```
In [ ]:
        import cirq
        import numpy as np
         import matplotlib.pyplot as plt
         # import cirq_google
        from numpy import linalg as LA
        from math import log, e
        # print the supremacy chip
        #print(cirq_google.Sycamore)
In [ ]: # These are three qubits on a line
        qubits = cirq.LineQubit.range(3)
        a = qubits[0]
         b = qubits[1]
        c = qubits[2]
        # This is a collection of operations
        # Each operation is a gate
        ops = [cirq.H(a), cirq.H(b), cirq.CNOT(b, c), cirq.H(b)]
        circuit = cirq.Circuit(ops)
        # print circuit diagram
        print(circuit)
In [ ]: def Bell(n):
             q_chain = cirq.LineQubit.range(2)
            bell = cirq.Circuit()
             bell.append(cirq.H(q chain[0]))
            bell.append(cirq.CNOT(q_chain[0],q_chain[1]))
             if n == 1:
                 return bell
            elif n==2:
                 bell.append(cirq.Z(q_chain[1]))
                 return bell
             elif n==3:
                 bell.append(cirq.X(q_chain[1]))
                 return bell
             elif n==4:
                 bell.append(cirq.X(q chain[1]))
                 bell.append(cirq.Z(q_chain[0]))
                 return bell
         bell = Bell(4)
         print(bell) #1:
        simulator = cirq.Simulator()
         result = simulator.simulate(bell)
         print('Bra-ket notation for the wavefunction:')
        print(result.dirac_notation())
```

Exercise 4

```
In []: import random
    from math import pi

    random.seed(41)

# qubit order:
    # Message
# Alice
# Bob
```

```
# creating basics of circuit
        qubits = 3
        q_chain = cirq.LineQubit.range(qubits)
        preparation = cirq.Circuit()
        teleportation = cirq.Circuit()
        message = q_chain[0]
        alice = q_{chain}[1]
        bob = q_chain[2]
        # preparing the random state on qubit 0 and the Bell state on qubit 2 and 3
        t = pi*random.uniform(0, 1)
        gate = cirq.ry(t)
        preparation.append(gate(message))
        teleportation.append(gate(message))
        teleportation.append(cirq.H(alice))
        teleportation.append(cirq.CNOT(alice,bob))
        # teleportation measurement and correction
        teleportation.append(cirq.CNOT(message,alice))
        teleportation.append(cirq.H(message))
        teleportation.append(cirq.measure(message,alice))
        teleportation.append(cirq.CZ(message,bob))
        teleportation.append(cirq.CNOT(alice,bob))
        print('\n', teleportation, '\n')
        simulator = cirq.Simulator()
        trials = 4
        for i in range(trials):
                 preparation_result = simulator.simulate(preparation)
                teleportation_result = simulator.simulate(teleportation)
                 print('First qubit state after preparation \t System state after teleportation')
                 print(preparation result.dirac notation(), '\t\t', teleportation result.dirac notat
In [ ]:
        random.seed(42)
        q_chain = cirq.LineQubit.range(1)
        message_test = cirq.Circuit()
        t = pi*random.uniform(0, 1)
        gate = cirq.ry(t)
        message_test.append(gate(message))
        print(message_test)
        simulator = cirq.Simulator()
        result = simulator.simulate(message test)
        print('Bra-ket notation for the wavefunction:')
        print(result.dirac_notation())
```

Exercise 6:

```
import matplotlib.pyplot as plt
def Population(w,wo,w1,t):
        Omega = sqrt(w1**2 + (w-wo)**2)
        population = (w1*sin(Omega*t/2)/Omega)**2
        return population
def U(w,wo,w1,t,dt):
        q_chain = cirq.LineQubit.range(1)
        troter = cirq.Circuit()
        N = int(t/dt)
        thetaz = -wo*dt
        thetax = 2*w1*dt
        for n in range(N):
                rx = cirq.rx(thetax*cos(w*n*dt))
                rz = cirq.rz(thetaz)
                troter.append(rz(q chain[0]))
                troter.append(rx(q chain[0]))
        return troter, q_chain
def Troterization(w,wo,w1,t,dt, mode='simulate', noise=''):
        reps = 10000
        troter, q_chain = U(w,wo,w1,t,dt)
        simulator = cirq.Simulator()
        if mode == 'tomography':
                sampler = cirq.DensityMatrixSimulator()
                result = cirq.experiments.state_tomography(sampler=sampler, qubits=q_chain, c
                rho = result._density_matrix
                population = rho[1,1].real
                return population
        if mode == 'run':
                troter.append(cirq.measure(q_chain[0], key='0'))
                result = simulator.run(troter, repetitions=reps)
                counts = result.histogram(key = '0')
                # population for state |1> := prob(measurement=1)
                population = counts[1]/reps
                return population
        elif mode == 'simulate':
                result = simulator.simulate(troter)
                #print('Bra-ket notation for the wavefunction:')
                #print(result.dirac_notation())
                #print('Population of state |1>')
                population = abs(result.final_state_vector[-1])**2
                #print(np.around(population, 3))
                return population
        elif mode == 'noise':
                troter.append(cirq.measure(q_chain[0], key='0'))
                result = cirq.sample(program=troter, noise=noise, repetitions=reps) # type(re
                # type(histogram) = <class 'collections.Counter'>
                counts = result.histogram(key = '0')
                # population for state |1> := prob(measurement=1)
                population = counts[1]/reps
                return population
        else:
                raise Exception('Not a valid mode: {}'.format(mode))
```

```
In [ ]: # ts = np.linspace(0.2,4,100)
        t = 4
        dt = 0.05
        ts = np.linspace(0, t, int(t/dt))
        pop_a = []
        theo_a = []
        for t in ts:
            W, W0, W1 = 25.5, 25, 2
            theo_a.append(Population(w,wo,w1,t))
             pop_a.append(Troterization(w,wo,w1,t,dt,'run'))
        ws = np.arange(10, 40, 0.2)
        w0, w1 = 25, 2
        delta = np.arange(10-w0, 40-w0, 0.2)
         pop_b = []
        theo_b = []
        for w in ws:
            t = pi/w1
            theo b.append(Population(w,wo,w1,t))
            pop_b.append(Troterization(w,wo,w1,t,dt,'run'))
        t = 4
        dt = 0.05
        timeIntervals2 = np.linspace(0, t, int(t/dt))
        pop_c = []
        theo_c = []
        W, W0, W1 = 2, 2, 2
        for t in timeIntervals2:
            theo_c.append(Population(w,w0,w1,t))
             pop_c.append(Troterization(w,w0,w1,t,dt,'run'))
        t = 1.4
        dt = 0.05
        timeIntervals3 = np.linspace(0, t, int(t/dt))
        pop_d = []
        theo_d = []
        for t in timeIntervals3:
            W, W0, W1 = 3, 1, 6
            theo_d.append(Population(w,w0,w1,t))
             pop_d.append(Troterization(w,w0,w1,t,dt,'run'))
In [ ]: plt.figure()
```

```
plt.xlabel(r'Time t')
plt.ylabel(r'Population $|\beta(t)|^2$')
plt.plot(ts, theo_a, 'gray', label='Analytical')
plt.plot(ts, pop_a, 'green', label='Trotterization')
plt.ylim(-0.02, 1.02)
plt.legend()
plt.savefig("exercise06_01.png")
plt.show()
plt.figure()
plt.xlabel(r'Detuning $\Delta$')
plt.ylabel(r'Population $|\beta(t)|^2$')
plt.plot(delta, theo_b, 'gray', label='Analytical')
plt.plot(delta, pop_b, 'royalblue', label='Trotterization')
plt.ylim(-0.02, 1.02)
plt.legend()
plt.savefig("exercise06_02.png")
plt.show()
plt.figure()
plt.xlabel(r'Time t')
plt.ylabel(r'Population $|\beta(t)|^2$')
plt.plot(timeIntervals2, theo_c, 'gray', label='Analytical')
```

```
plt.plot(timeIntervals2, pop_c, 'darkred', label='Trotterization')
plt.ylim(-0.02, 1.02)
plt.legend()
plt.savefig("exercise06_03.png")
plt.show()

plt.figure()
plt.xlabel(r'Time t')
plt.ylabel(r'Population $\beta(t)\|^2$')
plt.plot(timeIntervals3, theo_d, 'gray', label='Analytical')
plt.plot(timeIntervals3, pop_d, 'darkred', label='Trotterization')
plt.ylim(-0.02, 1.02)
plt.legend()
plt.savefig("exercise06_04.png")
plt.show()
```

Exercise 7

```
In [ ]: dt = 0.05
        t = 7
        W, W0, W1 = 25.5, 25, 2
        p = 0.9
        gamma = 0.02
        pop_a = []
        pop_noise = []
        noise = cirq.ConstantQubitNoiseModel(cirq.GeneralizedAmplitudeDampingChannel(p=p,gamma=gamma)
        ts new = np.linspace(0, t, int(t/dt))
        for t in ts new:
            pop noise.append(Troterization(w,wo,w1,t,dt,'noise',noise))
        for t in ts_new:
            theo_a.append(Population(w,wo,w1,t))
            pop_a.append(Troterization(w,wo,w1,t,dt,'run'))
        plt.figure()
In [ ]:
        plt.xlabel('Time t')
        plt.ylabel(r'Population $|\beta(t)|^2$')
        plt.plot(ts_new, pop_a, 'gray', label='no noise')
        plt.plot(ts_new, pop_noise, 'darkviolet', label='GAD')
        plt.ylim(-0.02, 1.02)
        plt.legend()
        plt.grid(linestyle='-', linewidth=0.2)
        plt.savefig("exercise07_01.pdf")
        plt.show()
In []: W, wo, w1 = 25.5, 25, 2
        ps = [1,0]
        gammas = [0.02]
        dt = 0.05
        t = 10
        ts_new = np.linspace(0, t, int(t/dt))
        populations1 = np.zeros((len(ps),len(ts_new),len(gammas)))
        for i, p in enumerate(ps):
            for j, gamma in enumerate(gammas):
                for k, t in enumerate(ts_new):
                     noise = cirq.ConstantQubitNoiseModel(cirq.GeneralizedAmplitudeDampingChannel(p=p,
                     populations1[i,k,j] = (Troterization(w,wo,w1,t,dt,'noise',noise))
        ps = [0.5]
```

```
dt = 0.05
        t = 10
        ts_new_2 = np.linspace(0, t, int(t/dt))
        populations2 = np.zeros((len(ps),len(ts_new_2),len(gammas)))
        for i, p in enumerate(ps):
            for j, gamma in enumerate(gammas):
                for k, t in enumerate(ts_new_2):
                     noise = cirq.ConstantQubitNoiseModel(cirq.GeneralizedAmplitudeDampingChannel(p=p,
                     populations2[i,k,j] = (Troterization(w,wo,w1,t,dt,'noise',noise))
        ps = [1,0]
        gammas = [0.02]
        w, wo, w1 = 25.5, 25, 0
        dt = 0.4
        t = 40
        ts new 3 = np.linspace(0, t, int(t/dt))
        populations3 = np.zeros((len(ps),len(ts_new_3),len(gammas)))
        for i, p in enumerate(ps):
            for j, gamma in enumerate(gammas):
                for k, t in enumerate(ts_new_3):
                     noise = cirq.ConstantQubitNoiseModel(cirq.GeneralizedAmplitudeDampingChannel(p=p,
                     populations3[i,k,j] = (Troterization(w,wo,w1,t,dt,'noise',noise))
In [ ]: plt.figure()
        plt.xlabel('Time t')
        plt.ylabel(r'Population $|\beta(t)|^2$')
         plt.plot(ts_new, populations1[0,:,0], 'violet', label='p = 1')
         plt.plot(ts_new, populations1[1,:,0], 'darkviolet', label='p = 0')
        plt.ylim(-0.02, 1.02)
        plt.legend()
        plt.grid(linestyle='-', linewidth=0.2)
         plt.savefig("exercise07_02.png")
        plt.show()
        plt.figure()
        plt.xlabel('Time t')
        plt.ylabel(r'Population $|\beta(t)|^2$')
        plt.plot(ts_new_2, populations2[0,:,0], 'darkviolet', label=r'$\gamma = 0.02$')
         plt.plot(ts_new_2, populations2[0,:,1], 'violet', label=r'$\gamma = 0.005$')
        plt.ylim(-0.02, 1.02)
        plt.legend()
        plt.grid(linestyle='-', linewidth=0.2)
        plt.savefig("exercise07_03.png")
        plt.show()
        plt.figure()
        plt.xlabel('Time t')
        plt.ylabel(r'Population $|\beta(t)|^2$')
        plt.plot(ts_new_3, populations3[0,:,0], 'violet', label='p = 1')
        plt.plot(ts_new_3, populations3[1,:,0], 'darkviolet', label='p = 0')
        plt.ylim(-0.02, 1.02)
        plt.legend()
        plt.grid(linestyle='-', linewidth=0.2)
        plt.savefig("exercise07_04.png")
        plt.show()
```

Notes:

gammas = [0.02, 0.005]

For p=1 the channel reduces to the normal amplitude damping channel \ For p=0 the channel behaves like an amplitude enhancing channel, driving the system to the |1><1| state

Exercise 8:

```
In [ ]:
        def Expectation(prepare='', ops='', reps=int(1e6), debug=False):
            prepare: gates to prepare initial state
            ops: operators to change the Z measurement basis into another basis
            examples:
             - for measurement in X basis: H

    for measurement in Y basis: Sdagger, H

            qubits = cirq.LineQubit.range(1)
            q0 = qubits[0]
             simulator = cirq.Simulator(seed=np.random.seed(2))
            measurement_op = cirq.Circuit()
             if prepare:
                 measurement_op.append(p(q0) for p in prepare)
            if ops:
                 measurement op.append(p(q0) for p in ops)
            measurement_op.append(cirq.measure(q0, key='0'))
            if debug:
                 print(measurement_op)
             result = simulator.run(measurement_op, repetitions = reps)
             counts = result.histogram(key = '0')
            print(counts)
            N0, N1 = counts[0], counts[1]
             return (NO-N1)/reps
         prepare = [cirq.H, cirq.T]
        ops_X = [cirq.H]
        ops_Y = [cirq.S^{**}-1, cirq.H]
        EZ = Expectation(prepare, debug=True)
        EX = Expectation(prepare, ops_X, debug=True)
        EY = Expectation(prepare, ops_Y, debug=True)
         print('Expectation value of Z gate: ', EZ)
         print('Expectation value of X gate: ', EX)
        print('Expectation value of Y gate: ', EY)
```

Exercise 9:

Idea: to get the state we can diagonalize the matrix, and obtain the first (and only) eigen state, that would be our initial pure state

```
print(rho_tomo)
        w_tomo, v_tomo = LA.eig(rho_tomo)
        print(abs(w_tomo[0])**2)
        v_tomo[:,0]
In [ ]:
        #circuit for state tomography
        sampler = cirq.DensityMatrixSimulator()
        qchain = cirq.LineQubit.range(1)
        q0 = qchain[0]
        ops = [cirq.H(q0), cirq.T(q0)]
        circuit9 = cirq.Circuit(ops)
        # print circuit diagram
        print(circuit9)
        simulator = cirq.Simulator()
        result = simulator.simulate(circuit9)
        print('Bra-ket notation for the wavefunction:')
        print(result.dirac_notation())
        result = cirq.experiments.state_tomography(sampler=sampler, qubits=qchain, circuit= circuit9,
        rho = result. density matrix
        print(rho)
        print(rho[0,0])
        w_function, v_function = LA.eig(rho)
        print(abs(w_function[0])**2)
        v_function[:,0]
In [ ]: # comparison:
```

Exercise 10

plt.legend()

print(LA.norm(rho_tomo - rho))

rho_tomo += c*paulis_matrices[i]

```
In []: dt = 0.05
    ts_10 = np.linspace(0.2,2,100)
    pop_10 = []
    theo_10 = []
    for t in ts_10:
        w, wo, w1 = 25.5, 25, 2
        theo_10.append(Population(w,wo,w1,t))
        pop_10.append(Troterization(w,wo,w1,t,dt,'tomography'))
In []: plt.figure()
    plt.xlabel('Time t')
    plt.ylabel(r'Population $|\beta(t)|^2$')
    plt.plot(ts_10, np.array(theo_10), 'gray', label='Analytical')
    plt.plot(ts_10, pop_10, 'mediumseagreen', label='Tomography')
    plt.ylim(-0.02, 1.02)
```

```
plt.savefig("exercise10.pdf")
plt.show()
```

Exercise 11:

```
"""Define a custom gate with a parameter."""
In [ ]:
        class RXZ(cirq.Gate):
            def __init__(self, b):
                super(RXZ, self)
                self.b = b
            def _num_qubits_(self):
                return 2
            def _unitary_(self):
                return np.array([
                     [np.cos(self.b), 0.0, 1j*np.sin(self.b), 0.0],
                     [0.0, np.cos(self.b), 0.0, -1j*np.sin(self.b)],
                     [1j*np.sin(self.b), 0.0, np.cos(self.b), 0.0],
                     [0.0, -1j*np.sin(self.b), 0.0, np.cos(self.b)]
                1)
            def circuit diagram info (self, args):
                return f"RXZ({self.b})", f"RXZ({self.b})"
        class RZX(cirq.Gate):
            def __init__(self, b):
                super(RZX, self)
                self.b = b
            def _num_qubits_(self):
                 return 2
            def _unitary_(self):
                return np.array([
                     [np.cos(self.b), 1j*np.sin(self.b), 0.0, 0.0],
                     [1j*np.sin(self.b), np.cos(self.b), 0.0, 0.0],
                     [0.0, 0.0, np.cos(self.b), -1j*np.sin(self.b)],
                     [0.0, 0.0, -1j*np.sin(self.b), np.cos(self.b)]
                1)
            def _circuit_diagram_info_(self, args):
                return f"RZX({self.b})", f"RZX({self.b})"
        def expXZ(b):
            " exp(ib(XZ)) acting on two qubits"
            q_chain = cirq.LineQubit.range(2)
            rotation = cirq.Circuit()
            rotation.append(RXZ(b=b).on(*q chain))
            return rotation
        def expZX(b):
            " exp(ib(ZX)) acting on two qubits"
            q_chain = cirq.LineQubit.range(2)
            rotation = cirq.Circuit()
            rotation.append(RZX(b=b).on(*q_chain))
            return rotation
        def ansatz(a,b):
            q_chain = cirq.LineQubit.range(2)
            circ = cirq.Circuit()
            rz = cirq.rz(-2*a)
            circ.append(expZX(b))
```

```
circ.append(expXZ(b))
            circ.append(rz(q_chain[1]))
            circ.append(rz(q_chain[0]))
            return circ
        # testing our implementation of the Ansatz circuit
        ansatz_state = ansatz(np.pi/2,np.pi/2)
        print(ansatz_state)
        result = simulator.simulate(ansatz state)
        print('Bra-ket notation for the wavefunction:')
        print(result.dirac_notation()) #this is correct
In [ ]: # Exercise 12
        alist = np.linspace(0,np.pi,20)
        blist = np.linspace(0,np.pi,20)
        energies = np.empty(shape=(len(alist),len(blist)))
        q_chain = cirq.LineQubit.range(2)
        q0 = q chain[0]
        q1 = q_{chain}[1]
        for i, a in enumerate(alist):
            for j, b in enumerate(blist):
                 energies[i,j] = simulator.simulate_expectation_values(ansatz(a,b), observables=-cirq.
In [ ]: pos = plt.imshow(energies, cmap='hot', interpolation='nearest', extent=[0,alist[-1],blist[-1]
        plt.title('Energies')
        plt.ylabel('a')
        plt.xlabel('b')
        tick_pos = [0, np.pi/4, np.pi/2, 3*np.pi/4, np.pi]
        labels = ['0', '\$\pi/4\$', '\$\pi/2\$', '\$\pi/4\$', '\$\pi']
        plt.xticks(tick_pos, labels)
        plt.yticks(tick pos, labels)
        plt.colorbar(pos)
        plt.savefig('exercise12.pdf')
        plt.show()
        print("The smallest energy is: ", np.amin(energies)) #should be -2 in theory
In [ ]:
        arg = np.argmin(energies)
        print(arg)
        bmin= arg%20
        amin= (arg)//20
        print("Indices for the arrays are: (i: {} and j: {}) ".format(amin,bmin) )
        print("Values for the indices are: (a: {} and b: {}) ".format(alist[amin],blist[bmin]))
In [ ]: print(np.where(energies == np.amin(energies))) #printing the pair of indices for which the mi
        # here this is (5,5) and (14,14)
In [ ]: test =simulator.simulate_expectation_values(ansatz(alist[14],alist[14]), observables=-cirq.Z(
        print('Energy of ground state:', round(test.real,3))
In [ ]: def entanglement_entropy(circ, simulator=cirq.Simulator()):
            result = simulator.simulate(circ)
            state = result.final_state_vector
            C = np.reshape(state, newshape=(2,2))
            u, s, vh = LA.svd(C)
```

```
for sigma in s:
                if sigma != 0:
                    ent -= sigma**2 * log(sigma**2,2)
            return ent
In [ ]:
        #13
        gs = ansatz(alist[amin],blist[bmin])
        print(gs)
        result = simulator.simulate(gs)
        state = result.final_state_vector
        print('======"')
        print('Bra-ket notation for the wavefunction:')
        print(result.dirac_notation())
        print('=======')
        #correct up to a global phase of e^(-ipi/2)
        gs_state = np.array([1,1,1,-1])*0.5
        print(gs_state)
        print('Including the global phase and taking only the real part: ')
        print((np.exp(-1j*np.pi/2)*state).real)
        print(cirq.equal_up_to_global_phase(gs_state, state, atol=1))
        entanglement_entropy(gs) #should be equal to 1 in theory
In [ ]: # Exercise 15:
        entanglement = np.empty(shape=(len(alist),len(blist)))
        simulator = cirq.Simulator()
        for i, a in enumerate(alist):
            for j, b in enumerate(blist):
                ansatz_state = ansatz(a,b)
                entanglement[i,j] = entanglement_entropy(ansatz_state, simulator)
        pos2 = plt.imshow(entanglement, cmap='hot', interpolation='nearest', extent=[0,alist[-1],blis
In [ ]:
        plt.title('Entanglement')
        plt.ylabel('a')
        plt.xlabel('b')
        tick_pos = [0, np.pi/4, np.pi/2, 3*np.pi/4, np.pi]
        labels = ['0', 'pi/4', 'pi/2', '3pi/4', 'pi']
        plt.xticks(tick_pos, labels)
        plt.yticks(tick_pos, labels)
        plt.colorbar(pos2)
        plt.savefig('exercise15.pdf')
        plt.show()
        print("The highest Entropy is: ", np.amax(entanglement))
In [ ]:
        arg = np.argmax(entanglement)
        print('at',arg)
        b = arg \% 20
        a = (arg-b)//20
        print("Indices in the arrays are: (i: {} and j: {}) ".format(a,b) )
        print("Values for the indices are: (a: {} and b: {}) ".format(alist[a],blist[b]) )
In [ ]: print(np.where(entanglement == np.amax(entanglement))) #printing the pair of indices for whic
```

ent = 0

```
In [ ]: test = ansatz(alist[14],blist[5])
    entanglement_entropy(test, simulator)
```

the two ground states encountered coincid with points where entanglement is highest

Not used:

```
In [ ]: # Test for entanglement entropy in a bell state
        q0,q1 = cirq.LineQubit.range(2)
        circ = cirq.Circuit(cirq.H(q0), cirq.CNOT(q0,q1))
        print(circ)
        result = simulator.simulate(circ)
        state = result.final state vector
        C = np.reshape(state, newshape=(2,2))
        print(C)
        u, s, vh = LA.svd(C)
        ent = 0
        for sigma in s:
            if sigma != 0:
                ent -= sigma**2 * log(sigma**2,2)
        ent #should be equal to 1 by definition
        #entanglement= cirq.von_neumann_entropy(density)
        #print(entanglement)
```