

**Qubit transfer with high fidelity in 1D
fermion spin chains with nearest-
neighbour interaction based on new
recurrence relations for Racah polynomials**

Aynura M. Jafarova

E-mail: aynure.jafarova@gmail.com

ANAS Institute of Mathematics and Mechanics, Baku, Azerbaijan

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Levels of interconnection



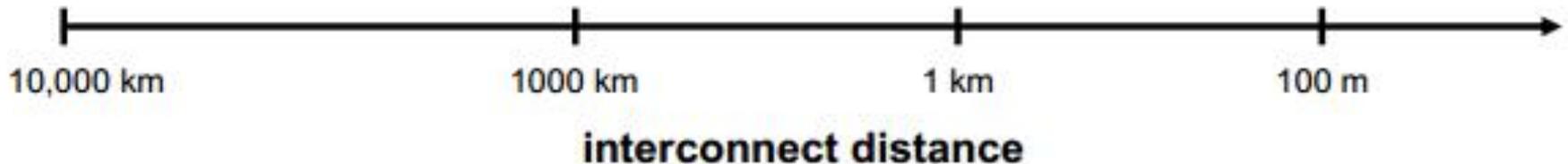
Telecommunications



Campus networks



LANs

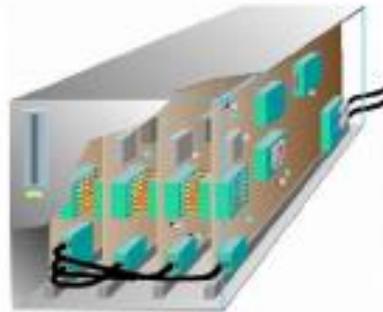


Optics currently dominates for long distance interconnects
Increasingly, optics is used in local area network applications

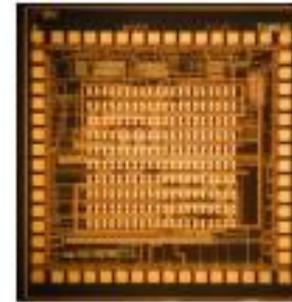
Levels of interconnection



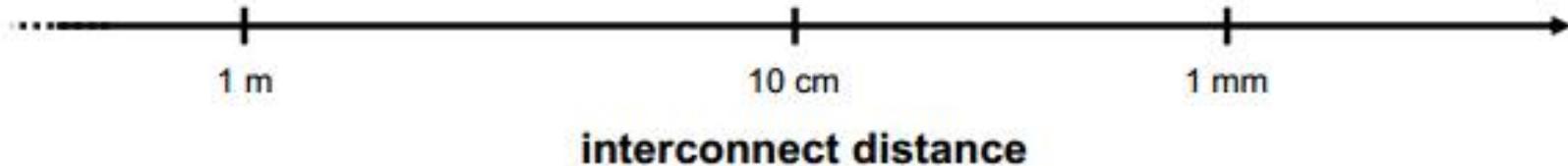
Backplanes & board-to-board



Chip-to-chip



On-chip



Electrical signaling within computers is encountering severe limitations - what can help at these $\ll 1\text{mm}$ length scales?



What is a quantum computer?

- A quantum computer is a machine that performs calculations based on the laws of quantum mechanics, which is the behavior of particles at the sub-atomic level.

Classical information

Stored as string of bits 0100110

Manipulation of (qu)bits (computation, dynamics)

Bit transformations (function computation)
All functions can be computed reversibly.

Bit states can be copied.

Transmission of (qu)bits (communication, dynamics)

Readout of (qu)bits (measurement)

Distinguishability of bit states

Quantum information

Stored as quantum state $|\psi\rangle$
of string of qubits

Unitary operations U (reversible)

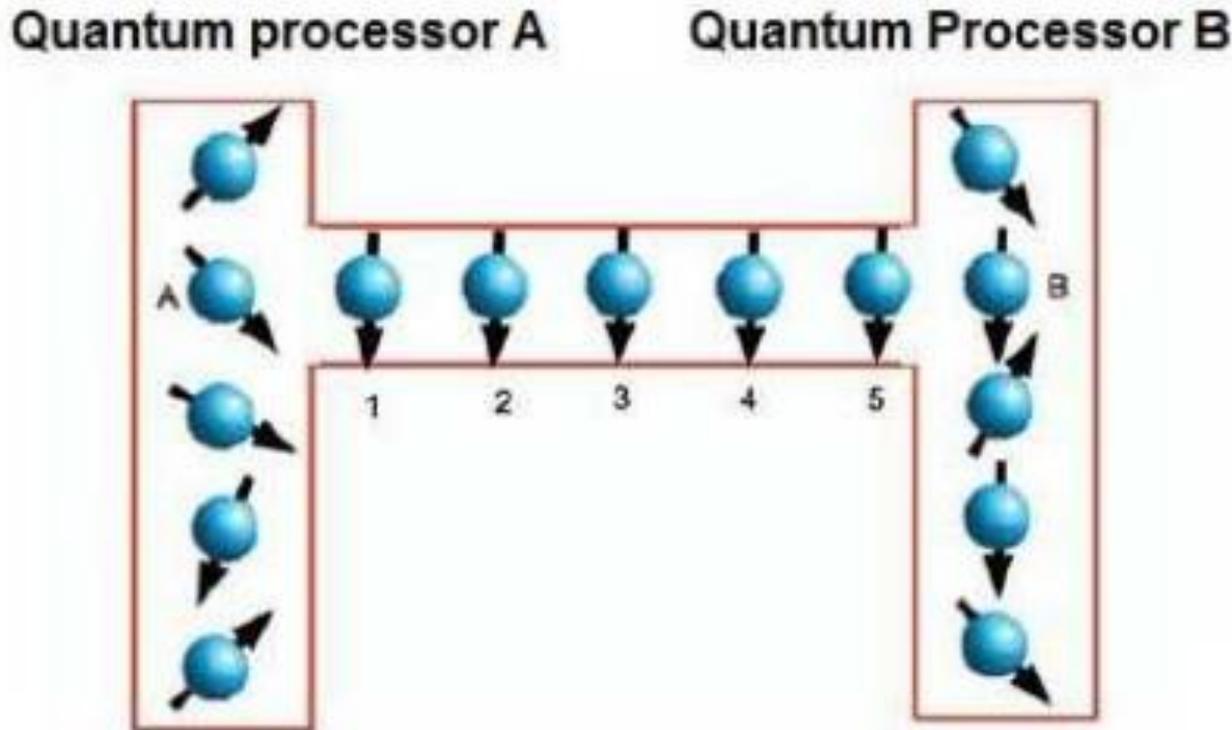
Qubit states cannot be copied, except for
orthogonal states $|\psi\rangle|0\rangle \xrightarrow{U} |\psi\rangle|\psi\rangle$
 $\langle\psi|\phi\rangle^2 = \langle\psi|\phi\rangle$

$$\begin{aligned} |\psi\rangle &= \alpha|0\rangle + \beta|1\rangle \\ p_0 &= |\alpha|^2 = |\langle 0|\psi\rangle|^2 \\ p_1 &= |\beta|^2 = |\langle 1|\psi\rangle|^2 \end{aligned}$$

Quantum states are not distinguishable, except for
orthogonal states

Quantum processors connected by spin chains

Realization is based on spin chain of the permanent fermions (spin-1/2 particles) with nearest-neighbour interaction: $|0\rangle$ - spin down, $|1\rangle$ - spin up.



Spin chains – recent development

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PHYSICAL REVIEW LETTERS

week ending
14 NOVEMBER 2003

Quantum Communication through an Unmodulated Spin Chain

Sougato Bose

*Institute for Quantum Information, MC 107-81, California Institute of Technology, Pasadena, California 91125-8100, USA
and Department of Physics and Astronomy, University College London, Gower Street, London WC1E 6BT, United Kingdom*

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We propose a scheme for using an unmodulated and unmeasured spin chain as a channel for short distance quantum communications. The state to be transmitted is placed on one spin of the chain and received later on a distant spin with some fidelity. We first obtain simple expressions for the fidelity of quantum state transfer and the amount of entanglement sharable between any two sites of an arbitrary Heisenberg ferromagnet using our scheme. We then apply this to the realizable case of an open ended chain with nearest neighbor interactions. The fidelity of quantum state transfer is obtained as an inverse discrete cosine transform and as a Bessel function series. We find that in a reasonable time, a qubit can be directly transmitted with better than classical fidelity across the full length of chains of up to 80 spins. Moreover, our channel allows distillable entanglement to be shared over arbitrary distances.

DOI: 10.1103/PhysRevLett.91.207901

PACS numbers: 03.67.Hk, 05.50.+q, 32.80.Lg

721 times cited

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Spin chains – recent development

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7 MAY 2004

Perfect State Transfer in Quantum Spin Networks

Matthias Christandl,^{1,*} Nilanjana Datta,² Artur Ekert,^{1,3} and Andrew J. Landahl^{4,5}

¹Centre for Quantum Computation, Centre for Mathematical Sciences, DAMTP, University of Cambridge, Wilberforce Road, Cambridge CB3 0WA, United Kingdom

²Statistical Laboratory, Centre for Mathematical Science, University of Cambridge, Wilberforce Road, Cambridge CB3 0WB, United Kingdom

³Department of Physics, National University of Singapore, Singapore 117 542, Singapore

⁴Center for Bits and Atoms, Massachusetts Institute of Technology, Cambridge, Massachusetts 02139, USA

⁵HP Labs, Palo Alto, California 94304-1126, USA

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We propose a class of qubit networks that admit the perfect state transfer of any quantum state in a fixed period of time. Unlike many other schemes for quantum computation and communication, these

589 times cited

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$$J_n = \begin{array}{cccccc} \sqrt{1 \cdot 5} & \sqrt{2 \cdot 4} & \sqrt{3 \cdot 3} & \sqrt{4 \cdot 2} & \sqrt{5 \cdot 1} & \\ \bullet & \bullet & \bullet & \bullet & \bullet & \bullet \\ n = & 1 & 2 & 3 & 4 & 5 & 6 \\ m = & -\frac{5}{2} & -\frac{3}{2} & -\frac{1}{2} & \frac{1}{2} & \frac{3}{2} & \frac{5}{2} \\ & A & & & & & B \end{array}$$

Spin chains – recent development

PRL 93, 230502 (2004)

PHYSICAL REVIEW LETTERS

week ending
3 DECEMBER 2004

Mirror Inversion of Quantum States in Linear Registers

Claudio Albanese,^{1,2,*} Matthias Christandl,^{3,†} Nilanjana Datta,^{4,‡} and Artur Ekert^{3,5,§}

¹*Department of Mathematics, Imperial College, London, SW7 2AZ, United Kingdom*

²*Department of Mathematics, National University of Singapore, Singapore 117543, Singapore*

³*Centre for Quantum Computation, DAMTP, University of Cambridge, Cambridge CB3 0WA, United Kingdom*

⁴*Statistical Laboratory, DPMMS, University of Cambridge, Cambridge CB3 0WB, United Kingdom*

⁵*Department of Physics, National University of Singapore, Singapore 117542, Singapore*

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Transfer of data in linear quantum registers can be significantly simplified with preengineered but not dynamically controlled interqubit couplings. We show how to implement a mirror inversion of the state of the register in each excitation subspace with respect to the center of the register. Our construction is especially appealing as it requires no dynamical control over individual interqubit interactions. If, however, individual control of the interactions is available then the mirror inversion operation can be performed on any substring of qubits in the register. In this case, a sequence of mirror inversions can generate any permutation of a quantum state of the involved qubits.

DOI: 10.1103/PhysRevLett.93.230502

PACS numbers: 03.67.Hk, 05.50.+q

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Spin chains with nearest-neighbor interaction

Hamiltonian of spin chain of $(N + 1)$ electrons coupled via the nearest-neighbour interaction is the following

Hamiltonian:

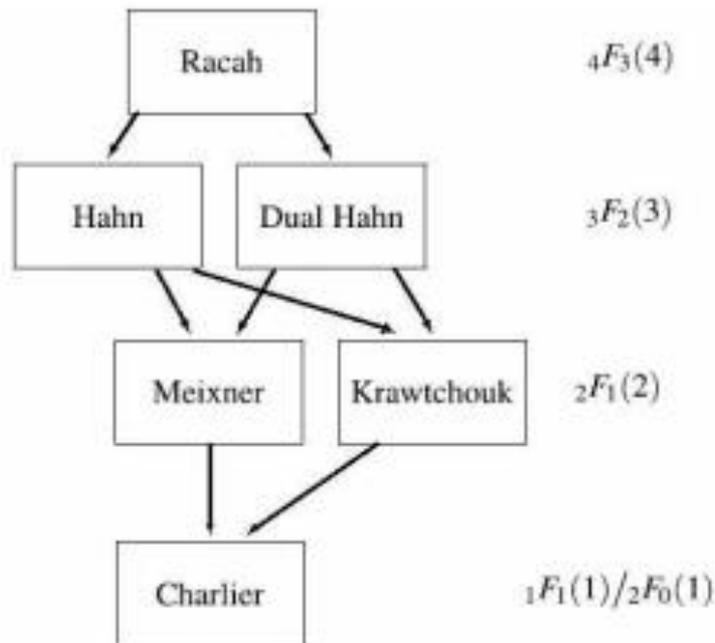
$$\hat{H} = \frac{1}{2} \sum_{k=0}^{N-1} J_k (\sigma_k^x \cdot \sigma_{k+1}^x + \sigma_{k+1}^y \cdot \sigma_k^y) + \frac{1}{2} \sum_{k=0}^N h_k (\sigma_k^z + 1)$$

J_k - coupling strength, h_k - Zeeman energy.

Pauli matrices:

$$\sigma^x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad \sigma^y = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \quad \sigma^z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$$

Askey scheme of the orthogonal polynomials



- R. Koekoek, P.A. Lesky, R. Swarttouw, Hypergeometric Orthogonal Polynomials and Their q-Analogues, Springer (2010)
-

Spin chains – simplest example of analytical solution



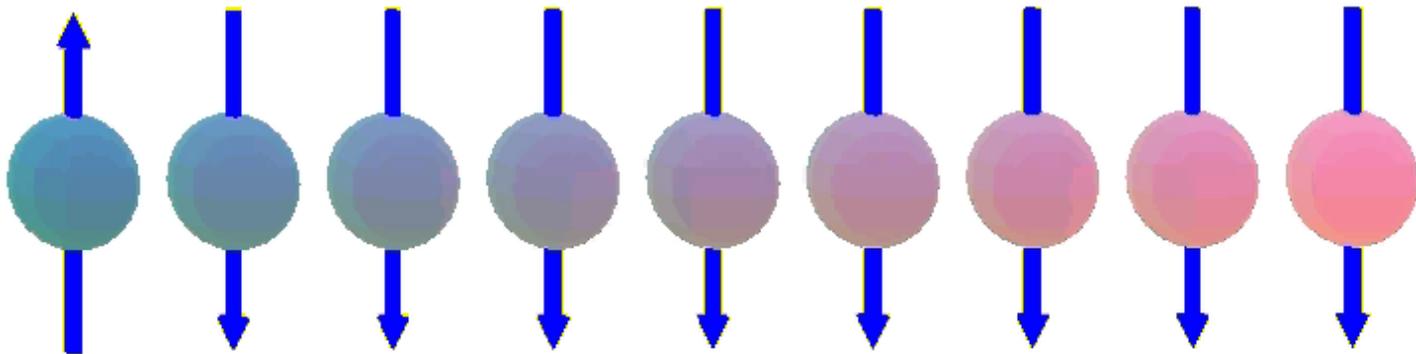
$$h_k = 0 \text{ and } J_k = 1.$$

$$\epsilon_j = 2 \cos \left(\frac{(j+1)\pi}{N+2} \right) \quad U_{ij} = \sqrt{\frac{2}{N+2}} \sin \left(\frac{(i+1)(j+1)\pi}{N+2} \right).$$

$$f_{r,s}(t) = \sum_{j=0}^N \frac{2}{N+2} \sin \left(\frac{(r+1)(j+1)\pi}{N+2} \right) \sin \left(\frac{(s+1)(j+1)\pi}{N+2} \right) \\ \times \exp \left(-2it \cos \left(\frac{(j+1)\pi}{N+2} \right) \right)$$

Simulation

TIME = 0.



Spin chains - another example - Krawtchouk polynomials

$$\triangleright \quad h_k = N/2 \text{ and } J_k = \frac{1}{2} \sqrt{(k+1)(N-k)}.$$

$$\epsilon_j = j \quad (j = 0, 1, \dots, N) \quad U_{ij} = \tilde{K}_i(j) \sim {}_2F_1 \left(\begin{matrix} -i, -j \\ -N \end{matrix}; \frac{1}{2} \right)$$

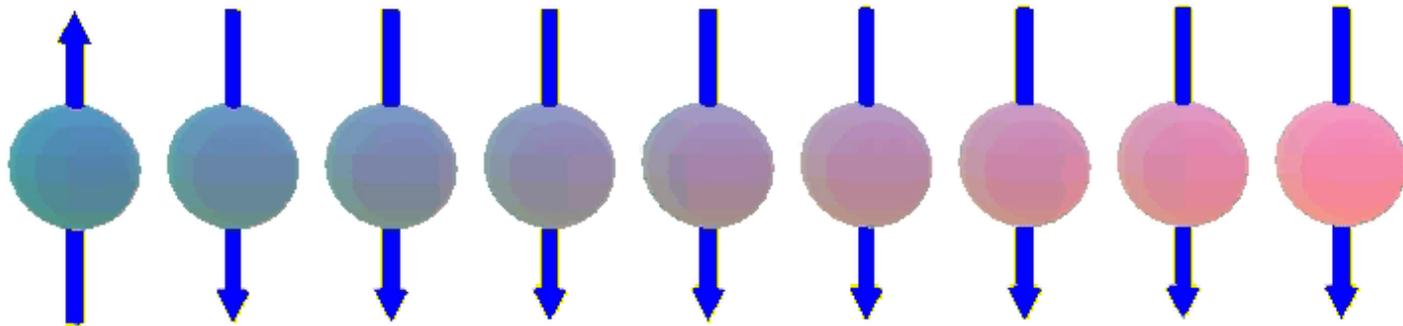
$$f_{r,s}(t) = \sqrt{\binom{N}{r} \binom{N}{s}} (\sqrt{p(1-p)})^{r+s} (1-z)^{r+s} (1-p+pz)^{N-r-s} \\ \times {}_2F_1 \left(\begin{matrix} -r, -s \\ -N \end{matrix}; \frac{-z}{p(1-p)(1-z)^2} \right).$$

$$f_{N,0}(t) = \left(\sqrt{p(1-p)} \right)^N (1 - e^{-it})^N$$

$$f_{N,0}(\pi) = 1 \quad p = 1/2.$$

Simulation

TIME = 0.



Spin chains - another example - q-Krawtchouk polynomials

Jacobi matrix elements

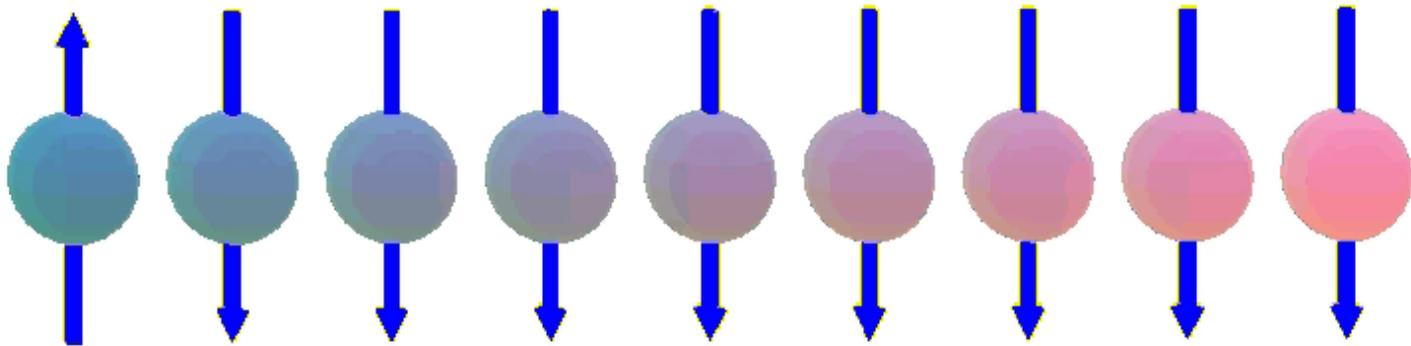
$$J_n = -\frac{A_n}{1-q} \sqrt{\frac{d_{n+1}}{d_n}}, \quad h_n = -\frac{A_n + C_n}{1-q},$$

$$U_{jk} = \tilde{K}_j(q^{-k}) \quad UU^T = U^T U = I.$$

$$\epsilon_j = -[-j] = -\frac{1 - q^{-j}}{1 - q} = q^{-1} + q^{-2} + \dots + q^{-j}.$$

Simulation

TIME = 0.



Hamiltonian of spin chain (N+1) electrons

$$\hat{H} = \sum_{k=0}^{N-1} J_k (a_k^+ a_{k+1} + a_{k+1}^+ a_k)$$

Pair of new recurrence relations for Racah polynomials

$$\begin{aligned} & \left[\frac{(x + \alpha + 2)(x + \gamma + \delta + 1)}{2x + \gamma + \delta + 2} e^{\partial_x} - \frac{(x + 1)(x - \alpha + \gamma + \delta)}{2x + \gamma + \delta + 2} \right] R_n(\lambda(x); \alpha + 1, \beta - 1, \gamma, \delta) \\ & \quad = (\alpha + 1) R_n(\lambda(x + 1); \alpha, \beta, \gamma, \delta - 1), \\ & \left[\frac{(\beta - \gamma - x - 1)(x + \delta)}{2x + \gamma + \delta + 1} + \frac{(x + \gamma + 1)(x + \beta + \delta)}{2x + \gamma + \delta + 1} e^{\partial_x} \right] R_n(\lambda(x); \alpha, \beta, \gamma, \delta - 1) \\ & \quad = \frac{(n + \alpha + 1)(n + \beta)}{\alpha + 1} R_n(\lambda(x + 1); \alpha + 1, \beta - 1, \gamma, \delta). \end{aligned}$$

Racah polynomials

$$R_n(\lambda(x); \alpha, \beta, \gamma, \delta)$$
$$= {}_4F_3 \left(\begin{matrix} -n, n + \alpha + \beta + 1, -x, x + \gamma + \delta + 1 \\ \alpha + 1, \beta + \delta + 1, \gamma + 1 \end{matrix} ; 1 \right),$$

$$\lambda(x) = x(x + \gamma + \delta + 1), \quad n = 0, 1, 2, \dots, m$$

Coupling strength

$$J_k = \begin{cases} \sqrt{(k+1)(m-k)f(\alpha, \beta, \delta)}; & k - \text{odd} \\ \sqrt{(k+2\alpha+2)(m-k+2\beta)g(\delta)}; & k - \text{even} \end{cases}$$

$$f(\alpha, \beta, \delta) = \frac{(k-2\alpha+2\delta-m)(k+2\beta+2\delta-1)}{(2k+2\delta-m-1)(2k+2\delta-m+1)}$$

$$g(\delta) = \frac{(k-m+2\delta-1)(k+2\delta)}{(2k+2\delta-m-1)(2k+2\delta-m+1)}$$

During computations, it is necessary to take into account satisfaction of the following three cases, under which the Racah polynomials hold finite-discrete orthogonality relation

$$\frac{(-\beta)_m(\gamma+\delta+2)_m}{(-\beta+\gamma+1)_m(\delta+1)_m},$$

$$\frac{(\alpha+\delta)_m(\gamma+\delta+2)_m}{(-\alpha+\gamma+\delta+1)_m(\delta+1)_m},$$

$$\frac{(\alpha+\beta+2)_m(-\delta)_m}{(\alpha-\delta+1)_m(\beta+1)_m},$$

$$\text{if } \alpha + 1 = -m,$$

$$\text{if } \beta + \delta + 1 = -m,$$

$$\text{if } \gamma + 1 = -m.$$

Transition from one end of the chain ($s = 0$) to the final end of the chain ($r = N = 2m + 1$), $t \equiv T = \pi/2$;

Time evolution correlation function:

$$f_{N,0}(\pi/2) = \sqrt{\frac{(\alpha - \delta + 2)_m (\alpha + \delta + 1)_m}{(\delta)_m (1 - \delta)_m}}$$

Conclusion

- Time-dependent correlation function becomes equal to 1, is called as perfect state transfer. One can easily check that expression of the correlation function, obtained in this work, never becomes equal to 1.
- Selection of nearest-neighbour coupling parameters J_k , corresponding to pair of the recurrence relations for the Racah polynomials, one can assess only the case of the qubit transfer with the high fidelity. This statement is very important, because, it allows observing the law of the breakdown of the perfect qubit transfer under the certain complication of the nearest-neighbour interaction parameter.

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Thank you for attention!