OPTIMIZATION OF A DERIVATIVE-FREE METHOD FOR SOLVING LARGE SCALE SYSTEMS OF NONLINEAR EQUATIONS

M.K. Dauda

Department of Mathematical Sciences, Kaduna State University, Nigeria.

Abstract

Mostly, research from the work of engineers, mathematicians, physicist, computer scientist, Astronomers and other scientist are equations, which are nonlinear in nature. Researchers made significant effort in solving equations of the form F(x) = $0, x \in \mathbb{R}^n$, yet some of the methods developed do have deficiency. In this article, an improved version of a derivative free method for Solving Nonlinear Equations via a derivative free line search is presented. Interestingly, without computing any derivative, the proposed method never fail to converge throughout the numerical experiments and it satisfied the descent condition. Different initial starting points were used on a benchmark problems and the output is based on number of iterations and CPU time. The approach yield a method of solving systems of nonlinear equations that is capable of significantly reducing the CPU time and number of iteration, as compared to its counterparts. Thus, suitable and achieved the objective. The efficiency of the proposed scheme has been confirmed by the numerical results presented.

Keywords: Optimization, Decent Direction, Derivative Free, Line Search, Nonlinear Equations

1. Introduction

Consider the global optimization problem of the form

 $minf(x), x \in \mathbb{R}^n$. (1) With condition $that(x_k + \alpha_k d_k) \le f(x_k)$, where α_k is the line search and d_k is the search direction. The search direction d_k is generally required to satisfy the descent condition $\nabla(x_k)^T d_k < 0$. Supposed f be a norm function defined by a function

$$f(x) = \frac{1}{2} \|F(x)\|^2$$

where, $|| \cdot ||$ stands for the Euclidian norm. Then the unconstrained optimization problem, (1) is equivalent systems to nonlinear equations problem

 $F(x) = 0, \qquad x \in \mathbb{R}^n$

where $F: \mathbb{R}^n \to \mathbb{R}^n$ is nonlinear mappingassumed to satisfy (i) There exists $x^* \in \mathbb{R}^n$ such that $F(x^*) = 0$, (ii) F is a continuously differentiable mapping in a neighborhood of x^* (iii) the Jacobian matrix of F at x given by J(x) = F'(x) is symmetric. There are various methods for solving nonlinear equations (3) [1-6]. Among them is the Newton's method, which plays an important role due to its rapid convergence rate and decreasing the function value of the sequence. However, the methods still suffers deficiency such as computation of the derivative F' and solution of some system of linear equations at each iteration. The iterative formula of a Newton method is given by $x_{k+1} = x_k + s_k$, $s_k = \alpha_k d_k$, $k = 0, 1, \dots$,

where α_k is a step length to be computed by a line search technique, x_{k+1} represents a new iterative point, x_k is the previous iteration, while d_k is the search direction to becalculated by solving the linear system of equations below,

 $F'(x_k)d_k = -F(x_k),$

Where $F'(x_k)$ is the Jacobian matrix of $F(x_k)$ at x_k .

The drawback of the technique (4) is the need to compute the Jacobian matrix $F'(x_k)$ at every iteration, which increases the difficulty in computation; this is due to the first-order derivative of the system. Sometimes the derivatives are not available or could not be obtained exactly especially for the large-scale problems [7, 8]. Hence these requires the need for derivative free methods for solving such problems. There are many scholars [9-13] that discussed derivative-free methods, yet some problems persist. Therefore, motivated by [14] the purpose of this article is to improve the derivative free method with decent direction for solving system of nonlinear equations via derivative free line search [16, 17] and the approximations $F_o(x_k) \approx \gamma_k I$. Where I is an identity matrix and γ_k is the new parameter introduced. The method [14] retained a norm descent property without computing the Jacobian matrix with less number of iterations and CPU time that is globally convergent.

Corresponding Author: Dauda M.K., Email: mkdfika@kasu.edu.ng, Tel: +2348038893151, +2348023582223

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The next section discussed the proposed method. Section 3 reports some numerical results of the proposed method while sections 4 gives the conclusion.

2. Derivation of the Method

This section discussed the Derivation of the proposed method for solving system of nonlinear equations (3). The aim is to improve the computational performance of the existing method [14] by the concept of a derivative free line search method, and introducing the approximations $F_o(x_k) \approx \gamma_k I$, where *I* is an identity matrix and γ_k is the new parameter.

$F(x_{k+1}) \approx F(x_k) + F^o(\delta)(x_{k+1} - x_k) $ (5) where the parameter δ fulfills the conditions $\delta \in [x_k, x_{k+1}]$, and that	
where the parameter δ fulfills the conditions $\delta \in [x_k, x_{k+1}]$, and that	
$\delta = x_k + \lambda(x_{k+1} - x_k) \qquad 0 \le \lambda \le 1. \tag{6}$	
Note that the distance between x_k , and x_{k+1} is small enough. By taking $\lambda = 1$ in (6), then $\delta = x_{k+1}$. Also, if $\lambda = 0$, then δ	S =
x_k . Therefore we have	
$F^{o}(\delta) \approx \gamma_{k+1} I. \tag{7}$	
Knowing this, the expression (5) becomes	
$F(x_{k+1}) - F(x_k) = \gamma_{k+1} I(x_{k+1} - x_k) = \gamma_{k+1} (x_{k+1} - x_k). $ (8)	
Now, (8) transformed to standard secant condition [3,4].	
$\gamma_{k+1}s_k = \gamma_k,\tag{9}$	
where, $y_k = F(x_{k+1}) - F(x_k)$ and $s_k = x_{k+1} - x_k$,	
In [1], Li and Fukushima used the term	
$g_k = \frac{F(x_k + \alpha_k F(x_k)) - F(x_k)}{2} \tag{10}$	

With the aid of (10), the scheme will maintain derivative free process and avoids computing exact gradient. It is clear that when $||F(x_k)||$ is small, then the approximate gradient $\nabla f(x_k)$, is given by $g_k \approx \nabla f(x_k)$.

Consequently, g_k^T is obtainable. Thus, pre-multiplying both side of (9) by g_k^T , the relation allows us to compute the parameter γ_{k+1} in the following way:

$\gamma_{k+1}g_k^T s_k = g_k^T \gamma_k$	(11)
Taking the transpose of both side of (11) gives	
$[\gamma_{k+1}g_k^T s_k]^T = [g_k^T \gamma_k]^T$	
$\gamma_{k+1}[g_k^T s_k]^T = [g_k^T \gamma_k]^T$	
Where γ_{k+1} is a parameter.	
$\gamma_{k+1}[s_k^T g_k] = [\gamma_k^T g_k]$	(12)
Multiply both side of (12)by $[s_k^T g_k]^{-1}$ yields	
$\gamma_{k+1}[s_k^T g_k]^{-1}[s_k^T g_k] = [s_k^T g_k]^{-1}[\gamma_k^T g_k]$	
$\gamma_{k+1} = [s_k^T g_k]^{-1} [y_k^T g_k]$	(13)
Thus, direction is given by	
$d_{k+1} = -\gamma_{k+1} F(x_k)$, for $k = 1, 2, 3, \cdots$	(14)
Finally, the general scheme is given by	
$x_{k+1} = x_k + \alpha_k d_k.$	(15)
where x_{k+1} represents a new iterative point, x_k is the previous iteration, while d_k is the search	direction to be obtain u

where x_{k+1} represents a new iterative point, x_k is the previous iteration, while d_k is the search direction to be obtain using (14). The basic requirement of the line search is to sufficiently decrease the function values i.e. $||F(x_k + \alpha_k d_k)|| \leq ||F(x_k)||$. Therefore, to compute the line search α_k in (15), the line search used in [16, 17] is the best choice, since it is derivative free in nature. This is given by $\alpha_k = max\{s, rs, r^2s, \cdots\}$ (16) satisfying $-g(x_k + \alpha_k d_k)^T d_k \geq \omega \alpha_k ||g(x_k + \alpha_k d_k)|| ||d_k||^2$

Remark

It is clear that the line search is well defined.

Algorithm 1.

Step 1: Given $\gamma_0 = 1, \epsilon = 10^{-4}, x_0, (r, r\omega) \in (0,1)$, set k = 0 and $d_0 = -g_0$. Step 2: Compute $F(x_k)$ and check the stopping conditions. If yes, then stop; otherwise continue with Step 3. Step 3: Compute search direction $d_{k+1} = \gamma_{k+1}F(x_k)$ using (14) Step 4: Compute step the length α_k (using (16). Step 5: Set $x_{k+1} = x_k + \alpha_k d_k$. Step 6: Compute $F(x_{k+1})$. Step 7: determine the parameter , γ_{k+1} from (13) Step 8: Repeatk = k + 1, and go to Step 3.

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3. **Numerical Results**

This section gives the computational performance of the proposed method named M1 and two other methods named M2[14] and M3[15]respectively. In the presentation, the proposed method [15] which is the improved version of [14], is compared with classical version of [14] and the method which was compared with [14] initially. To ascertain the efficiency of the result, all the three methods were tested using the same Benchmark test problems as indicated in the Appendix. The accuracy of the methods was confirmed using different initial points. The computational codes for all the methods was written in Matlab 7.9.0 (R2009b) [18], and run on a personal computer 2.00 GHz CPU processor and 3 GB RAM memory.

Table 1: Numerical Comparison of *M*1, *M*2 and *M*3 methods for problem 1 and 2.

			M1				M2		M3			
S/N	Dim	ISP	Iter	Time	Norm	Iter	Time	Norm	Iter	Time	Norm	
1	10	0.2	7	0.029974	5.78E-04	31	0.020822	8.12E-04	10	0.006045	6.84E-05	
	100		8	0.019030	7.31E-04	36	0.007647	8.42E-04	11	0.002768	8.49E-05	
	1000		9	0.004687	9.24E-04	41	0.020315	8.73E-04	13	0.008038	4.14E-05	
	10000		11	0.114284	4.67E-04	46	0.146886	9.04E-04	14	0.061264	5.14E-05	
	100000		12	0.307812	5.91E-04	51	1.732412	9.37E-04	15	0.634616	6.39E-05	
	10	0.9	5	0.001026	7.42E-04	27	0.005472	8.78E-04	11	0.002501	3.99E-05	
	100		6	0.001188	9.37E-04	32	0.007435	9.10E-04	12	0.002964	4.95E-05	
	1000		8	0.003464	4.74E-04	37	0.018337	9.44E-04	13	0.006808	6.15E-05	
	10000		9	0.023079	6.00E-04	42	0.139060	9.78E-04	14	0.057503	7.64E-05	
	100000		10	0.228708	7.58E-04	48	1.593948	8.11E-04	15	0.603471	9.49E-05	
2	10	0.2	5	0.006740	3.86E-04	31	0.025203	*	10	0.013997	4.70E-05	
	100		6	0.002321	1.47E-04	31	0.026398	*	11	0.003108	5.21E-05	
	1000		6	0.003298	4.64E-04	31	0.064822	*	12	0.007429	6.43E-05	
	10000		7	0.025496	1.76E-04	36	1.871130	*	13	0.051324	8.75E-05	
	100000		7	0.255355	5.56E-04	35	140.6098	*	15	0.594499	4.67E-05	
	10	0.9	5	0.001098	4.58E-04	124	0.021183	9.94E-04	9	0.002826	4.95E-05	
	100		6	0.001445	1.74E-04	139	0.027641	9.69E-04	10	0.002959	3.36E-05	
	1000		6	0.003226	5.50E-04	154	0.057261	9.44E-04	11	0.006268	2.41E-05	
	10000		7	0.027749	2.09E-04	168	0.394139	9.95E-04	11	0.040932	7.62E-05	
	100000		7	0.257811	6.60E-04	183	5.132788	9.70E-04	12	0.468681	2.62E-05	

Table 2: Numerical Comparison of M1, M2 and M3 methods for problem 3 and 4.

-			M1				М2		M3			
S/N	Dim	ISP	Iter	Time	Norm	Iter	Time	Norm	Iter	Time	Norm	
3	10	0.2	5	0.007880	3.51E-04	39	0.260591	*	9	0.007014	9.11E-05	
	100		6	0.001555	2.22E-04	37	0.033213	*	10	0.002861	9.56E-05	
	1000		6	0.002903	7.03E-04	37	0.086759	*	11	0.006741	7.62E-05	
	10000		7	0.023343	4.44E-04	37	2.112099	*	12	0.056222	7.99E-05	
	100000		8	0.273506	2.81E-04	37	142.4925	*	13	0.633558	6.37E-05	
	10	0.9	4	0.000924	6.76E-04	25	0.004853	8.74E-04	5	0.001823	5.77E-05	
	100		5	0.001744	4.27E-04	30	0.007191	9.06E-04	6	0.001918	8.15E-05	
	1000		6	0.003145	2.70E-04	35	0.017076	9.39E-04	7	0.004520	2.47E-05	
	10000		6	0.019402	8.55E-04	40	0.124208	9.72E-04	7	0.031964	7.82E-05	
	100000		7	0.260632	5.41E-04	46	1.999564	8.06E-04	9	0.452263	1.06E-05	
4	10	0.2	5	0.006751	3.51E-04	39	0.052702	*	9	0.005365	9.11E-05	
	100		6	0.001168	2.22E-04	37	0.028444	*	10	0.003112	9.56E-05	
	1000		6	0.002344	7.03E-04	37	0.071258	*	11	0.005913	7.62E-05	
	10000		7	0.019958	4.44E-04	37	1.873526	*	12	0.040622	7.99E-05	
	100000		8	0.210191	2.81E-04	37	140.1881	*	13	0.467590	6.37E-05	
	10	0.9	4	0.000875	6.76E-04	25	0.004130	8.74E-04	5	0.582910	5.77E-05	
	100]	5	0.001102	4.27E-04	30	0.007141	9.06E-04	6	0.012311	8.15E-05	
	1000		6	0.002571	2.70E-04	35	0.014490	9.39E-04	7	0.031729	2.47E-05	
	10000]	6	0.018040	8.55E-04	40	0.098346	9.72E-04	7	0.112399	7.82E-05	
	100000]	7	0.194157	5.41E-04	46	1.514717	8.06E-04	9	0.357280	1.06E-05	

According to the program, the iteration is terminated if the total number of iterations exceeds 1000 or $||F(x_k)|| \le 10^{-4}$. If the method fails, the symbol "*" is used and represents failure due to; (i) Memory requirement (ii) Number of iterations exceed 1000. (iii) If $||F(x_k)||$ is not a number.

For the both algorithms in [14,15], the following parameters $\omega_1 = \omega_2 = 10^{-4}$, r = 0.2 and $\eta_k = \frac{1}{(k+1)^2}$ are used.

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			M1				М2		М3		
S/N	Dim	ISP	Iter	Time	Norm	Iter	Time	Norm	Iter	Time	Norm
5	10	0.2	2000	0.238044	0.1974	112	0.027648	9.94E-04	8	0.013481	7.55E-05
	100		2000	0.277768	0.6243	127	0.035844	9.69E-04	10	0.002987	4.72E-07
	1000		2000	0.729098	1.9741	142	0.064516	9.44E-04	10	0.010842	1.49E-06
	10000		2000	5.708984	6.2426	156	0.469077	9.95E-04	10	0.051874	4.72E-06
	100000		2000	68.898187	19.7408	171	6.504744	9.70E-04	10	0.464678	1.49E-05
	10	0.9	2000	0.226144	0.1216	30	0.006853	9.13E-04	7	0.003896	2.49E-05
	100		2000	0.289621	0.3845	35	0.006963	9.47E-04	7	0.002062	7.86E-05
	1000		2000	0.810016	1.216	40	0.021452	9.81E-04	8	0.006757	8.99E-05
	10000		2000	5.595398	3.8452	46	0.140034	8.13E-04	10	0.046637	6.12E-05
	100000		2000	65.461812	12.1595	51	1.947417	8.43E-04	11	0.527849	1.09E-06
6	10	0.2	8	0.045419	7.53E-04	28	0.016831	8.26E-04	10	0.059030	5.05E-05
	100		10	0.003527	4.06E-04	33	0.007480	8.29E-04	10	0.002714	6.48E-05
	1000		11	0.003839	5.17E-04	38	0.019901	8.56E-04	11	0.007156	5.92E-05
	10000		12	0.037682	6.55E-04	43	0.179768	8.87E-04	12	0.074682	7.03E-05
	100000		13	0.364561	8.28E-04	48	2.121760	9.19E-04	13	0.652322	8.70E-05
	10	0.9	14	0.001905	4.29E-04	*	0.348713	1.1412	12	0.003284	6.58E-05
	100		16	0.002461	6.91E-04	*	0.419275	4.4653	13	0.003611	4.66E-05
	1000		19	0.007444	4.78E-04	*	1.136226	16.4158	14	0.010835	5.15E-05
	10000		21	0.070103	6.88E-04	*	9.293672	52.3308	15	0.071868	6.31E-05
	100000		23	0.829604	8.55E-04	*	96.63993	154.1355	16	0.814324	7.83E-05

Table 3: Numerical Comparison of M1, M2 and M3 methods for problem 5 and 6.

Tables 1 gives the numerical results of problems 1 and 2. Tables 2 gives the numerical results of problems 3 and 4 while Tables 3 gives the numerical results of problems 5 and 6. In all the Tables, the Columns for each methods is titled with "S/N", "Dim", "ISP", "Iter", "Time" and "Norm" which stands for the serial number of the problem, dimension of the problem, Initial Starting Point, total number of all iterations, the CPU time in solving the problems measured in seconds and residual at the stopping point ($||F(x_k)||$)respectively.

From the Table, it is observed that all the three methods.e.M1, M2 and M3 attempt to solve the systems of nonlinear equations (3), it is worth noting that the proposed method effectively solved the problems where the other two methods fails. For instance, in problems 2, 3 and 6, M2 fails where M1 successfully and effectively solved the problems with minimal number of iterations. The method M1considerably outperforms the M2 and M3 in almost all the tested problems, since it has the least number of iterations and CPU time as compared to the CPU time and the number of iterations for the methodM2 and M3 respectively. This is quite evident that the proposed method M1 improved the computational performance of M2. The efficiency and accuracy of the proposed algorithm was clear for it solves systems of nonlinear equations (3) where mostly the method M2 and M3 fails. This achievement is due to the derivative free line search used [16,17].



Figure 1: Performance profile of M1, M2 and M3 methods with respect to the number of iterations.



Figure 2: Performance profile of M1, M2 and M3 methods with respect to the CPU time.

Figures 1, 2, and 3 gives the graphical representation of the numerical results of all the three methods. The analysis was done using the profile by Dolan and More [19]. The proposed method operates with less number of iteration and CPU time denoted by M1 represented in pink color, while M2 and M3 performs below M1 with line represented in red color and blue color respectively. Obviously, the performance of the proposed method is better in terms of number of iterations, CPU time and residual function than the other two existing methods in comparison. Another merit of the proposed method is its ability to solve problems at shortest possible time. Generally, the results shows that the proposed method is an improvement of [14] with respect to matrix storage and computational time.

4. Conclusion

In this paper, an improved version of derivative-free decent method via acceleration parameter for solving systems of nonlinear equations is given. The aim of this research is to improve the computational performance and convergent rate of existing methods, thereby reducing the computational time at each iteration. The aim of this research has greatly achieved by doing away with a computation of the Jacobian inverse during the iteration process. It is a fully derivative-free iterative method, which retain the global convergence of the classical method under some appropriate conditions. The Numerical results reported using a set of large-scale test problems has shown that the proposed method is practically effective. The method is valid in terms of derivation, reliable in terms of number of iterations and accurate in terms of CPU time. Conclusively, the proposed method is recommended for solving large-scale system of nonlinear equations of the form $F(x) = 0, x \in \mathbb{R}^n$.

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Appendix.

The following test problems are used for the numerical experiments. The mapping *F* is taking as $(x) = (f_1(x); f_2(x); \dots; f_n(x))^T$; and $x = (x_1; x_2; \dots; x_n)^T$. The test problems 1,2,3,4,5 are from [20] while problem 6 is from [21].

Problem 1.

 $f_1(x) = (x_1 - x_2^2)(x_1 - \sin(x_2)),$ $f_2(x) = (\cos(x_2) - x_1)(x_2 - \cos(x_1)).$ **Problem 2.** $f_1(x) = -x_1 + 0.5x_2^2 - 1.5,$ $f_2(x) = -x_2 + 0.605 \exp(1 - x_1^2) + 0.395$ **Problem 3.**

$$f_1(x) = x_1^2 - 2x_1 + \frac{1}{3}x_2^3 + \frac{1}{3}$$

$$f_2(x) = x_1^3 - x_1x_2 - 2x_1 + 0.5x_2^2 + 1.5$$
Problem 4.
$$f_1(x) = x_1,$$

$$f_2(x) = \frac{10x_1}{x_1 + 0.1} + 2x_2^2.$$

Problem 5.

 $f_1(x) = -13 + x_1 + ((-x_2 + 5)x_2 - 2)x_2,$ $f_2(x) = -29 + x_1 + ((x_2 + 1)x_2 - 14)x_2.$

Problem 6.

 $F_1(x) = x_1(x_1^2 + x_2^2) - 1,$ $F_i(x) = x_i(x_{i-1}^2 + 2x_i^2 + x_{i+1}^2) - 1,$ $F_n(x) = x_n(x_{n-1}^2 + x_n^2).$ $i = 2,3, \dots n - 1.$

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