Key Performance Indicator for Topological Charge Diffusion on Lattice

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1 Introduction

Paper [1] studies the diffusion of topological charge on lattices with open and periodic boundary conditions in systematic way. Two quantities, *diffusion coefficient*, $D(t)$ and *tunneling rate*, τ_{tunn}, are used to measure how fast topological charge evolves during the Monte Carlo. These two quantities, which are defined in equation (1) of this document and in equation (4.2) of [1], could be used as a set of key performance indicators(KPI) for topological charge evolution. A baseline for comparison for future effort in fighting against the critical slowing down could be found in Table 3 in [1].

This document includes a manual of the package for measuring topological charge diffusion in the way described in [1]. We assume that the potential users have basic understanding of `bash` system and lattice QCD calculation. Statistical techniques used are discussed in the appendix since they are only briefly mentioned in the paper.

1.1 Original Paper

For an $SU(3)$ gauge theory *topological charge density* is defined as

$$\rho(\vec{x},t)=\frac{1}{32\pi^2}\epsilon_{\mu\nu\rho\lambda}\text{Tr}[F_{\mu\nu}(\vec{x},t)F_{\rho\lambda}(\vec{x},t)]$$

in the continuum limit. In the paper the topological charge density is measured using the "5Li" method[2] after wilson flow is run for a number of steps. This charge density is then summed on different time slices t and different configurations at molecular dynamic(MD) time τ,

$$Q(t,\tau)=\sum_{\vec{x}}\rho(\vec{x},t;\tau).$$

The next step is to construct the correlation function. The definition differs between lattices with open and periodic boundary condition.

1.1.1 Open Boundary Condition

The correlation function is defined as

$$C(t,t_0,\tau)\equiv\langle Q(t,\tau_0+\tau)Q(t_0,\tau_0)\rangle,$$

where the average is taken over all possible τ_0's.

1.1.2 Periodic Boundary Condition

The correlation function is defined as

$$C(\Delta t, \tau) \equiv \langle Q(t_0, \tau_0 + \tau) Q(t_0 + \Delta t, \tau_0) \rangle,$$

where the average is taken over all possible τ_0's and t_0's. This simplified version is used since we can utilize the translation invariance in the t-direction.

With the correlation function constructed a diffusion model is proposed,

$$\frac{\partial}{\partial \tau} C(t, t_0, \tau) = \frac{\partial}{\partial t} \left[D(t) \frac{\partial}{\partial t} C(t, t_0, \tau) \right] - \frac{1}{\tau_{\text{tunn}}} C(t, t_0, \tau) \tag{1}$$

for lattices with open boundary condition. The parameters in this model are $D(t)$ and τ_{tunn}. Intuitively $D(t)$ tells how fast the topological charge moves through the lattice and τ_{tunn} measures the rate at which topological charge is created.

For lattices with periodic boundary condition a replacment of $C(t, t_0, \tau)$ with $C(t, \tau)$ is understood and $D(t)$ is set to be a constant.

The paper finds this model fits well with the data. The fit is done by least square method and the error of the results are estimated with jackknife method[1].

2 Package

The package is available at https://github.com/RBC-UKQCD/DBW2-KPI. In order to run wilson flow and measure topological charge `CPS`(Columbia Physics System) is needed: https://github.com/RBC-UKQCD/cps-DBW2-KPI.

3 Install

3.1 CPS

We will need `gcc` and `MPI` to compile `CPS`. Make sure your `$PATH` includes `MPI` before compling.

1. Clone `CPS` from the git repository. Through out this document we will call this directory `CPS_REPO`.

2. Go to the top of the repository, `CPS_REPO`. `CPS` will be installed and placed under a directory specified as a variable `$cps`(line 4) in `CPS_REPO/conf.sh` so change that if you want. The default directory is `../cps-build/public/`. Through out this document we will call this directory `CPS_BUILD`.

3. Execute the `build.sh` script.

[1] See appendix for description of the jackknife method and the variance of the correlation function.

3.2 Analysis

We will need `g++`, `python`(version `2.7.5` tested) and `matlab`(version `R2013a (8.1.0.604)64-bit (glnxa64)` tested), later versions should be fine) to perform the whole analysis. Make sure these softwares are included in your `$PATH`.

1. Clone `DBW2-KPI` from the git repository. Through out this document we will call this directory `DBW2-KPI_REPO`.

2. Copy the file `CPS_BUILD/build-qmp/Makefile.users` to both `DBW2-KPI_REPO/ens_gen/` and `DBW2-KPI_REPO/wflow_tcharge_cps/` as `makefile`, i.e.

```
$ cp CPS_BUILD/build-qmp/Makefile.users DBW2-KPI_REPO/ens_gen/makefile
$ cp CPS_BUILD/build-qmp/Makefile.users DBW2-KPI_REPO/wflow_tcharge_cps/
    makefile
```

4 Usage

There are two stages when using this package. In the first stage `CPS` reads in the lattice configurations in standard *nersc* format, run wilson flow and measure the topological charge density. The second stage process the raw data produced by `CPS` and perform the fit.

4.1 First Stage

1. Go to `DBW2-KPI_REPO/wflow_tcharge_cps/binaries`.

2. Information of the incoming configurations are supposed to be given from line 100 to line 119. You might want to change these information to match your configurations. The comment there should be quite self-explaining.

3. `make`.

4. Use `mpirun -np $num_node NOARCH.x -qmp-geom $x $y $z $t` to run the job. You should change `$num_node` to match your environment. `$x`, `$y`, `$z` and `$t` are number of nodes to be allocated in each dimensions. The product of these four should be equal to `$num_node`. Examples are

```
mpirun -np 1 NOARCH.x -qmp-geom 1 1 1 1
mpirun -np 256 NOARCH.x -qmp-geom 4 4 4 4
```

5. The results will be placed under `DBW2-KPI_REPO/wflow_tcharge_cps/results/alg_wflow/`. The files would look like:

```
1    2.131295e+00
2    1.700994e+00    2.087967e+00    2.171048e+00    2.162269e+00    ...
3 AlgTcharge:
4 nleaf : 5
5    0 : 1x1
```

```
6    1 : 1x2
7    2 : 2x2
8    3 : 3x3
9    4 : 1x3
10 0 0 :    6.235056e-01
11 1 1 :    4.098351e-01
12 2 2 :    1.044206e-01
13 3 3 :    7.561517e-03
14 4 4 :    2.090085e-01
15 0 :    2.519808e-01   -1.512013e-01   -2.445761e-01    1.517083e-02    ...
16 1 :    8.713487e-02   -9.518725e-03   -2.284069e-02    3.067389e-03    ...
17 2 :    4.226348e-03   -4.861770e-03    9.862970e-03   -1.161517e-04    ...
18 3 :   -1.795864e-03    2.317363e-03   -9.817474e-04    1.049183e-03    ...
19 4 :    2.337879e-02    1.922419e-02    2.056840e-02   -6.668813e-03    ...
20 AlgWilsonFlow: dt = 5.000000e-02
```

The "5 Li" method can be found in [2]. Here in the outputfile line 5 shows number of leaves we use to calculate the topological charge density, line 5 through 9 show the size of the leaves, line 10 through 14 show the values of the different pieces of the total topological charge. line 15 through 19 show the same number but local on different time slices.

4.2 Second Stage

1. Go to `DBW2-KPI_REPO`.
2. The input information of this stage is specified in `DBW2-KPI_REPO/conf.sh`. Specify all the information in that file.
3. execute `DBW2-KPI_REPO/do.sh`.
4. The final result will be placed under `DBW2-KPI_REPO/correlation_to_fit/results/`.

4.3 Ensemble Generation

`DBW2-KPI_REPO/ens_gen/binaries` contains the code to generate open/periodic boundary condition pure gauge lattice with DBW2 action. Before compiling you should specify the information of the lattice you want to generate:

1. Lattice size is specified as `int x_sites`, `int y_sites`, `int z_sites` and `int t_sites` in `DBW2-KPI_REPO/ens_gen/vmls/do_arg.vml`.
2. t-direction boundary condition is specsified at line 94 of `DBW2-KPI_REPO/ens_gen/binaries/main.C`.
3. Number of steps in one trajectory and step size are specified as `int steps_per_traj` and `double step_size` in `DBW2-KPI_REPO/ens_gen/vmls/hmc_arg.vml`.

4. β and c_1 of the gauge action are specified as `double beta` and `c_1` in `DBW2-KPI_REPO/ens_gen/vmls/do_arg.vml`. The default value is $c_1 = -1.4088$ for DBW2 action.

5. You are probably safe to use all the other default values.

Now just go to `DBW2-KPI_REPO/ens_gen/binaries`, do `make` and execute the `NOARCH.x` in the same way mentioned in the first stage.

The topological charge density data is automatically generated and placed under `DBW2-KPI_REPO/ens_gen/results/alg_wflow/`. From here you can directly go to the second stage.

Appendices

A Statistics

A.1 Jackknife Method[3]

Let $[X_1, \cdots, X_n]$ be n (*independent and identically distributed*) random variables. We split this sample into g groups of size h each, $n = gh$[2]. Let T be an estimator of some parameter θ based on sample size n, T_{-i} be the corresponding estimator based on the sample of size $(g-1)h$, where the i-th group of size h has been excluded:

$$T = f[X_1, \cdots, X_n],$$

$$T_{-i} = f[X_1, \cdots, X_{(i-1)h}, X_{ih+1}, \cdots, X_n].$$

Write $\bar{T}_\bullet$ as the mean of the T_{-i}'s, the jackknife estimator is defined as

$$T_j = gT - (g-1)\bar{T}_\bullet, \ \bar{T}_\bullet = \langle T_{-i} \rangle_i = \frac{1}{g}\sum_{i=1}^{g} T_{-i},$$

and the estimator of the variance of the T_j is

$$\mathrm{var}[T_j] = \frac{g-1}{g}\sum_{i=1}^{g}(T_{-i} - \bar{T}_\bullet)^2.$$

Usually we use T instead of T_j, since they have the same expectation value.

A.2 Variance of the Correlation Function

A.2.1 Open Boundary Condition

Correlation function on lattices with open boundary condition is defined as

$$C(t, t_0, \tau) \equiv \langle Q(t, \tau_0 + \tau) Q(t_0, \tau_0) \rangle.$$

[2] If h is longer than the autocorrelation time of the random variables it seems the requirement that the random variables be *independent and identically distributed* can be skipped.

A natural estimator is:

$$\hat{C}(t, t_0, \tau) = \langle Q(t, \tau_0 + \tau) Q(t_0, \tau_0) \rangle_{\tau_0}.$$

Variance of $\hat{C}(t, t_0, \tau)$ is roughly[3]

$$\mathrm{var}[\hat{C}(t, t_0, \tau)] = \langle C(t, t, \tau_0) C(t_0, t_0, \tau_0) + C(t, t_0, \tau_0 + \tau) C(t_0, t, \tau_0 - \tau) \rangle_{\tau_0} .$$

A.2.2 Periodic Boundary Condition

Correlation function on lattices with periodic boundary condition is simplified as

$$C(\Delta t, \tau) \equiv \langle Q(t_0, \tau_0 + \tau) Q(t_0 + \Delta t, \tau_0) \rangle.$$

A natural estimator is:

$$\hat{C}(\Delta t, \tau) = \langle Q(t_0, \tau_0 + \tau) Q(t_0 + \Delta t, \tau_0) \rangle_{\tau_0, t_0}.$$

The corresponding variance is roughly

$$\mathrm{var}[\hat{C}(\Delta t, \tau)] = \left\langle C(t_0, \tau_0)^2 + C(t_0 + t, \tau_0 + \tau) C(t_0 - t, \tau_0 - \tau) \right\rangle_{\tau_0, t_0} .$$

References

[1] G. McGlynn, and R. D. Mawhinney, Phys. Rev. D **90**, 1 (2014), [arXiv:1406.4551].

[2] P. de Forcrand, M. García Pérez, and I.-O. Stamatescu, Nucl. Phys. B **499**, 409 (1997), [arXiv:hep-lat/9701012].

[3] R. G. Miller, Biometrika **61**, 1 (1974).

[4] N. Madras, and a. D. Sokal, J. Stat. Phys. **50**, 109 (1988).

[3] This is true if mean value of $Q(t, t_0, \tau)$ is zero and its the fourth joint cumulant κ_4 is also zero.[4] The mean value of the topological charge density is zero while we are assuming κ_4 to be zero.