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Stability Analysis of a More General Class of Systems with Delay-Dependent Coefficients

Chi Jin, Keqin Gu, IEEE Senior Member, Islam Boussaada, and Silviu-Iulian Niculescu, IEEE Fellow

Abstract—This paper presents a systematic method to analyse the stability of systems with single delay in which the coefficient polynomials of the characteristic equation depend on the delay. Such systems often arise in, for example, life science and engineering systems. A method to analyze such systems was presented by Beretta and Kuang in a 2002 paper, but with some very restrictive assumptions. This work extends their results to the general case with the exception of some degenerate cases. It is found that a much richer behavior is possible when the restrictive assumptions are removed. The interval of interest for the delay is partitioned into subintervals so that the magnitude condition generates a fixed number of frequencies as functions of the delay within each subinterval. The crossing conditions are expressed in a general form, and a simplified derivation for the first-order derivative criterion is obtained. Illustrative examples are also presented.

Index Terms—Delay Systems, Stability Analysis

I. INTRODUCTION

The presence of time-delay has been widely observed in physical and engineering systems, and it is often caused by the finite time needed to transfer materials, energy and information. Such systems may be modeled as delay differential equations, and have attracted significant attentions of scholars from mathematics, engineering, life science and economics for many years. See [3], [9], [10], [11] for some recent progress. For a linear system with constant coefficients and single delay or multiple commensurate delays, a number of effective methods have been proposed [1], [2], [5]. The methods are along the line of D-subdivision [6], [7], also known as the τ-decomposition method [17] as the parameter involved in this case is the delay τ. These methods roughly proceed as follows: starting with one value of delay τ0 that one knows the number of characteristic roots on the right-half plane (usually τ0 = 0), one sweeps through an interested interval (τ0, τf) of delays, and identify all delays τk, k = 1, 2, ..., N − 1 for which there are characteristic roots on the imaginary axis. By identifying the direction these roots cross the imaginary axis, the change of the number of right-half plane roots as τ goes through τk can be determined. Thus, the interval (τ0, τf) is divided into subintervals (τk−1, τk), and the number of right-half plane roots within each subinterval is constant and can be explicitly determined. Especially, the subintervals of delay for the systems to be stable can be identified.

There are, however, practical systems in, for example, life science and engineering, for which the coefficients of the system characteristic equation depend on the delay values. For example, in [15], the source and dissipative process of a stellar dynamo is described by the following equations

\[
\begin{align*}
B_\phi(t) &= c_1 e^{-c_2 T_0} A(t - T_0) - c_2 B_\phi(t), \\
A(t) &= c_3 e^{-c_2 T_1} B_\phi(t - T_1) - c_2 A(t),
\end{align*}
\]

where \( B_\phi \) is the strength of toroidal field, and \( A \) is the strength of poloidal field, and \( c_1, c_2, c_3, T_0, T_1 \) are positive constants. The characteristic equation of the above system can be easily obtained as the following with delay-dependent coefficients:

\[
\lambda^2 + 2c_2 \lambda + c_2^2 - c_1 c_3 e^{-c_2 \tau} e^{-\tau \lambda} = 0,
\]

where \( \tau = T_0 + T_1 \).

A model of hematopoietic stem cell dynamics is given in [16]. The model is nonlinear, and possesses two equilibria. The linearized equation in the neighborhood of the nonzero equilibrium has the following characteristic equation

\[
\lambda + A(\tau) - B(\tau) e^{-\lambda \tau} = 0,
\]

where \( A, B \) are nonlinear functions of \( \tau \). Therefore delay-dependent coefficients may result from the linearized dynamics of a nonlinear time-delay system.

Time-delay systems with delay-dependent coefficients can also arise from the analysis of partial differential equations. As an example, the modeling of cell density in a generic compartment in [27] suggests an advection or reaction-convection equation of the following form:

\[
\frac{\partial x(t, a)}{\partial t} + V \frac{\partial x(t, a)}{\partial a} = -\gamma(t, a) x(t, a).
\]

A time-delay system can be obtained from the above equation using the method of characteristics

\[
S(t) = 2\beta(S(t - \tau_0)) e^{-\gamma \tau} S(t - \tau_0) - [\beta(S(t)) + \delta] S(t).
\]

Detailed derivation and the meaning of the variables and functions in the above equation can be found in [27]. It is clear that the delay parameter \( \tau \) enters the system coefficients through the exponential term \( e^{-\gamma \tau} \).

Other examples of systems with delay-dependent coefficients include the sun flower model [26], control systems using a finite-difference scheme for stabilization [21] as well as various population dynamics models [14]. Indeed, it has been pointed out in [12] that the dynamics of a population that goes
through distinct life stages in general involves delay-dependent parameters.

While it is possible to use the existing methods mentioned above to determine the stability of such a system with a given delay value, they are no longer sufficient to determine the range of delays for the system to be stable. Beretta and Kuang [12] presented an effective method to carry out such a stability analysis for systems with a single delay. However, the authors of [12] made some very restrictive assumptions, and the main attention has been paid to the crossing direction. No procedure was given in [12] to identify all the pairs \((j\omega, \tau)\) that satisfy the characteristic equation. In general, the structure of the functions \(\omega(\tau)\) implicitly defined by \(F(\omega, \tau)\) has not been sufficiently described in [12] to systematically identify all such pairs.

The purpose of this paper is to extend the method to the general case with the exception of some degenerate cases. As we will see, the removal of such restrictive assumptions means that a much richer behavior is possible. More specifically, the interval of interest for delay needs to be divided into subintervals so that the number of continuous functions \(\omega(\tau)\) remains constant within each subinterval. The number of such functions may change as the delay moves from one subinterval to another. The dividing points of the interval are those delays for which two polynomial equations have a common real solution. Based on such a structure, the crossing delays and the corresponding crossing frequencies may be identified systematically. Furthermore, the delay intervals such that the system is stable may be determined based on the crossing directions of each critical delay-frequency pair.

The crossing direction in the general case may be determined numerically. With additional nondegeneracy assumption, the crossing direction may be conveniently determined analytically similar to the method given in [12], although we will show that a simplified derivation is possible.

A preliminary version of this paper was presented in [20].

The following notation will be used in this paper. For a polynomial, \(\text{ord}(\cdot)\) denotes its order. For any complex number \(c\), \(\Re(c)\), \(\Im(c)\) and \(\tau\) denote its real part, imaginary part and conjugate, respectively. \(\mathbb{R}\) stands for the set of real numbers and \(\mathbb{R}_+\) for non-negative reals. We will use \(\partial\) with a subscript to denote partial derivatives. For instance, \(\partial_\tau D(\lambda, \tau) := \frac{\partial D(\lambda, \tau)}{\partial \tau}\).

II. PROBLEM STATEMENT

Consider a time-delay system with characteristic equation of the form:

\[ D(\lambda, \tau) = P(\lambda, \tau) + Q(\lambda, \tau)e^{-\tau \lambda} = 0, \quad (2) \]

where \(P(\lambda, \tau)\) and \(Q(\lambda, \tau)\) are continuous in \(\tau\) and are polynomials of \(\lambda\) with real coefficients for each given \(\tau \in \mathcal{I}\), and \(\mathcal{I} = [\tau^*, \tau^*']\) is the range of delay parameters \(\tau\) of interest, \(0 \leq \tau < \tau^*\). In some context, we may write \(P_t(\lambda)\) and \(Q_t(\lambda)\) instead of \(P(\lambda, \tau)\) and \(Q(\lambda, \tau)\) in order to emphasize them as functions (polynomial in this case) of \(\lambda\) for a given \(\tau\).

The same convention is also used for other functions of two independent variables with \(\tau\) as one of them. For example, we may write \(D_\tau(\lambda)\) instead of \(D(\lambda, \tau)\) to emphasize that we are considering \(D\) as a function of \(\lambda\) for a given \(\tau\) even though it is no longer a polynomial.

As we will see later on, the solutions of (2) with \(\lambda\) on the imaginary axis plays an important role in stability analysis, in which case, (2) becomes

\[ D(j\omega, \tau) = 0, \quad (3) \]

where \(\omega\) is real. For this purpose, we define:

\[ F(\omega, \tau) = P(j\omega, \tau)P(-j\omega, \tau) - Q(j\omega, \tau)Q(-j\omega, \tau). \quad (4) \]

It is not difficult to see that a necessary but not sufficient condition for \((\omega, \tau)\) to satisfy (3) is

\[ F(\omega, \tau) = 0. \quad (5) \]

The equation (5) is known as the magnitude condition, which means that the norms of the two complex number \(P(j\omega, \tau)\) and \(Q(j\omega, \tau)\) are equal.

We will restrict ourselves to systems that satisfy the following four assumptions:

Assumption I. For all \(\tau \in \mathcal{I}\), \(P_\tau\) satisfies

\[ \text{ord}(P_\tau) = n. \quad (6) \]

Furthermore,

\[ \lim_{\omega \to 0} \left| \frac{Q_\tau(j\omega)}{P_\tau(j\omega)} \right| < 1. \quad (7) \]

Assumption II. No \((\omega, \tau) \in \mathbb{R}_+ \times \mathcal{I}\) satisfies

\[ P(j\omega, \tau) = 0, \]
\[ Q(j\omega, \tau) = 0, \]
simultaneously.

Assumption III. Any \((\omega^*, \tau^*) \in \mathbb{R}_+ \times \mathcal{I}\) that satisfies (3) must also satisfy

\[ \partial_\lambda D(\lambda, \tau) \big|_{\lambda=\omega^*} \neq 0. \quad (8) \]

Furthermore, let \(\lambda(\tau)\) be the function implicitly defined by (2) in a sufficiently small neighborhood of \((j\omega^*, \tau^*)\) within \(\mathbb{R}_+ \times \mathcal{I}\), then for all \(\tau \neq \tau^*, \tau \in \mathcal{I}, |\tau - \tau^*| \) sufficiently small, \(\Re(\lambda(\tau)) \neq 0\).

Assumption IV. There are only a finite number of \((j\omega, \tau)\) in \(\mathbb{R}_+ \times \mathcal{I}\) that simultaneously satisfy (5) and

\[ \partial_\omega F(\omega, \tau) = 0. \quad (9) \]

These four assumptions are less restrictive than typical in the literature either stated explicitly or implicitly. Assumption I above requires the leading coefficient of \(P_\tau\) not to vanish for all \(\tau \in \mathcal{I}\), and

\[ \text{ord}(Q_\tau) \leq n. \quad (10) \]

For time-delay systems of retarded type, (10) is satisfied with strict inequality. When (10) is an equality, the time-delay
system is of neutral type, and (7) requires the absolute value of the leading coefficient of $Q_\tau(\lambda)$ to be strictly less than that of $P_\tau(\lambda)$. Systems of neutral type involve some surprising subtleties. See [2] for an example for systems with single delay. For more comprehensive coverage see [4] and [9].

Assumption II is much less restrictive than the counterpart in [12] which is

$$P(j\omega, \tau) + Q(j\omega, \tau) \neq 0 \text{ for all } (\omega, \tau) \in \mathbb{R}^2. \quad (11)$$

Indeed, the two complex equations in Assumption II are equivalent to four real equations with two real “unknowns” $\omega$ and $\tau$. Obviously, cases that violate this assumption are degenerate and rare. On the other hand, the set

$$\{P(j\omega, \tau) + Q(j\omega, \tau) | (\omega, \tau) \in \mathbb{R}^2\}$$

is a region in the complex plane, and (11) requires this region not to include the origin, which is obviously much more restrictive. As will be presented later, the analysis is based on the phase condition on the set of parameters that satisfy the magnitude condition (5). The violation of this assumption makes the phase condition discontinuous at this point, and requires separate treatment which will not be pursued here.

In Assumption III, Condition (8) guarantees that $\lambda(\tau^*) = j\omega^*$ is a simple characteristic root, and consequently $\lambda(\tau)$ is well defined in a small neighborhood of $\tau = \tau^*$ by the implicit function theorem [22]. and $\lambda'(\tau)$ exists at $\tau^*$ if $D(\lambda, \tau)$ is differentiable with respect to $\tau$ at $(\lambda(\tau^*), \tau^*)$. The remaining part of the assumption means that the curve $\lambda(\tau)$ is on the imaginary axis only at one point $\lambda^* = \lambda(\tau^*)$ in this neighborhood. A more restrictive assumption is to assume $\Re(\lambda'(\tau)) \neq 0$, which is implicitly assumed in most works of this nature, including [12].

Assumption IV is also rather natural. It requires two real equations in two real variables to admit a finite number of solutions in the set $\mathbb{R}_+ \times \mathcal{J}$. This assumption holds for most systems with delay dependent coefficients in practice. This assumption allows the delay interval $\mathcal{J}$ to be divided into a finite number of sub-intervals such that the polynomial $F_\tau(\omega)$ has a constant number of simple positive roots within each subinterval.

In most cases, we may choose the lower limit $\tau'$ of $\mathcal{J}$ to be 0, and the upper limit $\tau''$ sufficiently large. We leave them in this general form so that the method we present here is still valid even if some of the assumptions are violated for some $\tau < \tau'$ or $\tau > \tau''$.

This paper provides extension of the analysis in [12] so that it is still applicable when the condition (11) is violated. In [12], it is also implicitly assumed that the number of real roots, $\pm \omega_k(\tau), k = 1, 2, \ldots, m,$ of $F_\tau(\omega)$ remains constant within the delay interval of interest $\mathcal{J}$, and they are continuously differentiable. With our relaxed assumptions, these are no longer true. Especially, the real roots may suddenly emerge or disappear as the delay $\tau$ increases within $\mathcal{J}$. It is therefore essential to understand the structure of this solution set in order to solve the stability problem. This will be discussed in the next section.

### III. Stability Analysis

The main idea for stability analysis here is along the line of $\tau$-decomposition method outlined in the introduction. The validity of the method is based on the fact that there exists a constant $c > 0$ for any closed interval of $\tau$ such that all roots of $D(\lambda, \tau)$ with $\Re(\lambda) > -c$ vary continuously as $\tau$ changes. This is true under Assumption I [2][4][9].

The critical aspects of the stability analysis are: (i) identifying the values of $\tau$ such that there is at least one root of $D_\tau(\lambda)$ on the imaginary axis, as well as the corresponding imaginary roots, and (ii) determining whether these imaginary roots move from the left-half plane to the right-half plane, or vice versa, or return to the original side as $\tau$ increases through these values. In this section, we will consider the first aspect, and describe the process of stability analysis assuming we know the answer to the second aspect. In the next section, we will describe some methods of accomplishing the second aspect.

To accomplish the first aspect stated in the last paragraph, it is useful to introduce the notation

$$S(\lambda, \tau) = \frac{P(\lambda, \tau)}{Q(\lambda, \tau)} e^{j\lambda}, \quad (12)$$

whenever

$$Q(\lambda, \tau) \neq 0. \quad (13)$$

Then

$$S(j\omega, \tau) = W(\omega, \tau)e^{j\theta(\omega, \tau)}, \quad (14)$$

where

$$W(\omega, \tau) = \frac{|P(j\omega, \tau)|}{|Q(j\omega, \tau)|}, \quad (15)$$

$$\theta(\omega, \tau) = \angle P(j\omega, \tau) - \angle Q(j\omega, \tau) + \omega \tau + \pi. \quad (16)$$

When $\lambda = j\omega$ is on the imaginary axis, we note that (3) is equivalent to the following two conditions

$$W(\omega, \tau) = 1, \quad (17)$$

$$\theta(\omega, \tau) = 2r\pi, \text{ for some integer } r, \quad (18)$$

provided that (13) holds.

Equation (17) is equivalent to (5), and represents the magnitude condition. Equation (18) is the phase condition.

To capture essentially the same phase relationship, in [12] a function different from $\theta(\omega, \tau)$ is introduced, which requires the more restrictive condition (11). Let

$$\mathcal{W} = \{(\tau, \omega) | \tau \in \mathcal{J}, \omega \in \mathbb{R}, F(\omega, \tau) = 0\}, \quad (19)$$

then $(\tau, \omega) \in \mathcal{W}$ if and only if $(\tau, \omega)$ satisfies (13) and (17) in view of Assumption II. Therefore, an effective approach to determine all $(\tau, \omega)$ satisfying (3) is to first determine the set $\mathcal{W}$, and then choose from $\mathcal{W}$ those $(\tau, \omega)$ that also satisfy (18).

To understand the structure of $\mathcal{W}$, we will examine the function $F(\omega, \tau) = F_\tau(\omega)$ more closely. For any given $\tau$, $F_\tau(\omega)$ is an $2n^{th}$ order polynomial with real coefficients in
view of Assumption I, and it is an even function. It can also be written as an $n^{th}$ order polynomial of $\omega^2$,

\[ F(a, \tau) = F(\omega, \tau), \]

\[ \alpha = \omega^2. \]

Therefore

\[ F(a, \tau) = 0 \]

will provide $n$ solutions $\omega_k$, $k = 1, 2, \ldots, n$. Without loss of generality, let $\omega_k$, $k = 1, 2, \ldots, n_p$, $n_p \leq n$, be the only real and non-negative solutions. Then, all the real solutions of (5) are $\pm \omega_k$, $k = 1, 2, \ldots, n_p$, where $\omega_k = \sqrt{\alpha_k}$. In general, the number of non-negative real roots $n_p$ depends on $\tau$. In order to understand this dependence, let $\tau^{(i)}$, $i = 1, 2, \ldots, K - 1$ be the set of all $\tau \in S$ such that $(\omega, \tau)$ simultaneously satisfies (5) and (9) for some $\omega \in \mathbb{R}_+$. (recall this set is indeed finite according to Assumption IV). We agree to order $\tau^{(i)}$ in ascending order

\[ \tau^{(1)} < \tau^{(2)} < \ldots < \tau^{(K-1)}. \]

We will also write $\tau^{(0)} = \tau^l$ and $\tau^{(K)} = \tau^u$. Then, we may partition $S$ into $K$ subintervals

\[ S^{(i)} = [\tau^{(i-1)}, \tau^{(i)}], \quad i = 1, 2, \ldots, K. \]

The interior of $S^{(i)}$ is denoted as $S^{(i)}_o = (\tau^{(i-1)}, \tau^{(i)})$. Then the structure of the set $W$ may be very clearly described in the following proposition.

**Proposition 1.** For a given $i$, the number of real roots of $F_\tau(\omega)$ are the same for all $\tau \in S^{(i)}_o$, and they are all simple. These real simple roots are continuous functions of $\tau$, and may be expressed as $\pm \omega_k^{(i)}(\tau)$, $k = 1, 2, \ldots, m^{(i)}$, where $m^{(i)} \leq n$, and $\omega_k^{(i)}(\tau) > 0$ for all $\tau \in S^{(i)}_o$.

**Proof.** For a fixed $i$, by definition, for all $\tau \in S^{(i)}_o$, any $\omega \in \mathbb{R}_+$ that satisfies

\[ F_\tau(\omega) = 0 \]

must satisfy

\[ F_\tau'(\omega) = \partial_\omega F(\omega, \tau) \neq 0, \]

from which we conclude that all real roots of $F_\tau(\omega)$ are simple. As $F_\tau(\omega)$ is an even function of $\omega$, we can also conclude that $-\omega$ is also a root if $\omega$ is a real root, and $\omega = 0$ is not a root (otherwise, it cannot be simple). To show the invariance of the number of real solutions within $S^{(i)}_o$, let $\tau^* \in S^{(i)}_o$, and let $\omega_k^{*}$, $k = 1, 2, \ldots, m$ be the only real roots of $F_\tau(\omega)$. By the continuity of roots with respect to coefficients[8], we may define $m$ continuous functions $\omega_k(\tau)$, $k = 1, 2, \ldots, m$ in $S^{(i)}_o$, $\omega_k(\tau^*) = \omega_k^{*}$, and each $\omega_k(\tau)$ is a root of $F_\tau(\omega)$. The proof is complete if we show that all $\omega_k(\tau)$ are real in $S^{(i)}_o$ as this also implies that $\omega_k(\tau)$ are simple roots of $F_\tau(\omega)$.

For a given $k$, let

\[ \tau_M = \sup\{\tau_l | \omega_k(\tau) \in \mathbb{R} \text{ for all } \tau \in [\tau^*, \tau_0]\}. \]

By continuity, $\omega_k(\tau_M)$ is real. We will show $\tau_M = \tau^{(i)}$. If not, for arbitrarily small $\epsilon > 0$, $\omega_k(\tau_M + \epsilon)$ is not real, which can be made arbitrarily close to $\omega_k(\tau_M)$ with sufficiently small $\epsilon$. But this means that its complex conjugate $\omega_k(\tau_M + \epsilon)$ is also a root of the polynomial with real coefficients $F_{\tau_M + \epsilon}(\omega)$ and arbitrarily close to $\omega_k(\tau_M)$. The continuity of roots with respect to the coefficients means that $\omega_k(\tau_M)$ cannot be a simple root of $F_{\tau_M}(\omega)$, which contradicts the first part of this proposition that we have already proven. We have thus shown that $\omega_k(\tau)$ is real for all $\tau \in \tau^{(i-1)}, \tau^{(i)}$. Similarly, we can show that $\omega_k(\tau)$ is real for all $\tau \in (\tau^{(i-1)}, \tau^l)$, and the proof is complete. □

As $\tau$ moves rightward from a point in $S^{(i)}_o$, some, say $m$, real roots, and $2l$ complex roots of $F_\tau(\omega)$ may merge to form a multiple root as $\tau$ reaches $\tau^{(i)}$, and some, say $2k$, become complex while $m + 2l - 2k$ roots remain real as $\tau$ enters $S^{(i+1)}_o$. The most common scenarios are either two real roots merge and become complex, or two complex roots merge and become real as $\tau$ moves from $S^{(i)}_o$ to $S^{(i+1)}_o$ through $\tau^{(i)}$.

A real root of $F_\tau(\omega)$ in $S^{(i)}_o$, say $\omega_k^{(i)}(\tau)$, that does not merge with other roots at $\tau^{(i)}$ remains real, and becomes $\omega_l^{(i+1)}$ for some $l \leq m^{(i+1)}$ as $\tau$ moves from $S^{(i)}_o$ to $S^{(i+1)}_o$ through $\tau^{(i)}$.

For a given $i$ and $k$, as $\omega_k^{(i)}$ depends on $\tau$ continuously in $S^{(i)}_o$, we will require $\angle P(j\omega_k^{(i)}(\tau), \tau)$ and $\angle Q(j\omega_k^{(i)}(\tau), \tau)$ to be continuous functions of $\tau$. This means that

\[ \theta_k^{(i)}(\tau) = \theta(\omega_k^{(i)}(\tau), \tau), \quad k = 1, 2, \ldots, m^{(i)} \]

are continuous functions of $\tau$ within $S^{(i)}_o$, and will be known as the phase functions. On the other hand, this continuity requirement means that the values of $\angle P(j\omega_k^{(i)}(\tau), \tau)$, $\angle Q(j\omega_k^{(i)}(\tau), \tau)$ and $\theta_k^{(i)}(\tau)$ may not be restricted to any $2\pi$ range. Furthermore, if $\omega_k^{(i)}(\tau)$ and $\omega_l^{(i)}(\tau)$ merge at, say, $\tau^{(i)}$, and we extend the definition of $\theta_k^{(i)}(\tau)$ and $\theta_l^{(i)}(\tau)$ to $\tau^{(i)}$ by continuity, then it is possible that

\[ \theta_k^{(i)}(\tau) - \theta_l^{(i)}(\tau) = 2\pi r, \]

for some integer $r \neq 0$ even though

\[ \omega_k^{(i)}(\tau^{(i)}) = \omega_l^{(i)}(\tau^{(i)}). \]

Going through each interval $S^{(i)}$ and each curve $\omega_k^{(i)}(\tau)$, we may identify all $\tau = \tau_i$ such that

\[ \theta_k^{(i)}(\tau_i) = 2\pi r, \quad r \text{ integer,} \]

for some $k$ if $\tau_i \in S^{(i)}$. Notice, the ends of the intervals, $\tau^{(i)}$, $i = 0, 1, \ldots, K$ should also be included. We will order such $\tau_i$ in an ascending order

\[ \tau^l \leq \tau_1 < \tau_2 < \ldots < \tau_k \leq \tau^u. \]

Each $\tau_i$ is known as a critical delay. For each given $\tau_i$, it is possible that more than one $k$ satisfies (28), and we denote the corresponding $\omega_k^{(i)}(\tau_i) \geq 0 = \omega_h, h = 1, 2, \ldots, H_i$. Therefore, we can identify all the pairs $(\omega_h, \tau_i)$, $h = 1, 2, \ldots, H_i; l = 1, 2, \ldots, L$, that satisfy (3). It should also be pointed out that it is possible that a simple root of $D_{\tau}(j\omega)$ may be a double root of $F_\tau(\omega)$. In other words, for some $\tau = \tau^{(i)}$, an $\omega$ that simultaneously satisfy (5) and (9) may satisfy (18) without violating Assumption III. Such
points pose special difficulty in determining crossing direction as will be shown in the next section.

Now we will describe the representation of the second aspect we mentioned at the beginning of this section, i.e., the movement of the imaginary roots. For a given pair \((\omega_h, \tau)\) that satisfies (3), a sufficiently small \(\varepsilon > 0\), and any \(\tau \in (\tau_i, \tau_i + \varepsilon)\), there is a unique solution \(\lambda^{\pm}_{lh}\) of (2) in the neighborhood of \(j\omega_h\). Assumption III and continuity means that \(\Re(\lambda^{\pm}_{lh})\) must be nonzero, and have the same sign for any \(\tau \in (\tau_i, \tau_i + \varepsilon)\). Similarly, let \(\lambda^{\pm}_{lh}\) be the unique solution of (2) in the neighborhood of \(j\omega_h\) corresponding to a given \(\tau \in (\tau_i - \varepsilon, \tau_i)\), then \(\Re(\lambda^{\pm}_{lh})\) must have the same sign for all such \(\tau\). We define

\[
\text{Inc}(\omega_h, \tau_i) = \frac{\text{sgn}(\Re(\lambda^{+}_{lh})) - \text{sgn}(\Re(\lambda^{-}_{lh}))}{2}.
\]

If \(\text{Inc}(\omega_h, \tau_i) = 1\), a root of \(D_{\tau}(\lambda)\) moves from the left-half plane to the right-half plane crossing the imaginary axis at \(j\omega_h\) as \(\tau\) increases from \(\tau_i - \varepsilon\) to \(\tau_i + \varepsilon\). On the other hand, if \(\text{Inc}(\omega_h, \tau_i) = -1\), then the root moves from the right-half plane to the left-half plane as \(\tau\) increases from \(\tau_i - \varepsilon\) to \(\tau_i + \varepsilon\). If \(\text{Inc}(\omega_h, \tau_i) = 0\), the root moves towards the imaginary axis, touching it at \(j\omega_h\), then return to the same half plane without crossing the imaginary axis. We also define

\[
\text{Inc}(\tau_i) = 2 \sum_{h=1}^{H_1} \text{Inc}(\omega_h, \tau_i).
\]

Then, as \(\tau\) increases from \(\tau_i - \varepsilon\) to \(\tau_i + \varepsilon\), there is a net increase of \(\text{Inc}(\tau_i)\) roots on the right-half plane. Notice, \(\omega_h > 0\), \(h = 1, 2, \ldots, H_1\) only accounts for the roots on the upper half of the imaginary axis, and the coefficient 2 in front of the summation sign in (30) accounts for the fact that the roots of \(D_{\tau}(\lambda)\) are symmetric to the real axis.

Let the number of right-half plane roots of \(D_{\tau}(\lambda)\) be \(N^{\mu}(\tau)\). Then, for any \(\tau \in \mathcal{S}, \tau \neq \tau_i, i = 1, 2, \ldots, L\), we have

\[
N^{\mu}(\tau) = N^{\mu}(\tau_i) + \sum_{l=1}^{L_2} \text{Inc}(\tau_i),
\]

where \(L_2 = \max(l \mid \tau_l < \tau)\).

If \(\tau = 0\), as \(D_{\tau}(\lambda)\) is a polynomial, \(N^{\mu}(\tau)\) is easily obtained. If \(\tau > 0\), \(N^{\mu}(\tau)\) may be obtained by a method covered in [5] or [1] (but notice the correction [2]). If there are imaginary roots for \(D_{\tau}(\lambda), N^{\mu}(\tau)\) should not count these imaginary roots, and \(\text{Inc}(\omega_h, \tau)\) should be defined as,

\[
\text{Inc}(\omega_h, \tau) = \begin{cases} 1, & \text{if } \text{sgn}(\Re(\lambda^{+}_{lh})) = 1, \\ 0, & \text{otherwise} \end{cases}
\]

(32)

instead. Obviously, \(N^{\mu}(\tau)\) remains the same in the interval \((\tau_i, \tau_i + 1)\) for any given \(l\). The system is stable if \(N^{\mu}(\tau) = 0\).

IV. CROSSING DIRECTION CONDITIONS

In the last section, a procedure of determining the range of \(\tau\) in \(\mathcal{S}\) such that \(D_{\tau}(\lambda)\) is stable has been developed, provided a method of determining \(\text{Inc}(\omega_h, \tau_i)\) is available. It is not difficult to determine \(\text{Inc}(\omega_h, \tau_i)\) according to the definition if a numerical method is used. Indeed, as the solution \((j\omega_h, \tau_i)\) is already known for \(D(\lambda, \tau)\), the Newton-Raphson method may be used to find the unique solution in the neighborhood of \(j\omega_h\) when \(\tau\) is very close to \(\tau_i\) and \(D(\lambda, \tau)\) is differentiable with respect to \(\tau\) in a neighborhood of \((\omega_h, \tau_i)\) [8]. In many cases, however, a simple analytical method can be used, which will be described as follows.

The simplest case is when

\[
\Re(\lambda^{+}_{lh}(\tau))_{\tau = \tau_i} \neq 0,
\]

where, \(\lambda^{+}_{lh}(\tau)\) is the implicit function defined by (2) in the neighborhood of \((j\omega_h, \tau_i)\) provided that \(\lambda^{+}_{lh}(\tau)\) is differentiable at \(\tau_i\). This can be guaranteed by requiring \(D(\lambda, \tau)\) to be differentiable w.r.t \(\tau\) at \((j\omega_h, \tau_i)\) [8]. Indeed, provided that (33) is satisfied, it is easy to see

\[
\text{Inc}(\omega_h, \tau_i) = \text{sgn}(\Re(\lambda^{+}_{lh}(\tau_i))),
\]

if \(\tau_i > \tau_i^l\). On the other hand, if \(\tau_i = \tau_i^l\), we have

\[
\text{Inc}(\omega_h, \tau_i) = \max\left\{0, \text{sgn}(\Re(\lambda^{+}_{lh}(\tau_i)))\right\}.
\]

(34)

If (33) is violated, and \(D(\lambda, \tau)\) is differentiable to a sufficiently high order at \((j\omega_h, \tau_i)\), then it follows from equation (8) in Assumption III and the implicit function theorem that the derivatives of \(\lambda(\tau)\) exist up to a sufficiently high order at the point \((j\omega_h, \tau_i)\) [8]. Consequently we may express \(\text{Inc}(\omega_h, \tau_i)\) using higher order derivatives. Suppose

\[
\Re\left(\frac{d^k\lambda(\tau)}{d\tau^k}\right)_{\tau = \tau_i} = 0, \quad k = 1, 2, \ldots, m - 1,
\]

\[
\Re\left(\frac{d^m\lambda(\tau)}{d\tau^m}\right)_{\tau = \tau_i} \neq 0.
\]

Then, if \(\tau_i > \tau_i^l\), then

\[
\text{Inc}(\omega_h, \tau_i) = \begin{cases} \text{sgn}(\Re\left(\frac{d^m\lambda(\tau)}{d\tau^m}\right)), & \text{if } m \text{ is odd}, \\ 0, & \text{if } m \text{ is even}. \end{cases}
\]

(36)

If \(\tau_i = \tau_i^l\), on the other hand, then

\[
\text{Inc}(\omega_h, \tau_i) = \max\left\{0, \text{sgn}(\Re\left(\frac{d^m\lambda(\tau)}{d\tau^m}\right))\right\}.
\]

(37)

If the condition (8) in Assumption III is violated for some imaginary characteristic root \(\lambda = j\omega^l\), we are then faced with a characteristic root with multiplicity and cannot regard it as a locally differentiable function of \(\tau\). In this case, the trajectory of characteristic roots parameterized by \(\tau\) may have several branches passing through the point \(j\omega^l\) on the imaginary axis. One may still determine the increment in the number of unstable roots based on these branches of curves, which can be locally characterized by the Newton-Puiseux series. Comprehensive analysis of this problem can be found in [23], [24] and [25]. An eigenvalue perturbation approach is taken in [23] and [24], which applies also to systems represented by the state-space matrices, whilst the analysis in [25] is based on the characteristic equations.

We will now give an explicit expression of \(\text{sgn}(\Re(\lambda^{+}_{lh}(\tau_i)))\) and leave the high-order analysis to future work. The expression is similar to that given in [12], but our derivation here is more succinct. For this purpose, we henceforth replace Assumption III by the following one:
Lemma 1. Any pair \((ω^*, τ^*)\) that satisfies Assumption IIIa must also satisfy (8).

Proof. At \((ω^*, τ^*)\)

\[
F = \bar{P}P - \bar{Q}Q = 0,
\]

\[
e^{-τλ} = -\bar{P}/P,
\]

\[
\frac{∂τ}{∂j}D = \frac{∂τ}{∂j}P + (\frac{∂j}{∂Q} - \frac{∂Q}{∂j})e^{-τλ} = 0.
\]

Therefore,

\[
\frac{∂ω}{∂ω}F = 2\Re\left(\frac{j}{ω}\frac{∂ω}{∂j}P - j\frac{∂j}{∂Q}Q\right) = -2\Re(\frac{∂ω}{∂Q}P - \frac{∂Q}{∂ω}P) = -2\Re\left(\frac{∂ω}{∂Q}P + \frac{∂j}{∂Q}Q + \frac{∂j}{∂Q}Pω\right) = -2\Re\left(\frac{∂ω}{∂Q}P + \frac{∂j}{∂Q}(Qe^{-τλ}) - τQe^{-τλ}\right) = -2\Re(\frac{∂ω}{∂Q}D).
\]

The above indicates that \(\frac{∂ω}{∂ω}F(ω^*, τ^*) \neq 0\) implies (8). \(\square\)

It should be pointed out that the converse is not necessarily true. Indeed, the proof above shows that \(\frac{∂ω}{∂ω}F(ω^*, τ^*) = 0\) only implies that \(\frac{∂ω}{∂Q}D(ω^*, τ^*)\) is parallel to \(P(jω^*, τ^*)\), which does not necessarily mean \(\frac{∂ω}{∂ω}D(ω^*, τ^*) = 0\).

Proposition 2. Let \((ω^*, τ^*) \in \Re × \mathcal{J}\) satisfy (3) and Assumption IIIa. Then (2) defines \(λ\) as a differentiable function of \(τ\) in a sufficiently small neighborhood of \((jω^*, τ^*)\), and

\[
\text{sgn}\left(\Re\left(\frac{dλ}{dτ}\right)\right) = \text{sgn}\left(\frac{∂ω}{∂ω}F(ω, τ)\right) = \frac{∂ω}{∂Q}(ω, τ) = \frac{∂ω}{∂Q}F\left(\begin{array}{c}
ω = \omega^* \\
τ = τ^*
\end{array}\right) = \frac{∂ω}{∂Q}F\left(\begin{array}{c}
ω = \omega^* \\
τ = τ^*
\end{array}\right),
\]

where

\[
\frac{dτ}{dτ} = \frac{∂ω}{∂Q}F\left(\begin{array}{c}
ω = \omega^* \\
τ = τ^*
\end{array}\right)
\]

is the total derivative of \(θ(ω, τ)\) with respect to \(τ\) when \(ω\) is considered as a function of \(τ\) defined implicitly by (5) in a sufficiently small neighborhood of \((ω^*, τ^*)\), and \(\frac{dτ}{dτ}\) is the derivative of the function \(ω(τ)\) so defined.

Proof. Lemma 1 and Assumption IIIa indicate that \(\frac{∂ω}{∂Q}D(λ, τ) \neq 0\) and \(\frac{∂τ}{∂j}D(λ, τ)\) exists in a neighborhood of \((jω^*, τ^*)\). Therefore, the equation (2), or equivalently

\[
S(λ, τ) = 1,
\]

defines \(λ\) as a differentiable function of \(τ\) in a small neighborhood of \(τ^*\) in view of the implicit function theorem. A differentiation of (40) yields

\[
\frac{dλ}{dτ} + \frac{∂τ}{∂j}S = 0,
\]

from which

\[
\frac{dλ}{dτ} = -\frac{∂τ}{∂j}S = -\frac{∂τ}{∂j}\frac{S}{|∂τS|^2}.
\]

But, at \(λ = jω^*\),

\[
\frac{∂τ}{∂j}S(λ, τ) = \frac{1}{j}\frac{∂ω}{∂j}S(jω, τ) = \frac{1}{j}\left(\Re\left(\frac{1}{W} - j\frac{∂ω}{∂j}W\right)\right) = \frac{1}{j}\left[\Re\left(\frac{1}{W} - j\frac{∂ω}{∂j}W\right)\right] = \frac{1}{j}\left[\Re\left(\frac{1}{W} - j\frac{∂ω}{∂j}W\right)\right].
\]

In the last step, (40) and (14) have been used. Similarly, we may obtain

\[
\frac{∂τ}{∂j}S = \frac{1}{W} - j\frac{∂ω}{∂j}W + \frac{∂ω}{∂j}W.
\]

Therefore,

\[
\text{sgn}\left(\Re\left(\frac{dλ}{dτ}\right)\right) = \frac{1}{W} - j\frac{∂ω}{∂j}W + \frac{∂ω}{∂j}W.
\]

When \(ω\) is a function of \(τ\) defined implicitly by (5), or equivalently by (17), we have:

\[
\frac{dτ}{dτ} = \frac{∂ω}{∂j}W = -\frac{∂ω}{∂j}W = -\frac{∂ω}{∂j}F(ω, τ).
\]

In view of \(|Q(ω^*, τ^*)| = |P(ω^*, τ^*)|\), it is easy to show that

\[
\left.\frac{∂ω}{∂j}W\right|_{ω=ω^*} = \frac{1}{|P|}\frac{∂ω}{∂j}F\left(\begin{array}{c}
ω = \omega^* \\
τ = τ^*
\end{array}\right).
\]

A substitution of (41) by (42) and (43) yields

\[
\text{sgn}\left(\Re\left(\frac{dλ}{dτ}\right)\right) = \text{sgn}\left(\Re\left(\frac{1}{|P|}\frac{∂ω}{∂j}F\left(\begin{array}{c}
ω = \omega^* \\
τ = τ^*
\end{array}\right)\right)\right),
\]

from which (39) can be easily derived. \(\square\)

We now make a useful observation about the first factor in (39).

Proposition 3. For any given \(i\) and \(k\), the quantity

\[
\text{sgn}\left(\Re\left(\frac{dλ}{dτ}\right)\right)_{ω=ω_i(τ^*)}
\]

remains constant for all \(τ \in \mathcal{J}_i^{(k)}\).

Proof. Due to the continuity of \(∂ωF(ω, τ)\), in order for \(∂ωF(ω_i(τ), τ)\) to change sign, it must first vanish, which violates the definition of \(\mathcal{J}_i^{(k)}\). \(\square\)

The above proposition indicates that the first factor in (39) only needs to be checked once for each curve \(ω_i(τ)\) within
the interval $\mathscr{I}_{\phi}^{(i)}$. Next, we will provide an explicit expression for the second factor.

**Proposition 4.** If $(\omega, \tau)$ satisfies (3),

$$\frac{dF}{d\tau} = \frac{1}{|P|^2} \left( P \frac{dF}{d\tau} \right) - P \frac{dF}{d\tau} - Q \frac{dF}{d\tau} + Q \frac{dF}{d\tau},$$

where the subscripts $r$ and $i$ represent the real and imaginary part of the quantities, respectively, and the total derivatives may be calculated by

$$\frac{dF}{d\tau} = \frac{dF}{d\tau} + \partial_x\phi,$$

where $\phi$ may be $P$, $P_r$, $Q$, or $Q_i$, and

$$\frac{dF}{d\tau} = -\partial_xF / \partial_\omega F.$$

**Proof.** Consider the identity

$$S = We^{\omega \tau} = -\frac{Pe^{i\omega \tau}}{Q},$$

(45)

By taking total derivative with respect to $\tau$, with $\omega(\tau)$ implicitly defined by (5), and noticing

$$W(\omega(\tau), \tau) = 1$$

for all $\tau$,

we obtain

$$j \frac{dF}{d\tau} We^{-j\omega \tau} = -\frac{dF}{d\tau} \left( \frac{P}{Q} \right) e^{i\omega \tau} - j \left( \frac{dF}{d\tau} + \frac{dF}{d\tau} \right) + \partial_x\phi.$$

Solving the above for $\frac{dF}{d\tau}$ and using (45), we obtain

$$\frac{dF}{d\tau} = \frac{1}{j} \left( \frac{1}{P} \frac{dF}{d\tau} - \frac{1}{Q} \frac{dF}{d\tau} \right) + \frac{dF}{d\tau} + \omega. $$

(46)

In view of $|P|^2 = |Q|^2$, the expression in the parentheses in (46) can be written as

$$\frac{1}{P} \frac{dF}{d\tau} - \frac{1}{Q} \frac{dF}{d\tau} = \frac{P}{PP} \frac{dF}{d\tau} - \frac{Q}{QQ} \frac{dF}{d\tau} = \frac{P}{PP} \frac{dF}{d\tau} - \frac{Q}{QQ} \frac{dF}{d\tau}.$$

A substitution of (46) by the above completes the proof. $\square$

While no explicit expression was given for $\frac{dF}{d\tau}$ in [12], an explicit expression of $S_n(\tau)$ in [12] could be obtained by going through the proof of Theorem 2.2 in [12]. Proposition 3 above can be considered as the consequence of Theorem 2.2 and Remark 2.2 in [12]. Indeed, it can be seen that $S_n(\tau)$ in Theorem 2.2 in [12] is equal to $(\theta(\tau) - 2n\pi)/\omega(\tau)$ here. Remark 2.2 in [12] indicates that the factor $\omega(\tau)$ does not affect the sign of the derivative at the crossing point.

It is interesting to apply the conclusions of Proposition 2 to the case of delay-independent coefficient polynomials, i.e., when $P(\lambda, \tau)$ and $Q(\lambda, \tau)$ are independent of $\tau$. In this case, $F(\omega, \tau)$ is independent of $\tau$, the curves $\phi(\tau)$ become constants, and $dF/\tau = \omega$. As a result, the crossing direction given in (39) is independent of delay. This fact is well-known in the literature on single or commensurate delay systems with delay-independent coefficients, and have been stated either implicitly [5] or explicitly [19] as the invariance property.

More generally, for systems with delay-dependent coefficient polynomials discussed in this paper, we may still identify delay intervals where the crossing direction is invariant provided $P(\lambda, \tau)$ and $Q(\lambda, \tau)$ are continuously differentiable with respect to $\tau$. Indeed, for a given subinterval $\mathscr{I}^{(i)} = (\tau^{(i-1)}, \tau^{(i)})$, and frequency curve $\omega(\tau)$, we may identify all the delay values $\tau_{ki} = 1, 2, \ldots, L - 1$, $\tau^{(i-1)} < \tau_{ki} < \tau^{(i)}$, $\tau^{(i)}$, such that $(dF/\tau)_{\tau = \tau^{(i)}} = 0$. Let $\xi_{ki} = \tau^{(i-1)}, \xi_{ki} = \tau^{(i)}$. Then, we may conclude, by continuity, that the crossing direction at the curve $\omega(\tau)$ remains invariant for all $\tau \in \left(\xi_{ki-1}, \xi_{ki}\right), l = 1, 2, \ldots, L$. Note that the intervals for invariant crossing direction $(\xi_{ki-1}, \xi_{ki})$ are different for different frequency curves in general.

V. NUMERICAL EXAMPLES

In this section, we present three examples to illustrate the method developed in this paper.

**Example 1.** We first consider the stellar dynamo model mentioned in the introduction. The system characteristic equation is given in (1). Therefore,

$$P(\lambda, \tau) = \lambda^2 + 2c_2\lambda + c_2^2,$$

$$Q(\lambda, \tau) = -c_1c_3e^{-c_2\tau}.$$  

The parameters are set as: $c_1 = -10$, $c_2 = 2$, $c_3 = 3$. We are concerned with the stability of the system for $\tau \in \mathscr{I} = [0, 2]$. Since $\text{ord}(P_2) = 2$ and $\text{ord}(Q_2) = 0$, Assumption I holds. Assumption II requires the following two equations do not hold simultaneously for real $\omega$ and $\tau \in \mathscr{I}:

$$-\omega^2 + 2jc_2\omega + c_2^2 = 0,$$

$$-c_1c_3e^{-c_2\tau} = 0,$$

which is obviously true. The other assumptions can be verified as we carry out the remaining analysis. The function $F$ in this case is

$$F(\omega, \tau) = \omega^4 + 2c_2^2\omega^2 + c_2^4 - c_1c_3^2e^{-c_2\tau}.$$  

(47)

Only one pair of parameters $(\omega, \tau) = (0, \tau^{(1)})$ simultaneously satisfies (5) and (9), where

$$\tau^{(1)} = -\frac{1}{2c_2} \ln\left(\frac{c_1^2}{c_2^2c_3^2}\right) \approx 1.006.$$  

Therefore, Assumption IV is satisfied. The interval $\mathscr{I}$ is thus partitioned into two subintervals $\mathscr{I}^{(1)} = [\tau^{(0)}, \tau^{(1)}], \mathscr{I}^{(2)} = [\tau^{(1)}, \tau^{(2)}]$, where $\tau^{(0)} = 0, \tau^{(2)} = 2$. There is one positive real root $\omega^{(1)}(\tau)$ of $F_1(\omega)$ for $\tau \in (0, \tau^{(1)})$. As $\tau$ reaches $\tau^{(1)}$, this solution merges with the negative solution $-\omega^{(1)}(\tau)$, and they become complex as $\tau$ enters $\mathscr{I}^{(2)}$, and $F_1(\omega)$ does not have any real solution for $\tau$ in $\mathscr{I}^{(2)}$. In this case, we have

$$\omega^{(1)}(\tau) = \sqrt{c_1c_3^2e^{-c_2\tau} - c_2^2}.$$

Corresponding to $\omega = \omega^{(1)}(\tau)$, $\theta^{(1)}(\tau)$ defined in (26) is plotted against $\tau$ in the top diagram of Figure 1. It can be seen that the curve intersects the horizontal line $2\pi$ at $\tau_1 \approx 0.2748$
and \( \tau_2 \approx 0.5314 \). Therefore, \( H_1 = 1, \omega_{h1} = \omega_1^{(1)}(\tau_1) \approx 3.6490 \), and \( H_2 = 1, \omega_{h2} = \omega_1^{(1)}(\tau_2) \approx 2.5228 \). Since both \( \tau_1 \) and \( \tau_2 \) are different from \( \tau^{(1)} \), it is easy to verify that Assumption IIIa holds because (9) does not hold for each \((\omega_h, \tau)\). Assumption III is further implied by Assumption IIIa.

Since \( \text{ord}(P_2) = 2 \) and \( \text{ord}(Q_2) = 1 \), Assumption I holds. Assumption II requires the following two equations do not hold simultaneously for real \( \omega \) and \( \tau \in \mathcal{S} \):

\[
\begin{align*}
-\omega^2 + aj \omega + c &= 0, \\
b(\tau)\omega + d(\tau) &= 0,
\end{align*}
\]

which can be easily verified to be true. The function \( F \) in this case is

\[
F(\omega, \tau) = \omega^4 + (2^2 - \omega^2 - 2c)\omega^2 + c^2 - d^2(\tau),
\]

Solving (5) and (9) together for \((\omega, \tau) \in \mathbb{R}_+ \times \mathcal{S}\), we obtain two pairs of solutions approximately equal to \((0, 1.981)\) and \((0.720, 2.391)\). The interval \( \mathcal{S} \) is thus partitioned into three subintervals \( \mathcal{S}^{(1)} = [\tau^{(0)}, \tau^{(1)}], \mathcal{S}^{(2)} = [\tau^{(1)}, \tau^{(2)}], \mathcal{S}^{(3)} = [\tau^{(2)}, \tau^{(3)}] \), where \( \tau^{(0)} = 0, \tau^{(1)} \approx 1.981, \tau^{(2)} \approx 2.391, \tau^{(3)} = 2.5 \). The polynomial \( F_2(\omega) \) has one positive real root, namely \( \omega_1^{(1)}(\tau) \), in the interval \( (\tau^{(0)}, \tau^{(1)}) \) and two positive roots, namely \( \omega_1^{(2)}(\tau) \) and \( \omega_2^{(2)}(\tau) \), in the interval \( (\tau^{(1)}, \tau^{(2)}) \). It has no real root for \( \tau \in (\tau^{(2)}, \tau^{(3)}) \). We have the following expressions:

\[
\begin{align*}
\omega_1^{(1)}(\tau) &= 2^{-1/2} \sqrt{(b^2(\tau) + 2c - a^2) + \Delta/2(\tau), \tau \in \mathcal{S}^{(1)},} \\
\omega_1^{(2)}(\tau) &= 2^{-1/2} \sqrt{(b^2(\tau) + 2c - a^2) + \Delta/2(\tau), \tau \in \mathcal{S}^{(2)},} \\
\omega_2^{(2)}(\tau) &= 2^{-1/2} \sqrt{(b^2(\tau) + 2c - a^2) - \Delta/2(\tau), \tau \in \mathcal{S}^{(2)},}
\end{align*}
\]

where \( \Delta(\tau) = (b^2(\tau) + 2c - a^2)^2 - 4(c^2 - d^2(\tau)) \). We observe that \( \pm \omega_2^{(2)}(\tau) \) emerge as a pair of real roots of \( F_2(\omega) \) at \( \tau = \tau^{(1)} \) and \( \omega_1^{(2)}(\tau) = 0 \). As \( \tau \) approaches \( \tau^{(2)} \) from the left, the solution \( \omega_1^{(2)}(\tau) \) merges with \( \omega_2^{(2)}(\tau) \). These two positive roots become complex as \( \tau \) increases beyond \( \tau^{(2)} \). The corresponding phase functions \( \theta_1^{(1)}(\tau), \theta_1^{(2)}(\tau), \theta_2^{(2)}(\tau) \) are plotted against \( \tau \) in the top diagram of Figure 2. These curves intersect the horizontal line \( 0 \) at \( \tau_1 \approx 0.7576 \) and \( \tau_2 \approx 2.1745 \). Therefore, \( H_1 = 1, \omega_{h1} = \omega_1^{(1)}(\tau_1) \approx 2.7556 \) and \( H_2 = 1, \omega_{h2} = \omega_1^{(2)}(\tau_2) \approx 1.1837 \). Since both \( \tau_1 \) and \( \tau_2 \) are different from either \( \tau^{(1)} \) or \( \tau^{(2)} \), it is easy to see that (9) does not hold for each \((\omega_h, \tau)\). Consequently we deduce that Assumption IIIa must hold, which also implies Assumption III.

It can be verified that \( \partial_\omega F(\omega_1^{(1)}(\tau), 1) > 0, \partial_\omega F(\omega_1^{(2)}(\tau), 2) > 0 \), therefore \( \partial_\omega F(\omega_1^{(1)}(\tau), \tau) > 0 \) for \( \tau \in (\tau^{(0)}, \tau^{(1)}) \) and \( \partial_\omega F(\omega_1^{(2)}(\tau), \tau) > 0 \) for \( \tau \in (\tau^{(1)}, \tau^{(2)}) \). Computation shows that

\[
\begin{align*}
\frac{d}{d\tau} \theta_1^{(1)}(\tau) > 0, \quad \frac{d}{d\tau} \theta_1^{(2)}(\tau) < 0,
\end{align*}
\]

which also follows from the graph of phase functions plotted in the top diagram of Fig.2. We deduce by using (39) that a pair of characteristic roots cross the imaginary axis from the left-half plane to the right-half plane as \( \tau \) increases through \( \tau_1 \). Another pair of characteristic roots cross the imaginary axis from the right-half plane to the left-half plane as \( \tau \) increases through \( \tau_2 \). Consequently, we have \( \text{Inc}(\omega_{h1}, \tau_1) = 1 \).
increases and goes through $\tau_1$ and $\tau_2$ respectively. The system has two unstable characteristic roots for $\tau = 0$, therefore it is asymptotically stable for $\tau \in (\tau_1, \tau_2)$ and unstable for $\tau \in [0, \tau_1) \cup (\tau_2, 1]$.

VI. CONCLUSION

A method of stability analysis for time-delay systems with coefficients depending on the delay has been developed. The method is an extension of the one given in [12] to a more general case. The method partitions the range of interest into sub-intervals so that the magnitude condition yields a fixed number of solutions of frequencies $\omega$ as functions of the delay $\tau$ within each subinterval. The crossing conditions is expressed in a general form, and a simplified derivation for the first order derivative crossing criterion is obtained.

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