

Fermionic fields in the pseudoparticle approach

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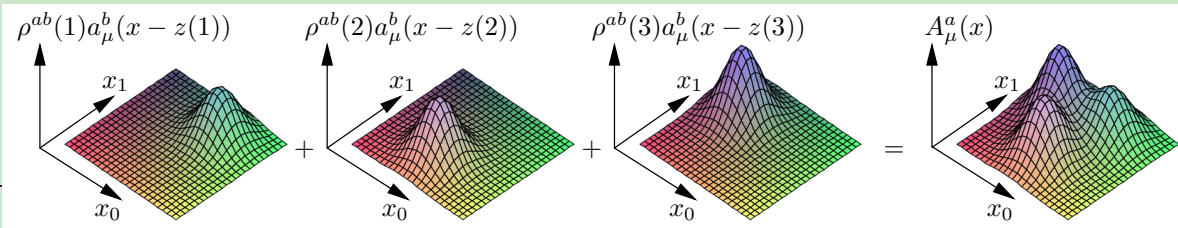
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Introduction (1)

- Models for SU(2) Yang-Mills theory with a small number of degrees of freedom:
 - Regular gauge instanton and meron models (F. Lenz, J. W. Negele, M. Thies, 2003).
 - Pseudoparticle approach (F. Lenz, M.W., 2005).
 - Calorons with non-trivial holonomy (P. Gerhold, E.-M. Ilgenfritz, M. Müller-Preussker, 2006).
 - Ensemble of dyons (D. Diakonov, V. Petrov, 2007).
- Basic principle: restrict the path integral to those gauge field configurations, which can be represented by a linear superposition of a small number of localized building blocks (instantons, merons, akryons, calorons, dyons, ...).



Introduction (2)

- Successes of these models:
 - Linear potential between two static charges at large separations (confinement).
 - Confinement-deconfinement phase transition.
 - String tension, topological susceptibility, critical temperature, ... in qualitative agreement with lattice results.
- Intention: get a better understanding of confining gauge field configurations and the mechanism of confinement.
- **How can fermions be included in such models?**
- In this talk: first steps in this direction.
 - Basic principle of the PP approach in fermionic theories.
 - Testing ground: the 1+1-dimensional Gross-Neveu model in the large- N -limit (phase diagram, chiral condensate).

Basic principle (1)

- Starting point: action and partition function of **any theory with quadratic fermion interaction**:

$$S[\psi, \bar{\psi}, \phi] = \int d^{d+1}x \left(\bar{\psi} Q(\phi) \psi + \mathcal{L}(\phi) \right)$$

$$Z = \int D\psi D\bar{\psi} D\phi e^{-S[\psi, \bar{\psi}, \phi]}$$

(Q : Dirac operator; ϕ : any type and number of bosonic fields, e.g. the non-Abelian gauge field in QCD).

Basic principle (2)

- Consider only those fermionic field configurations, which can be represented by a linear superposition a fixed number of localized building blocks:

$$\psi(x) = \sum_j \underbrace{\eta_j G_j(x)}_{j\text{-th PP}}$$

(η_j : Grassmann valued spinors; G_j : functions, which are localized in space as well as in time, i.e. PPs).

- Define the functional integration over all fermionic field configurations as an integration over the Grassmann valued spinors:

$$\int D\psi D\bar{\psi} \dots = \int \left(\prod_j d\eta_j d\bar{\eta}_j \right) \dots$$

Basic principle (3)

- Integrate out the fermions:

$$S_{\text{effective}}[\phi] = \int d^{d+1}x \mathcal{L}(\phi) - \ln \left(\det \left(\langle G_j | Q | G_{j'} \rangle \right) \right)$$

$$Z \propto \int D\phi e^{-S_{\text{effective}}[\phi]}$$

($\langle G_j | Q | G_{j'} \rangle$ is a finite matrix; **“Q-regularization”**).

- If $\det(Q)$ is real and positive, $\det(Q) = \sqrt{\det(Q^\dagger Q)}$. This suggests another PP regularization:

$$S_{\text{effective}}[\phi] = \int d^{d+1}x \mathcal{L}(\phi) - \frac{1}{2} \ln \left(\det \left(\langle G_j | Q^\dagger Q | G_{j'} \rangle \right) \right)$$

(**“Q[†]Q-regularization”**).

- **The “Q[†]Q-regularization” has significant advantages over the naive “Q-regularization”.**

Q versus $Q^\dagger Q$ (1)

- For the sake of simplicity: consider all PPs G_j to be orthonormal, i.e. $\langle G_j | G_{j'} \rangle = \delta_{jj'}$ (this is not a restriction!).
- The problem of the Q -regularization:
 - **Applying the Dirac operator Q to one of the PPs $G_{j'}$ in general yields a function, which is (partially) outside the PP function space $\text{span}\{G_n\}$, i.e.**

$$QG_{j'}(x) = \sum_k a_{j'k} G_k(x) + h_{j'} H_{j'}(x)$$

($H_{j'}$ normalized, $H_{j'} \perp \text{span}\{G_n\}$).

- If $|\sum_k a_{j'k} G_k| \gg |h_{j'}| \rightarrow$ no problem.
- If $|\sum_k a_{j'k} G_k| \lesssim |h_{j'}| \rightarrow$ when computing the fermionic matrix $\langle G_j | Q | G_{j'} \rangle$, a significant part of $QG_{j'}$ is simply ignored, just because $H_{j'}$ is perpendicular to the PP function space $\text{span}\{G_n\}$.

Q versus $Q^\dagger Q$ (2)

- The advantage of the $Q^\dagger Q$ -regularization:
 - **Both the left hand sides $\langle G_j | Q^\dagger$ and the right hand sides $Q | G_{j'} \rangle$ of the matrix elements $\langle G_j | Q^\dagger Q | G_{j'} \rangle$ might be outside the PP function space $\text{span}\{G_n\}$, but they form the same function space, $\text{span}\{QG_n\}$, in which their overlap is computed.**
- For “better arguments” cf. M.W., arXiv:0704.3023 [hep-lat].

Testing ground: Gross-Neveu model (1)

- Action and partition function of the 1+1-dimensional Gross-Neveu model:

$$S = \int d^2x \left(\sum_{n=1}^N \bar{\psi}^{(n)} \left(\gamma_0 (\partial_0 + \mu) + \gamma_1 \partial_1 \right) \psi^{(n)} - \frac{g^2}{2} \left(\sum_{n=1}^N \bar{\psi}^{(n)} \psi^{(n)} \right)^2 \right)$$

$$Z = \int \left(\prod_{n=1}^N D\psi^{(n)} D\bar{\psi}^{(n)} \right) e^{-S}$$

(N : number of flavors; μ : chemical potential; g : coupling constant).

Testing ground: Gross-Neveu model (2)

- Introduce a real scalar field σ and integrate out the fermions:

$$S_{\text{effective}} = N \left(\frac{1}{2\lambda} \int d^2x \sigma^2 - \ln \left(\det \left(\gamma_0(\partial_0 + \mu) + \gamma_1 \partial_1 + \sigma \right) \right) \right)$$

$$Z \propto \int D\sigma e^{-S_{\text{effective}}}$$

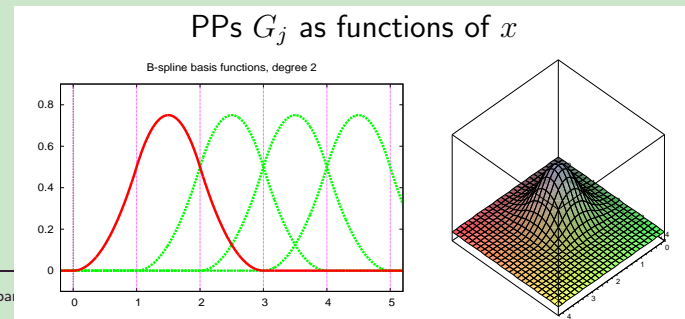
$$(\lambda = Ng^2).$$

- Large- N limit:

- $N \rightarrow \infty$, $\lambda = Ng^2 = \text{constant}$.
- There is no need to compute the σ -path integral anymore.
- It is sufficient to minimize $S_{\text{effective}}$ with respect to σ .
- $\sigma = -g^2 \sum_{n=1}^N \bar{\psi}^{(n)} \psi^{(n)}$ (chiral condensate).

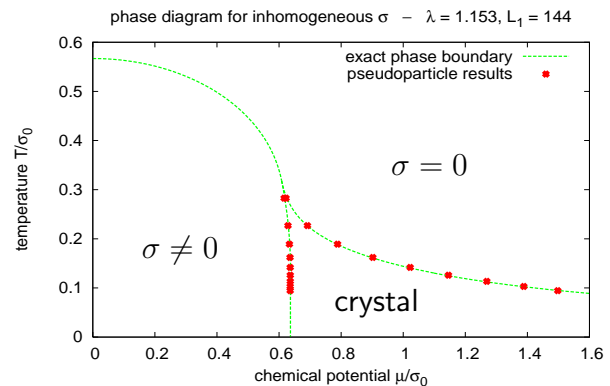
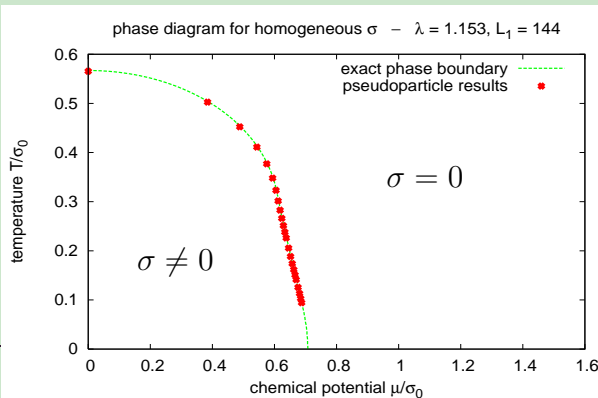
Fermionic PPs

- Fermionic PPs in this talk: a large number of uniformly distributed “hat functions”.
 - B-spline basis functions of degree 2.
 - 8 ... 56 PPs in time direction, 144 PPs in space direction.
 - Antiperiodic boundary conditions in time direction, periodic boundary conditions in space direction.
- “Sensible set of field configurations” (any not too heavily oscillating field configuration can be approximated)
 - we can expect to reproduce correct Gross-Neveu results.
- Piecewise polynomial functions
 - certain integrals can be calculated analytically.



Phase diagram

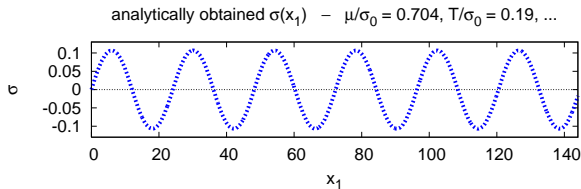
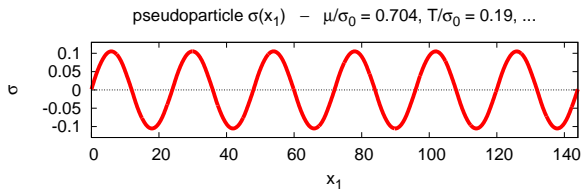
- **Q -regularization: completely wrong and useless results.**
 - No improvement, when using a larger number of PPs.
 - No improvement, when using a different type of PPs.
- **$Q^\dagger Q$ -regularization: excellent agreement with analytical results.**
 - Homogeneous chiral condensate: analytical results by U. Wolff, 1985.
 - Inhomogeneous chiral condensate: analytical results by O. Schnetz, M. Thies, K. Urlichs, 2004.



Chiral condensate

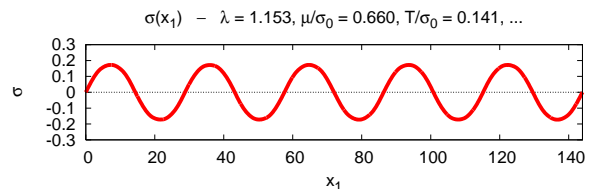
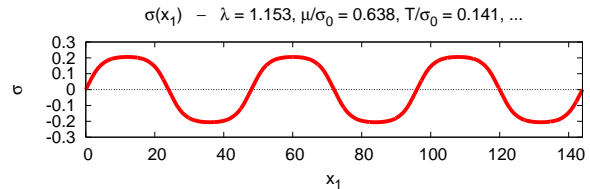
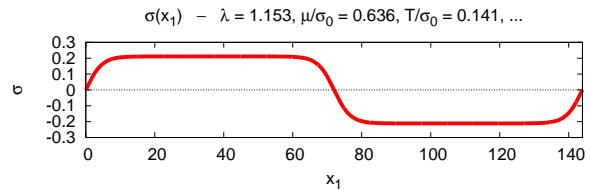
- $Q^\dagger Q$ -regularization: excellent agreement with analytical results.
 - The sin-like behavior “inside the crystal phase” changes to a kink-antikink structure, when approaching the left phase boundary.

PP result



analytical result

close to the left phase boundary



inside the crystal phase

Summary

- Always apply the $Q^\dagger Q$ -regularization and not the naive Q -regularization, when including fermionic fields in the pseudoparticle approach.
- The application of the PP approach to compute the phase diagram of the 1+1-dimensional Gross-Neveu model in the large- N -limit has been a successful test.

Outlook

- Next steps:
 - Apply the PP approach to QCD.
 - Try to identify a small number of physically relevant fermionic PPs (PPs, which are able to approximate typical low lying eigenmodes of the Dirac operator?).
- Current research:
 - Chiral symmetry breaking by computing the low lying eigenmodes of the Dirac operator in the quenched approximation (Banks-Casher relation).
- Goals:
 - **Obtain a model with a small number of degrees of freedom, which exhibits chiral symmetry breaking and a confinement deconfinement phase transition at the same time.**
 - Compute further fermionic observables: pion masses, ...