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# Lattice field theories

## Exercise sheet 6 – Wilson loop and string tension

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Please see the supplemental material at the end of this exercise sheet for further explanations and hints.

### Exercise 11: Connecting strong and weak coupling

Consider a 2-dimensional SU(2) Yang-Mills theory on the lattice, whose partition function reads

$$Z = \int \mathcal{D}U \exp \left\{ -\beta \sum_n \left[ 1 - \frac{1}{2} \text{ReTr} U_{01}(n) \right] \right\}, \quad (1)$$

where  $U_{01}$  is the plaquette at site  $n$  and directions 0 and 1. The Wilson loop defined in a contour  $\mathcal{C}$  of size  $T \times R$  lattice site  $n$  is given by

$$W(\mathcal{C}) \equiv W(T, R) = \frac{1}{2} \left\langle \text{Tr} \left[ \prod_{\mathcal{C}} U \right] \right\rangle. \quad (2)$$

Use the Metropolis algorithm to simulate this theory on a  $10 \times 10$  lattice with  $0.1 \leq \beta \leq 15.0$ , and compute the Wilson loop for square contours of different sizes.

You should observe that the  $1 \times 1$  Wilson loop has the following limiting behaviours:

- at strong coupling ( $\beta \ll 1$ ):

$$W(1, 1) \underset{\beta \rightarrow 0}{\sim} \frac{\beta}{4}; \quad (3)$$

- at weak coupling ( $\beta \gg 1$ ):

$$W(1, 1) \underset{\beta \rightarrow \infty}{\sim} 1 - \frac{3}{2\beta}. \quad (4)$$

You should also note that larger Wilson loops are suppressed at small  $\beta$ , and that getting good signals for them can be hard. You may want to run longer simulations than in the  $\phi^4$  case.

**Hint:** In order to improve statistics, consider computing  $W(1, 1)$  on all lattice points and then taking the volume average. Do not forget about the periodic boundary conditions.

**Hint:** For large  $\beta$  configurations where plaquette is close to the identity will become more and more important. Consider reducing the variance of your proposals in these situations, e.g., using a narrower Gaussian to generate random numbers, to achieve a higher acceptance rate.

### Exercise 12: Computing the string tension

From strong coupling arguments, we expect the above Wilson loop to have the form

$$W(T, R) = \exp[-\hat{\sigma}RT - \hat{\alpha}(R + T) + \hat{\gamma}], \quad (5)$$

where  $\hat{\sigma}$  is the *string tension* in lattice units, and  $\hat{\alpha}$  accounts for subleading effects proportional to the perimeter of the contour. Those subleading effects can be eliminated by studying the *Creutz ratio*,

$$\hat{\chi}(T, R) = -\ln \left( \frac{W(T, R) W(T-1, R-1)}{W(T-1, R) W(T, R-1)} \right). \quad (6)$$

If eq. (5) correctly describes the behaviour of the Wilson loop, then  $\hat{\chi}$  will coincide with the string tension.

In a way similar to the previous exercise, compute Wilson loops of square and rectangular shapes and construct Creutz ratios for various values of  $\beta$ . Like before, ratios computed at larger values of  $T$  or  $R$  will be noisier. Since we expect  $\hat{\chi}$  to be independent of both  $T$  and  $R$ , a constant fit (including errors, obviously!) will give the final result.

At Strong coupling, your string tension should behave as

$$\hat{\sigma} \underset{\beta \rightarrow 0}{\sim} -\ln \frac{\beta}{4}. \quad (7)$$

**Hint:** Due to larger Wilson loops being suppressed at small  $\beta$  their values may come out negative, but statistically compatible with zero and probably large relative errors. Since the string tension is a real number, taking the logarithm of negative numbers makes no sense. In such situations, you can either disregard the Creutz ratios where  $W$  is negative, or run your simulation for longer.

**Hint:** In order to improve statistics, consider computing  $W(T, R)$  on all lattice points and then taking the volume average. Do not forget about the periodic boundary conditions.

**Hint:** When computing  $\hat{\chi}(1, 1)$  use only  $W(1, 1)$  in the calculation, i.e., take the other loops to be equal to 1.

### Supplemental material

In this exercise we are dealing with a field whose elements belong to the group  $SU(2)$ , and the updating procedure should be such that update proposals also belong to the group.

Since every element of  $SU(2)$  group can be written as the exponential of a member of the  $\mathfrak{su}(2)$  Lie algebra, updating a particular link  $U$  can be written as

$$U \rightarrow RU, \quad R = \exp \left[ i \sum_{a=1}^3 \alpha^a \sigma^a \right], \quad (8)$$

with  $\sigma^a$  being the Pauli matrices and  $\alpha^a$  real numbers. Note that the update may equally well be written as  $U \rightarrow UR$ . Because of the commutation properties of the Pauli matrices the computation of  $R$  can be simplified as

$$R = \cos(|\vec{\alpha}|)\mathbb{1} + i \frac{\alpha^a}{|\vec{\alpha}|} \sigma^a \sin(|\vec{\alpha}|), \quad (9)$$

with  $\mathbb{1}$  being the  $2 \times 2$  identity matrix and  $|\vec{\alpha}| = \sqrt{(\alpha^1)^2 + (\alpha^2)^2 + (\alpha^3)^2}$ . Thus it requires only simple arithmetic and trigonometric functions.

```

matrix U[Nt] [Nx];
for n ← 0 to N_steps do
  for 0 ≤ t < N_t and 0 ≤ x < N_x do           // sweep over the lattice
    oldAction ← S(U);                          // save old action
    oldValue ← U[t] [x];                       // save old U
    R ← randomSU2();                            // random SU(2) matrix
    U[t] [x] ← R * U[t] [x];                  // random change to U within SU(2)
    ΔS = S(U) - oldAction;
    if ΔS > 0 then
      r ← UniformReal(0,1);
      if r > exp(-ΔS) then
        U[t] [x] ← oldValue;                   // reject proposal
      end
    end
  end
end
Obs ← Obs + O(U);                             // update computation of observable
end
Obs ← Obs / N_steps;

```

As always, you should skip a number of steps at the beginning, so the system can thermalise. And remember that consecutive configurations in the Markov chain are likely to be correlated, so it is also a good idea to skip some between consecutive measurements.

Below, you can find a plot of the string tension as a function of  $\beta$  to use for cross-checking.

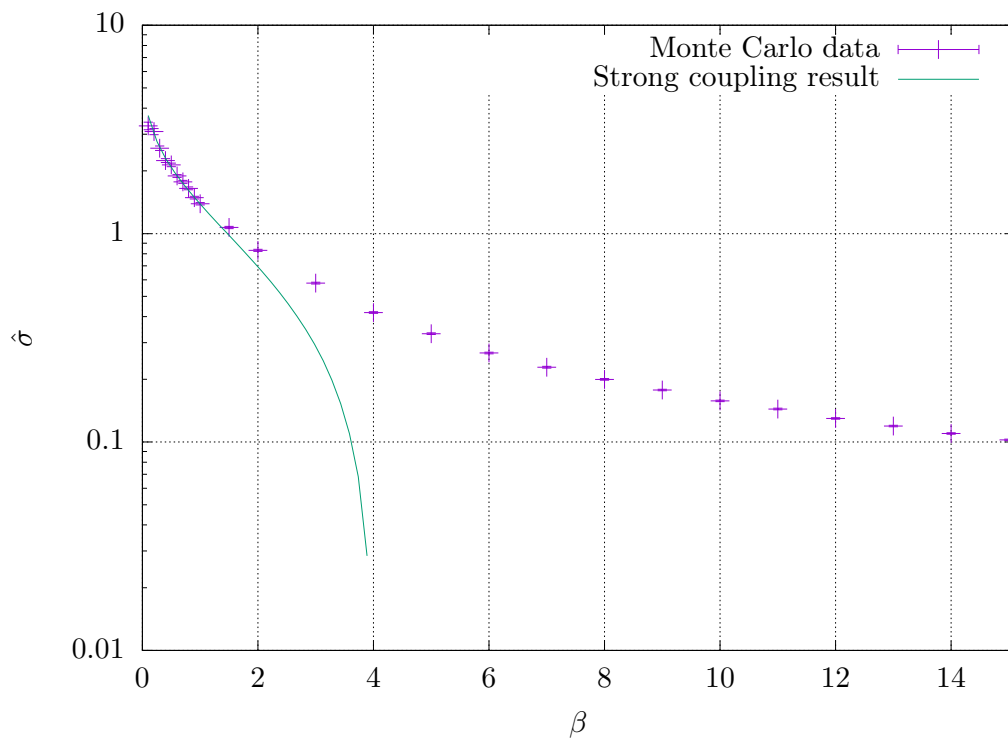


Figure 1: String tension, in lattice units, as a function of the (inverse) coupling. Computed on a  $10 \times 10$  lattice.