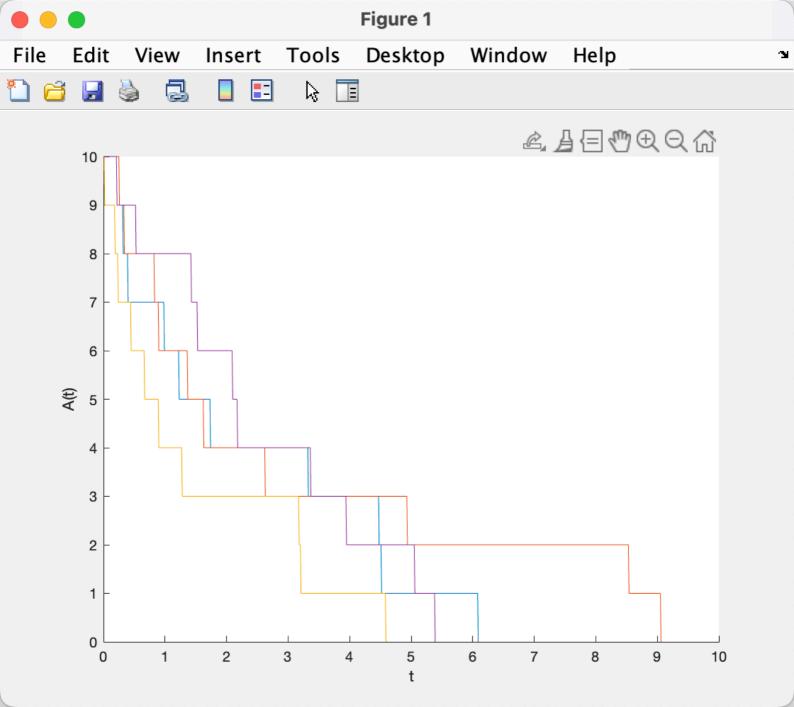
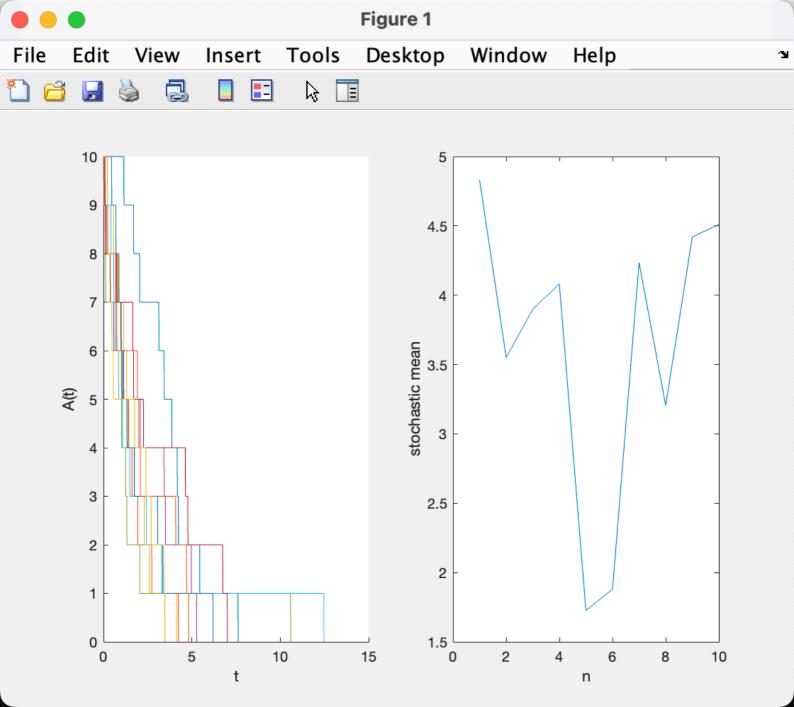
```
task3.m ×
  task1.m 🗶
               task2.m ×
                                           +
         clc
 1
 2
         clear all
 3
4
         close all
 5
 6
         n=4; % number of realizations
 7
         A0=10; % starting number of molecules
         % A=A0; % store the A
 8
 9
         detat=0.01; % time step
10
         k=0.5; % rate constant
11
12
         figure
13
         hold on
14
         for i=1:n
15
             A=A0; % store the A
16
             j=1;
17
             while A(j)>0
18
                  r=rand();
19
                  if r<A(end)*k*detat</pre>
20
                      A(j+1)=A(j)-1;
21
                  else
22
                      A(j+1)=A(j);
23
                  end
24
                  j=j+1;
25
             end
26
             plot(0:detat:(j-1)*detat,A)
27
              xlabel('t')
28
              ylabel('A(t)')
29
         end
30
```



```
task1.m
             X
                  task2.m × task3.m
                                                 +
 1
         clc
 2
         clear all
         close all
 3
 4
 5
 6
         n=10; % number of realizations
 7
         A0=10; % starting number of molecules
         % A=A0; % store the A
 8
9
         detat=0.01; % time step
         k=0.5; % rate constant
10
         mean_A=zeros(n,1);
11
12
13
         figure(1)
         subplot(121)
14
         hold on
15
         for i=1:n
16
              A=A0; % store the A
17
              j=1;
18
             while A(j)>0
19
20
                  r=rand();
21
                  tao=log(1/r)/A(j)/k;
22
                  count=ceil(tao/detat);
                  A(j+count)=A(j)-1;
23
24
                  j=j+count;
25
              end
26
              for ii=2:length(A)
27
                  if A(ii)==0
28
                      A(ii)=A(ii-1);
29
                  end
30
              end
31
              A(end)=0;
32
              mean_A(i)=mean(A);
33
              plot(0:detat:(length(A)-1)*detat,A)
34
              xlabel('t')
35
              ylabel('A(t)')
36
         end
         subplot(1,2,2)
37
         plot(1:n,mean_A)
38
39
         xlabel('n')
         ylabel('stochastic mean')
40
41
42
```



```
task1.m
                      task2.m
                                   🗶 task3.m 🗶
                 ×
 1
         clc
 2
          clear all
 3
          close all
 4
 5
 6
         n=100; % number of realizations
 7
         A0=10; % starting number of molecules
 8
         % A=A0; % store the A
 9
         detat=0.01; % time step
         k1=0.5; % rate constant of k1
10
11
         k2=0.2; % rate constant of k1
12
         mean_A=zeros(n,1);
13
14
         figure(1)
         subplot(121)
15
         hold on
16
17
         for i=1:n
     Ш
18
              A=A0; % store the A
19
              j=1;
     白
20
              while j<=100
21
                  r1=rand();
22
                  r2=rand();
23
                  alpha0=A(j)*k1+k2;
24
                  tao=log(1/r1)/alpha0;
25
                  count=ceil(tao/detat);
26
                  if r2<k2/alpha0
27
                      A(j+count)=A(j)+1;
28
                  else
29
                      A(j+count)=A(j)-1;
30
                  end
31
                  j=j+count;
32
              end
33
              for ii=2:length(A)
34
                  if A(ii)==0
35
                      A(ii)=A(ii-1);
36
                  end
37
              end
              A(end)=0;
38
39
              A=A(1:100); % 100 reactions
              mean_A(i)=mean(A);
40
41
                subplot(3,4,i)
42
              plot(0:detat:(length(A)-1)*detat,A)
              xlabel('t')
43
44
              ylabel('A(t)')
45
         end
          subplot(1,2,2)
46
47
         plot(1:n,mean_A)
         xlabel('n')
48
49
         ylabel('stochastic mean')
50
```

