

Homework 3

Kevin Yang - 50244152

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Link to repo:

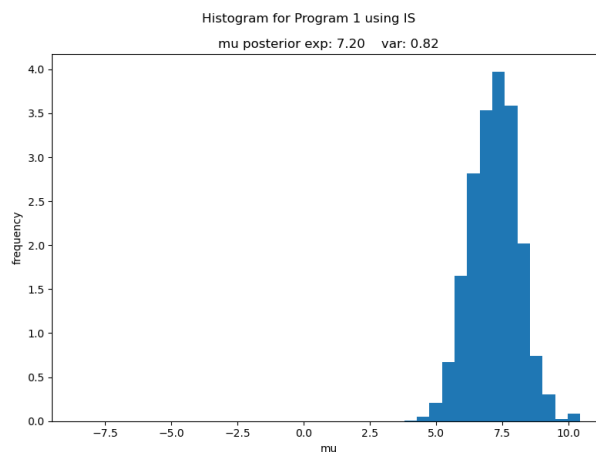
<https://github.com/keviny2/CPSC532W-Assignments/tree/main/FOPPL/HW>

1 Program 1

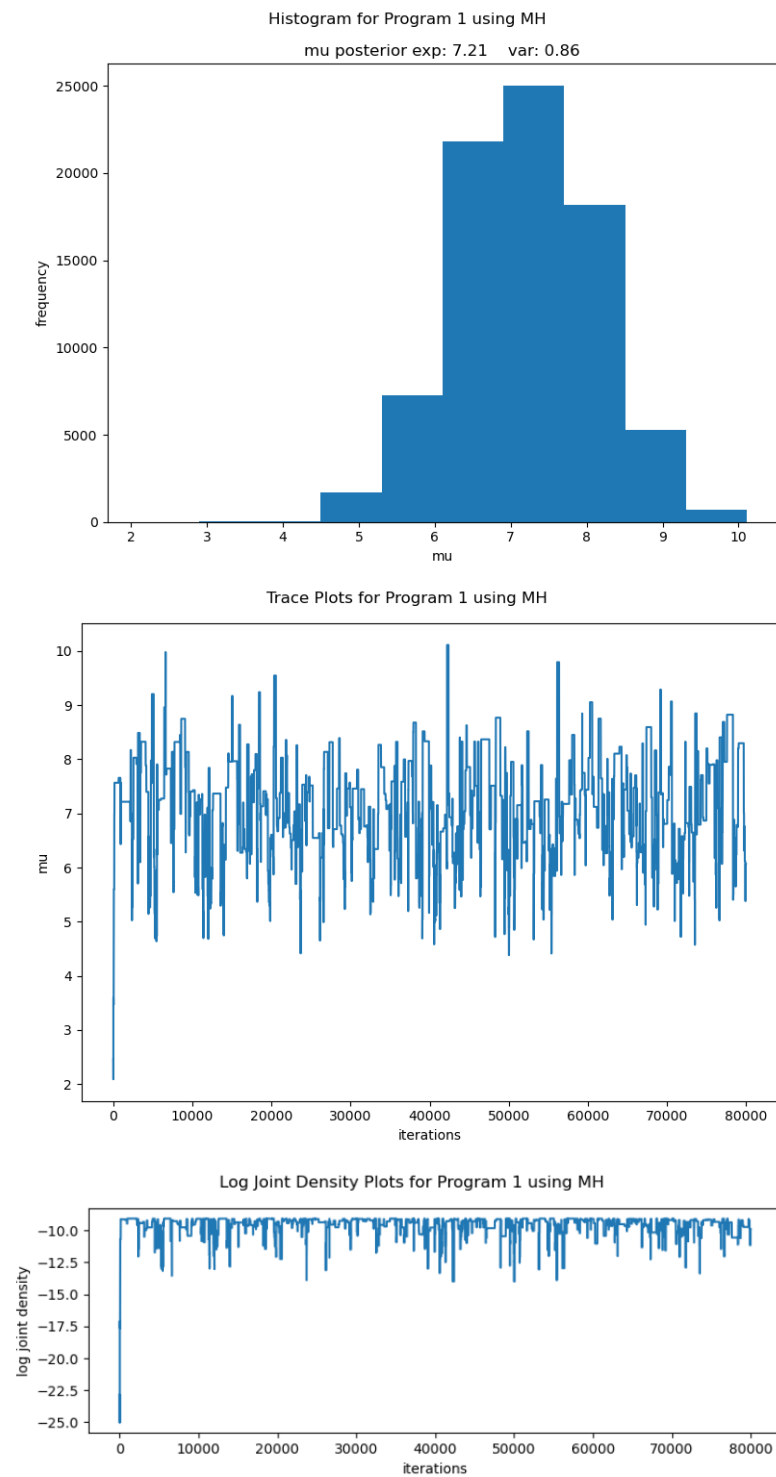
1.1 Wall-Clock Time

```
===== Likelihood Weighting =====
Took 71.83 seconds to finish Program 1
Posterior Expectation mu: tensor(7.1993)
Posterior Variance mu: tensor(0.8229)
===== Metropolis within Gibbs =====
Took 174.75 seconds to finish Program 1
Posterior Expectation mu: tensor(7.2083)
Posterior Variance mu: tensor(0.8607)
first attempt
===== Hamiltonian Monte Carlo =====
Took 323.58718633651733 seconds to finish Program 1
Posterior Expectation mu: tensor(7.2545)
Posterior Variance mu: tensor(0.7976)
```

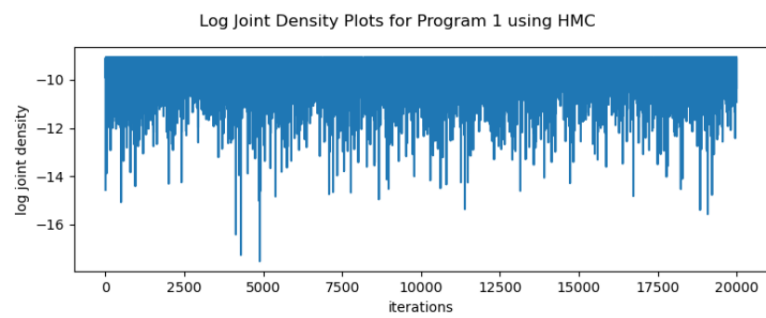
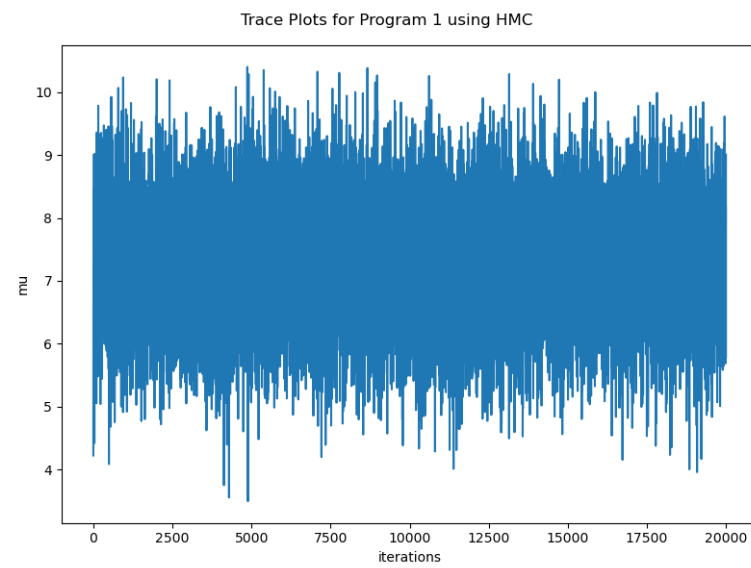
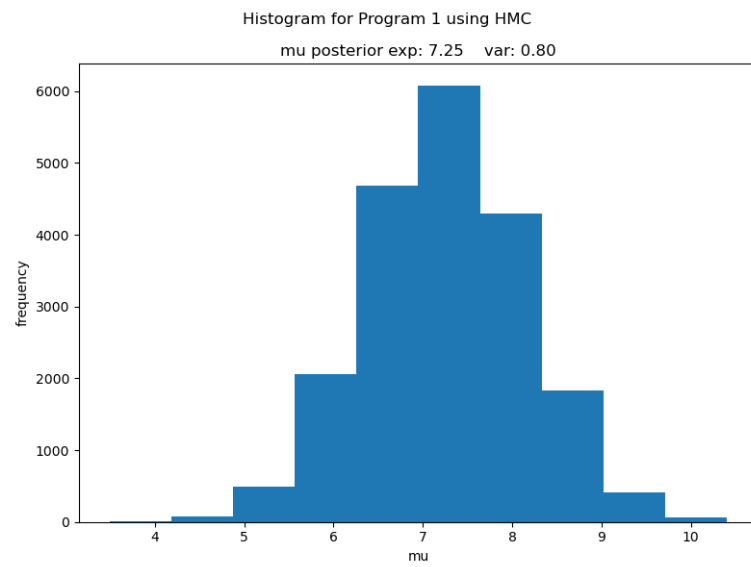
1.2 IS



1.3 MH-Gibbs



1.4 HMC

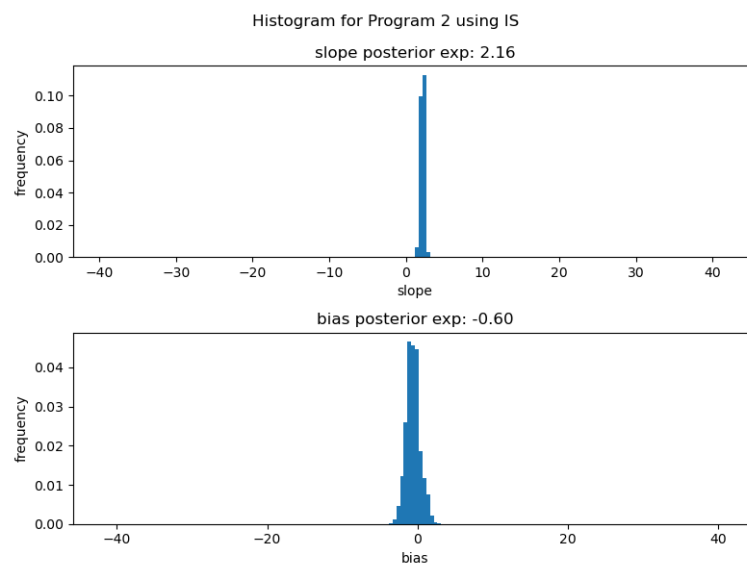


2 Program 2

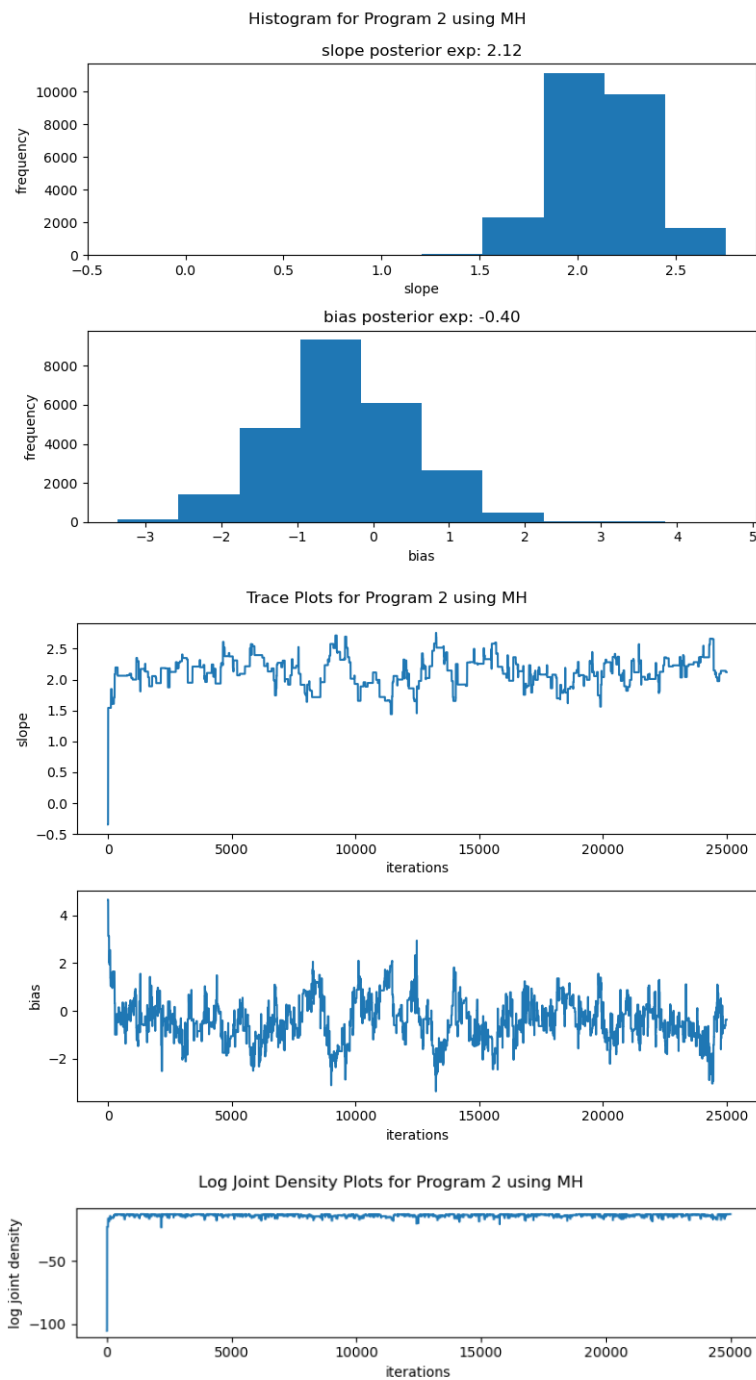
2.1 Wall-Clock Time

```
===== Likelihood Weighting =====
Took 345.71 seconds to finish Program 2
Posterior Expectation slope: tensor(2.1622)
posterior covariance of slope and bias:
[[ 0.05525479 -0.19291554]
 [-0.19291554  0.85292497]]
Posterior Expectation bias: tensor(-0.5960)
===== Metropolis within Gibbs =====
Took 231.19 seconds to finish Program 2
Posterior Expectation slope: tensor(2.1177)
posterior covariance of slope and bias:
[[ 0.04993265 -0.17395253]
 [-0.17395253  0.78755741]]
Posterior Expectation bias: tensor(-0.3951)
first attempt
===== Hamiltonian Monte Carlo =====
Took 420.389123916626 seconds to finish Program 2
Posterior Expectation slope: tensor(2.1442)
posterior covariance of slope and bias:
[[ 0.0529382 -0.19120865]
 [-0.19120865  0.82222667]]
Posterior Expectation bias: tensor(-0.4859)
```

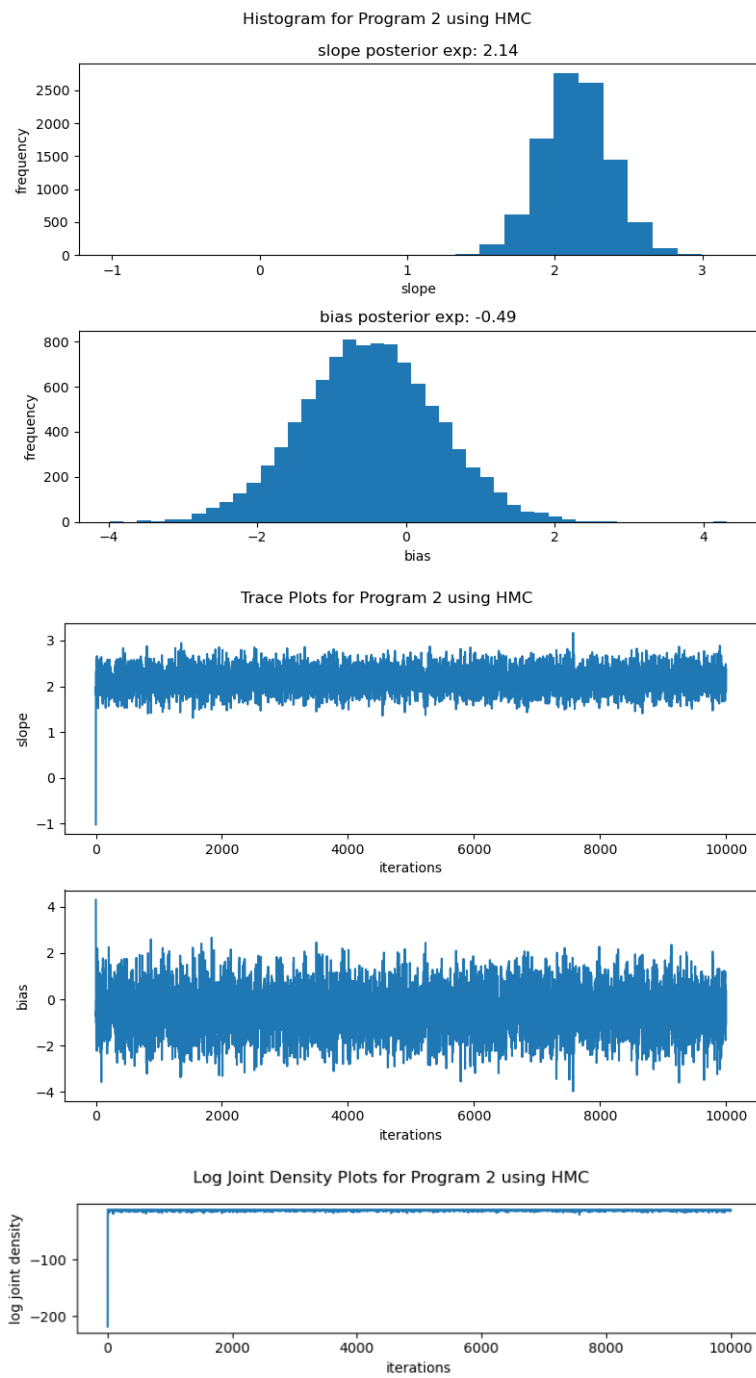
2.2 IS



2.3 MH-Gibbs



2.4 HMC

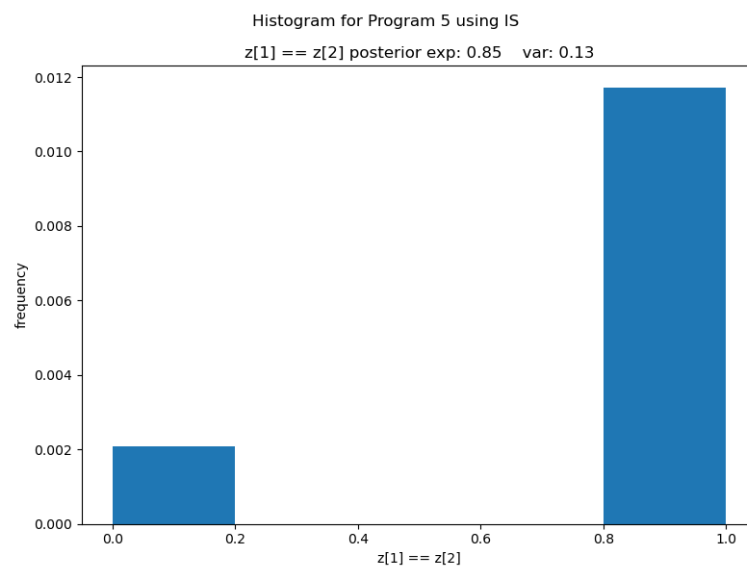


3 Program 3

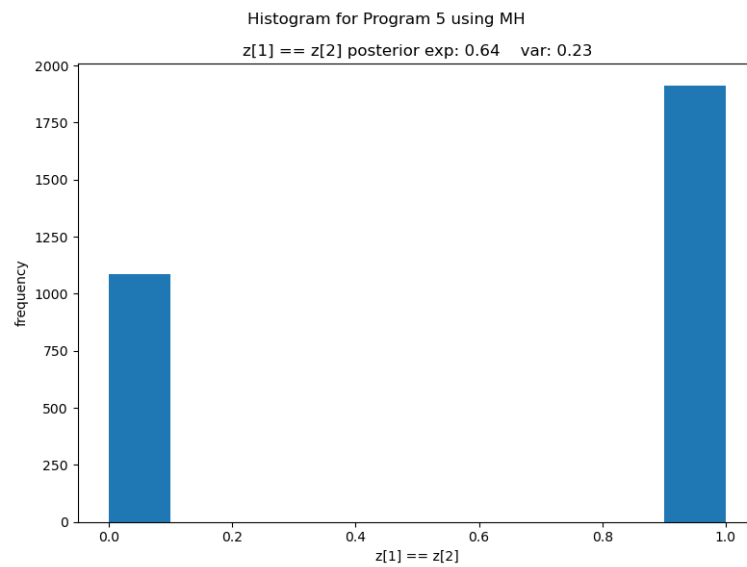
3.1 Wall-Clock Time

```
===== Likelihood Weighting =====  
Took 436.88 seconds to finish Program 5  
Posterior Expectation z[1] == z[2]: tensor(0.8492)  
Posterior Variance z[1] == z[2]: tensor(0.1281)  
===== Metropolis within Gibbs =====  
Took 250.12 seconds to finish Program 5  
Posterior Expectation z[1] == z[2]: tensor(0.6373)  
Posterior Variance z[1] == z[2]: tensor(0.2312)
```

3.2 IS



3.3 MH-Gibbs

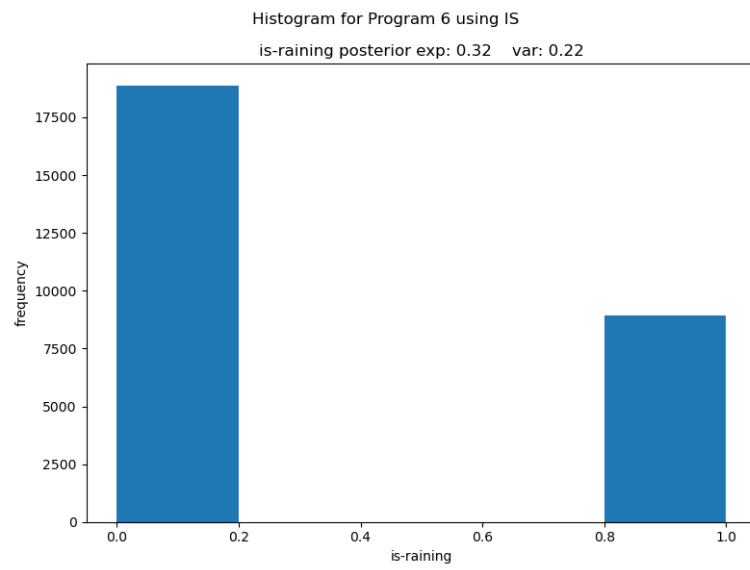


4 Program 4

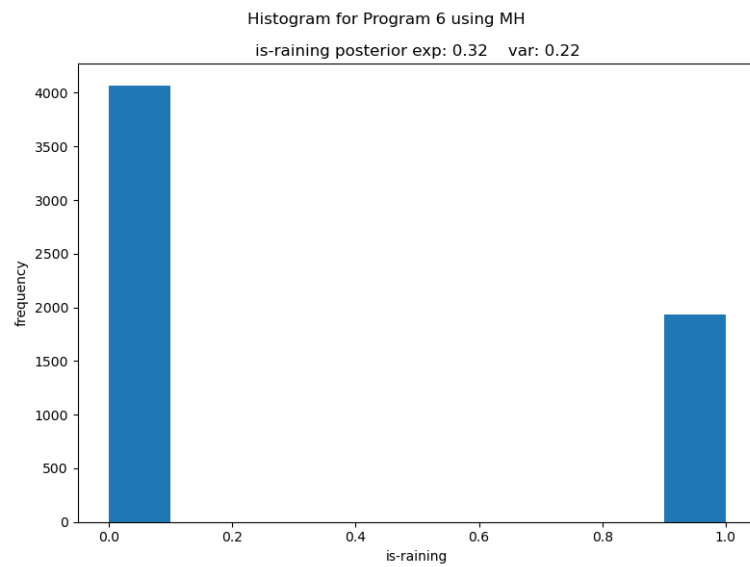
4.1 Wall-Clock Time

```
===== Likelihood Weighting =====  
Took 138.92 seconds to finish Program 6  
Posterior Expectation is-raining: tensor(0.3213)  
Posterior Variance is-raining: tensor(0.2180)  
===== Metropolis within Gibbs =====  
Took 65.43 seconds to finish Program 6  
Posterior Expectation is-raining: tensor(0.3222)  
Posterior Variance is-raining: tensor(0.2184)
```


4.2 IS



4.3 MH-Gibbs

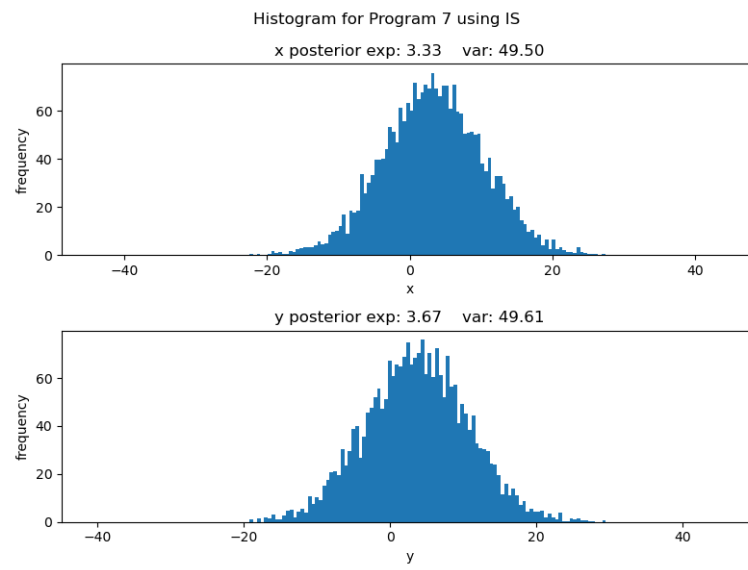


5 Program 5

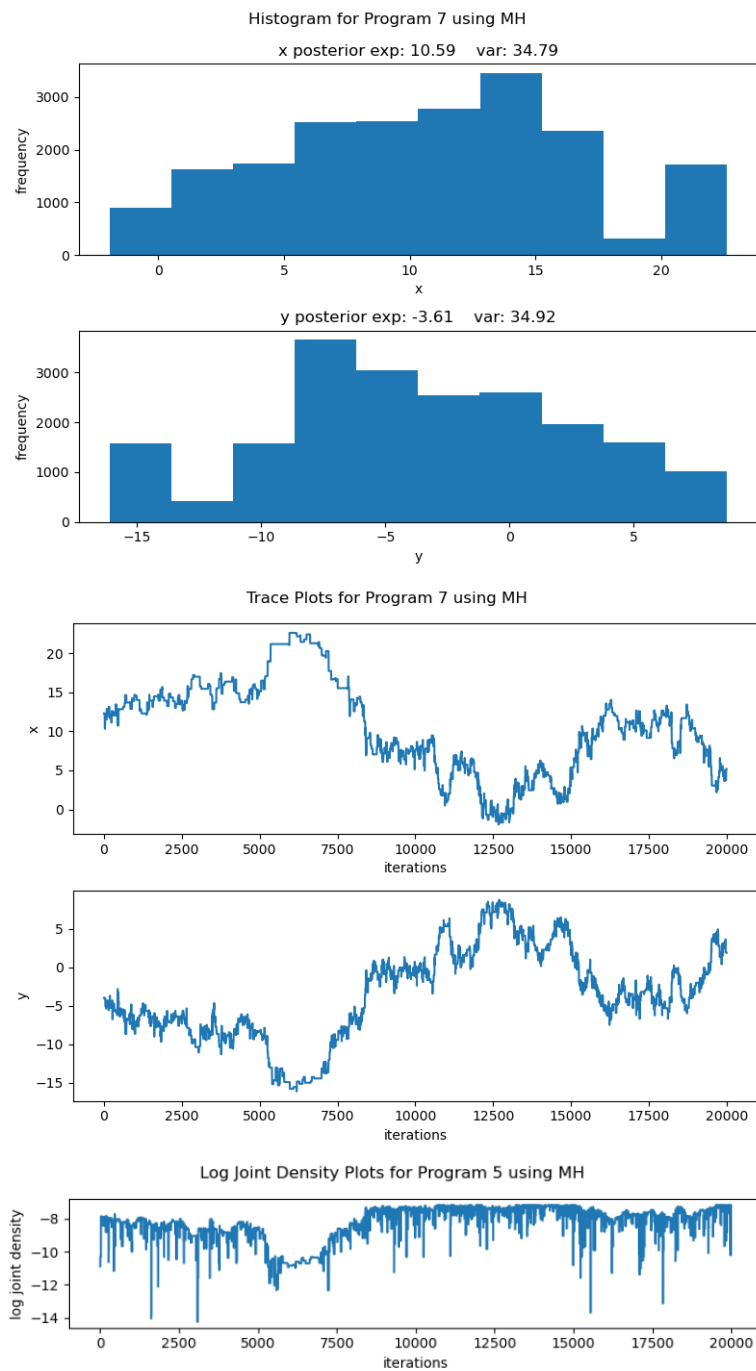
5.1 Wall-Clock Time

```
===== Likelihood Weighting =====
Took 71.46 seconds to finish Program 7
Posterior Expectation x: tensor(3.3320)
Posterior Variance x: tensor(49.5026)
Posterior Expectation y: tensor(3.6669)
Posterior Variance y: tensor(49.6122)
===== Metropolis within Gibbs =====
Took 61.62 seconds to finish Program 7
Posterior Expectation x: tensor(10.5921)
Posterior Variance x: tensor(34.7903)
Posterior Expectation y: tensor(-3.6118)
Posterior Variance y: tensor(34.9180)
C:\Users\kevin\OneDrive - The University Of British C
    fig, axs = plt.subplots(len(parameter_names), figsi
first attempt
===== Hamiltonian Monte Carlo =====
second attempt
===== Hamiltonian Monte Carlo =====
third attempt
===== Hamiltonian Monte Carlo =====
fourth attempt
===== Hamiltonian Monte Carlo =====
fifth attempt
===== Hamiltonian Monte Carlo =====
Took 279.50547552108765 seconds to finish Program 7
Posterior Expectation x: tensor(3.4514)
Posterior Variance x: tensor(38.6782)
Posterior Expectation y: tensor(3.5359)
Posterior Variance y: tensor(38.6634)
```

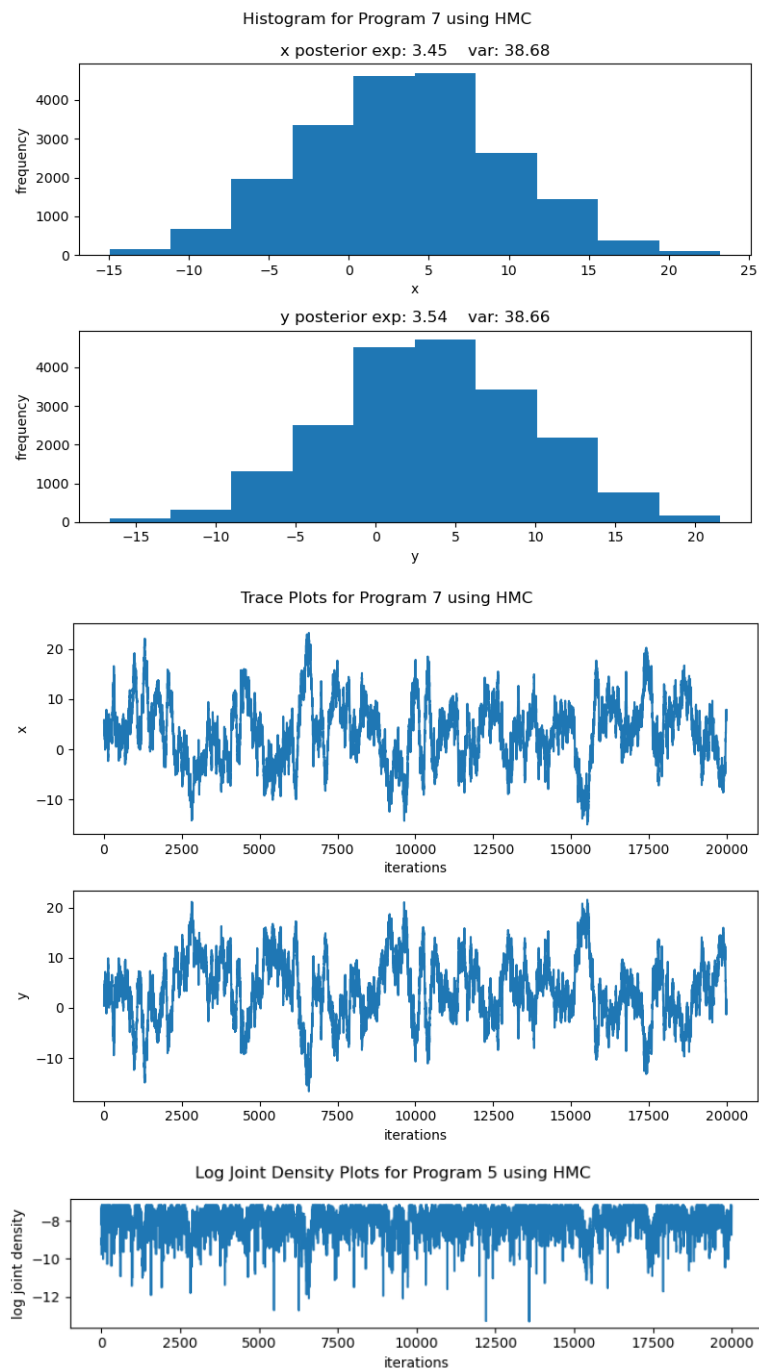
5.2 IS



5.3 MH-Gibbs



5.4 HMC



6 Code Snippets

```
def sample(self, num_samples, num):
    """
    importance sampling procedure
    """

    print('=' * 10, 'Likelihood Weighting', '=' * 10)

    ast = load_ast('programs/saved_asts/hw3/program{}.pkl'.format(num))
    sig = {'logW': 0}
    samples = []
    for i in range(num_samples):
        r_i, sig_i = evaluate_program(ast, sig, 'IS')
        logW_i = copy.deepcopy(sig['logW'])
        samples.append([r_i, logW_i])
        sig['logW'] = 0

    return samples


def gibbs_step(self, graph):
    """
    performs one entire gibbs sweep

    :param graph: current state of the graph
    :return:
    """

    # get a list of the latent variables
    regex = re.compile(r'observe*')
    latent_vars = [vertex for vertex in graph[1]['V'] if not regex.match(vertex)]

    # populate the proposal map self.Q
    self.populate_proposals(graph, latent_vars)

    for latent_var in latent_vars:
        # get the prior distribution for node latent_var
        d = deterministic_eval(substitute_sampled_vertices(self.Q[latent_var], graph[1]['V']))

        # make a copy of the graph (line 15 of algorithm 1 on pg.82 in the text)
        graph_propose = copy.deepcopy(graph)

        # sample a new latent_var from the proposal distribution (just the prior for now)
        graph_propose[1]['V'][latent_var] = d.sample()

        # compute acceptance ratio
        alpha = self.accept(latent_var, graph_propose, graph)

        # MH step
        u = torch.rand(1)
        if u < alpha:
            graph = copy.deepcopy(graph_propose)

    return graph


def accept(self, latent_var, graph_propose, graph):
    """
    compute acceptance probability

    :param latent_var: node of interest
    :param graph_propose: new proposed graph
    :param graph: original graph
    :return:
    """

    # both transition kernels are the same because we're just sampling from the prior
    d_old = deterministic_eval(substitute_sampled_vertices(self.Q[latent_var], graph[1]['V']))
    d_new = deterministic_eval(substitute_sampled_vertices(self.Q[latent_var], graph_propose[1]['V']))

    # create variables for the old and new parameter values
    old_value = graph[1]['V'][latent_var]
    new_value = graph_propose[1]['V'][latent_var]

    # first part of the log prob
    log_alpha = d_new.log_prob(old_value) - d_old.log_prob(new_value)

    # v_x contains all the children of node latent_var
    v_x = graph[1]['A'][latent_var]

    # compute the ratio of the joint likelihoods
    variable_bindings = graph[1]['V']
    variable_bindings_propose = graph_propose[1]['V']
```

```

for v in V_X:
    # substitute variable bindings
    raw_expression = graph_propose[1]['P'][v]
    expression = substitute_sampled_vertices(raw_expression, variable_bindings_propose)
    log_alpha += deterministic_eval(expression)

    # substitute variable bindings
    raw_expression = graph[1]['P'][v]
    expression = substitute_sampled_vertices(raw_expression, variable_bindings)
    log_alpha -= deterministic_eval(expression)

# NEED TO INCLUDE THE LATENT VARIABLE ITSELF TOO!!!!
# likelihood under new model
raw_expression = ['observe*', graph_propose[1]['P'][latent_var][1], graph_propose[1]['Y'][latent_var]]
expression = substitute_sampled_vertices(raw_expression, variable_bindings_propose)
log_alpha += deterministic_eval(expression)

# likelihood under old model
raw_expression = ['observe*', graph[1]['P'][latent_var][1], graph[1]['Y'][latent_var]]
expression = substitute_sampled_vertices(raw_expression, variable_bindings)
log_alpha -= deterministic_eval(expression)

# exponentiate to exit log-space
return torch.exp(log_alpha)

```

```

def hmc(self, X, Y, graph, num_samples):
    graph_old = copy.deepcopy(graph)
    X_old = copy.deepcopy(X)

    samples = []
    for s in range(num_samples):
        R_old = torch.distributions.multivariate_normal.MultivariateNormal(torch.zeros(len(self.latent_vars)),
                                                                            self.M).sample()

        # X_new is a dictionary containing the proposed new variable bindings
        X_new, R_new = self.leapfrog(X_old, Y, graph, R_old, graph)

        # MH step
        u = torch.rand(1)
        alpha = torch.exp(-self.H(X_new, Y, graph, R_new) + self.H(X_old, Y, graph_old, R_old))
        if u < alpha:
            X_old = X_new

        samples.append(deterministic_eval(substitute_sampled_vertices(graph[2], X_old)))

    return samples

```

```

def leapfrog(self, X, Y, R, graph):
    """
    perform leapfrog integration

    :param X: dictionary containing latent nodes and their corresponding values
    :param Y: dictionary containing observed nodes and their corresponding values
    :param graph: graphical model
    :param R: momentum
    :return: updated values for: X, R_half
    """

    R_half = R - 0.5 * self.epsilon * self.grad_U(X, Y, graph)

    for t in range(self.T - 1):
        # update latent variables
        for idx, node in enumerate(self.latent_vars):
            X[node] = X[node].detach() + (self.epsilon * R_half[idx])
            X[node].requires_grad = True

        R_half = R_half.detach() - self.epsilon * self.grad_U(X, Y, graph)

    # one last update
    for idx, node in enumerate(self.latent_vars):
        X[node] = X[node].detach() + (self.epsilon * R_half[idx])
        X[node].requires_grad = True

    R_half = R_half.detach() - 0.5 * self.epsilon * self.grad_U(X, Y, graph)

    return X, R_half

```

```

@staticmethod
def U(X, Y, graph):
    """
    returns potential energy U (log of the joint)

    :param graph: graphical model
    :return:
    """
    # iterate over every node in the graph to obtain the joint
    log_gamma = 0 # accumulator

    # compute log likelihood for latent nodes
    for node in list(X.keys()):
        # substitute variables with their values
        expression = substitute_sampled_vertices(graph[1]['P'][node], {**X, **Y})

        # if latent compute likelihood using prior distribution
        expression[0] = 'observe*'
        expression.append(X[node])
        log_gamma += deterministic_eval(expression)

    # compute log likelihood for observed nodes
    for node in list(Y.keys()):
        expression = substitute_sampled_vertices(graph[1]['P'][node], {**X, **Y})
        log_gamma += deterministic_eval(expression)

    return -log_gamma

def H(self, X, Y, graph, R):
    """
    use auxiliary variable technique

    :param graph: graphical model
    :param R: momentum
    :return:  $\exp(-U(X)) + 0.5 * R.T * \text{inv}(M) * R$ 
    """
    if self.M.size() == torch.Size([]):
        return torch.exp(-self.U(X, Y, graph) +
                        0.5 * torch.matmul(R.T,
                                           torch.FloatTensor([torch.matmul(torch.FloatTensor([self.M]), R)])))

    return torch.exp(-self.U(X, Y, graph) +
                    0.5 * torch.matmul(R.T,
                                        torch.matmul(torch.inverse(self.M), R)))

class Dirac:
    def __init__(self, param):
        """
        implementation of dirac function

        :param param:  $x + y$ 
        """
        self.param = param
        self.log_norm_const = torch.log(torch.tensor(2)) + torch.lgamma(torch.tensor(5/4))

    def log_prob(self, obs):
        log_prob = -(self.param - obs)**4 - self.log_norm_const
        return log_prob

```