

```

%Step 5 and 6

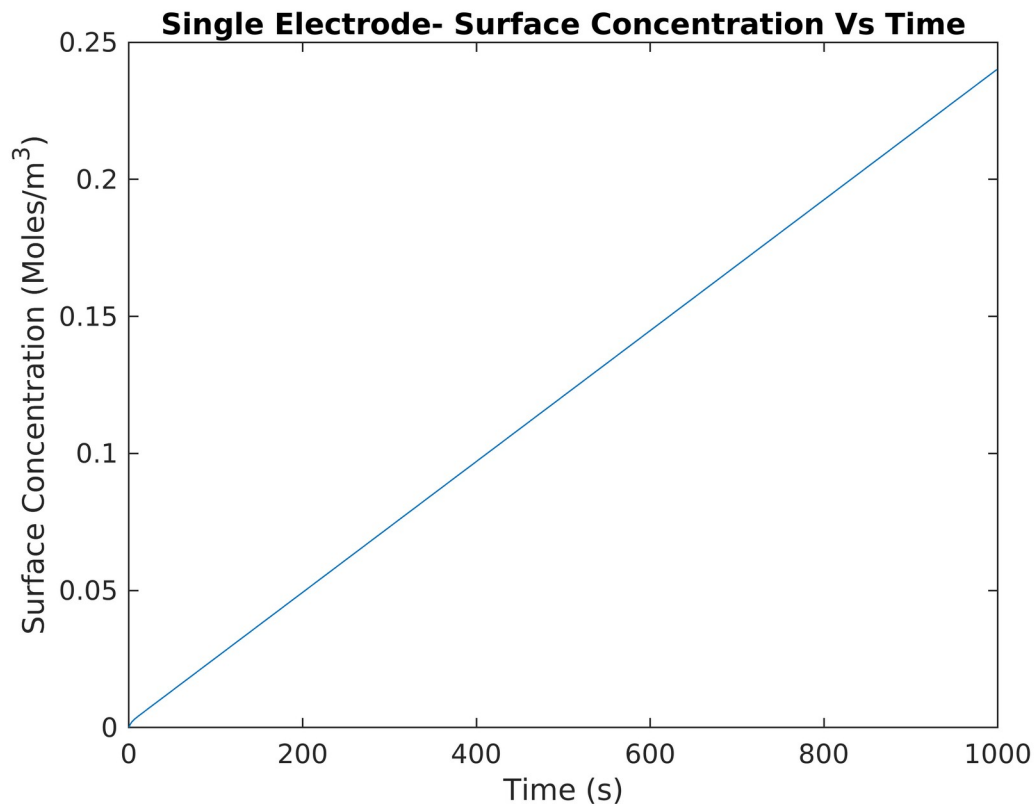
%Radius
R=10;
%Diffusivity
D=1;
nSteps=1000;
dt=1;

t=(0:dt:nSteps-1)';
u=ones(nSteps,1);
u_sin=sin(t/(10*pi));

%Initial Conditions
y1_0 = 0;
y2_0 =0;
sigmaY=0;

%Constant Input
y=lithium_model(R, D, u, y1_0, y2_0, dt, sigmaY);
plot(t, y)
title("Single Electrode- Surface Concentration Vs Time")
xlabel("Time (s)")
ylabel("Surface Concentration (Moles/m^3)")

```

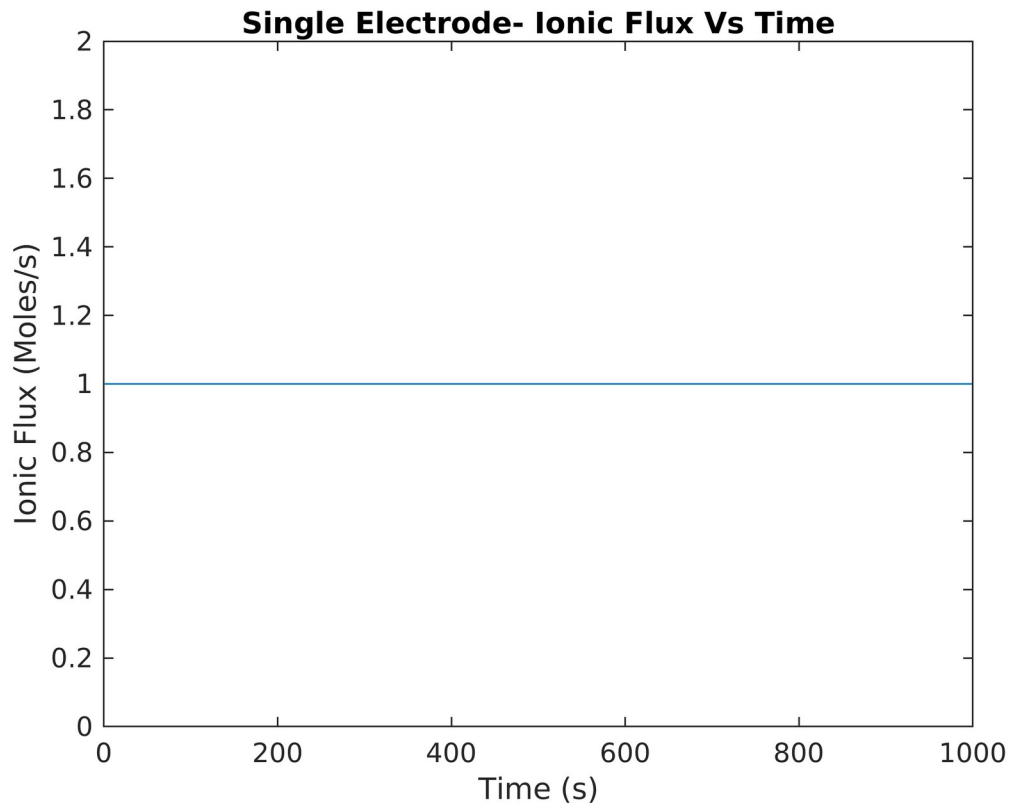


```

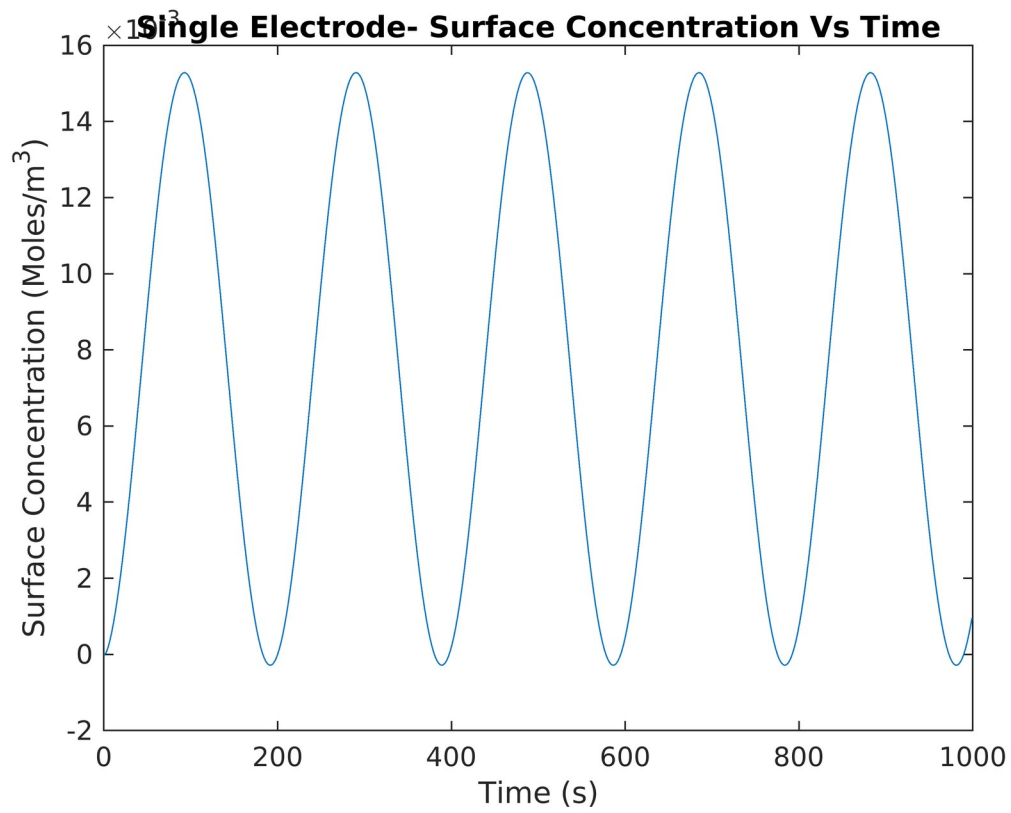
plot(t,u)
title("Single Electrode- Ionic Flux Vs Time")

```

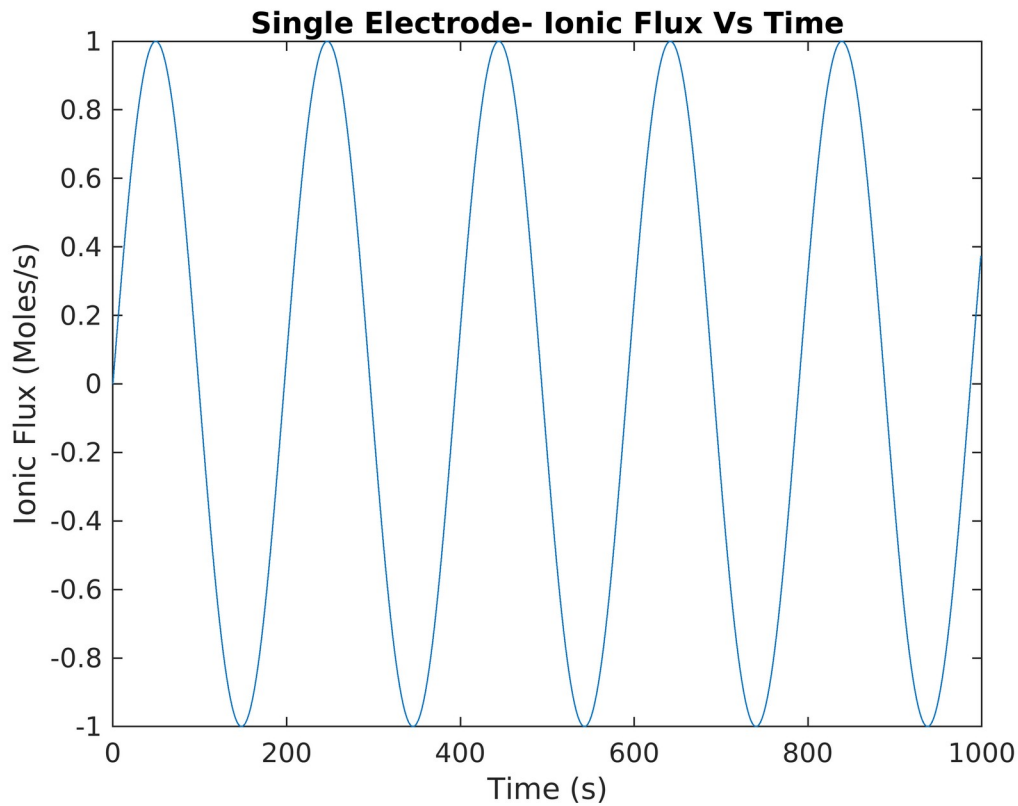
```
xlabel("Time (s)")
ylabel("Ionic Flux (Moles/s)")
```



```
%Sinusoidal Input
y=lithium_model(R, D, u_sin, y1_0, y2_0, dt, sigmaY);
plot(t, y)
title("Single Electrode- Surface Concentration Vs Time")
xlabel("Time (s)")
ylabel("Surface Concentration (Moles/m^3)")
```



```
plot(t,u_sin)
title("Single Electrode- Ionic Flux Vs Time")
xlabel("Time (s)")
ylabel("Ionic Flux (Moles/s)")
```



%Step 7:

As you insert charge into the battery, and electrons leave the anode, Lithium ions will convert to the ionic state and therefore increase the concentration of anode

%Step 8:

%Independent, Identically Distributed, Zero-mean, Gaussian Noise

%Constant Input

sigmaY = 0.1;

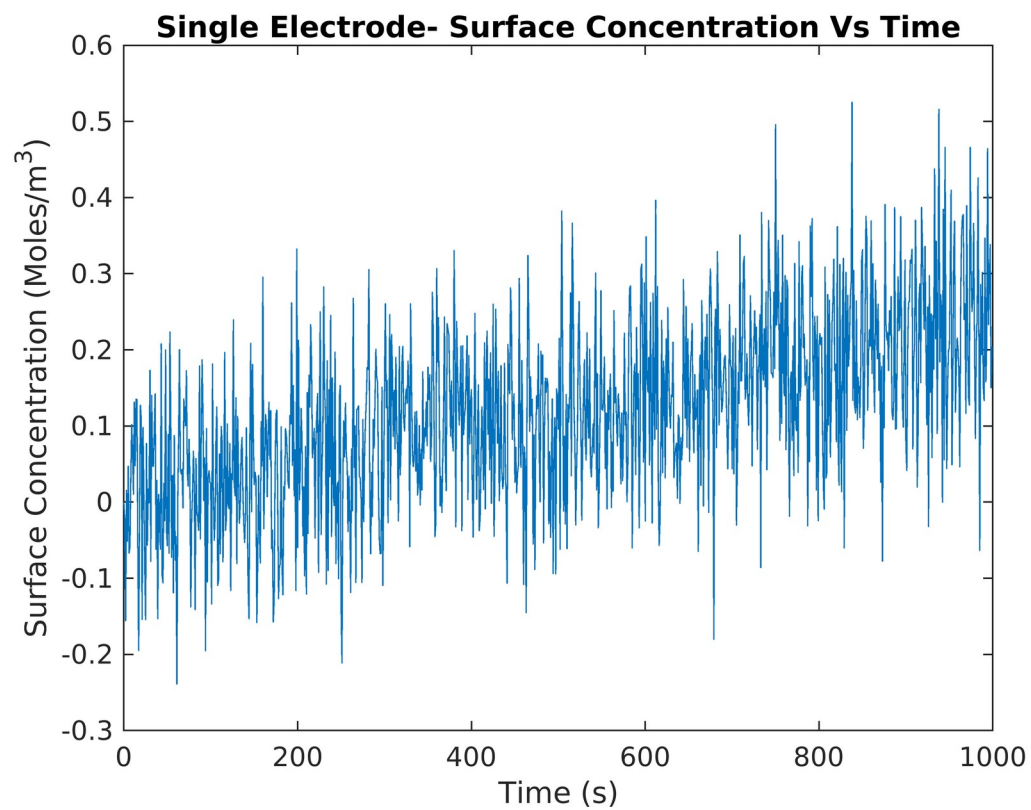
y_noise = lithium_model(R, D, u, y1_0, y2_0, dt, sigmaY);

plot(t,y_noise)

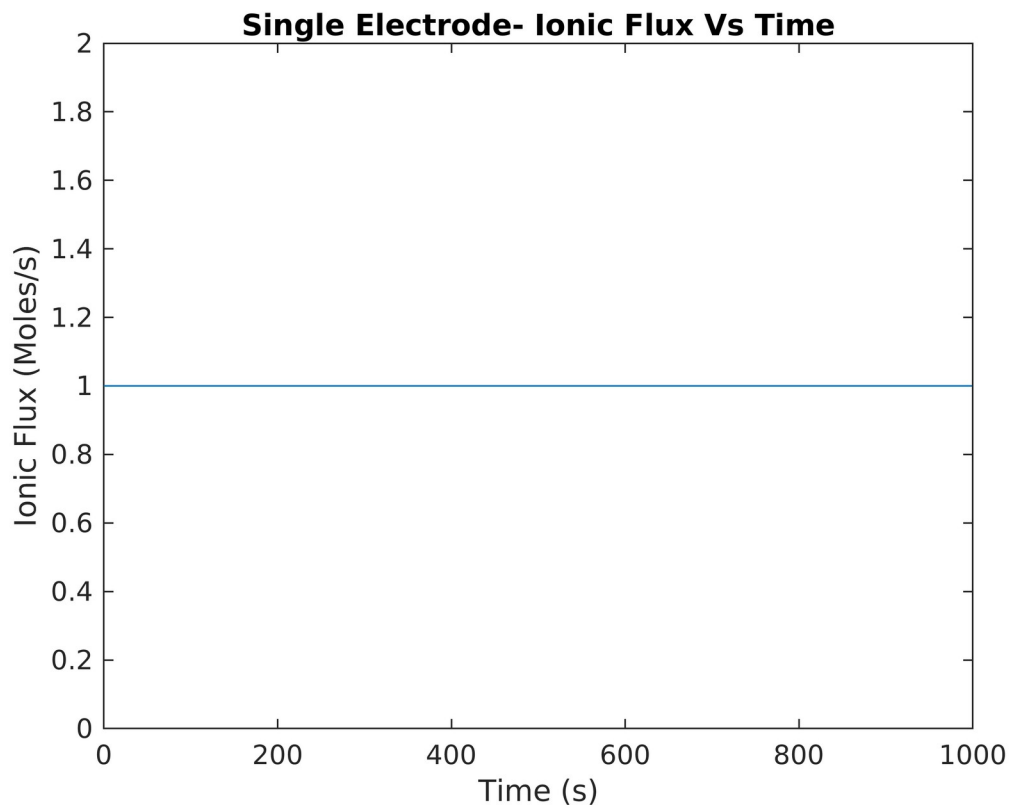
title("Single Electrode- Surface Concentration Vs Time")

xlabel("Time (s)")

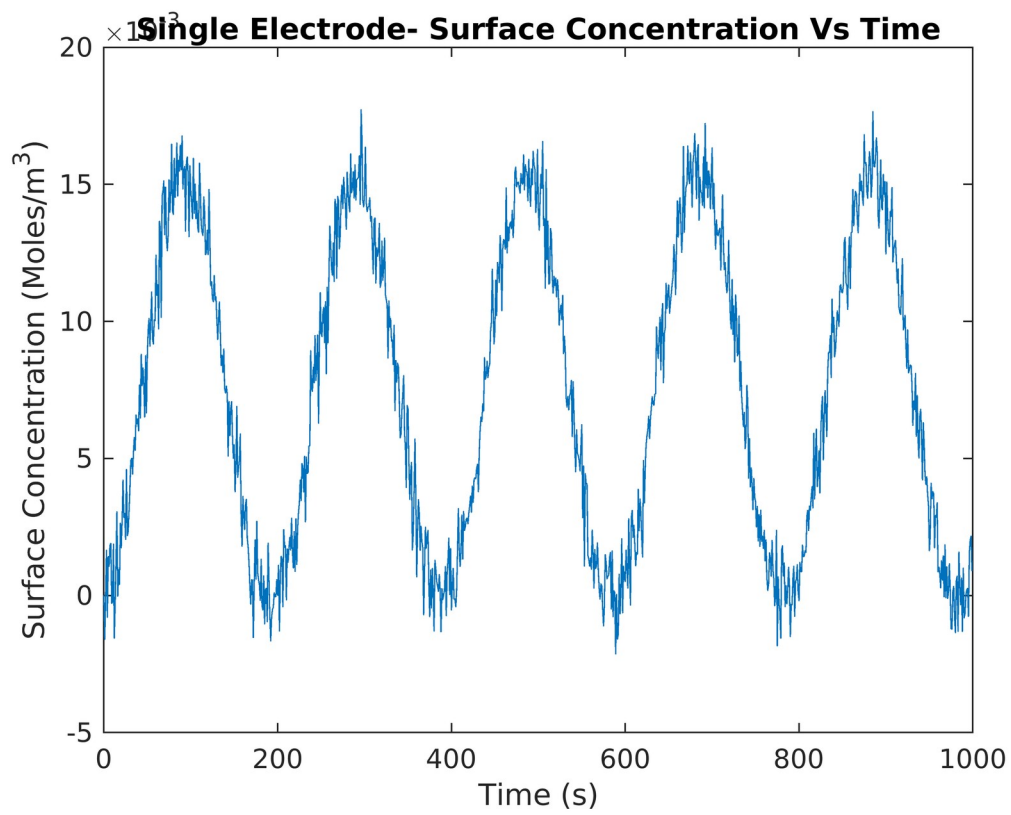
ylabel("Surface Concentration (Moles/m^3)")



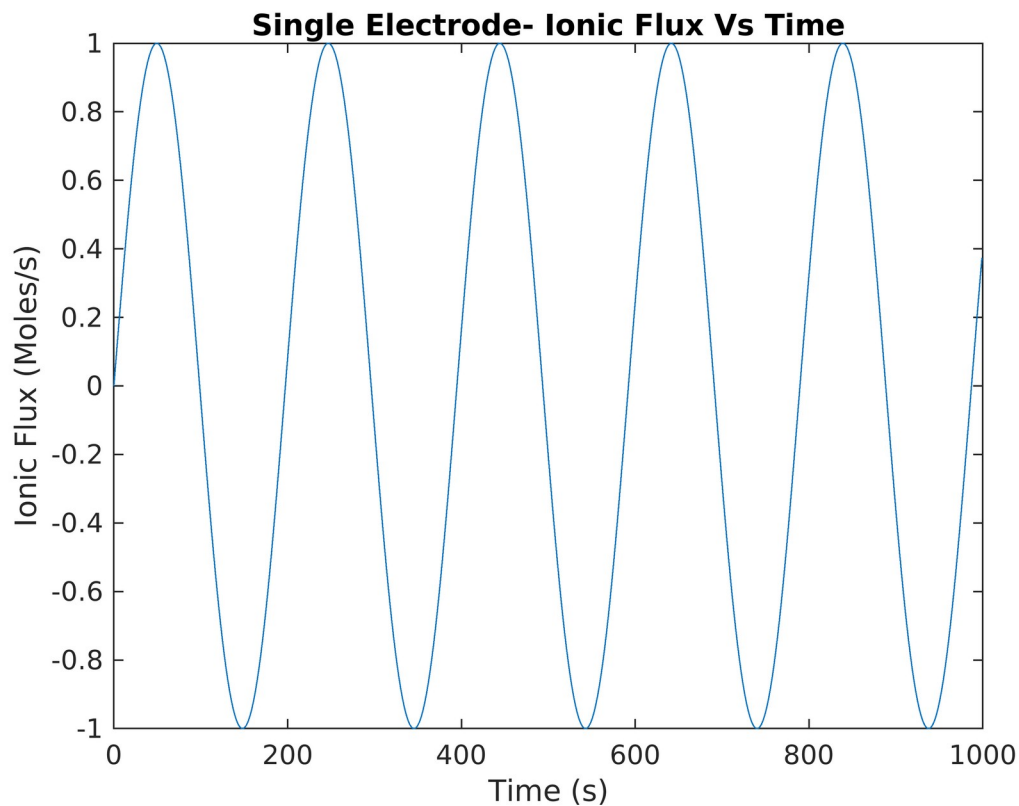
```
plot(t,u)
title("Single Electrode- Ionic Flux Vs Time")
xlabel("Time (s)")
ylabel("Ionic Flux (Moles/s)")
```



```
%Sinusoidal Input
sigmaY = 0.001;
y_noise = lithium_model(R, D, u_sin, y1_0, y2_0, dt, sigmaY);
plot(t,y_noise)
title("Single Electrode- Surface Concentration Vs Time")
xlabel("Time (s)")
ylabel("Surface Concentration (Moles/m^3)")
```



```
plot(t,u_sin)
title("Single Electrode- Ionic Flux Vs Time")
xlabel("Time (s)")
ylabel("Ionic Flux (Moles/s)")
```



```
%Step 9:
```

The `fminsearch` is able to converge to a reasonable value of R reliably, but D is unable to converge to the correct value.

```
clear all

Rtrue = 10;
Dtrue = 1;
N = 1000;
deltaT = 1;
sigmaY = 0.0005;
u = ones(N,1);
y1_0 = 0;
y2_0 = 0;

yhat = lithium_model(Rtrue, Dtrue, u, y1_0, y2_0, deltaT, sigmaY);

Rguess = 30;
Dguess = 8;
thetaGuess(1) = Rguess;
thetaGuess(2) = Dguess;

options = optimset('MaxFunEvals',100000,'TolX',1E-6,'TolFun',1E-6);
optimFunction = @(theta)(LSP(theta, Rtrue, Dtrue, u, N, deltaT, yhat, y1_0, y2_0, sigmaY));
```



```
varEstimation=fminsearch(optimFunction,thetaGuess,options)
```

```
Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 0.000561
varEstimation = 1x2
9.9730    4.5852
```

The fminsearch is able to converge to a reasonable value of R reliably, but D is unable to converge to the correct value. My parameter estimates appear to be biased as sigmaY is small I am unsatisfied with the estimated accuracy because D does not converge to the correct value

```
%Step 10:

nMonteCarlo= 500;

parfor i=1:nMonteCarlo
    optimFunction = @(theta) (LSP(theta, Rtrue, Dtrue, u, N, deltaT, yhat, y1_0, y2_0, s
    bestTheta=fminsearch(optimFunction,thetaGuess,options);
    RMonteCarlo(i) = bestTheta(1);
    DMonteCarlo(i) = bestTheta(2);
end
```

```
Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 8.170867
```

```
Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 3.974066
```

```
Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 0.000584
```

```
Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 6.846169
```

```
Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 0.000521
```

```
Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 0.000572
```

```
Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
```

Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 6.841349

Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 6.842117

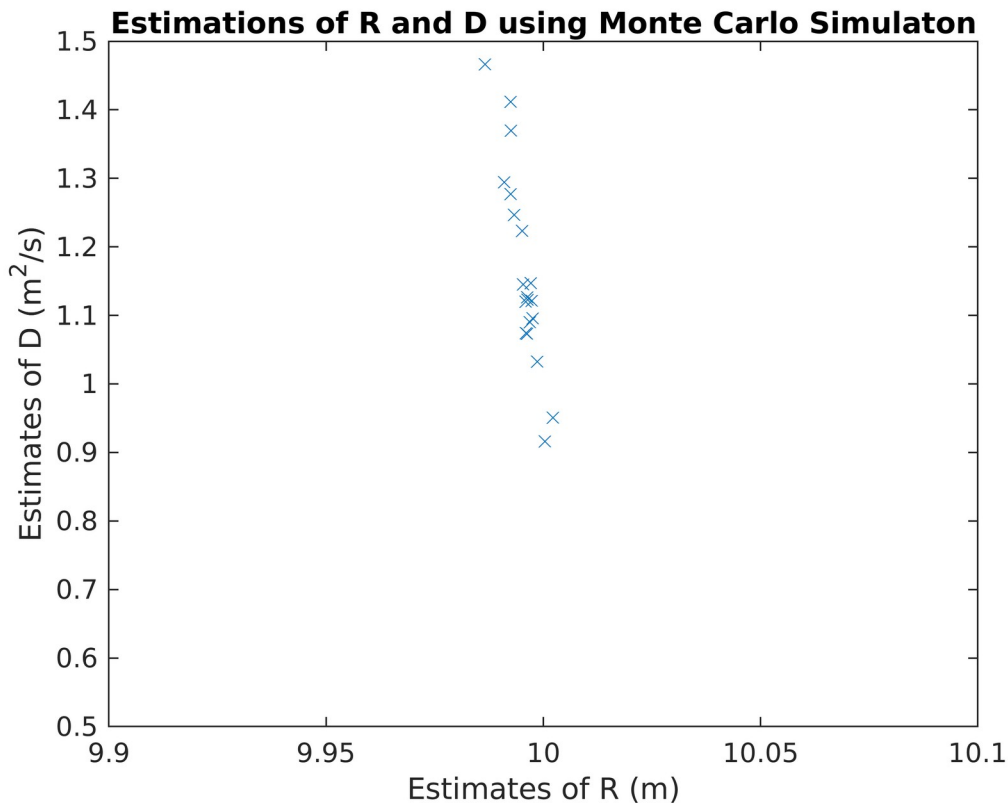
Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 0.000578

Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 0.000240

Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
Current function value: 0.000568

Exiting: Maximum number of iterations has been exceeded
- increase MaxIter option.
C...

```
plot(RMonteCarlo, DMonteCarlo, 'x')
title("Estimations of R and D using Monte Carlo Simulaton")
ylabel('Estimates of D (m^2/s)')
xlabel('Estimates of R (m)')
deltaD=.5;
deltaR=.1;
ylim([Dtrue-deltaD,Dtrue+deltaD]);
xlim([Rtrue-deltaR,Rtrue+deltaR]);
hold off
```



```

plot(RMonteCarlo, DMonteCarlo, 'x')
title("Estimations of R and D using Monte Carlo Simulaton")
xlabel('Estimates of R (m)')
ylabel('Estimates of D (m^2/s)')
hold on
yNominal = lithium_model(Rtrue, Dtrue ,u, y1_0, y2_0, deltaT, 0);
epsilon = 1e-3;

y1 = lithium_model(Rtrue*(1+epsilon), Dtrue, u, y1_0, y2_0, deltaT, 0);
y2 = lithium_model(Rtrue, Dtrue*(1+epsilon), u, y1_0, y2_0, deltaT, 0);

s1 = (y1-yNominal)/(Rtrue*epsilon);
s2 = (y2-yNominal)/(Rtrue*epsilon);

S= zeros(length(s1), 2);
S(:,1) = s1;
S(:,2) = s2;

F = (1/(sigmaY^2))*(S')*S;
CRLB = inv(F);
[V, D] = eig(CRLB);
lambda1 = D(1,1);
lambda2 = D(2,2);
v1 = V(:,1);
v2 = V(:,2);

xprime = linspace(-sqrt(lambda1),sqrt(lambda1));

```

```

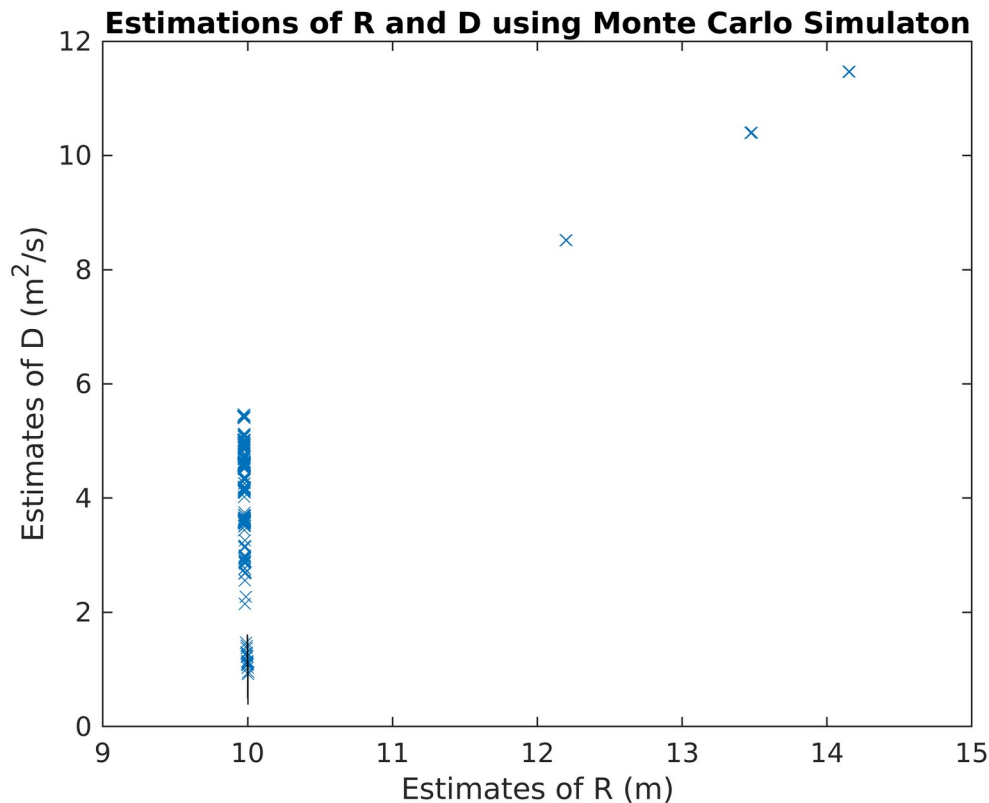
yprime = sqrt((1-(xprime.^2)/lambda1)*lambda2);

xypoints = [Rtrue;Dtrue]+v1*xprime+v2*yprime;
xypointsRef = [Rtrue;Dtrue]+v1*xprime-v2*yprime;
xy2points = [Rtrue;Dtrue]+2*v1*xprime+2*v2*yprime;
xy2pointsRef = [Rtrue;Dtrue]+2*v1*xprime-2*v2*yprime;
xy3points = [Rtrue;Dtrue]+3*v1*xprime+3*v2*yprime;
xy3pointsRef = [Rtrue;Dtrue]+3*v1*xprime-3*v2*yprime;

hold on

plot(xypoints(1,:), xypoints(2,:), 'k')
plot(xypointsRef(1,:), xypointsRef(2,:), 'k')
plot(xy2points(1,:), xy2points(2,:), 'k')
plot(xy2pointsRef(1,:), xy2pointsRef(2,:), 'k')
plot(xy3points(1,:), xy3points(2,:), 'k')
plot(xy3pointsRef(1,:), xy3pointsRef(2,:), 'k')

```



I am not satisfied with the results of the fisher information anylisis, as can be seen above, there is a portion of the estimations of R and D that converge to the correct value, and a portion that don't converge, then another portion that converge to arroneous values. the ellipsisies that are produced are unable to account for the outliers and the two distinct clusters of points.

```

function y = lithium_model(R, D, u, y1_0, y2_0, dt, sigmaY)
%Pre-Allocation

```

```

nSteps = length(u);
y1=zeros(nSteps,1);
y2=zeros(nSteps,1);
y=zeros(nSteps,1);

%Initial Conditions
y1(1)=y1_0;
y2(1)=y2_0;

%PADE Second Order Approximation
%Transfer Function Coefficients

m=-1/(4*pi*R^2*D);
a=(-3*D*m)/R;
b=-(2*m*R)/7;
c=R^2/(35*D);

%Foward Euler
for i=2:nSteps
    y1(i)=y1(i-1)+a*u(i-1)*dt;
    y2(i)=y2(i-1)+((b-c*a)/c)*u(i-1)*dt-(1/c)*y2(i-1)*dt;
    y(i)=(y1(i)+y2(i))+sigmaY*randn(1);
end
return
end

function error = LSP(theta, Rexact, Dexact, u, nSteps, dt, yhat, y1_0, y2_0, sigmaY)

nSteps = 1000;
R = theta(1);
D = theta(2);

yhat = lithium_model(Rexact, Dexact, u, y1_0, y2_0, dt, sigmaY);
y = lithium_model(R, D, u, y1_0, y2_0, dt, 0);

error = 0;

for i=1:nSteps
    error = error + (yhat(i)-y(i))^2;
end

end

```