

Parallel implementation of $CNOT^N$ and C_2NOT^2 gates via homonuclear and heteronuclear Förster interactions of Rydberg atoms

Ahmed M. Farouk^{1,2}, I.I. Beterov^{1,3}, Peng Xu⁴, S. Bergamini⁵, I.I. Ryabstev^{1,3}

¹ Novosibirsk State University, Russia.

² Al-Azhar University, Egypt.

³ Institute of Semiconductors, SB-RAS, Russia.

⁴ Wuhan Institute of Quantum Technology, China.

⁵ The Open University, United Kingdom.

ahmed.farouk@azhar.edu.eg

Физика Ультрахолодных Атомов – 2022
"Physics of ultracold atoms - 2022"

December 21, 2022.



Introduction

- Alkali Rydberg atoms
- C-NOT gate and Bell states

Parallel Implementation of C-NOT^N gates

- Physical system
- Calculation of interaction energy
- homonuclear and heteronuclear architectures
- Generation of GHZ-state

Implementation of C₂NOT² gates

Gate Error

Experimental demonstration of EIT

Conclusion and Outlook



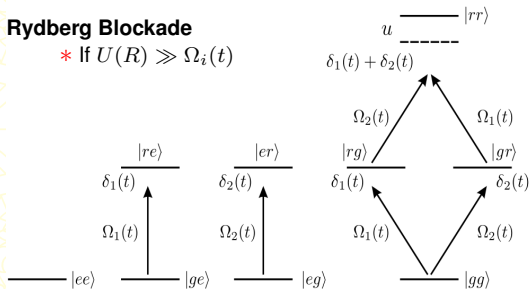
Alkali Rydberg atoms

Rydberg atom is a neutral atom with one or more electrons in a highly excited state with principal quantum number $n \gg 1$.

Alkali metals, particularly rubidium and cesium, are the species of atoms most commonly used for Rydberg atom experiments due to their single valence electrons.

Rydberg Blockade

* If $U(R) \gg \Omega_i(t)$



PRL 85, 2208 (2000).

Le Roy radius

$$R_{LR} \propto n^3$$

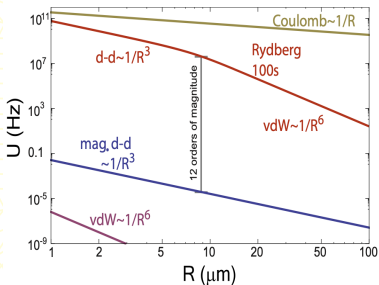
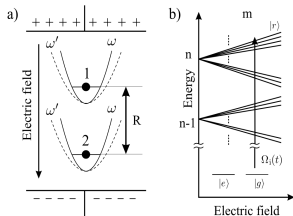
$$U_{d-d}(R) = \frac{C_3}{R^3}$$

$$C_3 \propto n^4$$

$$U_{vdW}(R) = \frac{C_6}{R^6}$$

$$C_6 \propto n^{11}$$

ARC, Comp. Phys. Com. 220, 319 (2017).



Rev. Mod. Phys. 82, 3 (2313) (2010).

C-NOT gate and Bell states

$$\text{CNOT}|x\rangle|y\rangle = |x\rangle|x \otimes y\rangle$$

$$\text{CNOT} = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 0 & 1 & 0 \end{pmatrix}$$

CNOT gate	
input	output
$ 00\rangle$	$ 00\rangle$
$ 01\rangle$	$ 01\rangle$
$ 10\rangle$	$ 11\rangle$
$ 11\rangle$	$ 10\rangle$



Nielsen and Chuang (2000)
[Quantum computation and quantum information]

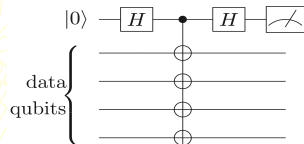
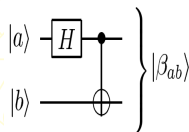
Since superposition of input states leads to a superposition of the corresponding output states

$$|00\rangle \xrightarrow{\text{Hadamard gate}} \frac{1}{\sqrt{2}} (|00\rangle + |10\rangle) \xrightarrow{\text{C-NOT gate}} \frac{1}{\sqrt{2}} (|00\rangle + |11\rangle)$$

Bell states

$$|\Phi^\pm\rangle = \frac{1}{\sqrt{2}} (|00\rangle \pm |11\rangle)$$

$$|\Psi^\pm\rangle = \frac{1}{\sqrt{2}} (|01\rangle \pm |10\rangle)$$



PRA 96, 052320 (2017)

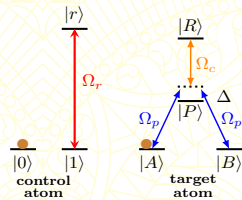
Hadamard gate

$$H|0\rangle = |+\rangle = \frac{1}{\sqrt{2}} (|0\rangle + |1\rangle)$$

$$H|1\rangle = |-\rangle = \frac{1}{\sqrt{2}} (|0\rangle - |1\rangle)$$

Parallel Implementation of C-NOT^N gates

(I) Blocking gate: $|0\rangle|A^N\rangle \rightarrow |0\rangle|A^N\rangle$
 $|0\rangle|B^N\rangle \rightarrow |0\rangle|B^N\rangle.$

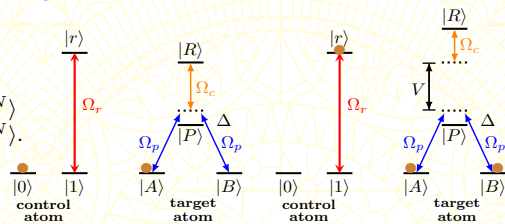


Blocking scheme

Parallel Implementation of C-NOT^N gates

(I) Blocking gate: $|0\rangle|A^N\rangle \rightarrow |0\rangle|A^N\rangle$
 $|0\rangle|B^N\rangle \rightarrow |0\rangle|B^N\rangle.$

(II) Transfer gate: $|1\rangle|A^N\rangle \rightarrow -(-1)^N|1\rangle|B^N\rangle$
 $|1\rangle|B^N\rangle \rightarrow -(-1)^N|1\rangle|A^N\rangle.$



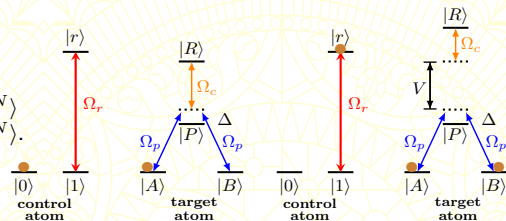
Blocking scheme

Transfer scheme

Parallel Implementation of C-NOT^N gates

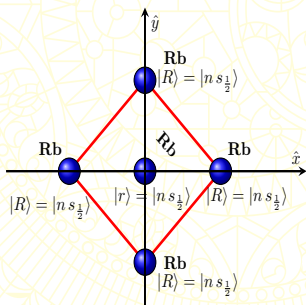
(I) Blocking gate: $|0\rangle|A^N\rangle \rightarrow |0\rangle|A^N\rangle$
 $|0\rangle|B^N\rangle \rightarrow |0\rangle|B^N\rangle.$

(II) Transfer gate: $|1\rangle|A^N\rangle \rightarrow -(-1)^N|1\rangle|B^N\rangle$
 $|1\rangle|B^N\rangle \rightarrow -(-1)^N|1\rangle|A^N\rangle.$



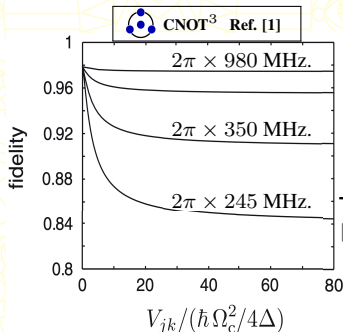
Blocking scheme

Transfer scheme



Homonuclear architecture

— Refs: [1-4]

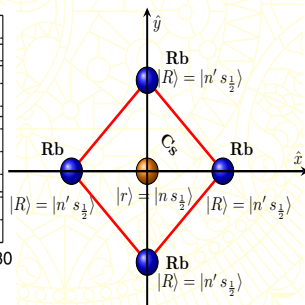


[1] PRL 102, 170502 (2009),

[3] New J. Phys. 16 053045 (2014),

[2] Nature Physics 6, 382 (2010),

[4] PRA 96, 052320 (2017).



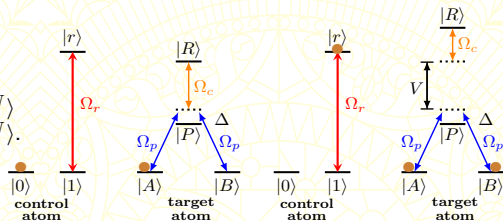
Heteronuclear architecture

— [This study]

Parallel Implementation of C-NOT^N gates

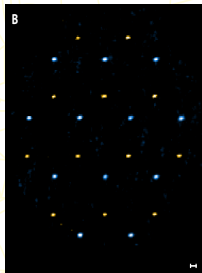
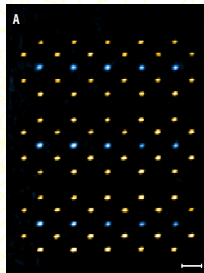
(I) Blocking gate: $|0\rangle|A^N\rangle \rightarrow |0\rangle|A^N\rangle$
 $|0\rangle|B^N\rangle \rightarrow |0\rangle|B^N\rangle.$

(II) Transfer gate: $|1\rangle|A^N\rangle \rightarrow -(-1)^N|1\rangle|B^N\rangle$
 $|1\rangle|B^N\rangle \rightarrow -(-1)^N|1\rangle|A^N\rangle.$



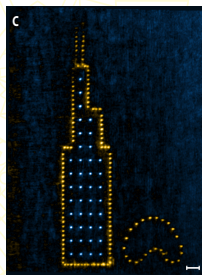
Blocking scheme

Transfer scheme



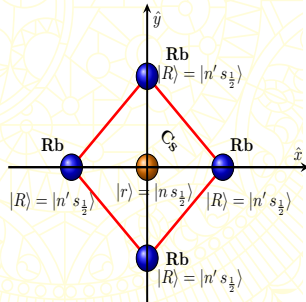
Heteronuclear architecture

- PRX 12, 011040 (2022)



Heteronuclear architecture

- [This study]



Parallel Implementation of C-NOT^N gates

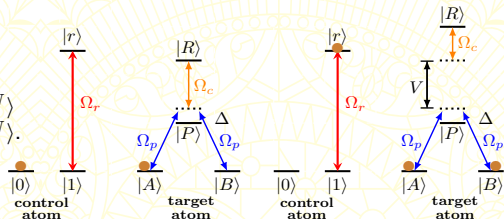
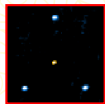
(I) Blocking gate: $|0\rangle|A^N\rangle \rightarrow |0\rangle|A^N\rangle$
 $|0\rangle|B^N\rangle \rightarrow |0\rangle|B^N\rangle.$

(II) Transfer gate: $|1\rangle|A^N\rangle \rightarrow -(-1)^N|1\rangle|B^N\rangle$
 $|1\rangle|B^N\rangle \rightarrow -(-1)^N|1\rangle|A^N\rangle.$

CNOT³

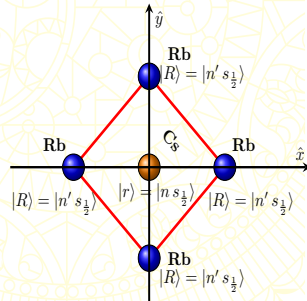
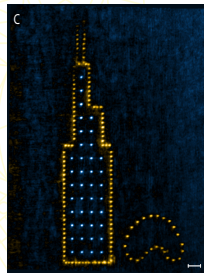
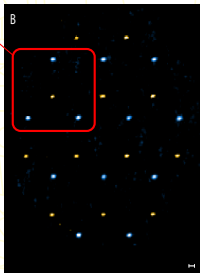
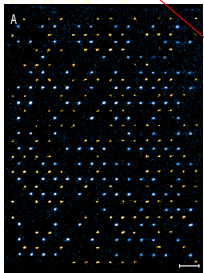
CNOT⁴

C₂NOT²



Blocking scheme

Transfer scheme



Heteronuclear architecture

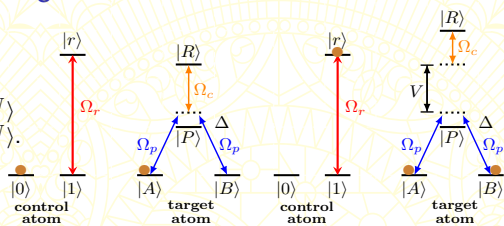
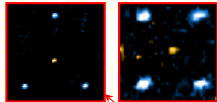
Heteronuclear architecture

Parallel Implementation of C-NOT^N gates

(I) Blocking gate: $|0\rangle|A^N\rangle \rightarrow |0\rangle|A^N\rangle$
 $|0\rangle|B^N\rangle \rightarrow |0\rangle|B^N\rangle.$

(II) Transfer gate: $|1\rangle|A^N\rangle \rightarrow -(-1)^N|1\rangle|B^N\rangle$
 $|1\rangle|B^N\rangle \rightarrow -(-1)^N|1\rangle|A^N\rangle.$

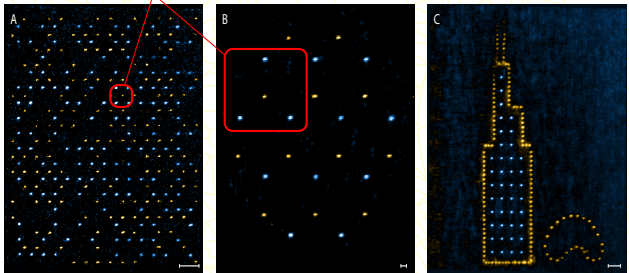
CNOT³ CNOT⁴ C₂NOT²



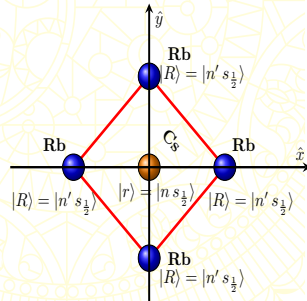
Blocking scheme

Transfer scheme

● ¹³³Cs ● ⁸⁷Rb



Heteronuclear architecture



Heteronuclear architecture

- PRX 12, 011040 (2022)

- [This study]

Parallel Implementation of C-NOT^N gates

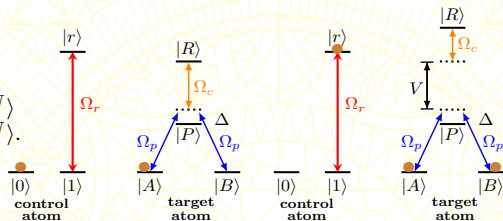
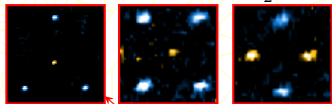
(I) Blocking gate: $|0\rangle|A^N\rangle \rightarrow |0\rangle|A^N\rangle$
 $|0\rangle|B^N\rangle \rightarrow |0\rangle|B^N\rangle.$

(II) Transfer gate: $|1\rangle|A^N\rangle \rightarrow -(-1)^N|1\rangle|B^N\rangle$
 $|1\rangle|B^N\rangle \rightarrow -(-1)^N|1\rangle|A^N\rangle.$

CNOT³

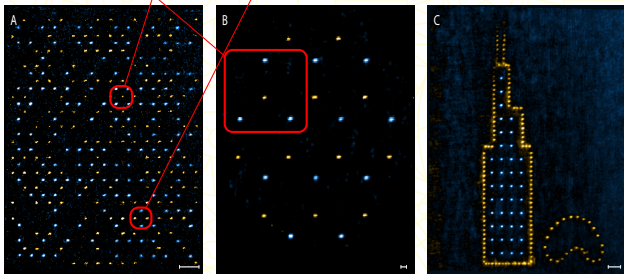
CNOT⁴

C₂NOT²



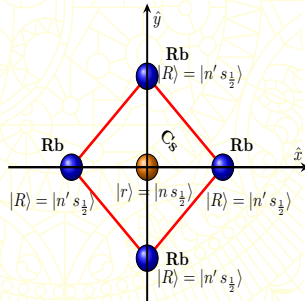
Blocking scheme

Transfer scheme



Heteronuclear architecture

- PRX 12, 011040 (2022)



Heteronuclear architecture

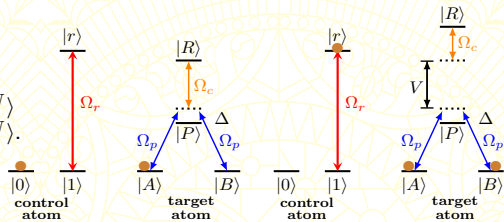
- [This study]

Parallel Implementation of C-NOT^N gates

(I) Blocking gate: $|0\rangle|A^N\rangle \rightarrow |0\rangle|A^N\rangle$
 $|0\rangle|B^N\rangle \rightarrow |0\rangle|B^N\rangle.$

(II) Transfer gate: $|1\rangle|A^N\rangle \rightarrow -(-1)^N|1\rangle|B^N\rangle$
 $|1\rangle|B^N\rangle \rightarrow -(-1)^N|1\rangle|A^N\rangle.$

PRL102, 170502 (2009)



Blocking scheme

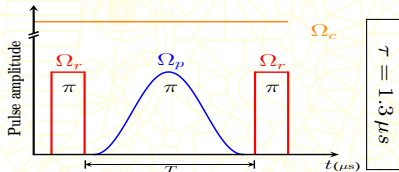
Transfer scheme

The Hamiltonian of control atom can be written as:

$$\hat{H}_C = \frac{\hbar}{2} \begin{pmatrix} 0 & 0 & 0 \\ 0 & 0 & \Omega_r \\ 0 & \Omega_r & 0 \end{pmatrix}.$$

The Hamiltonian of target atom can be written as:

$$\hat{H}_T = \frac{\hbar}{2} \begin{pmatrix} 0 & 0 & \Omega_p(t) & 0 \\ 0 & 0 & \Omega_p(t) & 0 \\ \Omega_p(t) & \Omega_p(t) & -2\Delta & \Omega_c \\ 0 & 0 & \Omega_c & 0 \end{pmatrix}.$$



$$\Omega_p(t) = \sqrt{\frac{16\pi\Delta}{3T}} \sin^2\left(\frac{\pi t}{T}\right)$$

$$\int_0^T \Omega_p^2(t) dt = 2\pi\Delta$$

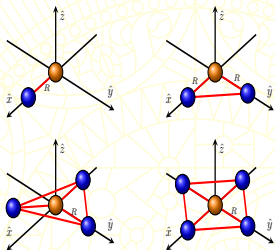
The Hamiltonian of the combined system for N target atom

$$\hat{H} = \hat{H}_C \otimes \hat{I} + \hat{I} \otimes \hat{H}_T + \sum_j^N V_{ct_j} |r\rangle\langle r| \otimes |R\rangle_j\langle R| + \sum_{j \neq k}^N V_{t_j t_k} |R\rangle_j\langle R| \otimes |R\rangle_k\langle R|.$$

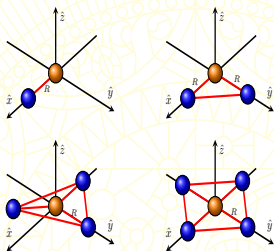
control-target interaction

target-target interaction

Spatial configurations of CNOT^N gates



Spatial configurations of CNOT^N gates



$$\hat{\mathcal{H}}_t = \hat{H}_{t_1} \otimes \hat{I}_{\mathcal{N}_t} \otimes \hat{I}_{\mathcal{N}_t} \otimes \hat{I}_{\mathcal{N}_t} + \hat{I}_{\mathcal{N}_t} \otimes \hat{H}_{t_2} \otimes \hat{I}_{\mathcal{N}_t} \otimes \hat{I}_{\mathcal{N}_t} + \hat{I}_{\mathcal{N}_t} \otimes \hat{I}_{\mathcal{N}_t} \otimes \hat{H}_{t_3} \otimes \hat{I}_{\mathcal{N}_t} + \hat{I}_{\mathcal{N}_t} \otimes \hat{I}_{\mathcal{N}_t} \otimes \hat{I}_{\mathcal{N}_t} \otimes \hat{H}_{t_4},$$

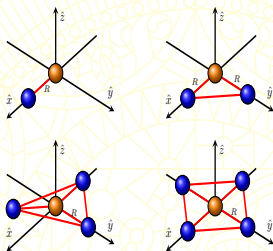
Control-target Hamiltonian:

$$|r\rangle\langle r| \otimes |R\rangle_1\langle R| = |r\rangle\langle r| \otimes \left(|R\rangle_1\langle R| \otimes \hat{I}_{\mathcal{N}_t} \otimes \hat{I}_{\mathcal{N}_t} \otimes \hat{I}_{\mathcal{N}_t} \right),$$

Target-target Hamiltonian:

$$|R\rangle_1\langle R| \otimes |R\rangle_2\langle R| = \hat{I}_{\mathcal{N}_c} \otimes \left(|R\rangle_1\langle R| \otimes |R\rangle_2\langle R| \otimes \hat{I}_{\mathcal{N}_t} \otimes \hat{I}_{\mathcal{N}_t} \right).$$

Spatial configurations of CNOT^N gates



$$\hat{\mathcal{H}}_t = \hat{H}_{t_1} \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} + \hat{I}_{N_t} \otimes \hat{H}_{t_2} \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} + \hat{I}_{N_t} \otimes \hat{I}_{N_t} \otimes \hat{H}_{t_3} \otimes \hat{I}_{N_t} + \hat{I}_{N_t} \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \otimes \hat{H}_{t_4},$$

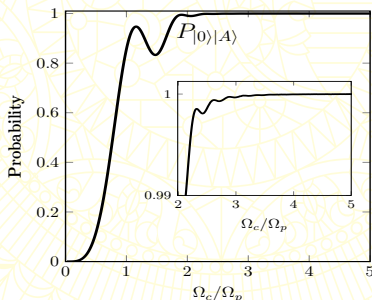
Control-target Hamiltonian:

$$|r\rangle\langle r| \otimes |R\rangle_1\langle R| = |r\rangle\langle r| \otimes \left(|R\rangle_1\langle R| \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \right),$$

Target-target Hamiltonian:

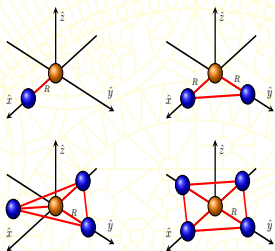
$$|R\rangle_1\langle R| \otimes |R\rangle_2\langle R| = \hat{I}_{N_c} \otimes \left(|R\rangle_1\langle R| \otimes |R\rangle_2\langle R| \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \right).$$

The behavior of blocking gate:-



• $\Delta = 2\pi \times 1.2$ GHz, • $\Omega_p = 2\pi \times 50$ MHz.

Spatial configurations of CNOT^N gates



$$\hat{\mathcal{H}}_t = \hat{H}_{t_1} \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} + \hat{I}_{N_t} \otimes \hat{H}_{t_2} \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \\ \otimes \hat{I}_{N_t} + \hat{I}_{N_t} \otimes \hat{I}_{N_t} \otimes \hat{H}_{t_3} \otimes \hat{I}_{N_t} + \hat{I}_{N_t} \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \otimes \hat{H}_{t_4},$$

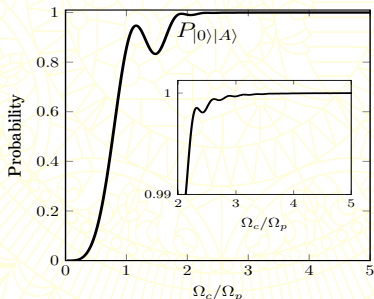
Control-target Hamiltonian:

$$|r\rangle\langle r| \otimes |R\rangle_1\langle R| = |r\rangle\langle r| \otimes \left(|R\rangle_1\langle R| \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \right),$$

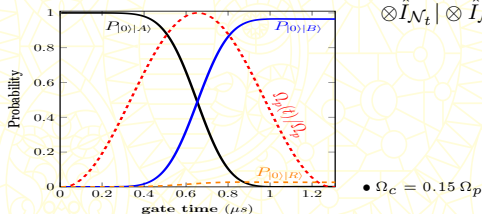
Target-target Hamiltonian:

$$|R\rangle_1\langle R| \otimes |R\rangle_2\langle R| = \hat{I}_{N_c} \otimes \left(|R\rangle_1\langle R| \otimes |R\rangle_2\langle R| \otimes \hat{I}_{N_t} \otimes \hat{I}_{N_t} \right).$$

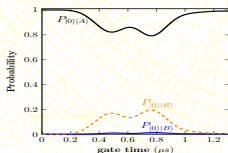
The behavior of blocking gate:-



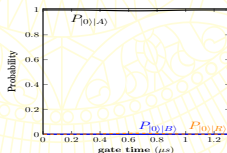
• $\Delta = 2\pi \times 1.2$ GHz, • $\Omega_p = 2\pi \times 50$ MHz.



• $\Omega_c = 0.15 \Omega_p$



• $\Omega_c = 2 \Omega_p$



• $\Omega_c = 14 \Omega_p$

Interaction energies and Fidelity

These values are calculated using ARC considering $\Delta n = 2$, $\Delta \ell = 2$, $\theta = \pi/2$,

$\phi = 0$, minimum state contribution = 0.2, and maximum energy difference

$\Delta E/h = 25$ GHz.

N ^o	Atom 1	Atom 2	R_{LR} (μm)	R_{vdW} (μm)	$C_3/2\pi$ (GHz $\cdot \mu\text{m}^3$)	$C_6/2\pi$ (GHz $\cdot \mu\text{m}^6$)
1	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	1.8	4.5	4.20	2036
2	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	2.0	9.5	1.92	2364
3	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	1.9	31.5	14.25	2484

[ARC, *Comp. Phys. Com.* **220**, 319 (2017).]

- R_{LR} is the interatomic distance where the theory of Le Roy-Bernstein is satisfied.
- R_{vdW} is crossover distance, marking boundary between dipole-dipole and vdW regimes.
- C_3 is the dispersive coefficient of d-d interaction energy.
- C_6 is the dispersive coefficient of vdW interaction energy.

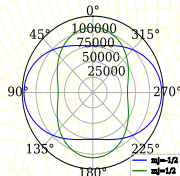
Interaction energies and Fidelity

These values are calculated using ARC considering $\Delta n = 2$, $\Delta \ell = 2$, $\theta = \pi/2$,

$\phi = 0$, minimum state contribution = 0.2, and maximum energy difference

$\Delta E/h = 25$ GHz.

N ^o	Atom 1	Atom 2	R_{LR} (μm)	R_{vdW} (μm)	$C_3/2\pi$ (GHz $\cdot \mu\text{m}^3$)	$C_6/2\pi$ (GHz $\cdot \mu\text{m}^6$)
1	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	1.8	4.5	4.20	2036
2	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	2.0	9.5	1.92	2364
3	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	1.9	31.5	14.25	2484



[ARC, Comp. Phys. Com. 220, 319 (2017).]

- R_{LR} is the interatomic distance where the theory of Le Roy-Bernstein is satisfied.
- R_{vdW} is crossover distance, marking boundary between dipole-dipole and vdW regimes.
- C_3 is the dispersive coefficient of d-d interaction energy.
- C_6 is the dispersive coefficient of vdW interaction energy.

Interaction energies and Fidelity

These values are calculated using ARC considering $\Delta n = 2$, $\Delta \ell = 2$, $\theta = \pi/2$,

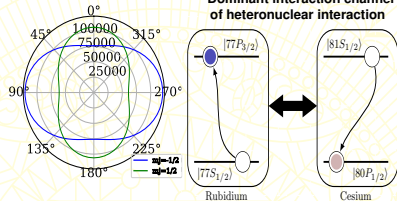
$\phi = 0$, minimum state contribution = 0.2, and maximum energy difference

$\Delta E/h = 25$ GHz.

N ^o	Atom 1	Atom 2	R_{LR} (μm)	R_{vdW} (μm)	$C_3/2\pi$ (GHz $\cdot \mu\text{m}^3$)	$C_6/2\pi$ (GHz $\cdot \mu\text{m}^6$)
1	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	1.8	4.5	4.20	2036
2	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	2.0	9.5	1.92	2364
3	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	1.9	31.5	14.25	2484

[ARC, Comp. Phys. Com. 220, 319 (2017).]

- R_{LR} is the interatomic distance where the theory of Le Roy-Bernstein is satisfied.
- R_{vdW} is crossover distance, marking boundary between dipole-dipole and vdW regimes.
- C_3 is the dispersive coefficient of d-d interaction energy.
- C_6 is the dispersive coefficient of vdW interaction energy.



Interaction energies and Fidelity

These values are calculated using ARC considering $\Delta n = 2$, $\Delta \ell = 2$, $\theta = \pi/2$.

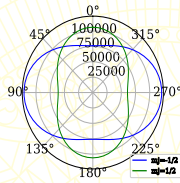
$\phi = 0$, minimum state contribution = 0.2, and maximum energy difference

$\Delta E/h = 25$ GHz.

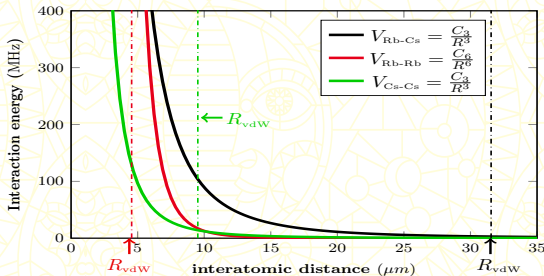
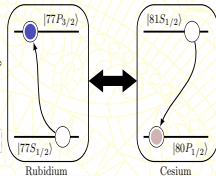
N ^o	Atom 1	Atom 2	R_{LR} (μm)	R_{vdW} (μm)	$C_3/2\pi$ (GHz $\cdot \mu\text{m}^3$)	$C_6/2\pi$ (GHz $\cdot \mu\text{m}^6$)
1	Rubidium 77S _{1/2} , m _j = 1/2⟩	Rubidium 77S _{1/2} , m _j = 1/2⟩	1.8	4.5	4.20	2036
2	Cesium 81S _{1/2} , m _j = -1/2⟩	Cesium 81S _{1/2} , m _j = -1/2⟩	2.0	9.5	1.92	2364
3	Rubidium 77S _{1/2} , m _j = 1/2⟩	Cesium 80P _{1/2} , m _j = -1/2⟩	1.9	31.5	14.25	2484

[ARC, Comp. Phys. Com. 220, 319 (2017).]

- R_{LR} is the interatomic distance where the theory of Le Roy-Bernstein is satisfied.
- R_{vdW} is crossover distance, marking boundary between dipole-dipole and vdW regimes.
- C_3 is the dispersive coefficient of d-d interaction energy.
- C_6 is the dispersive coefficient of vdW interaction energy.



Dominant interaction channel of heteronuclear interaction



Interaction energies and Fidelity

These values are calculated using ARC considering $\Delta n = 2$, $\Delta \ell = 2$, $\theta = \pi/2$.

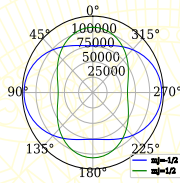
$\phi = 0$, minimum state contribution = 0.2, and maximum energy difference

$\Delta E/h = 25$ GHz.

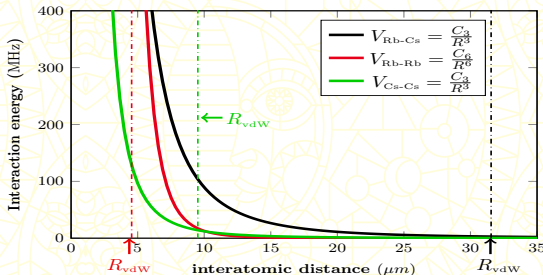
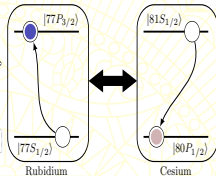
N ^o	Atom 1	Atom 2	R_{LR} (μm)	R_{vdW} (μm)	$C_3/2\pi$ (GHz $\cdot \mu\text{m}^3$)	$C_6/2\pi$ (GHz $\cdot \mu\text{m}^6$)
1	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	1.8	4.5	4.20	2036
2	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	2.0	9.5	1.92	2364
3	Rubidium $ 77S_{1/2}, m_j = 1/2\rangle$	Cesium $ 81S_{1/2}, m_j = -1/2\rangle$	1.9	31.5	14.25	2484

[ARC, Comp. Phys. Com. 220, 319 (2017).]

- R_{LR} is the interatomic distance where the theory of Le Roy-Bernstein is satisfied.
- R_{vdW} is crossover distance, marking boundary between dipole-dipole and vdW regimes.
- C_3 is the dispersive coefficient of d-d interaction energy.
- C_6 is the dispersive coefficient of vdW interaction energy.



Dominant interaction channel of heteronuclear interaction



Fidelity

$$F = \text{Tr} \left(\sqrt{\sqrt{\hat{\rho}} \hat{\sigma} \sqrt{\hat{\rho}}} \right)$$

$\hat{\rho}$ is the density matrix of the computational basis

$$\frac{1}{\sqrt{2}} (|0\rangle|A\rangle^{\otimes N} + |1\rangle|A\rangle^{\otimes N}) \rightarrow \frac{1}{\sqrt{2}} (|0\rangle|A\rangle^{\otimes N} + |1\rangle|B\rangle^{\otimes N})$$

$$\hat{\sigma} = |\Phi^+\rangle \langle \Phi^+|$$

$|\Phi^+\rangle$ is one of Bell states basis.

Interaction energies and Fidelity

These values are calculated using ARC considering $\Delta n = 2$, $\Delta \ell = 2$, $\theta = \pi/2$.

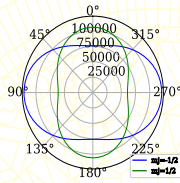
$\phi = 0$, minimum state contribution = 0.2, and maximum energy difference

$\Delta E/h = 25$ GHz.

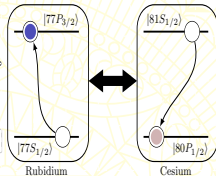
N ^o	Atom 1	Atom 2	R_{LR} (μm)	R_{vdW} (μm)	$C_3/2\pi$ (GHz $\cdot \mu\text{m}^3$)	$C_6/2\pi$ (GHz $\cdot \mu\text{m}^6$)
1	Rubidium [77S _{1/2} , m _j = 1/2]	Rubidium [77S _{1/2} , m _j = 1/2]	1.8	4.5	4.20	2036
2	Cesium [81S _{1/2} , m _j = -1/2]	Cesium [81S _{1/2} , m _j = -1/2]	2.0	9.5	1.92	2364
3	Rubidium [77S _{1/2} , m _j = 1/2]	Cesium [81S _{1/2} , m _j = -1/2]	1.9	31.5	14.25	2484

[ARC, Comp. Phys. Com. 220, 319 (2017).]

- R_{LR} is the interatomic distance where the theory of Le Roy-Bernstein is satisfied.
- R_{vdW} is crossover distance, marking boundary between dipole-dipole and vdW regimes.
- C_3 is the dispersive coefficient of d-d interaction energy.
- C_6 is the dispersive coefficient of vdW interaction energy.



Dominant interaction channel of heteronuclear interaction



Fidelity

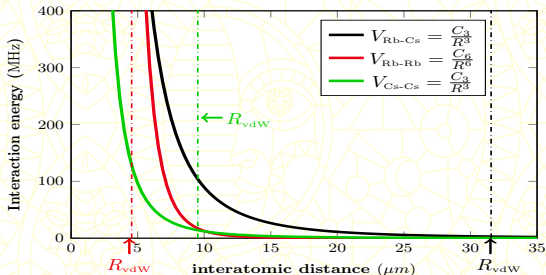
$$F = \text{Tr} \left(\sqrt{\sqrt{\hat{\rho}} \hat{\sigma} \sqrt{\hat{\rho}}} \right)$$

$\hat{\rho}$ is the density matrix of the computational basis

$$\frac{1}{\sqrt{2}} (|0\rangle|A\rangle \otimes^N + |1\rangle|A\rangle \otimes^N) \rightarrow \frac{1}{\sqrt{2}} (|0\rangle|A\rangle \otimes^N + |1\rangle|B\rangle \otimes^N)$$

$$\hat{\sigma} = |\Phi^+\rangle \langle \Phi^+|$$

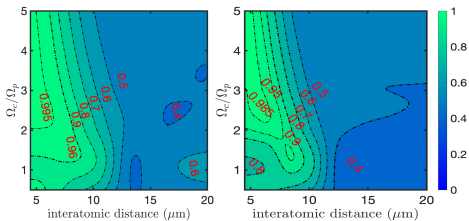
$|\Phi^+\rangle$ is one of Bell states basis.



- $\Delta = 2\pi \times 1.2$ GHz,
- $\Omega_p = 2\pi \times 50$ MHz,
- $\tau_{|r\rangle} = 505 - 548 \mu\text{s}$,
- $|P\rangle = |6 P_{3/2}\rangle$,
- $\tau_{|P\rangle} = 0.131 \mu\text{s}$.

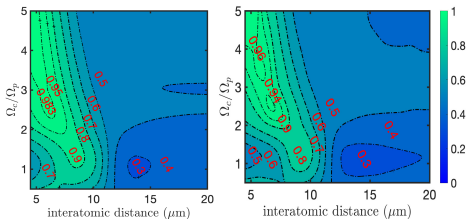
Fidelity contours

Homomuclear architectures of symmetric Rb atoms



$N = 1$

$N = 2$

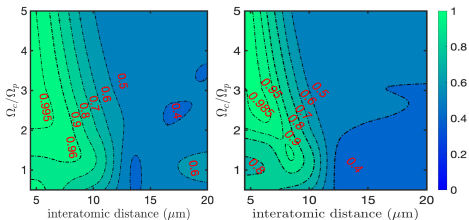


$N = 3$

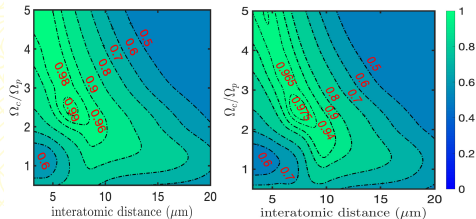
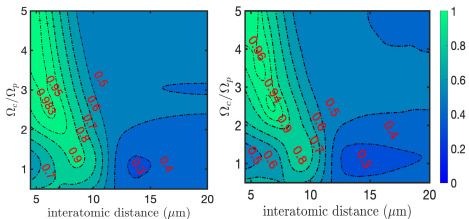
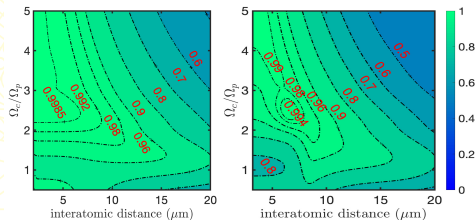
$N = 4$

Fidelity contours

Homonuclear architectures of symmetric Rb atoms



Heteronuclear architectures of Rb & Cs atoms



Fidelity contours

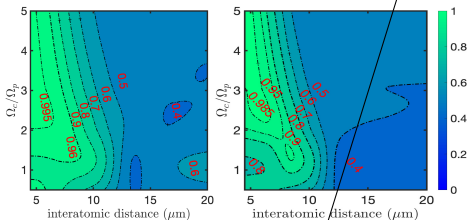
$$\mathcal{F}_{\text{Rb-Rb}}^{N=4} =$$



$$\mathcal{F}_{\text{Cs-Rb}}^{N=4} =$$

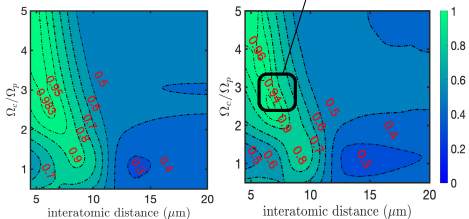


Homonuclear architectures of symmetric Rb atoms



$N = 1$

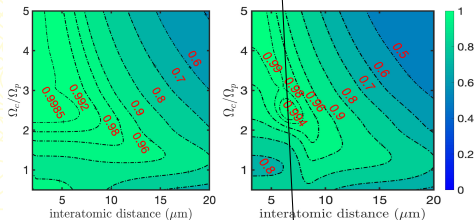
$N = 2$



$N = 3$

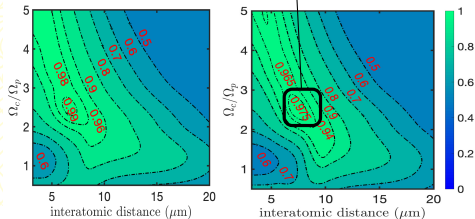
$N = 4$

Heteronuclear architectures of Rb & Cs atoms



$N = 1$

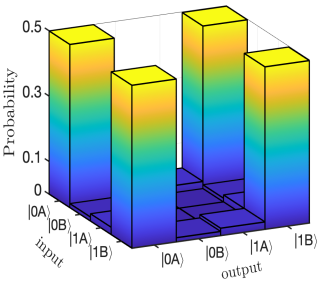
$N = 2$



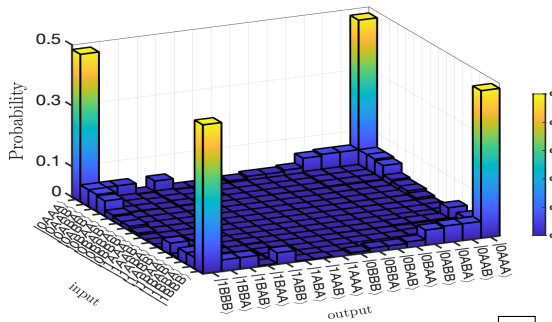
$N = 3$

$N = 4$

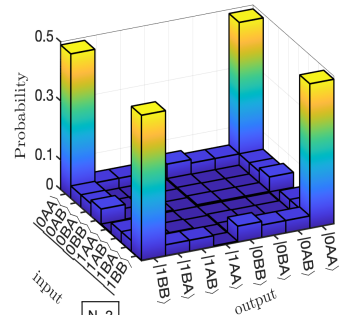
Generation of GHZ-state



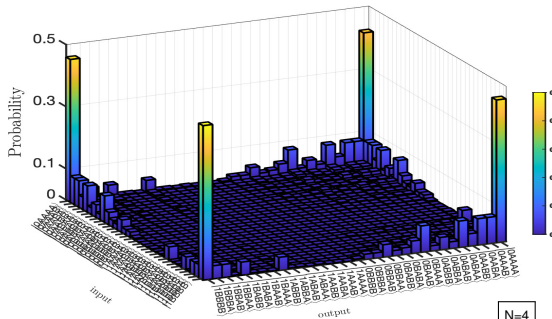
N=1



N=3

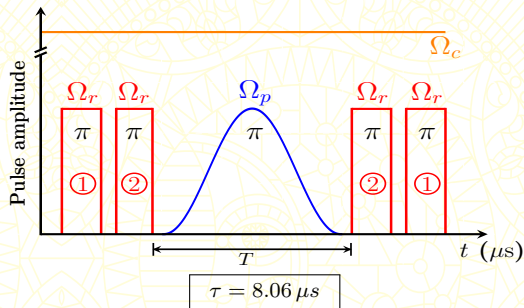
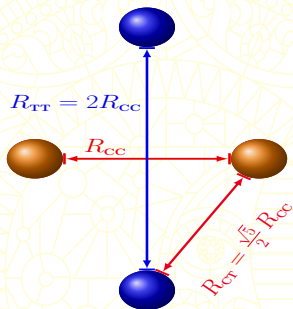


N=2



N=4

Demonstrating C_2NOT^2 gates

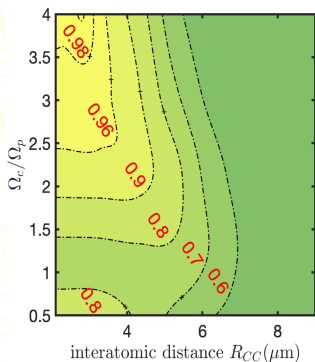


(I) Blocking: $|00\rangle|AA\rangle \rightarrow |00\rangle|AA\rangle$,
 $|00\rangle|BB\rangle \rightarrow |00\rangle|BB\rangle$,
 $|00\rangle|AB\rangle \rightarrow |00\rangle|AB\rangle$,
 $|00\rangle|BA\rangle \rightarrow |00\rangle|BA\rangle$,

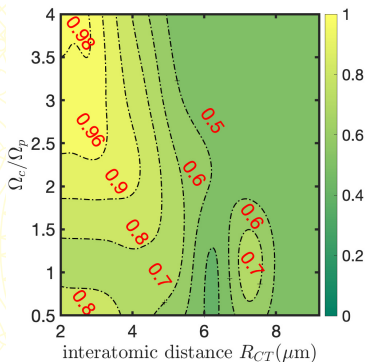
(II) Transfer: $|11\rangle|AA\rangle \rightarrow |11\rangle|BB\rangle$,
 $|11\rangle|BB\rangle \rightarrow |11\rangle|AA\rangle$,
 $|01\rangle|AB\rangle \rightarrow |01\rangle|BA\rangle$,
 $|01\rangle|BA\rangle \rightarrow |01\rangle|AB\rangle$,
 $|10\rangle|AB\rangle \rightarrow |10\rangle|BA\rangle$,
 $|10\rangle|BA\rangle \rightarrow |10\rangle|AB\rangle$.

Fidelity of C_2NOT^2

- Interaction of cesium atoms as control and rubidium atoms as targets.

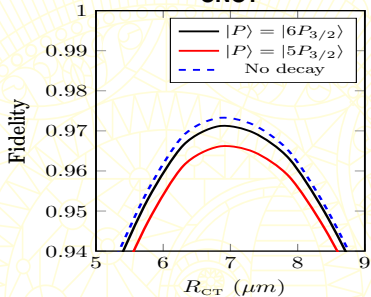


$$V_{CC} = 0.$$

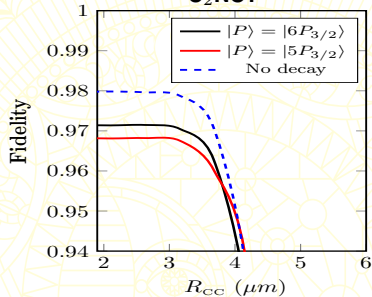


$$V_{CC} \neq 0.$$

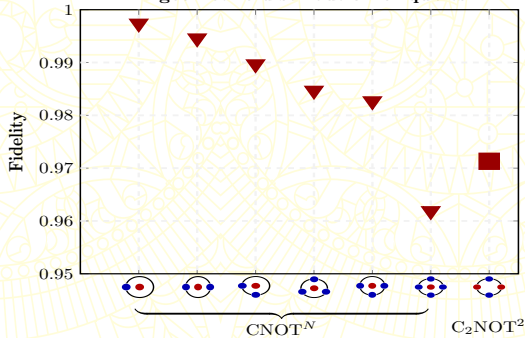
CNOT⁴



C₂NOT²

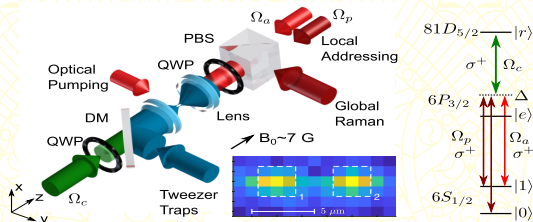


geometric distribution of qubits



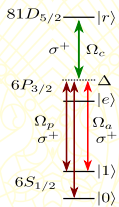
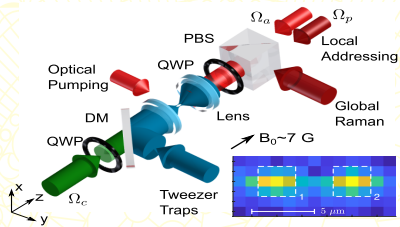
Foremost Experimental work demonstrating EIT

McDonnell et al, PRL 129, 200501 (2022)
ArXIV:2204.03733 (2022)

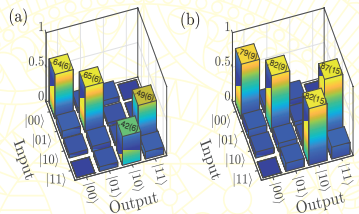


Experiment Setup

Foremost Experimental work demonstrating EIT McDonnell et al, PRL 129, 200501 (2022) ArXIV:2204.03733 (2022)



Experiment Setup

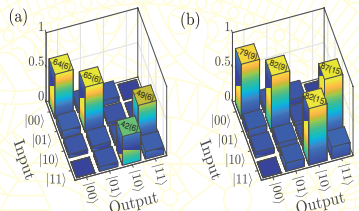
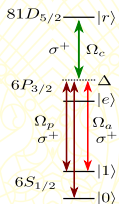
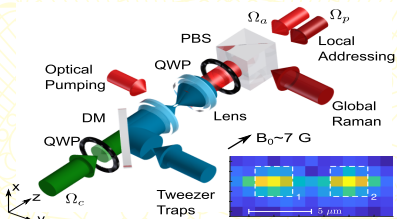


(a) $\mathcal{F}^{\text{RAW}} = 0.55$ (b) $\mathcal{F}^{\text{COR}} = 0.82$

Gate Measurement

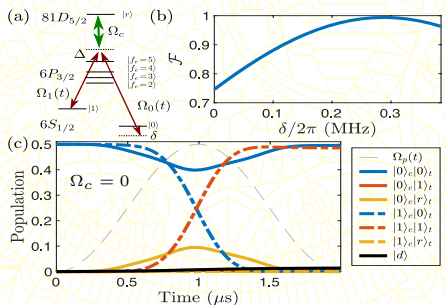
Foremost Experimental work demonstrating EIT

McDonnell et al, PRL 129, 200501 (2022)
ArXIV:2204.03733 (2022)



Experiment Setup

Gate Measurement

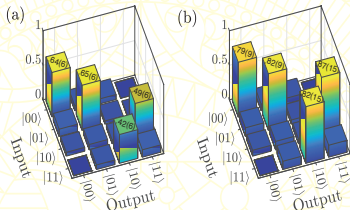
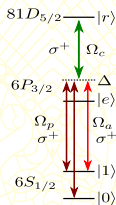
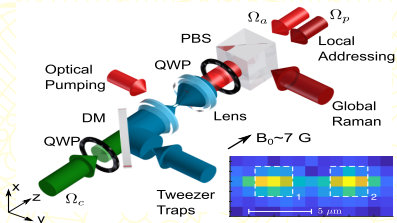


EIT Protocol via $6P_{3/2}$

$\Delta/2\pi = 1.05\text{GHz}$
 $F = 0.98$

Foremost Experimental work demonstrating EIT

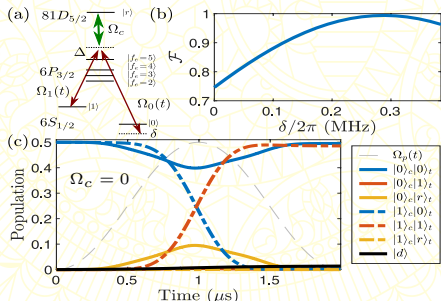
McDonnell et al, PRL 129, 200501 (2022)
ArXIV:2204.03733 (2022)



(a) $\mathcal{F}_{\text{RAW}} = 0.55$ (b) $\mathcal{F}_{\text{COR}} = 0.82$

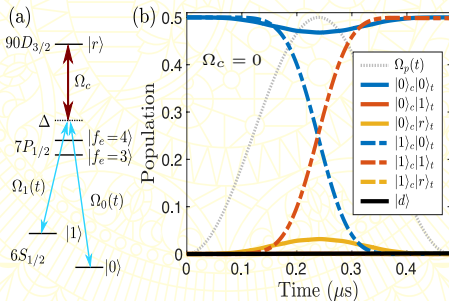
Experiment Setup

Gate Measurement



EIT Protocol via $6P_{3/2}$

$\Delta/2\pi = 1.05\text{GHz}$
 $\mathcal{F} = 0.98$

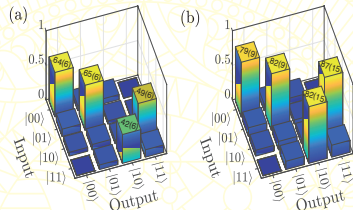
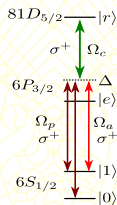
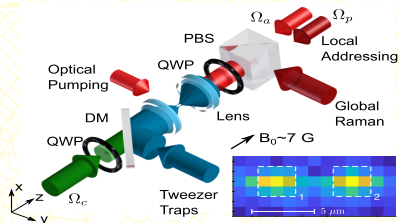


EIT Protocol via $7P_{1/2}$

$\Delta/2\pi = 5\text{GHz}$
 $\mathcal{F} > 0.998$

Foremost Experimental work demonstrating EIT

McDonnell et al. PRL 129, 200501 (2022)
ArXIV:2204.03733 (2022)

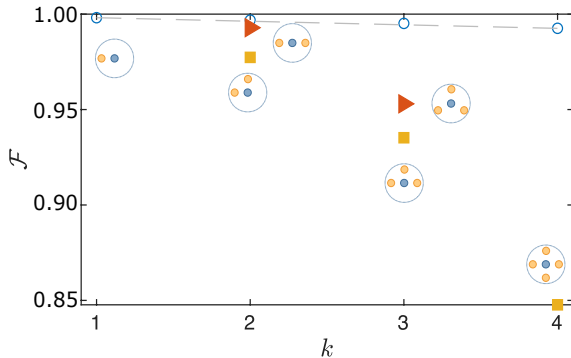


(a) $\mathcal{F}^{\text{RAW}} = 0.55$

(b) $\mathcal{F}^{\text{COR}} = 0.82$

Experiment Setup

Gate Measurement



Conclusion and Outlook

We can conclude the following:

- ▶ Heteronuclear architecture of alkali atoms is much better than the homonuclear architecture in terms of fidelity.
- ▶ The obtained Fidelity for CNOT⁴, with cesium control atom and rubidium target atoms, is $\simeq 97.5\%$ for fast gate implementation $\tau = 1.303 \mu\text{s}$ and moderate value of Rabi frequency $\Omega_c > 2\pi \times 125 \text{ MHz}$.
- ▶ The obtained fidelity is higher than the lower bound fidelity threshold for creating logic qubits of CNOT gates without the need to perform error correction surface codes.
 - IEEE Access, 7, 121501-121529, (2019).
- ▶ Using the 2nd resonance levels of target atoms (for Rb $5 S_{1/2} \rightarrow 6 P_{3/2} \rightarrow n S_{1/2}$, and for Cs $6 S_{1/2} \rightarrow 7 P_{3/2} \rightarrow n' S_{1/2}$) enhances the fidelity of the combined system.

We can conclude the following:

- ▶ Heteronuclear architecture of alkali atoms is much better than the homonuclear architecture in terms of fidelity.
- ▶ The obtained Fidelity for CNOT⁴, with cesium control atom and rubidium target atoms, is $\simeq 97.5\%$ for fast gate implementation $\tau = 1.303 \mu\text{s}$ and moderate value of Rabi frequency $\Omega_c > 2\pi \times 125 \text{ MHz}$.
- ▶ The obtained fidelity is higher than the lower bound fidelity threshold for creating logic qubits of CNOT gates without the need to perform error correction surface codes.
 - IEEE Access, 7, 121501-121529, (2019).
- ▶ Using the 2nd resonance levels of target atoms (for Rb $5 S_{1/2} \rightarrow 6 P_{3/2} \rightarrow n S_{1/2}$, and for Cs $6 S_{1/2} \rightarrow 7 P_{3/2} \rightarrow n' S_{1/2}$) enhances the fidelity of the combined system.

Future work

- ▶ Study the effect of coherent transport of a moving control atom implementing CNOT^N gate and the mutual information of the subsystems.

End of text!

Thanks!