Chaotic Convergence of Newton's method

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Abstract—Problem statement: In 1680 Newton proposed an algorithm for finding roots of polynomials. His method has since evolved, but the core concept remains intact. Here we briefly review this evolution, and consider the question of convergence.

Methods: Newton's method may be applied to any complex analytic function, thus holds for solutions of linear differential equations. The derivation is based on a Taylor series expansion in the Laplace frequency $s = \sigma + j\omega$. The convergence of Newton's method depends on the important concept known as the Region of convergence (RoC).

Findings: Under certain conditions, non-linear (NL) limit-cycles appear, resulting in a reduced rate of convergence to a root. Since Newton's method is inherently complex analytic (that is, linear and convergent), it is important to establish the source of this NL divergence. We show that this NL effect is due to violations of the Nyquist Sampling theorem, also known as aliasing. Aliasing is a well-known concept in discrete-time signal processing, due to the sampling a signal at less that twice its highest frequency. When a time signal is under-sampled, frequencies above the sampling frequency are shifted down in frequency. This is the definition of this nonlinear phenomena, known as aliasing, which follows from the reduced-sampling. Here the conditions and method for uniform convergence are explored.

Conclusions: The source of the nonlinear limit-cycle is explained in terms of aliasing. We numerically demonstrate that reducing the step-size always results in a more stable convergence. The down side is that it always results in a sub-optimal convergence. It follows that a dynamic step size would be ideal, by slowly increasing the step-size until it fails, and then decreasing it until it converges. A balance of the two methods seems like a potential solution, but this remains unproven.

I. INTRODUCTION

Newton's method (NM) is a venerable method for finding the roots of polynomials. However its utility has been questioned. First, and most important, does his method always converge? From numerical experiments, it does converge for the vast initial guesses. Thus a key and important question: does the convergence depend on this initial guess? This question was carefully evaluated by Willkinson, who studied conditions of sever divergence.¹

Assuming it does converge, what are the necessary conditions for convergence? In the following discussion we assume a monic polynomial of degree N. The fundamental theorem of algebra states that every polynomial of degree N has N roots, which are typically complex $s_n \in \mathbb{C}$.

Because Newton's algorithm converges within the RoC for any complex analytic function, it converges to one of the roots of the polynomial when the nearest root s_r ($s \in \mathbb{C}$) inside the RoC out to the nearest pole. This follows because every complex-analytic point on the complex plan has a region of convergence (Allen, 2020). We show that on the boundaries of the RoC regions, the method becomes hyper-sensitive to the initial condition (i.e., guess s_0).

We propose a complex *adaptive step-size* $\eta = ae^{j\phi} \in \mathbb{C}$ which we adaptively adjusted, greatly reducing, even removing the nonlinear effects of aliasing. Introducing this adaptive step-size (η) is known as the *damped Newton's method* (Galántai, 2000, p. 25). In the limit as $\eta \to 0$, the NL aliasing is avoided. This naturally follows from the complex-analytic properties of Newton's method.

Under some special conditions, a minor modification in the initial guess s_0 (n = 0) can result in the n + 1estimate of the root (s_{n+1}) to cross (or not cross) an RoC boundary, resulting in NM to divert its initial path (or not) to a alternate root. When this happens, the change in the step $\delta = s_{n+1} - s_n$ is unpredictable, and possibly even chaotic. It is this condition that is the source of a convergence instabilities due to aliasing, possibly leading to a limit cycle. This chaotic behavior is a main topic of this document.

The N regions of convergence (RoC) are investigated. From the definition of the RoC, these contiguous natu-

¹https://en.wikipedia.org/wiki/Wilkinson's_polynomial

rally existing regions are defined over all $s_0 \in \mathbb{C} < \infty$. That is, every possible s_0 comes from one of the N RoC regions. This naturally happens as $\eta \to 0$, since Newton's method is complex analytic (the step size is the ratio of two polynomials with different roots). The magnitude of $|\eta| = a$ may also be manually reduced to avoid crossing the boundary between two RoCs. As a is reduced, the trajectory naturally moves away from the poles and the smaller it is, the greater the effect. For this reason adaptively setting $|\eta|$ is likely the optimum balance, to minimize the computation while avoiding NL aliasing, which results from any crossing of RoC boundaries.

As an alternate to reducing $|\eta|$, one can modify its angle ϕ , redirecting the trajectory away from any RoC boundary, so to avoid crossings it. Note that we have not yet implemented this idea.

We show that when $\eta = 1$, depending on s_0 , the solution can cross an RoC boundary (i.e., diverge). In such cases the target root will change, resulting in a chaotic trajectory. Examples are provided. Depending critically on s_0 , as long as the RoC region remains the same, every iteration converges.

Newton's method (NM) a is a venerable complexanalytic mathematical algorithm for finding roots of any monic polynomial $P_N(s)$, where N is the degree and $s = \sigma + \omega j \in \mathbb{C}$ is the Laplace frequency. However the convergence properties of NM are controversial² (Stewart, 2012, p. 347). In this report we shall investigate why such a controversy developed, and discuss how to assure convergence. In our experience, given some care, the method always converges to a root.

Every initial guess s_0 on the plane of a complex analytic function is uniquely associated with one of the N roots of that function, which in turn are associated a unique region of convergence (RoC). This follows from the complex analytic property of a function (those that may be expanded in a complex-analytic Taylor series). When the trajectory of NM jumps to a different RoC, corresponding to a different root, it has been interpreted as a failure to convergence. What then happens in the examples presented here, the iteration still converges, but to a different root. NM contains properties that are similar to *dynamic analysis*, a mathematical science first introduced by Poincaré.

This question of the convergence of NM was recently explored in Allen (2020), where no instability or limitcycles were observed. An explanation is due: Newton's method was modified by applying an *adaptive stepsize* η , (Galántai, 2000, p. 25), a widely recognized contemporary technique in the engineering numerical analysis literature.³

A properly chosen adaptive step-size stabilizes the convergence, by forcing the convergence to remain in the target RoC. We show that its easy to detect when the divergence of the step, which should be monotonically decreasing. The onset of limit cycles are easily detected, and easily stabilized by modifying η .

Here we show that random jumps and limit-cycles can occur when $\eta = 1$. When the adaptive step-size η is sufficiently small, we show that the iteration always converges. The key here is to adaptively modify the magnitude of η , thereby constraining the trajectory to the initial RoC.

In Allen (2020, Fig. 3.2), two examples were provided using a fixed step-size ($\eta = 0.5$) and a random initial guess. The details of the step-size used by Allen (2020) was not discussed. One of these figures is presented in Fig. 1 (LEFT).

While most of the curves seem to converge to a root, there are some small percentage (e.g., 1%) of cases where the trajectories take huge jumps to random locations in the complex plane. We shall show that these jumps occur when the trajectory approaches any of the poles of Newton's method, that is, at the roots of $P'_N(s) = \frac{d}{ds}P_N(s)$. Near a pole the step can be arbitrarily large, depending on how close the step comes to the pole (Boas, 1987). We shall show that the poles are the source of the limit cycles, which are easily detected.

In Fig. 1 (LEFT), the five RoC regions are color coded, with each RoC region associated with one of the N roots. Due to the complex analytic nature of an RoCs, every point in the RoC is a valid initial condition. However this is limited by the numerical accuracy of the computer. Also the convergence depends on the size of the steps, defined as $s_{n+1} - s_n$, $s \in \mathbb{C}$, $n \in \mathbb{N}$, which typically decreases in magnitude with n. An exception occurs if s_{n+1} approach one of the N-1 poles of NM, causing the step to abruptly diverge. The properties of this small subset of initial conditions depends critically on the *step size* $|\eta|$, which is a key topic of this article.

For most initial guess $s_0 \in \mathbb{C}$ the iteration simply converges to a root, independent of $|\eta| \leq 1$. In fact for most starting values the solution converges for $|\eta| = 1$. However for s_0 values near the RoC boundary between two roots, the dependence is highly dependent on $|\eta|$, and can even be chaotic. This happens when s_0 defines a path that heads directly for a pole. In these cases the trajectory will be hypersensitive to both s_0 and $|\eta|$. The RoC regions are well defined non-overlapping complex-valued analytic regions. When s_0 is close to the RoC boundary, the convergence of NM critically depends on the magnitude and angle of the complex

²https://en.wikipedia.org/wiki/Newton's_method#Failure_of_the_method_to_ converge_to_the_root

³https://en.wikipedia.org/wiki/Adaptive_step_size



Fig. 1: Example LEFT: 1a This figure is taken from Allen (2020, p. 78). It is a plot of a thousand trajectories for Newton's method, with a random initial guess, taken from the complex plane, between [0,5] along the real axis and ± 1.5 along the imaginary axis, for a polynomial having N = 5 complex coefficients. Note the long straight lines that occasionally appear in the figure EXAMPLE **RIGHT:** 1b shows the poles and zeros of the polynomial having coefficients C = [1, 0, 0, 0, -1, -1] with random starting points, for 200 iterations of Newton's Method. In this case the roots (0) and poles (\times) are superimposed on top of the trajectories of Newton's method. An adaptive step-size of $\eta = 0.1$ is used to reduce the NL aliasing. **RIGHT:** AD ROOTS IN TITLE; LEFT: ADD ϕ_X .

Figure	$P_N(s) =$	$\Re s_r$	$\Im s_r$
Fig. 1a; LEFT	$s^5 - (13 + 0.5j)s^4 + (66.25 + 5j)s^3$	[4, 3, 3, 2, 1]	[1, -2, 2, -1, 1]/2
	$-(164.25+20.125j)s^2$		
	+(195+38.75j)s - 87.5 - 31.25j		
Fig. 1b; RIGHT	$s^5 - s - 1 \leftrightarrow [1, 0, 0, 0, -1, -1]$	[1.17, 0.18, 0.18, -0.76, -0.76]	[0, 1.08 -1.08, 0.35, -0.35]

TABLE I: Properties of the polynomials for the left and right panels.

step-size η . Even when $|\eta| \ll 1$, the convergence can become nonlinear (NL), resulting in a chaotic path. These observations are supported by several detailed numerical examples.

In Fig. 1 (LEFT), the red region, corresponding to the root at (2.0 - 0.5j), has a long narrow "RoC stream" for initial guesses s_0 south-east of (4.5 - 1j). There is a second green narrow neighboring related parallel stream, just north of the red stream, for initial values s_0 at (5 - 1j), which is in the RoC of root (3 + 1j).

While it may seem obvious given Fig. 1, I am not aware of any discussion of such distortion of the RoC's. The conditions for Fig. 1 are provided in Table I.

A. Convergence of Newton's method

Given the monic polynomial of degree $N \in \mathbb{N}$ $P_N(s) = s^N + c_{N-1}s^{N-1} + c_{N-2}s^{N-2} + \cdots + c_0$, and its derivative $P'_N(s) \equiv dP(s)/ds$ of degree N-1, we may express Newton's method as the ratio of monics

$$s_{n+1} - s_n = -\frac{\eta}{N} \frac{P_N(s_n)}{P'_N(s)}$$
 (I.1)

Our key idea is to define the ratio of monic polynomials as the *step-size* $S_N(s_n)$, thus modifying the syntax

of Newton's method. Equation A.5 then becomes

$$\frac{s_{n+1} - s_n}{\eta} = \frac{1}{N} S_N(s_n)$$

Scaling $P'_N(s)$ as a monic does not alter its roots.

Taking the limit $\eta \rightarrow 0$ results in the complex-analytic expression for NM

$$\frac{ds}{d\eta} \equiv \lim_{\eta \to 0} \left(\frac{s_{n+1} - s_n}{\eta} \right) = -\frac{1}{N} S_N(s_n).$$
(I.2)

The right hand side $S_N(s_n)$ is the reciprocal of the logderivative of $P_n(s)$ expressed as monics, and the left hand side is the slope of the Laplace frequency $(s = \sigma + j\omega)$ wrt η . Note that Eq. I.2 has no cellular structure along the RoC boundaries, due to its complex analytic properties.

B. What is going on?

In the limit as η goes to zero, the RoC boundaries are well defined analytic regions. But for small $\eta \neq 0$, no matter how small, the boundaries are cellular, becoming smooth only in the limit at zero. This cellular structure is always present even for the smallest nonzero values of η . Insight into how this happens is explained by example in Figs. 2, 3.

In Fig. 3 an infinitesimal change in s_0 leads to large jumps into different RoC. This is best shown by two starting values at $s_0 = 1 + \beta j$ for β just above 0.5 and again just below 1.0. The effect occurs only when the trajectory heads directly at a pole. Any starting value s_0 that goes directly at one of the poles can jump to a random RoC.

We have defined η as the *adaptive step-size*, because we can set η to modify the *step-size* S_N .— This result follows from a mathematical property cited by (Boas, 1987), that the entire plane may be found in the neighborhood of every pole.

Given this mapping, infinitesimal changes in the starting points which head directly at a pole, are reassigned to a random RoC, due to this analytic mapping.

a) In summary: Determining the RoCs for NM by analytic methods seems difficult, since the function $S_N(s_n)$ has poles, confounding the locations of the RoC boundaries. Based on Fig. 1, the RoC are complicated. If s_n approaches one of these the poles, the update can become arbitrarily large, depending on how close s_n is to the pole. If the *adaptive step-size* is within the RoC, this will not occur. When the value of s_{n+1} falls outside the RoC there can be an arbitrary increase in step-size. Normally this will not happen, since when s_n approaches a pole s_{n+1} is naturally "pushed" away from the pole, as may be seen in Fig. 2 (black circles).

II. EXAMPLES OF NEWTON'S METHOD

a) Example 1: We start with the monic polynomial of Example 1b, Fig. 1 (RIGHT), for N = 5,

$$P_5(s) = s^5 - s - 1. \tag{II.1}$$

In this case monic $\frac{1}{5}P'(s) = s^4 - 1/5$, has four roots, shown as bold \times symbols. These are the poles. The black bold circles are the desired zeros of $P_5(s)$.

As shown in Example 1b, Fig. 1 (RIGHT), given the initial guess s_0 and step-size η , as $n \to \infty$, s_n approaches a unique root s_r . Complex s_{n+1} is the n+1the estimate of the root given the *n* estimate s_n , defined by Eq. I.1.

The example shown in Fig. 2 is a zoomed-in version of Example 1b, Fig. 1 (RIGHT). When the step-size is $\eta = 0.5$ (red squares), s_{n+1} over-shoots the pole, resulting in a limit cycle. When the step-size is reduced to $\eta = 0.1$ (black circles), the trajectory is stable (avoids the pole) and converges smoothly to the complex root.

A. Discussion of Example 1b of Fig. 2

To study the convergence and limit-cycles it is helpful to use numerical methods and look at specific step-sizes and starting points, and explore for limit cycles.



Fig. 2: This is a zoomed-in version of Fig. 1b (RIGHT), presented as a colorized plot (Allen, 2020, p. 168) of $S_N(s)$, for $P_5(s) = s^5 - s - 1$ ([1, 0, 0, 0, -1, -1]). The magnitude of $S_N(s)$ is coded by the brightness, and the phase ($\angle(L_N(s))$) by the color (hue). Dark regions are zeros (roots of $P_n(s)$) while the white regions are poles of $S_N(s)$ (roots of P'(s)). Two trajectories of Newton's method are shown, as the black circles and red squares. The initial value for both cases is $s_0 = 1 + 0.75_2$. The black circles correspond to $\eta = 0.1$, while the red squares ($\eta = 0.5$) form a brief limit cycle. The vertical white lines are at {-1.0, 0, 1.0} and the horizontal lines are at {0, 1.0}. The polynomial coefficients are $P_5(s) = [1, 0, 0, 0, -1, -1]$, with roots $s_r = [1.167, 0.181 \pm 1.084_j], -0.765 \mp 0.3525_2$. The real poles (roots of $P'_5(s_r) = 0$) are $\pm 1/5^{0.25}$, while the imaginary poles are at $\pm 0.6687_j$.

Figure 2 shows two paths with the same initial condition $s_0 = 1 + 0.75j$ with different step-sizes. The utility of the reduced step-size is clear from the Figure.

In practice, once near a pole, it takes only a few additional steps to limit-cycle, recover and finally converge. Once near a zero, fewer than 10 steps typically give double-precision floating-point machine accuracy.

As discussed in the figure caption, if s_n is close to a root s_r of P'_N (i.e., a pole), the recursion dramatically fails, because the step becomes arbitrarily large, forcing the next trial to a random location in the *s* plane, denote \tilde{s}_r . In such cases the solution typically converges to a different root (RoC). It is not difficult to detect these large random steps by monitoring $|s_{n+1} - s_n|$, which must monotonically decrease.

If we start the iteration with the larger step size, the path develops into a NL limit-cycle near the pole at -0.669. It is a combination of the large steps and the proximity to the real pole that results in the nonlinear limit-cycle. On the 10 step it comes out of the limit cycle, and after 10 more steps, has converged to the root. When the step-size η is reduced from 0.5 to 0.1, as the path approaches the pole, it moves away, avoid the limit-cycle. With steps sizes of 0.2, 0.3 it becomes captured by the pole. The black circles show a smooth analytic trajectory, while the red-squares are chaotic.

With the step-size of 0.9 (not shown), the trajectory is similar to that of 0.5, but after 5 steps it is in well within a different RoC, corresponding to the zero at -.8 - .3j. After 20 steps, the error is less than 1%.

In summary, given a larger step size it still converges, but much more slowly, since the NL becomes greater. Thus the convergence time may be a crude metric of the NL. The smoothness of the trajectory may be more appropriate. This NL result is due to the reduced sample step size, also known as *aliasing*.

a) Example 2: This example is for $P_3(x) = x^3 - x + 1$ (C = [1, 0, -1, 1]), an example where Newton's method appears to fail. This example has two imaginary roots, $0.66236 \pm 0.56228j$, and a real root -1.32472. If the initial guess is taken to be $s_0 = 1$, the recursion proceeds using real arithmetic (Matlab and Octave). Due to the restriction that the computation is real, the solution is forced to the real line, where it limit cycles between 1.155 and 0.694. The iteration cannot converge if $s_r \in \mathbb{R}$ and $s_0 \in \mathbb{C}$ (Allen, 2020).

If $x_0 = j$, the solution converges in 3 steps to the upper complex root $x_3(3) = 0.639379 + 0.509792j$ If one starts the iteration with an imaginary component at $1 + j10^{-6}$, the iteration converges to the imaginary root in 13 steps.⁴

b) In summary: Roots $s_r \in \mathbb{C}$ may be found by a recursion that denotes a sequence $s_n \to s_r \in \mathbb{C}$, $n \in \mathbb{N}$, such that $P_N(s_n) \to 0$ as $n \to \infty$. As shown in Fig. 2, solving for s_{n+1} using Eq. A.5 always gives one of the roots, due to the analytic behavior of the complex logarithmic derivative $P'_N/P_N = d\ln(P_n(s))$.

When there are no limit cycles, each step (s_{n+1}) is always closer to the root, finally converging to the root for increased n. As it approaches the root, the linearity assumption becomes more accurate, resulting in a rapid convergence.

Even for cases where fractional derivatives are involved Newton's method will converge since the logderivative linearizes the equation (Allen, 2020, p. 197, #5).

B. Newton's method applied to functions other than polynomials

a) Example 2: Example of Plank's formula for Black Body radiation.

Planks famous BB radiation formula is (Kuhn, 1978; Allen, 2020)

$$S(\nu) = \frac{\nu^3}{e^{h\nu/kT} - 1}.$$
 (II.2)

In this historically important example, because the function is real (it is not complex analytic), the spectrum only has one pole, at $\nu = 0$. This formula is known to match the experimental data of the smoothed (nonanalytic) black-body power spectrum (Haar, 2016).

If we replace the real frequency ν with the negative Laplace frequency $-s = -\sigma - \omega j$, Eq. II.2 becomes

$$S(-s) = \frac{-s^3}{e^{-\hbar s/kT} - 1},$$
 (II.3)

⁴Octave program: ./DEMO-MATHINSIGHT/NEWTONEXP1.M

which is complex analytic, thus has a causal inverse Laplace transform. To use Newton's method we must compute NM update $L(s_n)$ where s_n is the present estimate of the root, defined as the the reciprocal of the logarithmic derivative (see derivation in Apdx. A). Taking the log followed by its derivative wrt $s \in \mathbb{C}$, gives

$$\frac{1}{L(s)} \equiv \frac{d}{ds} \ln S(-s)$$
$$= \frac{d}{ds} [-3\ln s + \ln(e^{-\hbar s/kT} - 1)],$$
$$= -\frac{3}{s} - \frac{\hbar}{kT} \cdot \frac{e^{-\hbar s/kT}}{e^{-\hbar s/kT} - 1}.$$

Thus there is a first order pole at s = 0 and poles at $h\nu_n/kT = 2\pi n$ for $n \in \mathbb{N}$. The discrete frequencies account for the eigen-modes in the black body radiation, as discussed by Kuhn, Plank and Einstein (Haar, 2016). Equation II.3 is causal, since it has a causal inverse \mathcal{LT} (Allen, 2020, p. 321)

$$-3u(t) - \frac{\hbar}{kT} \sum_{n=1}^{\infty} \delta(t - n\frac{\hbar}{kT})$$
(II.4)

The application of NM to Plank's famous formula can be used to make it complex analytic, by replace ν with the Laplace frequency $s = 2\pi\nu_n \jmath$ and h by \hbar . It is well established that complex analytic functions of the Laplace frequency s are *causal* (zero for negative time). In the case of Eq. II.3, S(-s) is causal, due to the Laplace transform relation of the exponent

$$\delta(t-\tau_o) \leftrightarrow e^{-s\tau_o}.$$

Here the time delay $\tau_o = \hbar/kT = (6.63/2\pi k) \cdot 10^{-11}$ [s], 6,280 [GHz], $\lambda \approx \frac{\pi}{2} 10^{-11}$ [m], or $\frac{\pi}{20}$ [Å], and T° [K] is the temperature.

Newton's method uses the reciprocal of L(s) to find the s_r ($S(s_r) = \infty$), given by

$$N(s_r) = 1 - e^{-\hbar s_r/kT} = 0.$$
(II.5)

There are an infinite number of such roots, since the roots are $\hbar s_r/kT \approx 2\pi j$. These poles are the missing discrete spectral lines (atomic resonances), required by quantum mechanics.

Applying Newton's method gives

X

$$x_{n+1} = x_n - \frac{e_n^x - 2}{e_n^x} = s_n - (1 - 2e^{-s_n}).$$

Since e^x is entire, there are no convergence issues.⁵ Since $x \in \mathbb{C}$, the imaginary part quickly decays to zero, and depending on the starting condition, approaches one of the infinite number of solutions, within a few steps.

⁵https://www.quantamagazine.org/how-mathematicians-make-sense-of-chaos-20220302/



Fig. 3: This numerical experiment for polynomial coefficients [1, 0, 0, 0, -1, -1] (the same as shown on the right panel of Fig. 1) having a step-size of 0.1), reveals the inner workings of Newton's method. We number the roots counter-clockwise from 1-5, with $s_1 =$ 1.2, $s_2 = 0.18123 + 1.08395\jmath$ and $s_5 = s_2^*$. Seventeen different starting values have been carefully chosen, to determine the RoCs. All the starting values are of the form $s_0 = 1 + j\beta$, where each β and the converged root are indexed in Table II. The root index goes from 1 to 5, counting counter clockwise from the eastern-most root. The scattering angle is determined by the residue of the scattering pole. Each curve is labeled twice, once at the starting point and again at another point on the trajectory. The carefully evaluated case is for starting points between 1 + 0.62j and 1 + 0.5999j, which converge to dramatically different RoCs, due to squarely hitting the positive real pole at $s_0 = 1 + j \, 0.6 \pm 0.001$.

TABLE II: Table of starting values $s_0 = 1 + \beta_j$ use in Fig. 3, along with the RoC targeted root index, defined as #1 for the real root at $0.2^{1/4}$. Root #1 converges from $s_0 = 0 + 1.25_j$, Root #2 is defined by counting counter-clockwise from #1, at $0.18 + 1.08_j$, starting fro $s_0 = 1 + 0.69_j$. Root #3 also converges from three values of β . Root #5 is the most carefully explored, starting from $1+\beta_j$. It is shown to converge to roots 1, 3, 4, 5, but not 2, which is reachable from very selective values of β . For other choices of β_0 , all 5 roots can be reached, as shown in Fig. 3 for $\eta = 1$. (For numeric values see _/M/ZvizDemo.m)

β	root
0.25, 0.4	#1
0.95, 0.99, 1.1	#2
0.69, 0.92, 0.93	#3
0.65, 0.632	#4
0.63, 0.631	#5

C. Example 3:

The impact of s_0 is shown in greater detail in Example 3, as shown in Fig. 3. When the value to s_0 is finely tuned, such that the trajectory intercepts a pole, a host of NL limit-cycles are exposed.

The Gauss-Lucas theorem⁶ comes into play at this point (Allen, 2020, p. 81). This theorem says that the convex hull of the roots of a polynomial bound the roots of its derivative. This theorem is relevant to the

convergence of Newton's method. Galántai (2000) has 75 relevant citations, many citing the same problems addressed here. The key to avoiding the troublesome limit-cycles is to detect them, and then reduce the stepsize.

The following quote is from Galántai (2000, p. 39):

The possibility that a small change in s_0 can cause a drastic change in convergence indicates the nasty nature of the convergence problem. The set of divergence points of the Newton method is best described for real polynomials.

As demonstrated in Fig. 3, we agree with Galántai's first point. His second seems vague: Is a "real polynomial" one with real coefficients, real roots, or both?

For example in Fig. 1, the red "stream" corresponding to the root near (2-0.5j) has an interesting long narrow "RoC-stream," converging from the lower-right quadrant, first seen at (4.5 - 1j). There is a second green RoCstream just north of the red stream, first seen near (4.5 - 0.9j). Thus a small change in the starting value s_0 robustly converges to a totally different root.

I am not aware of any discussion in the literature of this distortion of the RoC regions, bound to Newton's method. Presently I know of no way to predict the conformal remapping of the RoC regions for NM, other than tracking them, as done here. It seems likely that methods for doing must exist using modern analysis techniques (see Appendix III-A).

In the example of Fig. 3,

$$s_{n+1} = s_n - \frac{0.1}{5} \cdot \frac{s_n^5 - s_n - 1}{s_n^4 - 1/5},$$

for 17 carefully chosen initial guess $s_0 \in \mathbb{C}$. For readability, each trajectory is color-coded either red or blue.

a) Nonlinear limit cycles: It is well documented that limit cycles are nonlinear. Newton's method on the other hand is a linear recursion equation, with poles and zeros in the complex plane. The obvious research question is "Why does the complex-analytic linear equation become nonlinear?" We show how the these NL limitcycles may be easily avoided by removing (linearizing) the NL recursion once it is detected.

The suggested procedure will result in a net convergence speed-up, because the NL limit-cycle adds meandering NL steps to the recursion. If you experience a slowdown, try changing the step-size angle. This may be a panacea, since this is a 'local' modification that deals directly with the main problem of being on a RoC boundary. If you find an angle that reduces the chaos, then your moving in the right direction, away from the RoC boundary. This method seems obvious, yet unexplored. If your inside a "cell," it will fail and you need to reduce the magnitude as well.

⁶https://en.wikipedia.org/wiki/Gauss-Lucas_theorem



Fig. 4: Four colorized plots for $P_N = [1, 0, 0, 0, -1, -1]$ showing the N = 5 regions of convergence and two trajectories, for $s_0 = 1.8 - 1.5_J$ and $-1.95 - 0.1_J$. The four step-sizes are $\eta = \{1.0, 0.5, 0.2, 0.1\}$ (note that the imaginary axis is reversed). The fractal regions reside on the RoC boundaries, the sizes of which depend on the step-size, with the step-size of $\eta = 1.0$ (Upper-Left) resulting in large fractal regions. Reducing the step-size to $\eta = 1/2$ dramatically reduces the fractal regions. For $\eta = 0.1$ they almost disappear, except at $0.5 + 0_J$. In the dark RoC (purple) corresponding to root $-0.76 - 0.352_J$, two trajectories are shown. For the step-size of 1, a limit cycle is seen, for both trajectories. For the step-sizes [0.5, 0.2, 0.1], there are no limit cycles. As the trajectories approach the negative real pole, labeled as the red \times , they head for the root at $-0.76 - 0.352_J$. In summary: 1) limit cycle is easy because the path reverses (oscillates). An obvious method for avoiding limit cycles is to detect that the boundary has been crossed, corresponding to a different root, and restart with a reduced step-size, at step s_n or s_{n-1} .

b) Ratios of monics as NM: It can be notationally useful to define the *adaptive step-size* $S_N(s)$ as the ratio of monic polynomials

$$\frac{s_n^N + c_{N-1}s_n^{N-1} + \dots + c_0}{s_n^{N-1} + \frac{N-1}{N}c_{N-1}s_n^{N-2}\dots + \frac{1}{N}c_1} = \frac{1}{N}\frac{P_N(s_n)}{P'_N(s_n)}.$$
(II.6)

Using this trick we can absorb the factor of N into the definition of $\eta \equiv \frac{1}{N} e^{\phi j}$. Increasing N from 1 to 0.1 dramatically improves the convergence, while the poles (and zeros) of the step-size are unmodified.

Figure 4 quantifies the effect of reducing the step by up to 1/N ($|\eta| = [1, 1/2, 1/5, 1/10]$). For the largest step size, the trajectory of red squares in Fig. 2 limit cycle. This natural reduction in step-size by N, due to expressing the step-size as the ratio of monics, is dramatic. Given s_n , everything on the right is known; thus when s_n is within the RoC, s_{n+1} will converge to a unique root of $P_N(s)$ as $n \to \infty$. For sufficiently small step-size, the roots of Eq. II.6 are the solution to a linear difference equation, the simplest example being (Galántai, 2000)

$$s_{n+1} = s_n - \frac{\eta}{N} S_N(s_n).$$
 (II.7)

Introducing the *adaptive step-size* gain $|\eta| < 1 \in \mathbb{C}$, linearizes the iteration when s_n is in the neighborhood of a pole.

Near any pole, the step-size $|S_N(s_n)|$ can become arbitrary large, introducing aliasing (non-linearity) into the iteration.

III. SUMMARY AND DISCUSSION

A. The role of the step-size

In the derivation of NM we modified Eq. A.5 with the adaptive step-size $\eta < 1$, to obtain Eq. I.1. The effect of the reduced step-size is to force the trajectory to be more sensitive to the influence of the poles, rather than stepping over them. The modification of the step size S_N by η is an important modification to Newton's method. The smaller step-size can eliminate the nonlinear limit-cycles, as seen in the example of Fig. 4.

When the initial value for the iteration s_0 is close to the cross-over of two RoCs, $s_n \rightarrow s_{n+1}$ can cross over an RoC boundary, changing the limit point (root it converges to). A limit cycle can happen when s_n comes close to one of the poles of $S_N(s_n)$. At a pole, the value of S_N can become arbitrary large, causing the unmodified ($\eta = 1$) update $S_N = s_{n+1} - s_n$ to fail to satisfy the required RoC convergence condition (Eq. A.6).

One strategy for detecting the pole is to look at the magnitude of the step $(|\eta|)$. If $|\hat{s}_{n+1} - s_n| > 1$, the RoC condition has failed. The step must then be reverted back to s_n , and the step-size reduced, and s_{n+1} recomputed. This then repeated until the RoC condition $(|s_n| > |s_{n+1}|)$, thus avoiding a possible limit cycle.

Based on our numerical results, the addition of the convergence factor η seems unnecessary when the the initial value is well within the RoC, as required by Eq. A.6. The main question is when (and why) the limit-cycles are created with Newton's method. This question is at least partial explored in the example of Fig. 2. As long as the RoC condition is maintained, each step will progress closer to a root, and in the limit, as $n \to \infty$,

$$\frac{P_N(s_n)}{P'_N(s_n)} \to 0, \tag{III.1}$$

since $s_n \to s_r$ as $n \to \infty$.

We don't understand many observations in science (math and physics). But with some basic analysis, they are eventually explained. Einstein's 1905 analysis is the best known example. It is the *reductionist* method in science, and explains the success of the scientific method. This might be viewed as a form of evolution: success begets more success, while failure eventually dies off, perhaps slowly.

The process of systematically exploring these seemingly tiny discrepancy, almost always leads to new knowledge. Seeking out these idiosyncratic inconsistencies and trying to explain them is at the heart of the scientific method. When a problem is longstanding and considered fundamental, its resolution can even lead to a paradigm shift. Not surprisingly such deep insights are rarely welcomed by the scientific community, rather they are viewed with great skepticism. This can be good when if doesn't go on for 50 years.

The problem of finding roots using Newton's method is an excellent example. It is a case that can be explained only after careful thought and iterative analysis. I feel we are either close to that understanding, or it has been explained clearly enough that the debate can be stopped, and final conclusions may be reached. However, realize that there is no "final."

Limit cycles do exist in Newton's method, but in my view, they are due to under-sampling the complex plane. This is an example of aliasing, in the Nyquist sense, (Allen, 2020, p. 153,262). An under-sampled process becomes nonlinear when the "high frequencies" alias into the "base-band" frequencies. This nonlinear effect is easily removed by increasing the sampling rate above the Nyquist sampling frequency, defined as twice the highest frequency in the signal. While that concept is not clear in the context of Newton's method, it can explain limit-cycles, and slightly (2x-3x) increasing the computation, by decreasing the step-size η , the aliasing may be brought under control, and the problem becomes linear and well behaved. The onset of aliasing is easily detected. This leads to a well know method in signal processing called the *adaptive step-size*, which has been successfully applied in many engineering problems. It is, I believe, well understood and characterized in terms of aliasing (Allen and Sondhi, 1979; Rinzel and Miller, 1980, Sec. V, p. 126).

a) The linear prediction algorithm: An interesting alternative to stabilize NM is to use the linear prediction method, a causal recursion method invented in the 1940's (Vaidyanathan, 2007). It seems likely to me that the use of Linear Prediction (LP) could greatly improve the convergence properties of NM. The down side is that the LP method assume the step-size only has poles, which in our case is clearly not true. The zeros of $P_N(s)$ bias the estimate in a negative manner. However when the trajectory steps near a pole, the LP algorithm should fit the data extremely well, thus removing the influence of the pole. This approach could be especially effective if there were several poles in proximity.

APPENDIX

A. Derivation of Newton's method

Consider the monic polynomial $P_N(s)$, with $s, s_r, c_n \in \mathbb{C}$ and $n, k, N \in \mathbb{N}$:

$$P_N(s) = (s - s_r)^N + \sum_{k=1}^N c_{N-k} (s - s_r)^{N-k}, \quad (A.1)$$

where Taylor's formula is used to determine the coefficient vector $\boldsymbol{C} = [c_N, c_{N-1}, \cdots c_0]_{N \times 1}^T$

$$c_k = \left. \frac{1}{k!} \frac{d^k}{ds^k} P_N(s) \right|_{s=s_r}.$$
 (A.2)

Here $s = \sigma + j\omega$ is called the *Laplace frequency*, as defined by the Laplace transform (Allen, 2020). Depending on physical considerations, the coefficients c_k may be real or complex.

Assuming our initial guess for the root is s_0 is within the RoC (close to root s_r , we replace s_r with s_1 and swith s_0 , since $|(s_1 - s_0)^k| \ll |(s_r - s_0)|$ for $k \ge 2 \in \mathbb{N}$. Here we have assumed that within the RoC, the higher order terms may be ignored.

Iterating we increase n by 1. Thus $s_0 \rightarrow s_1$ and $s_1 \rightarrow s_2$, so the truncated Taylor series becomes

$$P_N(s_2) \approx (s_2 - s_1) \left. \frac{d}{ds} P_N(s) \right|_{s_2} + P_N(s_2).$$
 (A.3)

Generalizing this for $n \gg 1$ we find replace find $|(s_{n+1} - s_n)^k| \ll |(s_1 - s_0)|$ for $k \ge n \in \mathbb{N}$, (i.e., $\epsilon_n = s_1 - s_0$ is within its RoC), thus we may truncate Eq. A.1 to its linear term n = 1, resulting in the approximation Thus for large $n \to \infty$, $s_{n+1} \to s_r$, resulting in

$$\underline{P_N(s_{n+1})}^0 = (s_{n+1} - s_n)^N + \sum_{k=1}^N c'_k (s_{n+1})^{N-k}$$
(A.4)

Here c'_{n+1} is shorthand for $dP_N(s_{n+1})/ds$.

Solving for s_{n+1} gives Newton's method:

$$s_{n+1} = s_n - \frac{P_N(s_n)}{P'_N(s_n)}.$$
 (A.5)

Importantly, if s_n approaches a root of P'(s), the denominator can become arbitrarily large, resulting in a restart of the entire procedure.

On the other hand, if any guess of the root s_n is close to a root of (i.e., $P_N(s_r \pm \epsilon) \approx 0$) then for $n \ge 2 \in \mathbb{N}$, $\epsilon = s_n - s_r$ is within the RoC. Namely for all $k \in \mathbb{N} + 1$

$$(s_n - s_r)^k | \ll |(s_n - s_r)|.$$
 (A.6)

This complex analytic linearization step is the key to Newton's method. It will only be true if the difference equation remains linear, which requires Eq. A.6.

In summary: Newton's method is a linear approximation that critically depends on the RoC condition (Eq. A.6).

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