

# Infrared divergence of the color-Coulomb self-energy in Coulomb gauge QCD

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Collaboration with

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# 1 Introduction

# Confinement mechanism

- Hadrons are considered as bound states of quarks and gluons.
- Physical states are color singlet states.
  - ⇒ The energy of a color non-singlet state diverges.
  - ⇒ absence of isolated quarks and gluons in nature
- ▶ There are several approaches to understand the mechanism of color confinement where topological objects (collective infrared gluonic degrees of freedom) or unphysical degrees of freedom are responsible for the confinement.
  - color monopoles in maximal abelian gauge (dual superconducting scenario)
  - center vortices in maximal center gauge (center vortex model)
  - ghost in Landau gauge

In Coulomb gauge, what are the relevant degrees of freedom for the confinement?

## 2 Coulomb gauge QCD

# Coulomb gauge Hamiltonian

► The Coulomb gauge Hamiltonian can be expressed as the sum of the gluonic part and the instantaneous part:

$$H = \frac{1}{2} \int d^3x \{ (E_i^{\text{tr}})^2 + B_i^2 \} + \frac{1}{2} \int d^3y \int d^3z \rho^a(\vec{y}, t) \mathcal{V}^{ab}(\vec{y}, \vec{z}; A^{\text{tr}}) \rho^b(\vec{z}, t)$$

● color charge density

$$\rho^a = \underbrace{gf^{abc} A_i^{b,\text{tr}} E_i^{c,\text{tr}}}_{\text{gluons}} + \underbrace{igq^\dagger T^a q}_{\text{quarks}}$$

● kernel of the instantaneous interaction

$$\mathcal{V}^{ab}(\vec{y}, \vec{z}; A^{\text{tr}}) \equiv \int d^3x \mathcal{G}^{ac}(\vec{y}, \vec{x}; A^{\text{tr}}) (-\nabla_{\vec{x}}^2) \mathcal{G}^{cb}(\vec{x}, \vec{z}; A^{\text{tr}})$$

►  $\mathcal{G}$  is the Green's function of the Faddeev-Popov (FP) operator:

$$M^{ab} = -\partial_i D_i^{ab} = -\delta^{ab} \partial_i^2 - gf^{abc} A_i^{c,\text{tr}} \partial_i$$

# Color-Coulomb potential

(D. Zwanziger, NPB518, 237 (1998))

► We can show by the perturbative analysis that the instantaneous interaction causes an anti-screening effect.

⇒ We expect that the instantaneous interaction provides a confining force.

► The color-Coulomb potential is defined as **the instantaneous interaction energy between color charges**:

$$V_C(\vec{x} - \vec{y}) = g^2 \vec{T}^a \cdot \vec{T}^b \langle \mathcal{V}^{ab}(\vec{x}, \vec{y}; A^{\text{tr}}) \rangle$$

► At the tree level, the color-Coulomb potential is the  $1/R$  Coulomb potential.

# Coulomb gauge QCD

- Gribov ambiguity and instantaneous interaction  
⇒ conjecture: Instantaneous interaction provides a long-range attractive force?  
(V.N. Gribov, NPB138, 1 (1978))
- fundamental modular region and Gribov-Zwanziger scenario  
(D. Zwanziger, NPB412, 657 (1994))
- $g^2 D_{00}$  is a renormalization group invariant. (independent of the ultraviolet cutoff and the renormalization point and of the regularization and the renormalization scheme)  
⇒  $I$  and  $P$  are renormalization group invariants.  
(D. Zwanziger, NPB518, 237 (1998))
- Coulomb gauge is a renormalizable gauge (finite limit of the interpolating gauge).  
(L. Baulieu and D. Zwanziger, NPB548 (1999) 527)

- gluon propagators in SU(2) lattice simulations  
 $\implies \tilde{D}_{00}(\vec{k})$  is strongly enhanced at  $\vec{k} = \mathbf{0}$ , while  $\tilde{D}^{\text{tr}}(\vec{k})$  vanishes.  
 (A. Cucchieri and D. Zwanziger, PRD65 (2001) 014001)
- (scheme independent) invariant color charge  
 (A. Cucchieri and D. Zwanziger, PRD65 (2001) 014002)
- Zwanziger's inequality  $V_C(\mathbf{R}) \geq V(\mathbf{R})$   
 $\implies$  If the color-Coulomb potential is non-confining, then the physical potential is also non-confining, i.e., "No confinement without color-Coulomb confinement".  
 (D. Zwanziger, PRL90 (2003) 102001)
- gluon propagators and ghost propagator in SU(2) lattice simulations  
 $\implies$  The ghost form factor acquires an infrared singularity.  
 (K. Langfeld and L. Moyaerts, PRD70 (2004) 074507)

# Lattice QCD simulations — confinement phase

J. Greensite and S. Olejnik, PRD67, 094503 (2003)

A. Nakamura and T. Saito, PTP115, 189 (2006)

- ▶ The color-Coulomb potential behaves as a **linearly rising** potential at large distances.
- ▶ The color-Coulomb string tension is 2-3 times larger than that of the physical potential.  
⇒ Zwanziger's inequality
- ▶ **The instantaneous interaction provides a strong confining force.**

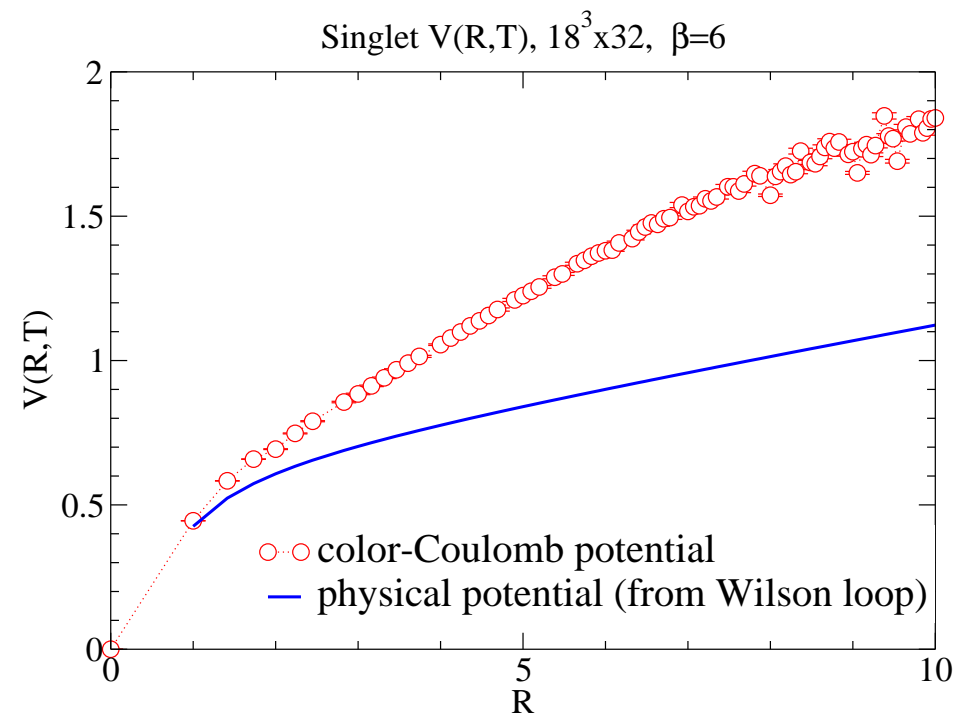


Figure 1: (from PTP115:189 (2006))

# Lattice QCD simulations — deconfinement phase

J. Greensite and S. Olejnik, PRD69, 074506 (2004)

A. Nakamura and T. Saito, PTP115, 189 (2006)

- The color-Coulomb potential is a confining potential even in the **deconfinement phase**.

⇒ The color-Coulomb string tension does not serve as an order parameter for confinement/deconfinement phase transition.

- In Coulomb gauge, the confinement is attributed to the **instantaneous interaction**, whereas the confinement/deconfinement phase transition will be caused by the **non-instantaneous interaction**.

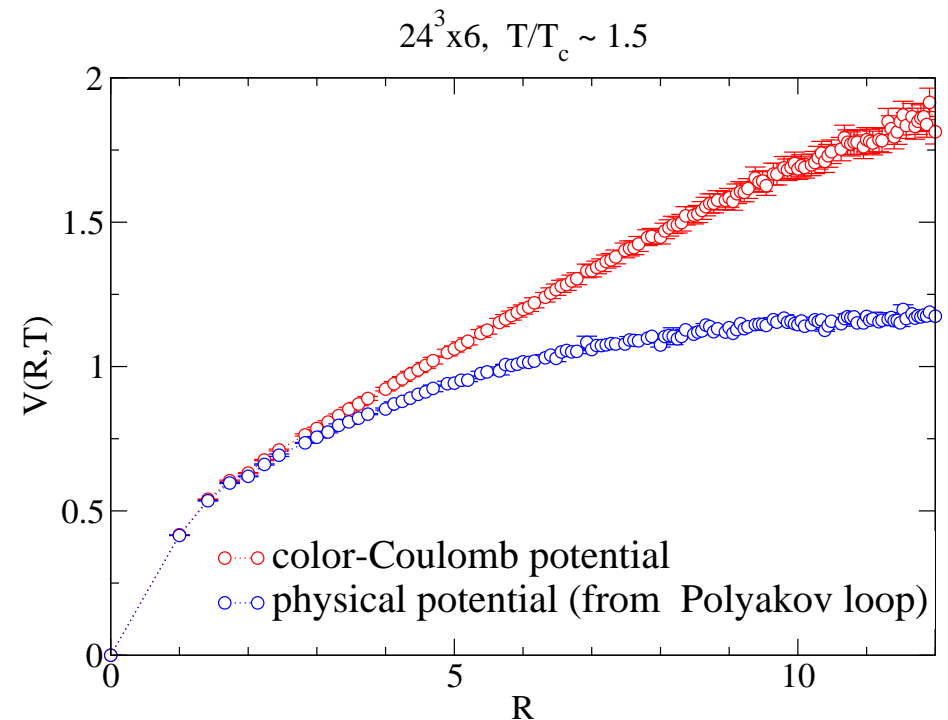


Figure 2: (from PTP115:189 (2006))

### **3 Eigenvalue distribution of the FP operator**

# Motivation for our study

Now, we have two questions:

- ▶ Why is the color-Coulomb potential confining in the confinement phase?
- ▶ Why is the color-Coulomb potential confining even in the deconfinement phase? In other words, why does not the color-Coulomb potential show a critical behavior?

⇒ To address these questions, we study the eigenvalue distribution of the FP operator.

# Color-Coulomb self-energy

(J. Greensite, S. Olejnik and D. Zwanziger, JHEP 05, 070 (2005))

- The color-Coulomb potential is defined as the instantaneous interaction energy between color charges:

$$\begin{aligned} V_C(\vec{x} - \vec{y}) &= g^2 \vec{T}^a \cdot \vec{T}^b \langle \mathcal{V}^{ab}(\vec{x}, \vec{y}; A^{\text{tr}}) \rangle \\ &= g^2 \vec{T}^a \cdot \vec{T}^b \left\langle \int d^3x \mathcal{G}^{ac}(\vec{x}, \vec{z}; A^{\text{tr}}) (-\nabla_{\vec{z}}^2) \mathcal{G}^{cb}(\vec{z}, \vec{y}; A^{\text{tr}}) \right\rangle \end{aligned}$$

- $\mathcal{G}$  is the Green's function of the Faddeev-Popov (FP) operator:

$$M^{ab} = -\partial_i D_i^{ab} = -\delta^{ab} \partial_i^2 - g f^{abc} A_i^{c,\text{tr}} \partial_i$$

- The Green's function of the FP operator can be expanded in terms of the eigenvectors  $\phi_n^a(\vec{x})$  and the eigenvalues  $\lambda_n$  of the FP operator:

$$\mathcal{G}^{ab}(\vec{x}, \vec{y}; A^{\text{tr}}) = \sum_{n=1}^{8V_3} \frac{\phi_n^{*a}(\vec{x}) \phi_n^b(\vec{y})}{\lambda_n}$$

- The color-Coulomb potential can be written as

$$V_C(\vec{x} - \vec{y}) = g^2 \vec{T}^a \cdot \vec{T}^b \left\langle \sum_{n,m=1}^{8V_3} \frac{\phi_n^{*a}(\vec{x}) \phi_m^b(\vec{y})}{\lambda_n \lambda_m} \int d^3 z \phi_n^c(\vec{z}) (-\nabla_{\vec{z}}^2) \phi_m^{*c}(\vec{z}) \right\rangle$$

- We set  $\vec{y} = \vec{x}$  and consider the color-Coulomb self-energy

$$V_C(\vec{x} - \vec{x}) = g^2 \frac{C_D}{N^2 - 1} \left\langle \sum_{n,m=1}^{8V_3} \frac{\phi_n^{*a}(\vec{x}) \phi_m^a(\vec{x})}{\lambda_n \lambda_m} \int d^3 z \phi_n^c(\vec{z}) (-\nabla_{\vec{z}}^2) \phi_m^{*c}(\vec{z}) \right\rangle$$

- By Fourier transforming  $V_C(\vec{x} - \vec{x})$  and using the orthonormal condition

$$\int d^3 x \phi_n^{*a}(\vec{x}) \phi_m^a(\vec{x}) = \delta_{nm},$$

we have

$$\int \frac{d^3 p}{(2\pi)^3} \tilde{V}_C(\vec{p}) = \frac{g^2 C_D}{8V_3} \left\langle \sum_{n=1}^{8V_3} \frac{F_n}{\lambda_n^2} \right\rangle$$

where  $F_n$  are the expectation values of the negative Laplacian in the FP eigenmodes

$$F_n = \int d^3x \phi_n^{*a}(\vec{x}) (-\nabla^2) \phi_n^a(\vec{x})$$

► We define the normalized density of the FP eigenvalues

$$\rho(\lambda) \equiv \frac{N(\lambda, \lambda + \Delta\lambda)}{8V_3\Delta\lambda}$$

where  $N(\lambda, \lambda + \Delta\lambda)$  is the number of eigenvalues in the range  $[\lambda, \lambda + \Delta\lambda]$ .

► Then, we have

$$\int_0^\Lambda \frac{d|\vec{p}|}{4\pi} |\vec{p}|^2 \tilde{V}_C(|\vec{p}|) = g^2 C_D \int_0^{\lambda_{max}} d\lambda \frac{\langle \rho(\lambda) F(\lambda) \rangle}{\lambda^2}$$

where  $\Lambda$  is the ultraviolet cutoff and  $\lambda_{max}$  is the corresponding maximum value of the FP eigenvalue. At the tree level,  $V_C$  is the Coulomb potential. Therefore at sufficiently large  $\vec{p}$ ,  $\tilde{V}_C(|\vec{p}|) \sim 1/|\vec{p}|^2$ , and the integration in the left-hand side is finite in the upper limit as long as we keep the cutoff finite.

$$\int_0^\Lambda \frac{d|\vec{p}|}{4\pi} |\vec{p}|^2 \tilde{V}_C(|\vec{p}|) = g^2 C_D \int_0^{\lambda_{max}} d\lambda \frac{\langle \rho(\lambda) F(\lambda) \rangle}{\lambda^2}$$

► If the condition

$$\lim_{\lambda \rightarrow 0} \frac{\langle \rho(\lambda) F(\lambda) \rangle}{\lambda} > 0$$

is satisfied in the infinite volume limit, the integration in the right-hand side diverges in the lower limit. In other words,

If the criterion is satisfied, the color-Coulomb potential is more singular than the Coulomb potential ( $\sim 1/|\vec{p}|^2$ ) in the infrared region.

This is the necessary condition for the color-Coulomb potential being a confining potential.

## Zero-th order in the coupling or an abelian theory

$$\left( M^{ab} = -\partial_i D_i^{ab} = -\delta^{ab} \partial_i^2 - g f^{abc} A_i^{c,\text{tr}} \partial_i \longrightarrow M = -\partial_i^2 \right)$$

- Eigenvalue equation:  $-\nabla^2 \phi(\vec{x}) = \lambda \phi(\vec{x})$
- Eigenvalue:  $\lambda = \vec{k}^2$
- Eigenvectors: plane waves
- Eigenvalue density:  $\rho(\lambda) = \sqrt{\lambda}/4\pi^2$
- expectation value of the negative Laplacian:  $F(\lambda) = \lambda$
- ▶ confinement criterion is not satisfied:

$$\lim_{\lambda \rightarrow 0} \frac{\rho(\lambda) F(\lambda)}{\lambda} = \lim_{\lambda \rightarrow 0} \frac{\sqrt{\lambda}}{4\pi^2} = 0$$

# Simulation details — Action and gauge fixing

- The gauge configurations are generated by the heat-bath Monte Carlo technique with the **Wilson plaquette action**:

$$S_W = \beta \sum_x \sum_{\mu < \nu} \left[ 1 - \frac{1}{3} \text{Tr}(U_{x;\mu\nu}) \right]$$

( $\beta = 6/g^2$  is the lattice coupling constant.)

- We adopt the **iterative method** to fix the gauge: we minimize the functional

$$F_U[g] = \sum_{i,x} \text{ReTr} \left( 1 - \frac{1}{3} g^\dagger(x) U_i(x) g(x + \hat{i}) \right)$$

with respect to the gauge rotation  $g(x)$  and find the local minimum of  $F_U[g]$ .

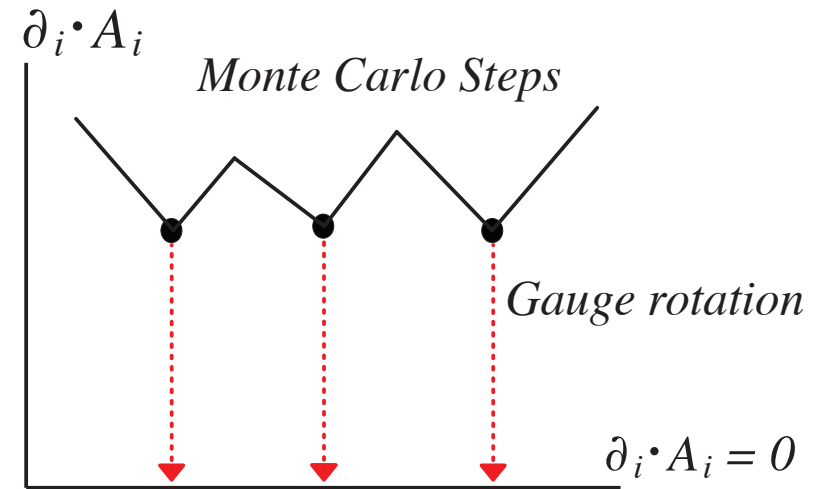


Figure 3: Wilson-Mandula method

# Simulation details — Eigenvalue equation

- ▶ Construct the lattice FP operator

$$M_{xy}^{ab} = \sum_i \Re \text{Tr} \left[ \{T^a, T^b\} \left( U_i(x) + U_i(x - \hat{i}) \right) \delta_{x,y} \right. \\ \left. - 2T^b T^a U_i(x) \delta_{y, x + \hat{i}} - 2T^a T^b U_i(x - \hat{i}) \delta_{y, x - \hat{i}} \right]$$

- ▶ Solve the eigenvalue equation

$$M\phi_n = \lambda_n\phi_n$$

using the Lanczos method and obtain the lowest 1000 eigenvalues and corresponding eigenvectors.

► Calculate

$$\rho(\lambda) = \frac{N(\lambda, \lambda + \Delta\lambda)}{8V_3\Delta\lambda} \quad \text{and} \quad F_\lambda = \int d^3x \phi_\lambda^{*a}(\vec{x}) (-\nabla^2) \phi_\lambda^a(\vec{x}),$$

and check whether the confinement criterion

$$\lim_{\lambda \rightarrow 0} \frac{\langle \rho(\lambda) F(\lambda) \rangle}{\lambda} > 0$$

is satisfied.

- Previous work: J. Greensite, S. Olejnik and D. Zwanziger, JHEP05, 070 (2005)  
⇒ FP eigenvalue density in SU(2) lattice simulation

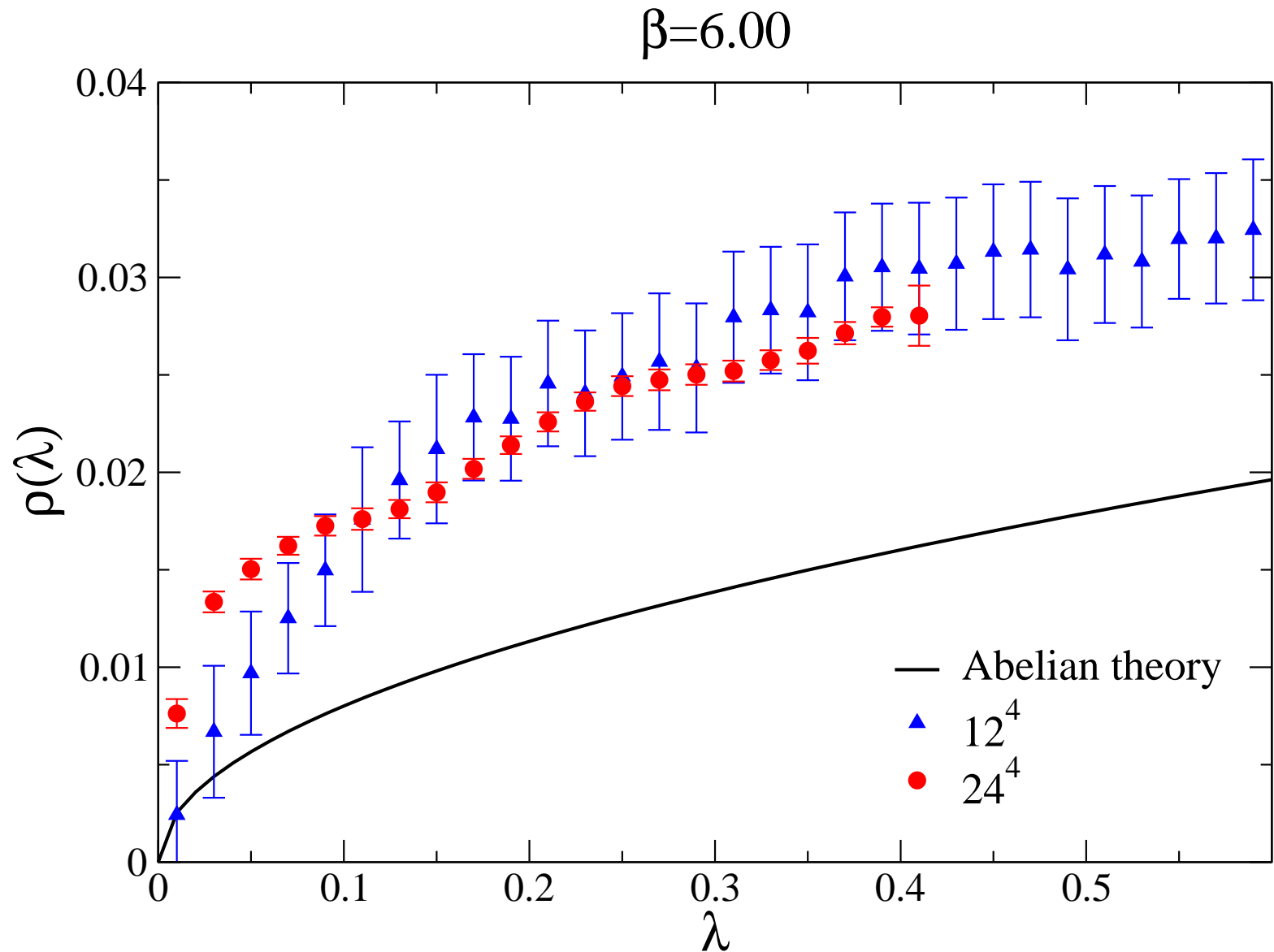


Figure 4:  $\rho(\lambda)$  in the confinement phase. The FP eigenvalue density near  $\lambda = 0$  increases at large lattice volume.

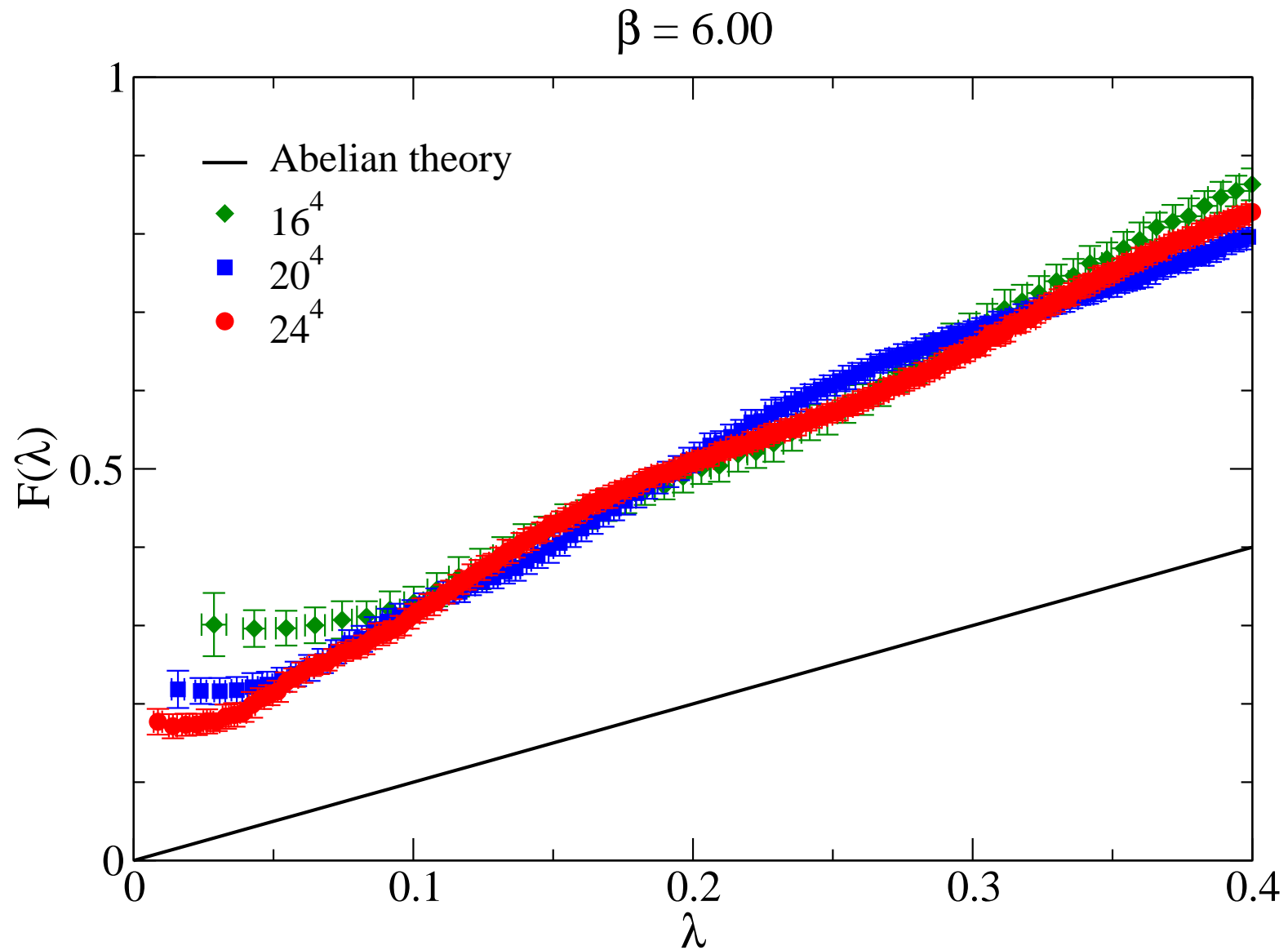


Figure 5:  $F(\lambda)$  in the confinement phase. We expect that  $F(\lambda)$  has a finite value at  $\lambda = 0$  in the infinite volume limit.

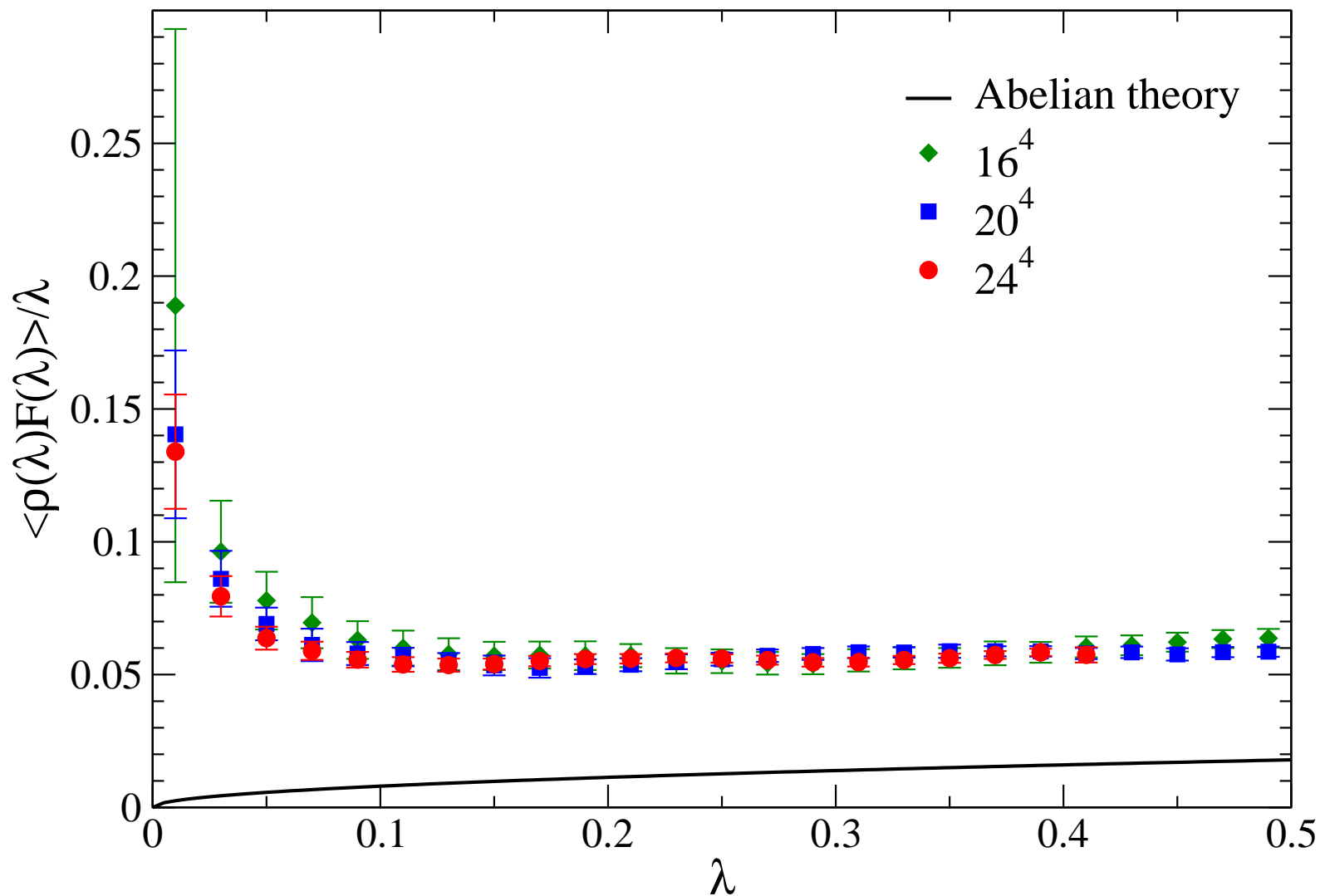
$\beta=6.00$ 

Figure 6:  $\rho(\lambda)F(\lambda)/\lambda$  in the confinement phase. We see the enhancement near  $\lambda = 0$  and we conclude that the necessary condition  $\lim_{\lambda \rightarrow 0} \rho(\lambda)F(\lambda)/\lambda > 0$  is satisfied in the confinement phase.

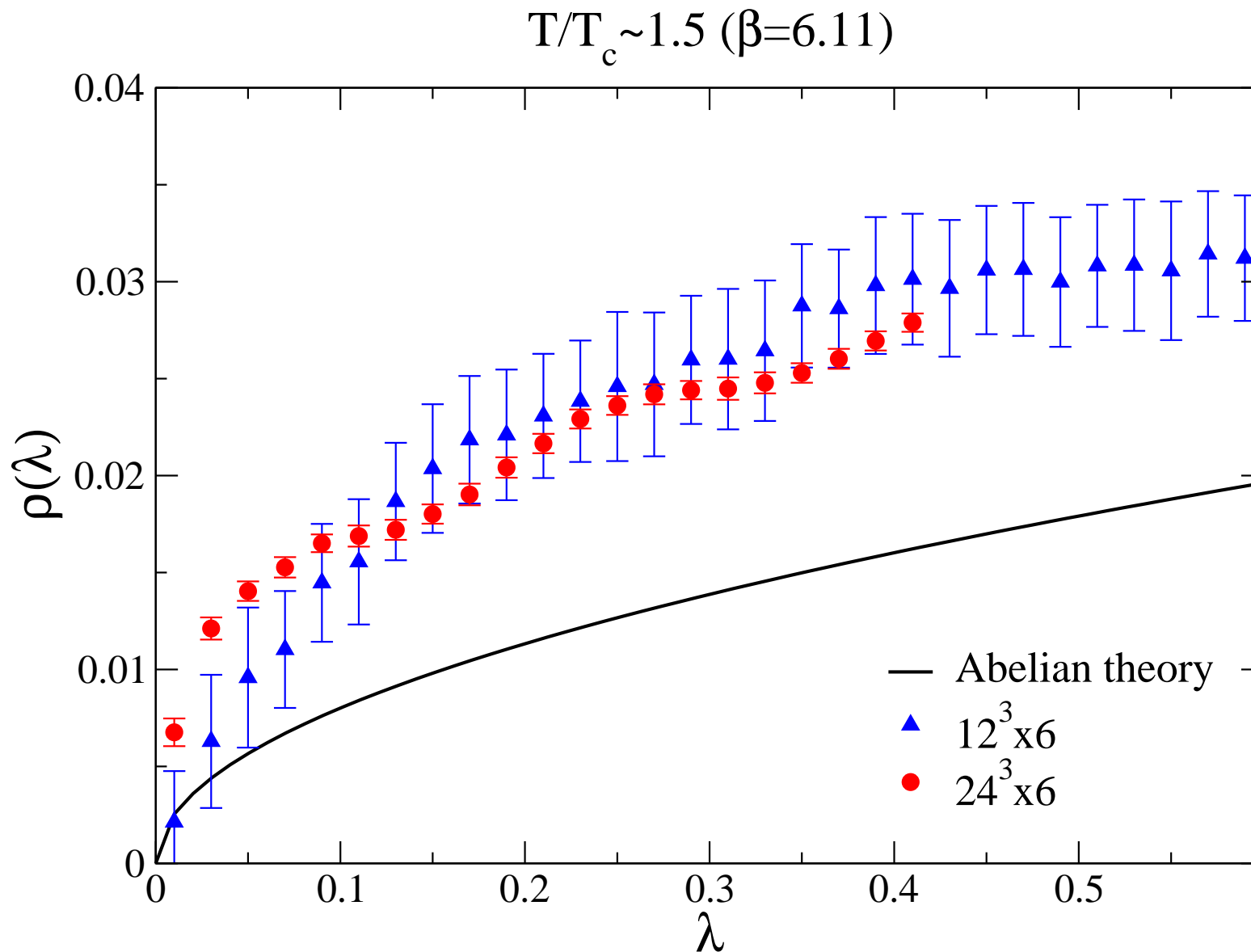


Figure 7:  $\rho(\lambda)$  in the deconfinement phase. The behavior of  $\rho(\lambda)$  in the deconfinement phase is qualitatively the same as in the confinement phase and the FP eigenvalue density does not show a critical behavior.

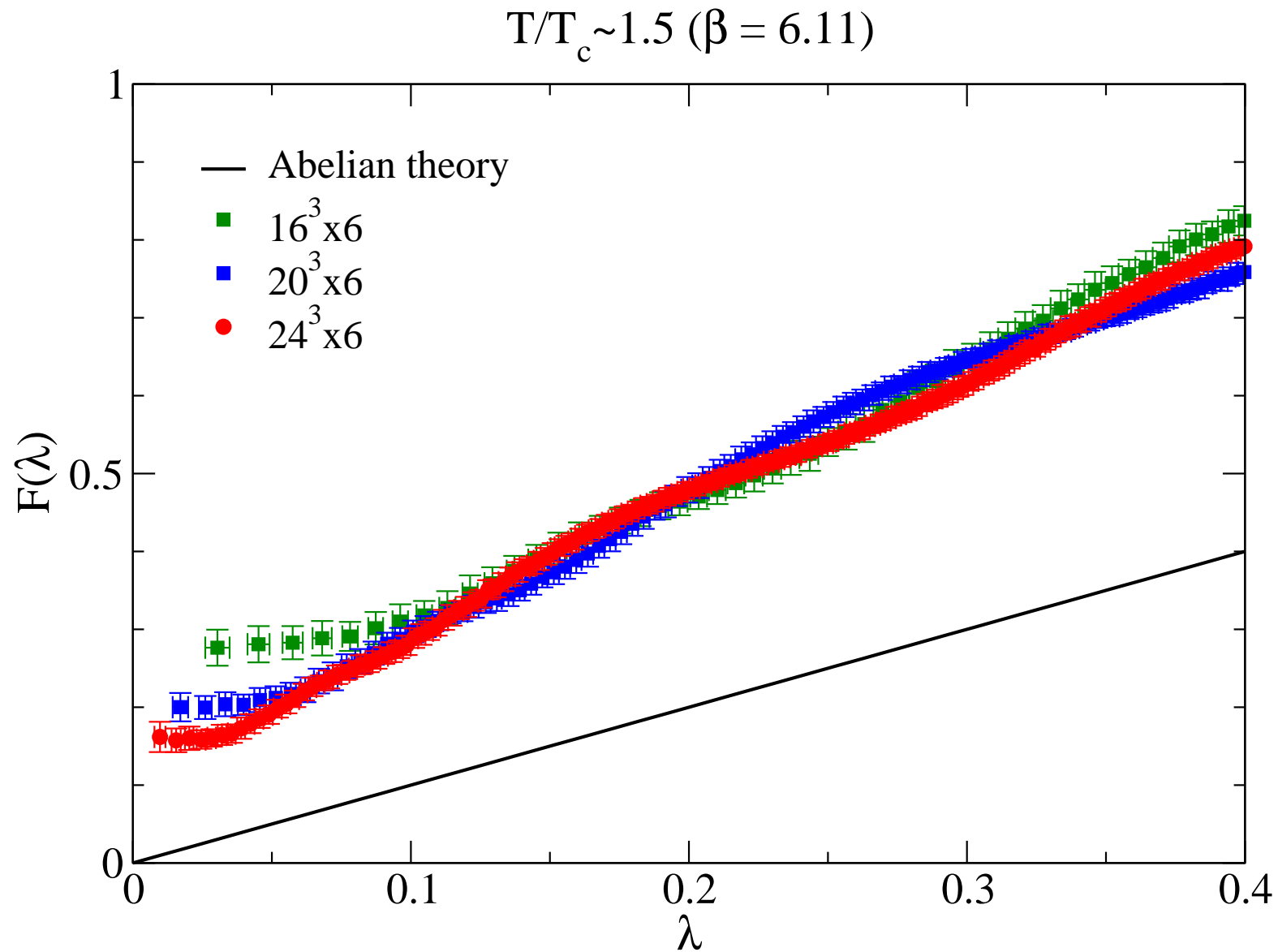


Figure 8:  $F(\lambda)$  in the deconfinement phase.  $F(\lambda)$  shows a similar behavior as in the confinement phase.

$$T/T_c \sim 1.5 \quad (\beta=6.11)$$

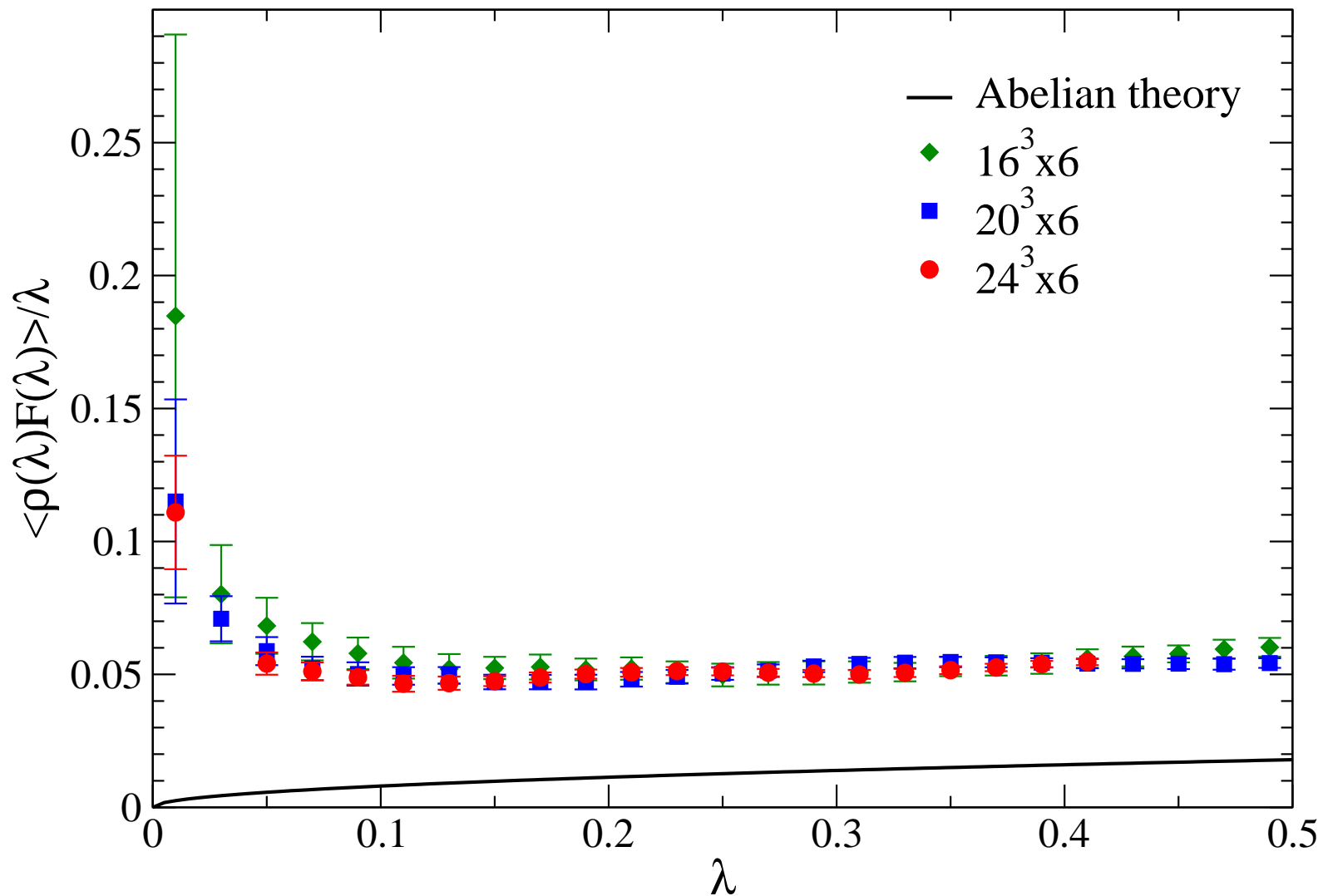


Figure 9:  $\rho(\lambda)F(\lambda)/\lambda$  in the deconfinement phase. The necessary condition  $\lim_{\lambda \rightarrow 0} \rho(\lambda)F(\lambda)/\lambda > 0$  is satisfied even in the deconfinement phase.

## **4 Summary**

# Summary

- ▶ We investigated the eigenvalue distribution of the FP operator and it is revealed that...
- The FP eigenvalue density of near-zero modes in  $SU(3)$  Yang-Mills theory is greater than that of the abelian theory and the necessary condition is satisfied in the confinement phase. Therefore, the color-Coulomb potential is more singular than the Coulomb potential in the infrared region.
- The near-zero modes of the FP operator survive above the critical temperature and the necessary condition is satisfied even in the deconfinement phase. This is the reason why the color-Coulomb potential is a confining potential in the deconfinement phase and does not show a critical behavior.