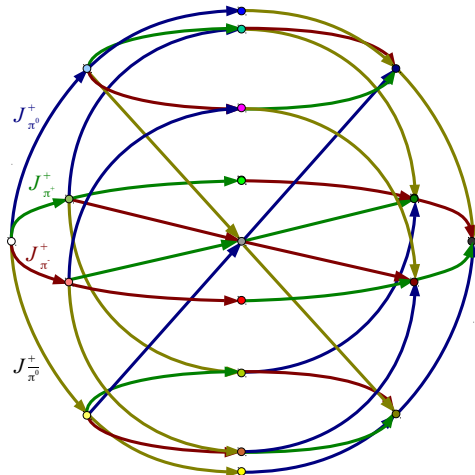
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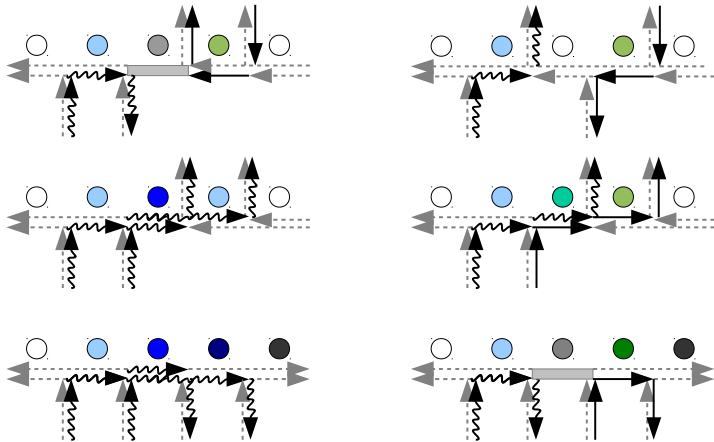
The Transition Rules Encoded in J^\pm



The Transition Rules Encoded in J^\pm



The Transition Rules Encoded in J^\pm



Why Study Strong Coupling QCD on the Lattice?

Two possible scenarios for the relation between SC-LQCD (back) and the (L)QCD phase diagram for four flavors (front):

