

Bipartite Entanglement Entropy

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Quasiperiodic Kicked Rotor

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- This has the drawback of increasing computational complexity of each individual step and the memory used at any given time is large.

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- This is better as the memory used is less but computation increases. Since fourier transforms are computationally cheap anyway, so it's fine.
- Peak memory required scales the same way but we have reduced it by a constant factor and it is not used in all calculations.

We use $\hbar = 2.85, \omega_2 = 2\pi\sqrt{5}, \omega_3 = 2\pi\sqrt{13}$, the momentum ranges from -10 to 10

$$H = \frac{p_1^2}{2} + p_2\omega_2 + p_3\omega_3 + K\cos(\theta_1)(1 + \alpha\cos(\theta_2)\cos(\theta_3)) \sum_n \delta(t - n)$$

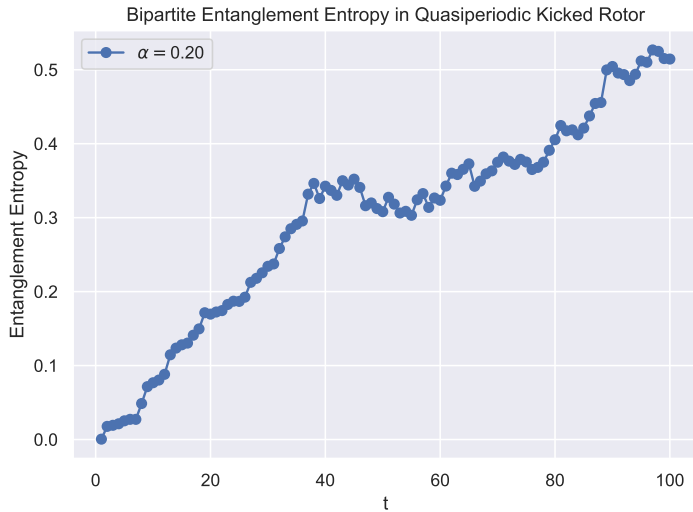


Figure 1: Precritical (Insulator): $K = 4, \alpha = 0.2$

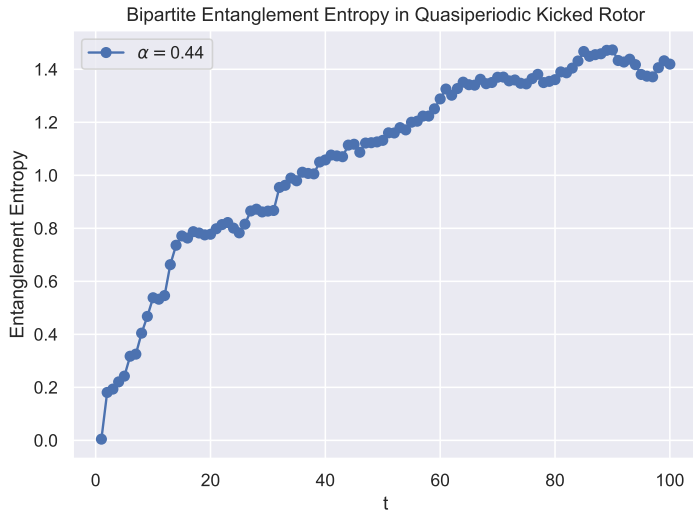


Figure 2: Critical: $K = 6.36, \alpha = 0.4375$

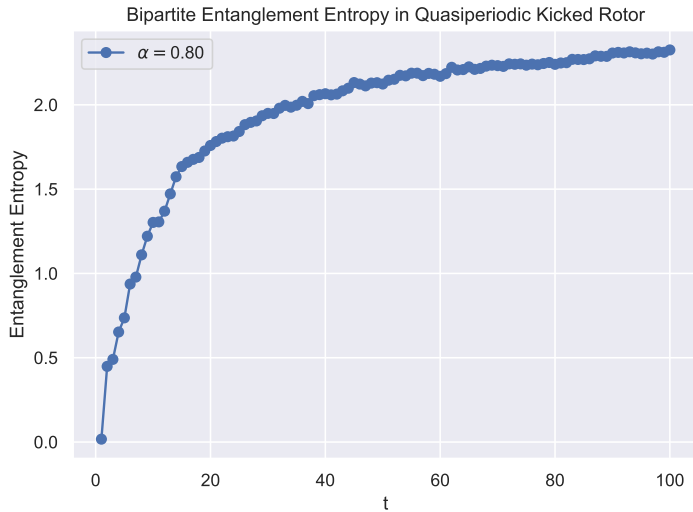


Figure 3: Post-critical (Metal): $K = 8, \alpha = 0.8$

- I don't see much of a trend here. The entanglement grows faster and higher with higher K values i.e. more diffusive the regime higher the entanglement for the same number of time steps but other than that, I don't see anything here.

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 - $S = -\rho_1 \ln(\rho_1)$

Momentum (p_1) distributions

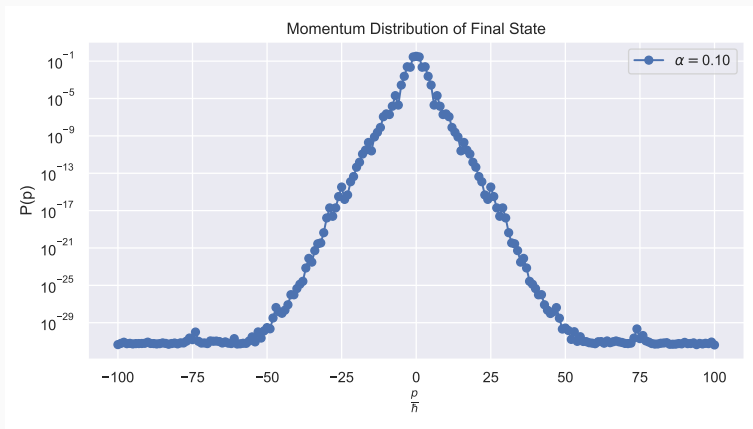


Figure 4: $K = 3$, $\alpha = 0.1$

Momentum (p_1) distributions

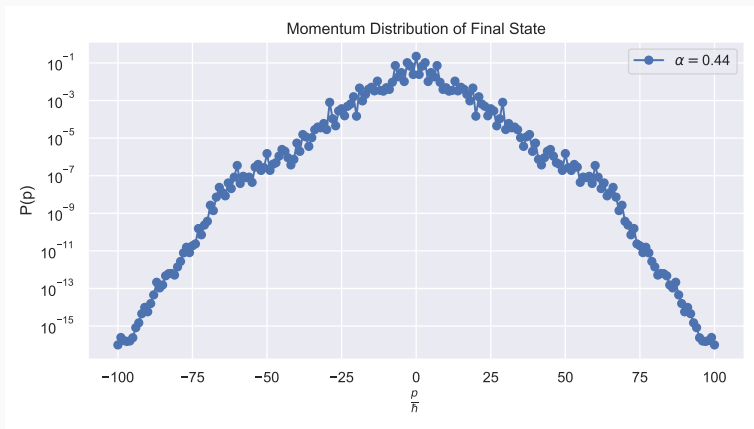


Figure 5: $K = 6.36$, $\alpha = 0.4375$

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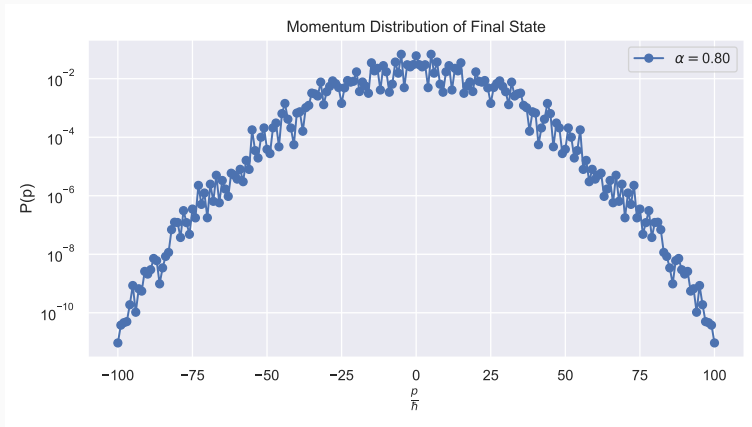


Figure 6: $K = 7$, $\alpha = 0.8$

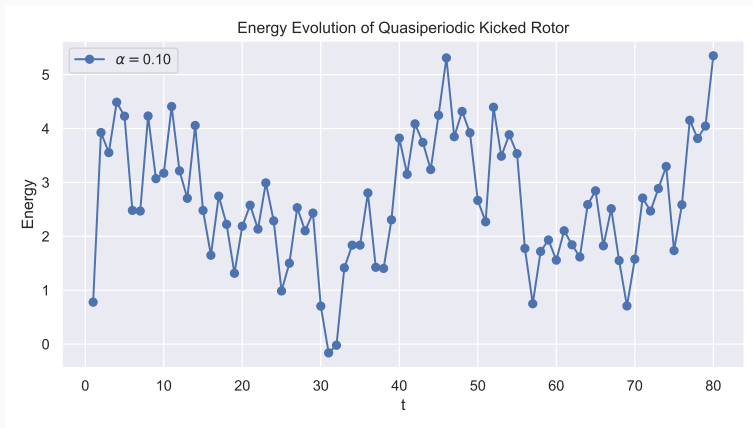


Figure 7: $K = 3$, $\alpha = 0.1$

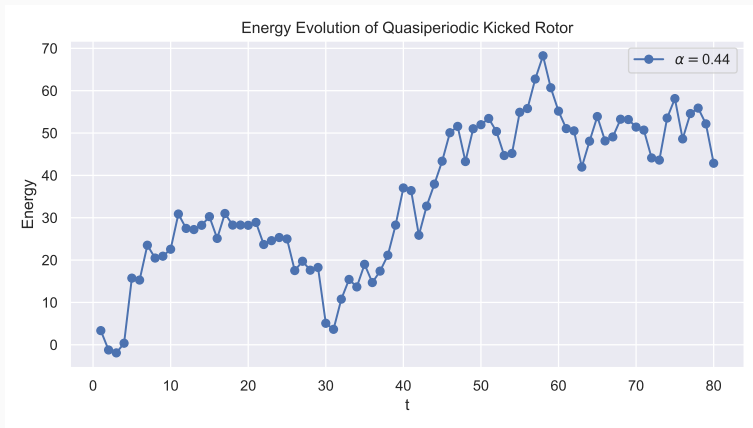


Figure 8: $K = 6.36$, $\alpha = 0.4375$

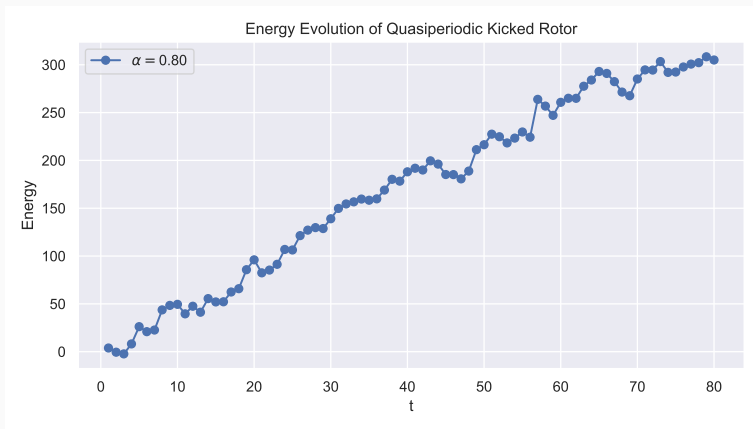


Figure 9: $K = 7$, $\alpha = 0.8$

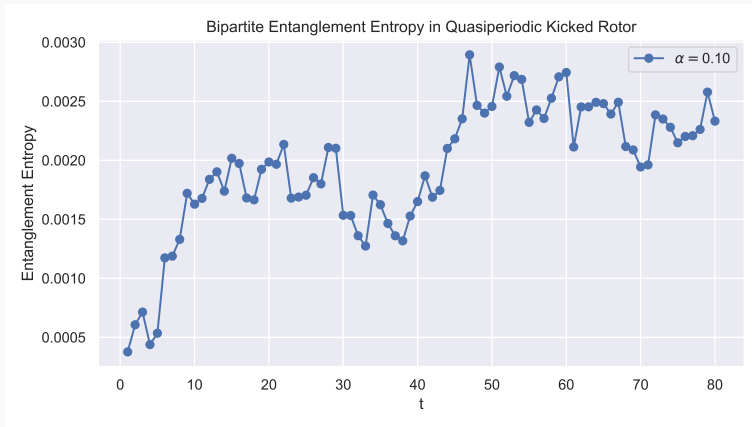


Figure 10: $K = 3$, $\alpha = 0.1$

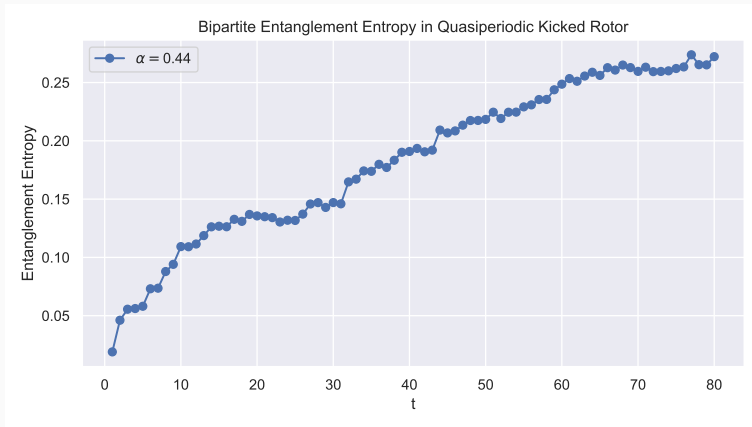


Figure 11: $K = 6.36$, $\alpha = 0.4375$

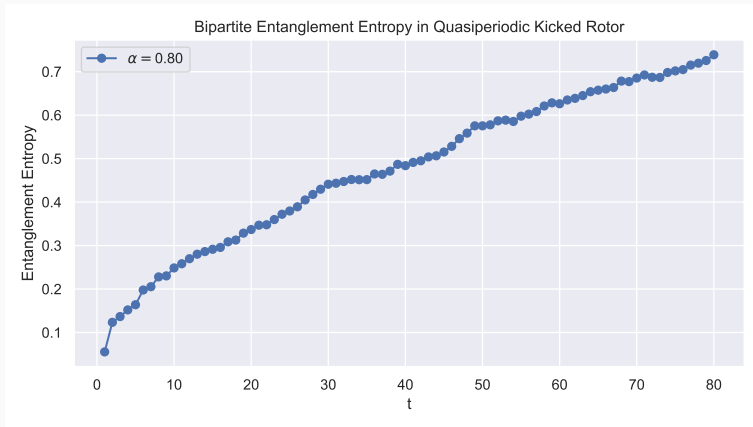


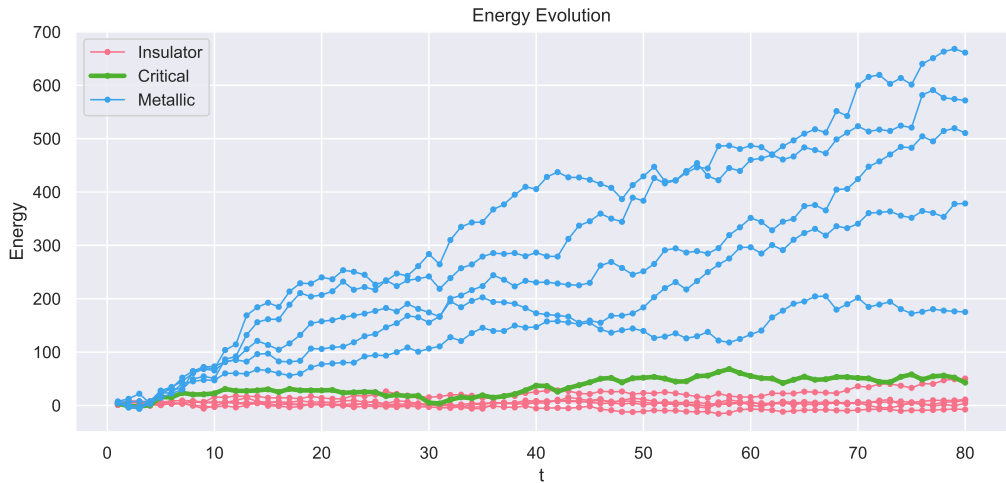
Figure 12: $K = 7$, $\alpha = 0.8$

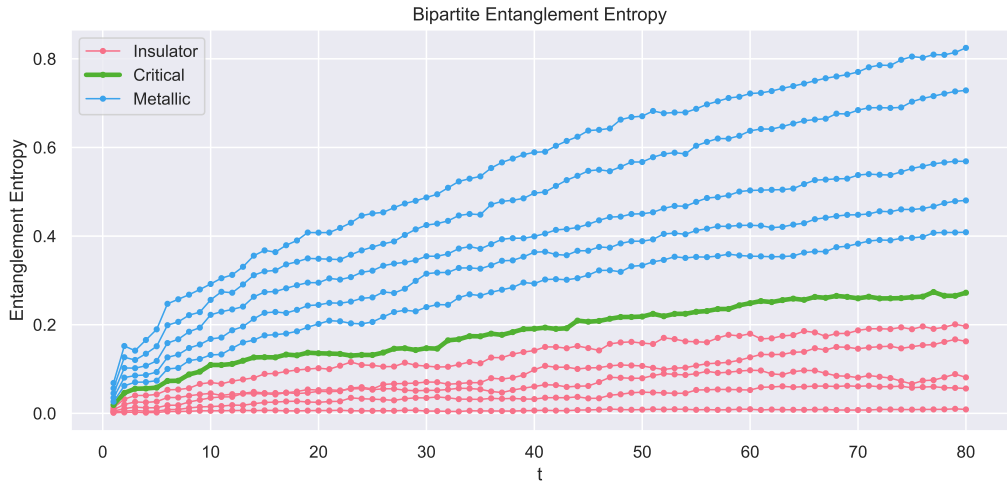
- We look at the following plots for the following at different K, α values from the insulator to the metallic regime:
 1. Energy Expectation Value
 2. Entanglement Entropy
 3. Momentum Distribution
- To study the changes in the energy and entropy values with K, we have plotted the following quantities:
 1. Entropy Difference: $S(K_{n+1}, \alpha_{n+1}) - S(K_n, \alpha_n)$ vs t
 2. Energy Difference: $E(K_{n+1}, \alpha_{n+1}) - E(K_n, \alpha_n)$ vs t

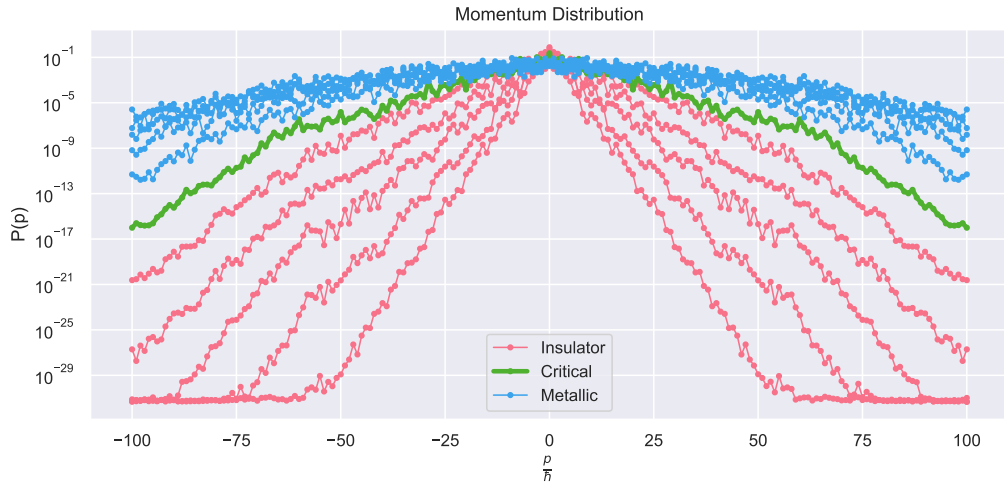
- First we take a big picture look with 11 values with $K \in [3.00, 9.72]$ and $\alpha \in [0.200, 0.6750]$. Both ranges are centred around the critical point¹. Momentum range is -100 to 100 with 80 timesteps.
- Then we look very close to the critical point with 11 values with $K \in [6.30, 6.42]$ and $\alpha \in [0.4000, 0.4750]$. Both ranges are centred around the critical point. Momentum range is -100 to 100 with 80 timesteps.

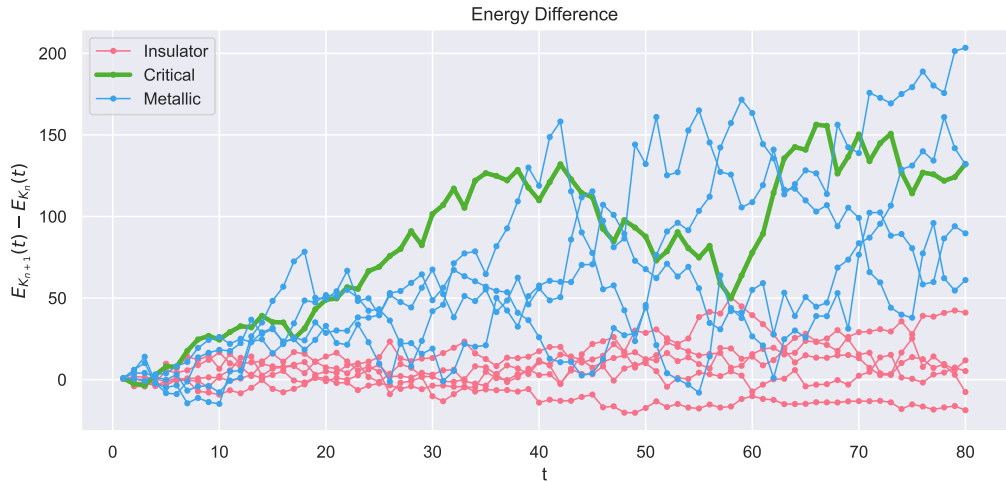
¹The (G. Lemarié, Grémaud, and Delande 2009) paper gives the value of K at critical point, but not of α .

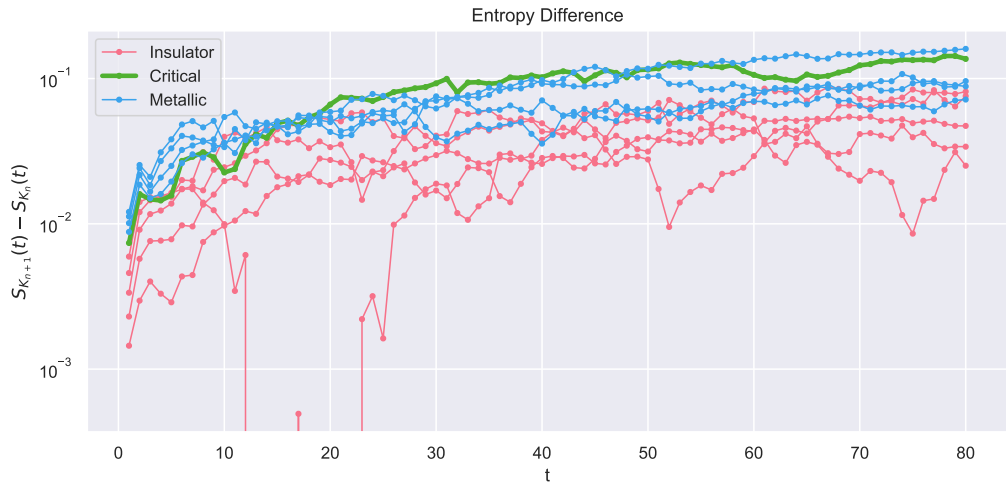
The Big Picture





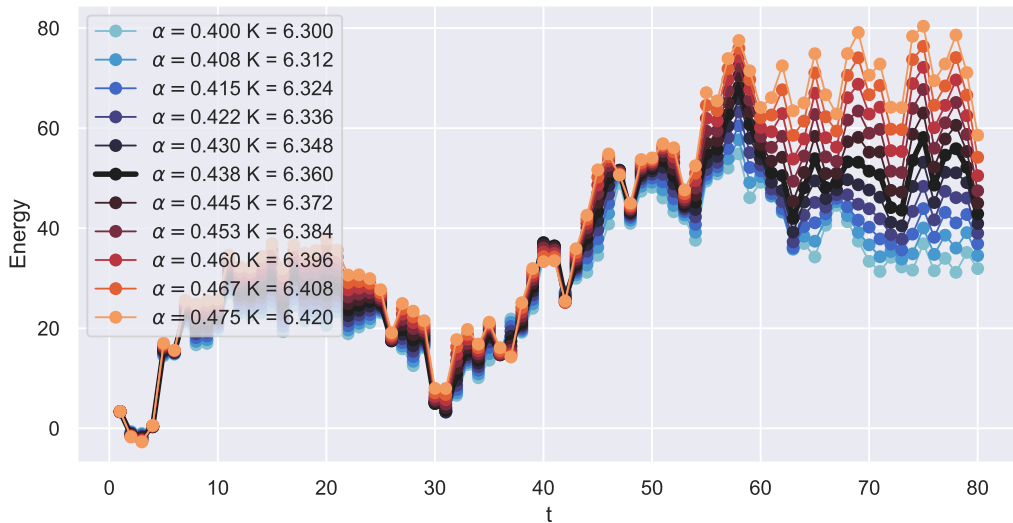




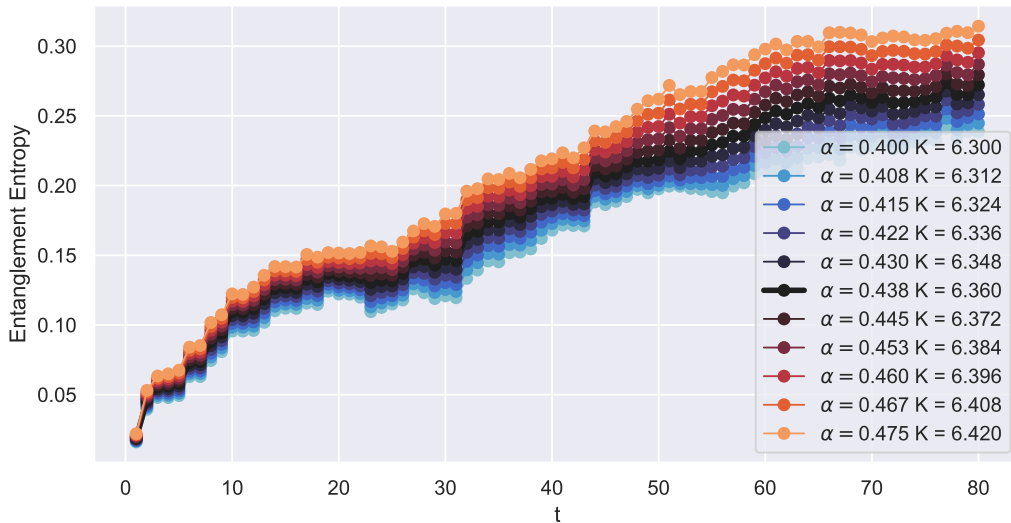


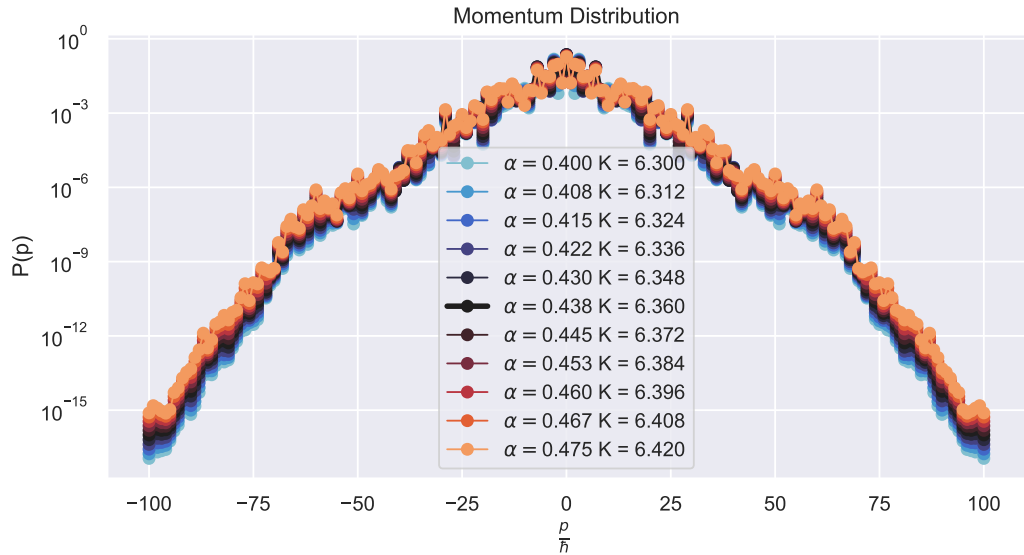
The Microscopic Picture

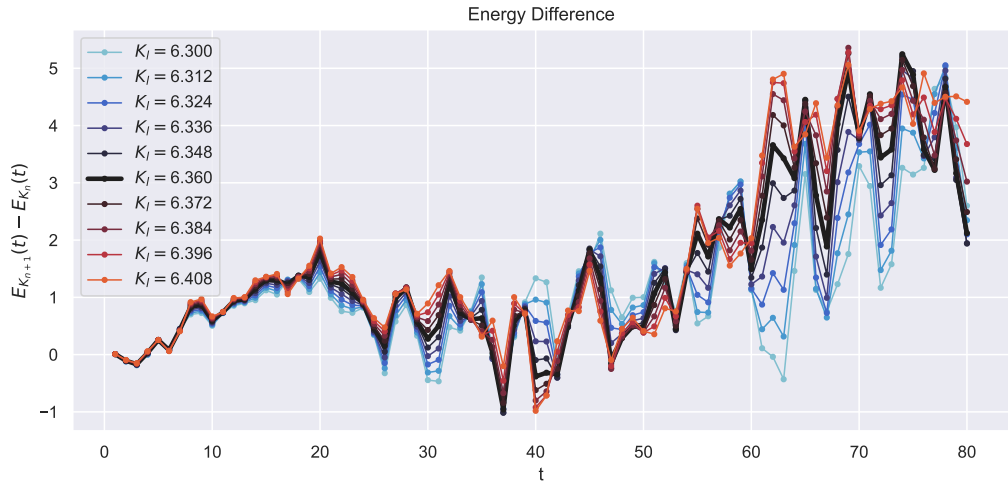
Energy Evolution

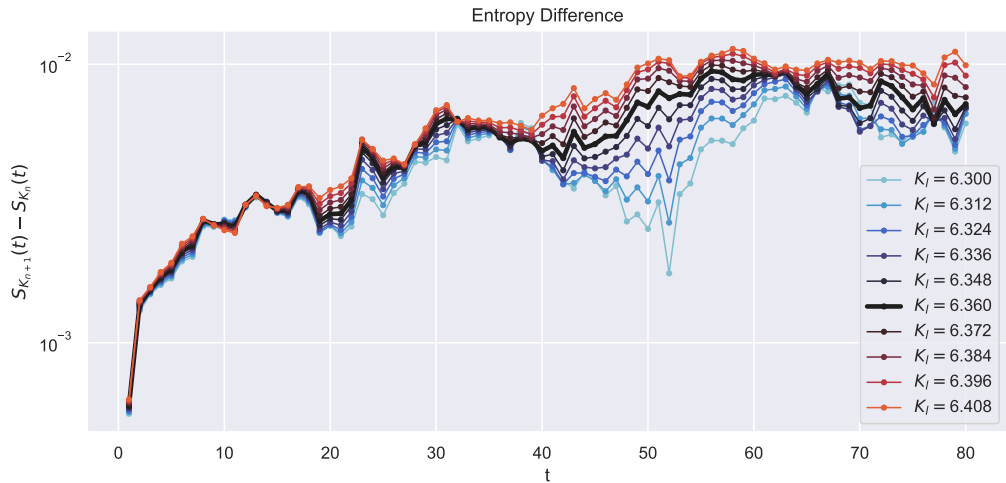


Bipartite Entanglement Entropy









- We use $\hbar = 1$, $\omega_2 = 2\pi\sqrt{5}$ and $\omega_3 = 2\pi\sqrt{13}$. The critical point of the metal insulator transition is known to occur at $K_c = 6.36 \pm 0.02$.

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- The value of α_c , however is not determined accurately, nor written in any literature as it doesn't form a part of the scaling function in a nice way, though it most certainly does matter as it generates the anisotropy. We determine the value of α_c roughly by using the fact that the authors of the cited papers traversed the $K - \alpha$ space along a straight line perpendicular to the transition parabola. (G. Lemarié, Grémaud, and Delande 2009) (Gabriel Lemarié et al. 2009)

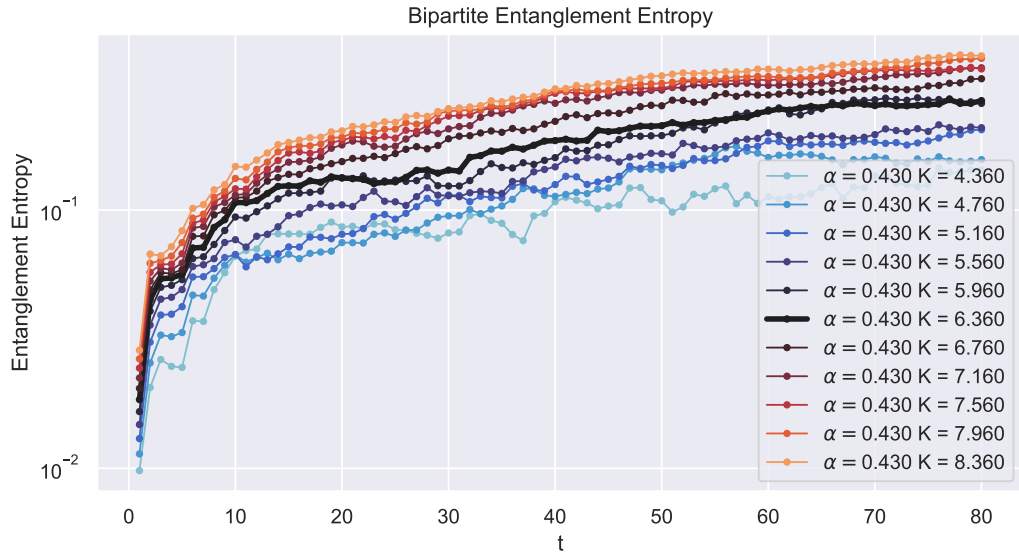
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- TL;DR we used $K_c = 6.36$ and $\alpha_c = 0.4303$.

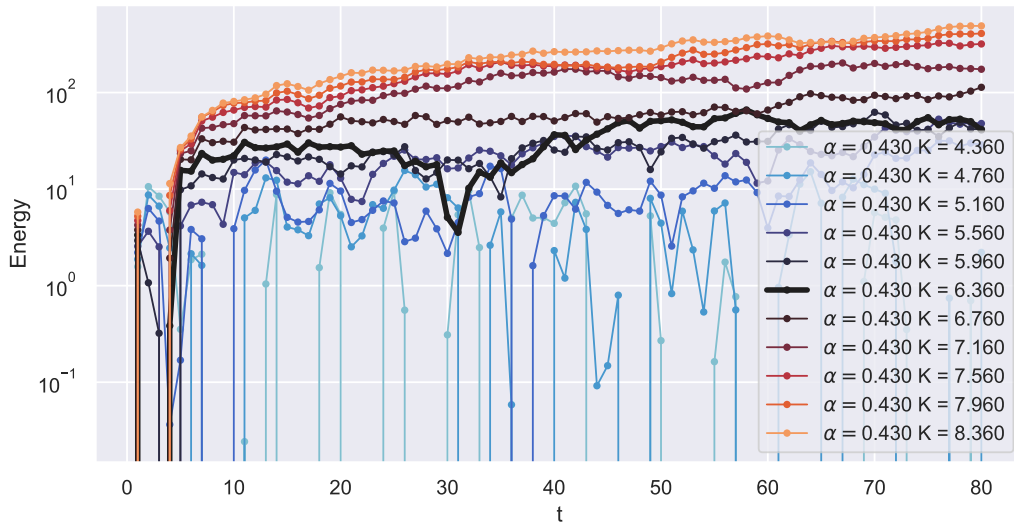
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- We used a basis size of 201 (-100 to 100) for each of the 3 coordinates and the simulations were done for 80 timesteps.

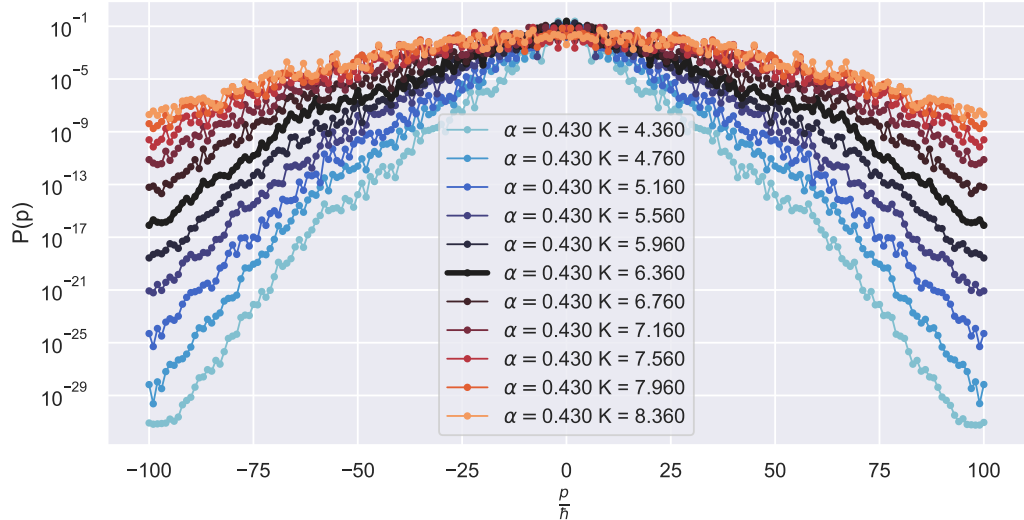
Varying K

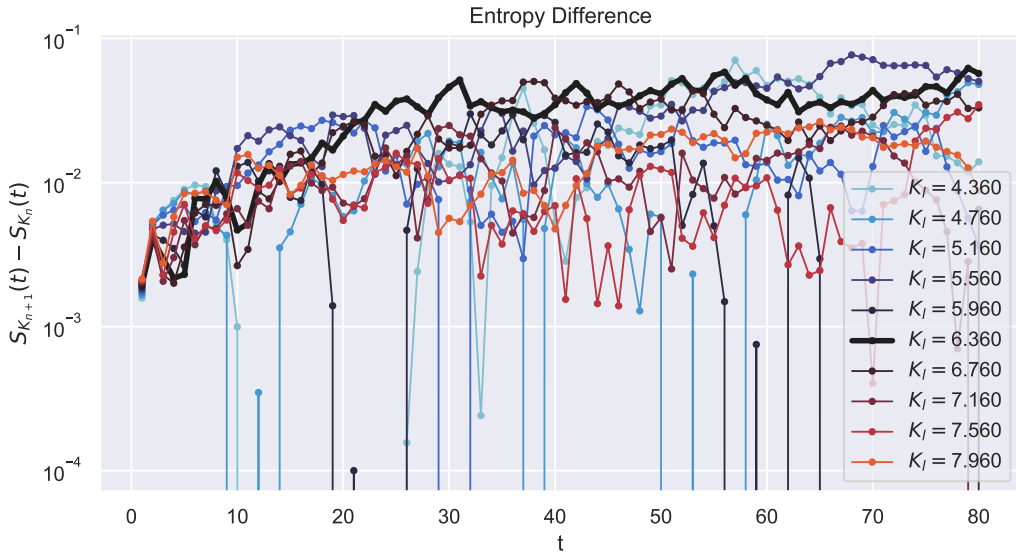


Energy Evolution

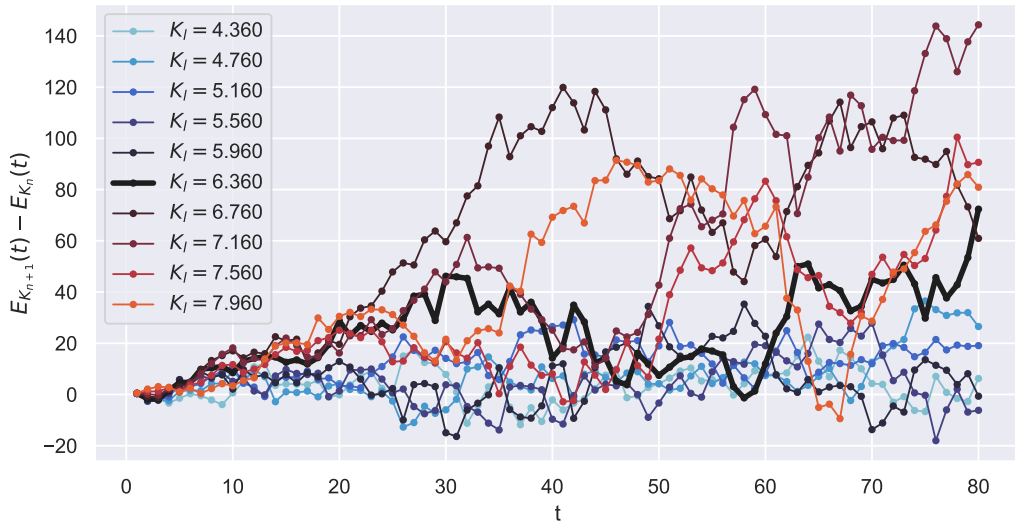


Momentum Distribution



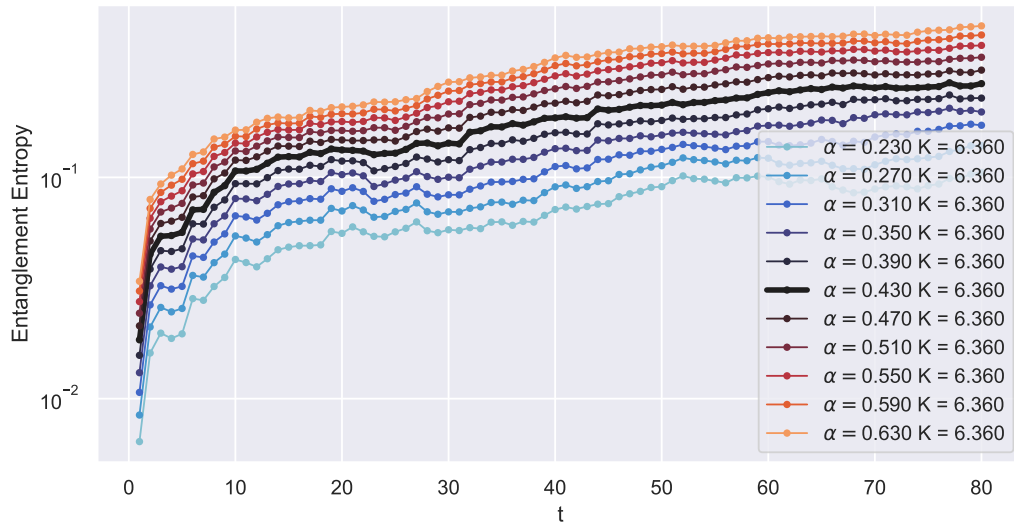


Energy Difference

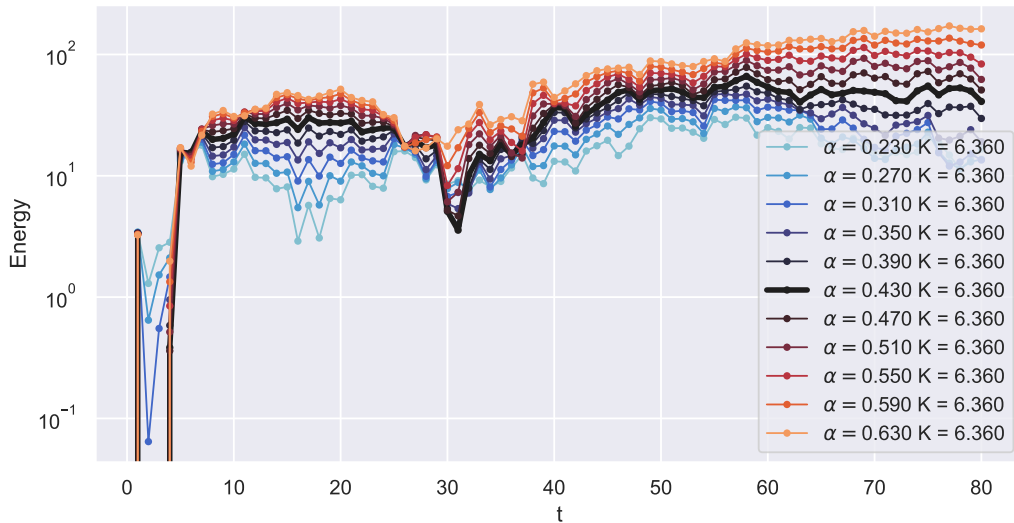


Varying α

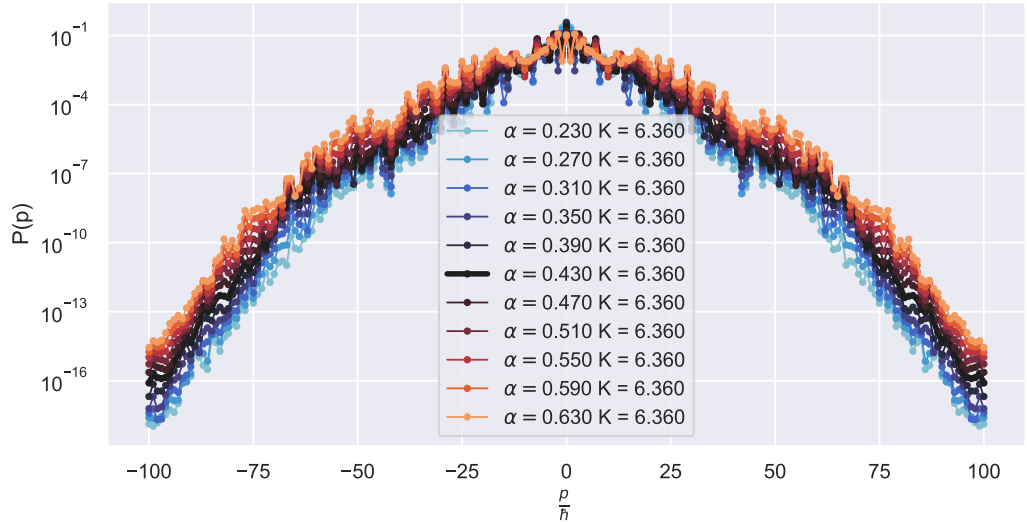
Bipartite Entanglement Entropy



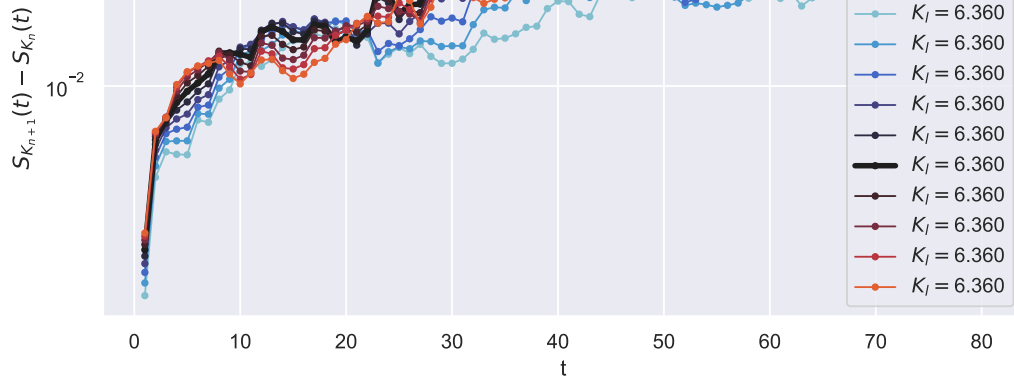
Energy Evolution



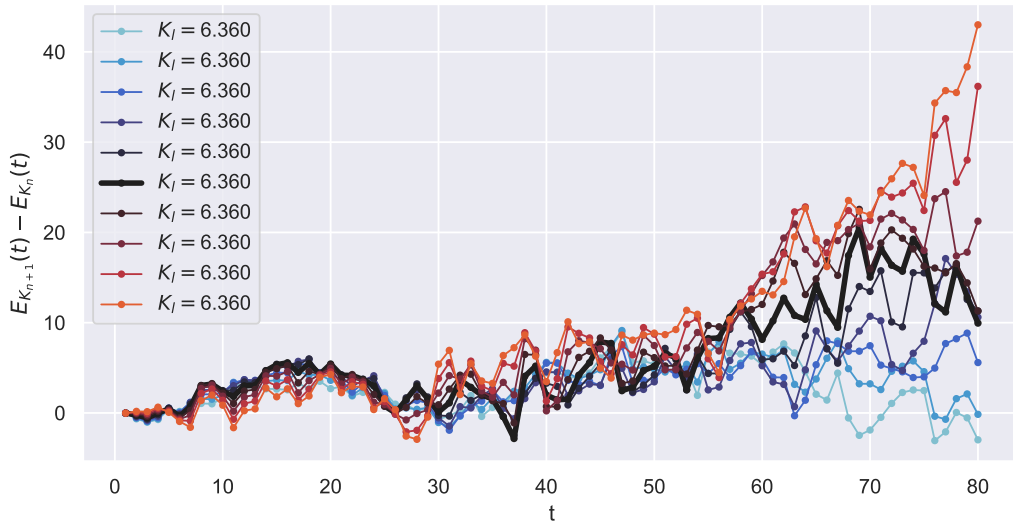
Momentum Distribution



Entropy Difference

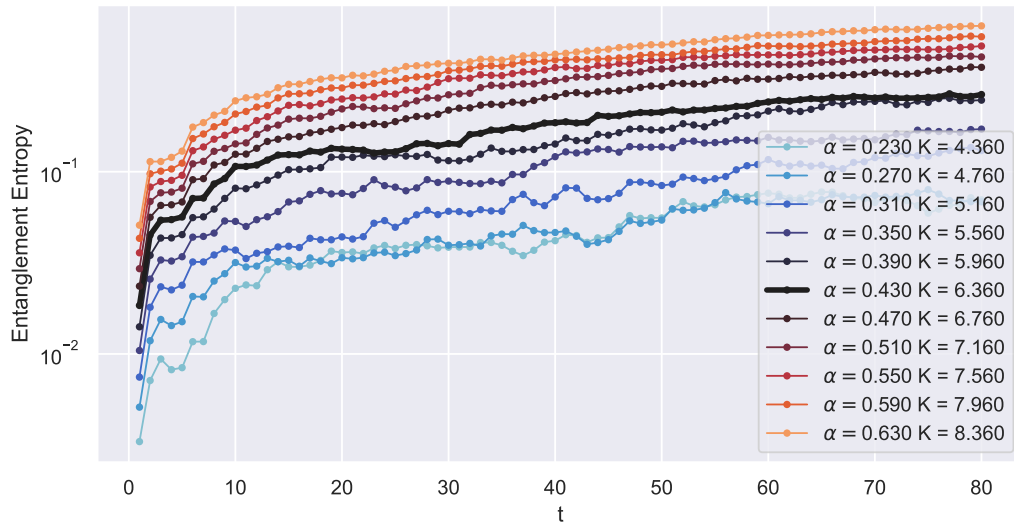


Energy Difference

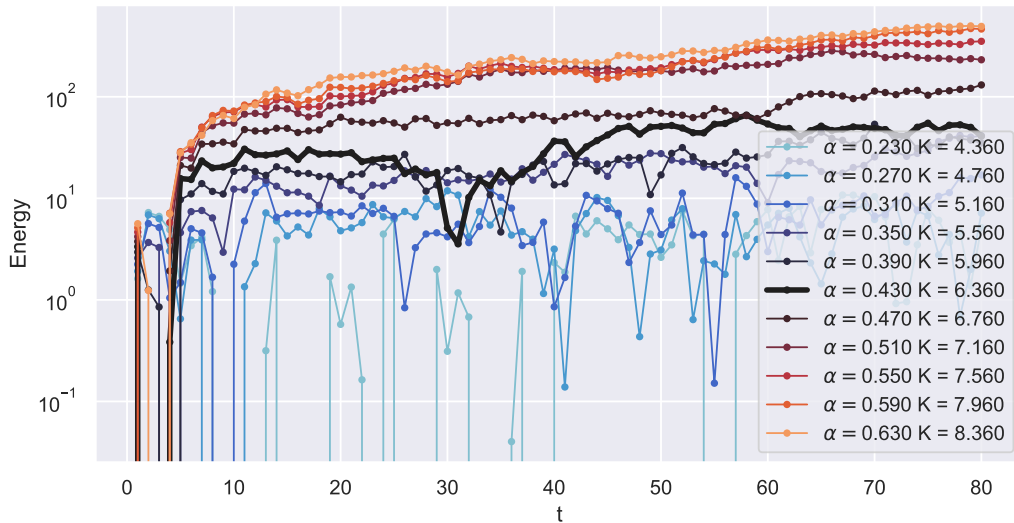


Varying K and α

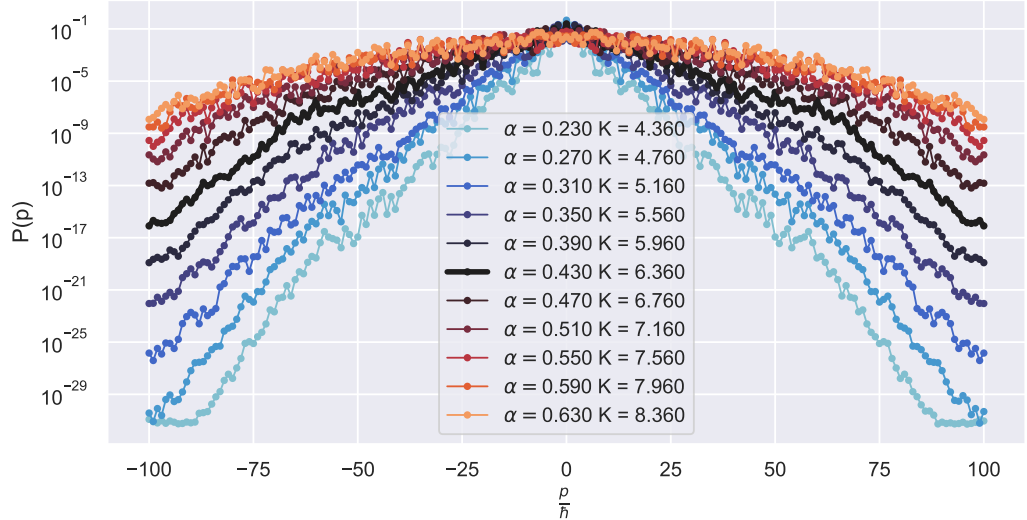
Bipartite Entanglement Entropy

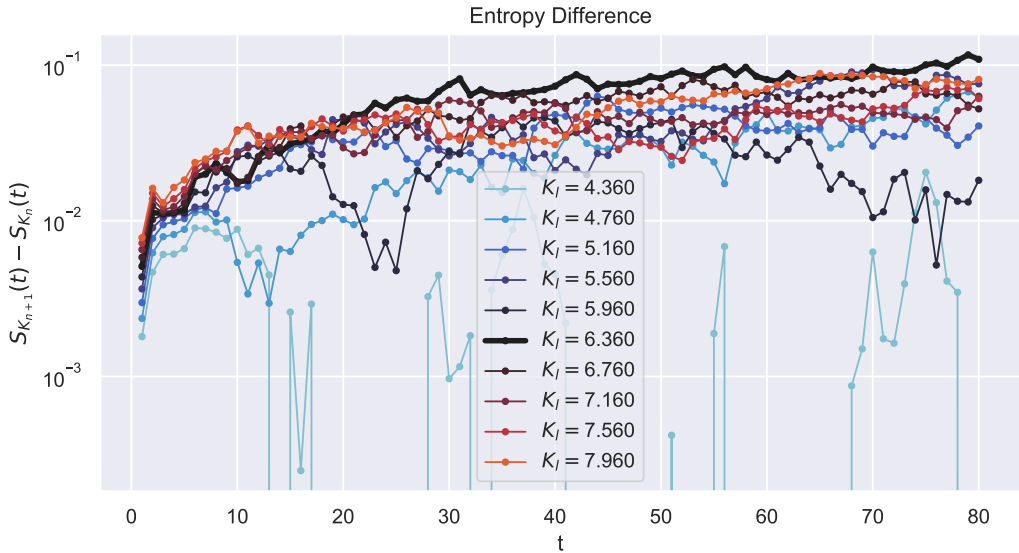


Energy Evolution

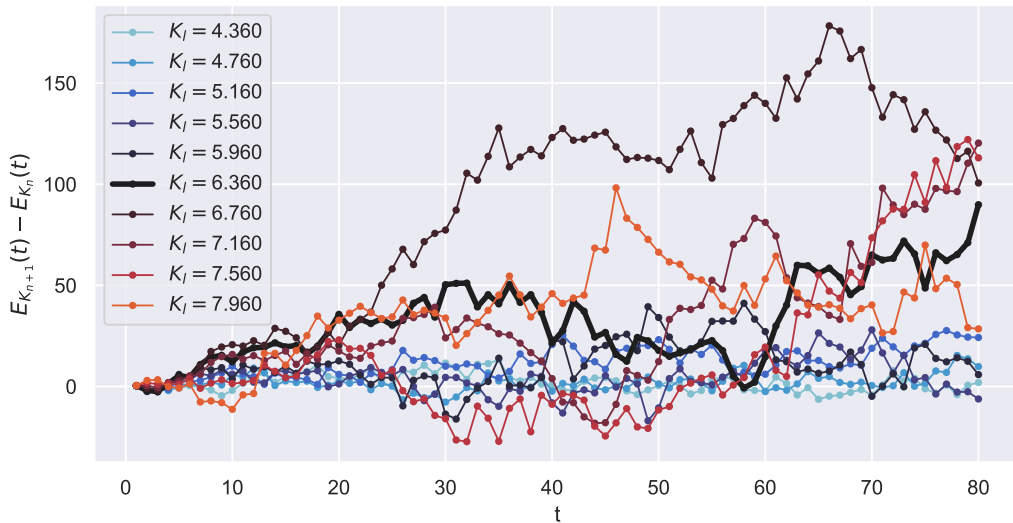


Momentum Distribution





Energy Difference



- Lemarié, Gabriel, Julien Chabé, Pascal Szriftgiser, Jean Claude Garreau, Benoît Grémaud, and Dominique Delande. 2009. "Observation of the Anderson Metal-Insulator Transition with Atomic Matter Waves: Theory and Experiment." *Phys. Rev. A* 80 (4): 043626.
<https://doi.org/10.1103/PhysRevA.80.043626>.
- Lemarié, G., B. Grémaud, and D. Delande. 2009. "Universality of the Anderson Transition with the Quasiperiodic Kicked Rotor." *Europhys. Lett.* 87 (3): 37007.
<https://doi.org/10.1209/0295-5075/87/37007>.