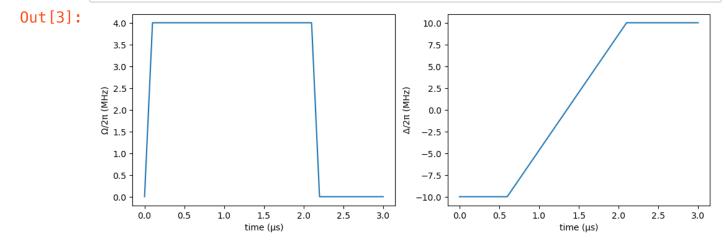
7/10/22, 8:13 PM task1 - Jupyter Notebook

```
In [2]: using Blogade
        using PythonCall
        using KrylovKit
        using SparseArrays
        plt = pyimport("matplotlib.pyplot");
```

Define time-dependent Rabi frequency, $\Omega(t)$, and detuning parameter, $\Delta(t)$

```
In [3]: |t_tot = 3.0;
          \Omega max = 2\pi * 4;
          \Omega = piecewise_linear(clocks = [0.0, 0.1, 2.1, 2.2, t_tot], values = [0.0, 0.1, 2.1, 2.2, t_tot]
          U = 2\pi * 10;
          \Delta = piecewise_linear(clocks = [0.0, 0.6, 2.1, t_tot], values = [-U, -U]
          fig, (ax1, ax2) = plt.subplots(ncols = 2, figsize = (12, 4))
          Blogade plot! (ax1, \Omega)
          ax1.set_ylabel("\Omega/2\pi (MHz)")
          Bloqade.plot!(ax2, \Delta)
          ax2.set_ylabel("\Delta/2\pi (MHz)")
          fiq
```



Create 1D chain with 9 atom sites

```
In [4]: nsites = 9
        atoms = generate_sites(ChainLattice(), nsites, scale = 5.48)
Out [4]:
```

0.0µm

Define time-dependent Rydberg Hamiltonian

```
In [5]: h = rydberg_h(atoms; \Delta, \Omega)

Out[5]: \sum \frac{2\pi \cdot 0.863 \times 10^{6.0}}{|r_i - r_j|^6} n_i n_j + 1 \cdot \Omega(t) \cdot \sum \sigma_i^x - \Delta(t) \cdot \sum n_i
```

Set all atoms to ground state in initial state and set up time-evolution problem

```
In [17]: reg = zero_state(9);
prob = SchrodingerProblem(reg, t_tot, h);
integrator = init(prob, Vern8());
```

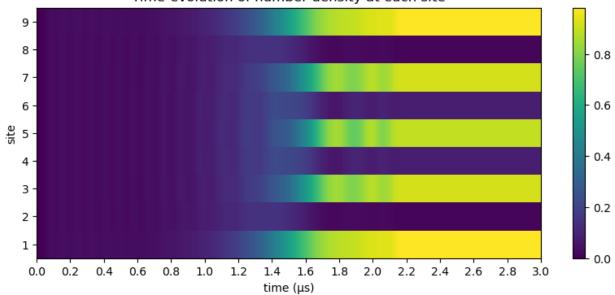
Calculate the site number densities, $\langle n_i(t) \rangle$, as the initial state adiabatically time-evolves

```
In [18]: densities = []
    for _ in TimeChoiceIterator(integrator, 0.0:1e-2:t_tot)
        push!(densities, rydberg_density(reg))
    end
    D = hcat(densities...);

    fig, ax = plt.subplots(figsize = (10, 4))
    shw = ax.imshow(real(D), interpolation = "nearest", aspect = "auto", e
    ax.set_title("Time-evolution of number density at each site")
    ax.set_xlabel("time (µs)")
    ax.set_ylabel("site")
    ax.set_yticks(0:0.2:t_tot)
    ax.set_yticks(1:nsites)
    bar = fig.colorbar(shw)
    fig
```

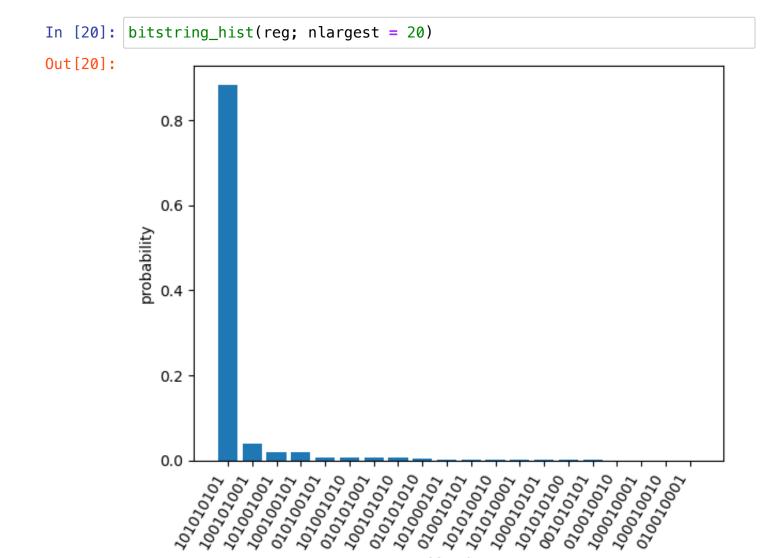
Out[18]:





In [19]: measure(reg)[]

Out[19]: 101010101 (2)

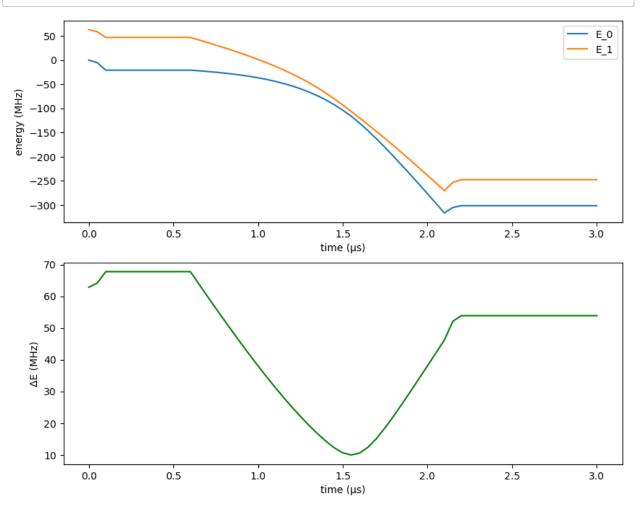


bitstring

Energy of ground and first excited states

```
In [9]: E0 = []; E1 = [];
        tseq = collect(0.0:0.05:t_tot)
        for t in tseq
            ht = h |> attime(t)
            h_m = mat(ht)
            vals, vecs, info = KrylovKit.eigsolve(h_m, 5, :SR)
            push!(E0, vals[1]);push!(E1, vals[2]);
        end
        fig, (ax1, ax2) = plt.subplots(nrows = 2, figsize = (10, 8))
        ax1.plot(tseq, E0, label = "E_0")
        ax1.plot(tseq, E1, label = "E_1")
        ax1.set_xlabel("time (µs)")
        ax1.set_ylabel("energy (MHz)")
        ax1.legend()
        ax2.plot(tseq, E1 - E0, 'g')
        ax2.set_xlabel("time (µs)")
        ax2.set_ylabel("ΔE (MHz)")
        fig
```

Out [9]:



```
In [11]: minimum(E1 - E0)
Out [11]: 9.95423632964821
          Expectation value, \langle \sigma_i^x \rangle, at each site
In [23]: measure(reg)[]
Out [23]: 101010101 (2)
In [22]: println("Expectation value of X at site")
          for i in 1:9
              println(i, ": ", expect(put(9,i=>X), reg))
          end
          Expectation value of X at site
          1: -0.0009154741016707858 + 0.0im
          2: -0.002741274442991784 + 0.0im
          3: 0.01565263209333362 + 0.0im
          4: -0.002034021969768495 - 1.0842021724855044e-19im
          5: 0.02624159404006243 + 0.0im
          6: -0.0020340219697753278 - 2.168404344971009e-19im
          7: 0.0156526320931435 + 6.938893903907228e-18im
          8: -0.0027412744429678765 + 4.336808689942018e-19im
          9: -0.0009154741018149708 + 0.0im
 In [ ]:
```