# SU(3) LATTICE MONTE CARLO CALCULATION OF THE GLUEBALL MASS SPECTRUM 

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We have calculated the glueball masses of various spins and parities in $\operatorname{SU}(3)$ gauge theory. Our first results give $m_{\mathrm{M}}\left(0^{++}\right)=(3.6 \pm 0.2) \Lambda_{\mathrm{mom}}, m_{\mathrm{E}}\left(0^{++}\right)=(4.3 \pm 0.3) \Lambda_{\mathrm{mom}}, m\left(0^{-+}\right)=\left(7.2_{-0.9}^{+1.6}\right) \Lambda_{\mathrm{mom}}, m_{\mathrm{M}}\left(2^{++}\right)=(8.1 \pm 1.1) \Lambda_{\mathrm{mom}}$ and $m_{\mathrm{E}}\left(2^{++}\right)=\left(8.3_{-1.0}^{+1.6}\right) \Lambda_{\mathrm{mom}}$ as well as information on the glueball wave functions.

In this letter we extend our recent calculation of the glueball mass spectrum [1] to the gauge group SU(3) (i.e., QCD).

We follow the calculational procedure outlined in our previous letter [1]. That is, we perform a variational calculation by choosing some à priori reasonable class of wave functions $\{\phi\}$ and vary $\phi$ within that class to maximize the expectation value
$\frac{\Gamma_{a}}{\Gamma_{0}}=\frac{\left\langle\phi_{\boldsymbol{P}=0}(t+a) \phi_{\boldsymbol{P}=0}(t)\right\rangle}{\left\langle\phi_{\boldsymbol{P}=0}(t) \phi_{\boldsymbol{P}=0}(t)\right\rangle}=\frac{\left\langle\phi_{\boldsymbol{P}=0}(t) \mathrm{e}^{\left.-H a_{\phi_{\boldsymbol{P}}=0}(t)\right\rangle}\right.}{\left\langle\phi_{\boldsymbol{P}=0}(t) \phi_{\boldsymbol{P}=0}(t)\right\rangle}$

$$
\begin{equation*}
=\frac{\left.\Sigma_{n=0}^{\infty} \mathrm{e}^{-m_{n} a}\left|\langle n| \phi_{P=0}(t)\right| \Omega\right\rangle\left.\right|^{2}}{\left.\Sigma_{n=0}^{\infty}\left|\langle n| \phi_{P=0}(t)\right| \Omega\right\rangle\left.\right|^{2}} \tag{1}
\end{equation*}
$$

where $a$ is the lattice spacing and $t$ the euclidean time. The resulting wave function is then used to calculate

$$
\begin{align*}
\frac{\Gamma_{2 a}}{\Gamma_{a}} & =\frac{\left\langle\phi_{P=0}(t+2 a) \phi_{P=0}(t)\right\rangle}{\left\langle\phi_{P=0}(t+a) \phi_{P=0}(t)\right\rangle}=\frac{\left\langle\phi_{P=0}(t) \mathrm{e}^{-H 2 a^{\prime}} \phi_{P=0}(t)\right\rangle}{\left\langle\phi_{P=0}(t) \mathrm{e}^{-H a} \phi_{P=0}(t)\right\rangle} \\
& =\frac{\left.\Sigma_{n=0}^{\infty} \mathrm{e}^{-m_{n} 2 a}\left|\langle n| \phi_{P=0}(t)\right| \Omega\right\rangle\left.\right|^{2}}{\left.\sum_{n=0}^{\infty} \mathrm{e}^{-m_{n} a}\left|\langle n| \phi_{P=0}(t)\right| \Omega\right\rangle\left.\right|^{2}}, \tag{2}
\end{align*}
$$

the logarithm of which will give us the lowest-lying
glueball mass ( $m=m_{0}$ ):
$m=-a^{-1} \ln \left(\Gamma_{2 a} / \Gamma_{a}\right)$.
The idea behind this procedure is that (i) increasing the theoretical accuracy involves searching for a larger signal and (ii) the higher mass states in (2) are double suppressed by the choice of the wave function and the exponential damping factor

It follows that (1) provides an upper bound on the mass:
$-a^{-1} \ln \left(\Gamma_{a} / \Gamma_{0}\right) \geqslant m$.
The pure variational calculations of the glueball mass(es) $[2,3]$ assume that this upper bound will be close to the actual mass(es). How much this is actually so for the rather small lattices and the limited class of wave functions considered can be checked [1] and will be discussed further along with the results.

We work on a $4^{3} \cdot 8$ lattice with 8 lattice points in the time direction. From our previous experience [1], where we have considered $4^{4}, 4^{3} \cdot 8,6^{4}$ and $8^{4}$ lattices, we infer that the $4^{3} \cdot 8$ lattice is large enough for our purposes provided that the coupling parameter $\beta$ is chosen appropriately. Working on a "rectangular" lattice has the further advantage that it lowers the physical temperature at nearly the same computer
cost and, hence, better approximates zero temperature QCD.

An appropriate value of $\beta$ is one for which
$a<$ glueball size $<2 a$.
To two loops the lattice spacing is given by ( $\beta=6 / \mathrm{g}^{2}$ )
$a=\left(83.5 / \Lambda_{\text {mom }}\right)\left(\frac{8}{33} \pi^{2} \beta\right)^{51 / 121} \exp \left(-\frac{4}{33} \pi^{2} \beta\right)$,
where the overall scale is set by the string tension ${ }^{\neq 1}$
[we "measure" $\Lambda_{\text {mom }}=(0.48 \pm 0.05) \sqrt{K}$ at $\beta=5.7$ ].
$\Lambda_{\mathrm{mom}} \approx 0.5 \sqrt{K}$,
which we believe is $\sqrt{K} \approx 400 \mathrm{MeV}$, corresponding to unit Regge slope. Assuming the the glueball size to be of the order of 0.5 fermi (which proves to be correct in the actual calculation), this constrains $\beta$ to the region $5 \lesssim \beta \leqslant 6$. We also require that $\beta$ lies in the continuum region. Leaning on the results of ref. [4] we infer that $5 \Sigma \beta$ is a suitable value.

The occurence of a sharp peak in the specific heat $\mathrm{at}^{\ddagger 2} \beta=5.45 \pm 0.1$ and a rapid increase in the correlation length near this peak, which has been attributed to nearby complex $\beta$ plane singularities [6], cuts the range of admissible values of $\beta$ further down to 5.6 $\zeta \beta$. We have taken "data" at $\beta=5.7$ and $\beta=5.9$ which we have checked lie in the scaling region. Incidentally, the authors in ref. [3] have taken most of their "data" below and at the peak in the specific heat which casts some doubt on their results.

We have constructed states $\phi$ with $J^{P C}=0^{++}, 0^{-+}$, $1^{--}, 1^{+-}$and $2^{++}$which we expect to be low-lying. Except for $0^{-+}$they can be composed of planar loops with purely space-like links $\left(0^{++}, 1^{+-}, 2^{++}\right)$or spaceand time-like links $\left(0^{++}, 1^{--}, 2^{++}\right)$as shown schematically in fig. 1. To project onto zero momentum states the sum is to be taken over all spacial lattice points. The loops that we have considered in our variational calculation are shown in fig. 2. We have confined ourselves to loops that extend only one lattice spacing in time ( $\mathrm{O}_{4}$ and $\mathrm{O}_{5}$, cf. fig. 2 ). That is to re-
${ }^{\ddagger 1}$ We haven taken the mean value of the string tensions "measured" by the authors in ref. [4].
$\neq 2$ This has been obtained by a straight-line interpolation $1 / g^{2} \sim N, N$ being the number of colours, of the $\operatorname{SU}(2)$, $\mathrm{SU}(4), \mathrm{SU}(5)$ and $\mathrm{SU}(6)$ peak/phase transition values reported in ref. [5].Note that they fall very nicely on a straight line.

Re

Im


Re

Re

Im


Fig. $1.0^{++} .2^{++}, 1^{+-}$and $1^{--}$wave functions composed of planar loops. The loops are understood to be summed over all space-like translations. The subscripts M and E stand for magnetic and electric, respectively.
duce the amount of time-like overlap in $\Gamma_{a}$ and to avoid time-like overlaps in $\Gamma_{2 a}$ at all which, otherwise, could lead us into conflict with physical positivity [7]. Note that for $0^{++}$and $2^{++}$two types of different operators arise: magnetic ( $\sim B^{2}$ ) with exclusively space-like extensions and electric ( $\sim E^{2}$ ) extending also in time


Fig. 2. Loops considered in our calculation ( $x, y$ stand for any space direction).
(cf. fig. 1). This doubling of glueball states has also been found in the bag model [8]. The most natural trial wave function for $0^{-+}$is $F \widetilde{F}$, the topological charge density, which is of the type $\boldsymbol{E} \cdot \boldsymbol{B}$. We have taken $\Sigma_{i, j, k= \pm 1}^{ \pm 3} \widetilde{\epsilon}_{0 i j k} F^{0 i} F^{j k}$ instead which extends only one lattice spacing in the time direction in contrast to $F \widetilde{F}$ (which does two) but else has the same quantum numbers and topology. The reason is to secure physical positivity [7] as mentioned before. There are various ways of implementing this operator on the lattice. We use the definition employed in ref. [9].

We have performed several independent lines of iterations of the lattice configuration using the heatbath method [10]. Before taking any "data" we have heated up the lattice starting from cold $(\beta=\infty)$ and hot $(\beta=0)$ starts. So far we have collected $\gtrsim 24000$ "events" at $\beta=5.7$ and $\gtrsim 7000$ "events" at $\beta=5.9$. Our first results for $\operatorname{SU}(3)$ glueball masses and wave functions are summarized below. A more detailed presentation will be given elsewhere.
$0^{++}$. As $0^{++}$has vacuum quantum numbers, it is necessary to replace $\phi$ by $\phi-\langle\phi\rangle$ in (1) and (2) to ensure that the vacuum state does not appear as intermediate state.

We first treat the magnetic and electric $0^{++}$glueball states independently, i.e., not accounting for any mixing between them.
(a) Magnetic. After maximization we obtain at $\beta=5.7$
$\Gamma_{a} / \Gamma_{0}=0.22 \pm 0.01$,
which, according to (4), (6), corresponds to the upper bound
$m_{\mathrm{M}}\left(0^{++}\right) \leqslant(5.52 \pm 0.18) \Lambda_{\mathrm{mom}}$
(where the subscript $M$ stands for magnetic). The accompanying best wave function is (cf. fig. 2)
$\phi_{\mathrm{M}}\left(0^{++}\right) \approx 0.33 O_{1}+0.40 O_{2}+0.85 O_{3}$,
where the operators on the right-hand side are understood to be summed over all orientations in accord with fig. 1 and over the spacial lattice for any fixed $t$. For ease of writing we have dropped the subscript $P$ $=0$.

In the next step we compute $\Gamma_{2 a} / \Gamma_{a}$ now using the wave function (10). We find
$\Gamma_{2 a} / \Gamma_{a}=0.37 \pm 0.02$,
which gives the mass
$m_{\mathrm{M}}\left(0^{++}\right)=(3.6 \pm 0.2) \Lambda_{\text {mom }}$.
We observe that the straightforward variational calculation would have yielded a considerably higher $0^{++}$ glueball mass (9) which disqualifies the pure variational approach. This is not surprising as the class of possible wave functions is rather limited for such a small lattice.
(b) Electric. Our best wave function is $(\beta=5.7$; cf . fig. 2)
$\phi_{\mathrm{E}}\left(0^{++}\right) \approx 0.42 \mathrm{O}_{4}+0.9 \mathrm{O}_{5}$
(where the subscript E stands for electric) which corresponds to
$\Gamma_{a} / \Gamma_{0}=0.16 \pm 0.01$
(note that $\Gamma_{a}$ is positive in line with physical positivity [7]). This results in the upper bound
$m_{\mathrm{E}}\left(0^{++}\right) S(6.68 \pm 0.22) \Lambda_{\text {mom }}$.
Using the optimized wave function (13), we then obtain
$\Gamma_{2 a} / \Gamma_{a}=0.31 \pm 0.02$,
which gives the mass
$m_{\mathrm{E}}\left(0^{++}\right)=(4.3 \pm 0.3) \Lambda_{\text {mom }}$.
Note that also in this case the pure variational calculation cannot be trusted.
(c) Mixing. In reality the magnetic and electric glueball operators will mix. Accordingly, we write
$\phi=\cos \Theta \phi_{M}+\sin \Theta \phi_{E}$.
The physical eigenstates and mass eigenvalues of the transfer matrix are then obtained by searching for two orthogonal wave functions $\phi_{1,2}$ that maximize and minimize, respectively, $\Gamma_{2 a} / \Gamma_{a}$. We have attempted this. But, due to the relatively large errors, we have no reliable numbers yet and even cannot tell with certainty that there are two states (or just one), not to mention the additional mixing with quark-antiquark states. We postpone the details to a forthcoming publication.

At $\beta=5.9$ our best wave functions have a rather small projection onto the lowest lying state which results in a large error of $\Gamma_{2 a} / \Gamma_{a}$.
$2^{+\boldsymbol{+}}$. We now go through the same analysis for $2^{++}$.
(a) Magnetic. At $\beta=5.7$ the best wave function is (cf. fig. 2)
$\phi_{\mathrm{M}}\left(2^{++}\right) \approx 0.18 \mathrm{O}_{1}+0.46 \mathrm{O}_{2}+0.87 \mathrm{O}_{3}$
with
$\Gamma_{a} / \Gamma_{0}=0.058 \pm 0.002$,
which gives the upper bound
$m_{\mathrm{M}}\left(2^{++}\right) \leqslant(10.3 \pm 0.2) \Lambda_{\text {mom }}$.
Using (19) we then obtain
$\Gamma_{2 a} / \Gamma_{a}=0.11 \pm 0.03$,
which gives

$$
\begin{equation*}
m_{\mathrm{M}}\left(2^{++}\right)=(8.0 \pm 1.2) \Lambda_{\mathrm{mom}} \tag{23}
\end{equation*}
$$

Our results at $\beta=5.9$ (with $\sim 1 / 4$ the statistics) agree very well with that. The optimal wave function is
$\phi_{\mathrm{M}}\left(2^{++}\right) \approx 0.14 O_{1}+0.34 O_{2}+0.93 O_{3}$
and gives only a slightly higher bound on $m$ while
$\Gamma_{2 a} / \Gamma_{a}=0.16 \pm 0.03$,
which results in
$m_{\mathrm{M}}\left(2^{++}\right)=(8.4 \pm 1.0) \Lambda_{\text {mom }}$.
This is to say that $\Gamma_{2 a} / \Gamma_{a}$ follows nicely the weak coupling renormalization group trajectory.

By taking the statistical average of (23) and (26) we finally obtain the mass
$m_{M}\left(2^{++}\right)=(8.1 \pm 1.1) \Lambda_{\text {mom }}$.
(b) Electric. At $\beta=5.7$ the best wave function is (cf. fig. 2)
$\phi_{\mathrm{E}}\left(2^{++}\right) \approx \mathcal{O}_{5}$
corresponding to
$\Gamma_{a} / \Gamma_{0}=0.022 \pm 0.002$
(note that $\Gamma_{a}$ is positive in line with physical positivity [7]) which leads to the bound
$m_{\mathrm{E}}\left(2^{++}\right) \leqslant(13.9 \pm 1.0) \Lambda_{\text {mom }}$.
From (28) we then obtain
$\Gamma_{2 a} / \Gamma_{a}=0.11 \pm 0.06$
and
of $\Gamma_{2 a} / \Gamma_{a} \approx\left|\Gamma_{a} / \Gamma_{0}\right|$, suggests that $\Sigma_{i, j, k= \pm 1}^{ \pm 3} \tilde{\epsilon}_{0 i j k}$ $\times F^{0 i F j k}$ is a good $0^{-+}$wave function.
$1^{+-}, l^{--}$. We have no useful bounds and numbers yet in this case.

So far we have expressed the glueball masses in units of $\Lambda_{\text {mom }}$. To make contact to experiment [11] we may use (7) now to fix the absolute scale (but note the obvious caveats). If we do so, we obtain the mass spectrum
$m_{\mathrm{M}}\left(0^{++}\right)=(720 \pm 40) \mathrm{MeV}$,
$m_{\mathrm{E}}\left(0^{++}\right)=(850 \pm 50) \mathrm{MeV}$,
$m\left(0^{-+}\right)=\left(1430_{-180}^{+320}\right) \mathrm{MeV}$,
$m_{\mathrm{M}}\left(2^{++}\right)=(1620 \pm 220) \mathrm{MeV}$,
$m_{\mathrm{E}}\left(2^{++}\right)=\left(1670_{-200}^{+320}\right) \mathrm{MeV}$
[where we have ignored the uncertainties in (7)].
To summarize, we have reported first results of a calculation of the low-lying glueball masses based on altogether $\gtrsim 31000$ Monte Carlo "events". Our aim for the future is to reduce the errors and to make a more systematic survey of the low-lying glueball spectrum, in particular: to look for "oddballs".

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