

FAST ITERATIVE SOLVERS

PROJECT 3: Finding Maximum Eigen Value

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1. Run the power method for 100, 500, 1000, and 5000, iterations and Record the error, and the runtime for each of those runs.

For: **nos6.mtx**, it converges to given tolerance of 10^{-8} in **781 iterations** and error is the difference of largest eigenvalue between successive iterations

ITERATIONS	ERROR	ELAPSED TIME (sec)
100	1.284×10^3	0.638790
500	5.1602×10^{-4}	0.665496
1000	9.31×10^{-10}	0.673188
5000	0	0.864531
1000(with tolerance limit)	7.45×10^{-9}	0.661298
5000(with tolerance limit)	7.45×10^{-9}	0.720034

(these are the values taken without setting the tolerance limit, if not mentioned otherwise)

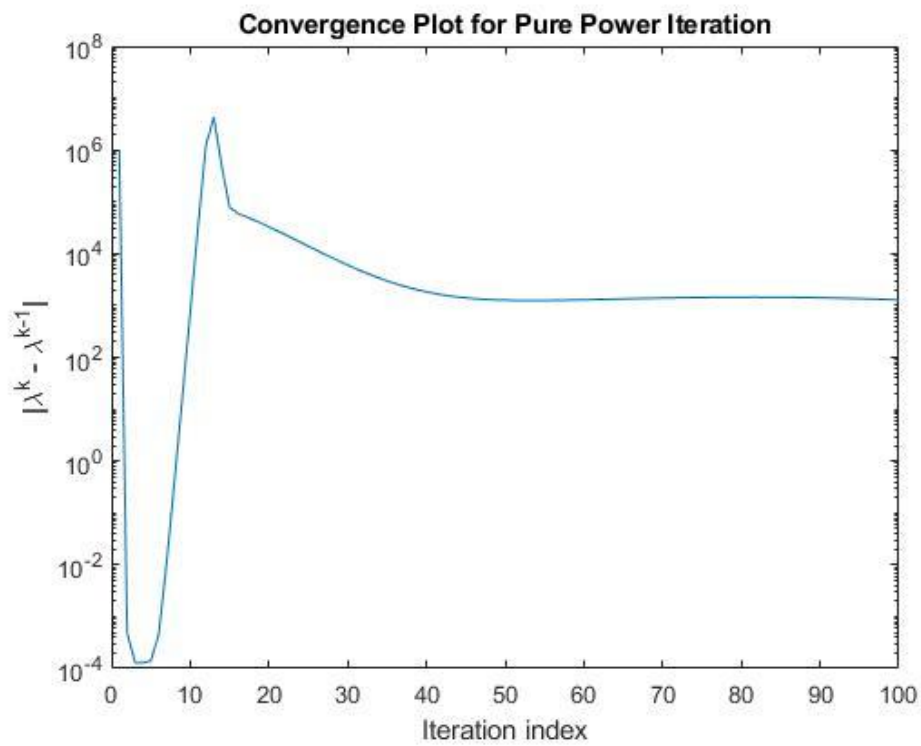
Recorded Eigen Value: 7.65×10^6

For: **s3rmt3m3.mtx**, it converges to given tolerance of 10^{-8} in **2145 iterations** and error is the difference of largest eigenvalue between successive iterations

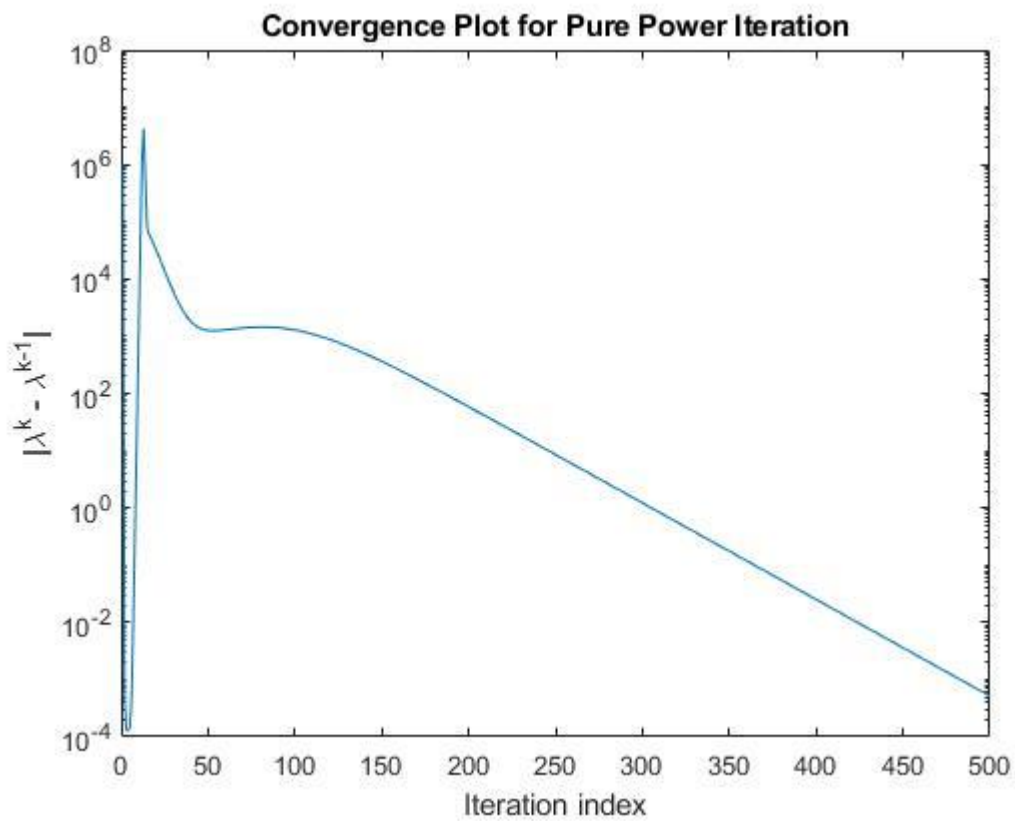
ITERATIONS	ERROR	ELAPSED TIME (sec)
100	0.9564	37.22
500	0.0041	37.93
1000	1.9078×10^{-5}	38.67
5000	1.8×10^{-12}	45.54
5000 (with tolerance limit)	9.98×10^{-9}	42.43

(these are the values taken without setting the tolerance limit, if not mentioned otherwise)

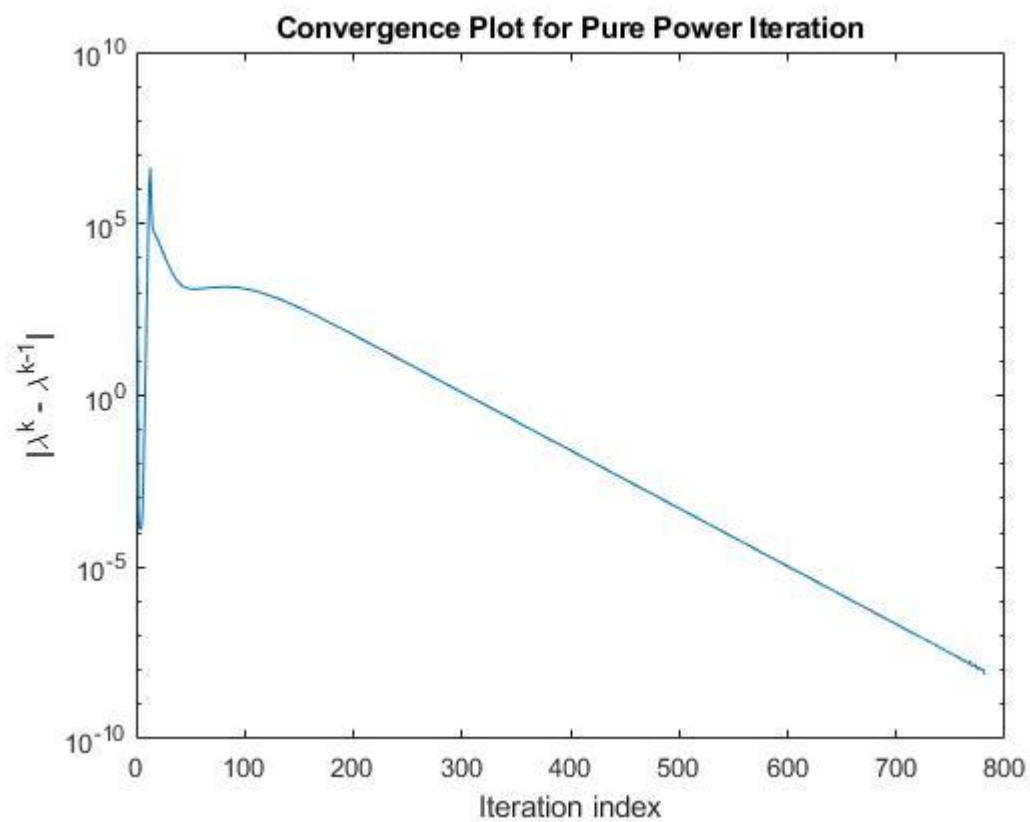
Convergence Plots for nos6.mtx for varying number of iterations



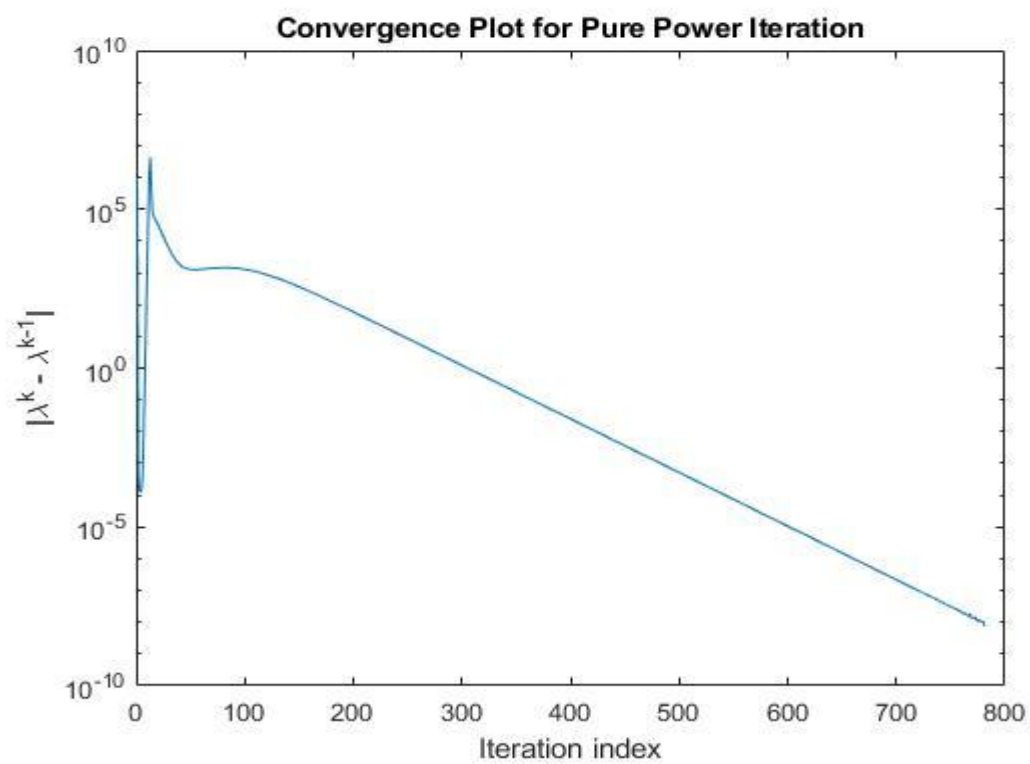
100 Iterations (tolerance of 10^{-8} is not met)



500 Iterations (tolerance of 10^{-8} is not met)



1000 Iterations(tolerance is met at 781 iterations)

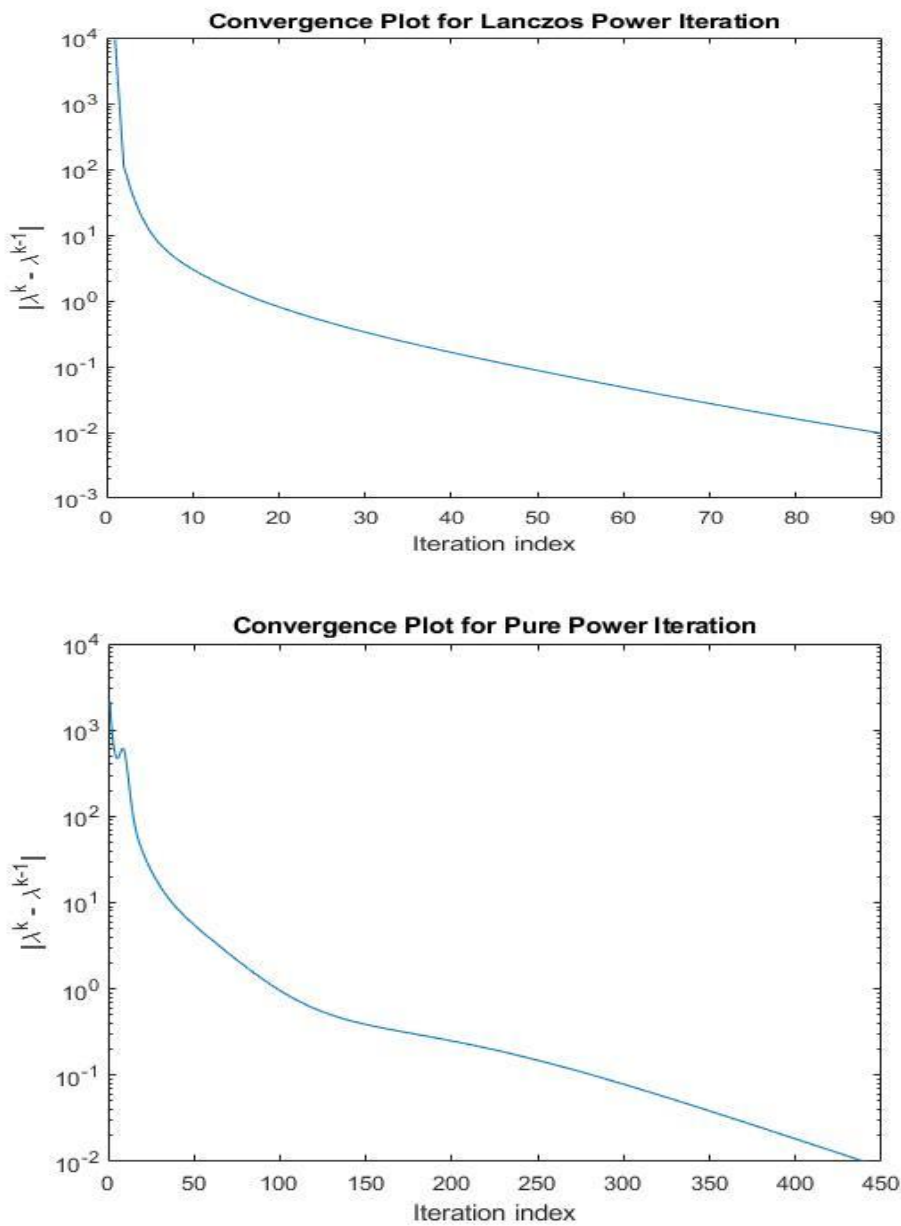


5000 iterations(tolerance is met at 781th iterations)

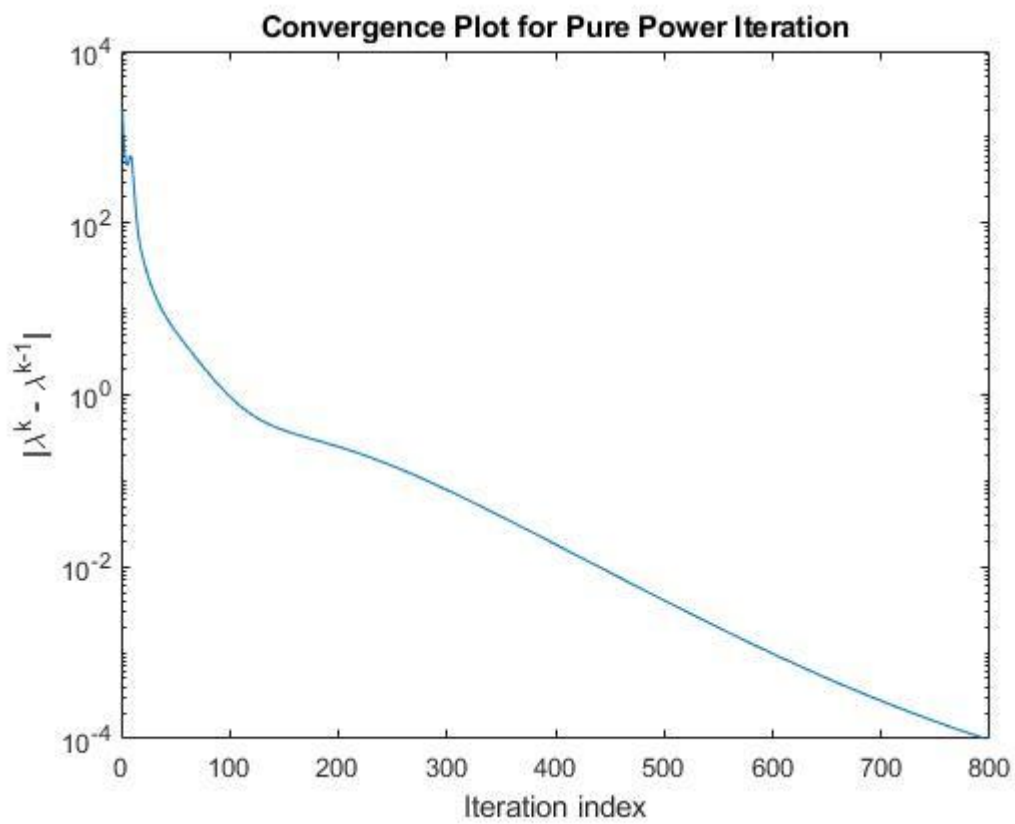
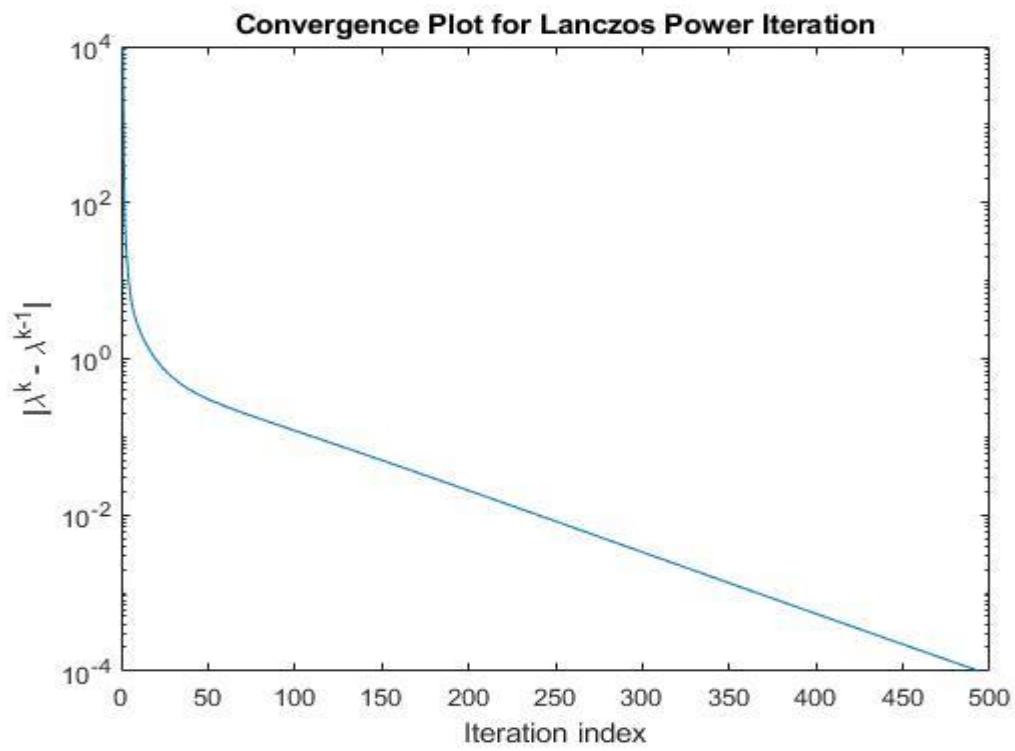
For: **s3rmt3m3.mtx** the maximum eigenvalue mentioned in matrix market is 9.5986080894852857E+03. And convergence plots for this matrix is shown in next section as comparison with Lanczos method.

2. Run the Lanczos method for $m = 30; 50; 75; 100$, where m is the dimension of the Krylov Space. For the power iteration you can use a convergence criterion error $< tol$, For the tolerance tol you can use 10^{-2} ($m = 30$), 10^{-4} ($m = 50$), 10^{-6} ($m = 75$), 10^{-10} ($m = 100$).

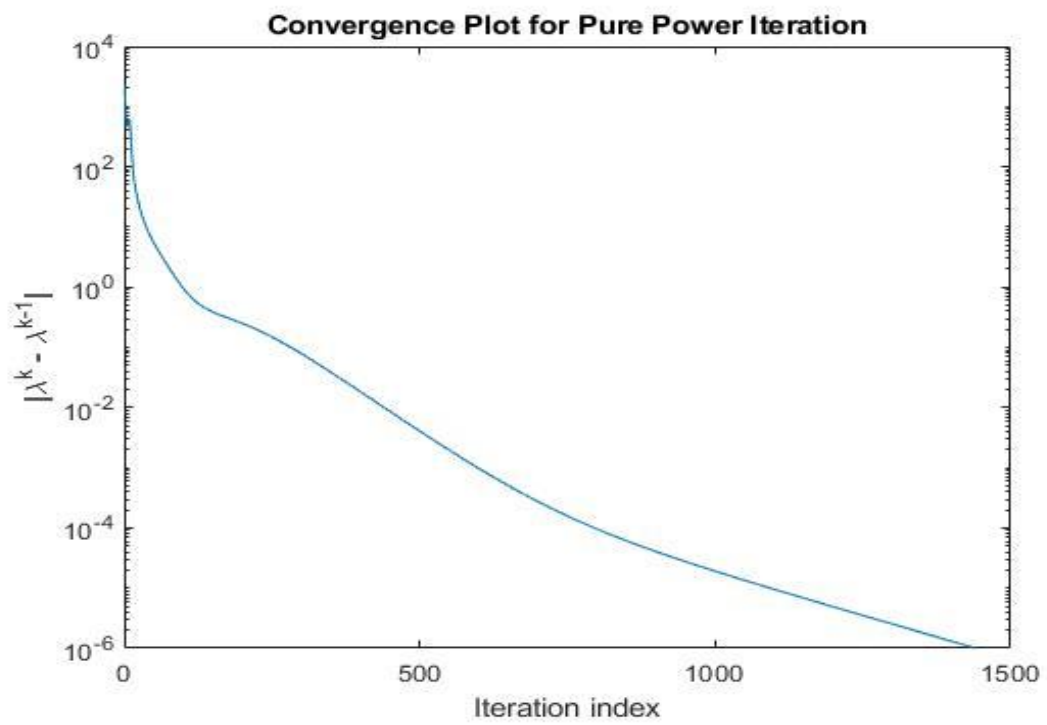
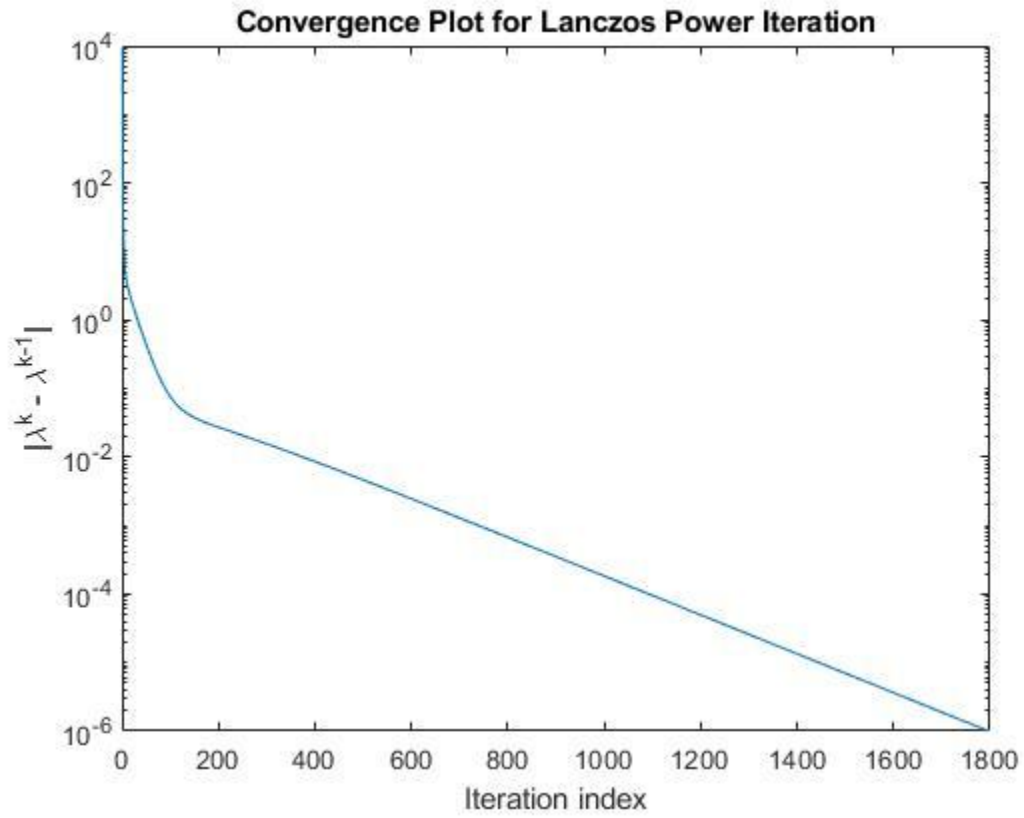
Maximum iterations are set to 5000, but in both methods tolerance is reached much before this. From here on we consider s3rmt matrix for further analysis. **Table in page 8** records needed data



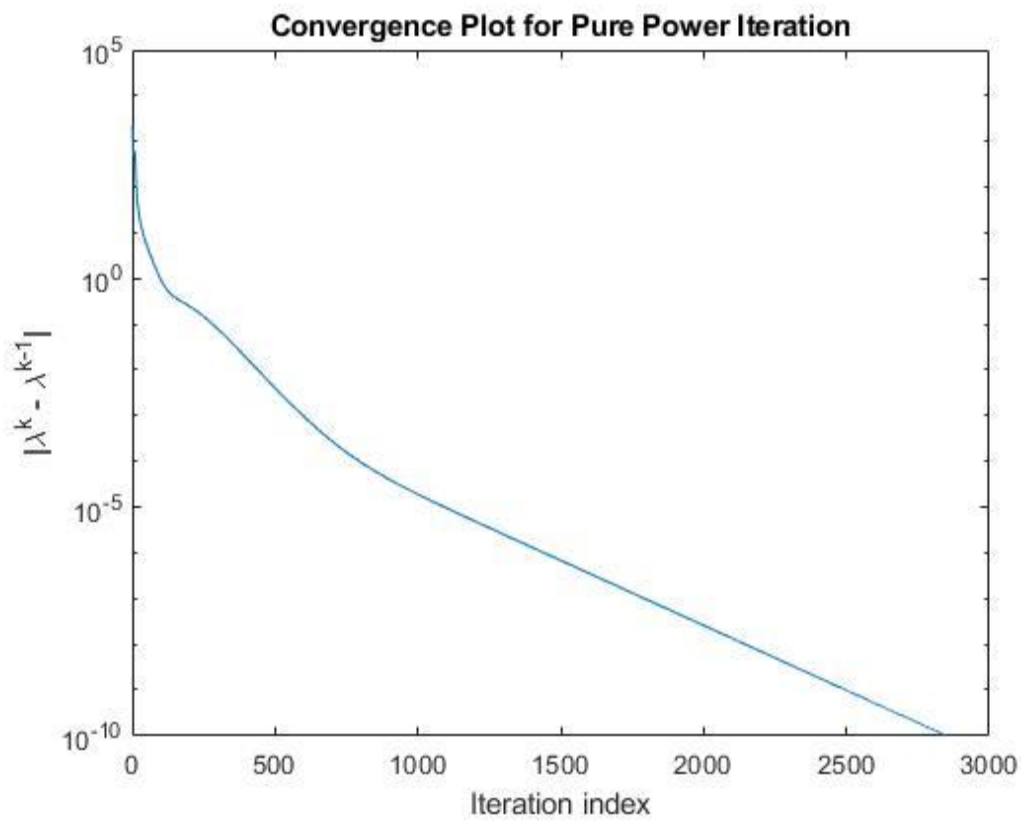
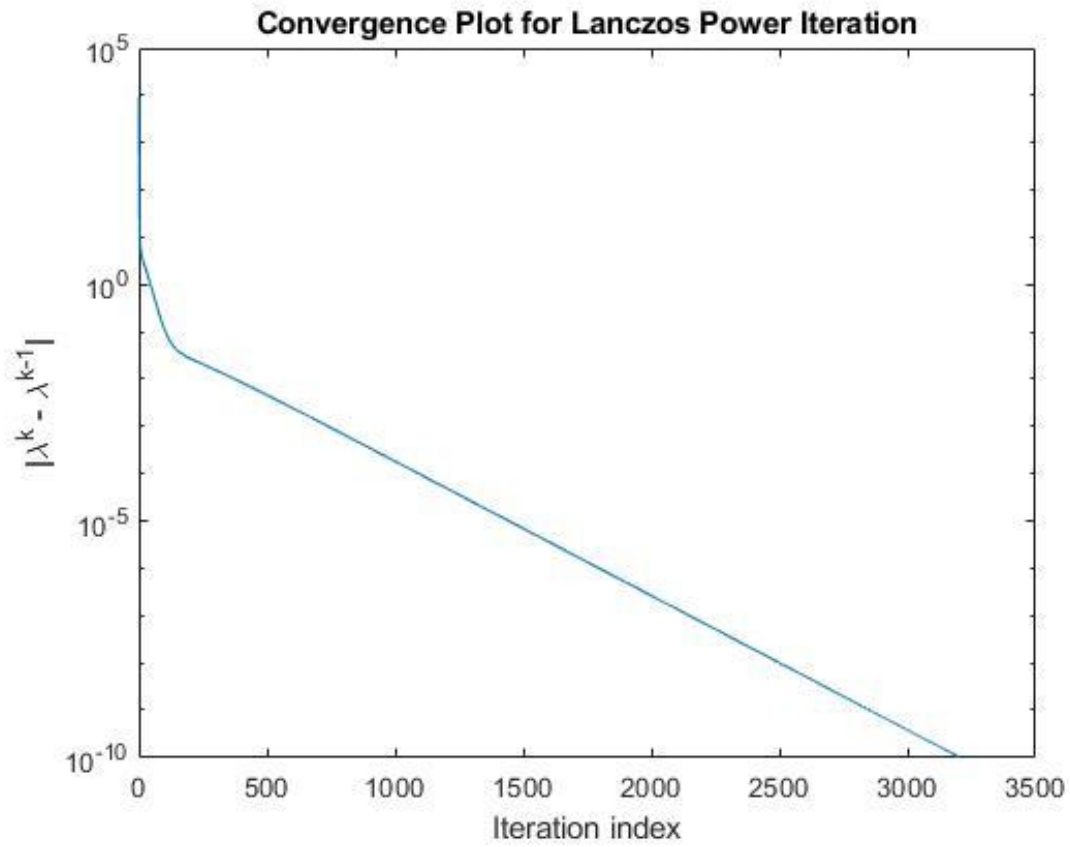
Plot 1 : with $m=30$, $tol = 10^{-2}$



Plot2: with $m=50$, $\text{tol} = 10^{-4}$



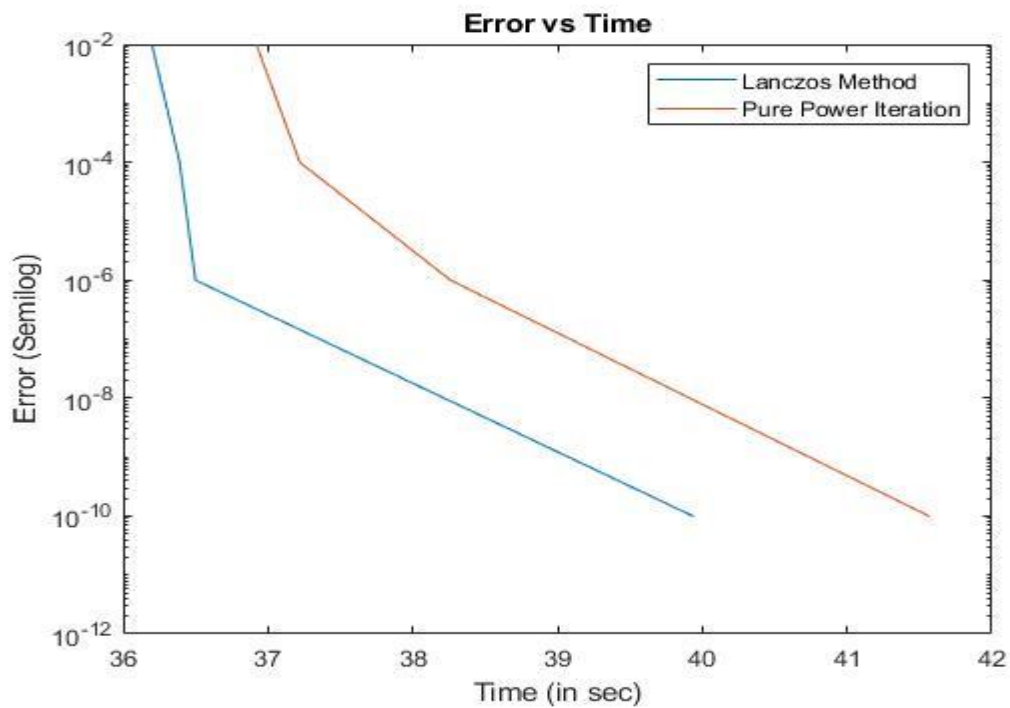
Plot 3: with $m=75$, $\text{tol} = 10^{-6}$



Plot 4: with $m=100$, $\text{tol} = 10^{-10}$

(m)	(tol)	Lanczos Method RunTime in Sec	Pure Power Iteration Method RunTime in sec	Lanczos Method Eigan Value	Pure Power Iteration Method Eigan Value	Lanczos Method Error	Pure Power Iteration Method Error
30	10^{-2}	36.20	36.92	9582.919746	9597.922393	0.0096722	0.0099617
50	10^{-4}	36.39	37.22	9598.531221	9598.595786	9.944×10^{-05}	9.953×10^{-05}
75	10^{-6}	36.52	38.26	9598.607937	9598.607937	9.970×10^{-07}	9.993×10^{-07}
100	10^{-10}	39.94	41.67	9598.608089	9598.608089	9.822×10^{-11}	9.822×10^{-11}

3. Plot the error on a semi-log scale for both schemes against execution time.



Plot 5: Error vs Run Time for Lanczos and Power Iteration method

We can observe that, to achieve the same convergence, Lanczos method takes lesser time than pure iteration method, because the smaller Hessenberg matrix of size $\sim(m \times m)$ is used to find the Eigen value, unlike in power iteration which uses very large sparse matrix.