
Lab 6

David Merrick

```
% 1.
% a.
lsq(1)
lsq(2)
lsq(3)

% Degree 1: Max error = 0.1548
% Degree 2: Max error = 0.0150
% Degree 3: Max error = 0.0011

% b.
lsq(4)
lsq(5)
lsq(6)

% Degree 4: Max error = 5.8444e-05
% Degree 5: Max error = 1.9226e-04
% Degree 6: Max error = 0.0015

% Max error stops decreasing at degree 5.

% c.
lsq(7)
lsq(8)
lsq(9)

% At degree 9, there are visible errors in the approximation.

% d.
lsq(10)
lsq(11)
lsq(12)

% maxerr on degree 10: 2.3753.
% maxerr on degree 11: 21.8856.
% maxerr on degree 12: 49.4773.
% The maximum error appears to be increasing exponentially.

% e.
% Polynomial approximations of  $e^x$  of degree 1-5 are more accurate than
% higher-degree approximations of this function. This does not necessarily
% mean that,
% in general, low-degree polynomials are better at approximating functions
% than high-degree polynomials. We have only tested this function.
% It might be possible to generate a table of values and use Chebyshev
% nodes to create a high-degree interpolating polynomial that is more
% accurate of an approximation and does not result in an ill-conditioned
% matrix.
```

```
% f.

% The 4x4 Hilbert matrix is not diagonally dominant, a prerequisite for
% Jacobian convergence, so this should not converge using Jacobian.

m=4
A=hilb(m+1);
N=diag(diag(A));
P=N-A;
norm(inv(N)*P)

m=10
A=hilb(m+1);
N=diag(diag(A));
P=N-A;
norm(inv(N)*P)

m=30
A=hilb(m+1);
N=diag(diag(A));
P=N-A;
norm(inv(N)*P)

m=40
A=hilb(m+1);
N=diag(diag(A));
P=N-A;
norm(inv(N)*P)

m=100
A=hilb(m+1);
N=diag(diag(A));
P=N-A;
norm(inv(N)*P)

m=200
A=hilb(m+1);
N=diag(diag(A));
P=N-A;
norm(inv(N)*P)

% For m=4, result is 4.6738. For m=10, result is 11.2735. For m=30, result
% is 33.2966. For m=40, result is 44.3076. For m=70, result is 77.395. For
% m=100, result is 110.3709. For m=200, result is 270.2244. As predicted,
% this does not converge using Jacobian iteration, so this method could not
% be applied to this linear system.

% g.
% Gauss-Seidel can be applied to any matrix with non-zero elements on the
% diagonals, but convergence is only guaranteed if the matrix is either
% diagonally dominant,
% or symmetric and positive definite. The Hilbert matrix is not
% diagonally dominant, but is symmetric, so it would be reasonable to
```

```
% predict that this will converge for the Hilbert matrix.

m=4
A=hilb(m+1);
N=tril(A);
P=N-A;
norm(inv(N)*P)

m=10
A=hilb(m+1);
N=tril(A);
P=N-A;
norm(inv(N)*P)

m=30
A=hilb(m+1);
N=tril(A);
P=N-A;
norm(inv(N)*P)

m=40
A=hilb(m+1);
N=tril(A);
P=N-A;
norm(inv(N)*P)

m=70
A=hilb(m+1);
N=tril(A);
P=N-A;
norm(inv(N)*P)

m=100
A=hilb(m+1);
N=tril(A);
P=N-A;
norm(inv(N)*P)

m=200
A=hilb(m+1);
N=tril(A);
P=N-A;
norm(inv(N)*P)

% For m=4, result is 1.0003. For m=10, result is 1.0015. For m=30, result
% is 1.0032. For m=40, result is 1.0035. For m=70, result is 1.0039. For
% m=100, result is 1.0041. For m=200, result is 1.0043. Clearly, as
% predicted, this does converge using Gauss-Seidel Iteration, so this
% method could be applied to this linear system.

% 2.
% a.
n=2
gaussweights(n)
```

```

n=4
gaussweights(n)

n=8
gaussweights(n)

% For n = 2,4,8, the calculated values match the values in Table 5.7 in the
% book.
% b.
n=32
gaussweights(n)

n=64
gaussweights(n)

% For n=32, the condition number is 1.2827e+12. For n=64, the condition
% number is 1.2338e+20. The condition number relates the reative accuracy
% of the input to the relative accuracy of the output. With a high
% condition number, small changes in the input can produce huge errors,
% and this is clearly demonstrated on the plot.

% c.
% Convergence for Jacobian iteration requires that the matrix is
% diagonally-dominant. With the Vandermonde matrix, this is not the case.
% So Jacobion iteration will not converge for this linear system.

% d.
% Convergence for Gauss-Seidel requires that the matrix is either
% diagonally-dominant, symmetrical, or positive definite. The Vandermonde
% matrix meets none of these criteria, so that method will not converge for
% this linear system.

maxerr =

    0.1548

ans =

    0

maxerr =

    0.0150

ans =

    1.1102e-16

```

maxerr =

0.0011

ans =

0

maxerr =

5.8444e-05

ans =

2.5438e-16

maxerr =

1.9226e-04

ans =

4.0792e-16

maxerr =

0.0015

ans =

3.2841e-16

maxerr =

0.0048

ans =

2.3009e-15

maxerr =

0.0051

ans =
 $1.9681e-15$

maxerr =
 0.1158

ans =
 $2.0337e-12$

maxerr =
 2.3753

ans =
 $1.7099e-10$

maxerr =
 21.8856

ans =
 $4.5044e-09$

maxerr =
 49.4773

ans =
 $8.5473e-08$

m =
 4

ans =
 4.6738

$m =$

10

$ans =$

11.2735

$m =$

30

$ans =$

33.2966

$m =$

40

$ans =$

44.3076

$m =$

100

$ans =$

110.3709

$m =$

200

$ans =$

220.4744

$m =$

4

ans =

1.0003

m =

10

ans =

1.0015

m =

30

ans =

1.0032

m =

40

ans =

1.0035

m =

70

ans =

1.0039

m =

100

ans =

1.0041

m =

200

ans =

1.0043

n =

2

ans =

2.7321

ans =

1 1

n =

4

ans =

21.9310

ans =

0.3479 0.6521 0.6521 0.3479

n =

8

ans =

798.9577

ans =

Columns 1 through 7

0.1012 0.2224 0.3137 0.3627 0.3627 0.3137 0.2224

Column 8

0.1012

n =

32

ans =

1.2827e+12

ans =

Columns 1 through 7

0.0070 0.0163 0.0254 0.0343 0.0428 0.0510 0.0587

Columns 8 through 14

0.0658 0.0723 0.0782 0.0833 0.0877 0.0912 0.0938

Columns 15 through 21

0.0956 0.0965 0.0965 0.0956 0.0938 0.0912 0.0877

Columns 22 through 28

0.0833 0.0782 0.0723 0.0658 0.0587 0.0510 0.0428

Columns 29 through 32

0.0343 0.0254 0.0163 0.0070

n =

64

*Warning: Matrix is close to singular or badly scaled. Results may be inaccurate.
RCOND = 7.132669e-21.*

ans =

1.2338e+20

ans =

Columns 1 through 7

0.0572 -0.1622 0.2423 -0.1803 -0.0105 0.4171 -0.8978

Columns 8 through 14

1.4791 -1.9049 2.2038 -2.1179 1.8182 -1.1832 0.6059

Columns 15 through 21

-0.0786 0.0529 -0.3540 1.1675 -1.9366 2.6211 -2.6597

Columns 22 through 28

2.3834 -1.7498 1.5161 -1.5566 2.0070 -1.8302 0.6451

Columns 29 through 35

2.2147 -5.8144 9.2520 -10.7129 9.7656 -6.2526 1.7344

Columns 36 through 42

2.6331 -5.2237 5.9092 -4.5707 2.3355 0.2944 -2.2814

Columns 43 through 49

3.6344 -3.9850 3.8568 -3.1967 2.5699 -1.8479 1.4037

Columns 50 through 56

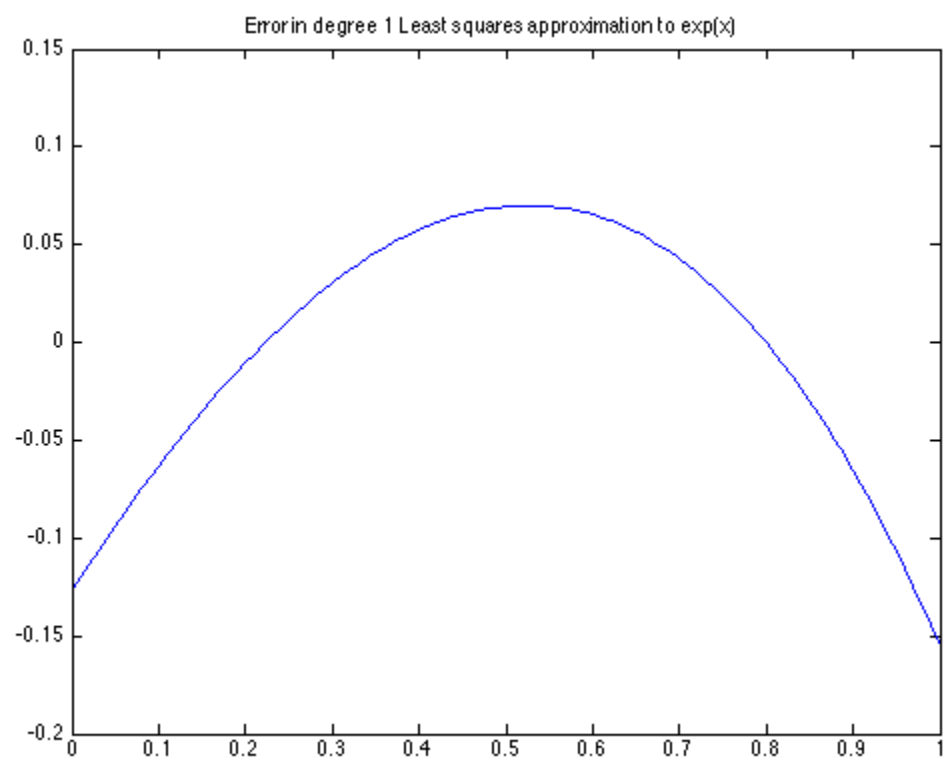
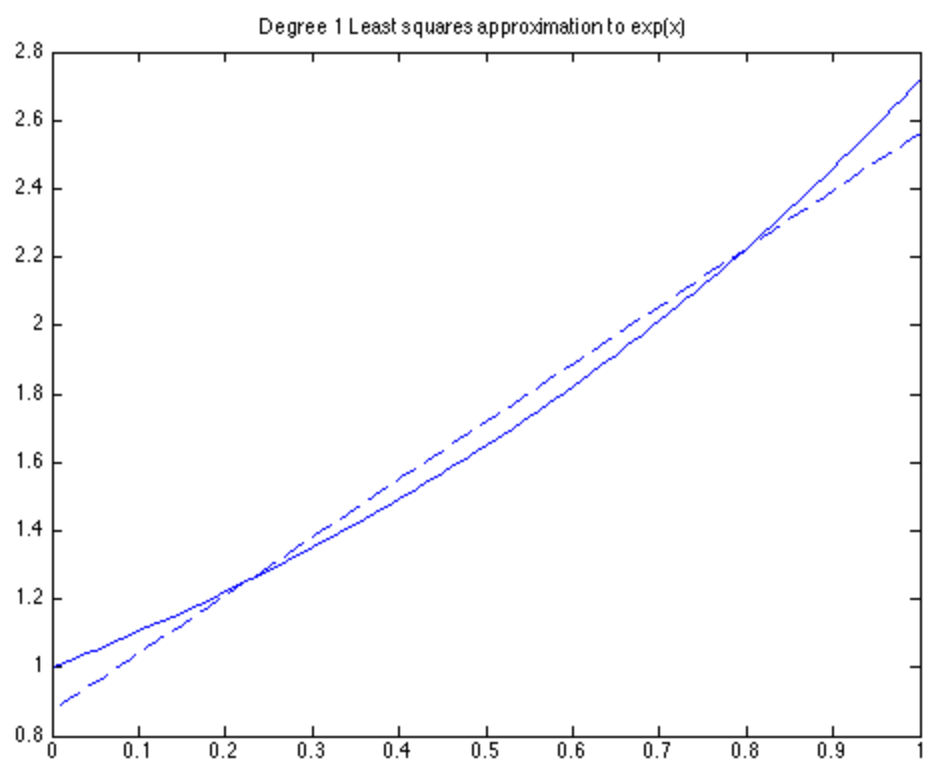
-0.9705 0.7783 -0.5307 0.4218 -0.2095 0.0916 0.1091

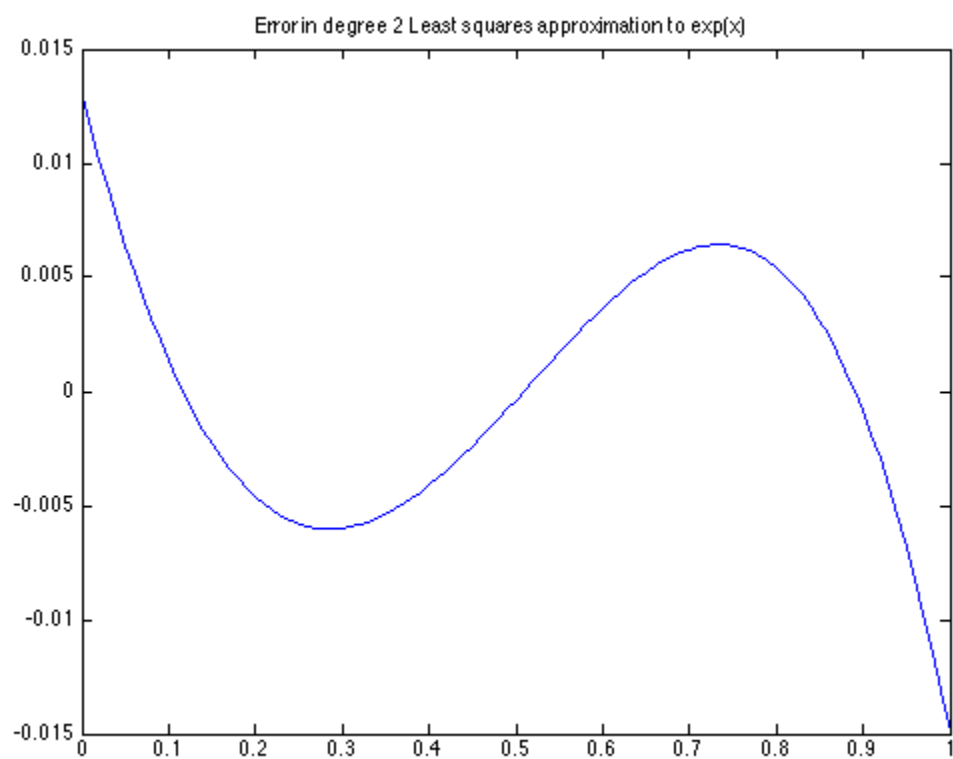
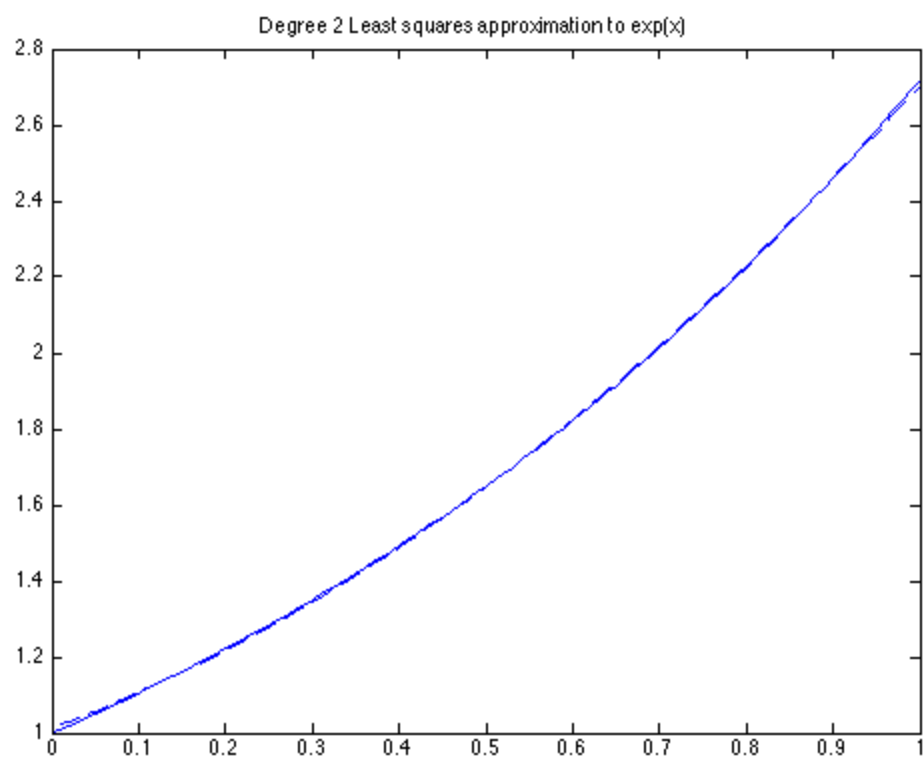
Columns 57 through 63

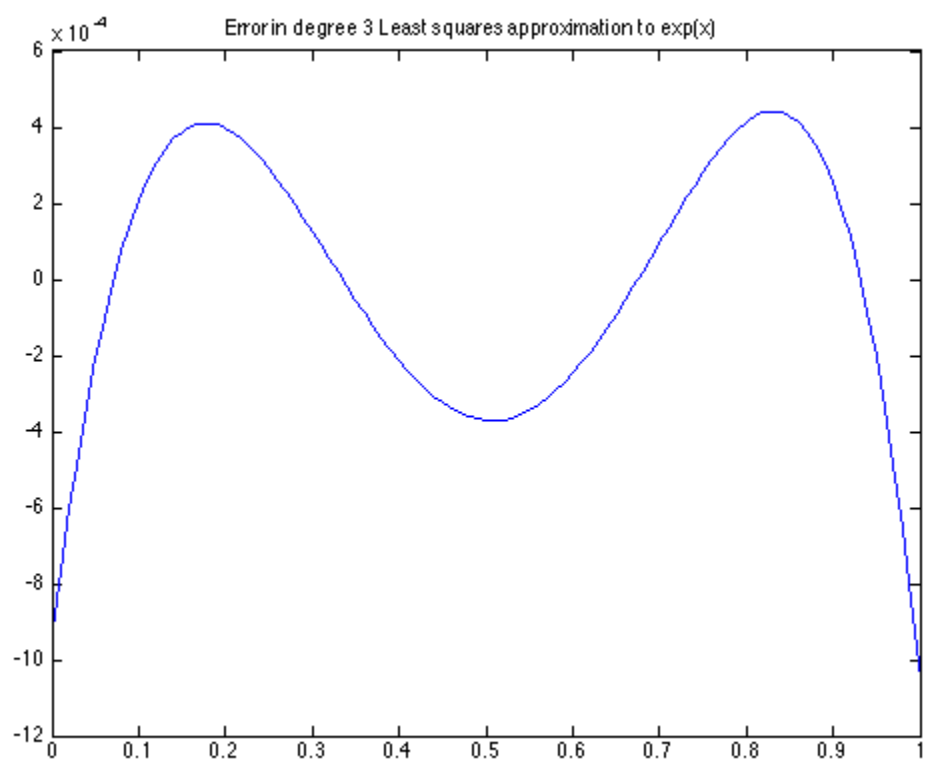
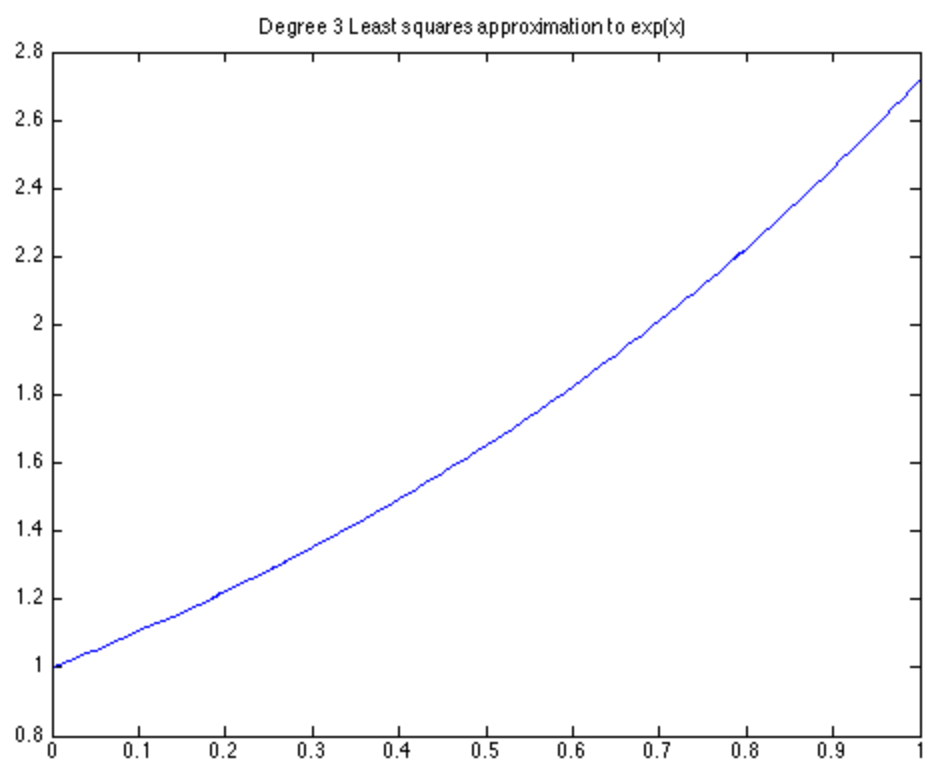
-0.2072 0.3386 -0.3575 0.3771 -0.3045 0.2334 -0.1227

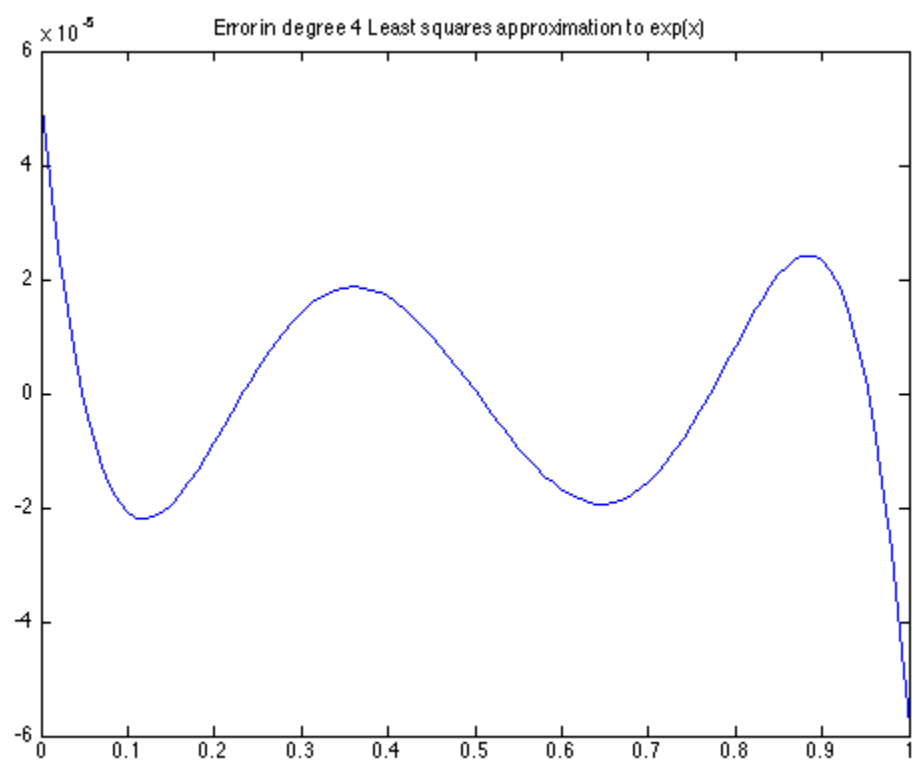
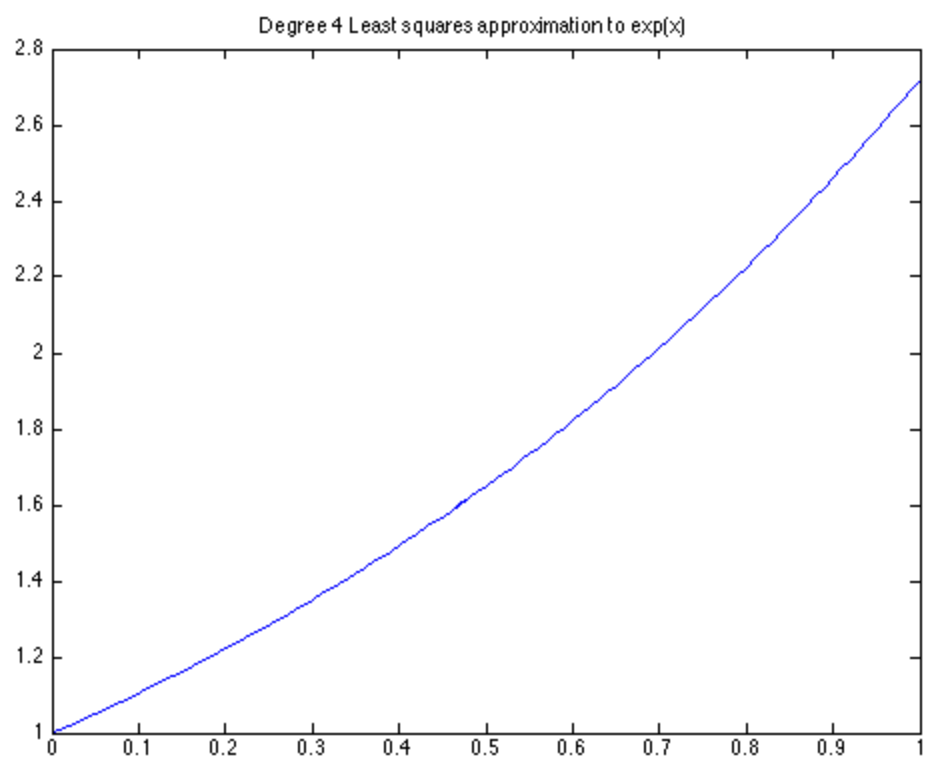
Column 64

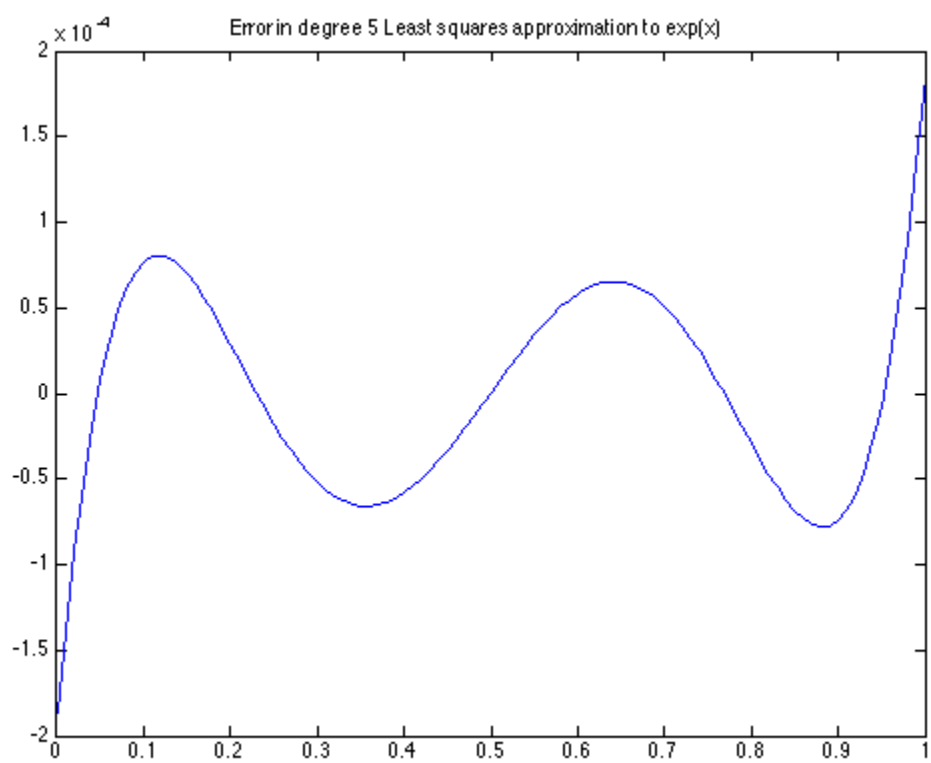
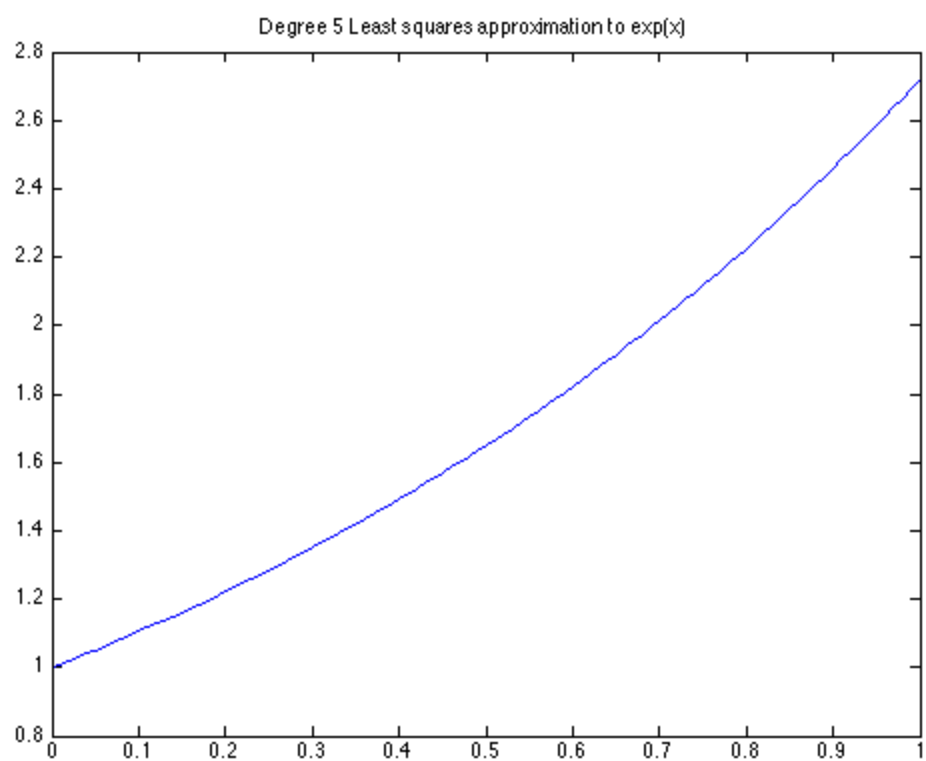
0.0399

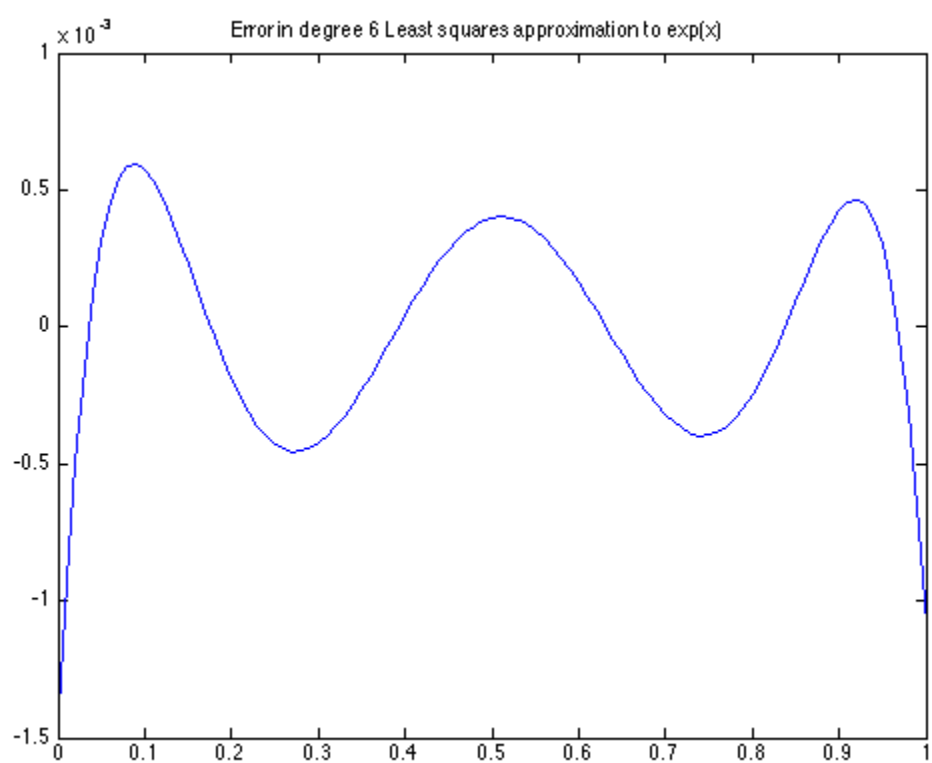
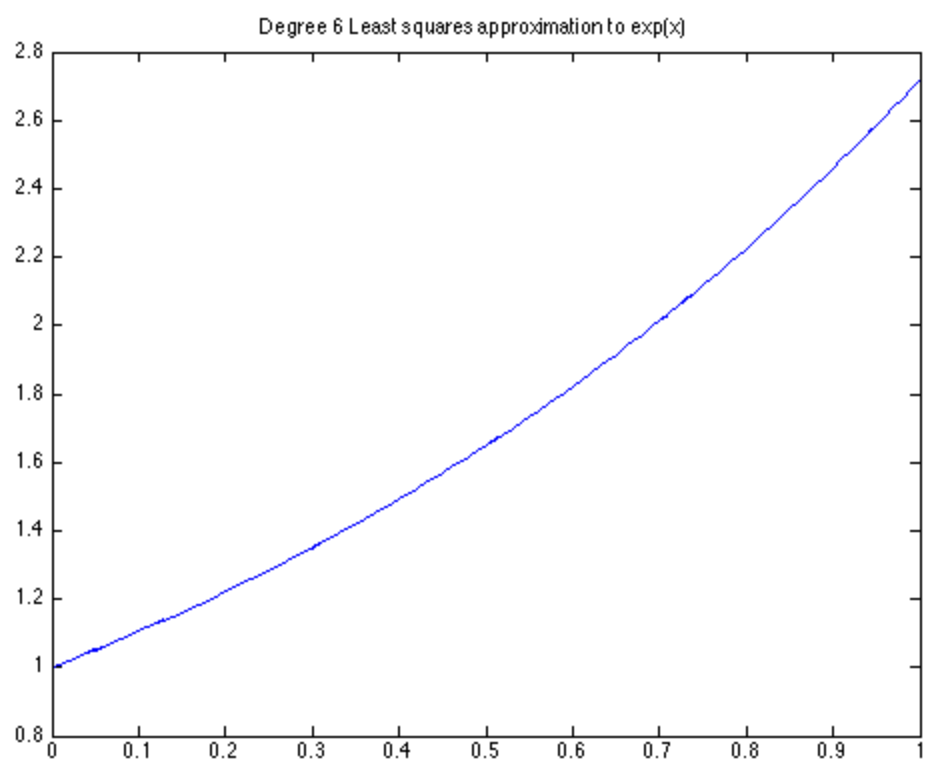


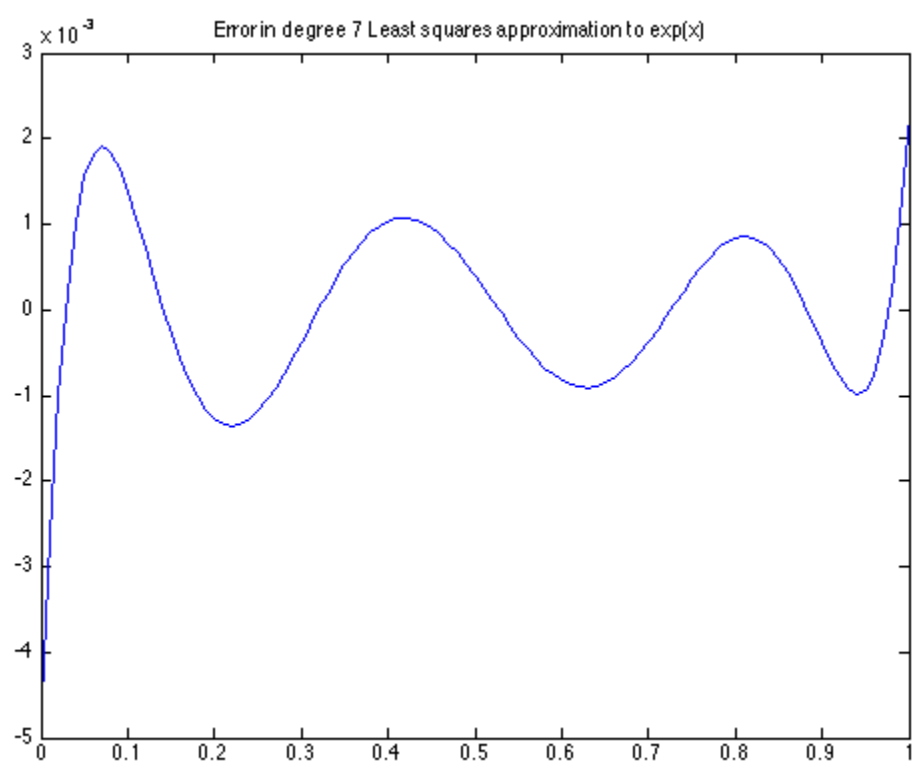
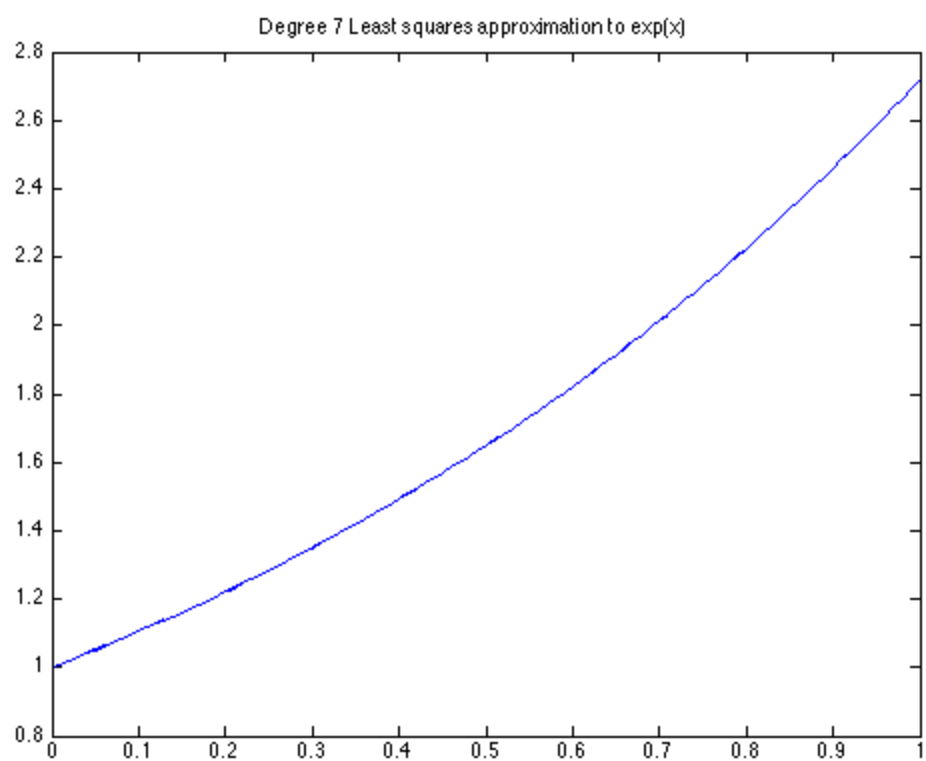


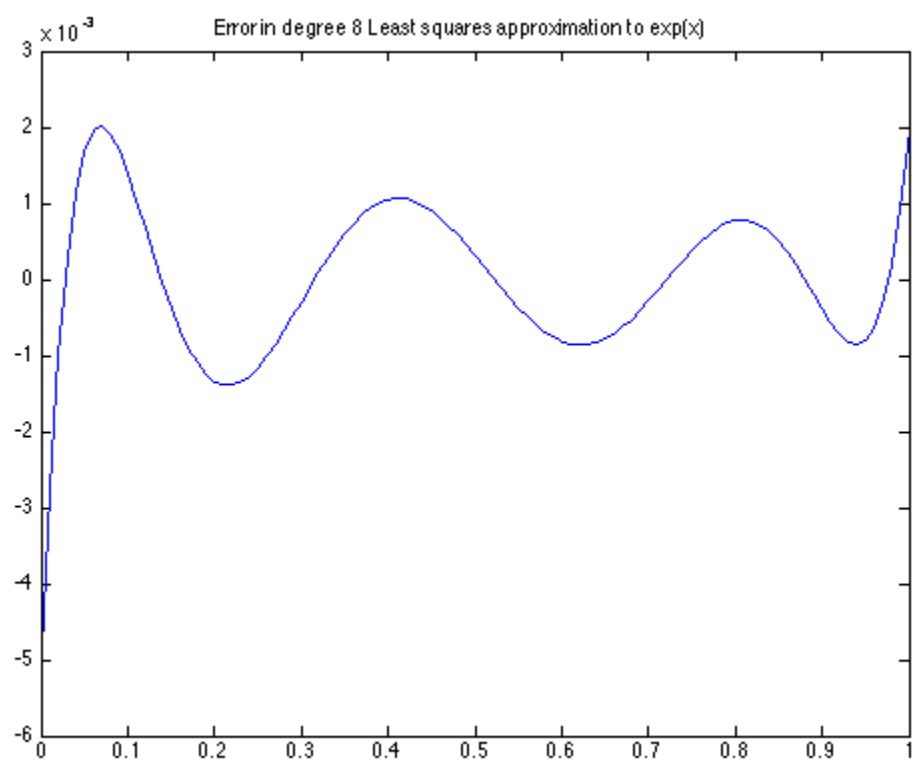
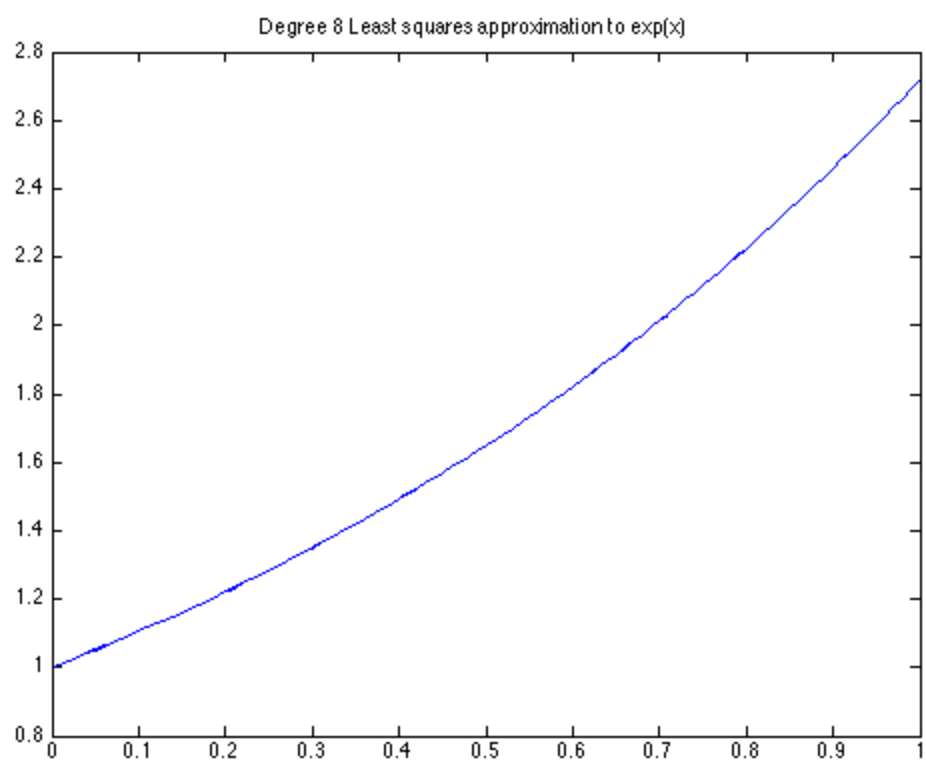


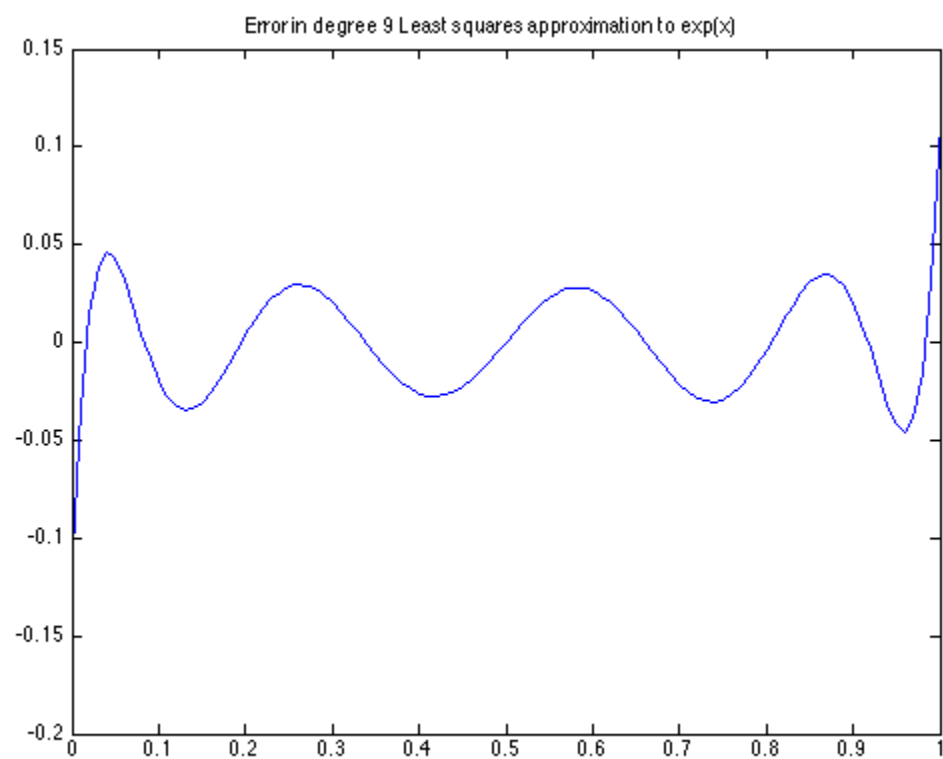
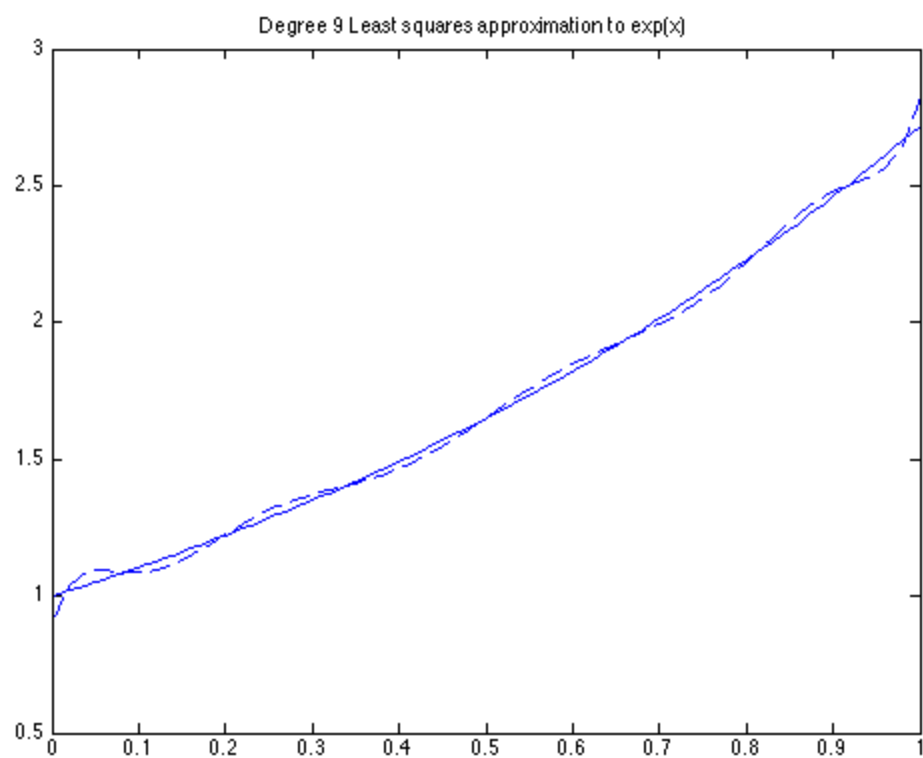


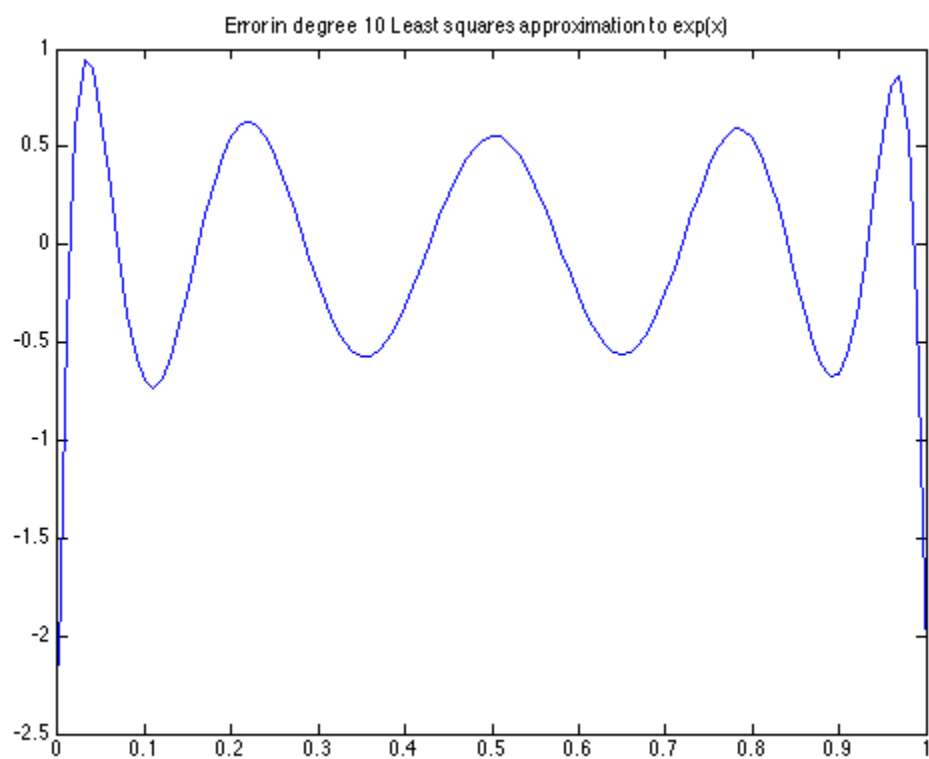
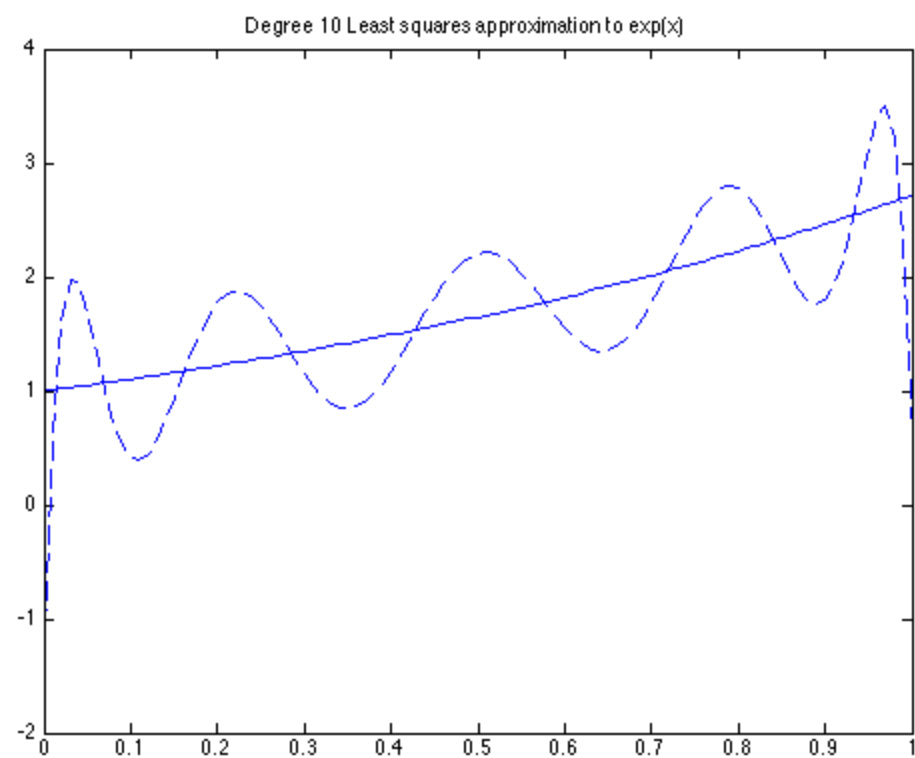


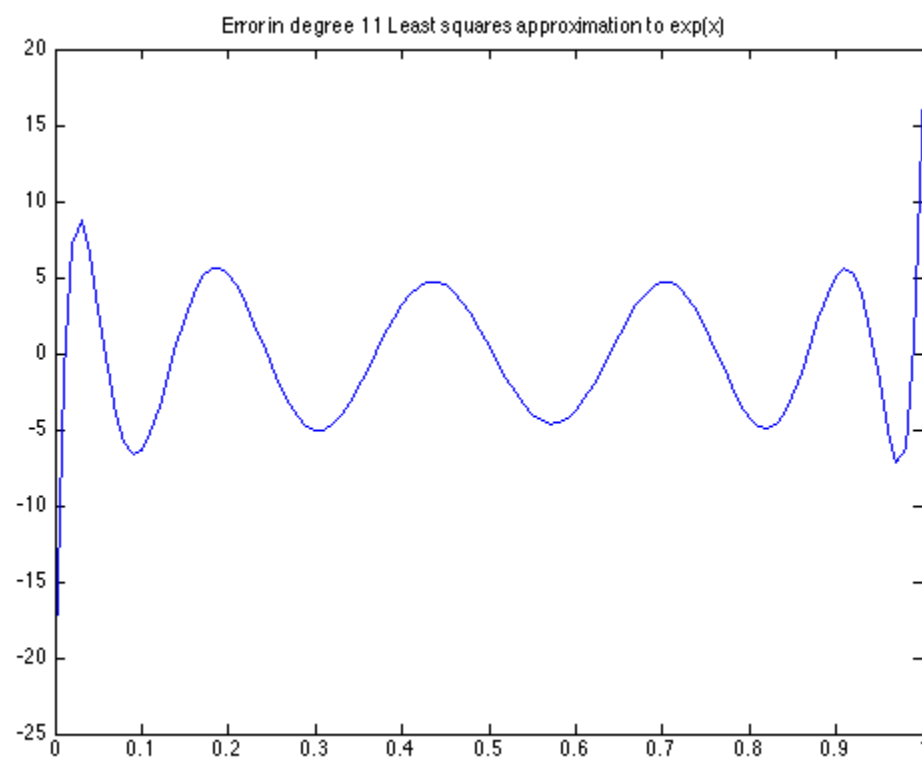
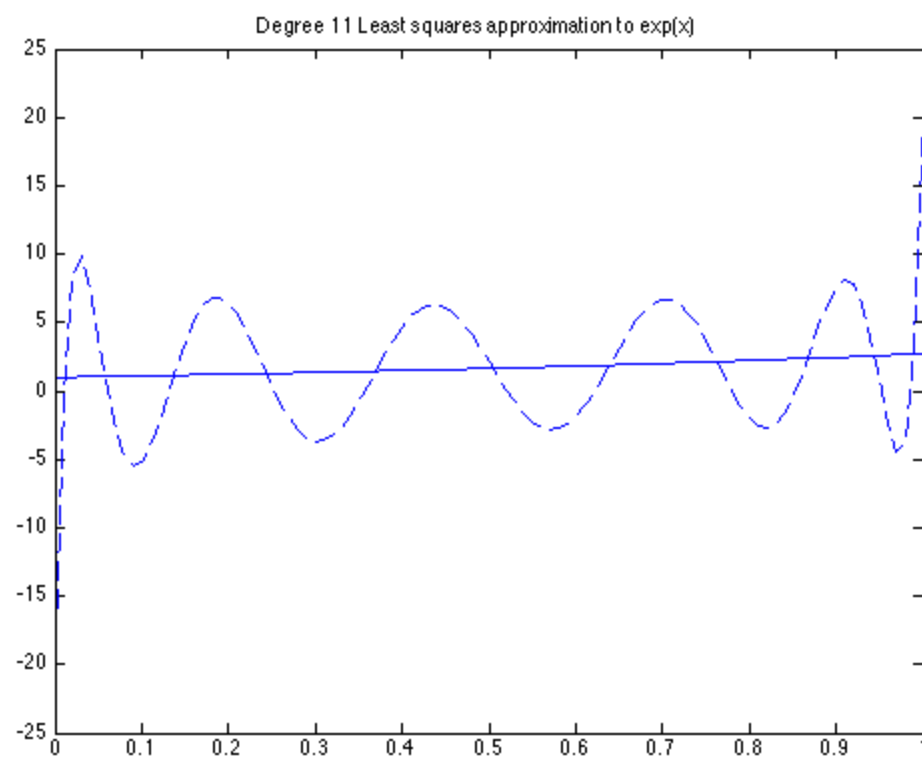


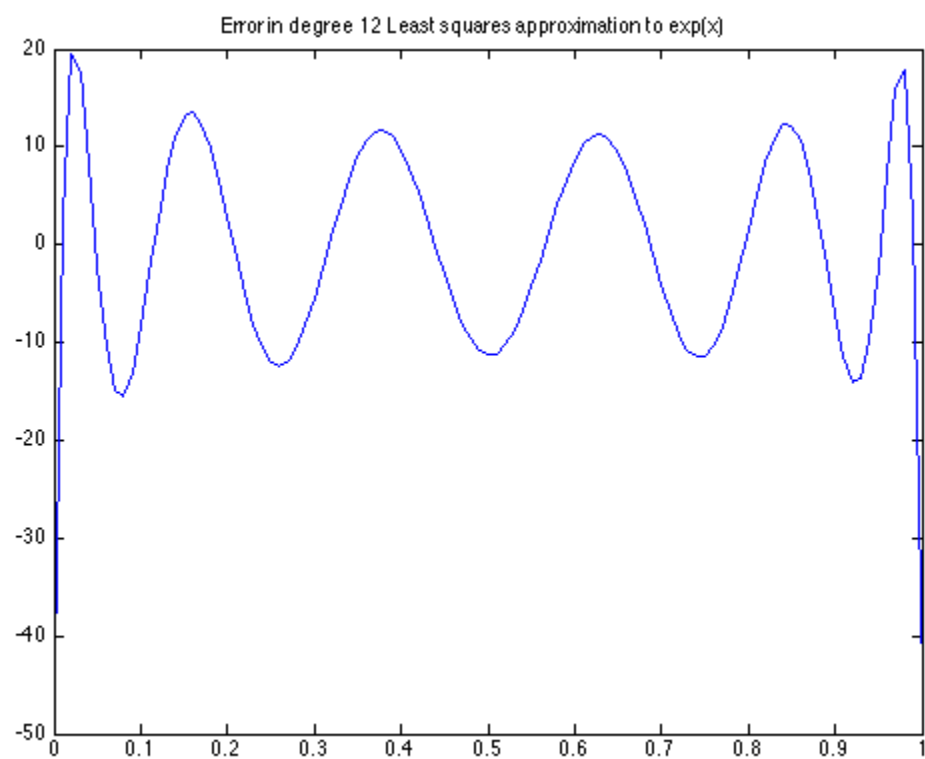
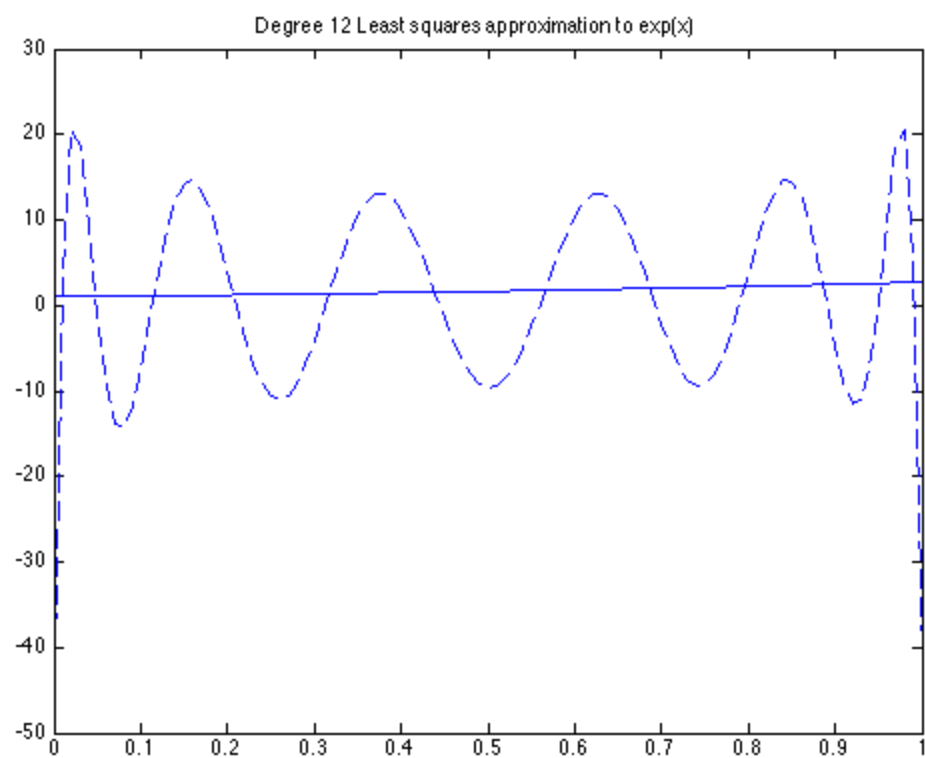


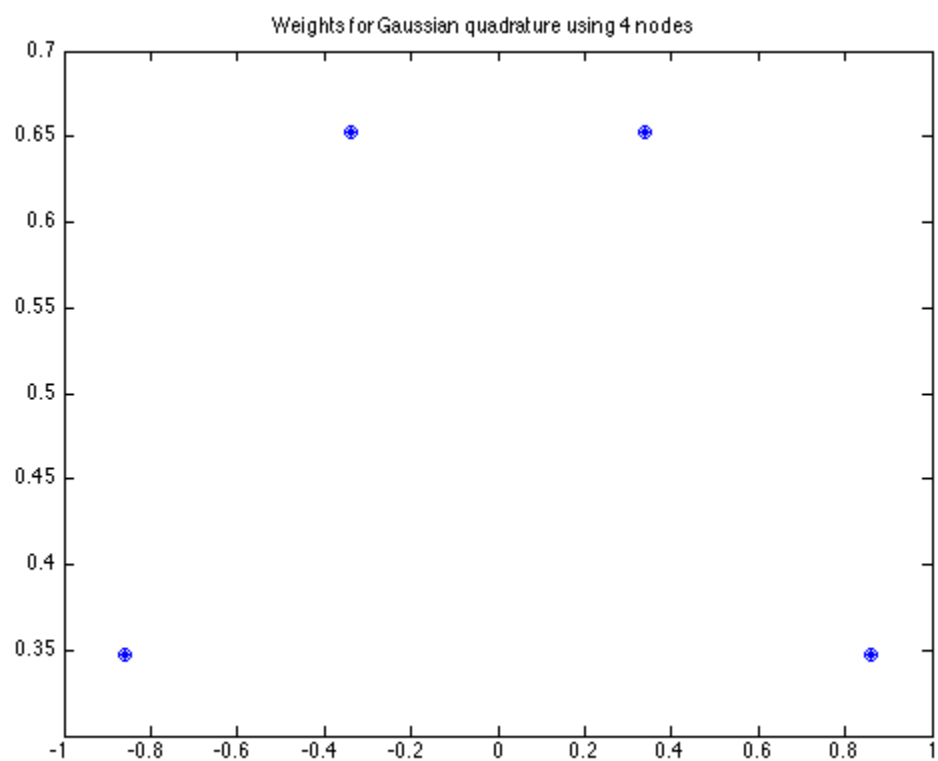
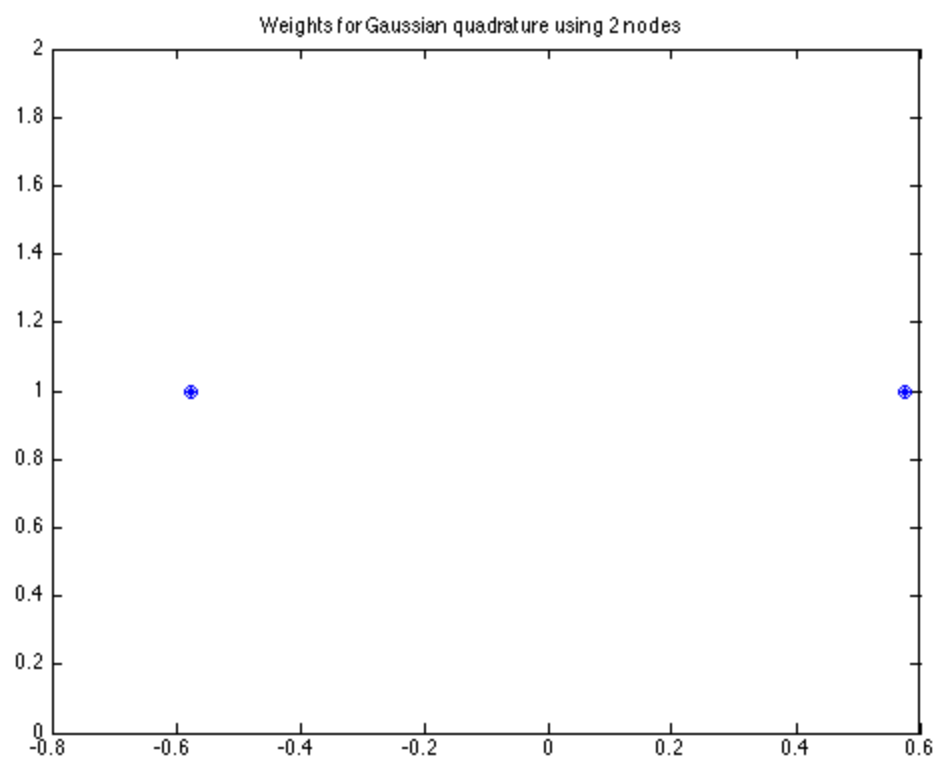


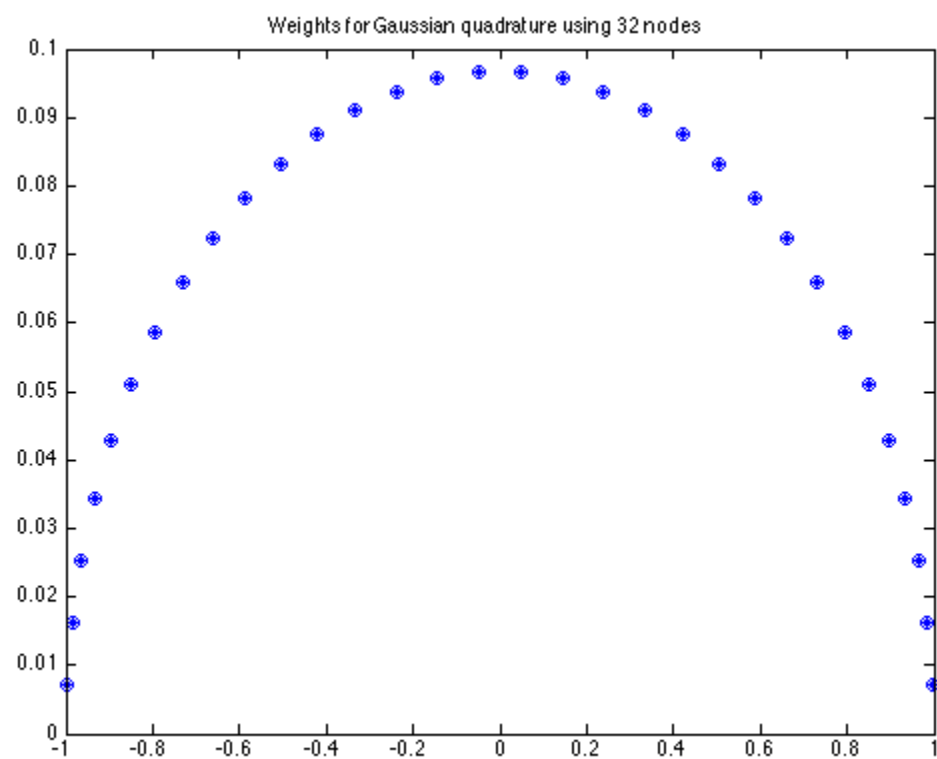
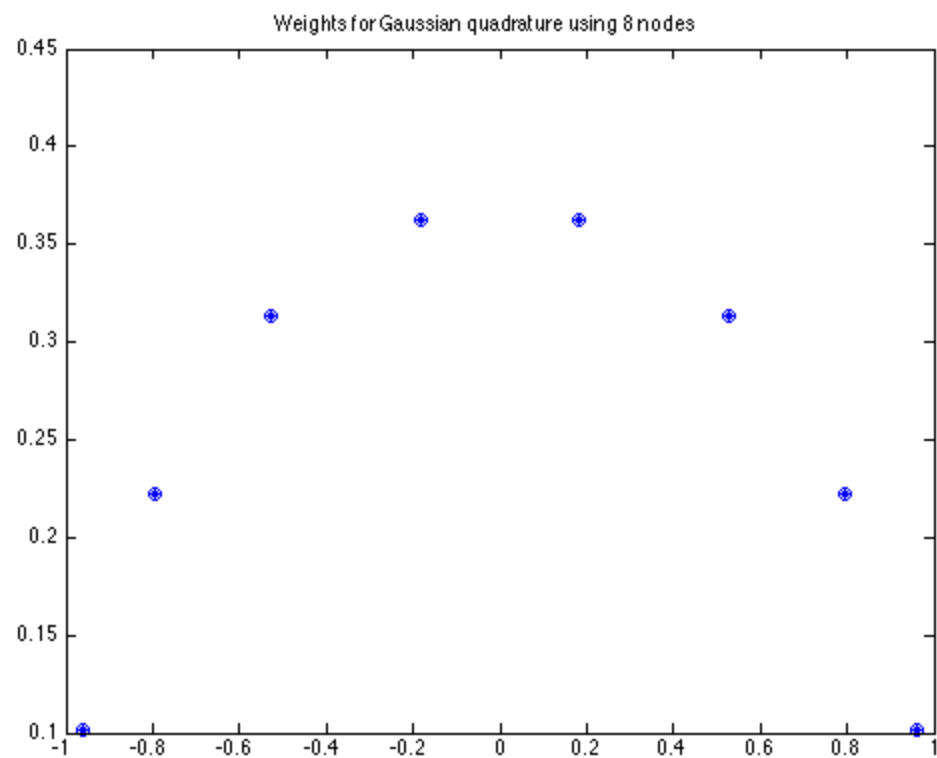


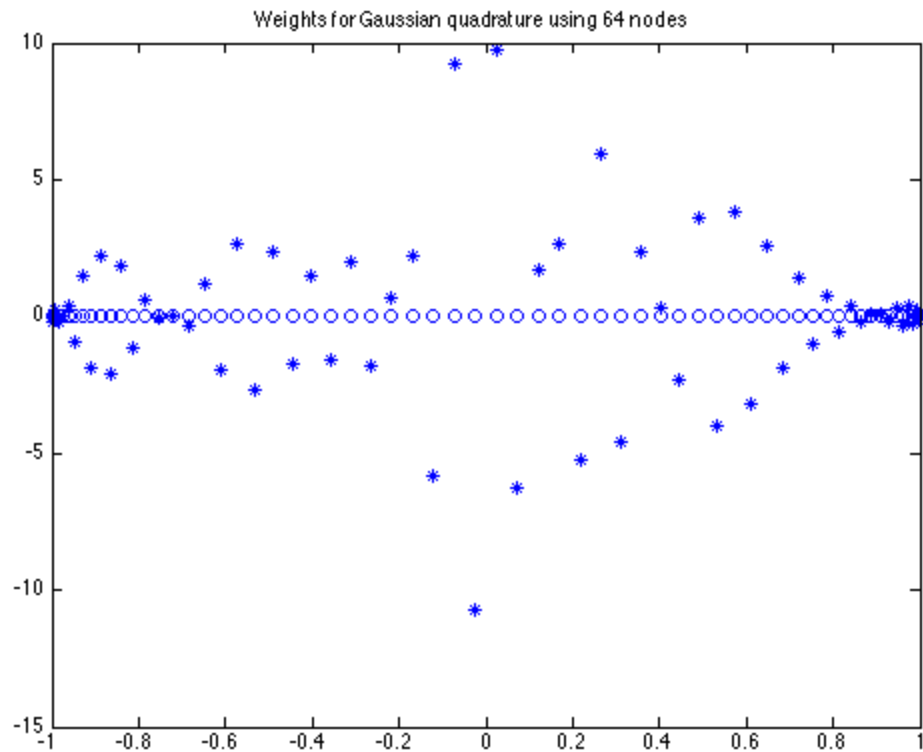












3.

```
cubspline(8)
cubspline(16)
```

```
% Run each of the methods 3 times, find the total, and divide by 3
% to display the average times.
```

```
% a. GEPivot
```

```
tot641=0;
tot1281=0;
tot2561=0;
for i=1:3
    [res64, time64, itnum64] = cubspline(64);
    [res128, time128, itnum128] = cubspline(128);
    [res256, time256, itnum256] = cubspline(256);
    tot641 = tot641 + time64;
    tot1281 = tot1281 + time128;
    tot2561 = tot2561 + time256;
end
```

```
% b. Jacobian
```

```
tot642=0;
tot1282=0;
tot2562=0;
for i=1:3
```

```

    [res64, time64, itnum64] = cubspline2(64);
    [res128, time128, itnum128] = cubspline2(128);
    [res256, time256, itnum256] = cubspline2(256);
    tot642 = tot642 + time64;
    tot1282 = tot1282 + time128;
    tot2562 = tot2562 + time256;
end

```

```

% c. Gauss-Seidel: ignore timings

```

```

tot643=0;
tot1283=0;
tot2563=0;
for i=1:3
    [res64, time64, itnum64] = cubspline3(64);
    [res128, time128, itnum128] = cubspline3(128);
    [res256, time256, itnum256] = cubspline3(256);
    tot643 = tot643 + time64;
    tot1283 = tot1283 + time128;
    tot2563 = tot2563 + time256;
end

```

```

% d. Conjugate-Gradient

```

```

tot644=0;
tot1284=0;
tot2564=0;
for i=1:3
    [res64, time64, itnum64] = cubspline4(64);
    [res128, time128, itnum128] = cubspline4(128);
    [res256, time256, itnum256] = cubspline4(256);
    tot644 = tot644 + time64;
    tot1284 = tot1284 + time128;
    tot2564 = tot2564 + time256;
end

```

```

% e. Tridiag

```

```

tot645=0;
tot1285=0;
tot2565=0;
for i=1:3
    [res64, time64, itnum64] = cubspline5(64);
    [res128, time128, itnum128] = cubspline5(128);
    [res256, time256, itnum256] = cubspline5(256);
    tot645 = tot645 + time64;
    tot1285 = tot1285 + time128;
    tot2565 = tot2565 + time256;
end

```

```

disp(sprintf('Runtimes for solving tridiagonal systems using \n'))
disp(sprintf('\n\t GEPivot\t Jacobi\t Conjugate-Gradient\t Tridiagonal\n'))
disp(sprintf('64\t %d\t %d\t %d\t %d\n', tot641/3, tot642/3, tot644/3, tot645/3));
disp(sprintf('128\t %d\t %d\t %d\t %d\n', tot1281/3, tot1282/3, tot1284/3, tot1285/3));
disp(sprintf('256\t %d\t %d\t %d\t %d\n', tot2561/3, tot2562/3, tot2564/3, tot2565/3));

```

```
% Tridiagonal had the lowest average runtimes, followed by Jacobi, then
% Conjugate-Gradient, and GEPivot was the slowest.
```

```
ans =
```

```
1.0923e-14
```

```
ans =
```

```
4.4702e-14
```

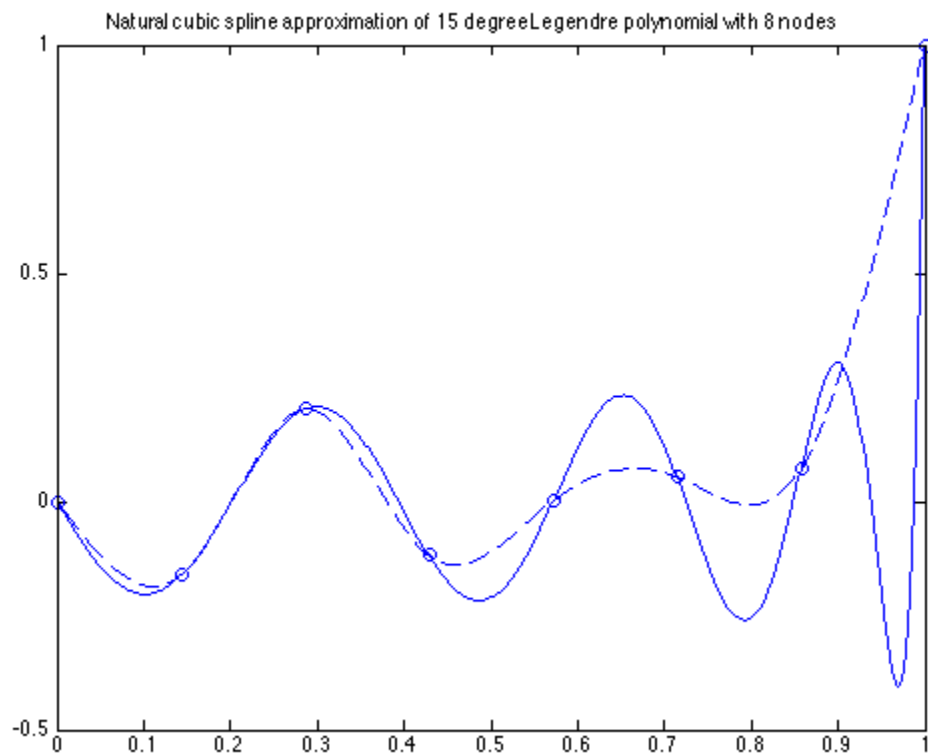
```
Runtimes for solving tridiagonal systems using
```

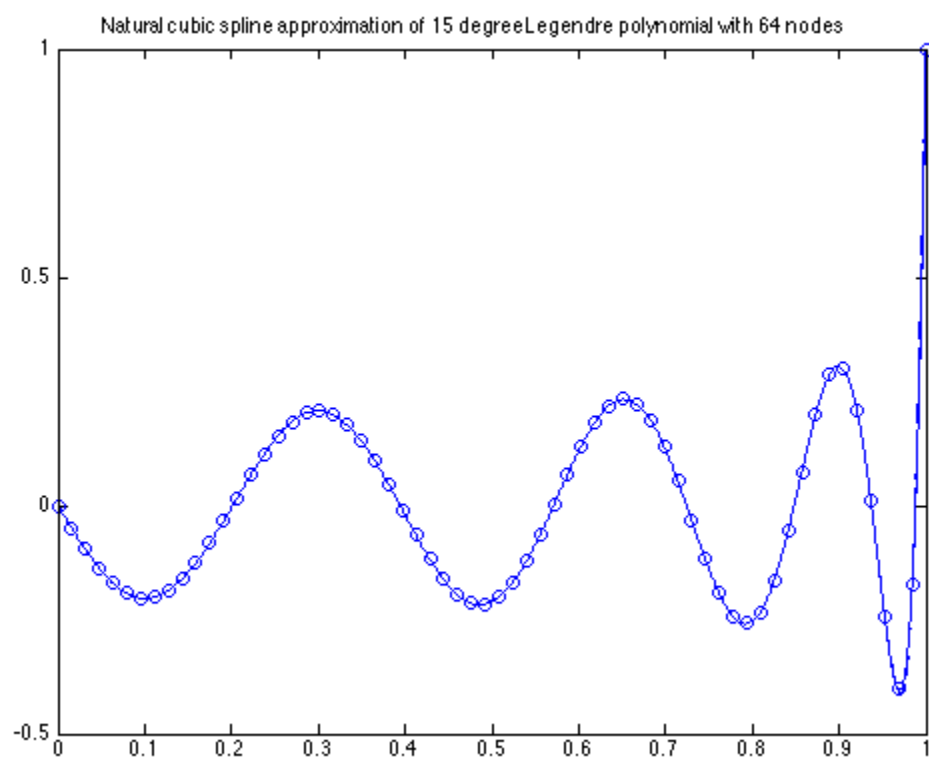
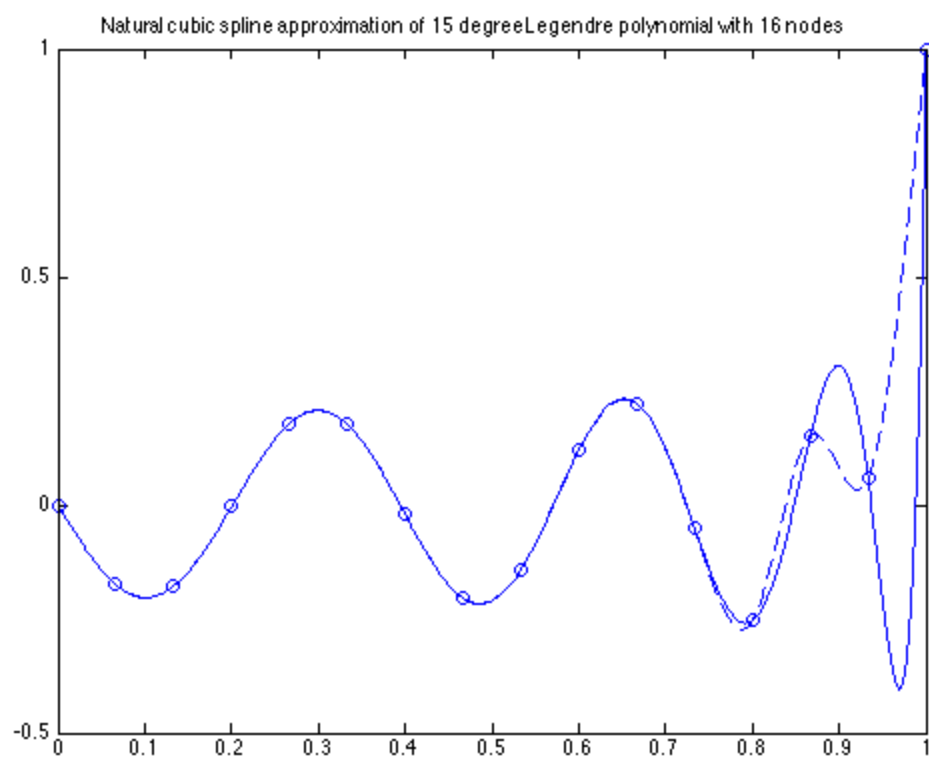
```
n GEPivot Jacobi Conjugate-Gradient Tridiagonal
```

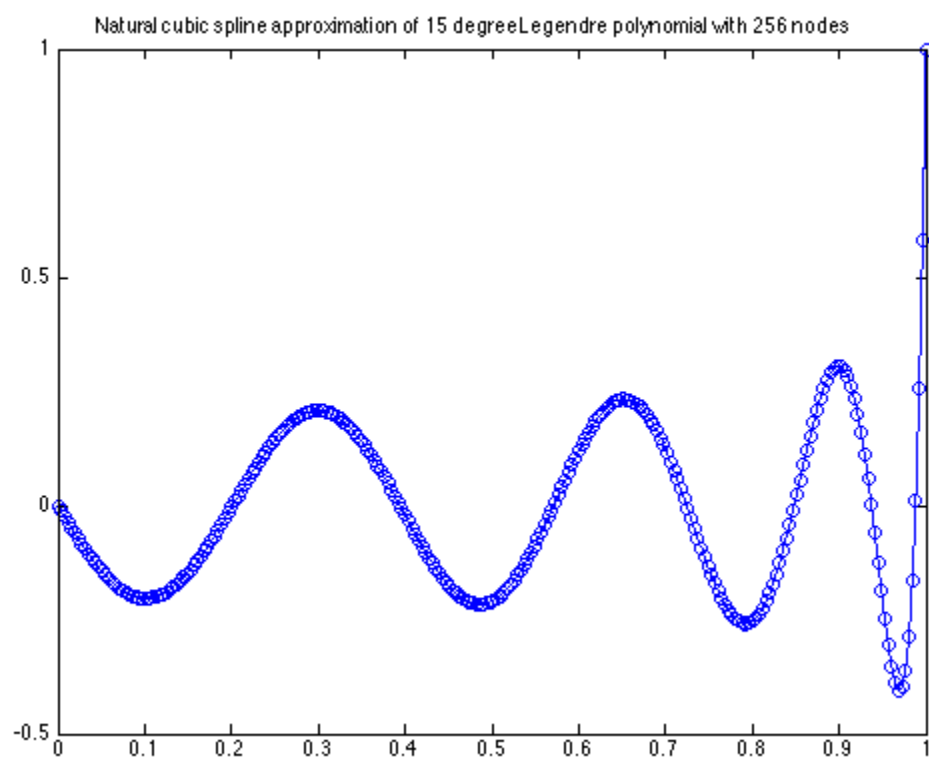
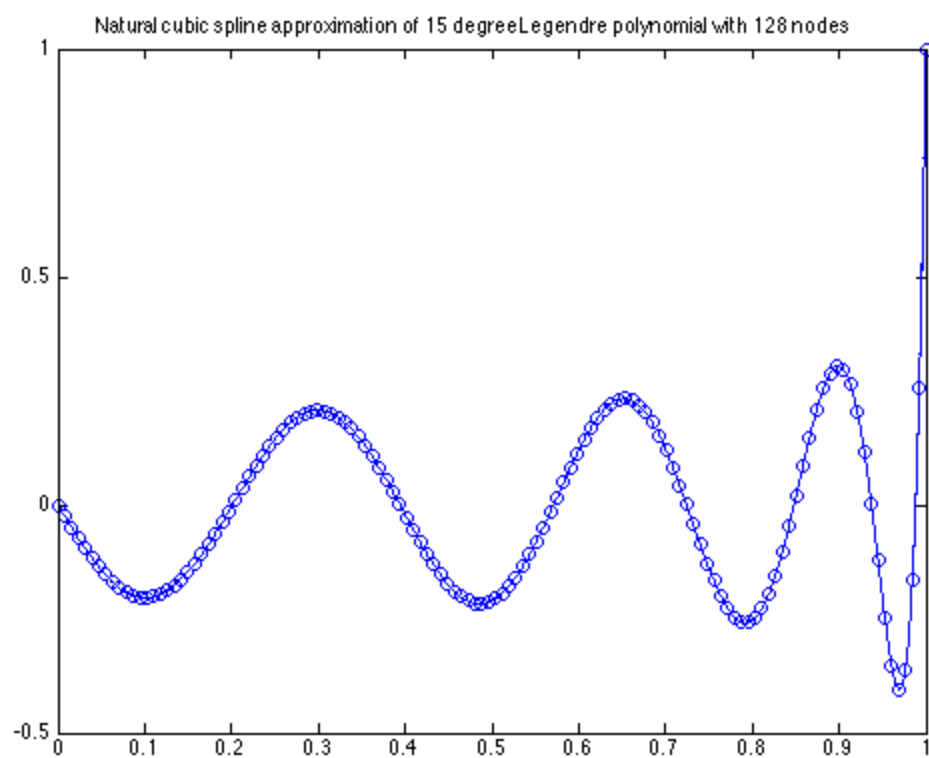
```
64 1.730935e-02 4.215353e-04 2.429209e-03 5.137500e-05
```

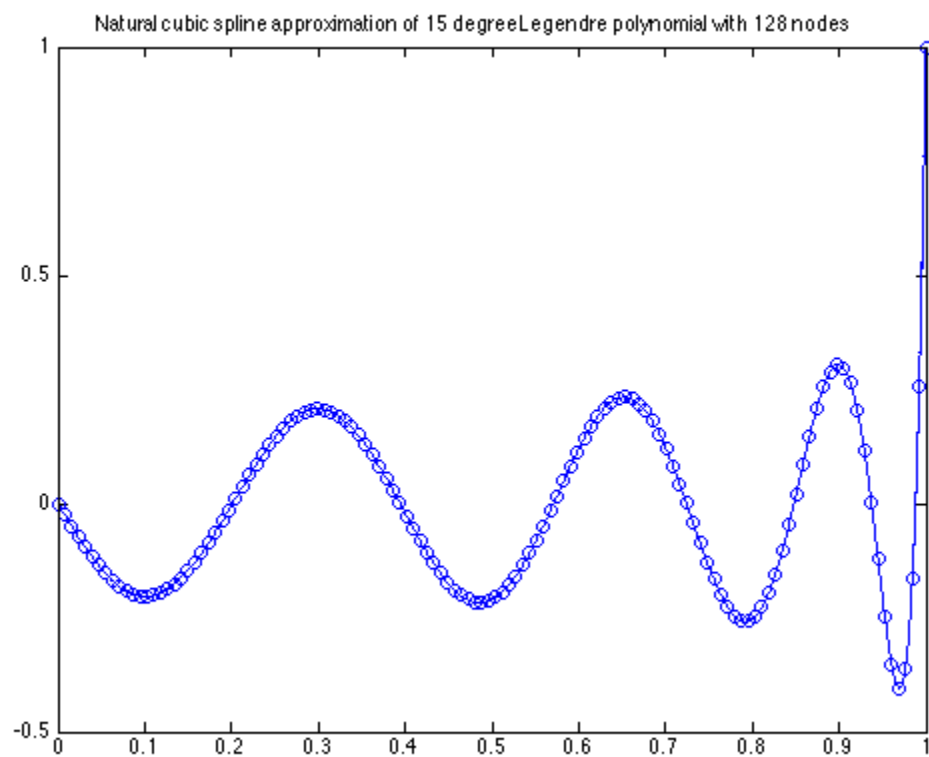
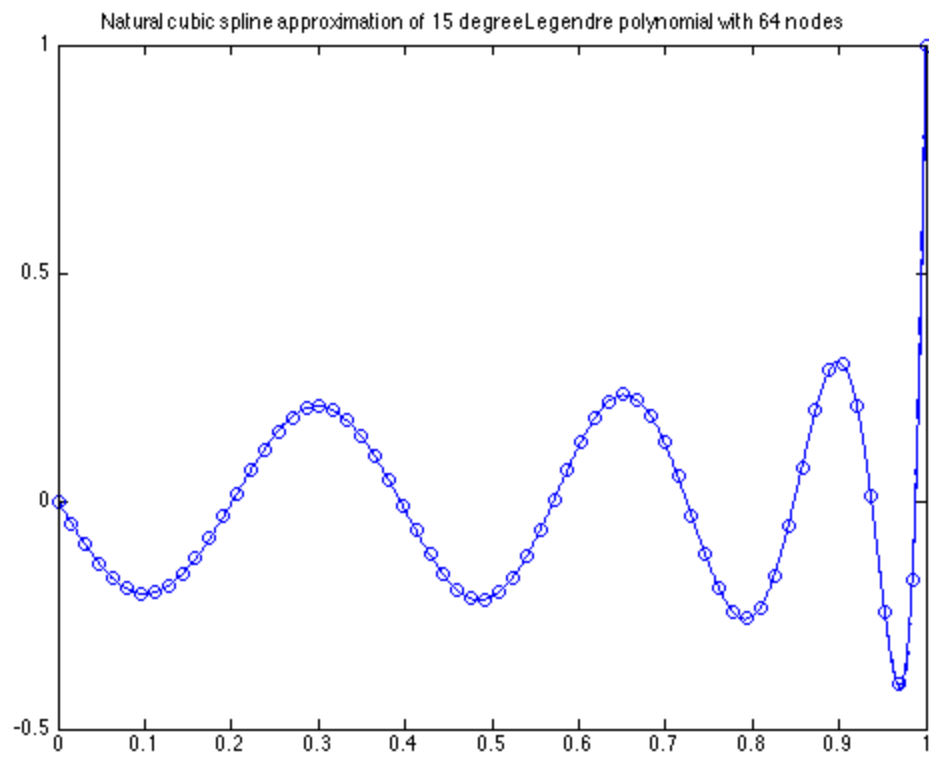
```
128 8.051737e-02 7.223910e-04 3.633219e-03 5.584033e-05
```

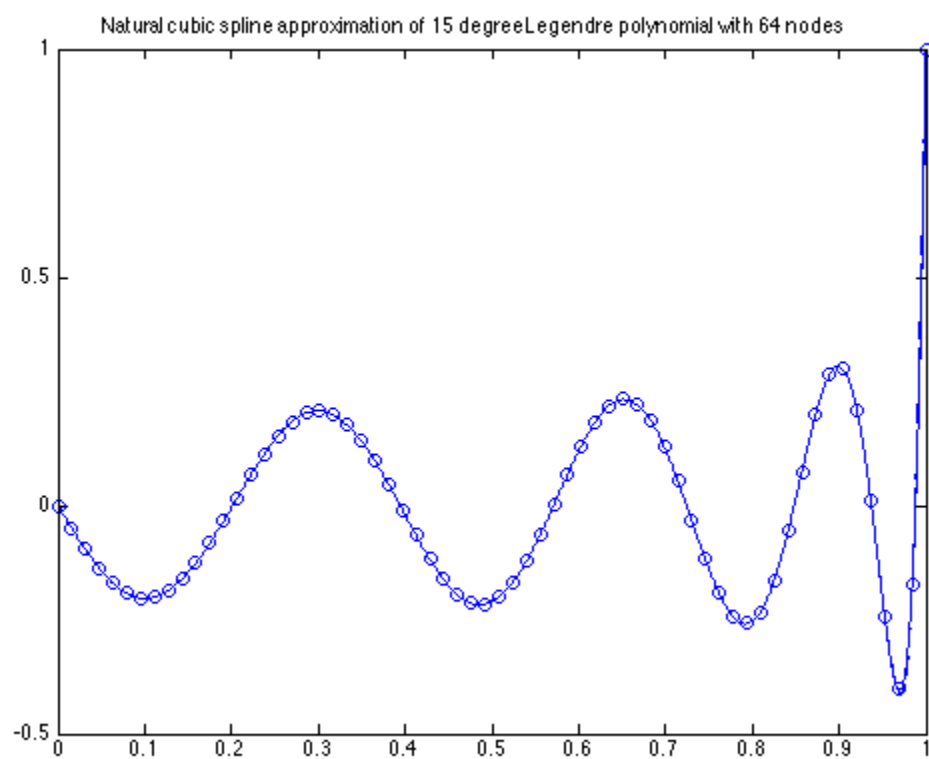
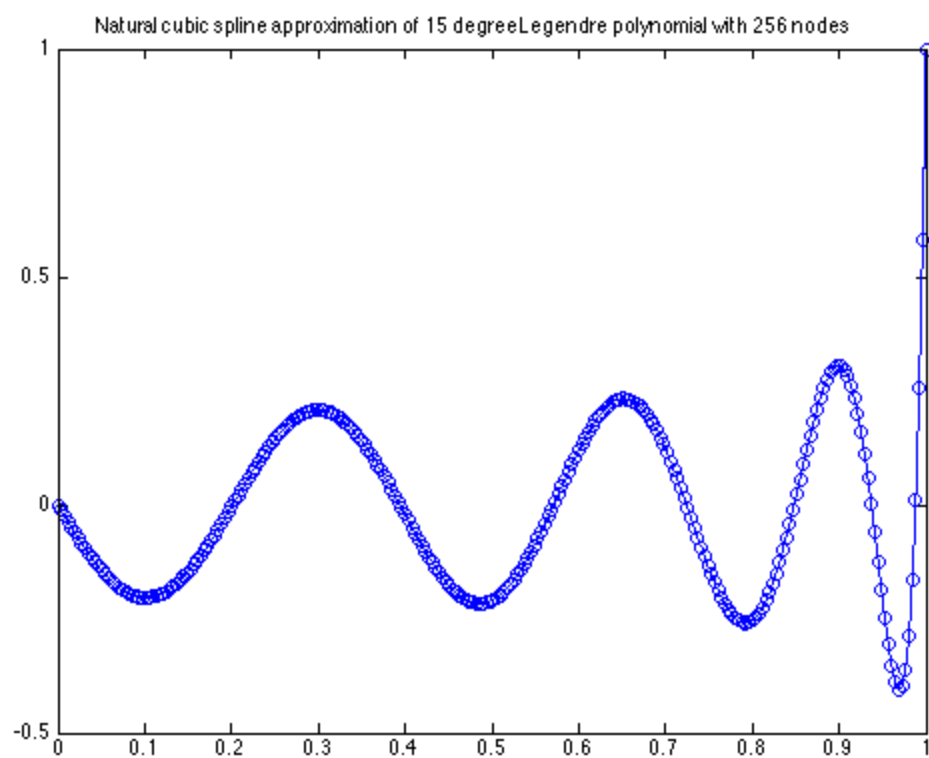
```
256 3.496016e-01 1.648796e-03 5.013996e-03 1.086917e-04
```

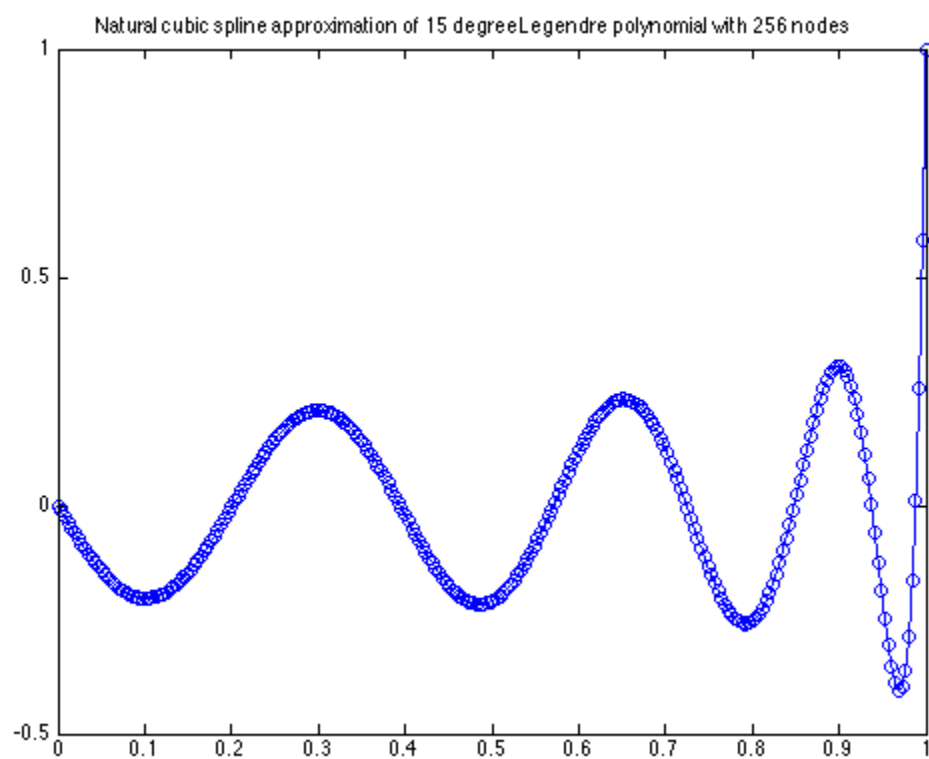
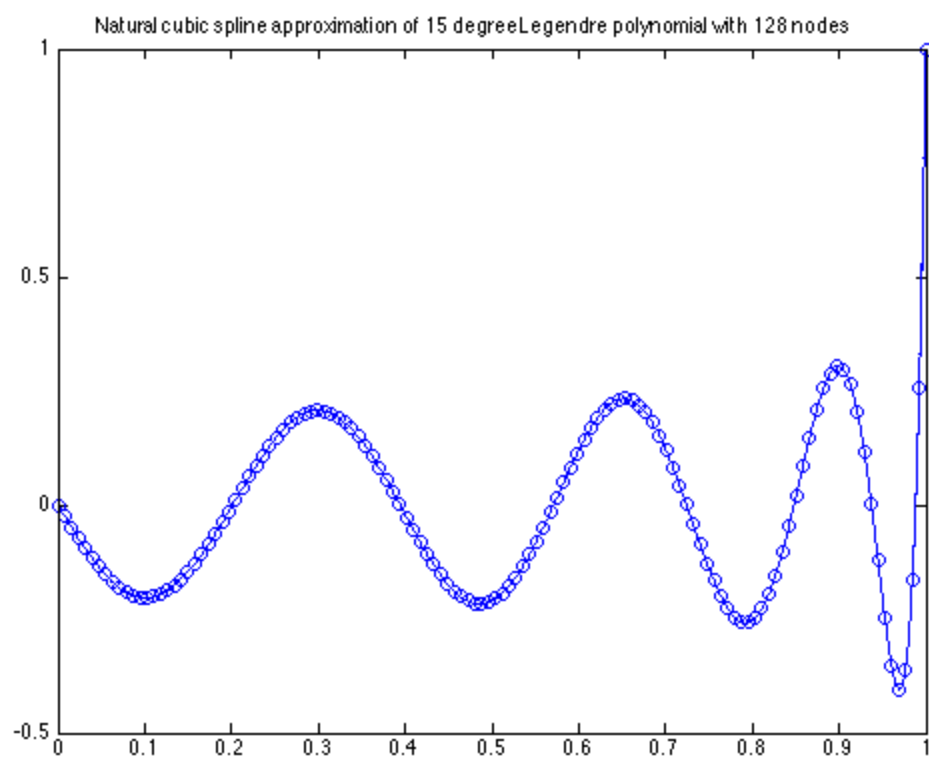












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