The Journey of Simulating Two Excitation Heisenberg Spin Chains

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Background: Heisenberg Spin Chains







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Research Goals – Finding Stability

1. Develop simulations that yield <u>qualitative</u> understanding and can handle many spins over long lifetimes.

2. Generate a new analytical model for <u>quantitatively</u> understanding two excitation stability.

Properties of Localized Excitations

for Systems with Two Excitations Propagating



Qualifying Localized States' Stability

Entanglement Propagation of Two Excitations on XXZ Heisenberg Spin Chains





QCTF – Analytically Propagating



Single Excitation QCTF Outcome

$$\tilde{\mathcal{Q}}_{q}(s) = \frac{1}{N^{4}} \sum_{\substack{m_{1},m_{2} \\ m_{3},m_{4} \\ p \neq q}} \left(s - \frac{i\mathbf{J}}{\hbar} \left(\cos\left(\frac{2\pi}{N}m_{2}\right) - \cos\left(\frac{2\pi}{N}m_{1}\right) - \cos\left(\frac{2\pi}{N}m_{4}\right) + \cos\left(\frac{2\pi}{N}m_{3}\right) \right) \right)^{-1} e^{\frac{2i\pi}{N} \left(q(m_{1}-m_{3})+p(m_{4}-m_{2})\right)}.$$

Two Excitation QCTF – Central Insights

- a) Maintain same entanglement measure as 1 excitation QCTF by parameterizing momenta eigenstates.
- b) Use 4 orthogonal structural



c) Adjust measure to recreate final cancellations.

Two Excitation QCTF – Analysis

Current Entanglement formulation

$$\tilde{Q}_{M}(s) = \sum_{1 \le \alpha < \beta}^{N} \sum_{1 \le \varsigma < \delta}^{N} \left(4x \sum_{\lambda_{1} \le \lambda_{2}}^{N-1} \right) \left[s - J \frac{i}{\hbar} (E_{1} - E_{2}' - E_{3}'' + E_{4}''') \right]$$
$$a(\alpha, \beta) a^{*}(\psi_{0}) a'^{*}(\varsigma, \delta) a'(\psi_{0}) a''^{*}(\alpha, \beta) a''(\psi_{0}) a'''(\varsigma, \delta) a'''(\psi_{0}) a'''(\varsigma, \delta) a'''(\psi_{0}) a'''(\varsigma, \delta) a'''(\varphi_{0}) a'''(\varsigma, \delta) a''(\varphi_{0}) a'''(\varphi_{0}) a''''(\varphi_{0}) a'''(\varphi_{0}) a'''(\varphi_{0}) a'''(\varphi_{0}) a'''(\varphi_{0}) a''''(\varphi_{0}) a'''(\varphi_{0}) a''''(\varphi_{0}) a''''(\varphi_{0}) a''''(\varphi_{0}) a''''(\varphi_{0}) a''''(\varphi_{0}) a'''''(\varphi_{0}) a''''(\varphi_{0}) a''''(\varphi_{0}) a'''''(\varphi_{0}) a'''''(\varphi_{0}) a'''''''(\varphi_{$$

Points of improvement for future formulations

- 1. High level analysis is required.
- 2. Advanced numerical programming is required.
- 3. Eigenstates are inseparable in this formulation.

Conclusion

- 1. Algorithmically solved for two excitation spin chain dynamics regardless of chain size.
- 2. Simulated spin chain evolution using numerical eigenstate solutions and the Schrödinger equation.
- 3. Found a QCTF formulation that reveals entanglement dynamics analytically

Future Work

- 1. Find an appropriate simplification of QCTF for analyzing bound states.
- 2. Analyze a two-particle correlation measure using QCTF methods.

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Citation

- (1) Ball, P. First 100-Qubit Quantum Computer Enters Crowded Race. *Nature* 2021, 599, 542. URL: https://www.nature.com/articles/d41586-021-03476-5 (accessed 2023-07-27)
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