

COMP5930M-Scientific Computation

Coursework 1

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1. A floating sphere

- (a). Nonlinear function and MATLAB implementation
- (b). The initial bracket plus justification for its validity
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2. Chemical engineering

- (a). Nonlinear system for the steady-state problem and MATLAB code
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- (b) Implement in MATLAB code
- (c) two correct solutions, figures
- (d) Implement in MATLAB code about the path and the tracing function, and the table of angles

1. A floating sphere

(a). Nonlinear function and MATLAB implementation

- The weight of water $W_w = \rho_w V_{cap} = \pi H^2(a - \frac{H}{3})$ and the weight of the sphere $W_s = \frac{4}{3} \pi \rho_s a^3$.
- Because of $W_w = W_s$, so the nonlinear equation $F_{a,\rho_s}(H) = W_w - W_s = 0$.
the nonlinear equation is down for given values of a and ρ_s :

$$F_{a,\rho_s}(H) = \pi H^2(a - \frac{H}{3}) - \frac{4}{3} \pi \rho_s a^3 = 0$$

- The implementation of MATLAB:

```

1. function f_h = fun_sphere(H, a, rhos)
2.     % The nonlinear equations Fa,phos (H) = 0
3.     % the weight of the water: W_w = pi*H.^2*(a - H/3)
4.     % the weight of the sphere: W_s = 4/3*pi*rhos*a.^3
5.     %
6.     % function f_h = fun_sphere(H, a, rhos)
7.     %
8.     % Input:
9.     %     H - a depth below the water surface
10.    %     a - the radius of sphere, constant value
11.    %     rhos- the density of sphere, constant value
12.    %
13.    % Output:
14.    %     f_h - final function value
15.    %
16.
17.    f_h = pi*H.^2*(a - H/3) - 4/3*pi*rhos*a.^3;
18.
19. end

```

(b). The initial bracket plus justification for its validity

- the initial bracket: $[0 + \delta, 3a - \delta], \delta = 10^{-6}$

Because the H_L and H_R must meet the following conditions:

- (1). $0 < H < 3a$, due to the $W_w > 0$
- (2). $F(x_L)F(x_R) \leq 0$
- (3). The $F(H)$ can be get a maximum at the point $2a$, So, the suitable value can be $[0 + \delta, 2a], \delta = 10^{-6}$ OR $[2a, 3a - \delta], \delta = 10^{-6}$ must be find a H to fit the $F(H) = 0$.

(c). Two correct solutions plus the number of iterations

- In given values $a = 1$ and $p_s = 0.45$,
- In Bisection method, the initial bracket $[10^{-6}, 3 - 10^{-6}]$, the solution is :
 $H = 0.9332$ and the number of iteration is 22.

- In Newton Method, the init guess $x_0 = 1.0$, the solution is : $H = 0.9332$ and the number of iteration is **3**.
- As the following figure:

```
>> bisection(@fun_sphere, 1e-6, 2-1e-6, 1e-6, 100)
  i   x_i          |F(x_i)|
  0   1.00000000    2.09e-01
  1   0.50000050    1.23e+00
  2   0.75000025    5.60e-01
  3   0.87500013    1.81e-01
  4   0.93750006    1.33e-02
  5   0.90625009    8.42e-02
  6   0.92187508    3.55e-02
  7   0.92968757    1.11e-02
  8   0.93359382    1.12e-03
  9   0.93164069    4.98e-03
 10   0.93261725    1.93e-03
 11   0.93310554    4.02e-04
 12   0.93334968    3.61e-04
 13   0.93322761    2.04e-05
 14   0.93328864    1.70e-04
 15   0.93325812    7.51e-05
 16   0.93324286    2.73e-05
 17   0.93323524    3.47e-06
 18   0.93323142    8.46e-06
 19   0.93323333    2.50e-06
 20   0.93323428    4.85e-07
 21   0.93323428    4.85e-07
```

ans =

列 1 至 8

1.0000	0.5000	0.7500	0.8750	0.9375	0.9063	0.9219	0.9297
--------	--------	--------	--------	--------	--------	--------	--------

列 9 至 16

0.9336	0.9316	0.9326	0.9331	0.9333	0.9332	0.9333	0.9333
--------	--------	--------	--------	--------	--------	--------	--------

列 17 至 21

0.9332	0.9332	0.9332	0.9332	0.9332
--------	--------	--------	--------	--------

 >>

```
>> newtonSys(@fun_sphere, @dfun_sphere, 1.0, 1e-6, 100)
  k   |F(x_k)|
  0     0.209
  1    0.00031
  2    2.06e-09
```

Converged

ans =

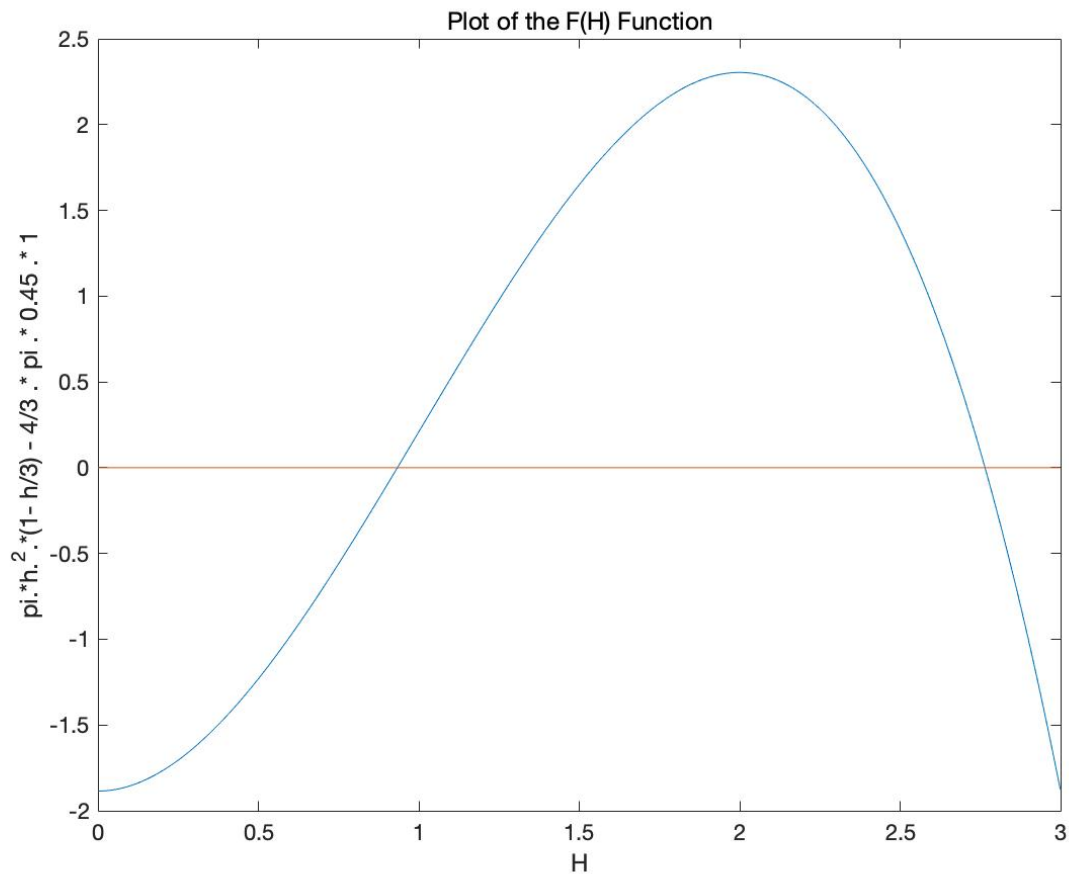
0.9332

 >>

(d). number of solutions and corresponding initial guesses

- By changing the initial guess x_0 , can be get **2** solutions, the first one is:

- $H = 0.9332$, the second one is $H = 2.7645$.



- if the initial guess $x_0 = 0.5$, the solution is : **0.9332**.

```
>> newtonSys(@fun_sphere, @dfun_sphere, 0.5, 1e-6, 100)
k   |F(x_k)|
0    1.23
1    0.279
2    0.000184
3    7.28e-10
Converged
ans =
    0.9332
```

```
>> |
```

- if the initial guess $x_0 = 2.5$, the solution is : **2.7645**.

```

>> newtonSys(@fun_sphere, @dfun_sphere, 2.5, 1e-6, 100)
k   |F(x_k)|
0   1.39
1   0.635
2   0.0395
3   0.000192
4   4.66e-09
Converged

ans =

    2.7645
fx >> |

```

2. Chemical engineering

(a). Nonlinear system for the steady-state problem and MATLAB code

For the case $n=5$ reactors, the nonlinear equation system $F(U) = 0$.

```

1. function f = fun_chemical(a, V, G, k, a0)
2.
3.     % The nonlinear equation system F(U)
4.     % function f = fun_chemical(a, V, G, k, a0)
5.     % Input:
6.     %     a - current solution
7.     %     V - constants
8.     %     G - constants
9.     %     k - constants
10.    %     a0 - constants
11.    % Output:
12.    %     f = final function value
13.    f = [(k*V/G)*a(1).^2 - a0 + a(1);
14.         (k*V/G)*a(2).^2 - a(1) + a(2);
15.         (k*V/G)*a(3).^2 - a(2) + a(3);
16.         (k*V/G)*a(4).^2 - a(3) + a(4);
17.         (k*V/G)*a(5).^2 - a(4) + a(5)
18.         ];
19. end

```

(b). Jacobian matrix for the steady-state problem and MATLAB code

the Jacobian matrix

```
1. function jf = Jfun_chemical(a, V, G, k, a0)
2.     % The nonlinear equation system F(U)
3.     % function f = fun_chemical(a, V, G, k, a0)
4.     % Input:
5.     %     a - current solution
6.     %     V - constants
7.     %     G - constants
8.     %     k - constants
9.     %     a0 - constants
10.    % Output:
11.    %     jf = Jacobian matrix
12.    jf = [
13.        (k*V/G)*2*a(1) + 1, 0, 0, 0, 0;
14.        -1, (k*V/G)*2*a(2) + 1, 0, 0, 0;
15.        0, -1, (k*V/G)*2*a(3) + 1, 0, 0;
16.        0, 0, -1, (k*V/G)*2*a(4) + 1, 0;
17.        0, 0, 0, -1, (k*V/G)*2*a(5) + 1
18.    ];
19. end
```

(c). Numerical solution and values of $|F|$

- In case the $n = 5$ with the following the parameters: $V = 1.0$, $G = 35.0$, $k=0.6$, $a0=6.0$.
- The values of $|F|$ at each iterations:

iterations: $|f(x)|$

0 6.000000

1 1.379973

2 0.078314

3 0.000231

4 0.000000

5 0.000000

the solutions:

5.4844

5.0476

4.6732

4.3490

4.0656

(d). Discretised time-dependent equations

- the time-dependent concentrations $a_i(t)$.
- (1). $\frac{da_i}{dt} = -\beta a_i^2 + a_{i-1} - a_i$, for $t > 0$; and $a_i = a(t_i)$
- Implicit Euler approximation for (2). $\frac{da_i}{dt} = f(u)$
- and the backward Euler(implicit) method is: (3). $\frac{a_{i+1} - a_i}{\Delta t} = f(u_{i+1})$
- So we can get this formal:
- $F_i(U) = \frac{a_{i+1} - a_i}{\Delta t} + \beta a_{i+1}^2 + a_{i+1} - a_i = 0$ given an initial $a_i(0) = a_i^0$ for $i = 1, \dots, n$.

(e). Jacobian matrix for the time-dependent problem, behaviour for limit values of times step

- the Jacobian matrix is :

$$\begin{bmatrix} \frac{\partial F_i}{\partial a_i} \\ \frac{\partial F_i}{\partial a_{i+1}} \end{bmatrix} = \begin{bmatrix} -1 - \frac{1}{\Delta t} \\ 1 + \frac{1}{\Delta t} + 2\beta a_{i+1} \end{bmatrix}$$

- when $(\Delta t \rightarrow 0)$, the Jacobian matrix is trend to $[-1; \infty]$, this reactor is unsteady-state.
- when $(\Delta t \rightarrow \infty)$, and this reactor is steady-state.

3. Control of a robot arm

(a) System of equations

- The system of nonlinear equations in the form $\mathbf{F}(\mathbf{x}) = \mathbf{0}$.
- Base on the equations of the location of the free end (loc_x, loc_y) and $(x_1, x_2) = (\theta, \phi)$:

$$loc_x = \cos(\theta) + \cos(\phi)$$

$$loc_y = \sin(\theta) + \sin(\phi)$$

- So, \mathbf{x} is the vector $\{x_1, x_2\}$, \mathbf{F} is a set $\{F_1(\mathbf{x}), F_2(\mathbf{x})\}$ nonlinear equations:

$$F_1(x_1, x_2) = \cos(x_1) + \cos(x_2) - loc_x$$

$$F_2(x_1, x_2) = \sin(x_1) + \sin(x_2) - loc_y$$

- Find (x_1^*, x_2^*) such that $F_1(x_1^*, x_2^*) = 0$ and $F_2(x_1^*, x_2^*) = 0$.

$$\cos(x_1) + \cos(x_2) - loc_x = 0$$

$$\sin(x_1) + \sin(x_2) - loc_y = 0$$

(b) Implement in MATLAB code

the matlab filename: **fun_arm.m**

```

1. function f = fun_arm(x, locx, locy)
2.     % Systems of nonlinear equations for control of a ro
    bot arm
3.     % system of 2 nonlinear equations
4.     % function f_x = fun_arm(x, locx, locy)
5.     %
6.     % Input: x - current solution
7.     %         locx - current location x
8.     %         locy - current location y
9.     %
10.    % Output: f - final function value
11.
12.    f = [cos(x(1)) + cos(x(2)) - locx;
13.         sin(x(1)) + sin(x(2)) - locy];
14. end

```

The Jacobian function: **Jfun_arm(x)**:

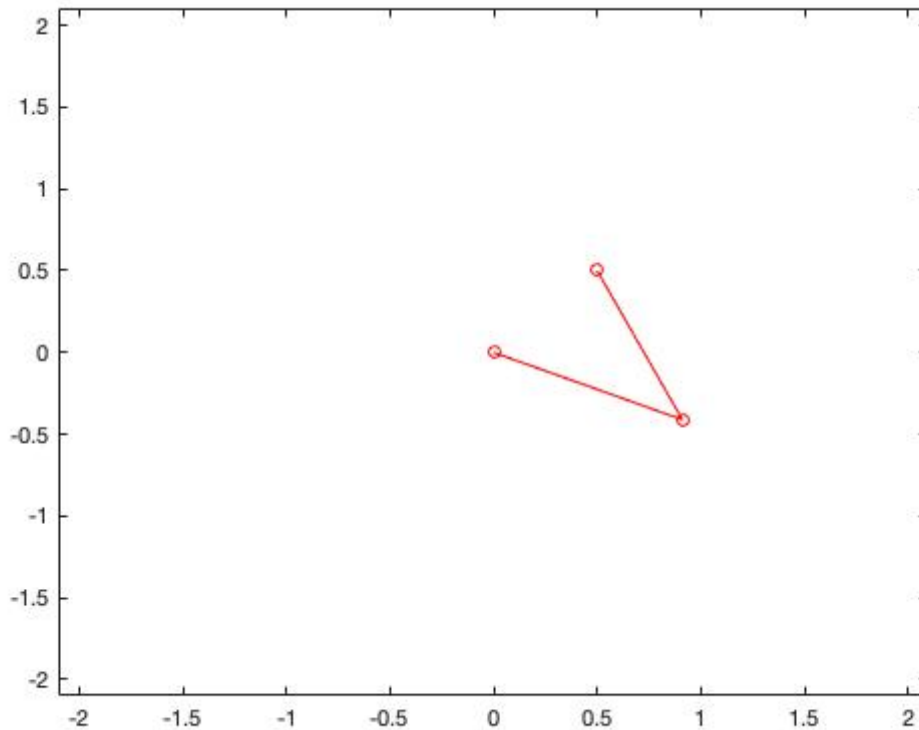
```

1. function jf = Jfun_arm(x)
2.     % Analytical Jacobian
3.     % function jf = trueJacobian(x)
4.     %
5.     % Input:
6.     %         x - current solution
7.     %
8.     %
9.     % Output: jf - Jacobian matrix
10.
11.    jf = [-sin(x(1)), -sin(x(2)); cos(x(1)), cos(x(2))];
12.
13. end

```

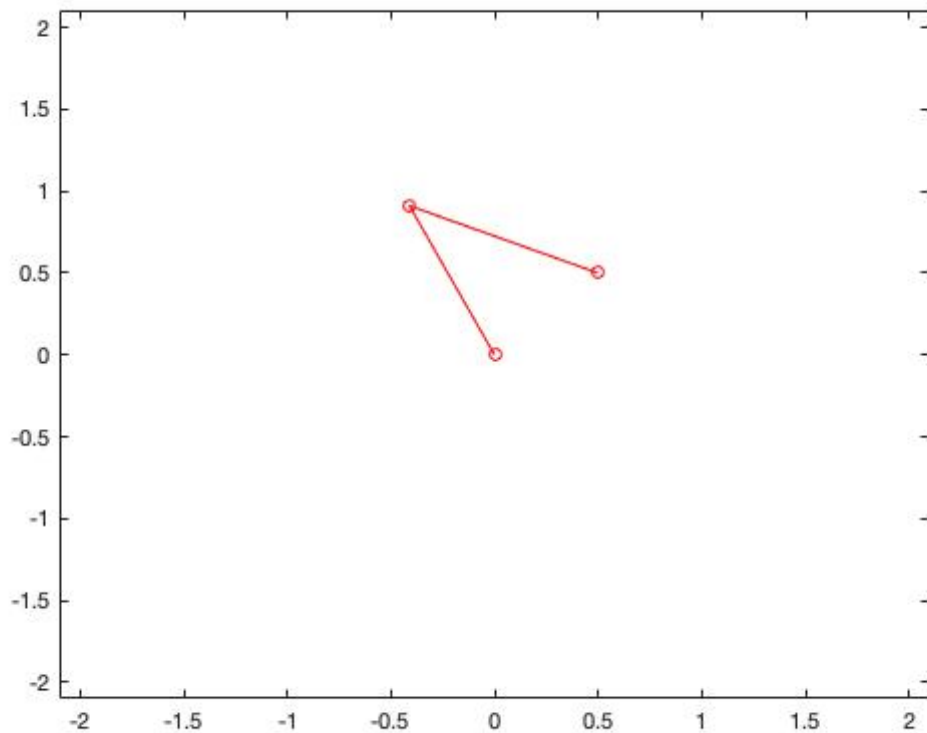

(c) two correct solutions, figures

- **Case i**, the initial x_0 for Newton method, when x_0 at location $(x_1, x_2) = (-1, 1)$
the solution is: $(\theta, \phi) = (x_1, x_2) = (-0.4240, 1.9948)$
the figure:



- **Case ii**, the initial x_0 for Newton method, when x_0 at location $(x_1, x_2) = (2, 0)$
the solution is: $(\theta, \phi) = (x_1, x_2) = (1.9948, -0.4240)$

the figure:



(d) Implement in MATLAB code about the path and the tracing function, and the table of angles

1. the traceFn is the function defining the path. the code: filename: **traceFn.m**

```
1. function [locx, locy] = traceFn(t)
2.     % traceFn is the function defining path
3.     % of the free end of the arm.
4.     % trace path defined by the following function
5.     % (x,y)=(t, 0.1+sin(2t-0.5)), and (-1<=t<= 1)
6.     %
7.     % function [locx, locy] = traceFn(t)
8.     % Input: t - param variables, from t=-1 to t=1
9.     %
10.    % Output: locx - the location point x
11.    %          locy - the location point y
12.
13.    locx = t;
14.    locy = 0.1 + sin(2*t - 0.5);
15.
```

16. **end**

2. The matlab filename: “**traceArm.m**”


```
1. function t_out = traceArm(traceFn,nSteps, x0)
2.
3.     % Trace path of the free end of arm at initial x0
4.     % function t_out = traceArm(traceFn,nSteps, x0)
5.     % Input:
6.     %     traceFn -to create the location (x,y)
7.     %     nSteps -the number of steps for t splitting the
      t
8.     %     x0      -the init guess value for Newton method
9.     % Output:
10.    %     t_out   - a table to recording the angles
11.    %     theta and phi at each step.
12.    step = 2 / (nSteps -1);
13.    i = 1;
14.
15.    step_list = [];
16.    t = [];
17.    x_list = [];
18.    y_list = [];
19.    theta_list = [];
20.    phi_list = [];
21.    for st = -1:step:1
22.        [locx, locy] = feval(traceFn, st)
23.        [xx, f] = newtonSys2(@fun_arm, @Jfun_arm, x0, 1e
-10, 100,locx,locy);
24.        t = [t st];
25.        step_list = [step_list i];
26.        x_list = [x_list locx];
27.        y_list = [y_list locy];
28.        theta_list = [theta_list xx(1,:)];
29.        phi_list = [phi_list xx(2,:)];
30.        i = i + 1;
31.    end
32.    nSteps = step_list';
33.    nT = t';
34.    nLocx = x_list';
```

```

35.     nLocy = y_list';
36.     Theta = theta_list';
37.     Phi    = phi_list';
38.     t_out = table(nSteps,nT, nLocx, nLocy,Theta,Phi);
39.
40. end

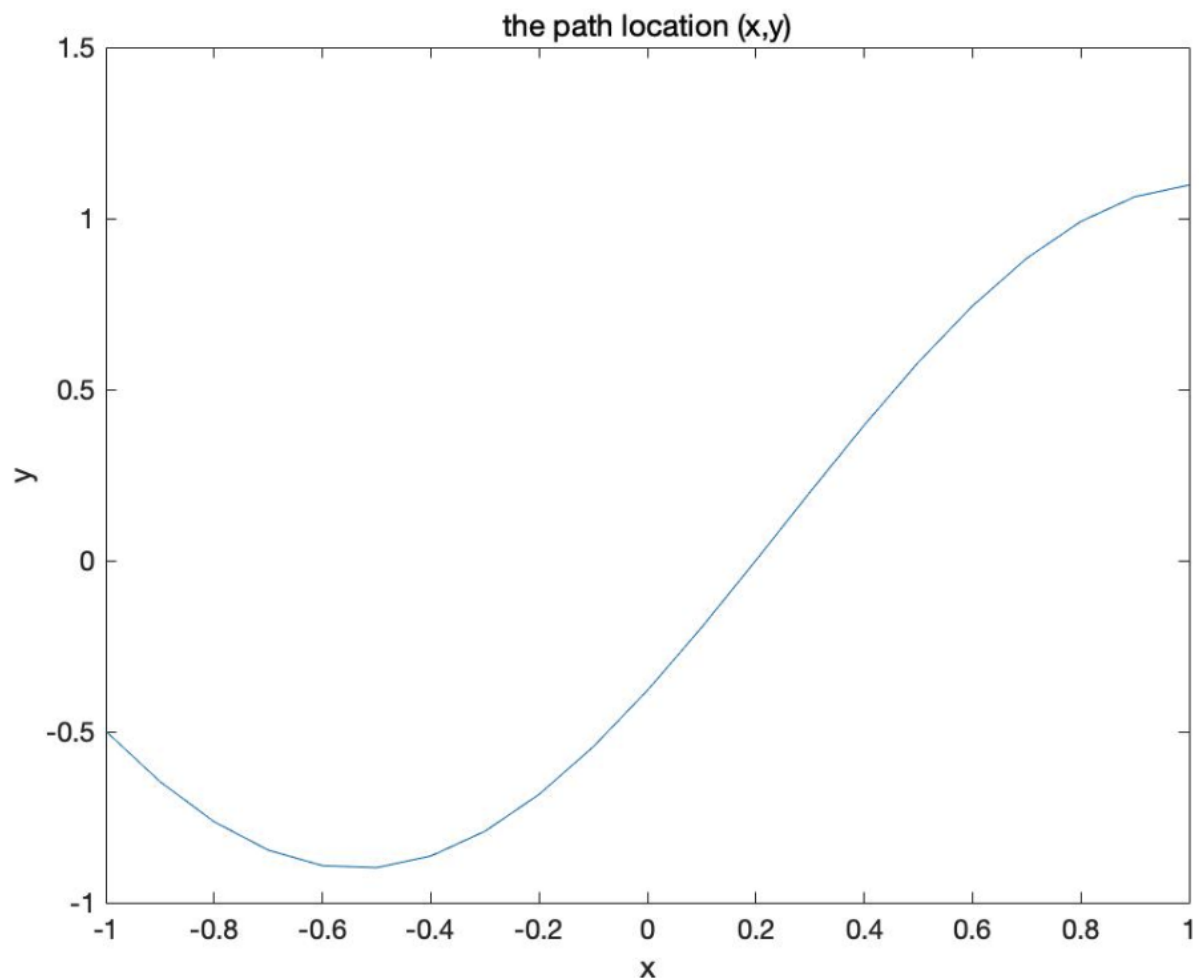
```

3. In case the \mathbf{x}_0 : (0.5, 0.5), the θ and ϕ at each step, following down:

 21x6 table

	1 nSteps	2 nT	3 nLocx	4 nLocy	5 Theta	6 Phi	7
1	1	-1	-1	-0.4985	-1.7012	2.6260	
2	2	-0.9000	-0.9000	-0.6457	-1.5354	2.7801	
3	3	-0.8000	-0.8000	-0.7632	48.8708	53.1839	
4	4	-0.7000	-0.7000	-0.8463	-3.2513	-1.2724	
5	5	-0.6000	-0.6000	-0.8917	-1.1596	3.1166	
6	6	-0.5000	-0.5000	-0.8975	3.1728	5.2354	
7	7	-0.4000	-0.4000	-0.8636	-7.2129	-3.0794	
8	8	-0.3000	-0.3000	-0.7912	-0.7993	3.2160	
9	9	-0.2000	-0.2000	-0.6833	-0.6487	3.2208	
10	10	-0.1000	-0.1000	-0.5442	-0.4620	3.2402	
11	11	0	0	-0.3794	3.3325	6.0923	
12	12	0.1000	0.1000	-0.1955	28.8571	31.7787	
13	13	0.2000	0.2000	1.6658e-04	-1.4698	1.4715	
14	14	0.3000	0.3000	0.1998	-0.8020	1.9772	
15	15	0.4000	0.4000	0.3955	-0.5059	2.0655	
16	16	0.5000	0.5000	0.5794	-0.3193	2.0370	
17	17	0.6000	0.6000	0.7442	-0.1802	1.9647	
18	18	0.7000	0.7000	0.8833	-0.0715	1.8728	
19	19	0.8000	0.8000	0.9912	0.0114	1.7721	
20	20	0.9000	0.9000	1.0636	0.0684	1.6686	
21	21	1	1	1.0975	0.0977	1.5660	
22							

The path of free end of arm.



The newtonSys2 function

```
1. function [ x,f ] = newtonSys2( fnon, fjac, x0, tol, maxI
   t,locx,locy)
2.
3. % Basic Newton algorithm for systems of nonlinear equati
   ons
4. % function [ x,f ] = newtonSys( fnon, fjac, x0, tol, max
   it )
5. % Input: fnon - function handle for nonlinear system
6. %         fjac - function handle for Jacobian matrix
7. %         x0 - initial state (column vector)
8. %         tol - convergence tolerance
9. %         maxIt - maximum allowed number of iterations
10. % Output: x - final point
11. %         f - final function value
12.
13. fprintf(' x      |f(x)|\n')
```

```

14.
15. n = length(x0);
16.
17. x = x0; % initial point
18. f = feval(fnon,x,locx, locy); % initial function values
19. normf = norm(f);
20. it = 0;
21. fprintf(' %d %12.6f\n',it,norm(f));
22.
23. while (normf>tol) && (it<maxIt)
24.
25.     J = feval(fjac, x); % build Jacobian
26.     delta = -J\f; % solve linear system
27.
28.     xkp = x;
29.     xk = x + delta;
30.
31.     x = x + delta; % update x
32.     f = feval(fnon,x, locx, locy); % new function values
33.     normf = norm(f);
34.
35.     it = it + 1;
36.     % Print the new estimate and function value.
37.     fprintf(' %d %12.6f\n',it,normf)
38.
39.     %plot([xkp(1) xk(1)], [xkp(2) xk(2)], 'ok-', 'LineWidth
    h', 2)
40.     %pause
41. end
42.
43. if( it==maxIt)
44.     fprintf(' WARNING: Not converged\n')
45. else
46.     fprintf(' SUCCESS: Converged\n')
47. end

```