

# Saturated Nuclear Matter in the Large $N_c$ and Heavy Quark Limits of Quantum Chromodynamics

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- Background
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  - What is Quantum Chromodynamics (QCD)?
  - QCD Intractability
- $N_c$  and  $M_q$  Limits: A Tractable Approach
- Saturated Nuclear Matter
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  - New considerations: Sub-leading order
- Calculations and correction terms
  - Nearest-neighbor distance
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# Subatomic Particles

## Hadrons and Their Constituents

- Nuclei consist of protons and neutrons, which are hadrons
- Hadrons are particles whose dominant interaction is the **strong** nuclear force

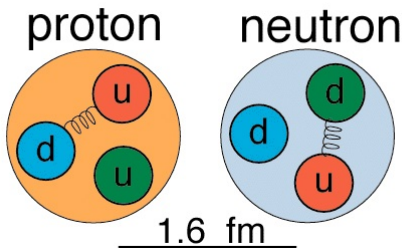
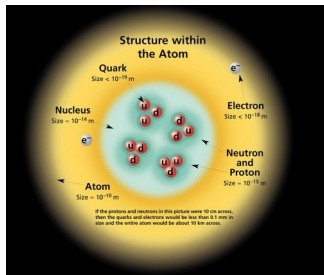
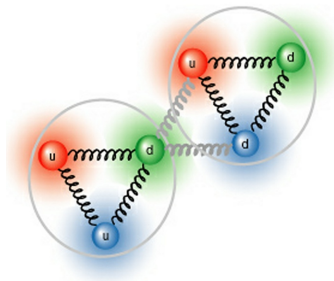


Figure : Left: Subatomic structure from AP Chemistry Wikispaces  
Right: PBS portrayal of up and down quarks that make up nucleons

# Quantum Chromodynamics (QCD)

- Quantum field theory of the strong interaction
  - Governs nuclear physics
- QCD is a theory of quarks and gluons
  - Quarks are sub-nuclear particles carrying “color”
    - Colors are red, green, and blue
    - Objects must be in a color-singlet state, i.e. color-neutral
  - “Gluons” mediate the strong interaction

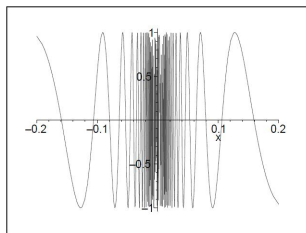


**Figure :** Gluons are what keep protons in a nucleus together.  
Image courtesy of ScienceBlogs.com.

# QCD Intractability

In principle, one could extract all useful information from the QCD Lagrangian. In practice:

- Perturbative expansions fail to converge
- Numerical techniques
  - Lattice QCD numericizes path integrals
  - Rapidly oscillating functions prevent accurate results



**Figure :**  $\sin(\frac{1}{x})$  is an example of an oscillatory function that would be impossible to integrate numerically.

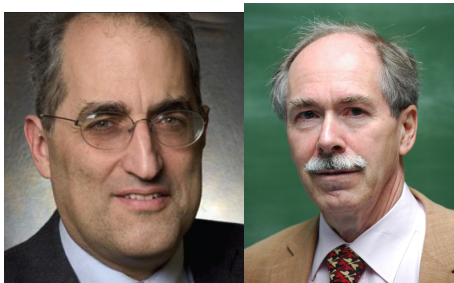
# The $N_c$ and $M_q$ Limits of QCD

$N_c$  : number of colors.

$M_q$  : quark mass.

When  $N_c$  and  $M_q$  grow arbitrarily large, QCD simplifies greatly.

This model is based on expansions in powers of  $\frac{1}{N_c}$ .



**Figure** : Edward Witten (left) and Gerard 't Hooft (right), pioneers of the double limit of QCD. Images from the Institute for Advanced Study.

Important to note:

- Hadrons have two sub-groups
  - Baryons (three quarks) and mesons (quark-antiquark)
  - Mesons are unstable and thus unimportant to this research

Simplifications of the double limit:

- Reduces a field theory to a quantum mechanical theory to describe baryons
- Particles move nonrelativistically
- The strong force reduces to a color-Coulomb interaction

# The Problem: Saturated Nuclear Matter

- Does nuclear matter bind to itself in the double limit?
  - Self-binds, i.e. saturates in the real world
  - Previous research: repulsive at leading order
- Sub-leading order considerations
  - Sub-leading order usually unimportant
  - “Glueballs,” conglomerate particles of color-carrying gluons, provide an attractive force at sub-leading order

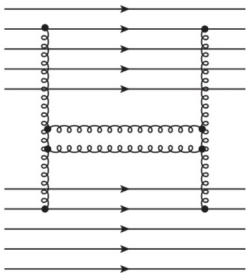


Figure : Two interacting gluons must exchange at least two quarks to preserve a color singlet, giving rise to the “glueball.”



- Analytic and numerical components
  - Analytic work completed by hand
  - Wolfram Mathematica software used to generate numerical solutions
- The two approaches produced results that agree closely

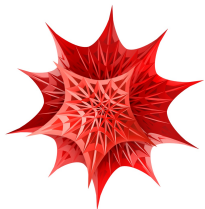


Figure : Mathematica Icon, courtesy of Wolfram.com.

# Leading Order

At leading order, the interaction energy per baryon is entirely Pauli repulsion:

$$E_{\text{Pauli}} = c_1 N_c M_q \tilde{\alpha}_s^2 \tilde{d}^p \exp(-c_2 \tilde{d}), \quad (1)$$

where  $c_1$ ,  $c_2$ , and  $p$  are known numerical constants and  $\tilde{d} = (\tilde{\alpha}_s m_q d)$ . The variable  $d$  is the distance between nearest-neighbor baryons.

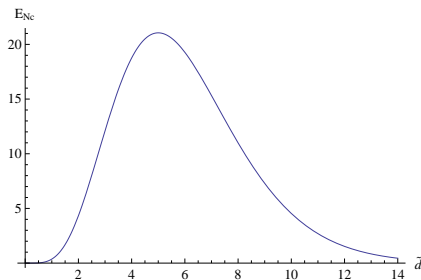


Figure : Leading order interaction energy vs. separation distance.

## Sub-leading order

According to quantum field theory, point sources have a Yukawa potential:

$$V(r) = -a \frac{e^{-kmr}}{r}, \quad (2)$$

where  $r$  is the distance to the particle,  $m$  is the mass of the particle, and  $a$  and  $k$  are constants.

The potential for a system of  $n_{gb}$  glueballs is

$$V_{\text{glueball}} = \sum_{i=1}^{n_{gb}} -a \frac{\exp(-km_i r)}{r} \quad (3)$$

Taking  $m_1$  to be the least massive glueball and factoring:

$$V = -a \frac{e^{-km_1 r}}{r} \left[ 1 + \sum_{j=1}^{n_{gb}} \frac{\exp(-km_j r)}{\exp(-km_1 r)} \right] \quad (4)$$

So all but the lightest glueball is exponentially suppressed.

# Sub-leading order, continued

Glueball potential is re-expressed as

$$E_{gb} = -\tilde{g}_{gb} \frac{\exp(-\tilde{m}_{gb}\tilde{d})}{\tilde{d}} \quad (5)$$

where  $\tilde{m}_{gb} \equiv \frac{m_{gb}}{\tilde{\alpha}_s M_q}$  and  $\tilde{g}_{gb}$  is a dimensionless constant.

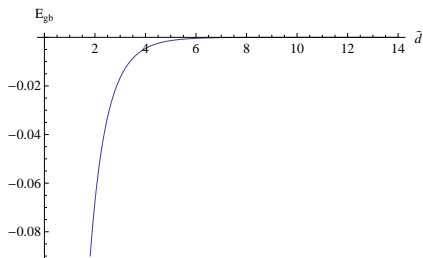


Figure : Glueball interaction energy vs. separation distance.

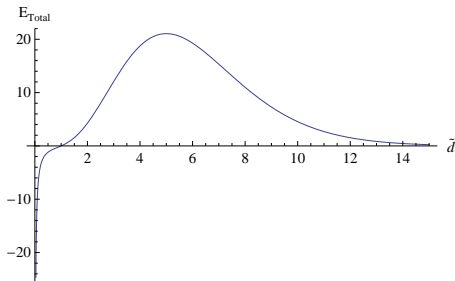
# Nuclear Matter Saturates in the Double Limit

Interaction energy at sub-leading order in  $N_c$ :

$$E_{\text{Interaction}} = E_{\text{Pauli}} + E_{\text{gb}} \quad (6)$$

Differentiating yields separation distance between baryons:

$$\tilde{d} \approx \frac{1}{(c_2 - \tilde{m}_{\text{gb}})} \ln\left(\frac{c_1 N_c M_q \tilde{\alpha}_s}{\tilde{g}_{\text{gb}}}\right) \quad (7)$$



**Figure :** Total interaction energy vs. separation distance. Saturation occurs at the x-intercept.

The derivation of the nearest-neighbor distance used the approximation  $\ln \tilde{d} \ll \tilde{d}$ .

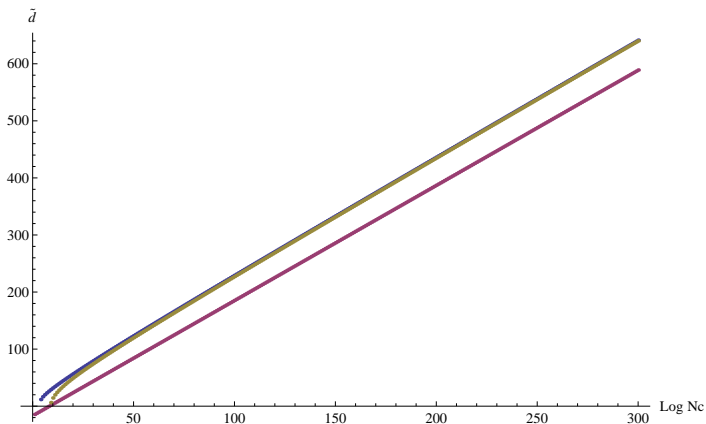
Error terms are of the form:

$$\ln(\ln(\frac{c_1 N_c M_q \tilde{\alpha}_s}{\tilde{g}_{gb}})) \quad (8)$$

Note that  $\ln \tilde{d}$  converges so slowly as to be a constant for all intents and purposes.

# Analytic and Numerical Agreement

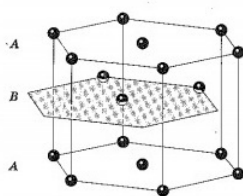
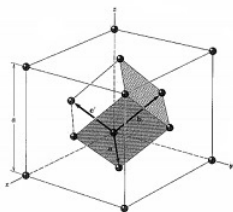
Analytic and numerical values for  $\tilde{d}$  agree in the large  $N_c$  limit.



**Figure :** The analytic approximation (red) is approximately linear in  $\log N_c$  and parallel to the numerical calculation (blue). The analytic approximation with the correction term (green) is indistinguishable from the numerical calculation for large  $N_c$ .

# Crystal Structure

- Baryons form a crystal in the  $M_q$  limit
- Lowest energy configuration corresponds to maximal atomic packing fraction
- Greatest atomic packing fraction:  $\frac{\pi}{3\sqrt{2}} \approx .74$ 
  - Face-Centered Cubic
  - Hexagonal Close-Packed



**Figure :** Face-Centered Cubic (left) and Hexagonal Close-Packed (right) crystal structures. Images are from Kittel's "Introduction to Solid State Physics."



- Primary focus: baryon interactions at sub-leading order
  - Phenomena at different orders in  $N_c$  typically do not affect each other
- Nuclear matter saturates
  - Nearest-neighbor distance  $\tilde{d} = \frac{1}{c_2 - \tilde{m}_{\text{gb}}} \ln\left(\frac{c_1 N_c M_q \tilde{\alpha}_s}{\tilde{g}_{\text{gb}}}\right)$
  - Result depends solely upon quantities that are either known or can be computed via lattice QCD
  - Correction term to  $\tilde{d}$ :  $(p + 1) \ln \tilde{d}$
  - FCC or HCP Crystal Structure

This research lays a foundation from which more realistic physical situations in QCD may be studied.

- Calculation of nuclear matter interaction energy
- Identifying whether FCC or HCP is more energetically favorable
  - Purely a theoretical question, as nuclear matter does not generally crystallize
- Lowering  $N_c$  and  $M_q$  into more realistic regimes
  - A phase transition is expected upon lowering  $M_q$

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