$$m_1 \ddot{q}_1 + k_1 q_1 - k_1 q_2 = f(t) + \omega(t)$$
 $m_2 \ddot{q}_2 = -k_1 q_1 + (k_1 + k_2) q_2 - k_2 q_3 = 0$ 
 $m_3 \ddot{q}_6 - k_2 q_2 + k_2 q_3 = 0$ 

$$\begin{bmatrix} \dot{v}_{i} \\ \dot{q}_{i} \\ \dot{q}_{i$$

$$A_{p} = \begin{bmatrix} 0 & 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 0 & 1 \\ \frac{1}{1} \frac{$$

General form: Re = Apap + Bpu + Dpw 1 = ( R x p + By, 4 + Dy w Yz= Crxp + Byzu Z= Moxp + Dzw Z = Mp xp > Z= xp (full state feedback) Comparing with earlier notation { it = Ap xp + Bp u + Dp w Ap = Ap Bp = Bp we only consider leg from w to y. Dp - Op as to minimize Pep, and ensure CL system مع = دا<sub>ع</sub> minimize Pep and ensure CL system By2 = 0 Shability by placing the CL poles in LHP Dy =0 such lat Re(z) < -1. destred region for CL-Poles used in "Iming". Lumped system representation B = [DP BP] C = ( Cp. Cp.) r = size (Byz) = [size (Jz) size(u)] obj = [0 0 0 1] ⇒ min [cp ⇒ optimal H2 control design min (of Fee + P Fep) s.t. Tec < 4 Tep < V2 [Yi, Yz, Gi, Sei, X] = msfayn (S. T. obj., region) Fee of Fep of static state Clayston feelback contra gain (4=4 x) Acc = Apt Bp.G.

```
Design validation:
                                       norm2 (ScL)
             where AP+PAT+BBT = 0 (Lyapunov equation)
            ( Con a grant the worked to a grant to
   [co < 12
                        Re(z) < 0 & Stability (CL) gueranteed
· Stability
CL poles: Re ( ): (Au) = Re(2) < -1 - pok placement (LL) in
(2)
                        in today didded date illate
 System simulation:
                                                (surrichay in book proug)
   simulate open (SOL) and closed (SCL) loop systems using
   "Isim" for t= 0 to t= 10 seconds with At = 0.01 ser.
                        plot (be, because) if the be with (is
Uncertain system:
  lot uncertainty in kz (nominal kz = 4000 N/m)
  i.e. 3600 x kz $ 4400 N/m
  for kz = 3600: 1: 4400
    hz-norm = norm2 (See) 4 using Ci
  plot (k2, h2-norm) - Per us. k2 with G,
Robust state-feedback controller design:
 Affine system representation
  A= Ao+ k2 A1
```

 $B_0 = B$ ,  $C_0 = C$ ,  $D_0 = D$   $P = [k_1]; k_1 \in [3600, 4400]$ 

So = ( Hisys ( A., G., C., D., 1) : milled in . S1 = 1 hays (A1, 0, 0, 0, 0, 0) affays = psys (P, [s., s.]) T = Size (Byz) = [Size (Je) size (4)] obj = [0 0 0 1] + optimal H2 control design (min [cp) region = liveg - LHP with Re(2) <-1 [ 4, , y, , G2, See, X] = msfsyn (affsys, r, obj, region) Tep of CL system representation obtained

Tep of CL system

Robust Static state feedback control gain (u = Gz 4p) (guaranteed Hz parformane) Acc = Ap+ Bp Gz ( Au = A + 6k) For k2 = 3600 : 1: 4400 die 1 of and of the hannorm = norma (Sec) + using Ga plot (kz, hanorm) + [top us. kz with G2 in in the state of the state of (min soon : I he known) is to provide them old south of the states the

in poise the feet account a recovery

it when it wight to Conon so in take

- 0 0 0 0 1 . N. To 0 1 0 0 2 1

inglish relations of suchasticities del

neither the stay of might

12 x 201 6

# Problem 1

#### CODE:

```
% PROBLEM 1
close all
clear
clc
% Define the system parameters
m1 = 20; % kg
m2 = 40; % kg
m3 = 30; % kg
k1 = 2000; % N/m
k2 = 4000; % N/m
% Define the system matrices
Ap = [0 \ 0 \ 0 \ 1 \ 0 \ 0];
     000010;
      000001;
      -k1/m1 k1/m1 0 0 0 0;
      k1/m2 (-k1-k2)/m2 k2/m2 0 0 0;
      0 k2/m3 -k2/m3 0 0 0]
Bp = [0; 0; 0; 1/m1; 0; 0]
Dp = [0; 0; 0; 1/m1; 0; 0]
Cp2 = [0 \ 0 \ 1 \ 0 \ 0]
By2 = 0
Dy = 0
```

```
% Lumped system representation
A = Ap;
B = [Dp Bp];
C = [zeros(size(Cp2)); Cp2];
D = [Dy zeros(size(By2)); zeros(size(Dy)) By2];
S = ltisys(A, B, C, D)

% State-feedback controller G1 to minimize the energy-to-peak gain Γep
r = size(By2);
obj = [0 0 0 1]; % Optimal H2 control design objective
region = lmireg; % h (half-plane) --> l (LHP, i.e., Re(z) < x0) --> -1 (x0 = -1)
[glopt, g2opt, G1, Sc1, X] = msfsyn(S, r, obj, region)
```

```
% Closed-loop system
Acl = Ap + Bp*G1
h2_norm = norm2(Scl)
% Scl = ss(Acl, [Dp zeros(size(Bp))], C, D)
% h2_norm = norm(Scl, 2)

% Verification
if((h2_norm < g2opt) && all(real(eig(Acl)) < -1.0))
    disp('Verification of H2 norm and pole location constraints successful!')
else
    disp('Verification of H2 norm and pole location constraints failed!')
end</pre>
```

```
% Define the lumped (disturbance + control) input
t = 0:0.01:10;
w amplitude = 100.0;
w_duration = 2;
w_pulse = w_amplitude * (t >= 0 & t <= w_duration);</pre>
u = zeros(size(w_pulse));
w = [w_pulse; u];
% Simulate the open-loop system response
Sol = ss(A, B, C, D)
[y_ol, t_out, x_ol] = lsim(Sol, w', t);
% Plot the results
figure;
sgtitle('Open-Loop System Response');
subplot(3, 1, 1);
plot(t, w(1, :), t, w(2, :));
legend('W', 'U');
subplot(3, 1, 2);
plot(t, x_ol(:, 1), ...
     t, x_ol(:, 2), ...
     t, x_ol(:, 3), ...
    t, x_ol(:, 4), ...
     t, x_ol(:, 5), ...
     t, x_ol(:, 6) ...
legend('X1', 'X2', 'X3', 'X4', 'X5', 'X6');
subplot(3, 1, 3);
plot(t, y_ol(:, 2));
```

```
legend('Y1');
% Simulate the closed-loop system response
Scl = ss(Acl, [Dp zeros(size(Bp))], C, D)
[y_cl, t_out, x_cl] = lsim(Scl, w', t);
% Plot the results
figure;
sgtitle('Closed-Loop System Response');
subplot(3, 1, 1);
plot(t, w(1, :), t, w(2, :));
legend('W', 'U');
subplot(3, 1, 2);
plot(t, x_cl(:, 1), ...
    t, x_cl(:, 2), ...
     t, x_cl(:, 3), ...
    t, x_cl(:, 4), ...
     t, x_cl(:, 5), ...
     t, x_cl(:, 6) ...
     );
legend('X1', 'X2', 'X3', 'X4', 'X5', 'X6');
subplot(3, 1, 3);
plot(t, y_cl(:, 2));
legend('Y1');
```

```
% Define uncertain system parameters
m1 = 20; % kg
m2 = 40; % kg
m3 = 30; % kg
k1 = 2000; % N/m
i = 1;
for k2 = 3600:1:4400 % N/m
    % Define the system matrices
    Ap = [0 \ 0 \ 0 \ 1 \ 0 \ 0]
          000010;
          000001;
          -k1/m1 k1/m1 0 0 0 0;
          k1/m2 (-k1-k2)/m2 k2/m2 0 0 0;
          0 k2/m3 -k2/m3 0 0 0];
    Bp = [0; 0; 0; 1/m1; 0; 0];
    Dp = [0; 0; 0; 1/m1; 0; 0];
    Cp2 = [0 \ 0 \ 1 \ 0 \ 0];
    By2 = 0;
    Dy = 0;
```

```
% Lumped system representation
    A = Ap;
    B = [Dp Bp];
    C = [zeros(size(Cp2)); Cp2];
    D = [Dy zeros(size(By2)); zeros(size(Dy)) By2];
    % Closed-loop system
    Acl = Ap + Bp*G1;
    Scl = ltisys(Acl, [Dp zeros(size(Bp))], C, D);
    h2 norm(i) = norm2(Scl);
    % Scl = ss(Acl, [Dp zeros(size(Bp))], C, D);
    % h2_norm(i) = norm(Scl, 2);
    i = i+1;
end
figure;
sgtitle('Uncertain System with Controller G_1');
plot(3600:1:4400, h2_norm)
xlabel('Spring Constant k_2')
ylabel('Energy to Peak Gain Γ_ep')
```

```
% Define the system parameters
m1 = 20; % kg
m2 = 40; % kg
m3 = 30; % kg
k1 = 2000; % N/m
% System matrices in affine form
A0 = [0 \ 0 \ 0 \ 1 \ 0 \ 0;
     000010;
      000001;
      -k1/m1 k1/m1 0 0 0 0;
      k1/m2 -k1/m2 0 0 0 0;
      0 0 0 0 0 0]
A1 = [0 \ 0 \ 0 \ 0 \ 0];
     000000;
     000000;
     000000;
     0 -1/m2 1/m2 0 0 0;
     0 1/m3 -1/m3 0 0 0]
B0 = B
C0 = C
D0 = D
% Uncertain LTI system in affine form
```

```
S0 = ltisys(A0, B0, C0, D0, 1)
S1 = ltisys(A1, zeros(size(B0)), zeros(size(C0)), zeros(size(D0)), 0)

% Parameter vector
P = pvec('box', [3600, 4400]) % k2

% Affine system
affsys = psys(P, [S0, S1])

% State-feedback controller G2 to minimize the energy-to-peak gain Fep
r = size(By2);
obj = [0 0 0 1]; % Optimal H2 control design objective
region = lmireg; % h (half-plane) --> 1 (LHP, i.e., Re(z) < x0) --> -1 (x0 = -1)
[g1opt, g2opt, G2, Sc1, X] = msfsyn(affsys, r, obj, region)

disp('Optimal guaranteed energy-to-peak gain for the uncertain system is:')
disp(g2opt)
```

```
% Define uncertain system parameters
m1 = 20; % kg
m2 = 40; % kg
m3 = 30; % kg
k1 = 2000; % N/m
i = 1;
for k2 = 3600:1:4400 \% N/m
    % Define the system matrices
    Ap = [0 \ 0 \ 0 \ 1 \ 0 \ 0;
          000010;
          000001;
          -k1/m1 k1/m1 0 0 0 0;
          k1/m2 (-k1-k2)/m2 k2/m2 0 0 0;
          0 k2/m3 -k2/m3 0 0 0];
    Bp = [0; 0; 0; 1/m1; 0; 0];
    Dp = [0; 0; 0; 1/m1; 0; 0];
    Cp2 = [0 \ 0 \ 1 \ 0 \ 0];
    By2 = 0;
    Dy = 0;
    % Lumped system representation
    A = Ap;
    B = [Dp Bp];
    C = [zeros(size(Cp2)); Cp2];
    D = [Dy zeros(size(By2)); zeros(size(Dy)) By2];
    % Closed-loop system
```

```
Acl = Ap + Bp*G2;
Scl = ltisys(Acl, [Dp zeros(size(Bp))], C, D);
h2_norm(i) = norm2(Scl);
% Scl = ss(Acl, [Dp zeros(size(Bp))], C, D);
% h2_norm(i) = norm(Scl, 2);
i = i+1;
end

figure;
sgtitle('Uncertain System with Controller G_2');
plot(3600:1:4400, h2_norm)
xlabel('Spring Constant k_2')
ylabel('Energy to Peak Gain Γ_ep')
```

#### **OUTPUT:**

```
Ap = 6 \times 6
          0
                      0
                                  0
                                        1.0000
                                                          0
                                                                       0
                      0
                                                     1.0000
                                                                       0
          0
                                  0
                                              0
                                                                 1.0000
           0
                      0
                                  0
                                               0
                                                          0
 -100.0000 100.0000
                                  0
                                              0
                                                           0
                                                                      0
   50.0000 -150.0000 100.0000
                                              0
                                                          0
                                                                       0
                                              0
                                                           0
          0 133.3333 -133.3333
                                                                       0
Bp = 6 \times 1
           0
           0
           0
    0.0500
           0
           0
Dp = 6 \times 1
           0
           0
           0
    0.0500
           0
           0
Cp2 = 1 \times 6
                    1
                           0
                                  0
                                          0
      0
             0
By2 = 0
Dy = 0
S = 9 \times 9
          0
                      0
                                  0
                                        1.0000
                                                          0
                                                                      0
                                                                                   0
0
      6.0000
           0
                      0
                                  0
                                              0
                                                    1.0000
                                                                                   0
0
            0
          0
                      0
                                  0
                                              0
                                                           0
                                                                1.0000
                                                                                   0
 -100.0000 100.0000
                                  0
                                              0
                                                           0
                                                                      0
                                                                            0.0500
0.0500
                  0
   50.0000 -150.0000 100.0000
                                              0
                                                           0
                                                                      0
                                                                                   0
            0
```

	0 133	3.3333 -13	33.3333	0	0	0	0
0	0	0	0	0	0	0	0
Ο	0	U	U	U	U	U	U
O	0	0	1.0000	0	0	0	0
0	0						
	0	0	0	0	0	0	0
0	-Inf						

Select a region among the following:

- h) Half-plane
- d) Disk
- c) Conic sector
- e) Ellipsoid
- p) Parabola
- s) Horizontal strip
- m) Matrix description of the LMI region
- q) Quit

Select a region among the following:

- h) Half-plane
- d) Disk
- c) Conic sector
- e) Ellipsoid
- p) Parabola
- s) Horizontal strip
- m) Matrix description of the LMI region
- q) Quit

Optimization of 0.000 \* G^2 + 1.000 \* H^2 :

Solver for linear objective minimization under LMI constraints

Iterations : Best objective value so far

2 3 5 6 7 8 0.123781 9 0.101980 10 0.101980 11 0.028857 12 0.028857 13 0.028857 14 0.012000 15 0.012000 16 0.010372 17 0.010372 18 0.010372 6.093835e-03 19 20 6.093835e-03

```
21
                  6.093835e-03
22
                  3.043563e-03
23
                  3.043563e-03
24
                  3.043563e-03
25
                  1.240950e-03
26
                  1.240950e-03
27
                  1.231455e-03
                  3.550594e-04
28
29
                  2.305720e-04
30
                  2.305720e-04
31
                  1.078822e-04
32
                  1.078822e-04
33
                  1.078822e-04
34
                  5.183917e-05
35
                  5.183917e-05
36
                  5.183917e-05
37
                  2.614067e-05
38
                  2.614067e-05
39
                  2.614067e-05
40
                  1.606369e-05
41
                  1.606369e-05
42
                  1.596526e-05
43
                  3.082070e-06
44
                  3.082070e-06
45
                  2.599298e-06
                  2.599298e-06
46
                  2.599298e-06
47
48
                  9.165308e-07
```

Result: reached the target for the objective value

best objective value: 9.165308e-07

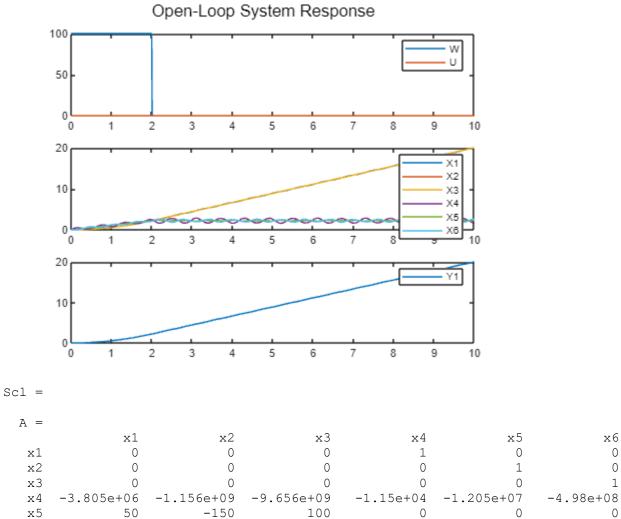
f-radius saturation: 0.520% of R = 1.00e+10

Guaranteed H2 performance: 9.57e-04

```
glopt =
    []
g2opt = 9.5736e-04
G1 = 1 \times 6
10^{11} \times
   -0.0008
              -0.2312
                         -1.9312
                                  -0.0000 -0.0024
                                                         -0.0996
Scl = 9 \times 8
10<sup>9</sup> ×
                                0
          0
                     0
                                     0.0000
                                                0
                                                                 0
                                                                             0
0.0000
                                                0.0000
          0
                     0
                                0
                                           0
                                                                 0
                                                                             0
0
          0
                     0
                                0
                                           0
                                                      0
                                                           0.0000
                                                                             0
0
   -0.0038
              -1.1562
                         -9.6558
                                    -0.0000
                                               -0.0120
                                                           -0.4980
                                                                       0.0000
0
   0.0000
              -0.0000
                         0.0000
                                          0
                                                      0
                                                                 0
                                                                             0
0
             0.0000
                         -0.0000
                                          0
                                                     0
          0
                                                                 0
                                                                             0
0
```

```
0
                 0
                       0
                                      0
                                                0
                                                           0
                                                                    0
0
                        0.0000
                                                 0
         0
                   0
                                       0
                                                           0
                                                                      0
0
                   0
                            0
                                                 0
                                                           0
         0
                                       0
                                                                      0
Inf
X = 6 \times 6
10^{6} \times
                       -0.0000
                                 -0.0080
                                           -0.0002
    0.0003
              0.0000
                                                     -0.0000
                                           -0.0000
   0.0000
              0.0000
                       0.0000
                                 0.0000
                                                     -0.0000
                                -0.0000
                                           0.0000
   -0.0000
              0.0000
                       0.0000
                                                     -0.0000
  -0.0080
              0.0000
                      -0.0000
                                 1.3121
                                          -0.0003
                                                     -0.0000
   -0.0002
             -0.0000
                       0.0000
                                 -0.0003
                                          0.0002
                                                     0.0000
   -0.0000
             -0.0000
                       -0.0000
                                 -0.0000
                                          0.0000
                                                      0.0000
Acl = 6 \times 6
10<sup>9</sup> ×
         0
                  0
                           0
                                 0.0000
                                                 0
                                            0.0000
         0
                  0
                             0
                                       0
                                                           0
         0
                  0
                           0
                                       0
                                                 0
                                                      0.0000
   -0.0038
             -1.1562
                       -9.6558
                                 -0.0000
                                           -0.0120
                                                     -0.4980
    0.0000
           -0.0000
                       0.0000
                                       0
                                                 0
             0.0000
                       -0.0000
                                       0
                                                 0
                                                           0
        0
h2 norm = 1.8822e-11
Verification of H2 norm and pole location constraints successful!
Sol =
  A =
                  x2
                          xЗ
                                           x5
           x1
                                   \times 4
                                                   x6
   x1
           0
                  0
                           0
                                   1
                                            0
                                                    0
  x2
           0
                   0
                           0
                                    0
                                            1
                                                    0
  xЗ
           0
                   0
                            0
                                    0
                                            0
                                                    1
                                                    0
  x4
         -100
                 100
                           0
                                   0
                                            0
           50
                -150
   x5
                          100
                                    0
                                            0
                                                    0
          0
              133.3 -133.3
                                    0
                                            0
                                                    0
  х6
  B =
               u2
         u1
   x1
         0
               0
                0
   x2
          0
                0
   xЗ
          0
       0.05 0.05
   x4
                0
  x5
         0
          0
                0
   х6
  C =
       x1 x2
               xЗ
                  \times 4
                       x5 x6
       0 0
              0
                  0
                      0
                          0
   у1
          0
               1
                    0
                        0
                            0
  у2
       0
  D =
      u1 u2
   у1
       0
          0
        0
            0
   y2
```

Continuous-time state-space model.



-133.3

x6

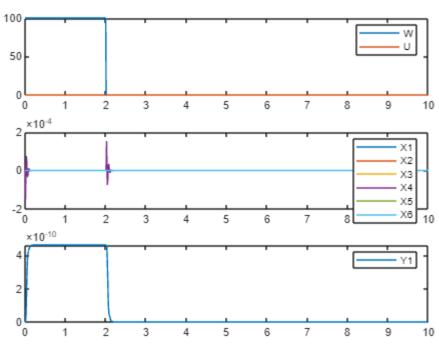
х6

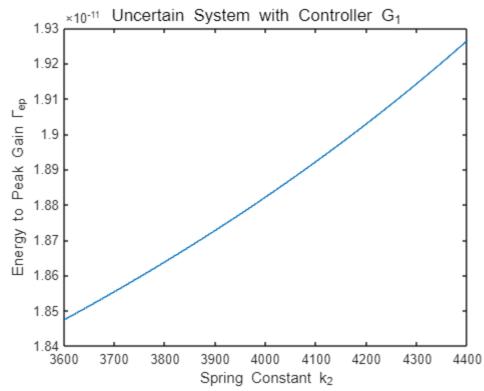
133.3

Continuous-time state-space model.

#### Model Properties

## Closed-Loop System Response





ΑO	=	6×6					
		0	0	0	1	0	0
		0	0	0	0	1	0
		0	0	0	0	0	1

```
-100
      100 0 0
                         0
    50 -50
              0
                     0
                                 0
                           0
                0
                      0
                                 0
    0
         0
                            0
A1 = 6 \times 6
                 0
                            0
                                      0
                                               0
                                                         0
         0
                 0
                                      0
                                                0
         0
                            0
                                                         0
         0
                  0
                            0
                                      0
                                                0
         0
                  0
                            0
                                      0
                                                0
                                                         0
         0
            -0.0250
                      0.0250
                                     0
                                               0
                                                         0
         0
             0.0333
                      -0.0333
                                     0
                                               0
                                                         0
B0 = 6 \times 2
        0
                  0
        0
                  0
        0
                  0
    0.0500
             0.0500
        0
                0
        0
                  0
C0 = 2 \times 6
     0
          0
                0
                      0
                          0
                                  0
     0
         0
               1
                      Ω
                           0
D0 = 2 \times 2
     0
         0
     0
         0
S0 = 9 \times 9
                           0
                                 1.0000
        0
                  0
                                           0
                                                         0
                                                                   0
    6.0000
0
                            0
                                          1.0000
        0
                  0
                                     0
                                                         0
                                                                   0
         0
        0
                            0
                                      0
                                                0
                                                   1.0000
                                                                   0
        0
-100.0000 100.0000
                            0
                                                0
                                                         0
                                                              0.0500
0.0500 0
   50.0000 -50.0000
                            0
                                      0
                                                0
                                                         0
                                                                   0
      0
0
        0
                            0
                                      0
                                                0
                                                         0
                                                                   0
                  0
0
         0
                  0
                            0
                                      0
                                                0
                                                         0
                                                                   0
         0
         0
        0
                  0
                       1.0000
                                      0
                                                0
                                                         0
                                                                   0
0
         0
        0
                  0
                          0
                                                0
0
      -Inf
S1 = 9 \times 9 \text{ complex}
   0.0000 - 1.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
6.0000 + 0.0000i
   0.0000 + 0.0000i 0.0000 - 1.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i
   0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 - 1.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i
  0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 - 1.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i
```

```
0.0000 + 0.0000i -0.0250 + 0.0000i 0.0250 + 0.0000i 0.0000 + 0.0000i
0.0000 - 1.0000i
              0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i
  0.0000 + 0.0000i 0.0333 + 0.0000i -0.0333 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 - 1.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i
  0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i
  0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i
  0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
-Inf + 0.0000i
P = 2 \times 5
            3600 4400
0 0
                                     0
                                      0
        1
affsys = 9 \times 26 complex
    -Inf + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0010 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
              0.0000 + 0.0000i
              0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 - 0.0010i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i
  0.0020 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0010 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 - 0.0010i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0010 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i
  0.0020 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0010 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i
  0.0060 + 0.0000i 0.0000 + 0.0000i -0.1000 + 0.0000i 0.1000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
              0.0001 + 0.0000i
                                              0.0000 + 0.0000i
                              0.0000 + 0.0000i
0.0001 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 - 0.0010i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i
  0.0020 + 0.0000i 0.0000 + 0.0000i 0.0500 + 0.0000i -0.0500 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
                              0.0000 + 0.0000i
              0.0000 + 0.0000i
0.0000 + 0.0000i
                                              0.0000 + 0.0000i
0.0000 + 0.0000i -0.0000 + 0.0000i 0.0000 + 0.0000i
                                              0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i
  0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
```

```
0.0000 + 0.0000i
              0.0000 + 0.0000i 0.0000 + 0.0000i
                                            0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i -0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i
  0.0100 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i
  0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0010 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i
  0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i
                              -Inf + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
-Inf + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i 0.0000 + 0.0000i
0.0000 + 0.0000i 0.0000 + 0.0000i
Select a region among the following:
   Half-plane
h)
  Disk
d)
C)
  Conic sector
  Ellipsoid
e)
p) Parabola
s) Horizontal strip
m) Matrix description of the LMI region
   Quit
q)
Select a region among the following:
  Half-plane
h)
  Disk
d)
  Conic sector
C)
  Ellipsoid
e)
  Parabola
p)
s) Horizontal strip
m) Matrix description of the LMI region
q)
Optimization of 0.000 * G^2 + 1.000 * H^2 :
 ______
Solver for linear objective minimization under LMI constraints
Iterations : Best objective value so far
```

1 2

3 4 5 6 7 8 9	
11 12 13 14 15 16 17 18 19 20 21 22 23	0.174617 0.084142 0.084142 0.084142 0.035606 0.035606 0.035606 0.019613 0.019613 0.019613 0.011597 0.011597
24 25 26 27 28 29 30 31 32 33 34 35	7.440620e-03 7.440620e-03 7.440620e-03 4.761071e-03 4.761071e-03 3.057379e-03 3.057379e-03 2.094639e-03 2.094639e-03 1.728979e-03
37 38 39 40 41 42 43 44 45 46 47 48	1.728979e-03 1.728979e-03 1.401557e-03 1.401557e-03 1.401557e-03 7.898234e-04 7.898234e-04 7.898234e-04 3.528895e-04 3.528895e-04 1.357855e-04 1.357855e-04
50 51 52 53 54 55 56 57 58	1.357855e-04 5.390543e-05 5.390543e-05 5.390543e-05 2.140225e-05 2.140225e-05 1.012459e-05 1.012459e-05

```
60
                   5.006723e-06
   61
                   5.006723e-06
   62
                   5.006723e-06
   63
                   2.216772e-06
   64
                   2.216772e-06
   65
                   2.216772e-06
   66
                   9.611413e-07
Result: reached the target for the objective value
         best objective value: 9.611413e-07
         f-radius saturation: 0.243\% of R = 1.00e+10
Guaranteed H2 performance: 9.80e-04
glopt =
   []
g2opt = 9.8038e-04
G2 = 1 \times 6
10^{11} \times
  -0.0009 -0.2480
                   -1.9988 -0.0000 -0.0028 -0.0997
Scl = 9 \times 19
10<sup>9</sup> ×
                         0
                                    0
                                             0
                                                 0.0000
     -Inf
                 0
      0
             0.0000
                          0
                                    0
                                             0
                                                  0
                                                          0.0000
        0
                0
                       0.0000
   0.0000
                 0
                       0
                                    0
                                                           0.0000
                                             0
                                                       0
                  0
                           0
                                    0
                                             0
                                                       0
0.0000
             0
                       0
                                0
   0.0000
                 0
                          0
                                    0
                                             0
                                                       0
0.0000
                       0
                                0
                                         0
   0.0000
                 0
                          0
                     -0.0046
                              -1.2400 -9.9938 -0.0000 -0.0138 -
   0.0000
                 0
                      0
                                0 -0.0046 -1.2400 -9.9938 -0.0000
0.4987 0.0000
-0.0138 -0.4987
                 0.0000
                                0
   0.0000
                 0
                     0.0000
                              -0.0000
                                        0.0000
                                                      0
                                                                0
                                       -0.0000
     0
                 0
                       0
                               0.0000
                                                  0.0000
                                                                0
         0
                  0
                           0
   0.0000
                          0
                               0.0000
                                       -0.0000
                                                     0
                 0
      0
                 0
                                 0
                                       0.0000
                                                 -0.0000
                          0
                                                                Ω
         0
                 0
                           0
   0.0000
                 0
                          0
                                    0
                                             0
                                                       0
                                                                0
                 0
                          0
                                    0
                                             0
                                                                0
         0
                                                       0
                 0
                           0
         0
        0
                 0
                          0
                                    0
                                         0.0000
                                                      0
                                                                0
        Ω
                 0
                          0
                                    0
                                          0
                                                   0.0000
                                                                0
        0
                 0
                           0
        0
                 0
                          0
                                    0
                                             0
                                                      0
                                                                0
        0
                                    0
                                             0
                                                                 0
               -Inf
                           0
         0
                0
                        -Inf
X = 6 \times 6
10^{5} \times
                                       -0.0008
   0.0013
            0.0000
                     0.0000
                             -0.0411
                                                 -0.0000
            0.0000
                     0.0000
                             0.0000 -0.0000
   0.0000
                                                 -0.0000
```

0 0

0

0

0

0

0

0

0

Ω

0

0

0.0000

-0.0411

-0.0008 -0.0000

0.0000

0.0000

0.0000

-0.0000

0.0000

-0.0000

5.9055

0.0058

0.0000

0.0058

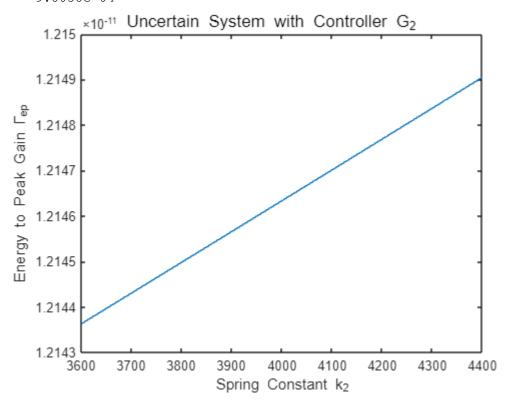
0.0007

-0.0000

0.0001

0.0000

-0.0000 -0.0000 -0.0000 0.0001 0.0000 0.0000 Optimal guaranteed energy-to-peak gain for the uncertain system is:  $9.8038e{-}04$ 



### **SCREENSHOT:**

