

Developing theoretical and experimental tools for a hybrid quantum simulation using trapped ions

Sainath Motlakunta

Quantum Information with Trapped Ions (QITI) Group

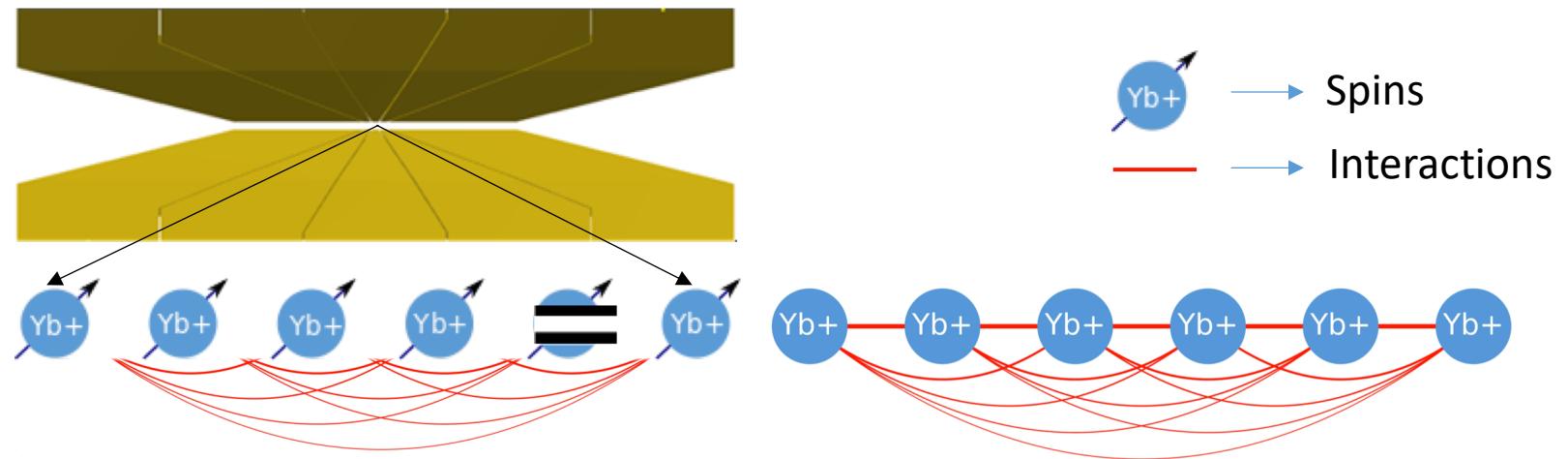


UNIVERSITY OF
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Quantum Computing

Talk outline

- **Part1: Theoretical tools:** A scheme to engineer programmable spin interaction graphs in a trapped ion quantum simulator.
- **Part2: Experimental tools:** Laser frequency stabilization for ion trap experiments

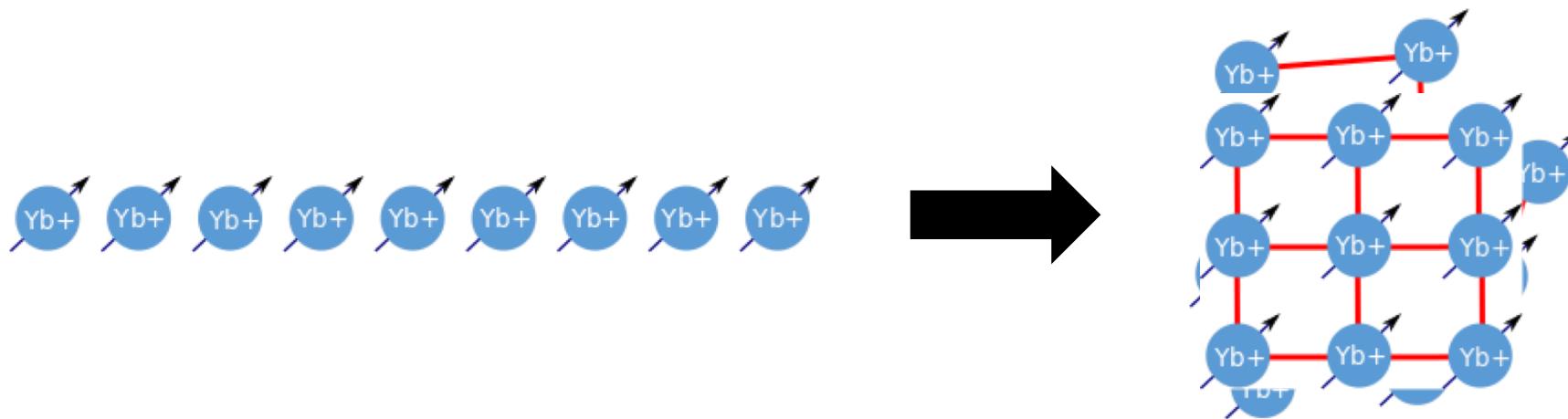
Trapped ions as Quantum Simulator



- Long coherence times (>10 min single qubit coherence time*)
- Ease of state preparation and isolation
- Controlled long range interactions
- Flexible fully connected graphs
- Useful to simulate systems with larger number of long range interactions

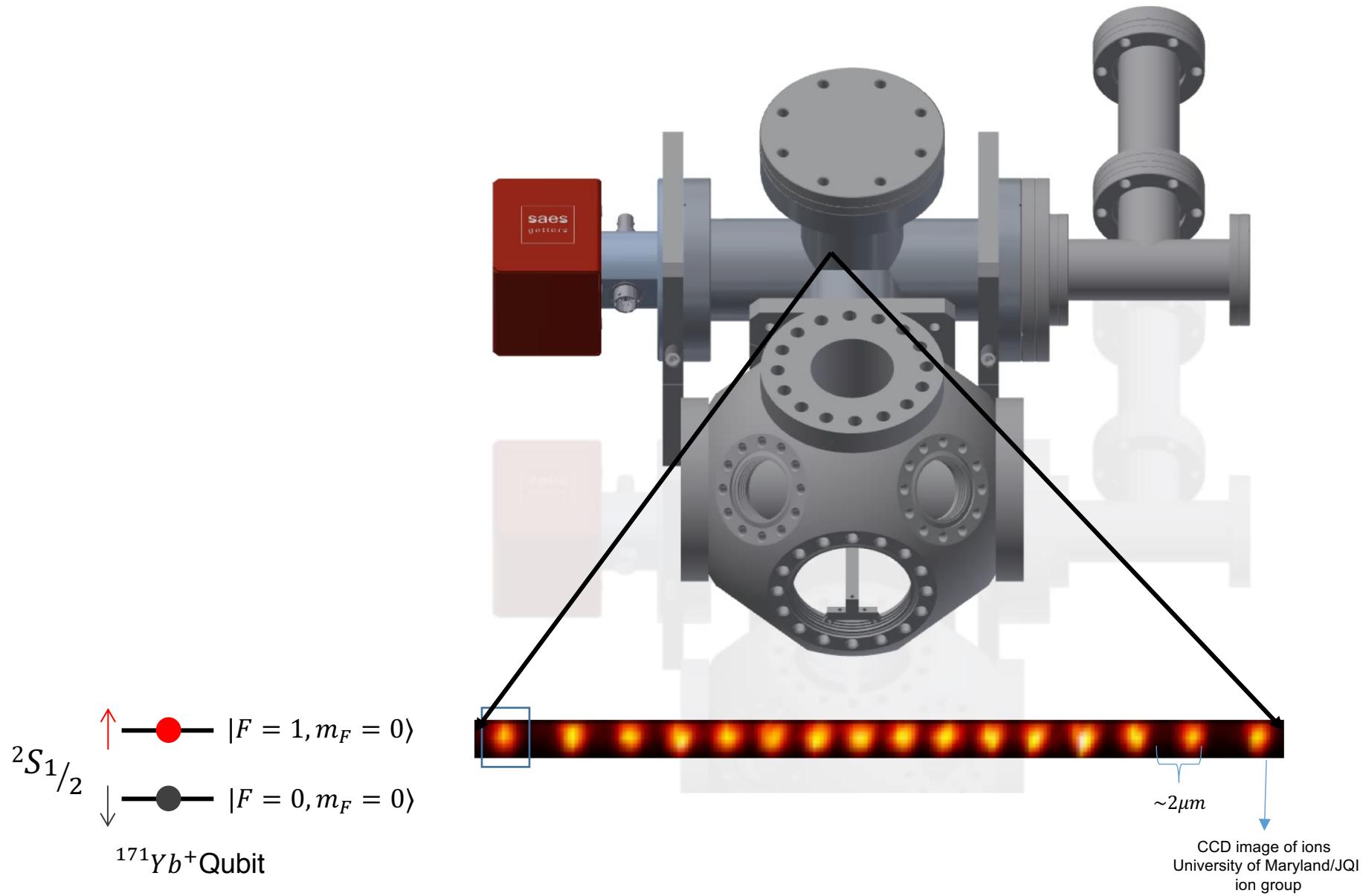
*Ye Wang et al *Nature Photonics* **11**, 646–650 (2017)

Engineer programmable spin interaction graphs

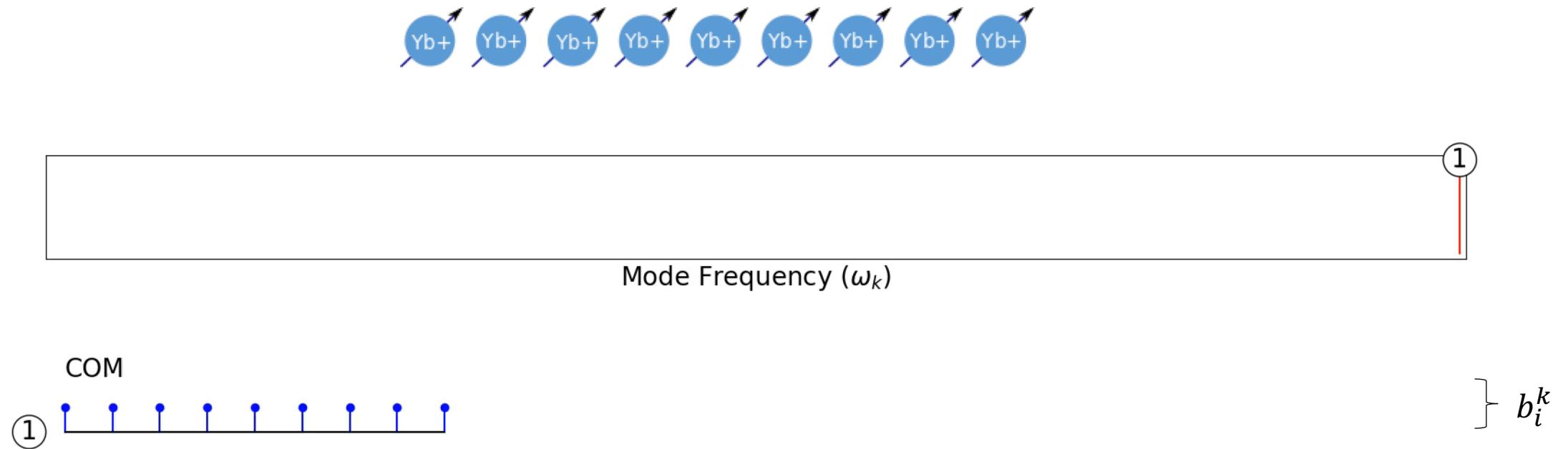


we propose a quantum simulation protocol to engineer XY spin dynamics on a 2D lattice with 1D chain of ions.

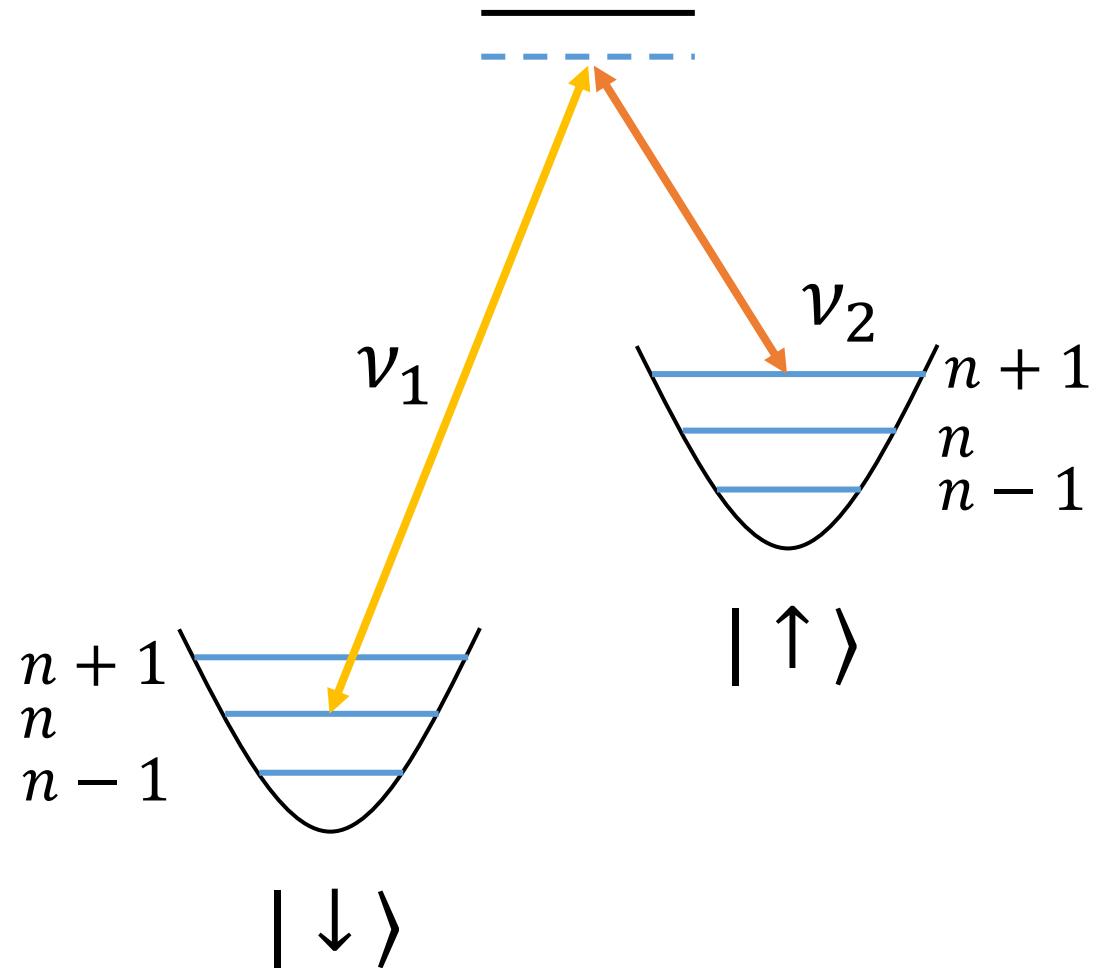
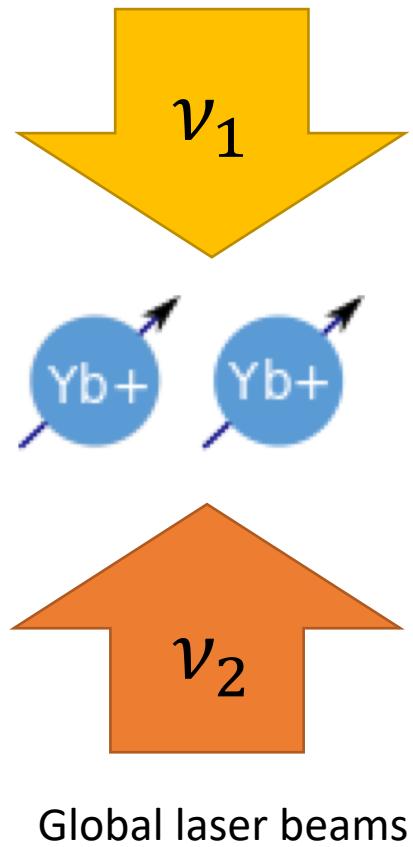
$^{171}Yb^+$ Ion Quantum Simulator



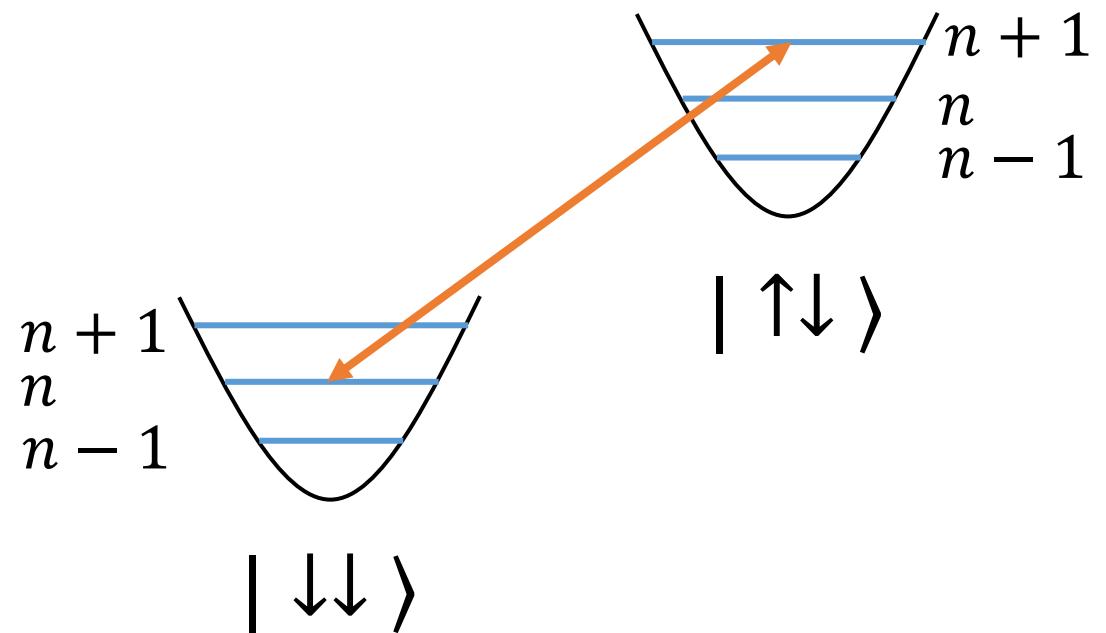
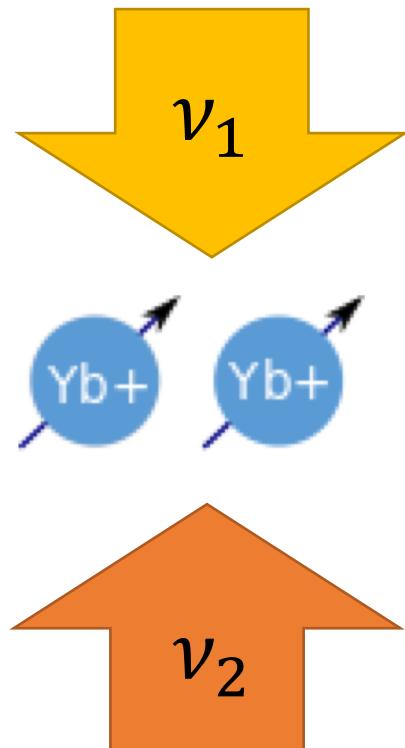
Normal modes of a chain of ions



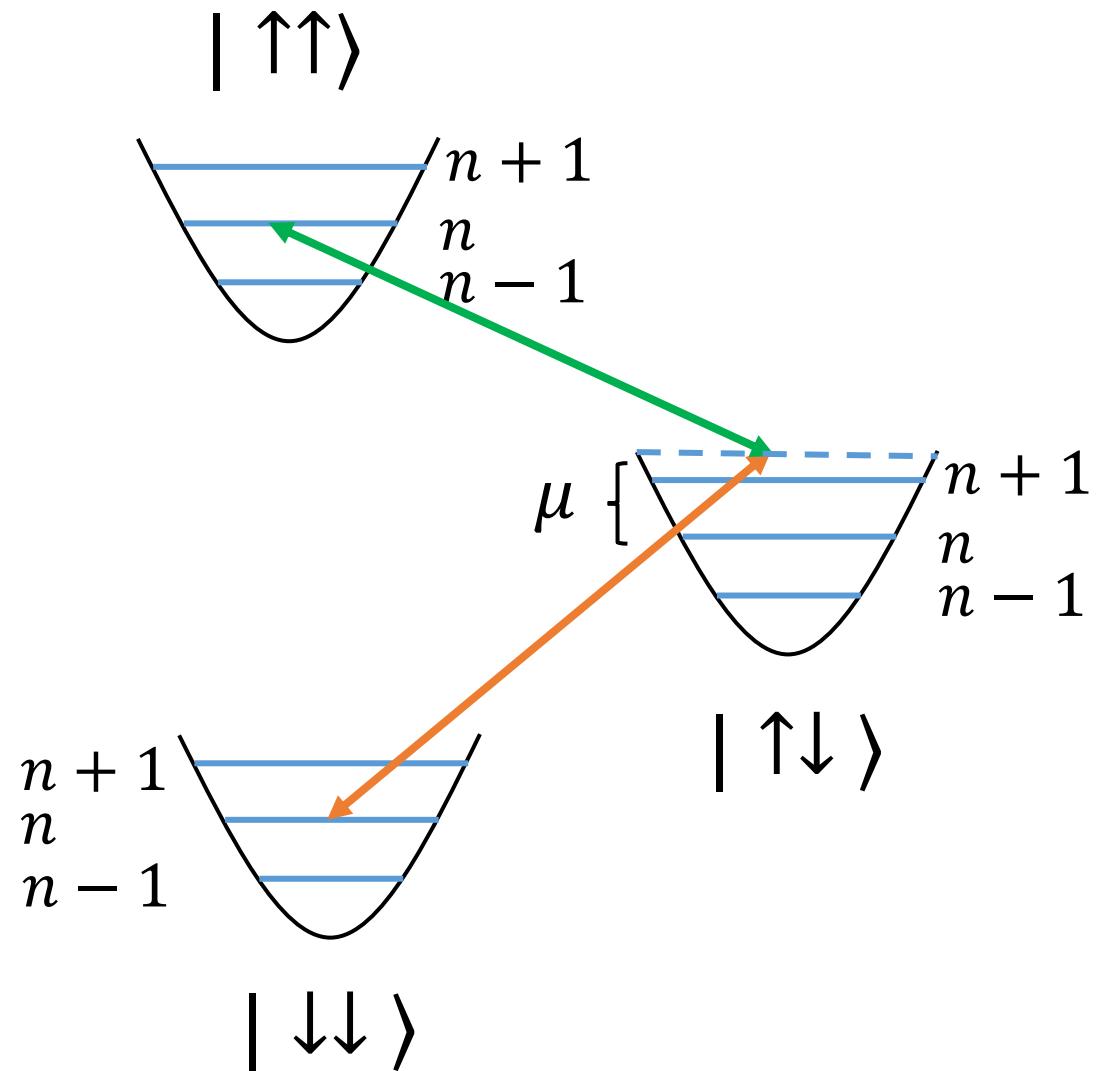
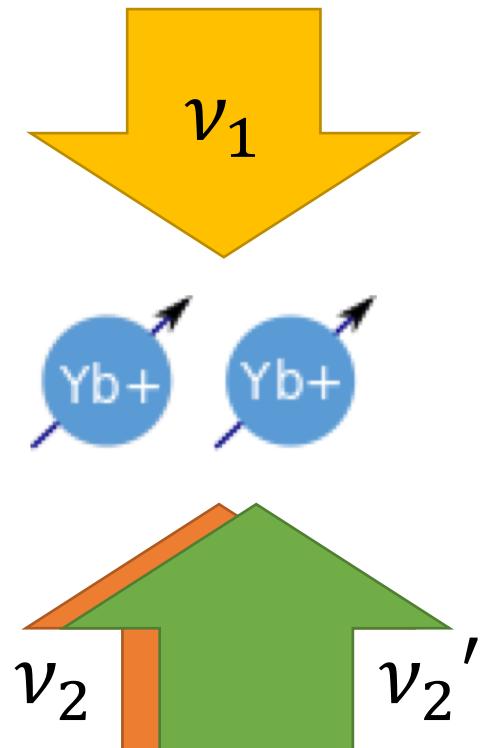
Phonon mediated spin-spin interactions



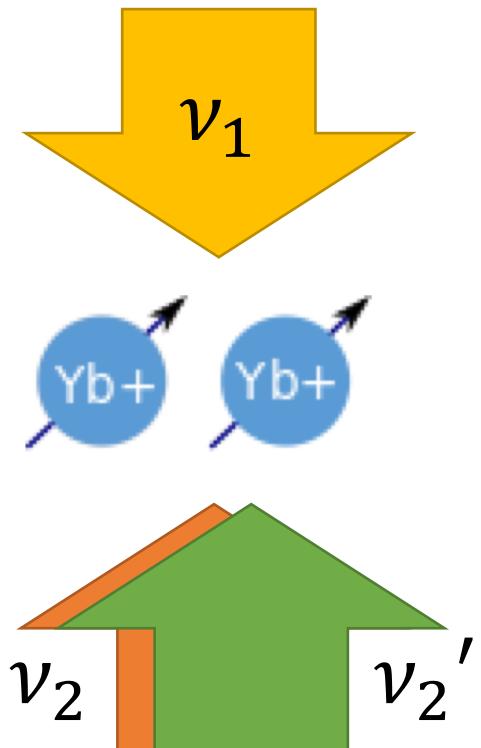
Phonon mediated spin-spin interactions



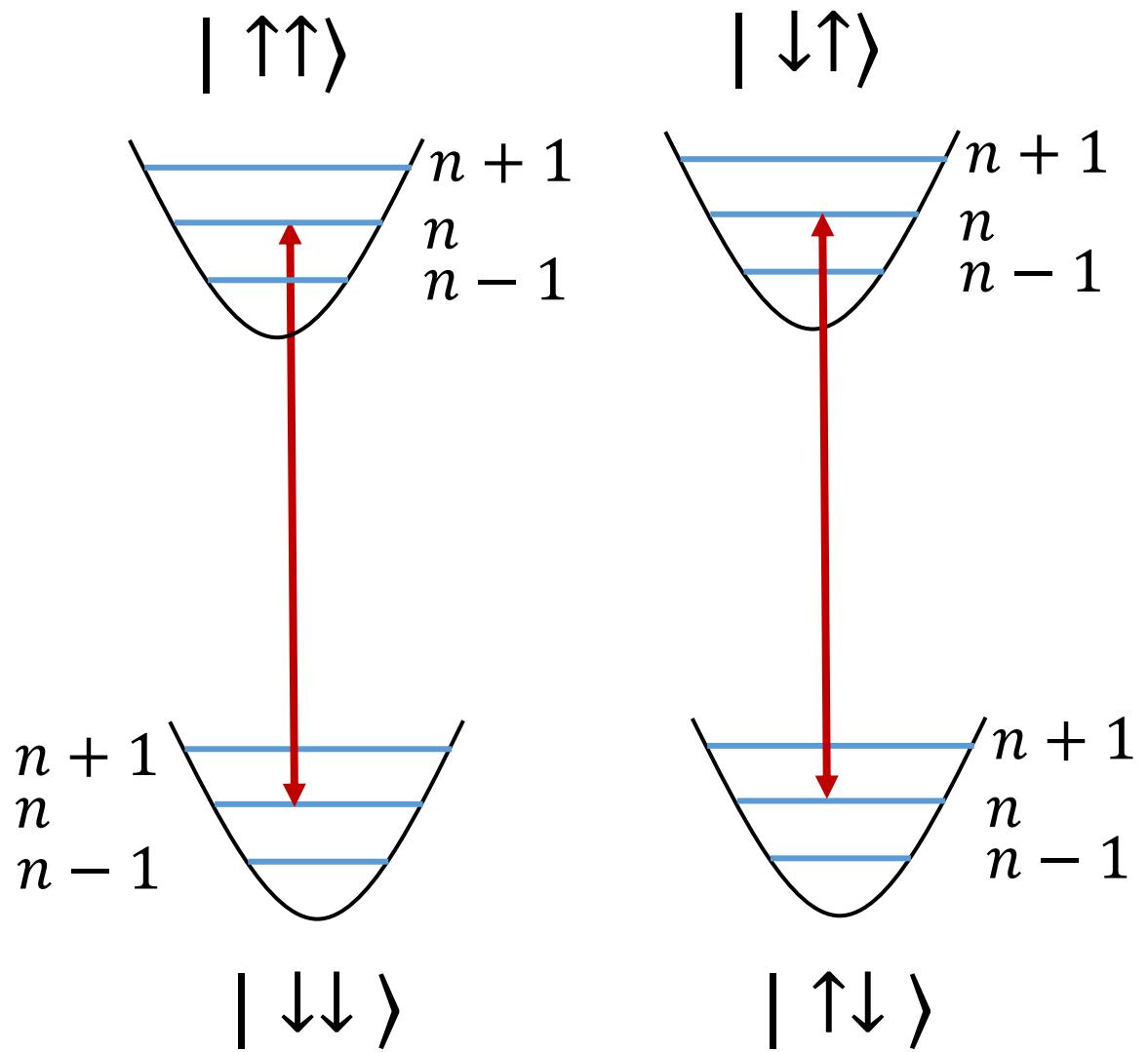
Phonon mediated spin-spin interactions



Phonon mediated spin-spin interactions

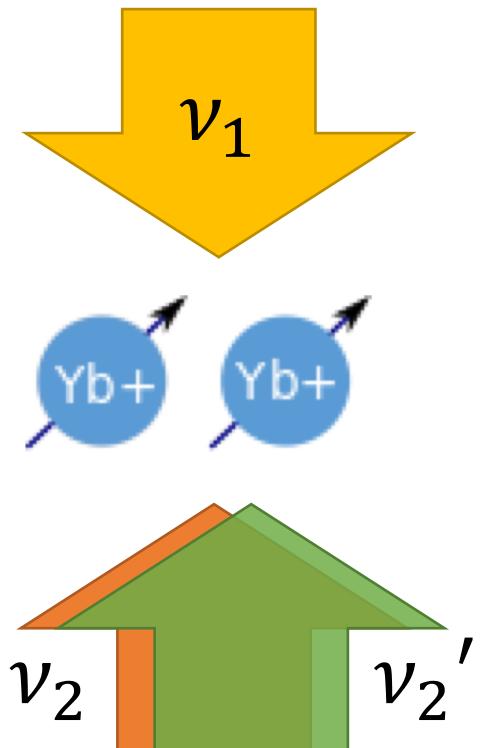


$$H_{eff} = H_{XX} = \sum_{i>j} J_{i,j} S_i^x S_j^x$$



- Cirac, Zoller **PRL** 4091 (1995)
- Molmer, Sorensen **PRL** 1835 (1999)

Phonon mediated spin-spin interactions



$$H_{XX} + B_z \sum_i S_i^z \xrightarrow{|B_z| \gg |J_{i,j}|} H_{XY}$$

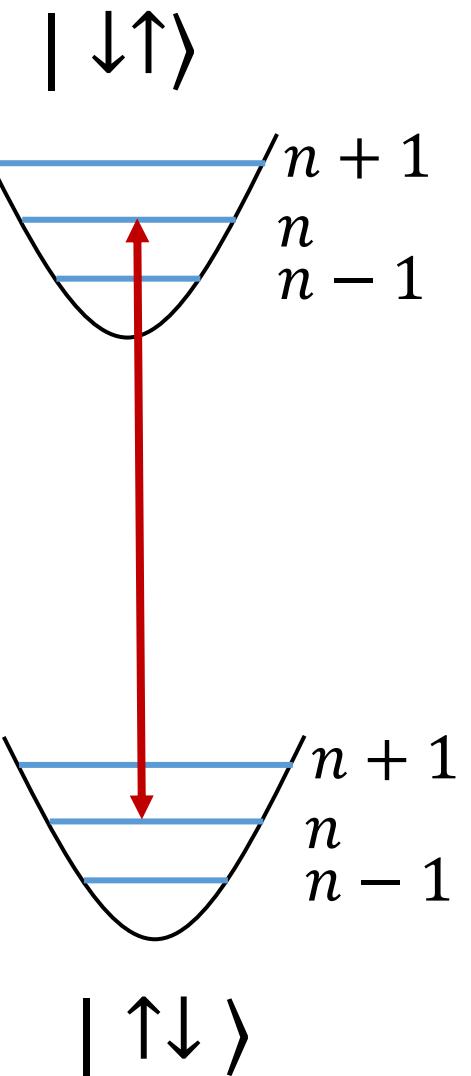
$$H_{XY} = \sum_{i < j} J_{ij} S_i^+ S_j^- + h.c.$$

$$J_{i,j} = \Omega_i \Omega_j \left(\frac{\hbar \Delta k^2}{2m} \right) \sum_k \frac{b_i^k b_j^k}{\mu^2 - \omega_k^2}$$

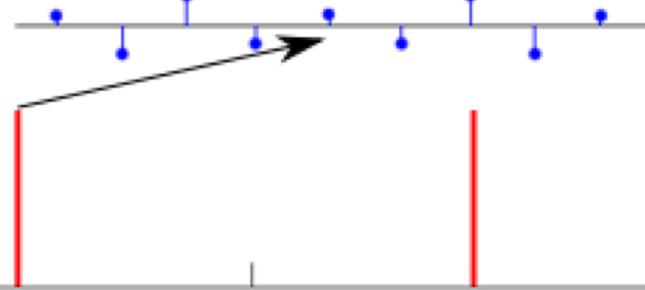
↑
Laser intensity over each ion

Mølmer-Sørensen detuning

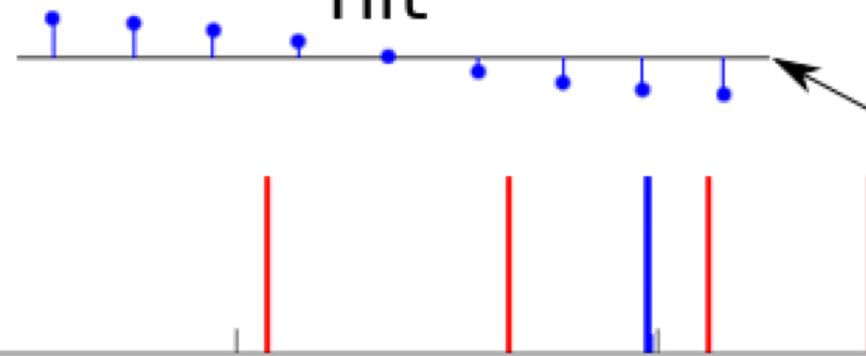
Normal mode eigenvector components



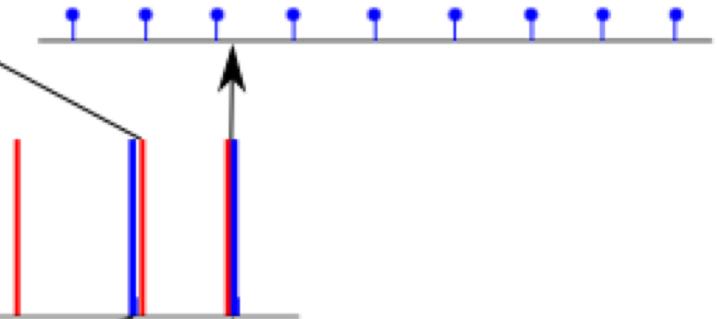
Zig-Zag



Tilt

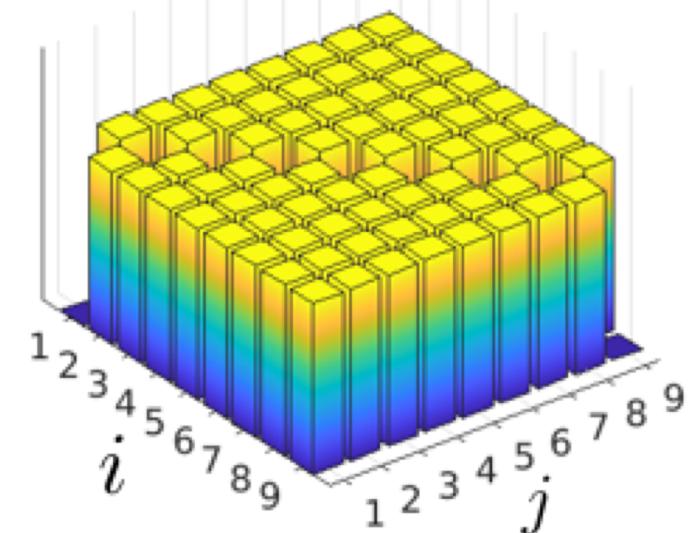
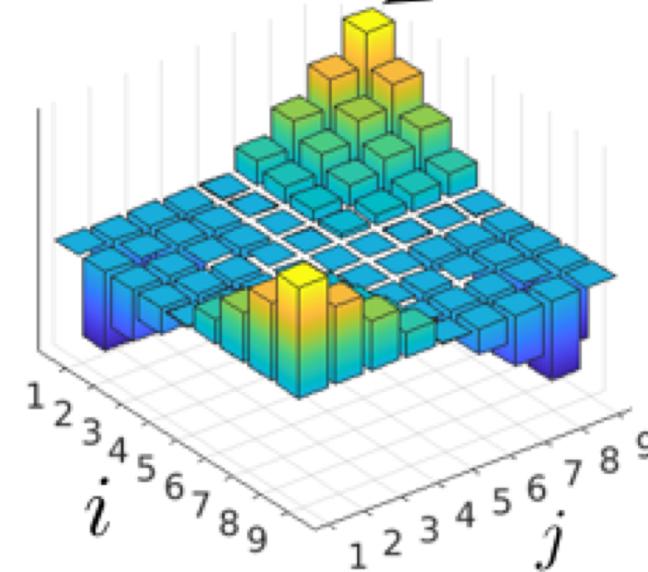
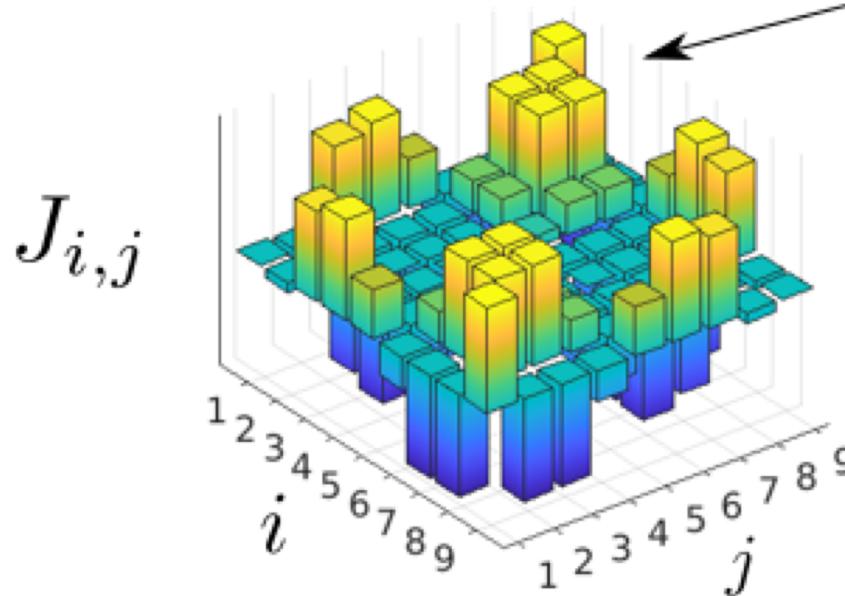


COM



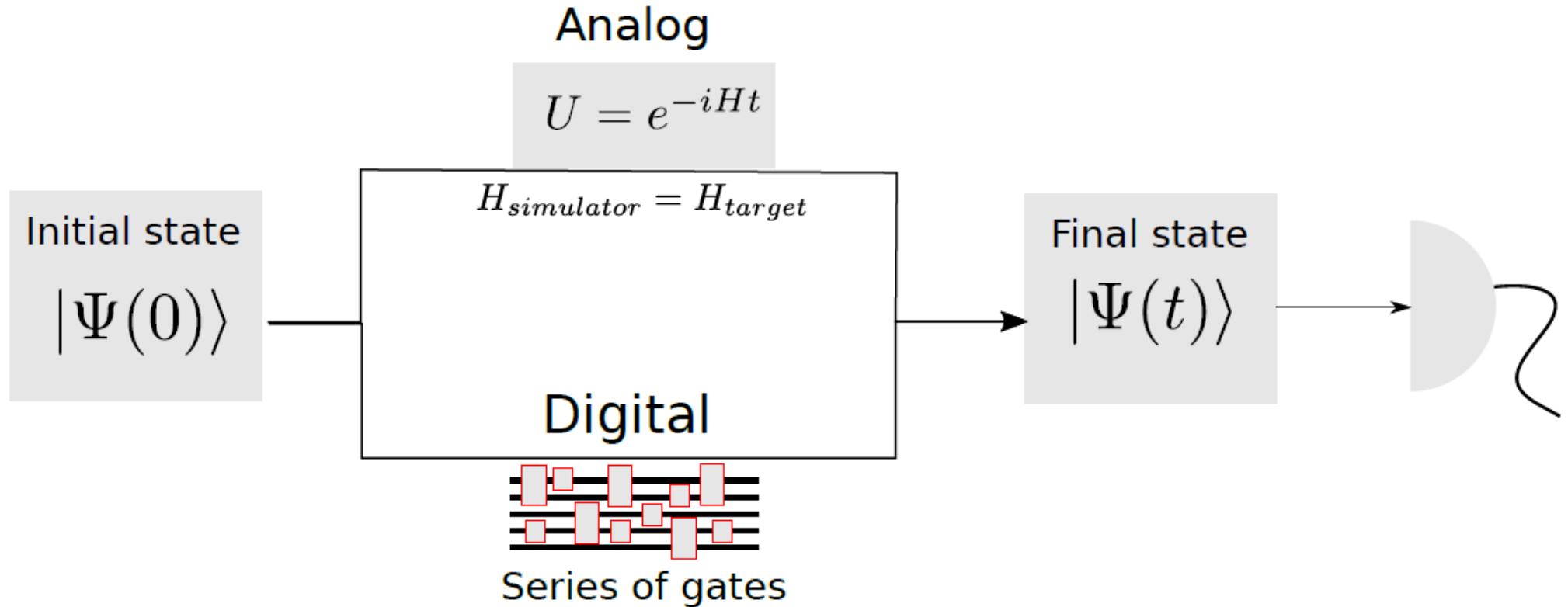
Mode freq

Detuning (μ)



Simulate various interaction profiles by changing the detuning – restricted to only certain types of profiles

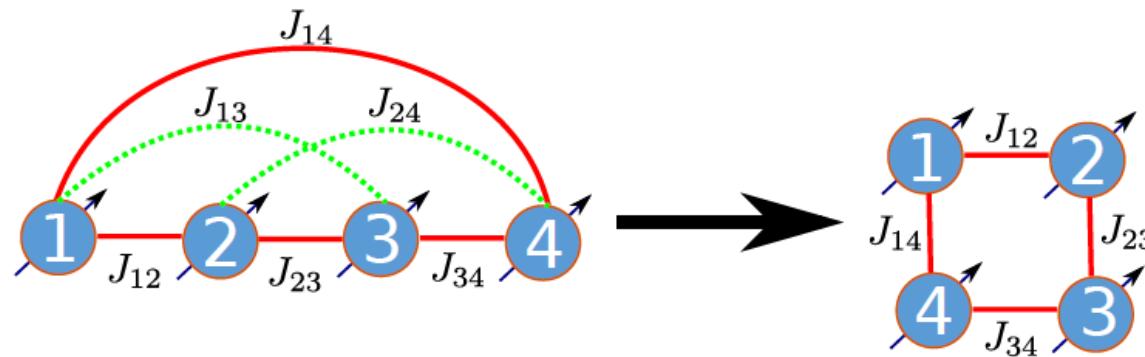
Types of Quantum simulation



Analog : Not Universal, restricted to simulation of certain types of Hamiltonians

Digital : Universal, but errors in simulation due to digitization

Hybrid simulation – Analog + Digital



Analog: Choose an interaction profile that can be engineered with global Raman beams. (Molmer-Sorensen scheme)

Digital: Suppress unwanted interactions and scale other spin interactions by using single spin operations only
e.g., using laser induced site dependent AC stark shift

Hybrid simulation – Analog + Digital

Consider a chain of ions with internal flip-flop Hamiltonian and external AC Stark shift

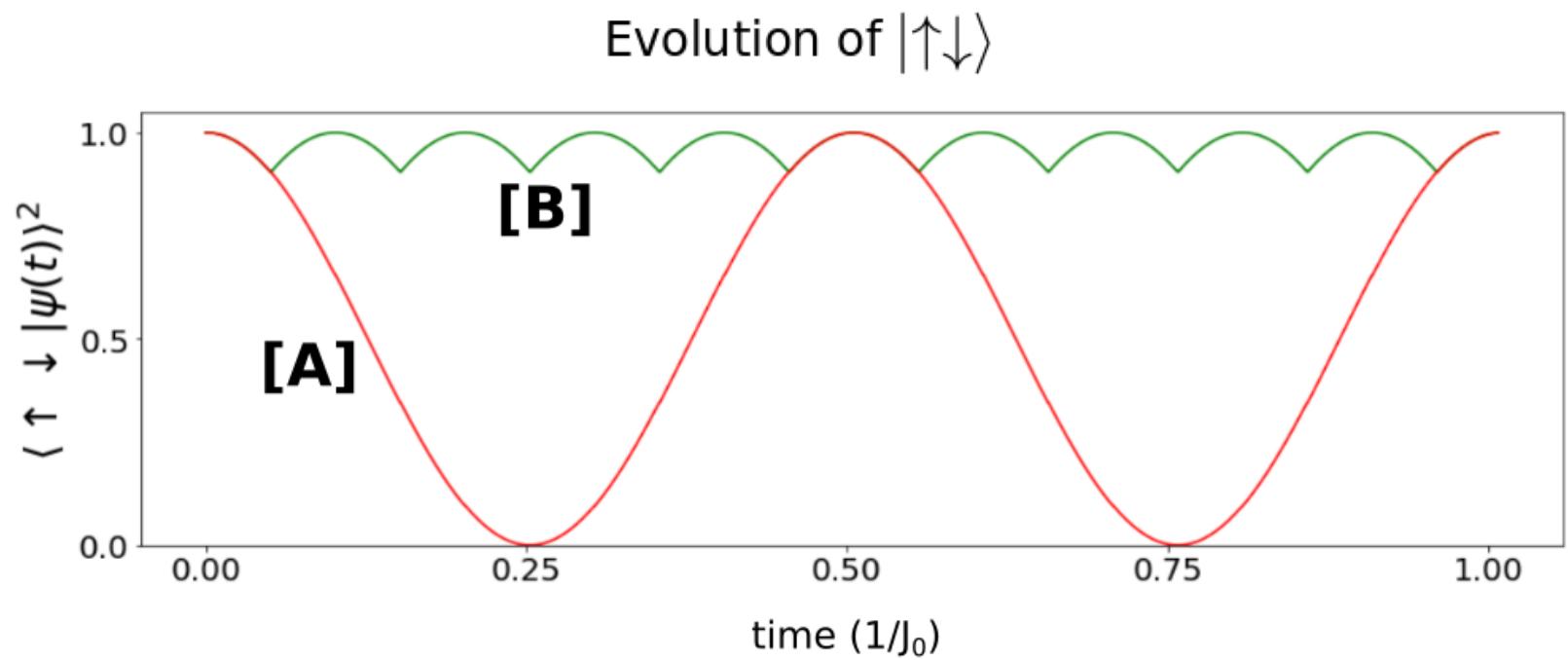
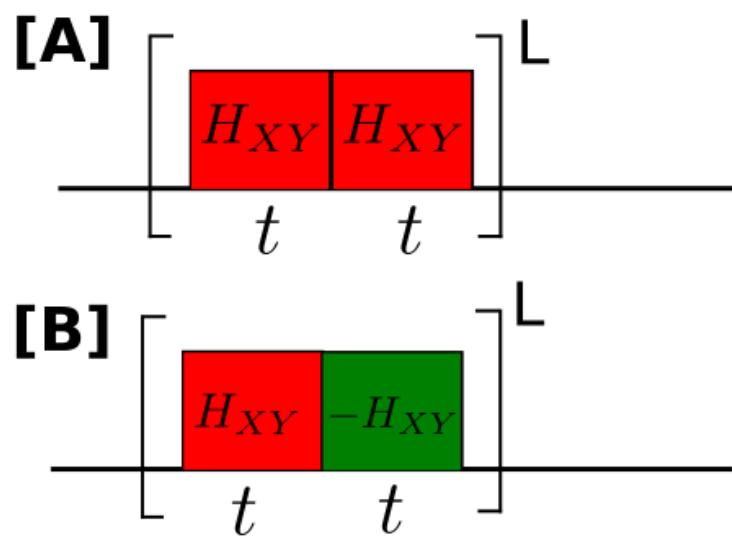
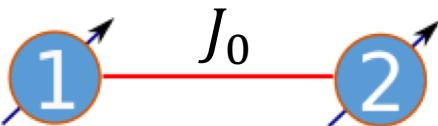
$$H_{XY} = \sum_{i < j} J_{ij} S_i^+ S_j^- + h.c. \quad J_{i,j} = \frac{J_0}{|i-j|^\alpha} \quad \alpha = (0, 3)$$

Created using a global Molmer-Sorensen scheme

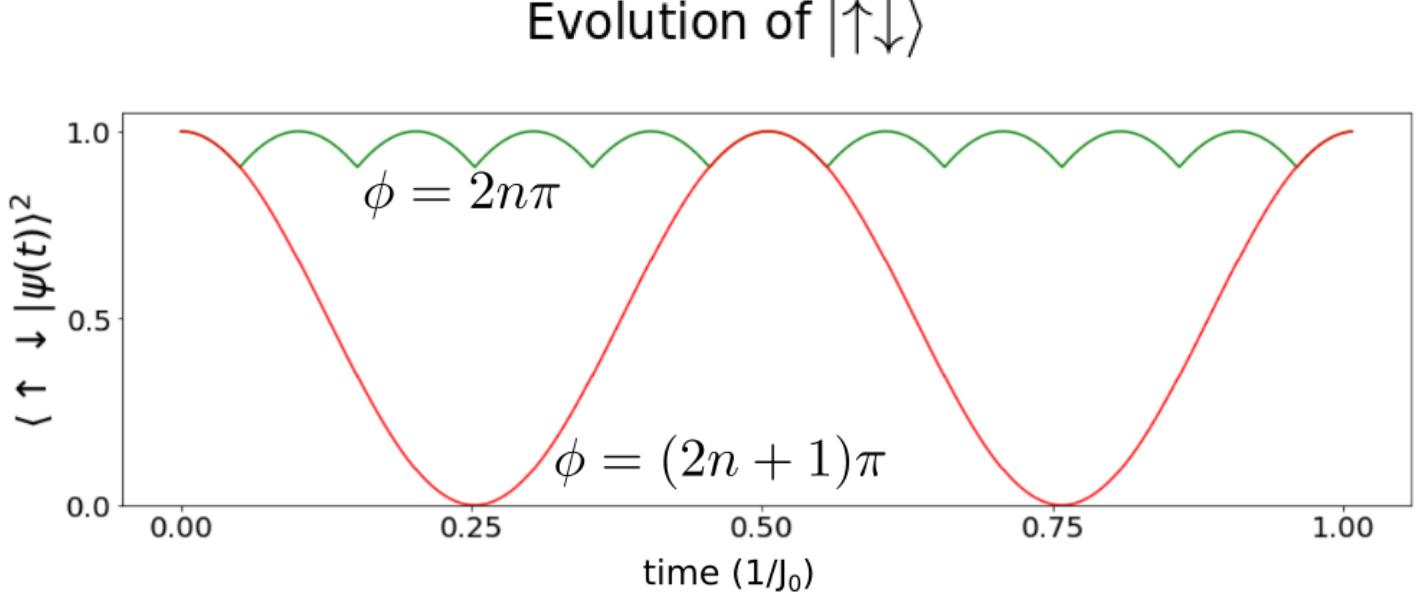
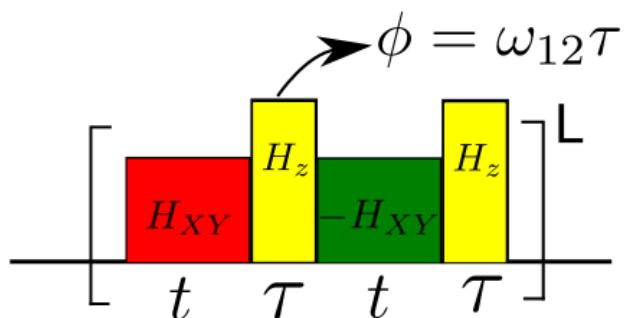
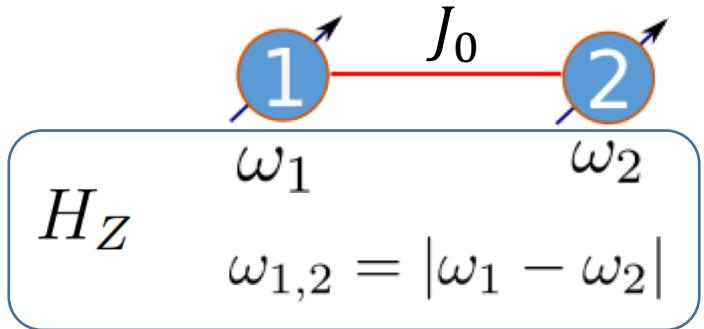
$$H_{stark} = H_Z = \sum_{i=1}^N \omega_i S_i^z$$

Created using an off-resonant laser beam with site dependent intensity inducing a differential AC Stark shift

Case of 2 ions



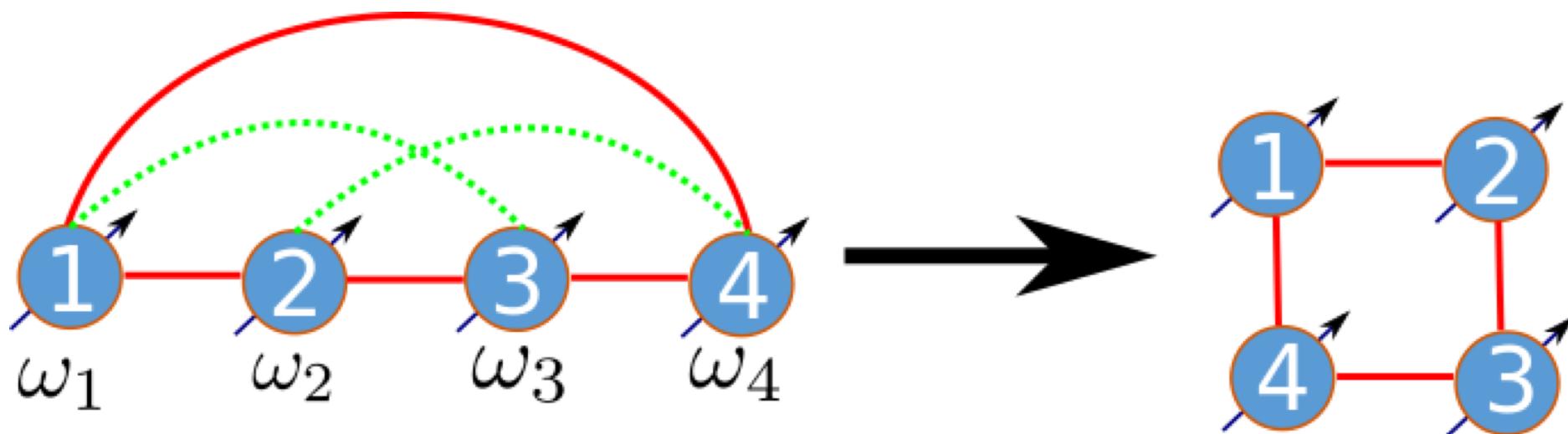
Case of 2 ions



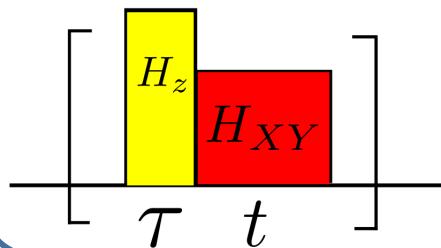
Pulse sequence cancels interaction if $\omega_{i,j}\tau = 2n\pi$

retains interaction if $\omega_{i,j}\tau = (2n+1)\pi$

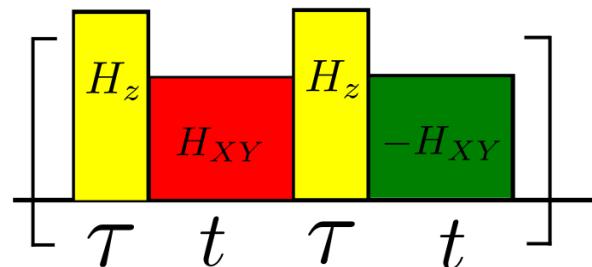
Goal



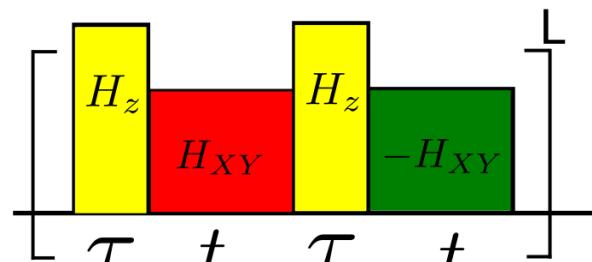
Average Hamiltonian for a sequence of pulses



$$\tilde{H}_{ff} = e^{-iH_Z\tau} H_{XY} e^{iH_Z\tau} = \sum_{i>j} J_{i,j} S_i^+ S_j^- e^{i\omega_{ij}\tau} + h.c$$



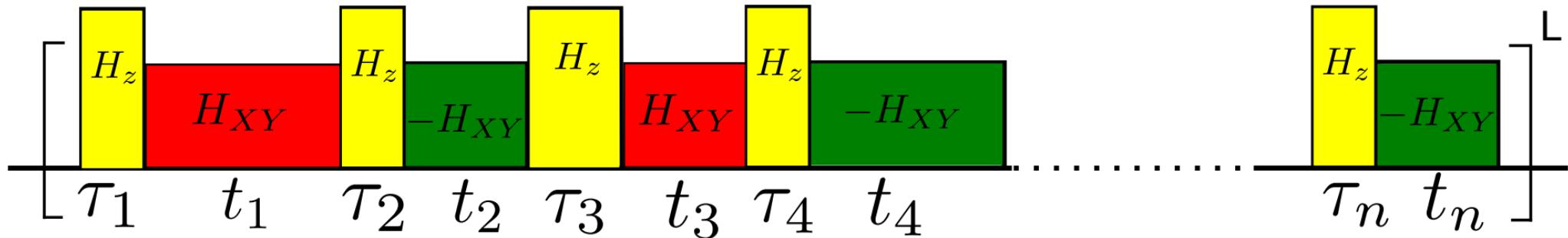
$$H_{avg} = \sum_{i>j} J_{i,j} S_i^+ S_j^- (e^{i\omega_{ij}\tau} - e^{2i\omega_{ij}\tau}) + h.c$$



$$H_{avg} = \sum_{i>j} J'_{i,j} S_i^+ S_j^- + h.c \quad \text{Valid if } J_0 * T_{cycle} \ll 1$$

$$J'_{i,j} = \frac{J_{ij}}{L} * (e^{i\omega_{ij}\tau_{tot}} - e^{2i\omega_{ij}\tau_{tot}} + \dots e^{Li\omega_{ij}\tau_{tot}})$$

Suppresses the interactions (i,j) where $\omega_{ij}\tau_{tot} = 2n\pi$



$$H_{avg} = \sum_{i>j} J''_{i,j} S_i^+ S_j^- + h.c$$

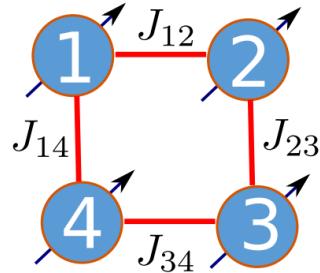
$$J''_{i,j} = \frac{J'_{i,j}}{T_{cycle}} * (t_1 e^{i\omega_{ij}\tau_1} - t_2 e^{i\omega_{ij}(\tau_1+\tau_2)} \dots)$$

Use this to scale (also cancel) the remaining interactions

Task : Find a solution for $\{\omega_i, \tau_k, t_k\}$ to simulate 2D lattices

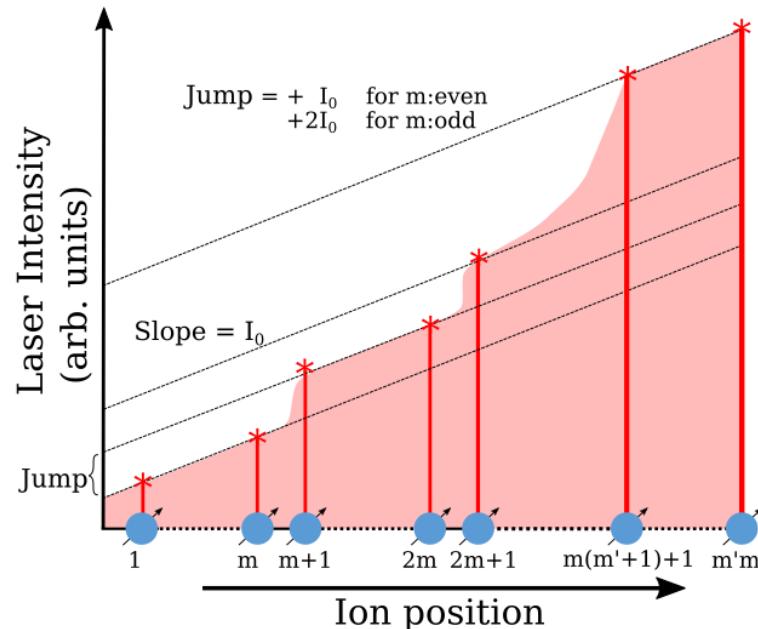
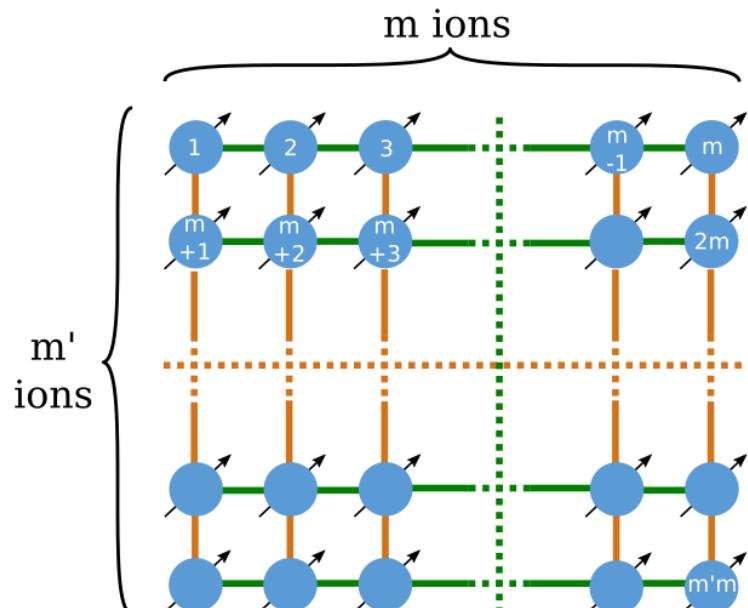
Stark shift ω_i pattern to simulate 2D lattices

$$\begin{aligned}\text{Bond class} &= \{(1, 2), (2, 3), (3, 4), (1, 4)\} \\ \text{NoBond class} &= \{(1, 3), (2, 4)\}\end{aligned}$$



Find ω_i such that

$$\{\omega_{ij} \mid (i, j) \in Bond\} \cap \{\omega_{ij} \mid (i, j) \in NoBond\} = \emptyset$$

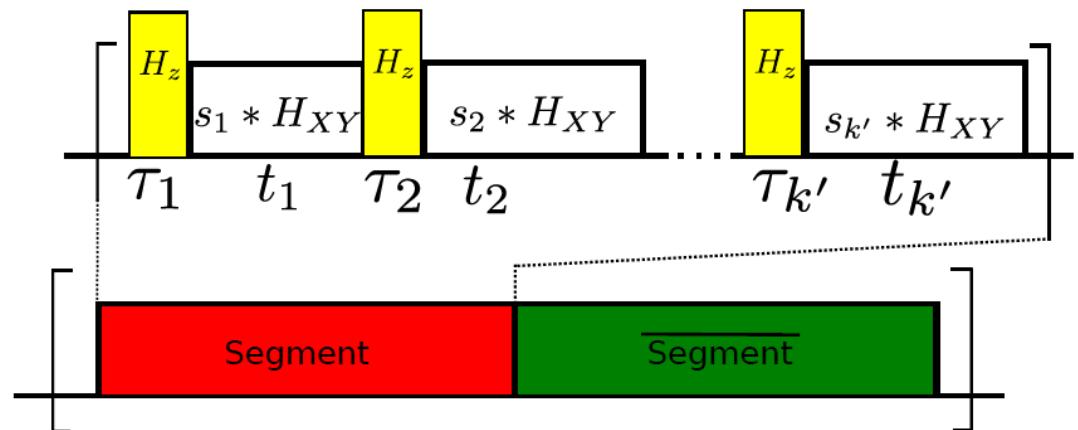


Construction of pulse sequence

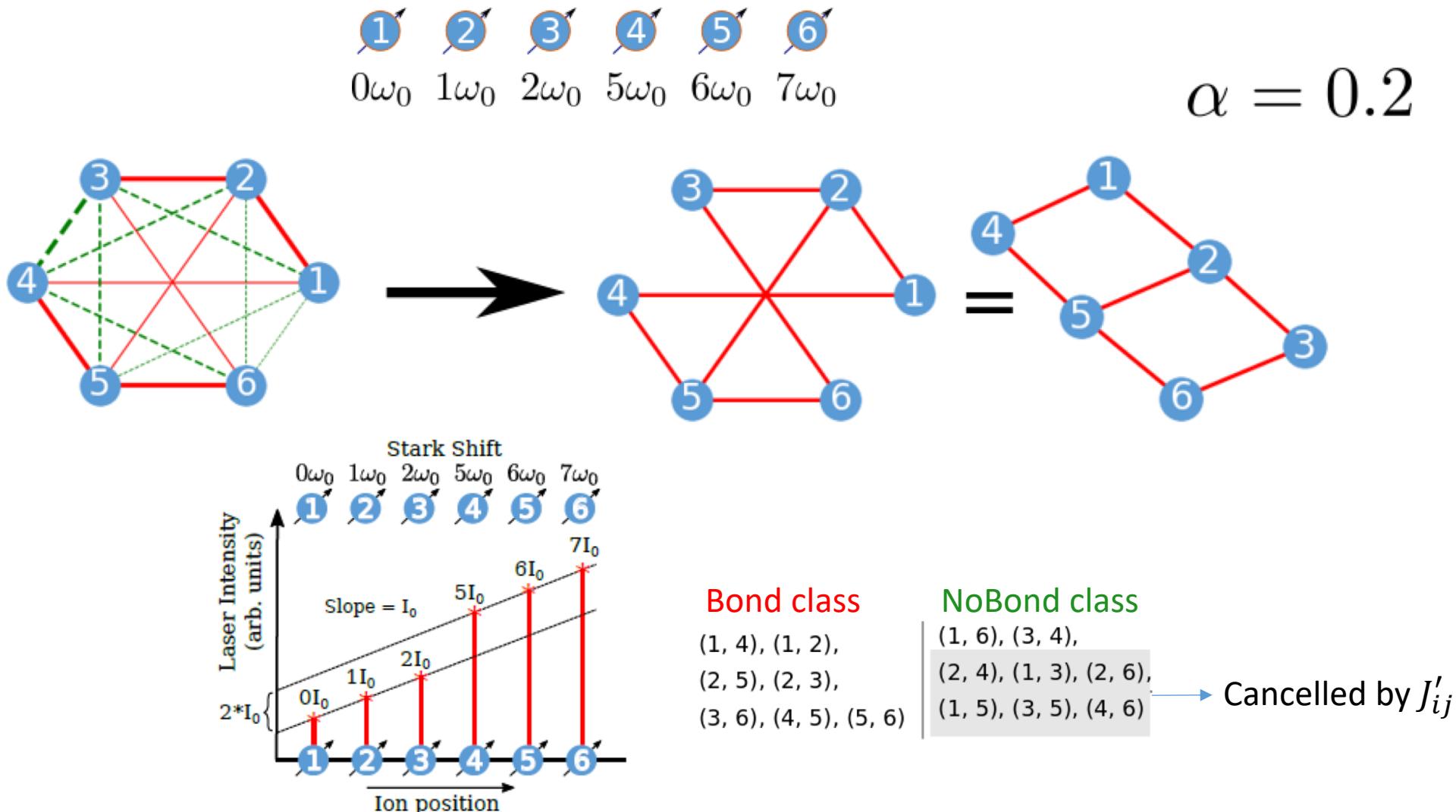
$$\begin{aligned}
 J''_{ij}(\{t_k, \tau_k\}) = F(\omega_{ij} * \tau_{tot}) &= 0 && \text{for } (i, j) \in \text{Nobond class} \\
 &= 1/m^\alpha && \text{for } (i, j) \in \text{Bond class and } \omega_{ij} = \omega_0 \\
 &= 1 && \text{for } (i, j) \in \text{Bond class and } \omega_{ij} \neq \omega_0
 \end{aligned}$$

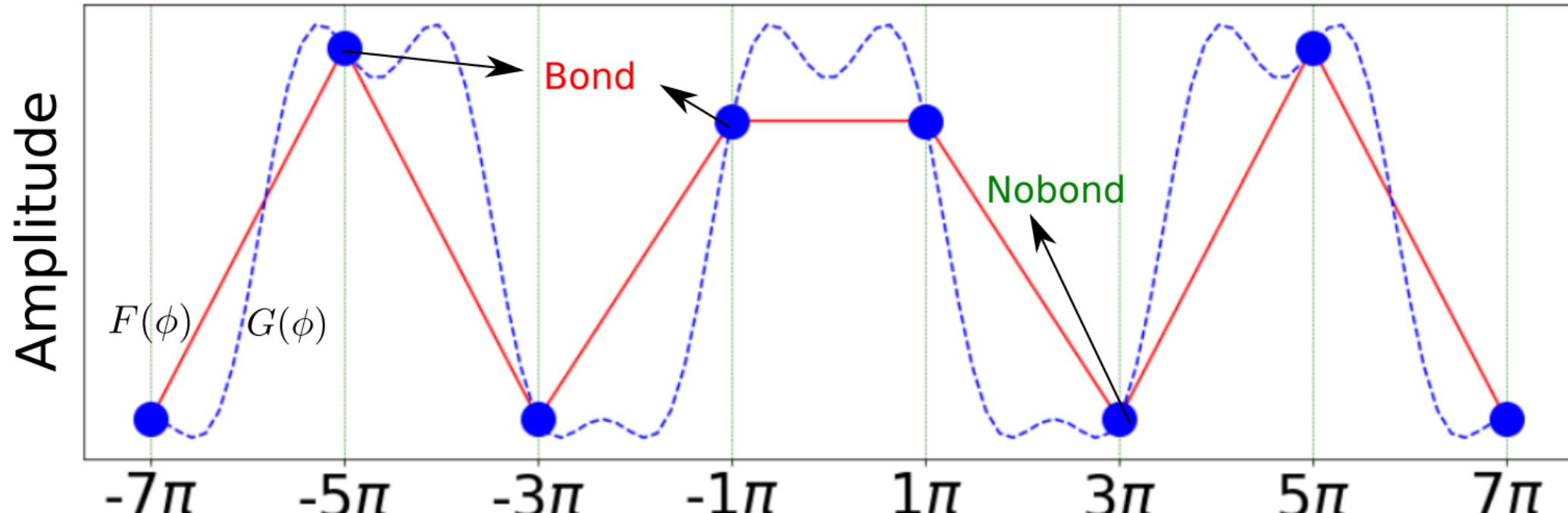
Fit with function $G(\phi) = \sum_{l=0}^{l'} a_l * \cos(iW\phi)$ obtain optimum values of $\{a_l, W\}$

$$\begin{aligned}
 t_k &= |a_k/2| && \text{for } k = (1, 2, \dots, l') \\
 &= t_{k-l'} && \text{for } k = (l' + 1, \dots, 2l') \\
 &= |a_0| && \text{for } k = 2l' + 1 \\
 \\
 \tau_k &= \tau_{tot} * W && \text{for } k \neq (l' + 1) \\
 &= \tau_{tot} - (l' - 1)\tau_1 && \text{for } k = l' + 1 \\
 \\
 s_k &= \text{sign}(a_k) && \text{for } k = (1, 2, \dots, l') \\
 &= -s_{k-l'} && \text{for } k = (l' + 1, \dots, 2l') \\
 &= -\text{sign}(a_0) && \text{for } k = 2l' + 1
 \end{aligned}$$



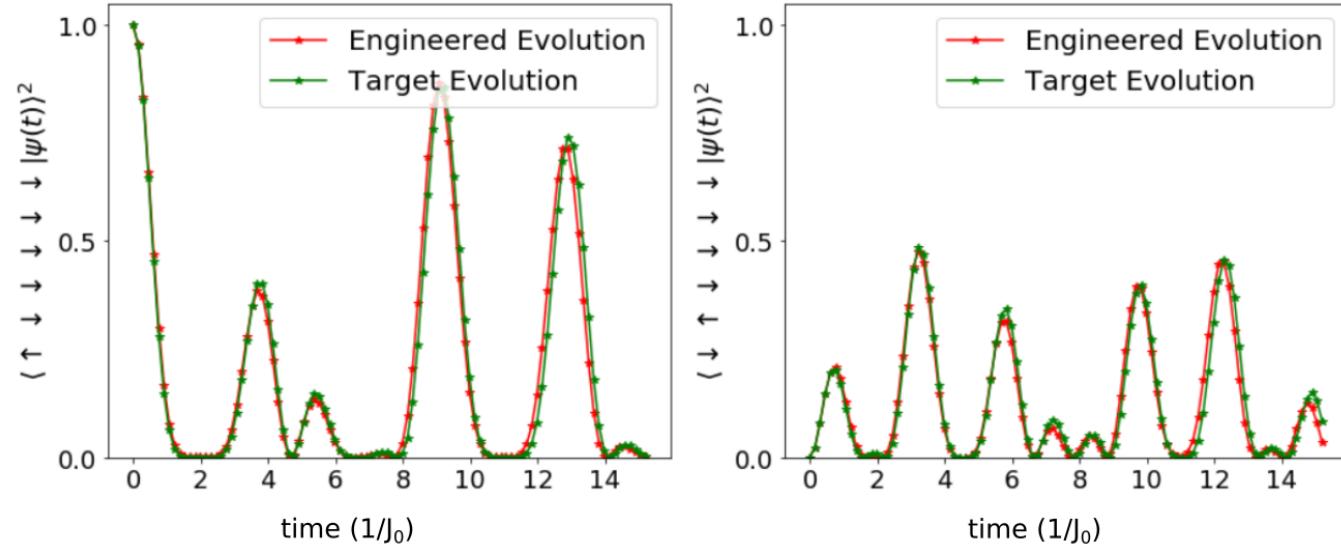
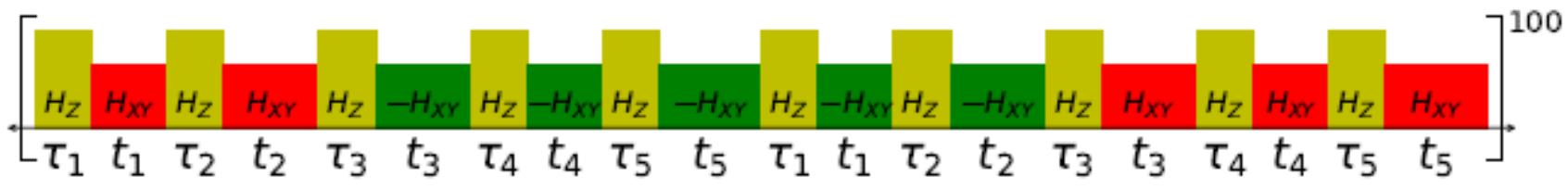
Simulating 3x2 lattice with 6 ion chain

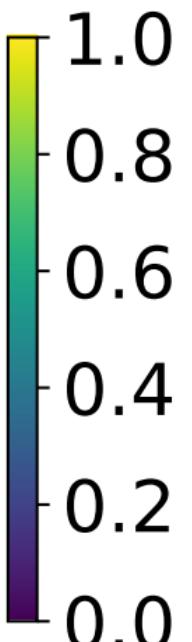
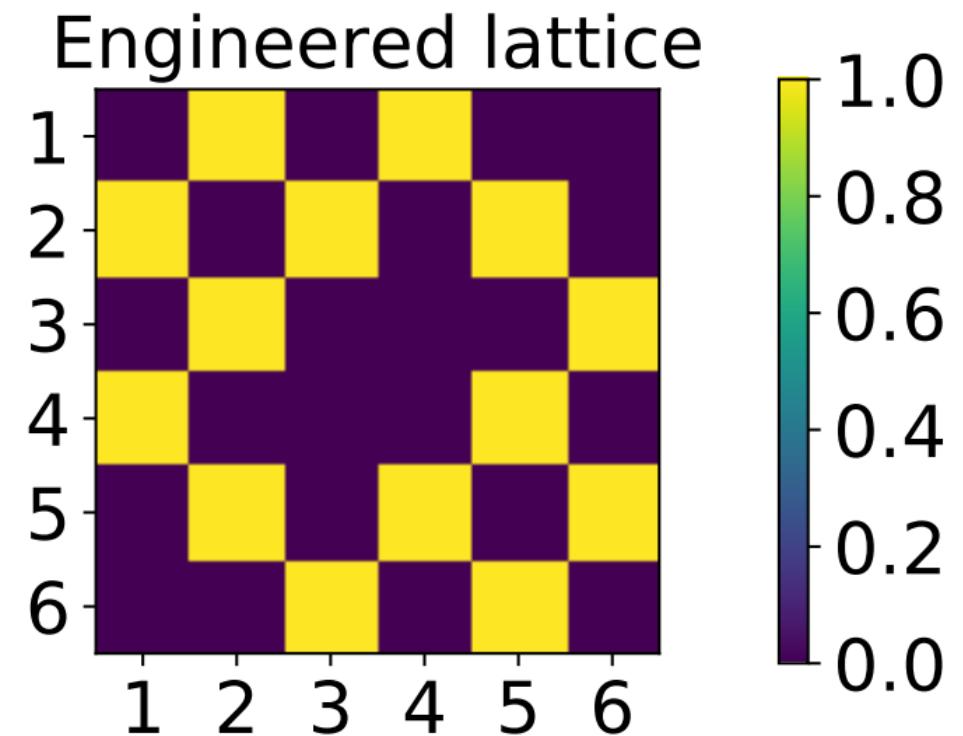
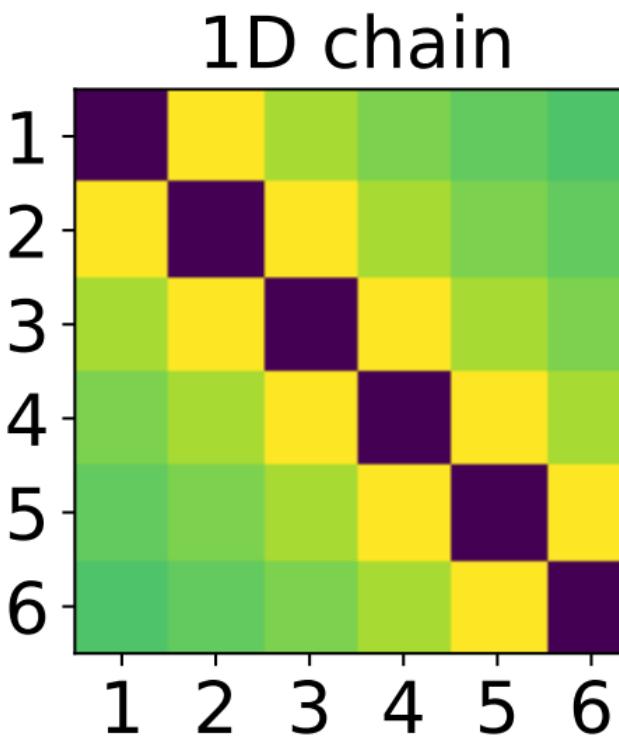
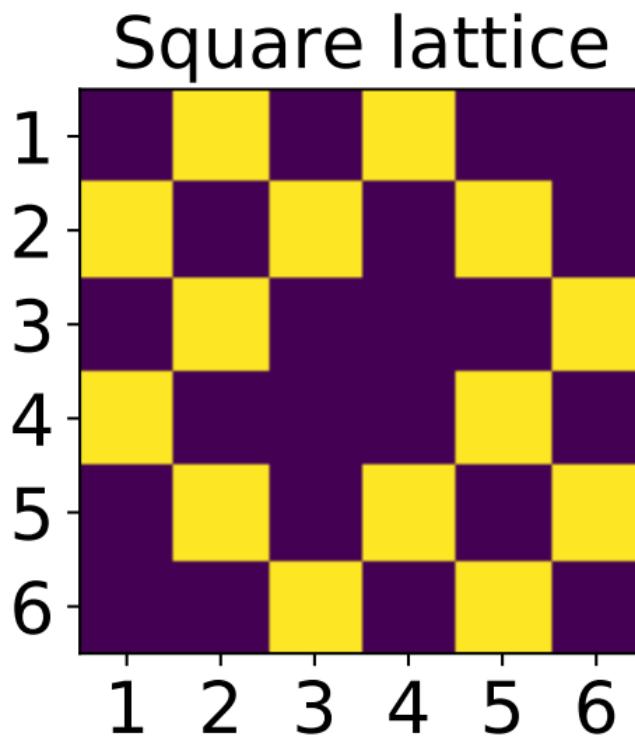




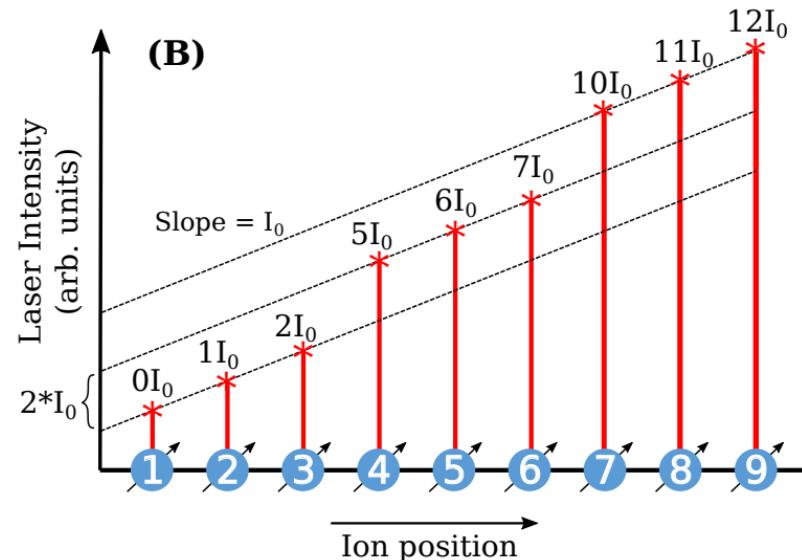
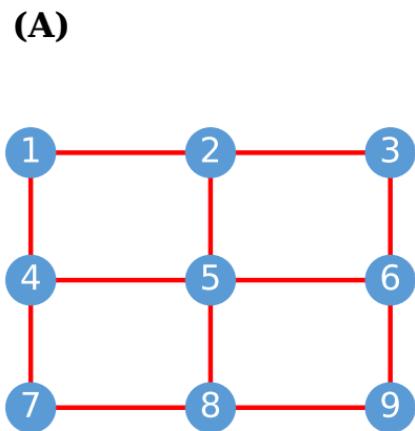
$$\phi = \omega_{ij}\tau$$

	1	2	3	4	5
$\tau(\mu s)$	6.283	6.283	6.283	6.283	6.283
$t(\mu s)$	2.989	14.961	14.961	2.989	26.055
<i>sign</i>	+1	+1	-1	-1	-1



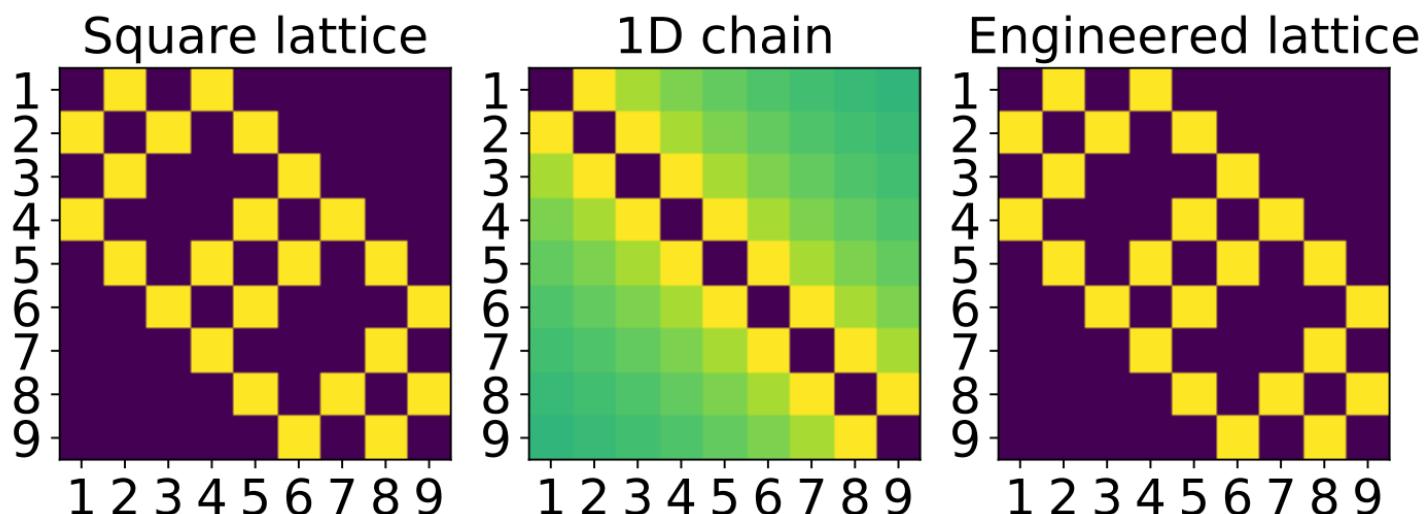


3x3 lattice with 9 ions



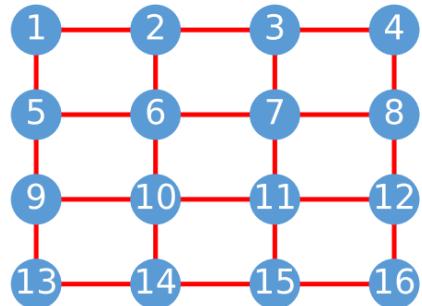
Fit parameters

W	a_0	a_1	a_2	a_3	a_4
0.1	0.360	0.305	-0.140	0.188	0.499

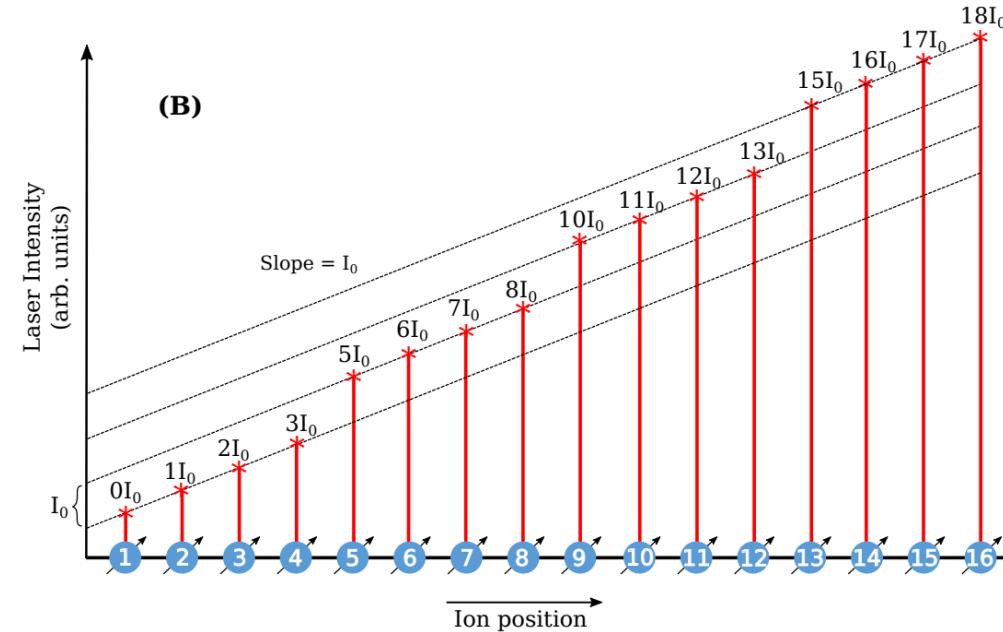


4x4 lattice with 16 ions

(A)

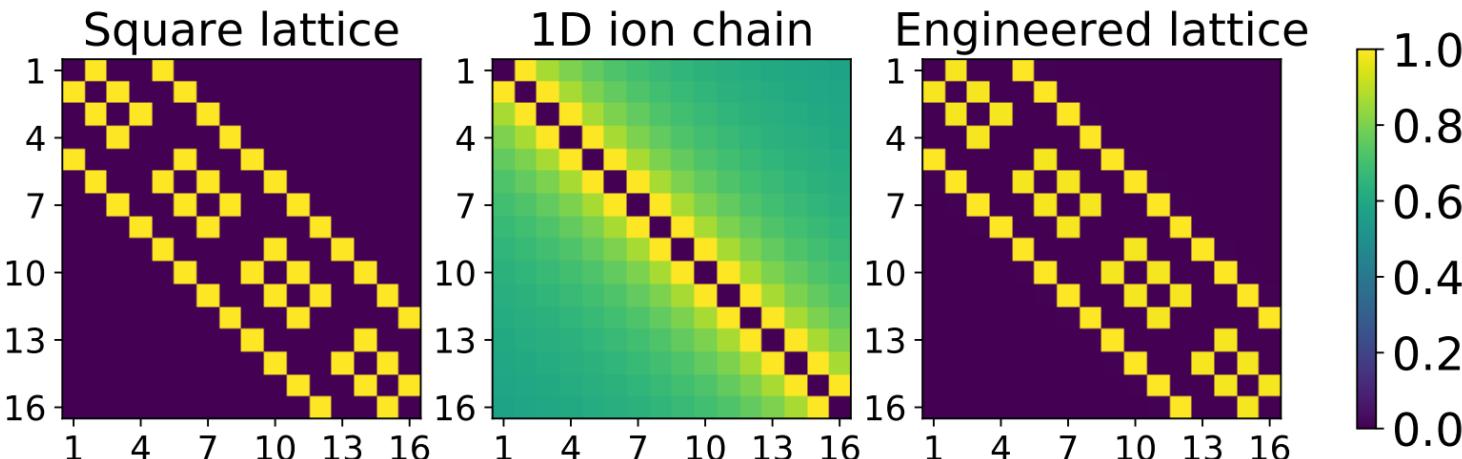


(B)

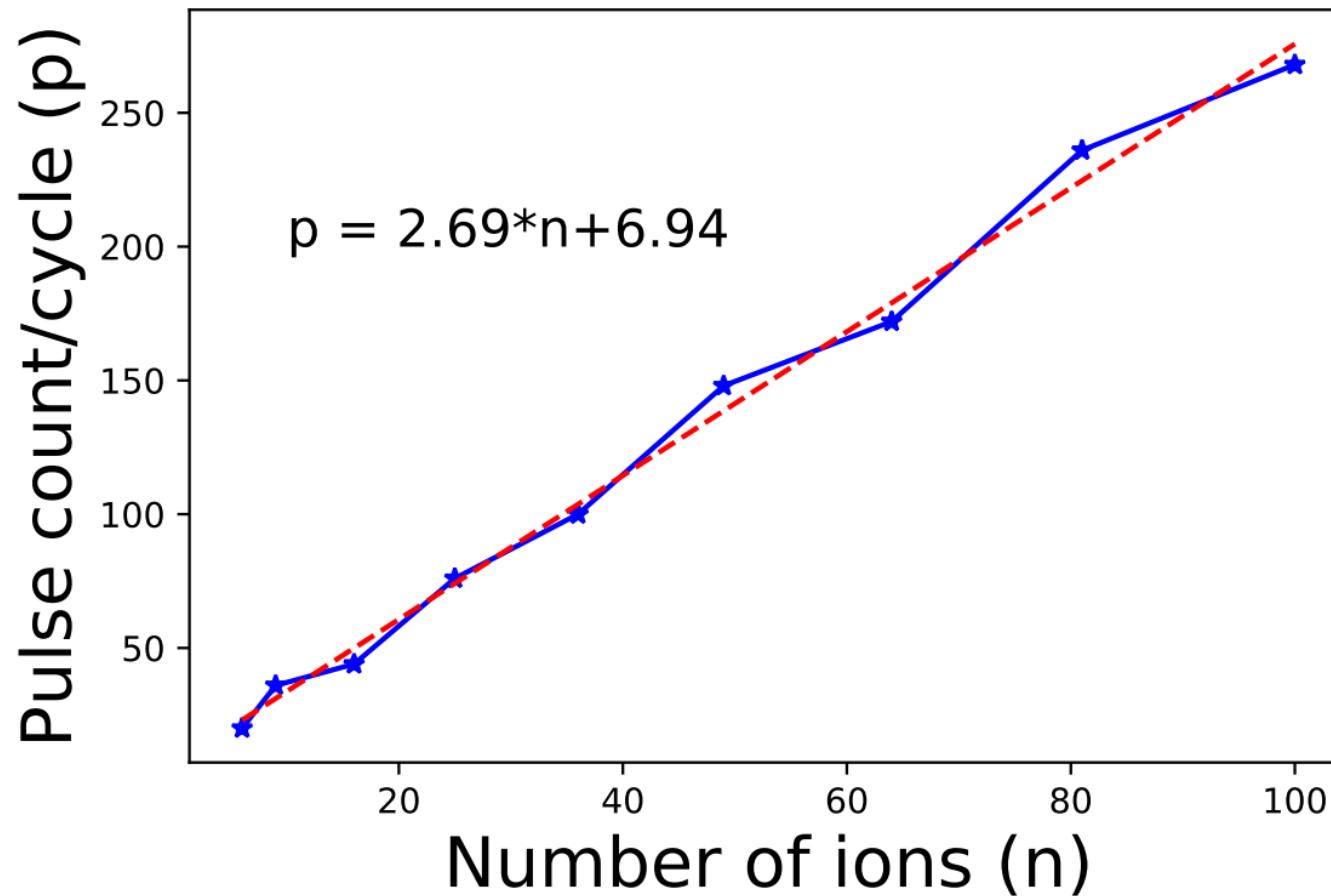


Fit parameters

W	a_0	a_1	a_2	a_3	a_4	a_5
0.083	0.293	0.33	-0.07	-0.057	0.293	0.387



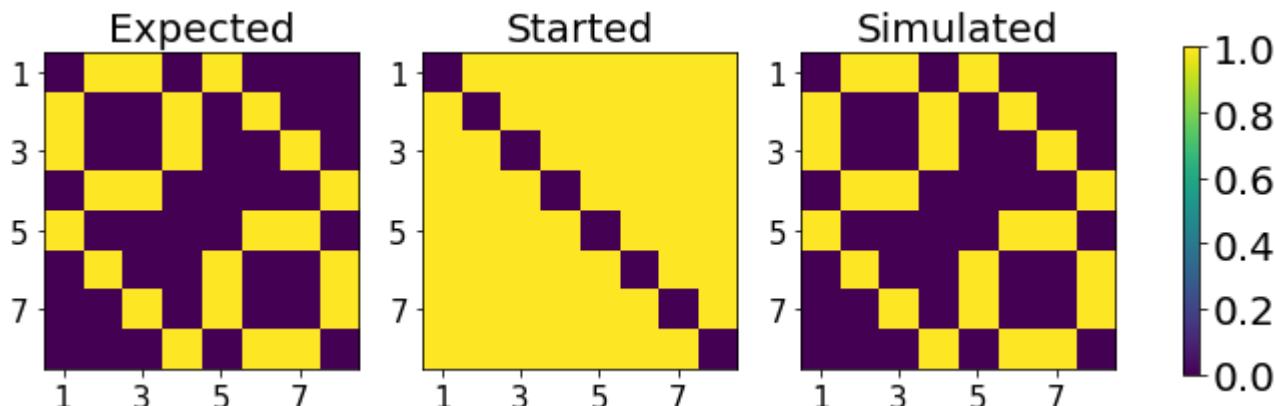
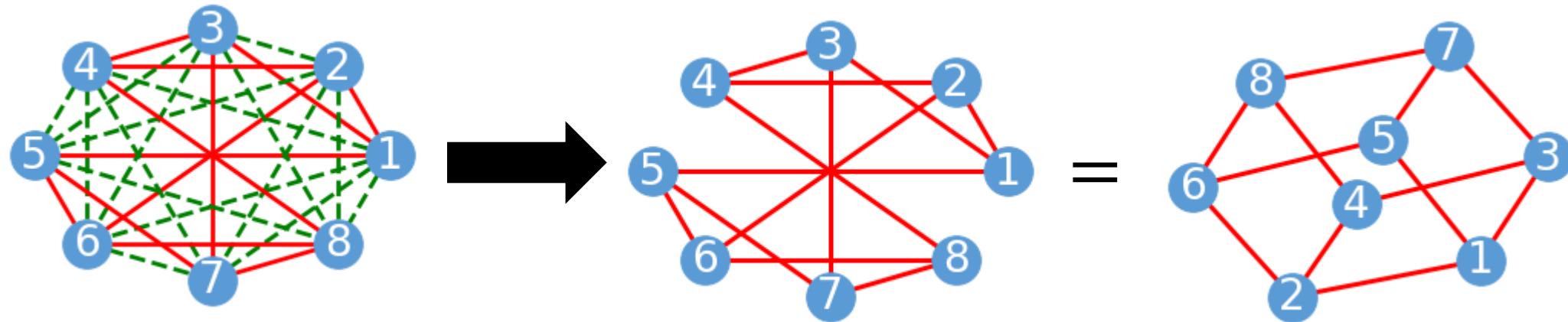
Scaling with the number of ions



For a target 2D square lattice

Extension to 3D lattice

2x2x2 lattice with 8 ions $\alpha = 0$

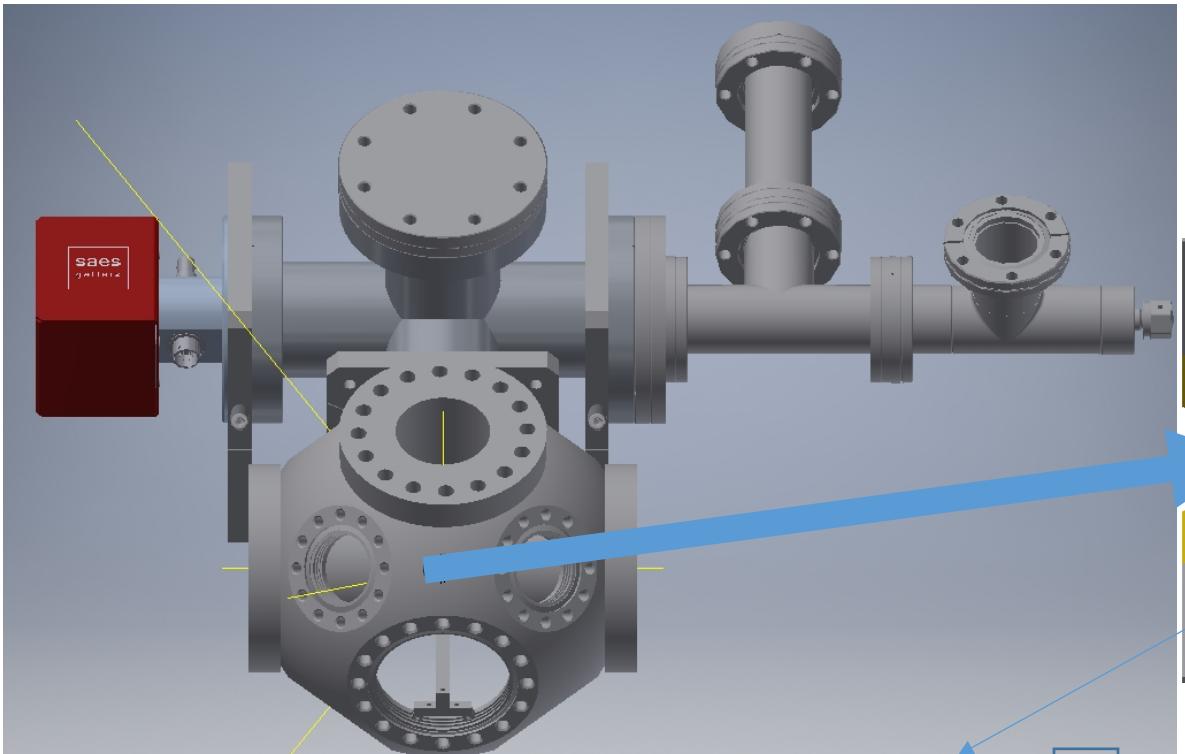


Summary and Outlook

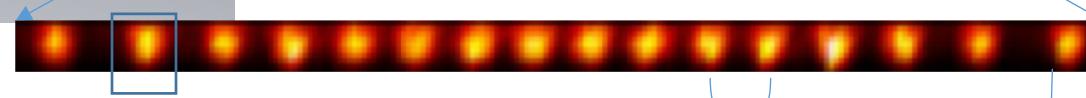
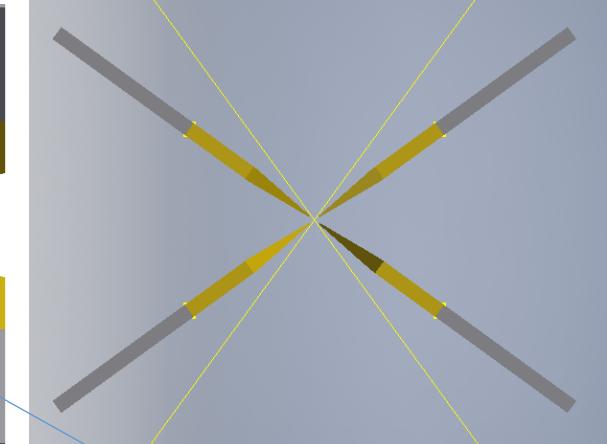
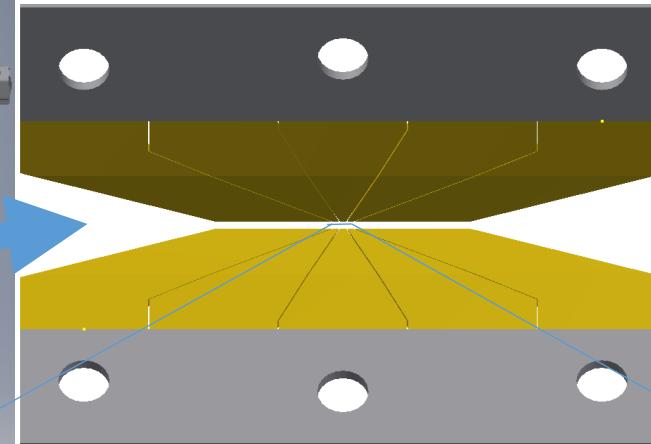
- Hybrid quantum simulation protocol to simulate 2D lattices for the case of 6, 9, 16 ions
- This protocol is scales linearly in number of pulses and stark shift gradient
- Can be used to simulate different geometries of spin graphs
- Explore additional control by dynamically changing Stark gradients

Part -2

Laser frequency stabilization



$^{171}\text{Yb}^+$ Ion Quantum Simulator



CCD image of ions
University of Maryland/JQI
ion group

$^2S_{1/2}$

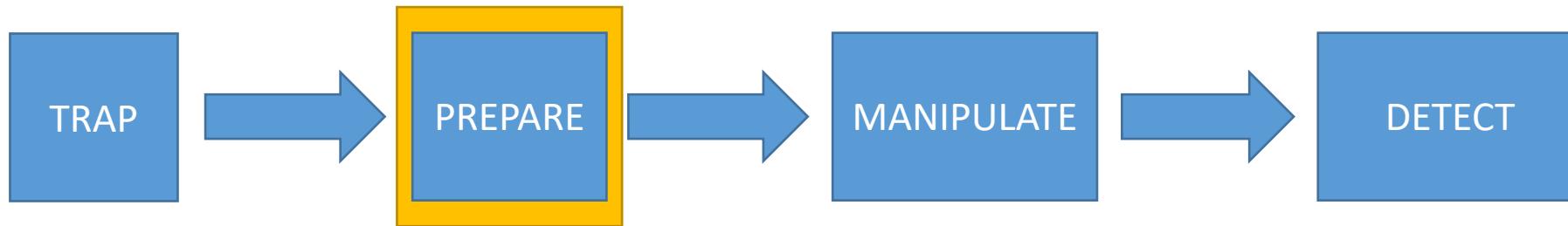
↑ —●— $|F = 1, m_F = 0\rangle$
 ↓ —●— $|F = 0, m_F = 0\rangle$

$^{171}\text{Yb}^+$ Qubit

Long coherence times > 10 min

Ye Wang et al *Nature Photonics* **11**, 646–650 (2017)

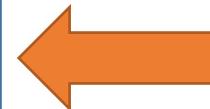
Trapped ion Quantum Simulator



- Cool the ions to motional ground state using Laser cooling
- Optically pump ions to ground state

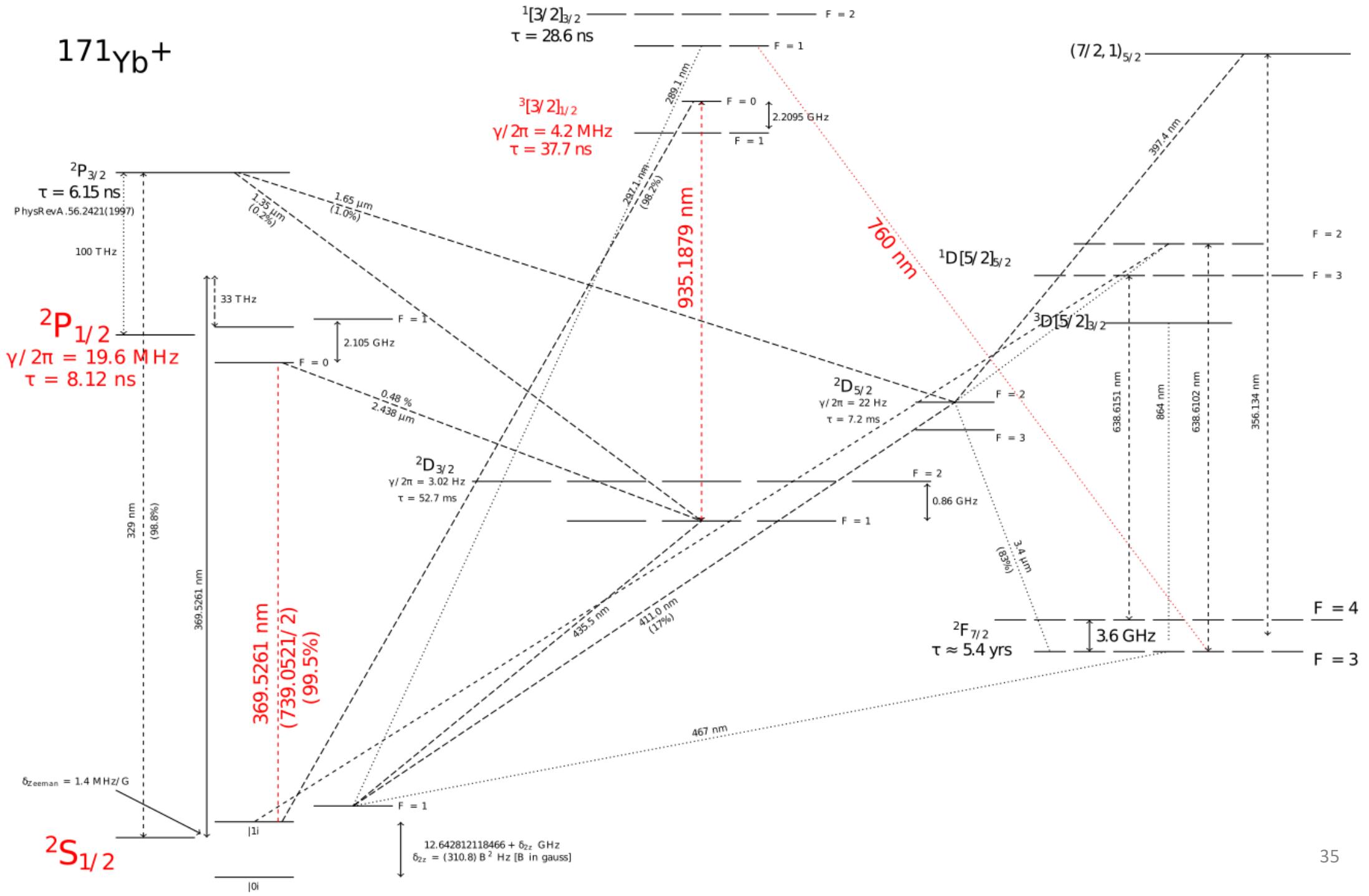
Challenge

- Frequency stabilize multiple lasers



Time

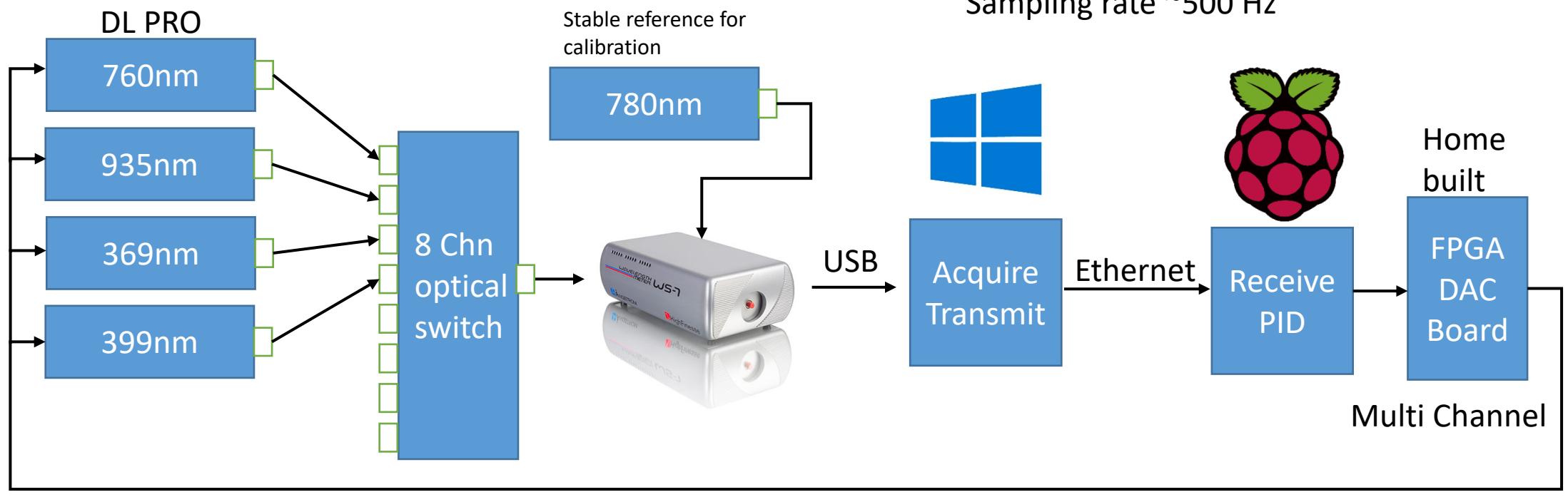
$^{171}\text{Yb}^+$



Frequency stabilization of lasers

- Stabilize lasers against drift due to ambient thermal and pressure fluctuations
- Line width of transitions order of few MHz
- No good atomic frequency reference for ions
- Solutions:
 - Laser lock based on Commercial Wavemeter
 - Lock based on in-house scanning Fabry Perot cavity

Wavemeter lock - scheme



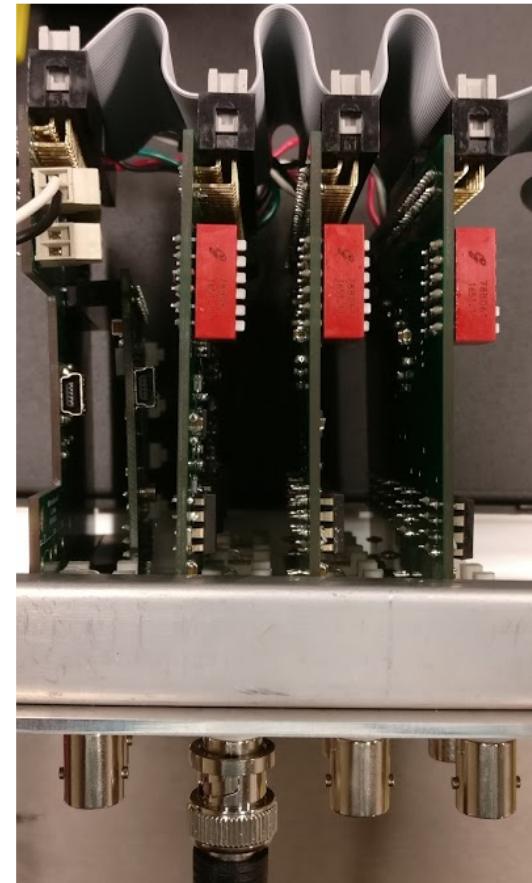
Analog Voltage 0-10V

Analog output board

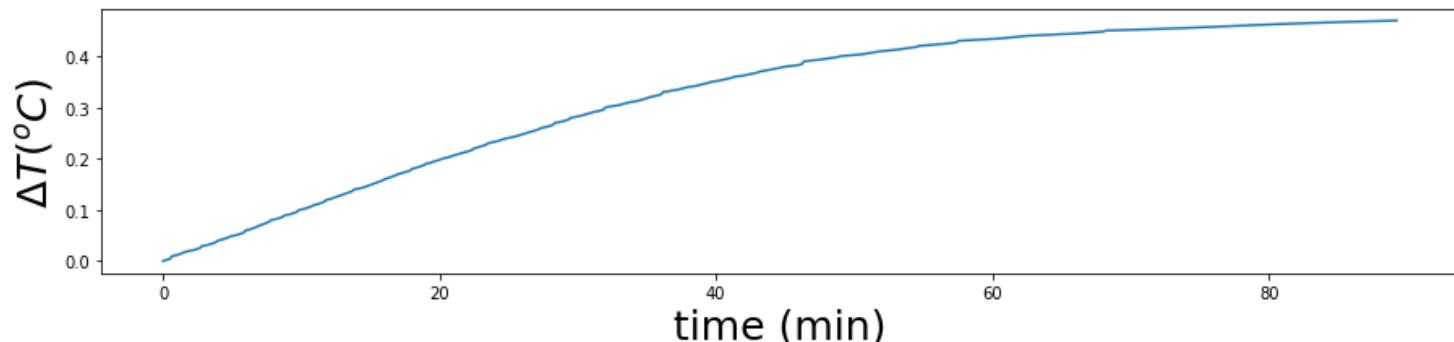
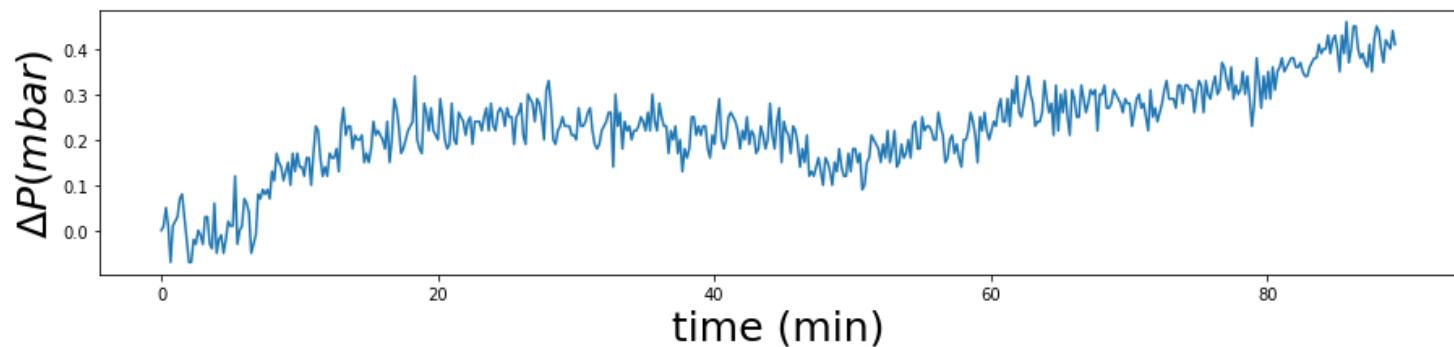
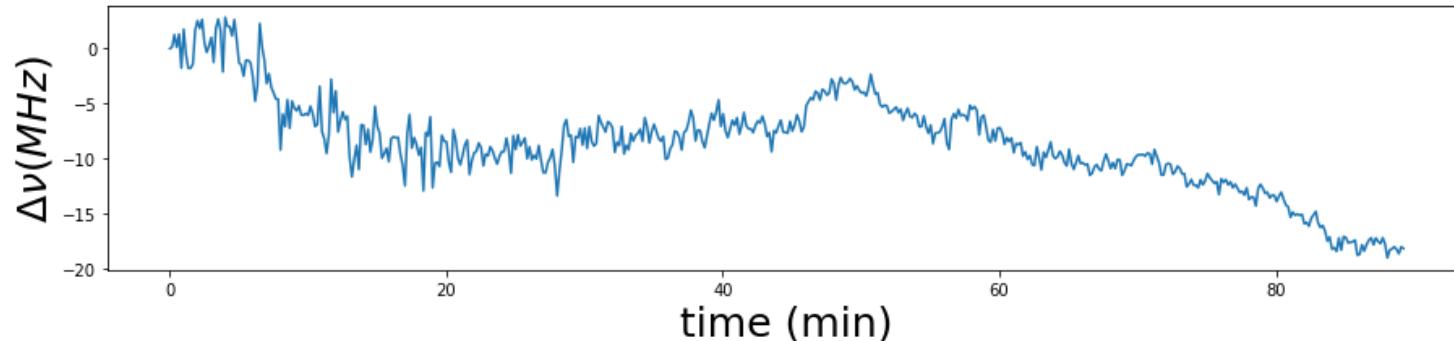
DAC7744: 4 channel 100 ksps 16bit

Numato Mimas FPGA board: Xilinx Spartan 6 LX9

Each board has 4 Analog output channels and has unique 6 bit address set by a DIP switch

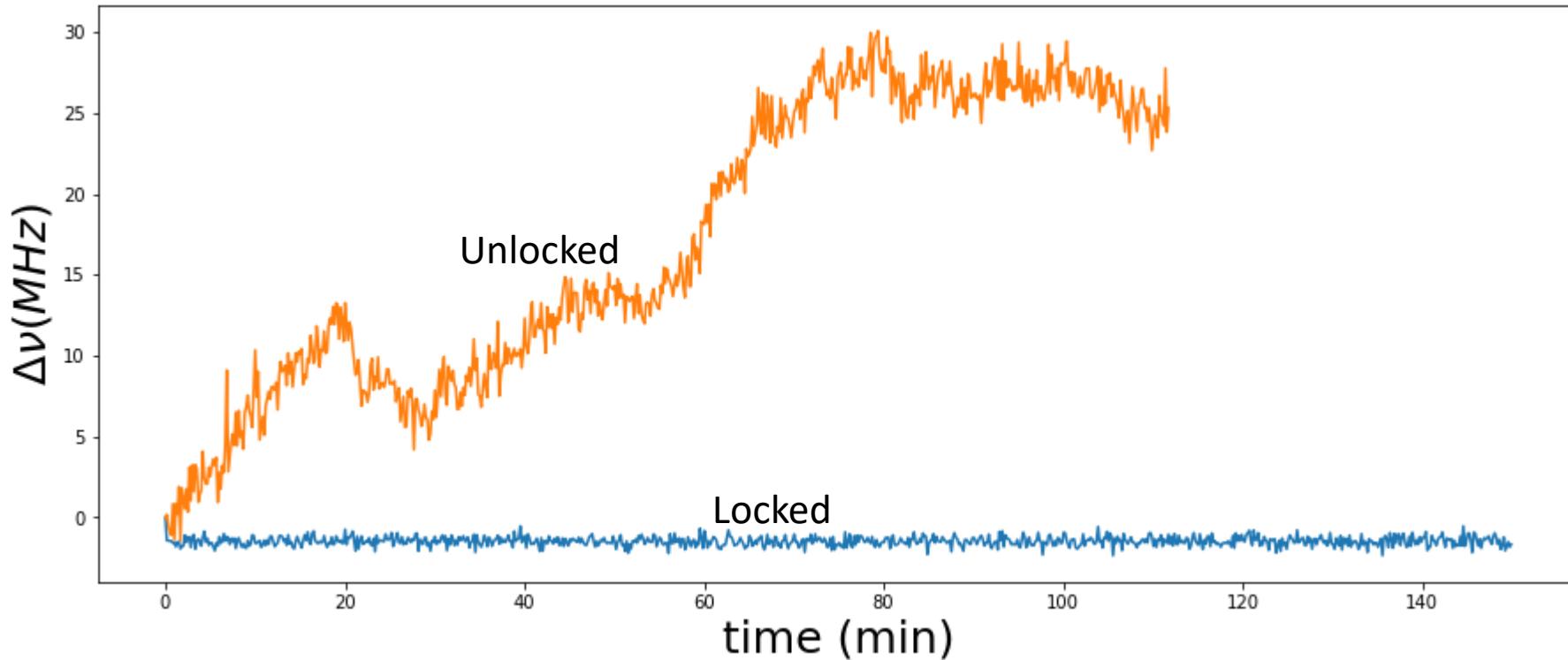


Measurements



Pressure and temperature
measured from wavemeter

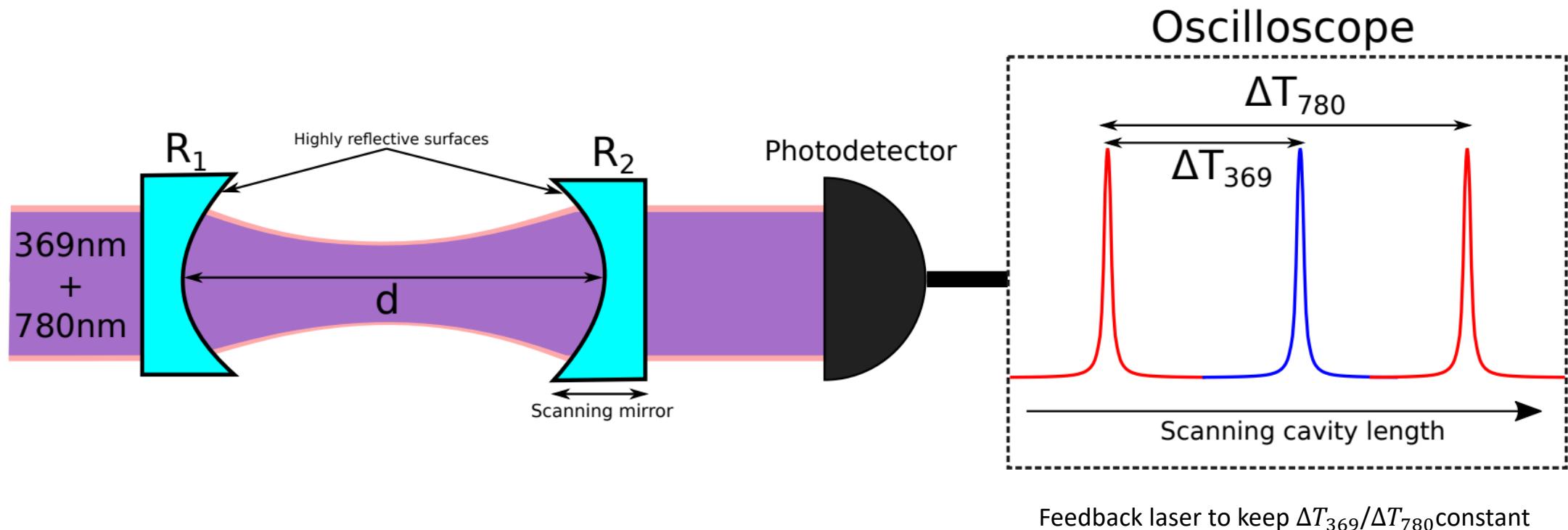
Measurements



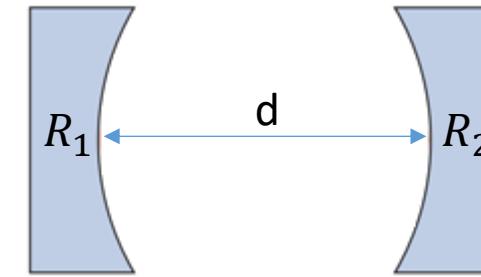
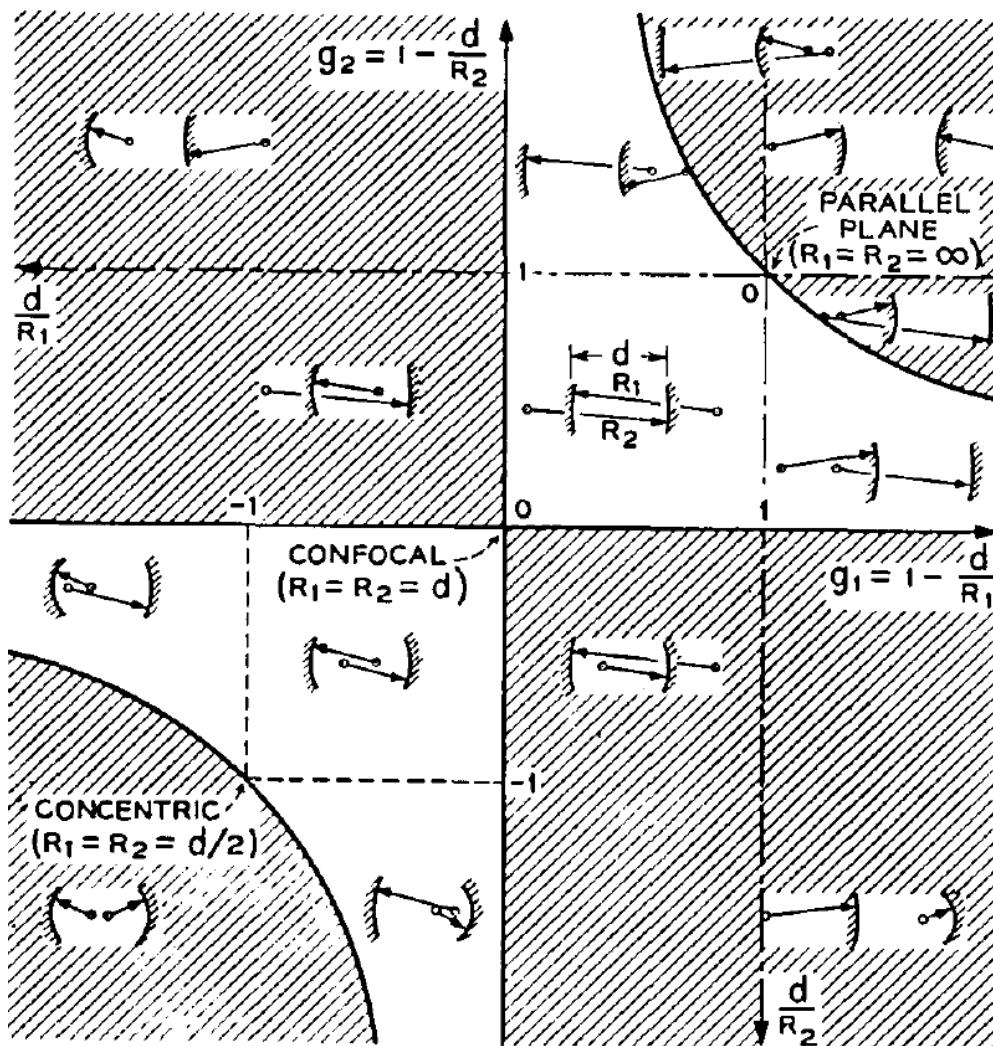
DLPRO 935nm laser

Limited by the resolution,
accuracy and response time
Of the wavemeter

Cavity based frequency lock



Stability condition for the cavity



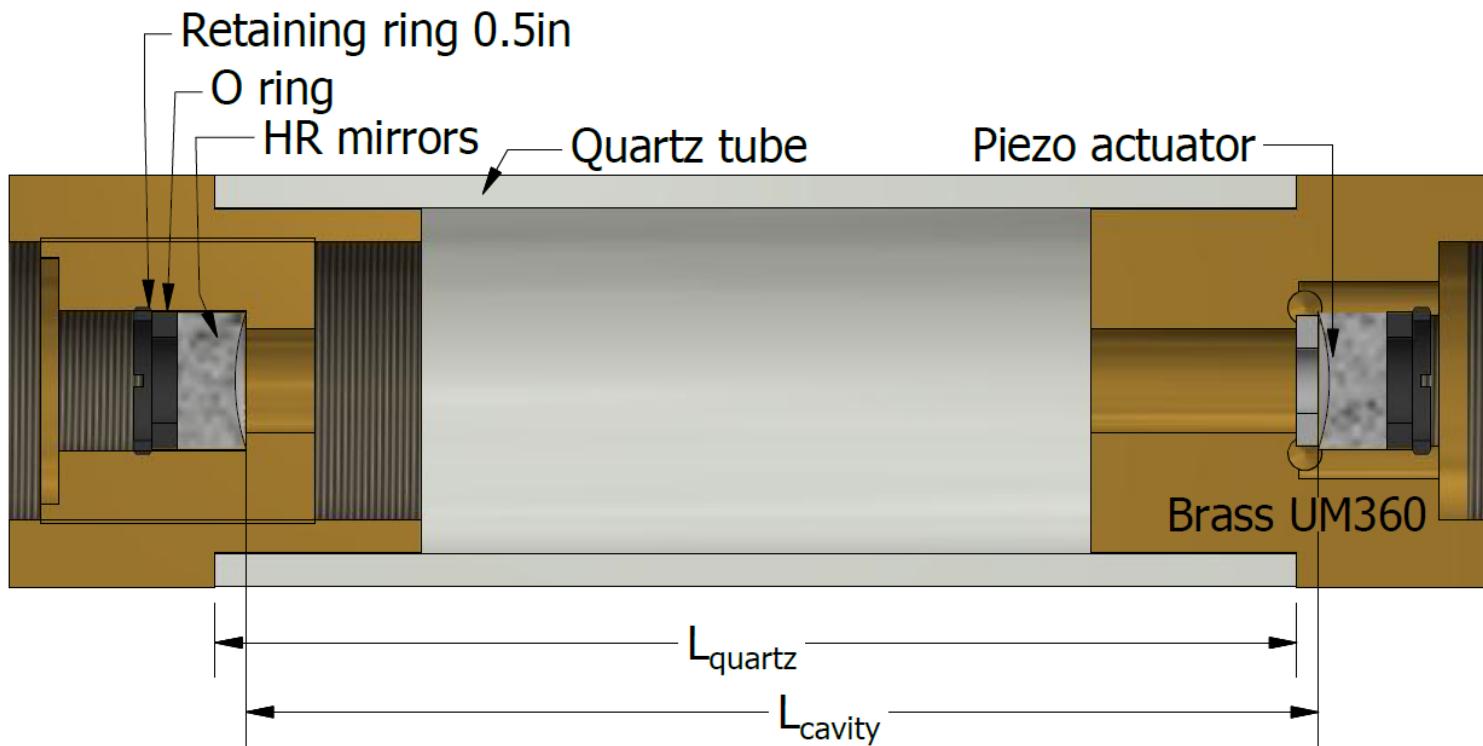
$$R_1 = R_2 = 150\text{mm}$$

$$\text{Define } g_i = \left(1 - \frac{d}{R_i}\right) i = 1,2$$

Cavity stability condition: $0 \leq g_1 * g_2 \leq 1$

We set $d \sim 100\text{ mm}$ $FSR \sim 1.5\text{GHz}$

Cavity construction



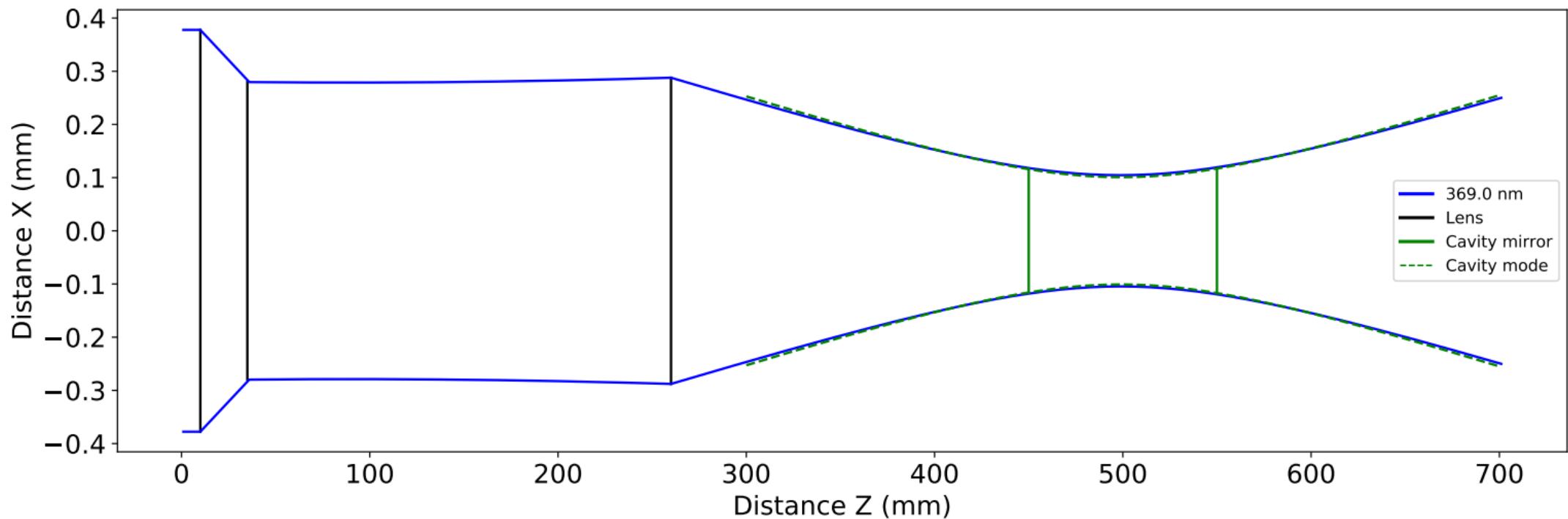
Concave mirrors ROC=150 mm
HR Mirrors R>99.5% @ 369nm - 925nm

Expected Finesse > 600

Endcaps: Brass (360 UM) (CTE: ~20ppm/ $^{\circ}\text{C}$)
Spacer : Quartz (CTE: ~0.5ppm/ $^{\circ}\text{C}$)

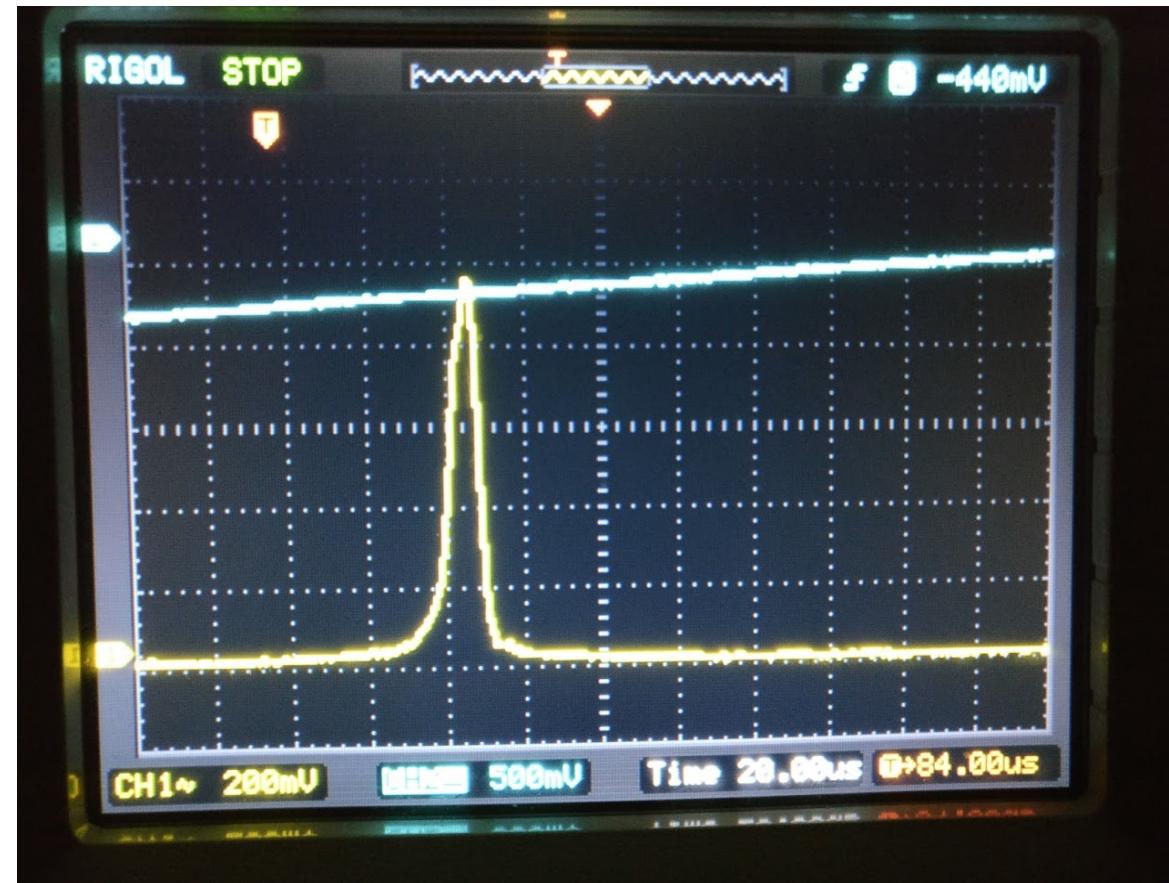
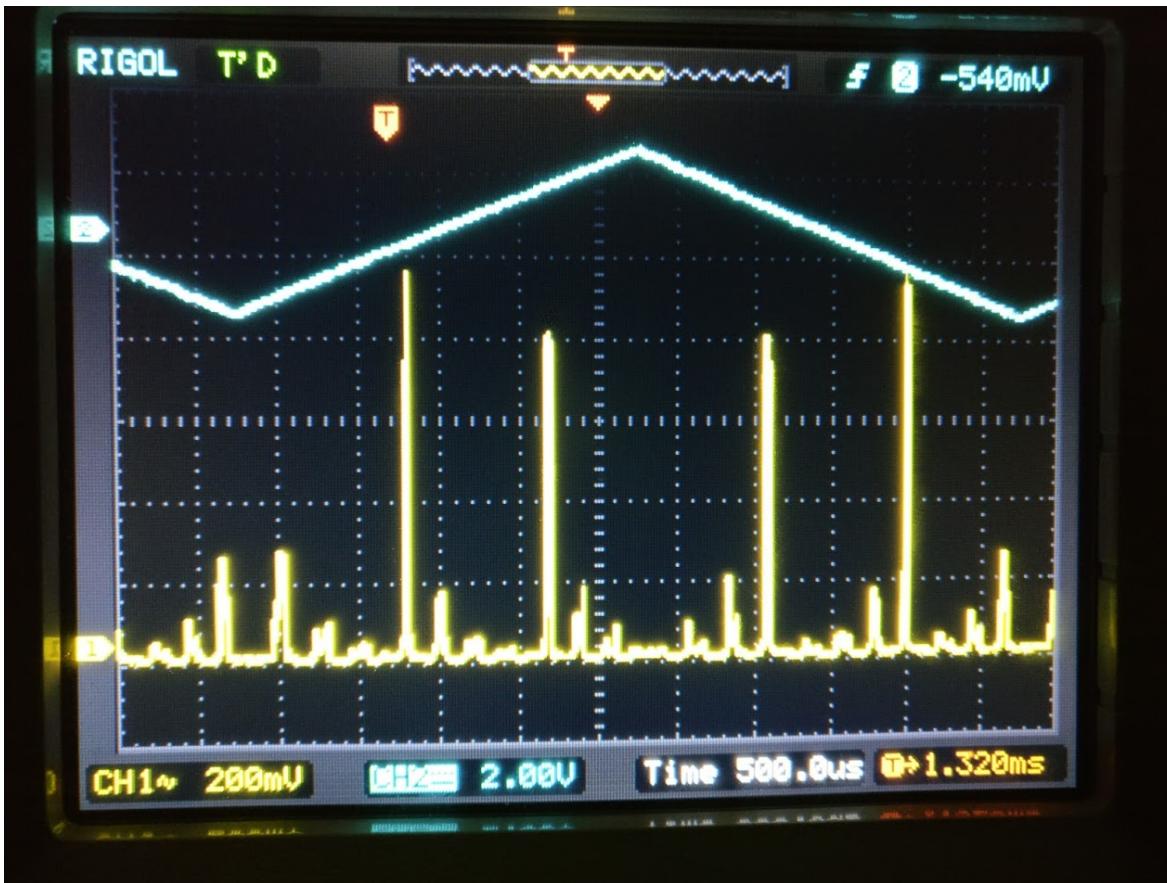
Dimensions are chosen to minimize thermal drifts

Cavity mode matching



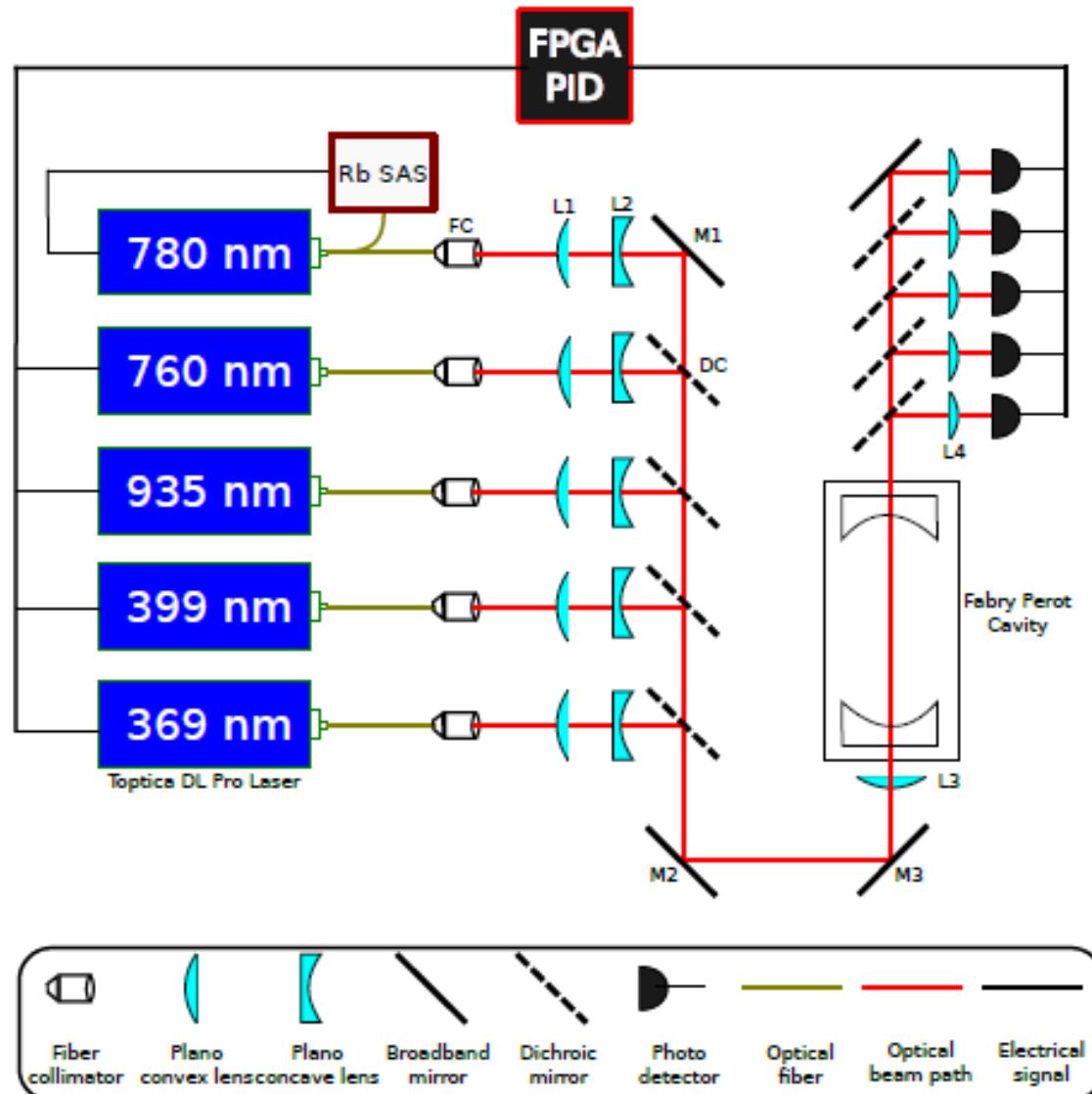
wavelength(nm)	f_{FC} (mm)	f_1 (mm)	f_2 (mm)	Ld(mm)	f_3
369.0		4.02	150.0	-75.0	75.0 150.0
399.0		4.02	150.0	-75.0	75.0 150.0
760.0		8.00	125.0	-50.0	74.0 150.0
780.0		8.00	125.0	-50.0	74.0 150.0
935.0		8.00	150.0	-50.0	99.0 150.0

Cavity measurements



Lorentzian finesse ~ 400

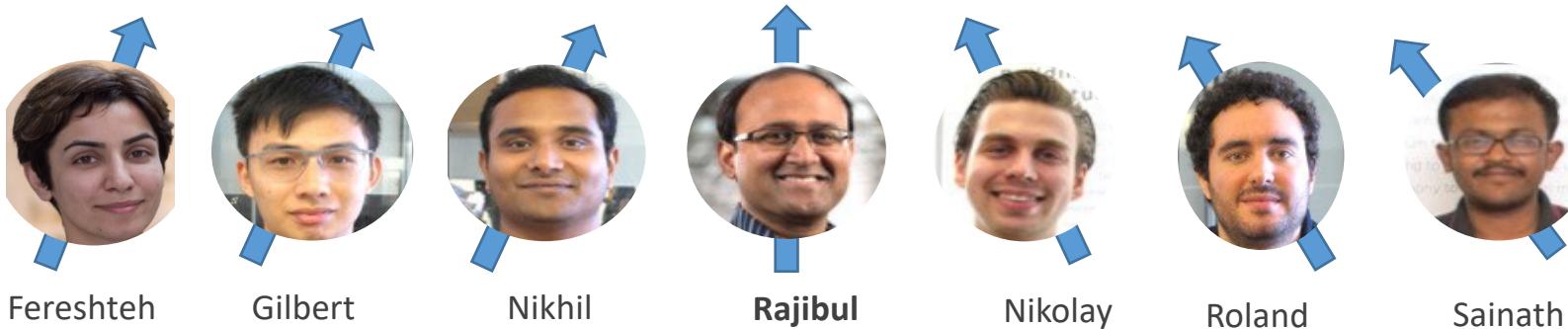
Single transfer cavity to lock multiple lasers



Summary and Outlook

- Alternate schemes for laser locking using a wavemeter and a transfer cavity
- Wavemeter lock has advantages of ease and scaling, but is limited by the resolution, accuracy and the sampling rate the instrument
- Cavity based lock has more resolution and sampling rate but cannot make absolute frequency measurements
- Combine both the schemes to obtain the advantages of both the systems.

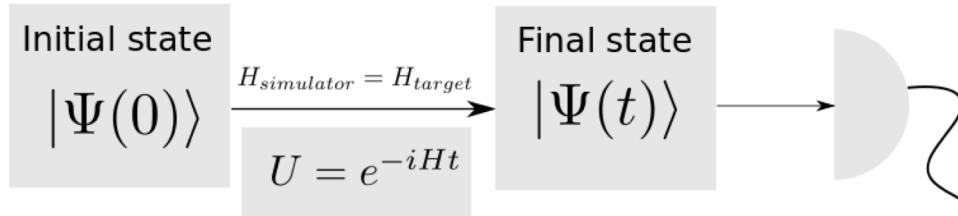
References



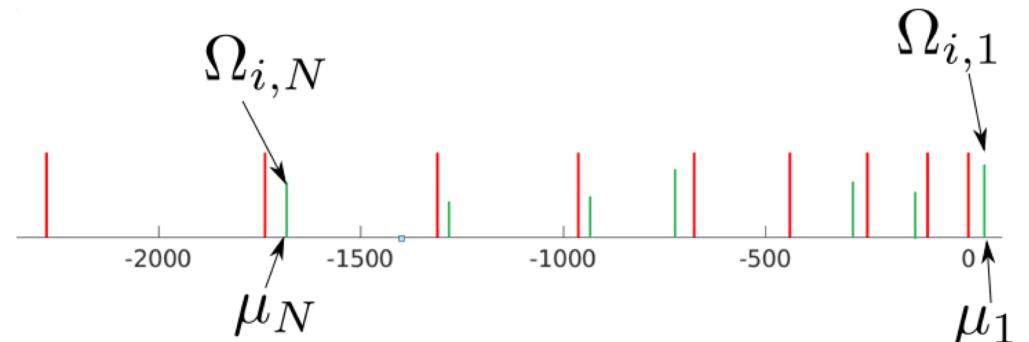
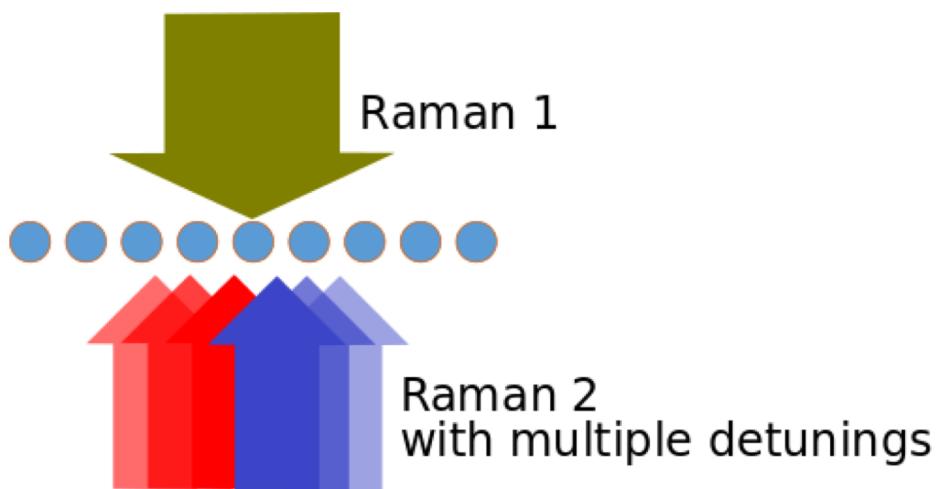
THANK YOU

Extra slides

Analog Quantum Simulation



Multi frequency extension of
Mølmer-Sørensen scheme

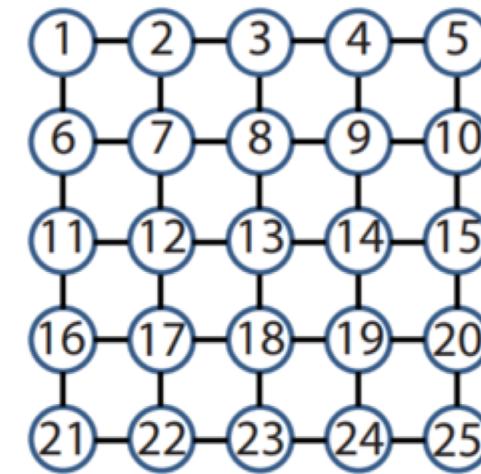
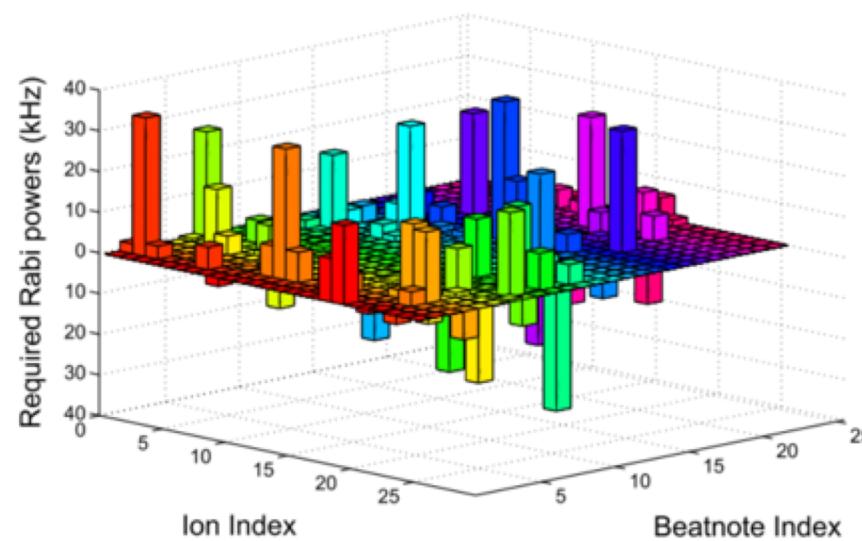


$$J_{i,j} = \sum_l \Omega_{i,l} \Omega_{j,l} \left(\frac{\hbar \Delta k^2}{2m} \right) \sum_k \frac{b_i^k b_j^k}{\mu_l^2 - \omega_k^2}$$

sum over various
Mølmer-Sørensen detunings

Optimize the values of $\{\mu_l, \Omega_{i,l}\}$ to create arbitrary J_{ij}

One such optimized solution to simulate 2D lattices*



Requires individual ion addressability with various Mølmer-Sørensen beams

Experimental parameters

