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# HIGHER-ORDER SENSITIVITY TO IMPERFECTIONS IN BIFURCATION BUCKLING **ANALYSIS**

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Abstract  $\cdots$  An alternative to the usual procedure of determining the maximum loads and corresponding coordinates as expansions in terms of an imperfection parameter is presented. The present approach is direct and does not require the development of inverse expansions. The procedure is based on the concept of least-degeneracy, which has also been used as the basis for the analysis in a wide variety of other applications. Use of the method is demonstrated by applying it to two examples : one symmetric bifurcation and one asymmetric bifurcation. It is shown that the present method very efficiently reproduces known results. A brief discussion of the consistent retention of terms in the various expansions is also presented.

#### **L. INTRODUCTION**

The general theory of elastic stability was originally formulated for continuous systems by Koiter (1945), and a number of researchers have contributed to a similar development for discrete systems. An account of the achievements in the theory in terms of generalized coordinates may be found in the books by Croll and Walker (1972). Thompson and Hunt (1973), Huseyin (1975) and, more recently, El Naschie (1990). Distinct critical states in the theory are found to be of two kinds: limit points and bifurcations points, and special attention is given to the sensitivity of critical states with respect to structural imperfections. An account of imperfection sensitivity may be found in the literature mentioned above and in the many references cited therein.

For critical states that are characterized as bifurcations, the usual approach to imperfection sensitivity is to define a new parameter  $\varepsilon$ , which serves as a measure of the amplitude of a given imperfection, and write the total potential energy  $V$  of the system in terms of the usual control  $\lambda$  and response Q, parameters plus the new imperfection parameter. A critical state satisfies two conditions: equilibrium (associated with the zero of the first variation of V) and critical stability (associated with the zero of the second variation of V) (Croll and Walker, 1972; Thompson and Hunt, 1973; Huseyin, 1975; El Naschie, 1990). Imperfection sensitivity is obtained by perturbing these two equations in order to find the critical state as a function of the imperfection amplitude. This dependence is usually written as  $\lambda_c = \lambda_c(\varepsilon)$ . The problem is that in bifurcation analysis such a relation has an infinite slope at  $\varepsilon = 0$  and the expansion cannot be made from that point by the use of regular perturbation methods.

One solution to this problem is described by Thompson and Hunt (1973). The perturbation is carried out with respect to another parameter, say  $Q_1$ , and series are found in the form  $\lambda_c = \lambda_c(Q_1)$ ;  $\epsilon = \epsilon(Q_1)$ :  $Q_{\kappa} = Q_{\kappa}(Q_1)$ , where c denotes critical values and  $j > 1$ . Then the middle expansion is inverted to provide  $Q_+ = Q_+(v)$ , and the result is substituted

into the two remaining series to produce the desired expansions. The three primary expansions are regular and can be obtained in the usual straightforward way. The final expansions are in terms of fractional powers of  $\varepsilon$  and have infinite slopes at  $\varepsilon = 0$ .

In this paper, we describe an alternative to this procedure, which does not require intermediate expansions. The relationship between  $\lambda_c$  and  $\varepsilon$  as well as those between the  $Q_k$ and  $\varepsilon$  are determined directly. The straightforward procedure of assuming expansions of the zeros in integral powers of  $\varepsilon$  does not produce results because the zeros of the first and second variations of the potential energy when  $\varepsilon = 0$  are always repeated roots. A consequence of repeated roots is that the expansions do not go in integral powers of  $\varepsilon$ . Indeed the powers of  $\varepsilon$  cannot be assumed *a priori* and must be determined as part of the solution. Similar problems arise in the analysis of dynamic systems when two or more equilibrium points coalesce. Moreover. in analysing boundary-layer phenomena, whether in plates. electrical conductors. or high Reynolds number flows, one must "stretch" one or more coordinates. Typically, the stretching goes according to a nonintegral power of a naturally occurring. small parameter (such as the square root of the reciprocal of the Reynolds number in the case of flows). Determining the proper stretching in these singularperturbation problems is a key element of the analysis, and that procedure is quite similar to the one employed here. The problem is discussed in the contexts of single algebraic equations and boundary layers in the text by Nayfeh (1981).

### 2. A SUMMARY OF THE AVAILABLE RESULTS

We shall follow the perturbation of Thompson and Hunt (1972). because it develops the formulation not only for first, but also for higher order terms in the imperfectionsensitivity perturbation analysis. The total potential energy of the system is given by  $V = V(Q_i, \lambda)$ ; the conditions of equilibrium and critical stability are written as

$$
V_{\parallel} = \frac{\partial V}{\partial Q_i} = 0 \tag{1}
$$

$$
V_{ij}x_j = \frac{\partial^2 V}{\partial Q_i \partial Q_j} x_j = 0,
$$
\n(2)

where eqn (2) is an eigenvalue problem in terms of the eigenvalue  $\lambda$  and the eigenvector  $x_i$ . Typically. eqns (I) and (2) have multiple roots: this is demonstrated in the section on applications below.

Following the general theory of elastic stability, one studies the critical state for  $\varepsilon = 0$ by writing the perturbation expansions

$$
\lambda(s) = \lambda_s + \lambda^{(1)}s + \frac{1}{2}\lambda^{(2)}s^2 + \cdots
$$
 (3)

$$
Q_i(s) = Q_i^s + Q_i^{s+1} s + \frac{1}{2} Q_i^{s+2} s^2 + \cdots,
$$
 (4)

where *s* is a suitable perturbation parameter (say  $s = Q_1$ ) and the coefficients are the derivatives

$$
(\ )^{(n)} \equiv \frac{\partial^n (\ )}{\partial s^n}\bigg|_{c}
$$

evaluated at the critical state.

For the imperfection-sensitivity analysis, we include the new parameter  $\varepsilon$  (a measure of the amplitude of the imperfection considered), so that  $V = V(Q_i, \lambda; \varepsilon)$ . With reference to



Fig. 1. (a) Equilibrium paths for different imperfection amplitudes. (b) Imperfection-sensitivity of the maximum load.  $\lambda^M = \lambda_{\text{max}}$ 

Fig. 1, eqns (1) and (2) represent the conditions of the maximum in each curve, for a specific imperfection. The particular example presented in Fig. 1 is an unstable symmetric bifurcation when  $\varepsilon = 0$ , but the imperfection breaks the bifurcation and leads to a nonlinear path with a maximum  $\lambda_{\text{max}}(\varepsilon)$ . At  $\varepsilon = 0$  in Fig. 1b, the curve is tangent to the  $\lambda$  axis; this means that a regular perturbation analysis cannot be carried out to represent the curve. Hunt (1971) proposed an alternative perturbation scheme by choosing an intermediate variable as the perturbation parameter, namely  $s = Q_1$ . Thus, the following equations are obtained:

$$
\lambda_{\max} = \lambda_c + \lambda_1 Q_+ + \frac{1}{2} \lambda_2 Q_+^2 + \cdots
$$
 (5)

$$
Q_{\text{max}} = Q_{\kappa} + Q_{\ell 1} Q_{\ell} + \frac{1}{2} Q_{\ell 2} Q_{\ell 1}^2 + \cdots
$$
 (6)

$$
\varepsilon = \varepsilon_0 + \varepsilon_1 Q_1 + \frac{1}{2} \varepsilon_2 Q_1^2 + \frac{1}{3!} \varepsilon_3 Q_1^3, \tag{7}
$$

where  $\lambda_1$ ,  $Q_{j1}$  and  $\varepsilon_1$  are the first derivatives with respect to  $Q_1$ : and  $\lambda_2$ ,  $Q_{j2}$  and  $\varepsilon_2$  are the second derivatives. The difference between eqns  $(5)$   $(7)$  and  $(3)$  and  $(4)$  is that the latter are computed by perturbing eqn (1) for  $\varepsilon = 0$ ; while the former are calculated from perturbations of (1) and (2) for  $\varepsilon \neq 0$ . Once the coefficients of eqns (5)–(7) are calculated, the series are inverted to obtain  $\lambda_{\text{max}} = \lambda(\varepsilon)$ . The explicit form for these series, for the case of asymmetric bifurcation, is

$$
\lambda = \lambda_c + \alpha(c)^{1/2} + \beta(c) + \gamma(c)^{-1/2} + \cdots. \tag{8}
$$

where

$$
\alpha = \pm \left(\frac{2}{\epsilon_2}\right)^{1/2}
$$
 (9a)

$$
\beta = \frac{\lambda_2}{\varepsilon_2} - \frac{\lambda_1 \varepsilon_3}{3(\varepsilon_2)^2}
$$
 (9b)

$$
y = \pm \left(\frac{2}{\epsilon_2}\right)^{1/2} \left\{\frac{\lambda_3}{3\lambda_2} - \frac{\lambda_2 \epsilon_3}{3(\epsilon_2)^2} + \frac{\lambda}{36(\epsilon_2)}\right\} \left[5(\epsilon_3)^2 - 3\epsilon_2 \epsilon_4\right] \bigg\}.
$$
 (9c)

In eqns (8) and (9), the notation  $1.2 +$  means that only the positive square root is considered, and the sign of  $\gamma$  depends upon the sign of  $\alpha$ . For the symmetric bifurcation the sensitivity equation is

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$$
\lambda_{\max} = \lambda_c + \alpha \varepsilon^{2/3} + \beta \varepsilon + \gamma \varepsilon^{4/3} + \cdots, \qquad (10)
$$

where

$$
\alpha = \frac{1}{2}\lambda_2 \left(\frac{6}{\varepsilon_2}\right)^{2/3} \tag{11a}
$$

$$
\beta = \frac{1}{2} \left[ \frac{2\lambda_3}{\epsilon_3} - \frac{\lambda_2 \epsilon_4}{\epsilon_3} \right]
$$
(11b)

$$
\gamma = \left(\frac{6}{\epsilon_3}\right)^{1/3} \left[\frac{\lambda_4}{4\epsilon_3} - \frac{\lambda_3 \epsilon_4}{4(\epsilon_3)^2} - \frac{\lambda_2 \epsilon_5}{10(\epsilon_3)^2} + \frac{7}{48} \frac{\lambda_2 (\epsilon_4)^2}{(\epsilon_3)^3}\right].
$$
 (11c)

## 3. PROPOSED SOLUTION

Rather than using an intermediate perturbation parameter, as in eqns  $(5)-(7)$ , we employ  $s = \varepsilon$ . We assume that the expansions for the maximum load,  $\lambda_{\text{max}}$ , and the corresponding coordinate,  $Q_{\text{max}}$ , have the following forms

$$
\lambda_{\text{max}} = \lambda_c + \lambda_1 \varepsilon^M + \lambda_2 \varepsilon^R + \lambda_3 \varepsilon^T + \cdots \tag{12a}
$$

$$
Q_{jmax} = Q_{jk} + Q_{j1} \varepsilon^N + Q_{j2} \varepsilon^P + Q_{j3} \varepsilon^S + \cdots, \qquad (12b)
$$

where  $j = 1, 2, \cdots K$ ; and the positive exponents M, N, P, R, S, and T as well as the coefficients  $\lambda_c$ ,  $\lambda_1$ ,  $\lambda_2$ , ... and  $Q_k$ ,  $Q_{i1}$ , ... are unknown. *K* is the order (i.e. number of coordinates needed to describe the state) of the system. We note that the coefficients  $\lambda_1$ ,  $\lambda_2$ ,  $\cdots$ ,  $Q_{i1}$ ,  $Q_{i2}$ ,  $\cdots$  in eqns (12a, b) are not related to those in eqns (5)–(7).

First, we substitute eqns (12a, b) into eqns (1) and (2); then we set  $\varepsilon = 0$  and solve for  $\lambda_c$  and  $Q_s$ . This is precisely the stability problem for no imperfections. Next, we include  $\lambda_1$ and the  $Q_i$  as well as  $\lambda_c$  and the  $Q_{\kappa}$ , but consider the remaining  $\lambda_n$  and  $Q_{i\eta}$  to be zero; then we select the heretofore arbitrary exponents *M* and *N* so as to obtain the "least-degenerate" two-term approximation (we shall explain and illustrate the concept of least degeneracy in the examples that follow); finally, we solve for  $\lambda_1$  and the  $Q_{i1}$ . Next we add  $\lambda_2$  and the  $Q_{i2}$ to the expansion; and repeat the procedure. Any desired number of terms may be included.

The proposed technique is applied in the next two sections to problems of symmetric and asymmetric bifurcation with one degree of freedom, in order to highlight the main features of the procedure.

#### 4. SYMMETRIC BIFLRCATION

The first example is a one-degree-of-freedom problem discussed by Croll and Walker (1972 : Chap. 5) and illustrated in Fig. 2. The exact form of the total potential energy is



Fig. 2. Example considered for symmetric bifurcation.

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$$
V(Q, \lambda; \varepsilon) = (\sin Q - \sin \varepsilon)^2 - 2\lambda(\cos \varepsilon - \cos Q), \tag{13}
$$

where  $\lambda \equiv F/KL$ , and Q and  $\epsilon$  have the meaning of angles, as shown in Fig. 2. Here we do not need a subscript on *Q.*

We begin by expanding the trigonometric functions as follows:

$$
\sin Q = Q - \frac{1}{6}Q^3 + \frac{1}{120}Q^5
$$
 (14a)

$$
\cos Q \doteq 1 - \frac{1}{2}Q^2 + \frac{1}{24}Q^4 - \frac{1}{720}Q^6 \tag{14b}
$$

$$
\sin \varepsilon = \varepsilon \tag{14c}
$$

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$$
\cos \varepsilon \doteq 1 \tag{14d}
$$

Next we substitute eqns (14a-d) into (13), retain terms up to and including fourth order, and obtain the following approximate expression for *V:*

$$
V(Q, \lambda; \varepsilon) = Q^2 - \frac{1}{3}Q^4 - \lambda \left(Q^2 - \frac{1}{12}Q^4\right) - \varepsilon \left(2Q - \frac{1}{3}Q^3\right).
$$
 (15a)

The equilibrium condition is given by

$$
\frac{1}{2}\frac{dV}{dQ} = Q_{\text{max}} - \frac{2}{3}Q_{\text{max}}^3 - \lambda_{\text{max}}\left(Q_{\text{max}} - \frac{1}{6}Q_{\text{max}}^3\right) - \varepsilon\left(1 - \frac{1}{2}Q_{\text{max}}^2\right) = 0
$$
 (15b)

and the condition of critical stability is given by

$$
\frac{1}{2}\frac{d^2V}{dQ^2} = 1 - 2Q_{\text{max}}^2 - \lambda_{\text{max}}\left(1 - \frac{1}{2}Q_{\text{max}}^2\right) - \varepsilon Q_{\text{max}} = 0. \tag{15c}
$$

When  $\varepsilon = 0$ , it follows from eqns (12a, b) that  $\lambda_{\text{max}} = \lambda_c$  and  $Q_{\text{max}} = Q_c$ ; then eqns (15b, c) reduce to

$$
6(1 - \lambda_c)Q_c - (4 - \lambda_c)Q_c^3 = 0
$$
 (16a)

and

$$
2(1 - \lambda_c) - (4 - \lambda_c)Q_c^2 = 0.
$$
 (16b)

The roots are  $\lambda_c = 1$  and  $Q_c = 0$ . We note the repeated roots.

To determine the influence of the imperfection  $\varepsilon$ , we begin by including the second terms in eqns (12a, b)

$$
\lambda_{\text{max}} = 1 + \lambda_1 \varepsilon^M \tag{17a}
$$

$$
Q_{\max} = Q_1 \varepsilon^N, \tag{17b}
$$

where *M* and *N* are arbitrary, but greater than zero, and  $\lambda_1$  and  $Q_1$  are unknown. Substituting eqns (I7a, b) into (ISb, c), we obtain

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$$
\frac{1}{2}\frac{dV}{dQ} = Q_1^2 \varepsilon^{2M+1} - Q_1^3 \varepsilon^{3M} - 2\varepsilon - \frac{1}{12}Q_1^4 \varepsilon^{4M+1} - 2\lambda_1 Q_1 \varepsilon^{M+N} + \frac{1}{3}\lambda_1 Q_1^3 \varepsilon^{3M+N} = 0 \quad (18a)
$$

$$
\frac{1}{2}\frac{d^2V}{dQ^2} = 2Q_1\varepsilon^{M+1} - 3Q_1^2\varepsilon^{2M} - \frac{1}{3}Q_1^3\varepsilon^{3M-1} - 2\lambda_1\varepsilon^N + \lambda_1Q_1^2\varepsilon^{2M+N} = 0.
$$
 (18b)

Some reduction in the number of terms can be made immediately: for small *s* and positive M and *N*,  $\varepsilon^{4M+1} \ll \varepsilon^{2M+1}$ ,  $\varepsilon^{3M+N} \ll \varepsilon^{3M}$ ,  $\varepsilon^{2M+N} \ll \varepsilon^{2M}$ , and  $\varepsilon^{3M+1} \ll \varepsilon^{M+1}$  regardless of the choice of M and N, and the corresponding terms can be ignored in eqns (18a, b). The result is

$$
Q_1^2 e^{2M+1} - Q_1^3 e^{3M} - 2\varepsilon - 2\lambda_1 Q_1 e^{M-N} = 0
$$
 (18c)

$$
2Q_{+}\varepsilon^{M+1} - 3Q_{+}^{2}\varepsilon^{2M} - 2\lambda_{+}\varepsilon^{N} = 0.
$$
 (18d)

To obtain the least-degenerate approximation from eqns (18c, d), we select  $M$  and  $N$ so that these equations retain the maximum number of terms and, hence, produce the maximum amount of information. Ideally, if it were possible, we would choose *M* and *N* such that

$$
2M + 1 = 3M = 1 = M + N.
$$
 (19a)

and

$$
M+1 = 2M = N.\t(19b)
$$

Generally, this is not possible and the proper selection of  $M$  and  $N$  will only satisfy some of the relationships given in eqns (19a, b). The idea is to choose the "'least-degenerate" combination. Next we consider the various possibilities one by one.

If we choose  $2M + 1 = 3M$  in eqn (19a), then  $M = 1$ , and it follows from eqn (19b) that  $N = 2$ . There is a single dominant term in eqn (18c), and it is  $-2\varepsilon$ , which leads to the inconsistent conclusion  $2 = 0$  if  $\varepsilon \neq 0$ . Hence, this is not the proper choice for M and N.

If we choose  $2M+1 = 1$ , then  $M = N = 0$ . This choice will not produce any new information. It places  $Q_1$  and  $\lambda_1$  on the same level with the solution corresponding to  $\varepsilon = 0$ ; hence,  $Q_1$  and  $\lambda_1$  are zero. This also is not the proper choice for *M* and *N*.

If we choose  $3M = 1$ , then  $M = 1/3$  and  $N = 2/3$ . Equations (18c, d) reduce to

$$
\varepsilon(-Q_1^3 - 2 - 2\lambda_1 Q_1) + O(\varepsilon^{3/3}) = 0 \tag{20a}
$$

and

$$
\varepsilon^{2/3}(-3Q_1^2 - 2\lambda_1) + O(\varepsilon^{4/3}) = 0.
$$
 (20b)

From which we find

$$
Q_1 = 1 \tag{21a}
$$

and

$$
\lambda_1 = -3/2. \tag{21b}
$$

This is the only choice for *M* and *N* that does not lead to inconsistencies and produces new information. Equations (20a. b) are called the least-degenerate form of eqns (l8c, d), and the results in eqns  $(21a, b)$  are called the least-degenerate approximation. So far the expansions have the following forms:

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$$
Q_{\text{max}} = \varepsilon^{1/3} \tag{22a}
$$

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and

$$
\lambda_{\max} = 1 - \frac{3}{2} \varepsilon^{2/3}.
$$
 (22b)

These results agree with eqn  $(10)$ .

To continue the process, we now add more terms to the expansions in eqns  $(12a, b)$ :

$$
Q_{s} = \varepsilon^{1/3} + Q_{2} \varepsilon^{p}
$$
 (23a)

$$
\lambda_{s} = 1 - \frac{3}{2}e^{2/s} + \lambda_{2}e^{R}.
$$
 (23b)

We can follow the procedure outlined above to determine  $P$  and  $R$ , or we can be guided by eqns (20) and (22) and choose  $P = 1$  and  $R = 4/3$ . (One can readily verify that the other choices for P and R lead to inconsistencies.) Then substituting eqns (23a, b) into eqns (l5b, c) leads to

$$
\varepsilon^{s-1}(1-4\lambda_2) + 0(\varepsilon^{-1}) = 0 \tag{24a}
$$

$$
\varepsilon^{4/3} (1 - 4\lambda_2 - 12Q_2) + O(\varepsilon^2) = 0.
$$
 (24b)

From which we find that

$$
Q_2 = 0 \tag{25a}
$$

and

$$
\lambda_2 = \frac{1}{4} \tag{25b}
$$

and that so far the expansions have the following forms:

$$
Q_{\text{max}} = \varepsilon^{1/3} + O(\varepsilon^{3/3})
$$
\n(26a)

$$
\lambda_{\text{max}} = 1 - \frac{3}{2} \varepsilon^{2/3} + \frac{1}{4} \varepsilon^{4/3} + O(\varepsilon^2). \tag{26b}
$$

The order of the truncation error in eqn (26b) is  $\vec{c}$ ; thus. if more terms in the expansions for  $Q_{\text{max}}$  and  $\lambda_{\text{max}}$  are to be found, then to be consistent we must include the second term in the expansion of  $\cos \varepsilon$  given in eqn (14d). A related question is how many terms in the expansions of  $Q_{\text{max}}$  and  $\lambda_{\text{max}}$  are consistent with the fourth-order expansion of *V* given in eqn (15a). To answer this question, we include more terms in the expansion of  $V$  and write

$$
V(Q, \lambda, \varepsilon) = Q^2 - \frac{1}{3}Q^4 + \frac{2}{45}Q^5 - \lambda \left(Q^3 - \frac{1}{12}Q^4 + \frac{1}{360}Q^6\right) - \varepsilon \left(2Q - \frac{1}{3}Q^3 + \frac{1}{60}Q^5\right) + \cdots
$$
\n(27)

When we repeat the procedure described above. we tind that when we include only fifth order terms in the expression for  $V$  the results are the same as those in eqns (26). But when sixth (and higher) order terms are included, we obtain

$\epsilon$ . $\lambda_{\text{max}}$	First order Eqn $(20)$	Second order	
		Eqn $(24)$	Eqn (30)
0.001	0.985		
0.005	0.956		
0.010	0.930		
0.050	0.796	0.800	0.803
0.100	0.677	0.689	0.694
0.200	0.487	0.516	0.531
0.300	0.328	0.378	0.403
0.400	0.186	0.259	0.296

Table 1. Maximum load.  $\lambda_{\text{max}}$ , for different values of the imperfection parameter,  $\varepsilon$ , for symmetric bifuration

$$
Q_{\max} = \varepsilon^{1/3} + \frac{1}{6}\varepsilon + O(\varepsilon^{5/3})
$$
 (28a)

$$
\lambda_{\max} = 1 - \frac{3}{2} \varepsilon^{3/2} + \frac{3}{8} \varepsilon^{4/3} + O(\varepsilon^2). \tag{28b}
$$

Comparing eqns (26) and (28), we find the last terms in the two sets of expansions are different; those in eqns (28a, b) are correct. Hence, if a three-term expansion for  $\lambda_{\text{max}}$  is desired, then *V* must contain sixth order terms.

For the present example, the results of first and second order perturbation analyses are compared in Table 1. As expected, the differences between first-order and second-order solutions in  $\lambda_{\text{max}}(\varepsilon)$  grow with  $\varepsilon$ . For example, for a value  $\varepsilon = 0.2$ , a first order solution would predict  $\lambda_{\text{max}} = 0.487$ , while a second order solution computed from a quartic functional leads to  $0.516$ . If the approximations in the function  $V$  are improved so that it includes terms with Q to the sixth power, the value of  $\lambda_{\text{max}} = 0.531$ . The quadratic approximation is conservative in the sense that it predicts values that are lower than the correct ones. In this case the first order perturbation solution is conservative with respect to the second order one.

#### 5. ASYMMETRIC BIFURCATION

As a second example, we consider the one-degree-of-freedom model studied in Croll and Walker (1972: Chap. 6) and shown in Fig. 3. The exact expression for the potential energy in this case is given by

$$
V(\Delta_1, \Delta_2, F; \varepsilon) = \frac{1}{2} K \Delta_2^2 - F \Delta_1,
$$
 (29a)

where



Fig. 3. Example considered for asymmetric bifurcation.

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$$
\Delta_1 = 2L(\cos \varepsilon - \cos Q) \tag{29b}
$$

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$$
\Delta_2 = \sqrt{2L[(1-\sin\epsilon)^{1/2} - (1-\sin Q)^{1/2}]}
$$
 (29c)

$$
F = \frac{1}{4} KL\lambda.
$$
 (29d)

Next we substitute the following expansions

$$
(1 - \sin Q)^{1/2} = 1 - \frac{1}{2}Q - \frac{1}{8}Q^2 + \frac{1}{48}Q^3 + \frac{1}{384}Q^4 - \frac{1}{3840}Q^5
$$
  

$$
\cos Q = 1 - \frac{1}{2}Q^2 + \frac{1}{24}Q^4
$$
  

$$
(1 - \sin \varepsilon)^{1/2} = 1 - \frac{1}{2}\varepsilon
$$
  

$$
\cos \varepsilon = 1
$$
 (30)

into eqns (29) and obtain the following approximate expression for  $V$ :

$$
V(Q_1, \lambda; \varepsilon) = Q^2 + \frac{1}{2}Q^3 - \frac{1}{48}Q^4 - \lambda \left(Q^2 - \frac{1}{12}Q^4\right) - \varepsilon \left(2Q + \frac{1}{2}Q^2 - \frac{1}{12}Q^3 - \frac{1}{96}Q^4\right).
$$
 (31a)

The equilibrium condition is given by

$$
\frac{1}{2}\frac{dV}{dQ} = Q_{\text{max}} + \frac{3}{4}Q_{\text{max}}^2 - \frac{1}{24}Q_{\text{max}}^3 + \lambda_{\text{max}}\left(Q_{\text{max}} - \frac{1}{6}Q_{\text{max}}^3\right) - \varepsilon\left(1 + \frac{1}{2}Q_{\text{max}} - \frac{1}{8}Q_{\text{max}}^2 - \frac{1}{48}Q_{\text{max}}^3\right) = 0 \quad (31b)
$$

and condition of critical stability is given by

$$
\frac{1}{2}\frac{d^2V}{dQ^2} = 1 + \frac{3}{2}Q_{\text{max}} - \frac{1}{8}Q_{\text{max}}^2 - \lambda_{\text{max}}\left(1 - \frac{1}{2}Q_{\text{max}}^2\right) - \varepsilon\left(\frac{1}{2} - \frac{1}{4}Q_{\text{max}} - \frac{1}{16}Q_{\text{max}}^2\right) = 0. \quad (31c)
$$

Next we substitute eqns (12a, b) into eqns (31b, *c*), set  $\varepsilon = 0$ , and solve for  $\lambda_c$  and  $Q_c$ . The solution is  $\lambda_c = 1$  and  $Q_c = 0$ . To determine the influence of the imperfection, we substitute  $\lambda_c$  and  $Q_c$  into egns (12a, b), retain the first two terms, substitute the result into eqns (31b, c), and obtain

$$
\frac{1}{2}\frac{dV}{dQ} = -\epsilon - \frac{1}{2}Q_{\perp}\epsilon^{M+1} + \frac{3}{4}Q_{\perp}^{2}\epsilon^{2M} - \lambda_{\perp}Q_{\perp}\epsilon^{M+N} = 0
$$
\n(32a)

$$
\frac{1}{2}\frac{d^2V}{dQ^2} = -\varepsilon + 3Q_1\varepsilon^M - 2\lambda_1\varepsilon^N = 0.
$$
 (32b)

Next, we choose the values M, N.  $\lambda_1$  and  $Q_1$  that produce the least-degenerate approximation. Following the procedure described in the previous section, we find that  $M = N = 1/2$ ,  $Q_1^2 = -4/3$ , and  $\lambda = 3.2Q_1$ . Thus, so far we have

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$$
\lambda_{\max} = 1 \pm \sqrt{-3\varepsilon} \tag{33a}
$$

$$
Q_{\text{max}} = \pm \frac{2}{3} \sqrt{-3\varepsilon}.
$$
 (33b)

This is the same result obtained by Croll and Walker (1972). For values of  $\varepsilon > 0$ ,  $\lambda_c$  and  $Q<sub>c</sub>$  are imaginary: hence, the system is only sensitive to imperfections that correspond to negative values of  $\varepsilon$ . The result is known as the "one-half-power law".

To extend the expansions. we write

$$
\lambda_{\text{max}} = 1 \pm \sqrt{-3\varepsilon + \lambda_2 \varepsilon^R} \tag{34a}
$$

$$
Q_{\text{max}} = \pm \frac{2}{3} \sqrt{-3\varepsilon + Q_2 \varepsilon^P}.
$$
 (34b)

Then we substitute eqns (34a, b) into eqns (31b, c) and find that  $P = R = 1$ ,  $\lambda_2 = -2/3$ , and  $Q_2 = 2(1 + \lambda_2)/3$ . Hence, so far we have

$$
\lambda_{\max} = 1 + \sqrt{-\varepsilon} - \frac{2}{9}(\varepsilon')
$$
 (35a)

$$
Q_{\max} = \frac{2}{3} \sqrt{-\epsilon} + \frac{2}{27} (\epsilon'), \qquad (35b)
$$

where  $\varepsilon' = 3\varepsilon$  and only the positive root is considered.

#### 6. DISCUSSION AND CONCLUSIONS

We have described a new. direct procedure for obtaining the relationships among the critical loads. the critical coordinates. and a measure of the amplitude of the imperfection. Difficulties arise in obtaining these expansions because the roots of the conditions for equilibrium and critical stability are repeated in the absence of an imperfection. Thus, the expansions of the critical loads and coordinates do not go in integral powers of the imperfection describing parameter. and regular perturbation procedures are not applicable. Instead. we develop a direct approach that is similar to the one used in the analysis of singular perturbation problems to ohtain the correct coordinate stretching. In the present approach. we assume the expansions for the critical loads and corresponding coordinates in terms of arbitrary powers of the imperfection describing parameter. Then, by using the principle of least-degeneracy. we determine the coefficients and exponents in the various terms.

Here we demonstrated the procedure by applying it for the first time to two wellknown problems. each having a single degree of freedom. In future work, the procedure will be extended to systems having many degrees of freedom.

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