

By acceptance of this article, the publisher or recipient acknowledges the U.S. Government's right to retain a nonexclusive, royalty-free license in and to any copyright covering the article.

CONF-830198--1

CONF-830198--1

DE86 003082

RADIAL MODE STRUCTURE OF CURVATURE-DRIVEN
INSTABILITIES IN EBT

D. A. Spong

Oak Ridge National Laboratory

MASTER

DISCLAIMER

This report was prepared as an account of work sponsored by an agency of the United States Government. Neither the United States Government nor any agency thereof, nor any of their employees, makes any warranty, express or implied, or assumes any legal liability or responsibility for the accuracy, completeness, or usefulness of any information, apparatus, product, or process disclosed, or represents that its use would not infringe privately owned rights. Reference herein to any specific commercial product, process, or service by trade name, trademark, manufacturer, or otherwise does not necessarily constitute or imply its endorsement, recommendation, or favoring by the United States Government or any agency thereof. The views and opinions of authors expressed herein do not necessarily state or reflect those of the United States Government or any agency thereof.

* Research sponsored by the Office of Fusion Energy, U.S. Department of Energy, under contract DE-AC05-84OR21400 with Martin Marietta Energy Systems, Incorporated.

RADIAL MODE STRUCTURE

OF CURVATURE-DRIVEN

INSTABILITIES IN EBT

D.A. SPONG

Oak Ridge National Lab.

Hot Electron Physics Minisymposium

Jan. 11, 1983

FEATURES OF THIS CALCULATION:

- RETAINS NONLOCAL STRUCTURE OF MODES.
- CONNECTS INNER AND OUTER RING REGIONS TOGETHER IN ONE TREATMENT.
- A SELF-CONSISTENT FINITE B EQUILIBRIUM B FIELD IS USED.
(including dB/dr and d^2B/dr^2)
- A WIDE RANGE OF EBT PARAMETERS HAVE BEEN EXAMINED.
- RELATIVISTIC EFFECTS ARE INCLUDED FOR THE HOT ELECTRON RING.

COLLABORATORS : H. L. BERK , J. W. VAN DAM,
M. N. ROSENBLUTH

ASSUMPTIONS OF PRESENT CALCULATION

- FINITE LARMOR RADIUS EFFECTS NEGLECTED.
- BALLOONING EFFECTS NEGLECTED AND
 $\vec{B} \cdot \vec{\nabla} (\text{EQUILIBRIUM QUANTITIES}) = 0$
- Z - PINCH GEOMETRY LOCALIZED TO RING
REGION USED WITH OUTGOING ENERGY BOUNDARY
CONDITIONS (NATURAL CURVATURE DRIFT).
- DELTA - FUNCTION HOT ELECTRON DISTRIBUTION:
$$F_{hot} = \frac{\delta P_{\perp n}}{\mu_0 B^2} \delta(P_{||}) \delta(u - u_0)$$
- WARM CORE ELECTRONS, COLD IONS.

BASIC EQUATIONS

● MOMENTUM BALANCE:

$$p_i \dot{x}_i = (\nabla \times B) \times B - \nabla P$$

$$P = \sum_i \int F_i \times p \, d^3p$$

$$\rho = m_e T \chi$$

● THE HOT ELECTRON PRESSURE TENSOR MAY BE

WRITTEN AS:

$$P = \sum_i \frac{\int dH \, d\mu \, B}{|p_{||}|} F(r, H, \mu) [4B(I - b_b) + p_{||}^2 b_b]$$

where $c^2 p_{||}^2 = H^2 - 24Bc^2 - m_e c^4$

$$\mu = \frac{p_{||}^2}{2B}$$

- F is obtained from the drift kinetic equation:

$$\frac{\partial F}{\partial t} + v_{\parallel} b \cdot \nabla F + v_0 \cdot \nabla F + H \frac{\partial F}{\partial H} = 0$$

where $v_{\parallel} = \frac{p_{\parallel}}{T} = \pm \frac{[H^2 - 2\mu B^2 c^2 - m^2 c^4]^{1/2}}{T c^2}$

$$v_0 = \frac{\mu b \times \nabla B}{q_j m_j B T} + \frac{p_{\parallel}^2}{q_j m_j B T} b \times (b \cdot \nabla) b + \frac{E_{\parallel} b}{B}$$

$$H = q_j E_{\parallel} v_{\parallel} + q_j E_{\perp} \cdot v_0 + \frac{\mu}{T m_j} \frac{\partial B}{\partial t}$$

- These equations are then combined, transformed to z-pinch geometry, and linearized about perturbed fields E_{\perp} and B_{\parallel} . These may be characterized by a "displacement" ξ :

$$\xi = i \frac{E_{\parallel} b}{\omega B} \exp[-i\omega t + ikz]$$

-14-

THIS RESULTS IN A SECOND ORDER EQUATION FOR ξ_T :

$$\frac{d}{dr} \left(r P \frac{d\xi_T}{dr} \right) - Q \xi_T = 0$$

where $P = \frac{\lambda B^2 (1 + G_1)}{D v_A^2}$

$$Q = \frac{r B^2}{v_A^2} \left[\frac{k^2 \lambda (1 + G_1)}{D} - \frac{\omega^2 \lambda}{v_A^2 D} \right]$$

$$- \frac{k^2 \mu_0 v_A^2}{r B^2} \frac{d}{dr} (P_H + P_C) - \frac{k^2 v_A^2}{r^2 D} (\sigma + G_3)$$

$$+ \frac{2 \lambda \omega k (1 - G_2)}{r \omega_{ci} D} + \frac{k^2 v_A^2 (1 - G_2)^2}{r^2 D}$$

$$+ \frac{v_A^2}{r B^2} \frac{d}{dr} \left[\frac{r \lambda B^2 S}{v_A^2 D} \right]$$

$$S = \frac{1 - G_2}{r} + \frac{\omega k (1 + G_1)}{\omega_{ci}}$$

$$D = 1 + G_1 - \frac{\lambda}{k^2 v_A^2}$$

$$\lambda = \frac{\omega^2 \omega_{ci}^2}{\omega_{ci}^2 - \omega^2}$$

G_1, G_2, G_3 ARE KINETIC INTEGRALS WHICH INVOLVE THE HOT ELECTRON DISTRIBUTION FUNCTION:

$$G_1 = -B \sum_i \int \frac{dp_* d\mu \mu^2}{T m_i} \left[\frac{1}{B} \frac{\partial F}{\partial \mu} + \frac{L_i^* F}{T \Omega} \right]$$

$$G_2 = - \sum_i \int \frac{dp_* d\mu \mu}{T m_i} p_{||}^2 \left[\frac{1}{B} \frac{\partial F}{\partial \mu} + \frac{L_i^* F}{T \Omega} \right]$$

$$G_3 = - \sum_i \int \frac{dp_* d\mu p_{||}^4}{B T m_i} \left[\frac{1}{p_{||}} \frac{\partial F}{\partial p_{||}} + \frac{L_i^* F}{T \Omega} \right]$$

where $\Omega = \omega - \omega_{DB} - \omega_{cv}$

$$L_i^* = \left(\frac{k}{m_i q_i B} \frac{\partial F}{\partial r} + \frac{T \omega_{cv}}{p_{||}} \frac{\partial F}{\partial p_{||}} \right)$$

THE INTEGRALS G_1 , G_2 , G_3 DEPEND ON THE MODEL USED FOR THE HOT ELECTRON DISTRIBUTION.

FOR THE CASE OF A DELTA FUNCTION WITH NO PARALLEL ENERGY:

$$F_{\text{Hot}} = \frac{\tau_0 p_{\perp H}}{\mu_0 B^2} \delta(p_{\perp}) \delta(\omega - \omega_0)$$

$$1 + G_1 = (\omega - \omega_{DB})^{-1} \left\{ \omega [1 + B_1 - B_{\perp H} \frac{\mu_0 B}{2 \tau_0^2 m_e^2 c^2}] \right.$$

$$\left. - \omega_{cv1} [1 + r \frac{d}{dr} (\frac{p_{\perp c}}{B^2})] \right\}$$

where $\omega_{cv1} = - \frac{k \mu_0}{e n_e \tau_0 r}$

$$\omega_{DB} = \frac{k \mu_0}{e n_e \tau_0 B} \frac{dB}{dr}$$

$$G_2 = \frac{B_{\perp c}}{2}$$

$$G_3 = \frac{3 B_{\perp c}}{2}$$

Using these forms for G_1, G_2, G_3 which are accurate for arbitrary Δ_B/R_c and ω/ω_{db} one can make a local approximation:

$$\frac{d^2 \xi_r}{dr^2} = -k_r^2 \xi_r, \quad k_r = \frac{\pi}{\Delta}$$

$$\frac{d\xi_r}{dr} = 0$$

$$\frac{1}{n_c} \frac{dn_c}{dr} = -\frac{1}{\Delta}$$

To obtain a fifth order dispersion relation:

$$y^5 + A y^4 + B y^3 + C y^2 + D y + E = 0$$

where $y = \frac{\omega}{\omega_{cav}}$

A. High Frequency Modes

Compressional Alfvén

Hot Electron interchange ($q_0 \gg 1$)

Retaining A, B and C terms in the 5th order eqn., $B^2 = 4AC$ gives the marginal stability boundary:

$$P \left\{ P^2 \left(1 + \frac{2\tilde{\beta}_c}{\tilde{\beta}_H} \right)^2 - 2P \left[\left(1 + \frac{1}{q} + \tilde{\beta}_c \right) \left(1 + \frac{2\tilde{\beta}_c}{\tilde{\beta}_H} \right) - \frac{4\tilde{\beta}_c}{\tilde{\beta}_H} \right] + \left(1 - \frac{1}{q} - \tilde{\beta}_c \right)^2 \right\} = 4 \left(\frac{\Delta}{R} \right)^2 (\tilde{\beta} - 1) \tilde{\beta}_H \left[P \tilde{\beta}_c \left(1 + \frac{2\tilde{\beta}_c}{\tilde{\beta}_H} - \frac{2}{\tilde{\beta}_H} \right) + \frac{1}{q} (1 - \tilde{\beta}_c) \right]$$

→ Cubic in $P = n_H/n_i$

2 roots near $P \approx 1 \rightarrow$ hot electron interchange

1 root at $P < (\Delta/R)^2 \ll 1 \rightarrow$ compressional Alfvén

$$\left(\tilde{\beta} = \frac{R_c}{2\Delta} \beta, \quad P = \frac{n_H}{n_i}, \quad q = \left(\frac{k}{k_\perp} \right)^2 \frac{v_{cv\perp}}{\Delta \omega_{ci}} \right)$$

Hot Electron Interchange

$P \approx 1 \rightarrow$ neglect R.H.S. \rightarrow quadratic
in P : only one physical
solution (i.e. with $P < 1$):

$$P < P_1 = \left[1 - \left(\frac{1}{q} + \tilde{\beta}_c \right)^{1/2} \right]^2$$

note: if $q < (4 - \tilde{\beta}_c)^{-1} \rightarrow P > 1$
and this mode stabilizes
(actually one goes over to
low freq. hot interchange)

Compressional Alfvén Mode

$P \ll 1$ root \Rightarrow neglect P^2 and P^3
terms

$$P > P_2 = \frac{1}{q} \left(\frac{2\Delta}{R} \right)^2 \tilde{\beta}_H (\tilde{\beta} - 1) (1 - \tilde{\beta}_c) \left(1 - \frac{1}{q} - \tilde{\beta} \right)$$

note: when $\tilde{\beta}_c = 1$ mode stabilizes

$$\text{also, when } 1 - \frac{1}{q} - \tilde{\beta}_c = 0 \quad P_2 \rightarrow \infty$$
$$(q = \left(\frac{k}{k_L} \right)^2 \frac{V_{ce}}{\Delta \omega_{ci}}) \quad P_1 \rightarrow 0$$

Low Frequency Modes ($\omega < \omega_{ci}$)

A. Low-frequency hot electron interchange
($q \ll 1$)

Keep B, C, D terms

(justified when $q \ll 1$ and $\beta_c \ll \beta_h$)

$C^2 = 4BD$ gives:

$$P < P_3 = \frac{1}{4} (k_\perp \Delta)^2 q_0 (1 - \tilde{\beta}_c)^2$$

B. Low-frequency background interchange

Keep C, D, and E terms ($\beta_c \ll \frac{2\Delta}{R_c}$)

Stability achieved in 2 ways:

$$\beta_h > \frac{4\Delta}{R_c} - 2\beta_c$$

(drift reversal)

or

$$\frac{n_h}{n_i} \gtrsim 8q_0 (k_\perp \Delta)^2 \left(\frac{\beta_c}{\beta_h} \right)$$

for $\tilde{\beta}_h \gg 1$,
 $\beta_c / \beta_h < 1$

(charge uncovering, $n_{ce} \neq n_{ci}$
frequency shift $\propto (n_{ci} - n_{ce})/n_{ce} = p$)

C. Interacting Background Interchange
 $(\beta_c \sim 2\Delta/R_c)$

Keeping B, C, D and E terms:

$$[1 + q(1 - \tilde{\beta}_c)]y^3 - (1 - \tilde{\beta}_c)y^2$$

$$+ \tau \left(1 + \frac{2\tilde{\beta}_c}{\tilde{\beta}_H}\right) y + \tau \tilde{\beta}_c \left(1 + \frac{2\tilde{\beta}_c}{\tilde{\beta}_H} - \frac{2}{\tilde{\beta}_H}\right) = 0$$

$$\text{where } \tau = \rho/q_0(k_\perp \Delta)^2$$

note: $\tau < 1$ is stability condition for low-freq.
 hot electron interchange

assuming $\tau \ll (1 - \tilde{\beta}_c)^2/3$ one finds
 that for $(1 - \tilde{\beta}_c)$ small, 2 roots
 of the cubic coalesce and lead
 to instability if:

$$1 - \tilde{\beta}_c < 3 \left[\frac{\tau}{4} \left(1 + \frac{2\tilde{\beta}_c}{\tilde{\beta}_H} - \frac{2}{\tilde{\beta}_H}\right) \right]^{1/3}$$

NUMERICAL SOLUTION

METHOD

- Radial equation: $\frac{1}{r} \frac{d}{dr} (rP \frac{ds}{dr}) - Qs = 0$
- Equivalent to 2 coupled first order equations:

$$\left\{ \begin{array}{l} \frac{dy_1}{dr} = \frac{y_2}{r \cdot P} \\ \frac{dy_2}{dr} = rQy_1 \end{array} \right.$$

where $y_1 = s$, $y_2 = r \cdot P \frac{ds}{dr}$

- Solve as 2-point boundary value problems on intervals r_{min} to r_m and r_m to r_{max} using the SUPPORT code.
- Powell's hybrid method is used to solve for ω by matching:

$$\left. \frac{dy_1}{dr} \right|_{r=r_m^+} = \left. \frac{dy_1}{dr} \right|_{r=r_m^-}$$

PLASMA AND HOT ELECTRON PROFILES
ARE CHOSEN WHICH ARE FLAT INSIDE
AND OUTSIDE THE RING AND WHICH
HAVE CONTINUOUS DERIVATIVES.

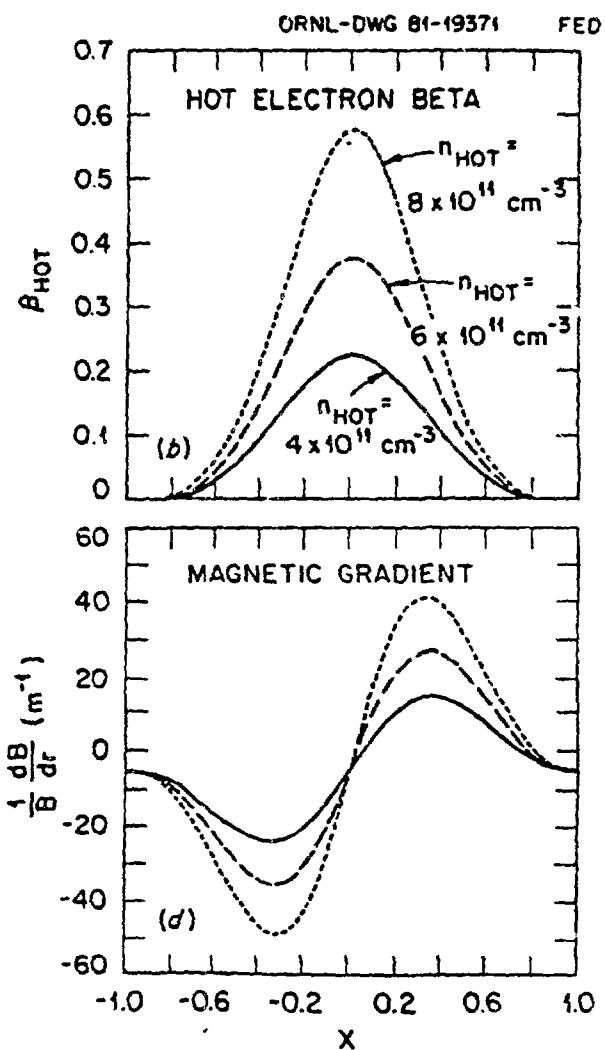
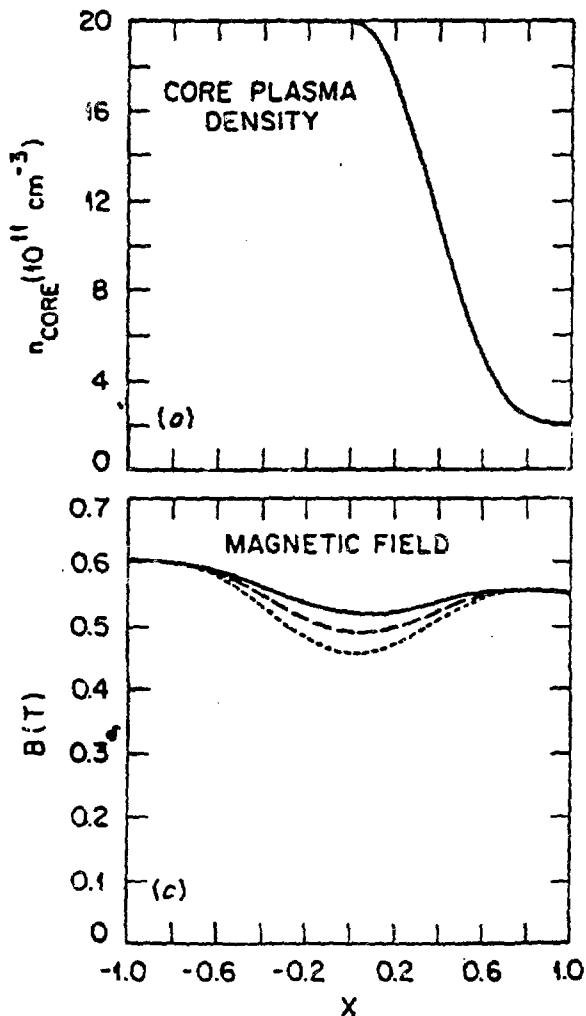
$$P_{\perp H}(r) = P_{\perp HO} \begin{cases} 0 & x < -x_0 \\ \frac{(x-x_0)^4(x+x_0)^4}{x_0^8} & -x_0 < x < x_0 \\ 0 & x > x_0 \end{cases}$$

$$P_c(r) = P_{co} \begin{cases} 1 & x < 0 \\ (1-s) \frac{(x-x_0)^4(x+x_0)^4}{x_0^8} + s & 0 < x < x_0 \\ s & x > x_0 \end{cases}$$

where $x = \frac{r-r_0}{\Delta}$

Δ = hot electron annulus
half-width

x_0 = parameter which controls
extent of pressure profile
 s = density shelf factor



A MAGNETIC EQUILIBRIUM MODEL
IS USED WITH FIELD LINE
CURVATURE APPROPRIATE TO
Z-PINCH GEOMETRY ($\kappa = -1/r$).

FROM:

$$\vec{r} \left(p_{\perp} + \frac{B^2}{2\mu_0} \right) = \vec{k} B^2 \left[1 - \frac{\mu_0(p_{\parallel H} - p_{\perp})}{B^2} \right]$$

ONE HAS:

$$\cdot \frac{1}{r^2} \frac{d}{dr} \left(r^2 \frac{B^2}{2\mu_0} \right) = - \frac{dp_{\perp}}{dr} - \frac{p_{\perp H}}{r}$$

WHICH CAN BE INTEGRATED:

$$B = \frac{1}{r} \left\{ r_i^2 B_i^2 - 2r^2 \mu_0 (p_{\perp H} + p_{\perp C}) \right. \\ \left. - 2\mu_0 \int_{r_i}^r r' dr' (p_{\perp H} + 2p_{\perp C}) \right\}^{1/2}$$

OUTGOING ENERGY / EVANESCENT
BOUNDARY CONDITIONS ARE USED
AT INSIDE AND OUTSIDE OF
ANNULUS.

$$\frac{d}{dx} \left(P \frac{d\psi}{dx} \right) - Q\psi = 0$$

OUTSIDE ANNULUS REGION,

$$P, Q \approx \text{constant to } O(\frac{1}{r})$$

∴ solution is matched onto
plane waves $\psi \sim e^{ik_x x}$

$$k_x = \pm i \sqrt{Q/P}$$

± sign determined by:

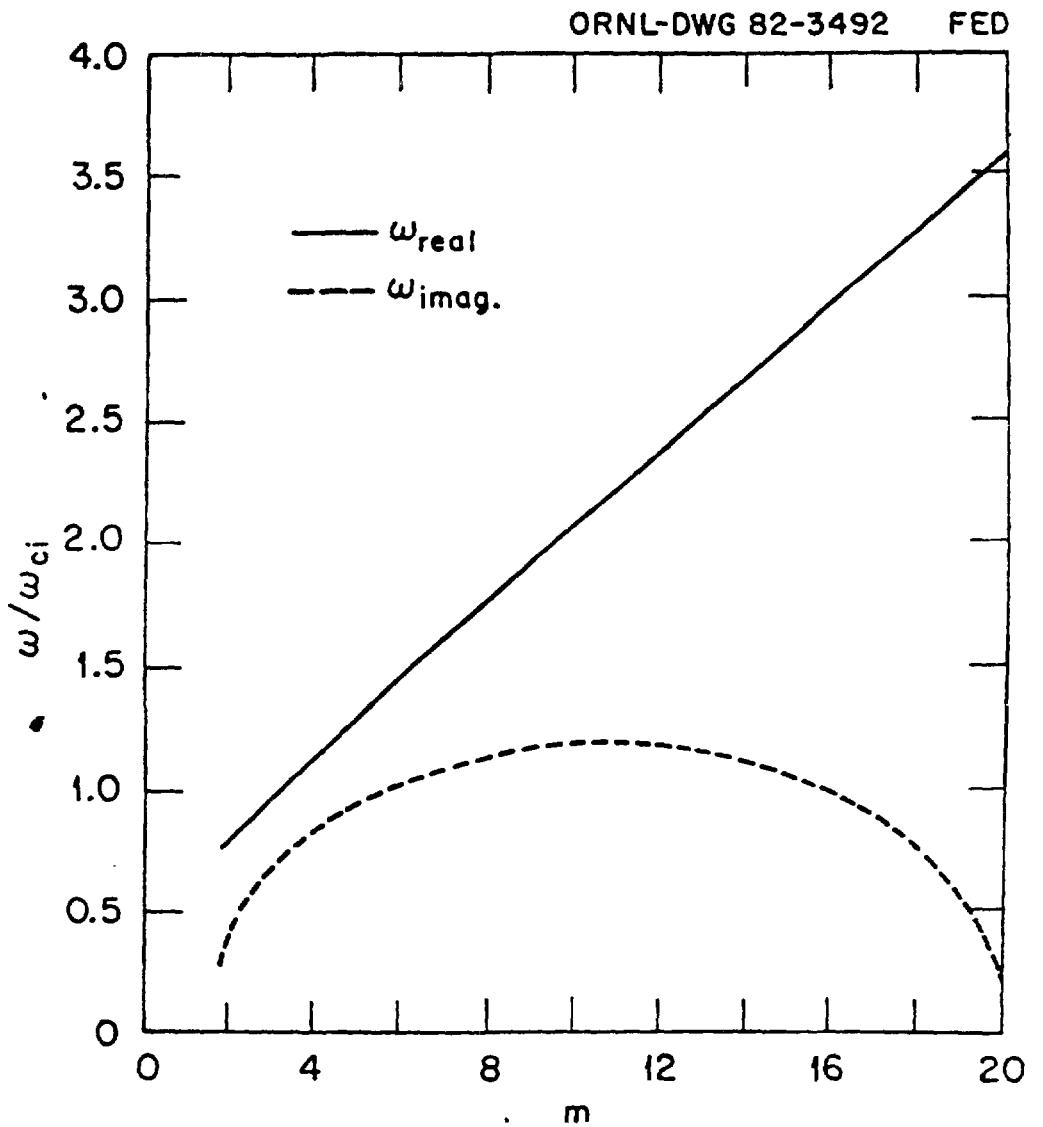
(1) $k_r > k_i$ outgoing energy

$$\frac{\partial w}{\partial k} \geq 0 \quad \text{for } x \geq 0$$

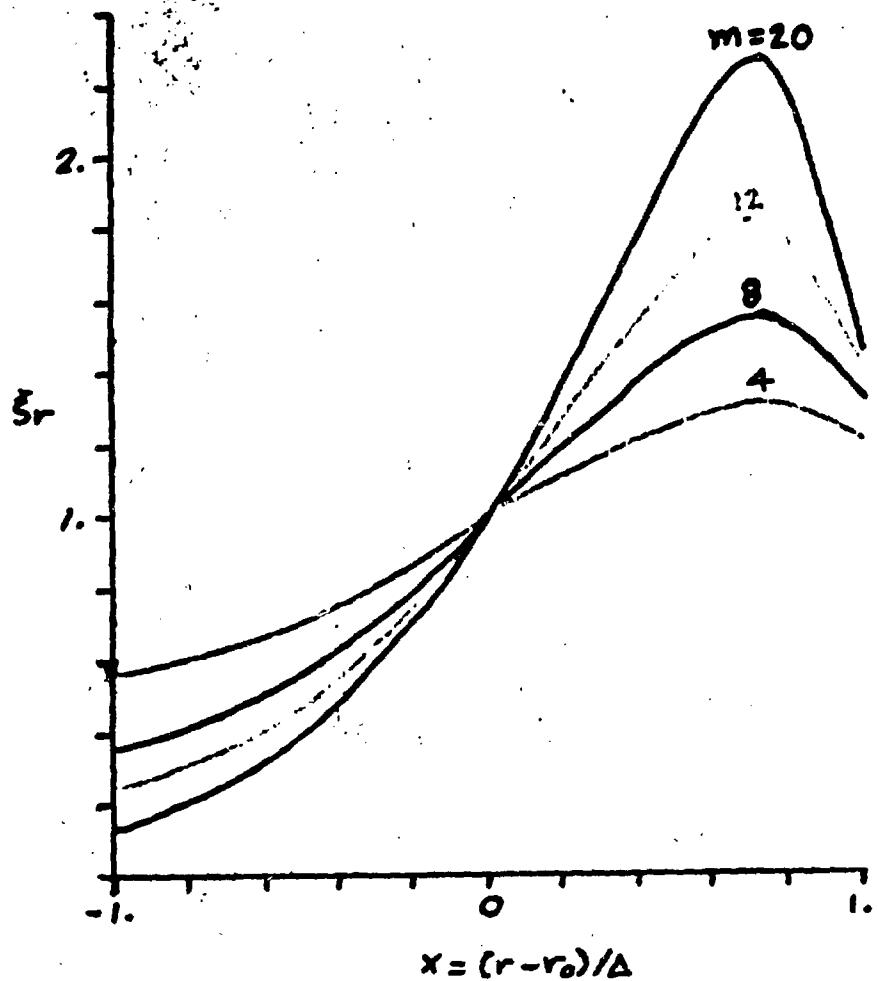
(2) $k_i > k_r$ evanescent

$$k_i \geq 0 \quad \text{for } x \geq 0$$

DEPENDENCE OF FREQUENCY ON AZIMUTHAL MODE NUMBER
FOR HOT ELECTRON INTERCHANGE ($N_{\text{core, elec.}} = 5 \times 10^{11} \text{ cm}^{-3}$,
 $N_{\text{hot}} = 5 \times 10^{11} \text{ cm}^{-3}$, $T_{\text{core}} = 0$, $T_{\text{hot}} = 500 \text{ keV}$)

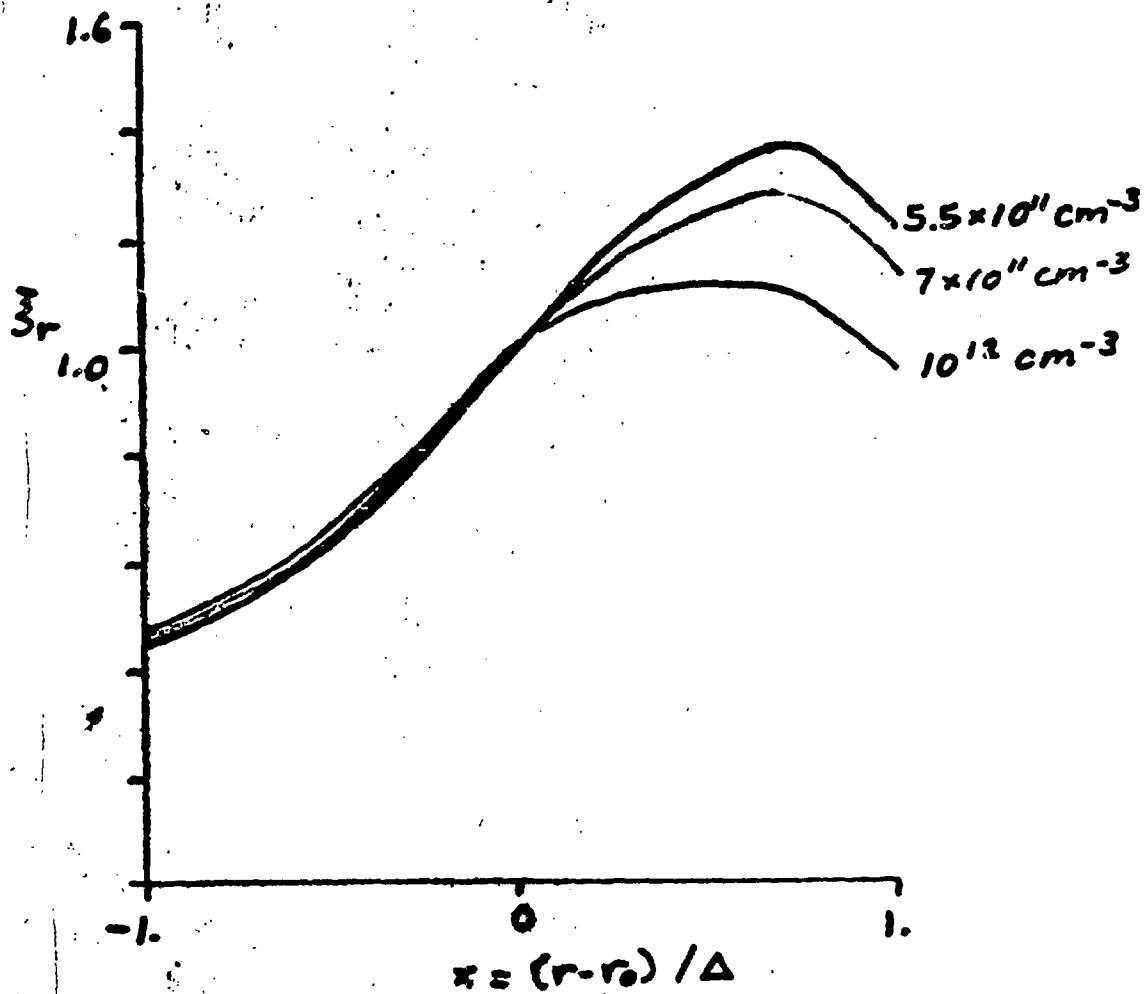


STRUCTURE OF HOT ELECTRON
INTERCHANGE VS. MODE NO.



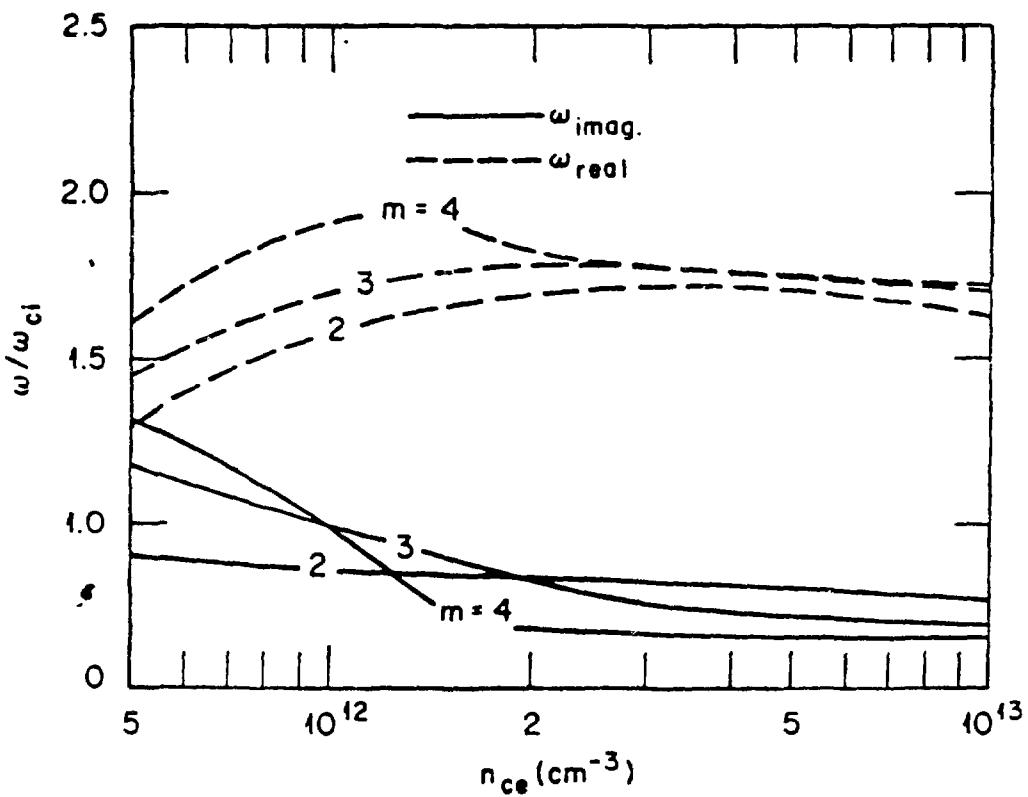
HOT ELECTRON INTERCHANGE MODE
STRUCTURE VS. CORE DENSITY

($m = 6$)

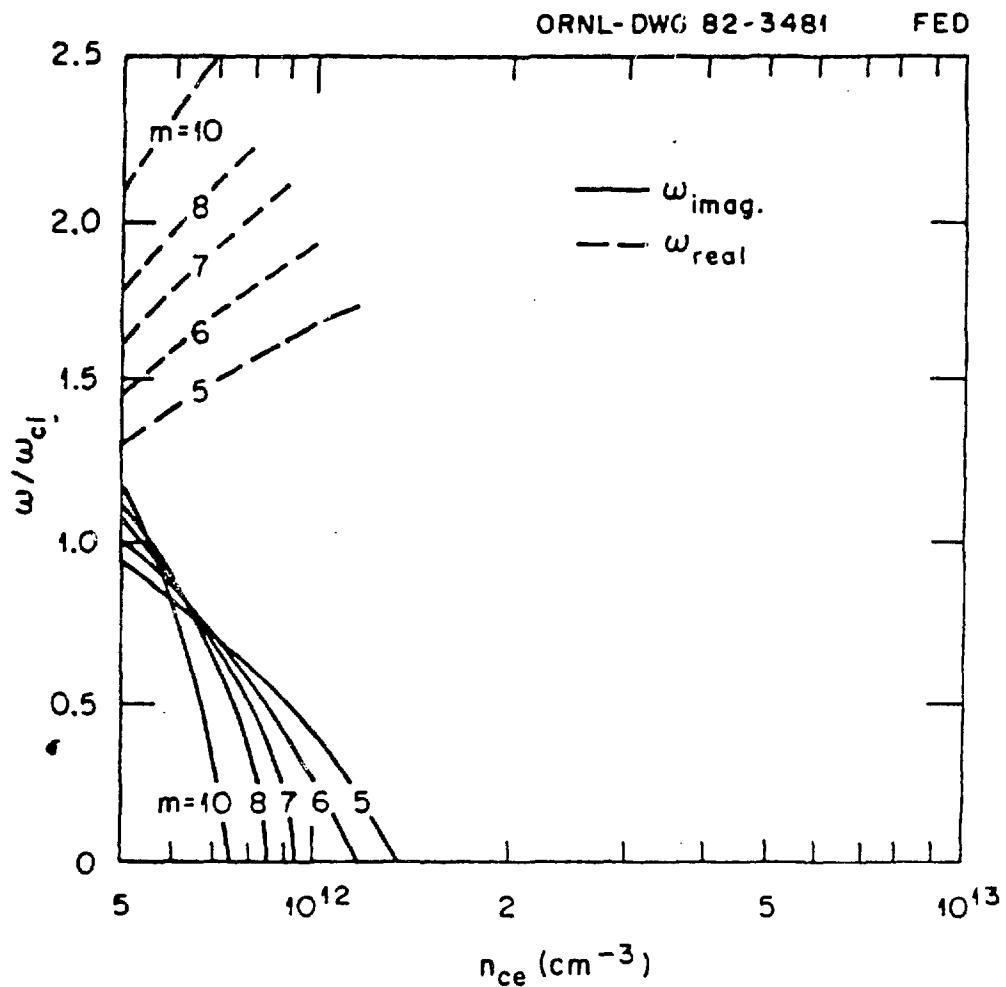


HOT ELECTRON INTERCHANGE
($m = 2-4$)

ORNL-DWG 82-3477 FED

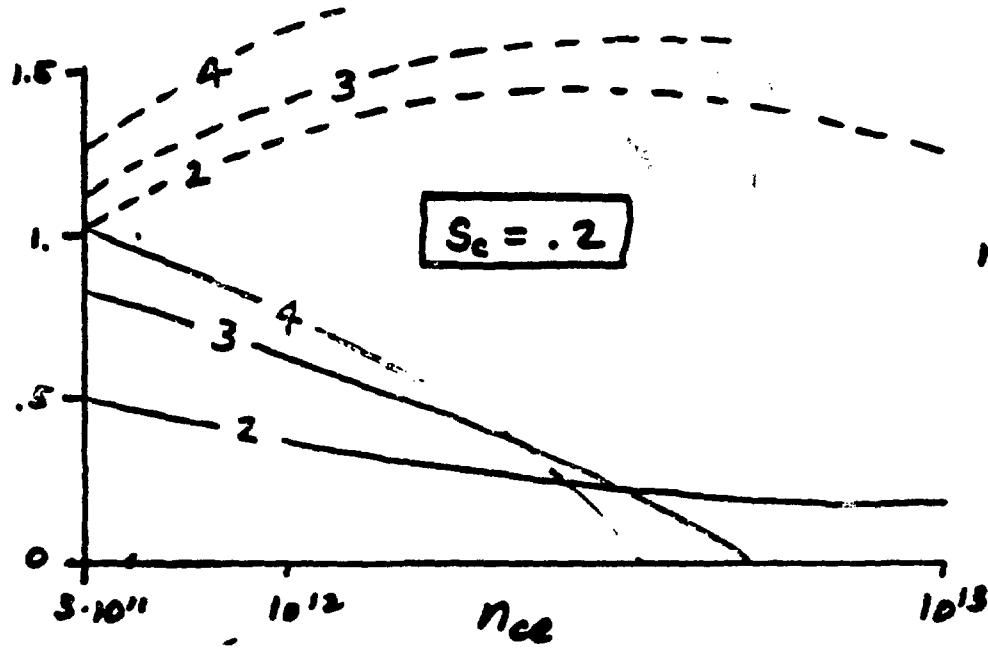
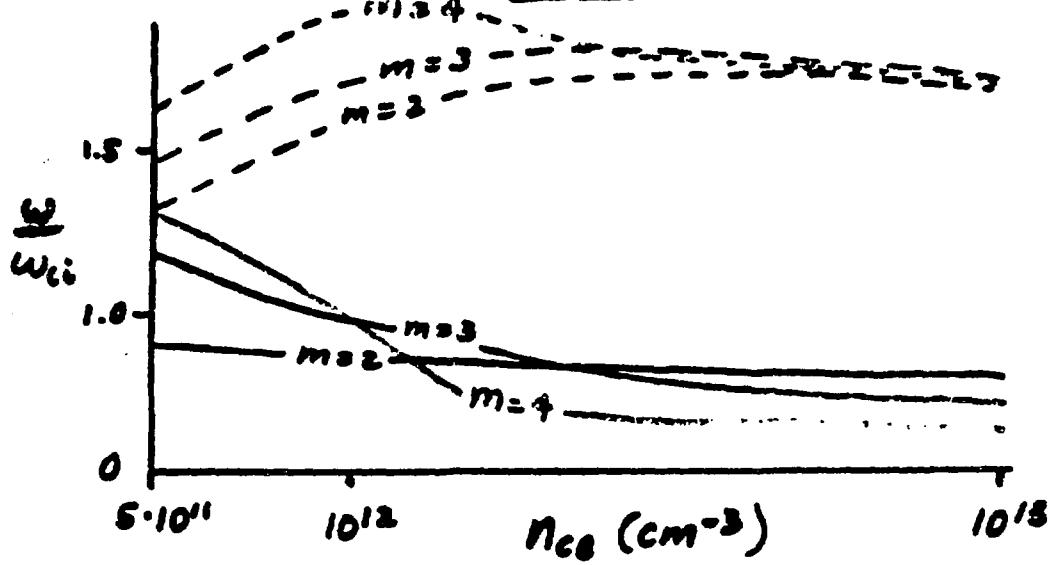


HOT ELECTRON INTERCHANGE
($m = 5-10$)

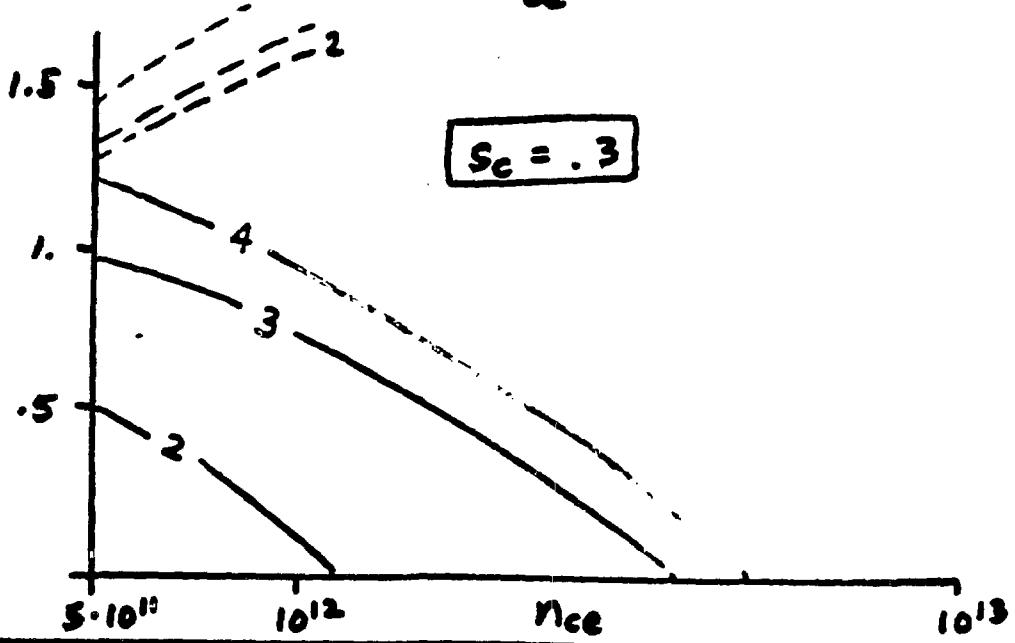


$S_c = .1$

-191-

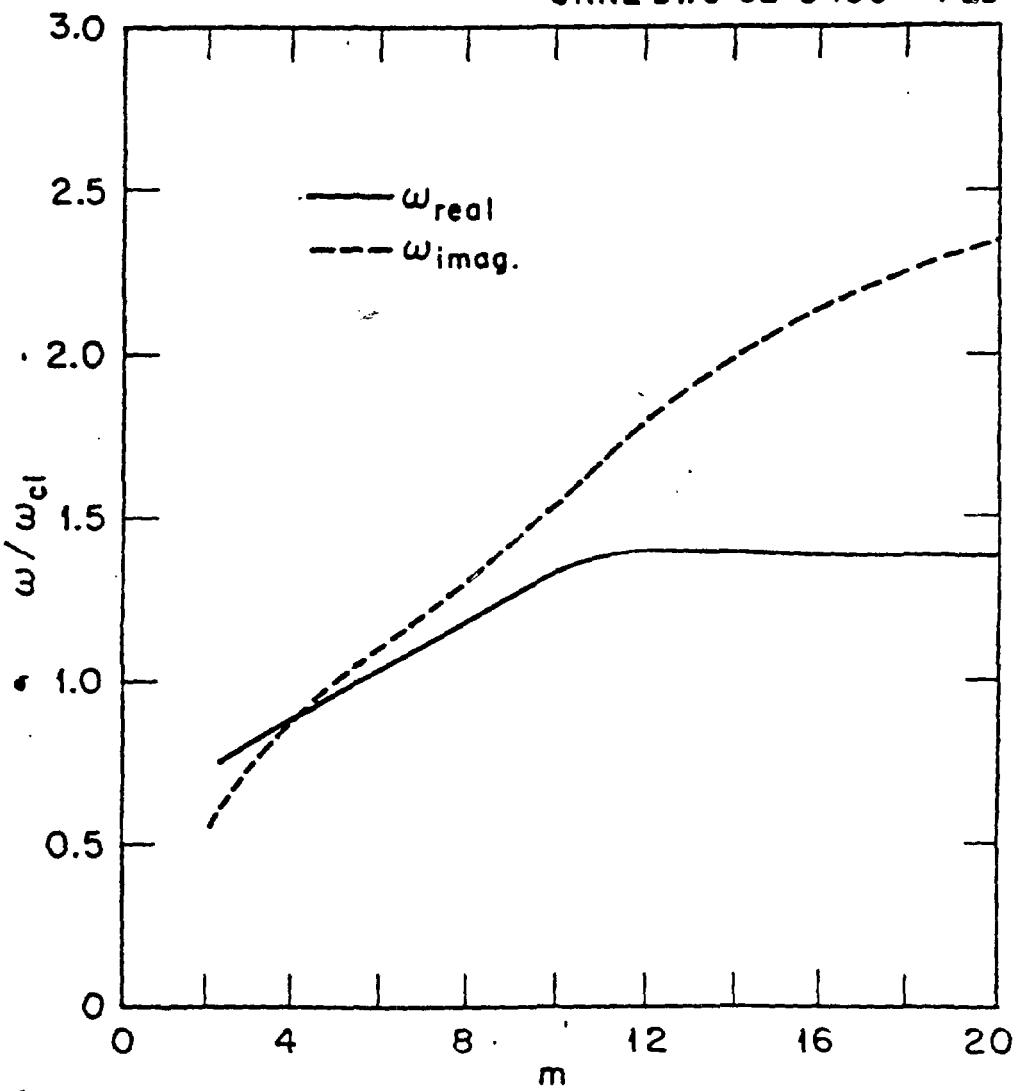


$S_c = .3$

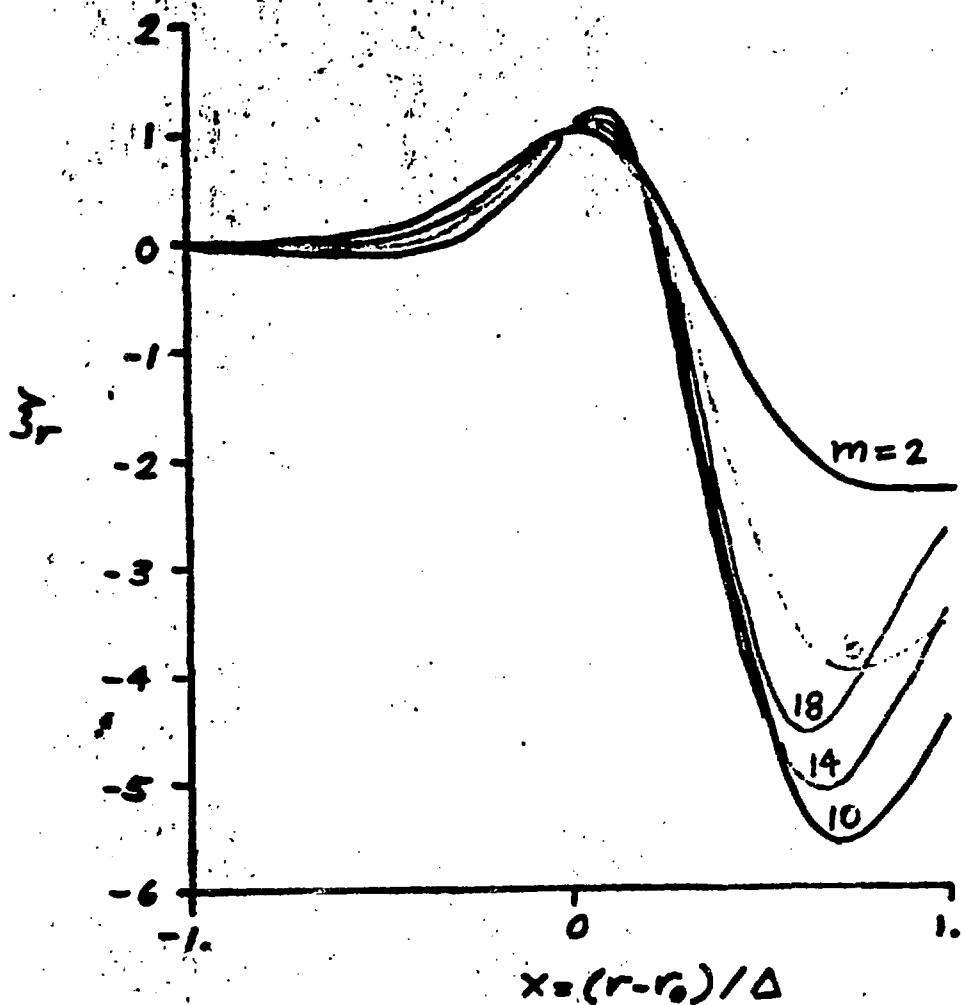


DEPENDENCE OF FREQUENCY ON AZIMUTHAL MODE NUMBER
FOR COMPRESSORIAL ALFVÉN MODE ($N_{\text{core}} = 7 \times 10^{14} \text{ cm}^{-3}$,
 $T_{\text{core}} = 0$, $N_{\text{hot}} = 5 \times 10^{11} \text{ cm}^{-3}$, $T_{\text{hot}} = 500 \text{ keV}$)

ORNL-DWG 82-3490 FED

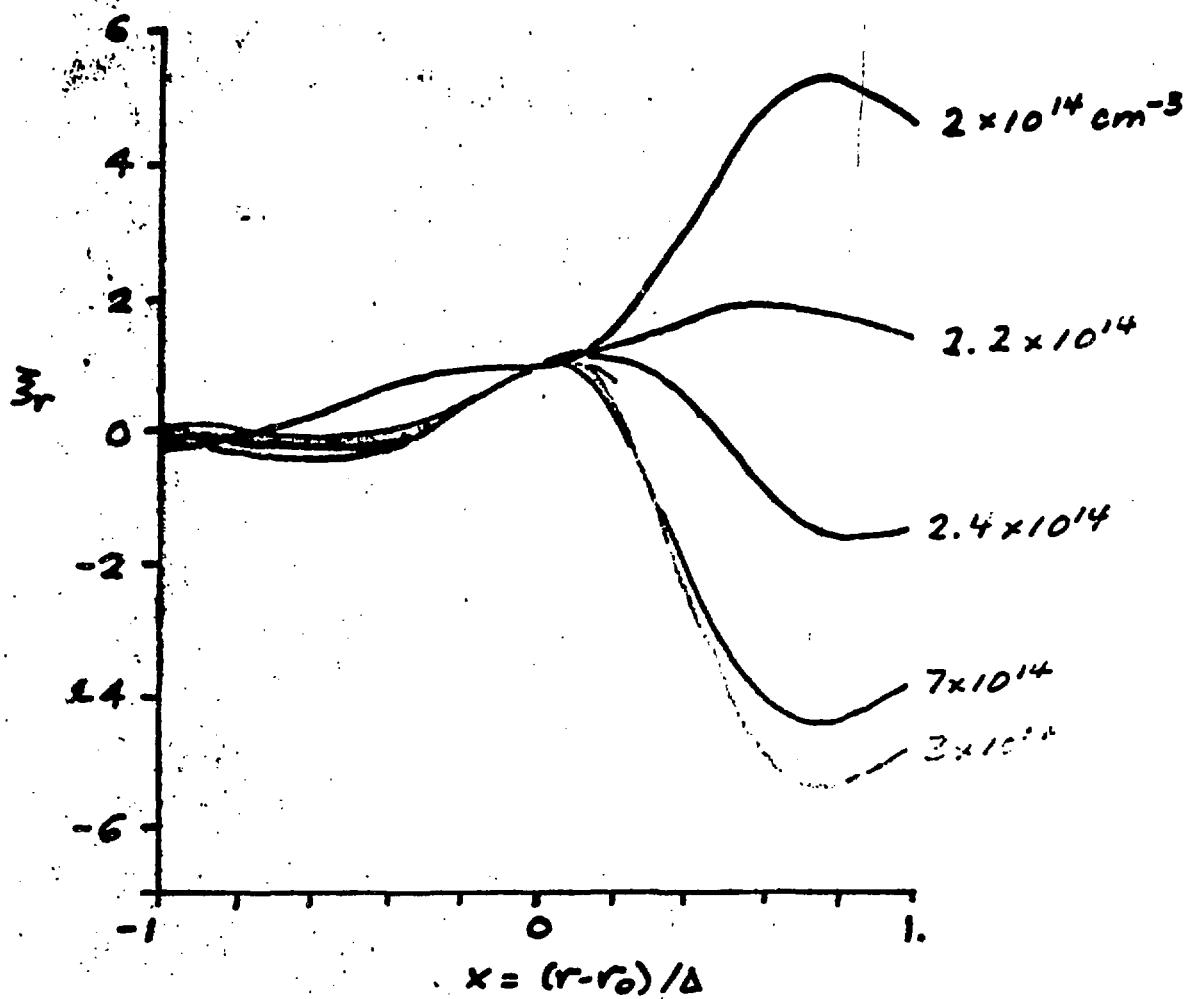


COMPRESSIVE ALFVEN MODE
STRUCTURE VS. MODE NO.



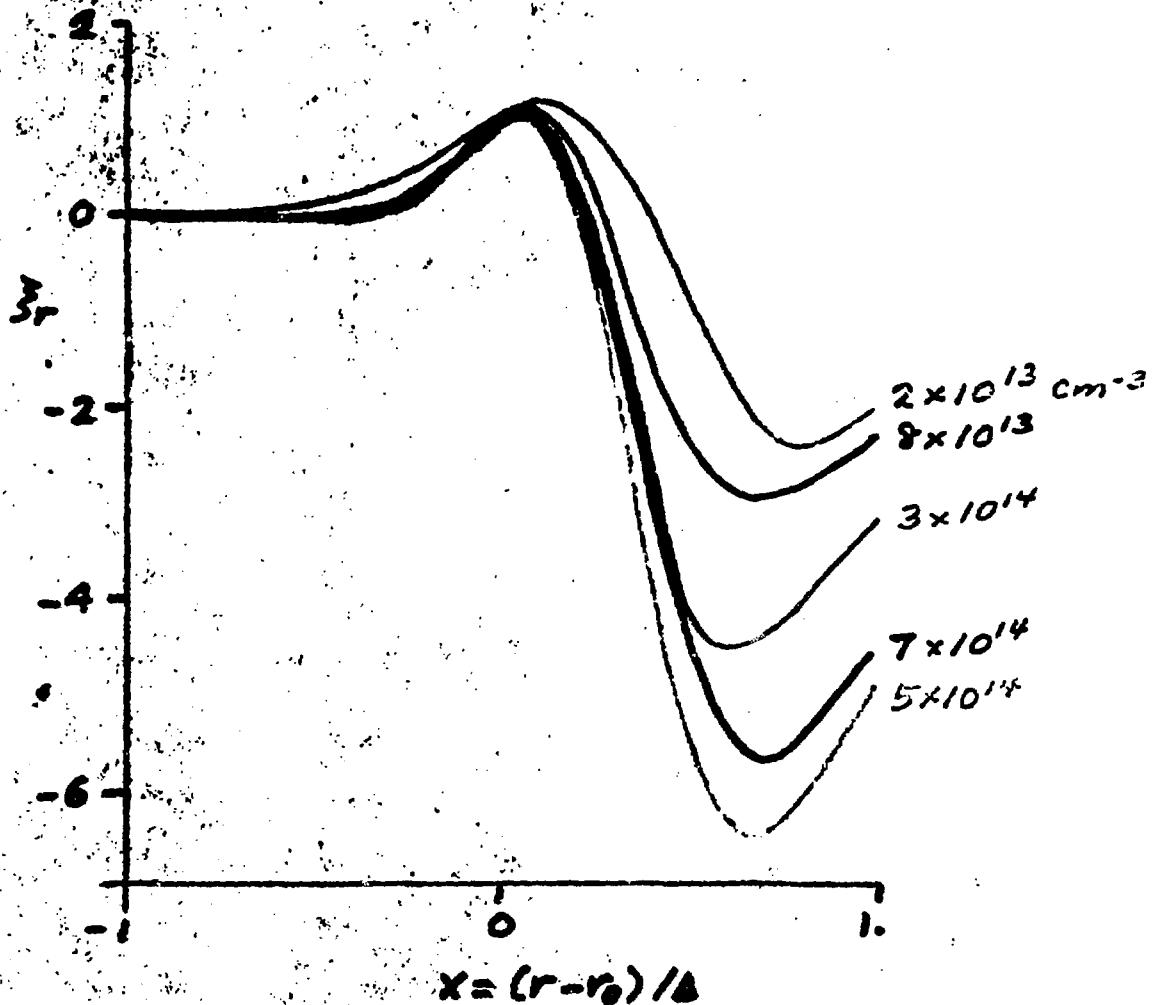
COMPRESSIVE ALFVEN MODE
STRUCTURE VS. CORE DENSITY

($m=7$)



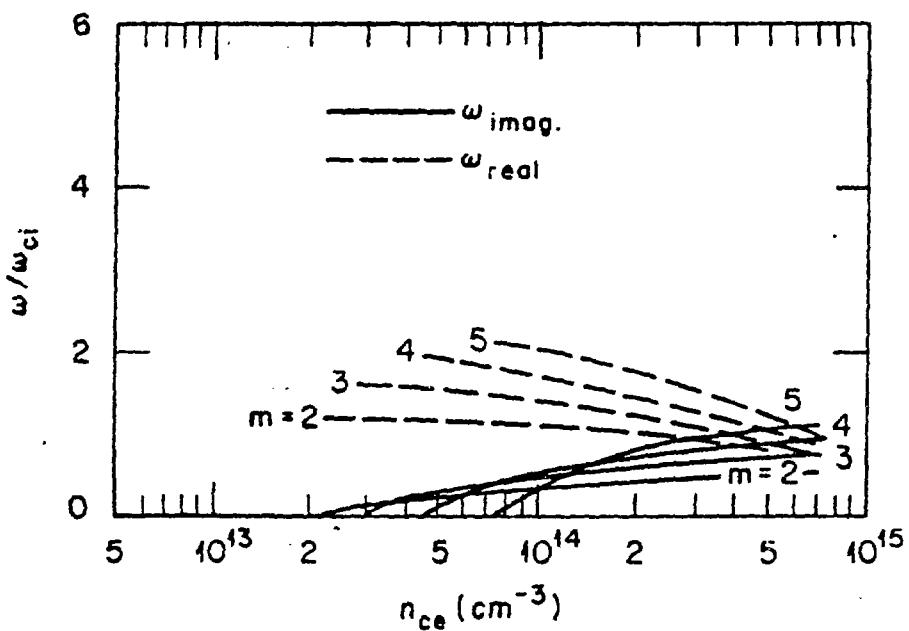
COMPRESSATIONAL ALFVEN MODE
STRUCTURE VS. CORE DENSITY

($m = 10$)



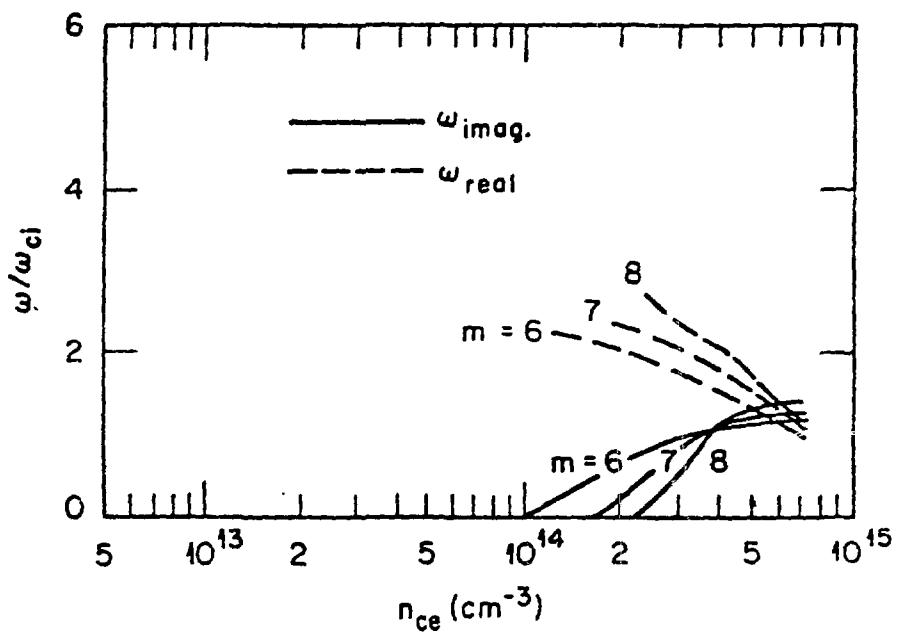
COMPRESSORIAL ALFVÉN WAVE
($m = 2-5$)

ORNL-DWG 82-3478 FED

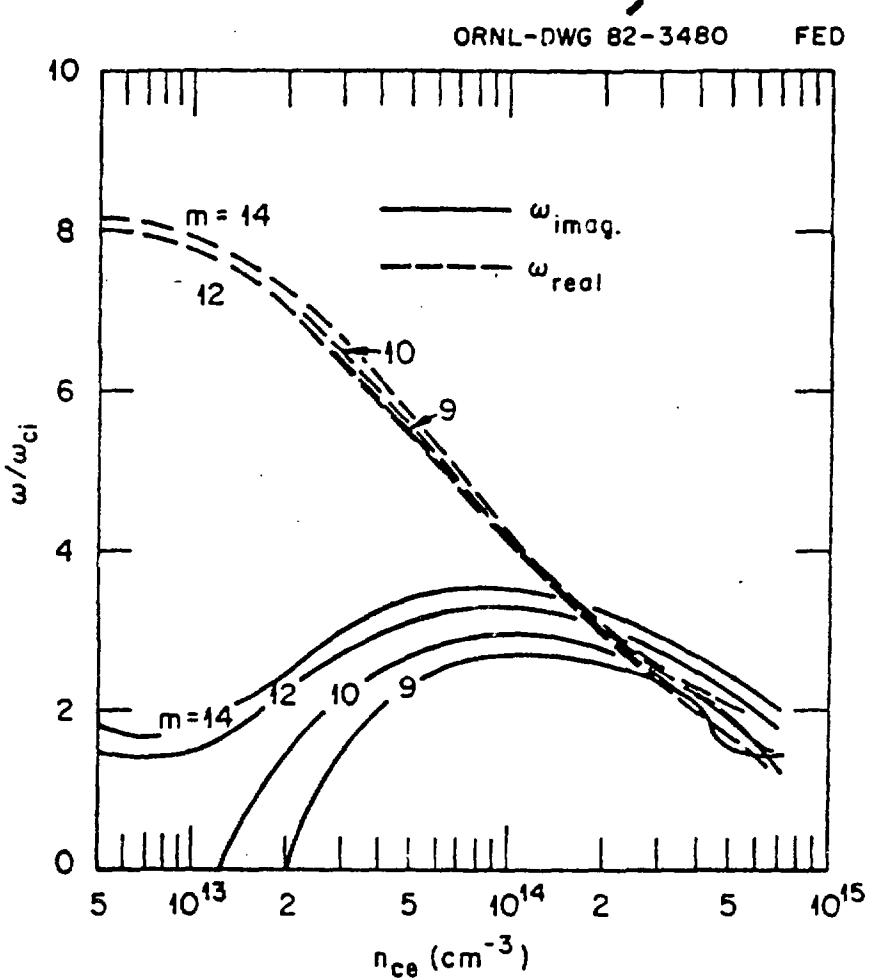


COMPRESSORIAL ALFVÉN WAVE
($m = 6-8$)

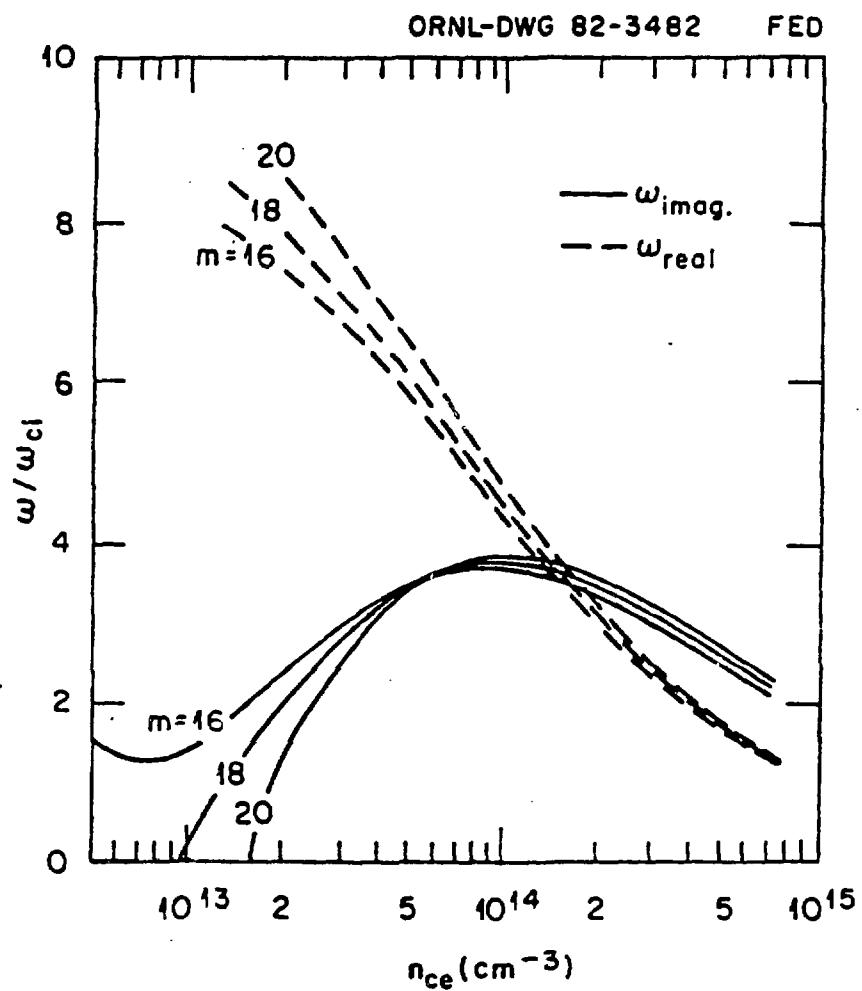
ORNL-DWG 82-3479 FED



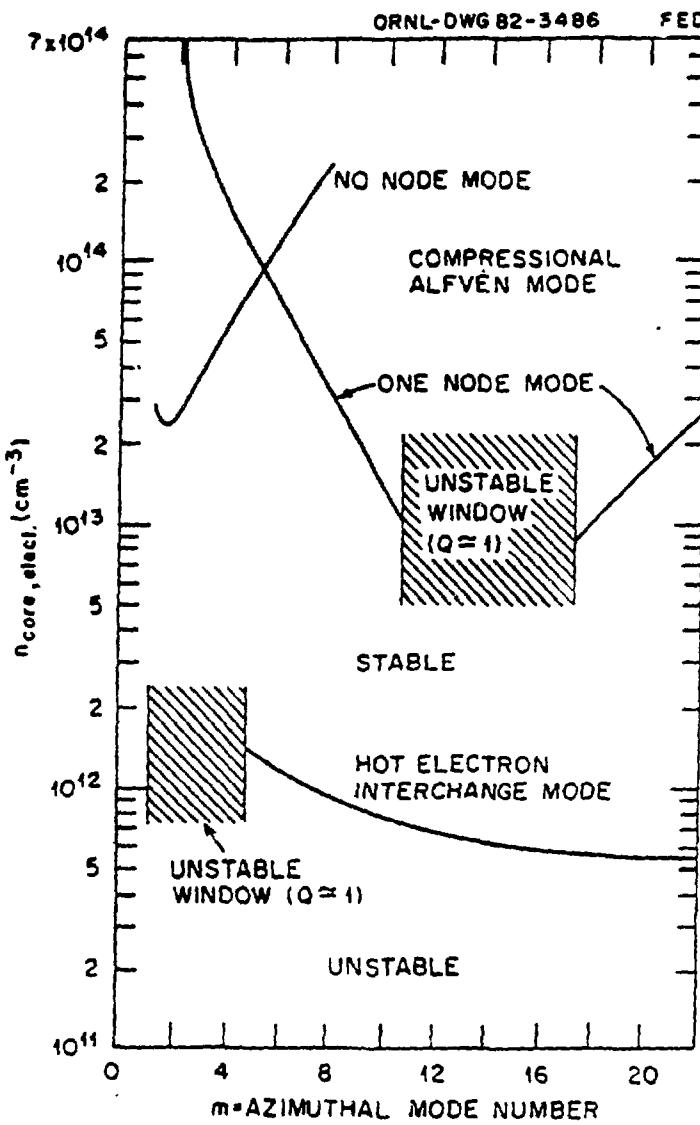
COMPRESSORIAL ALFVÉN WAVE
($m = 9-14$)



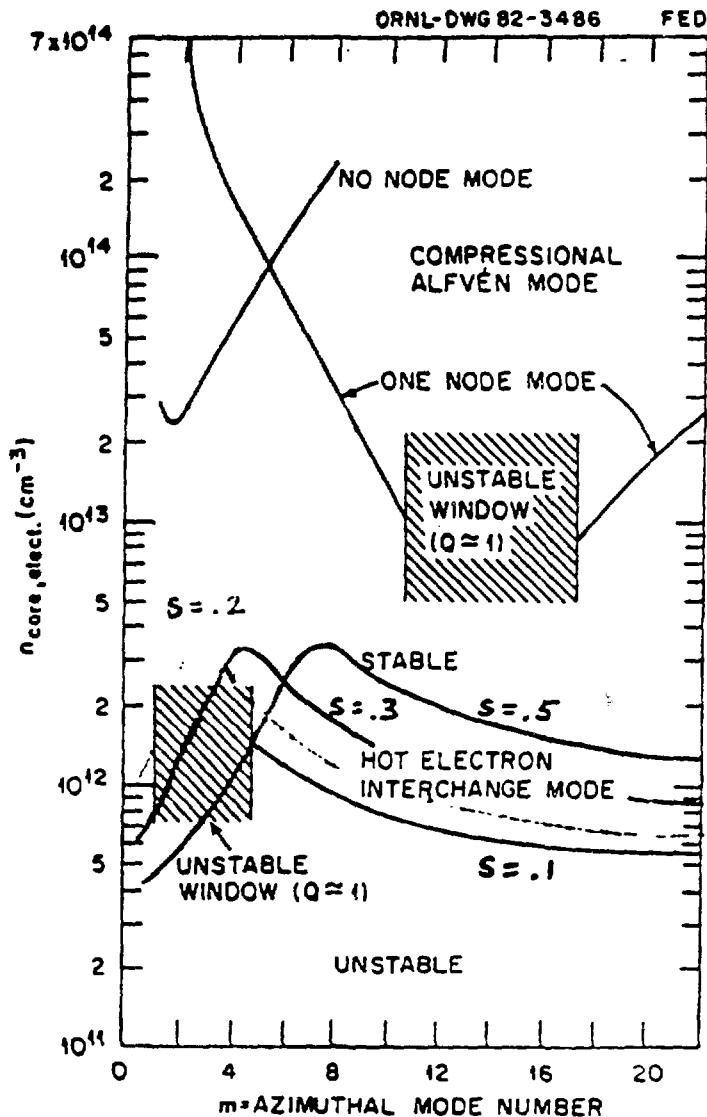
COMPRESSORIAL ALFVÉN WAVE
($m = 16-20$)



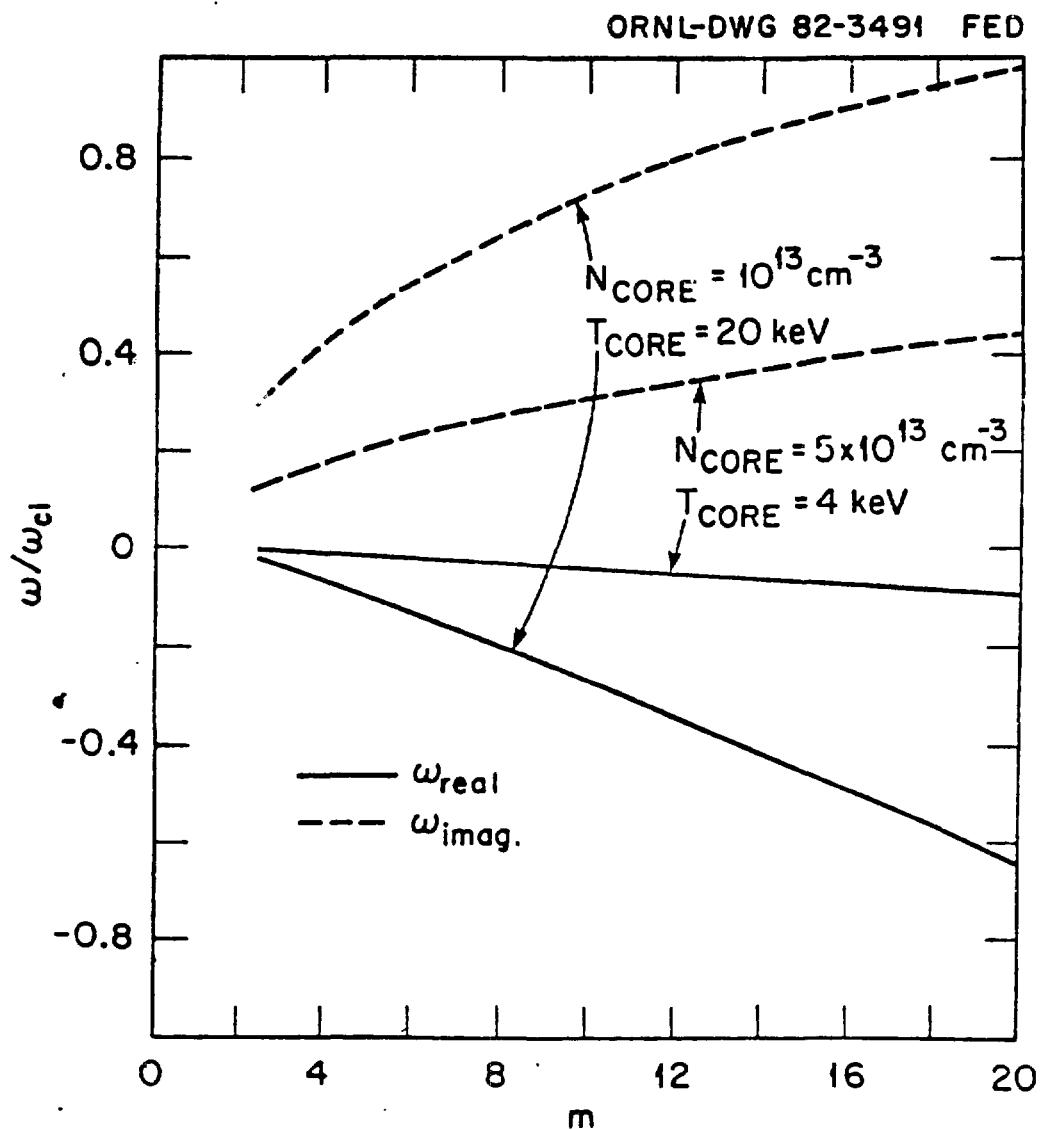
HIGH FREQUENCY MODE STABILITY BOUNDARIES
AS A FUNCTION OF m FOR $T_{\text{core}} = 0$,
 $N_{\text{hot}} = 5 \times 10^{11} \text{ cm}^{-3}$, $T_{\text{hot}} = 500 \text{ keV}$



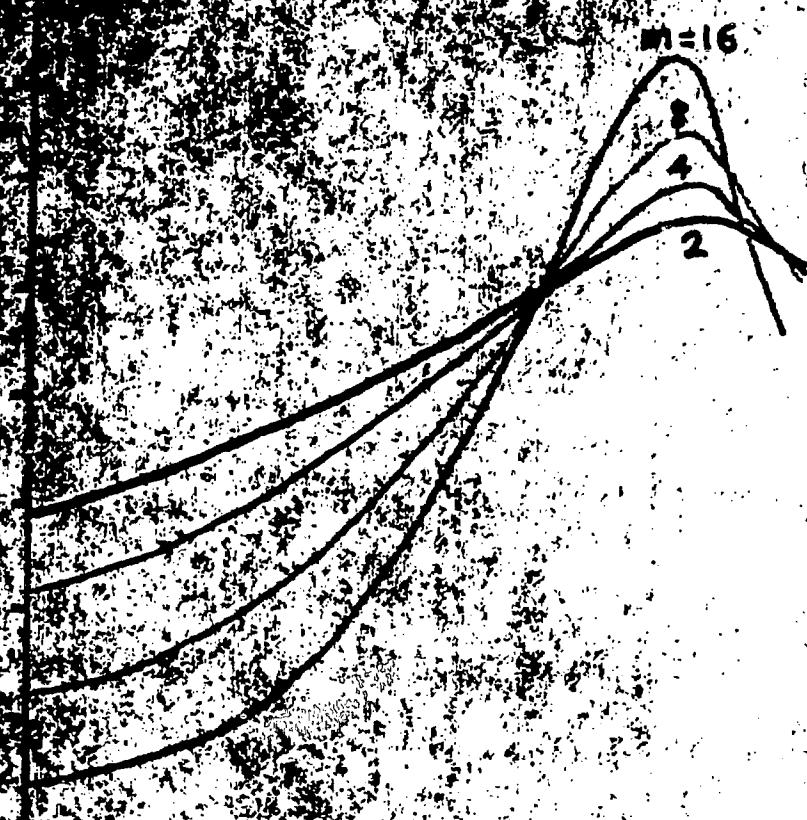
HIGH FREQUENCY MODE STABILITY BOUNDARIES
 AS A FUNCTION OF m FOR $T_{core} = 0$,
 $N_{hot} = 5 \times 10^{11} \text{ cm}^{-3}$, $T_{hot} = 500 \text{ keV}$



DEPENDENCE OF FREQUENCY ON AZIMUTHAL MODE NUMBER
FOR INTERACTING CORE INTERCHANGE
($N_{\text{hot}} = 5 \times 10^{11} \text{ cm}^{-3}$, $T_{\text{hot}} = 500 \text{ keV}$)

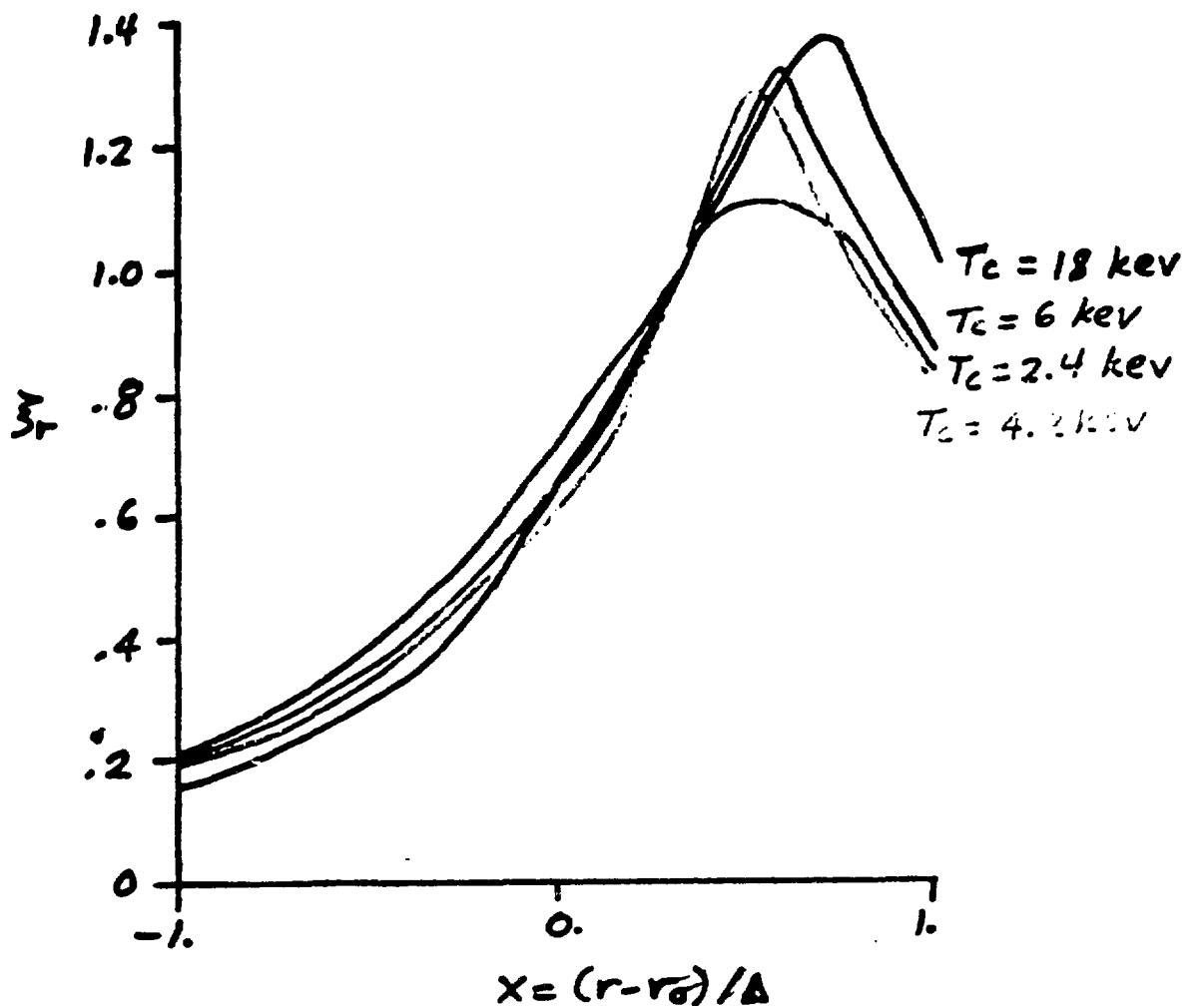


INTRINSIC PRESSURE-DRIVEN INTERCHANGE
STRUCTURE VS. M



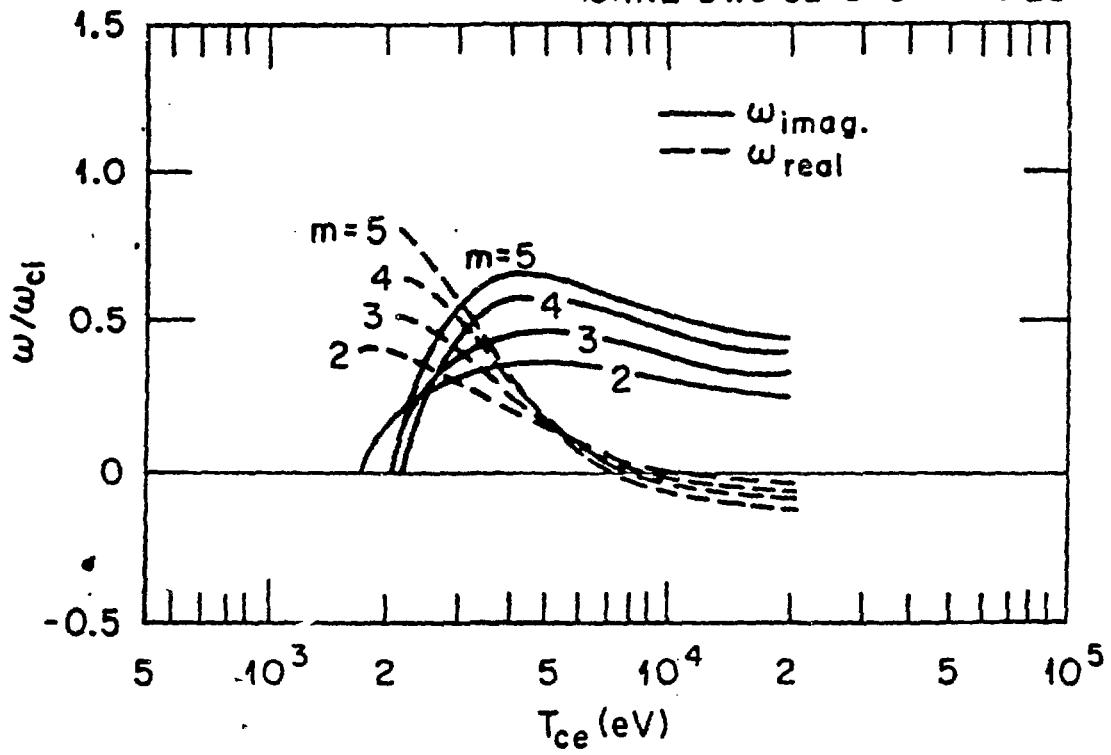
$$x = (r - r_0) / \Delta$$

INTERACTING PRESSURE -DRIVEN INTERCHANGE
MODE STRUCTURE VS. CORE TEMPERATURE
($m=11$)



INTERACTING CORE INTERCHANGE MODE
 $(N_{\text{core}} = 10^{13} \text{ cm}^{-3}, m = 2-5)$

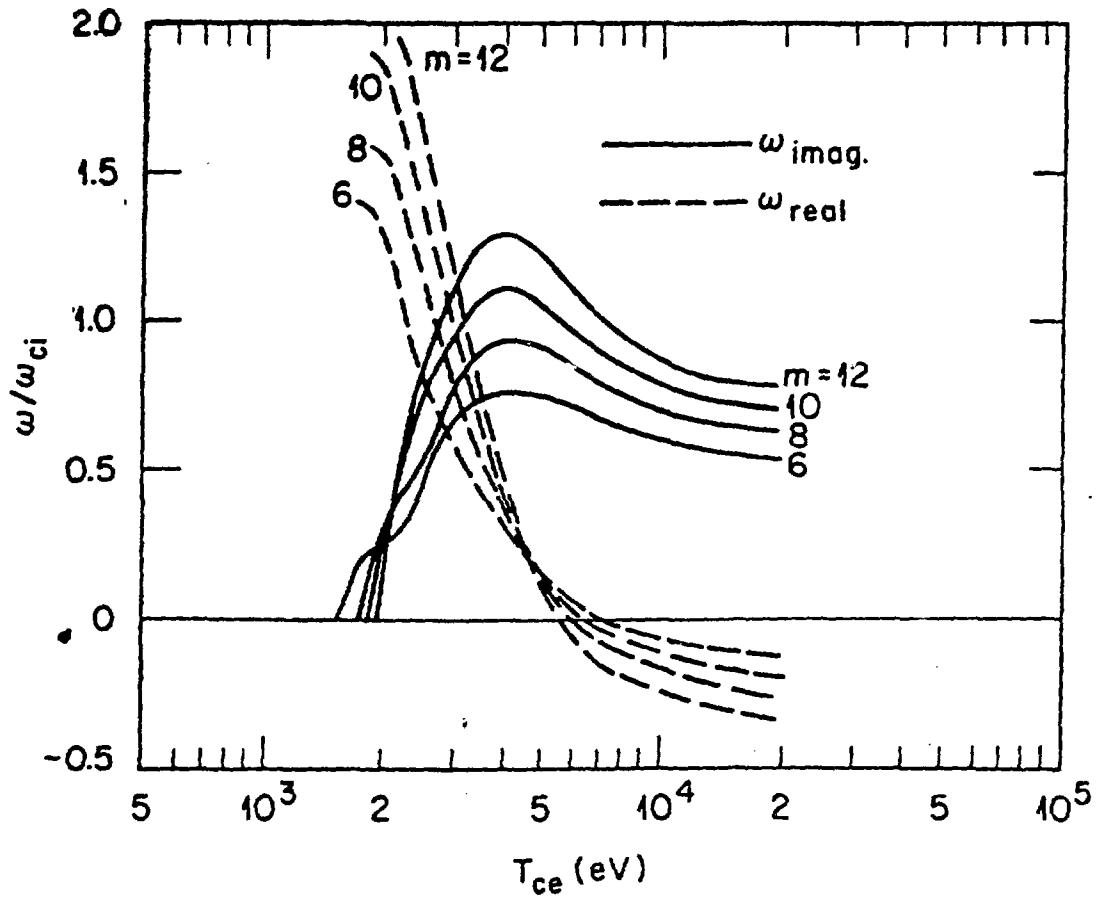
ORNL-DWG 82-3487 FED



INTERACTING CORE INTERCHANGE MODE
($N_{\text{core}} = 10^{13} \text{ cm}^{-3}$, $m = 6-12$)

ORNL-DWG 82-3489

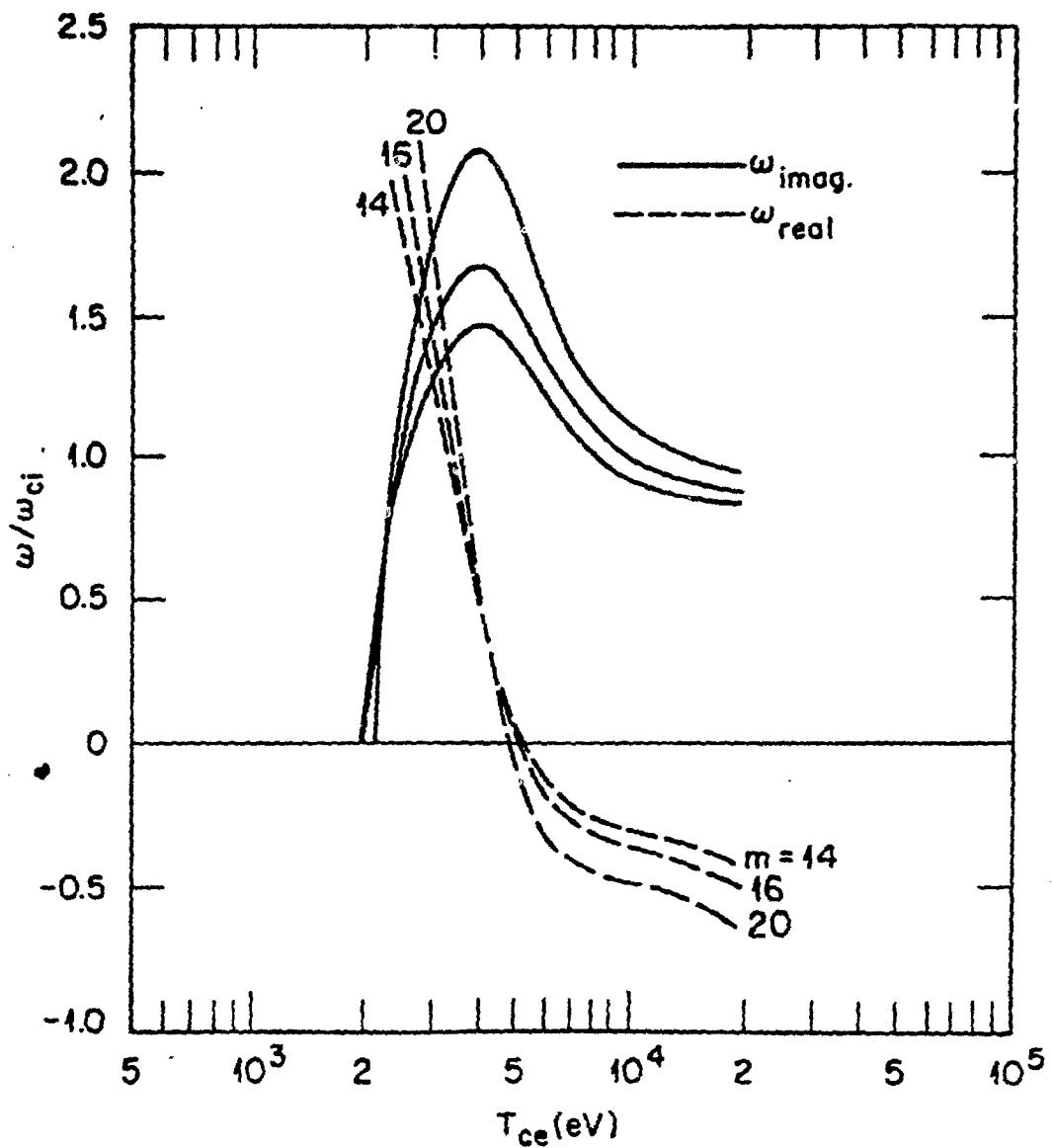
FED



INTERACTING CORE INTERCHANGE MODE
($N_{\text{core}} = 10^{13} \text{ cm}^{-3}$, $m = 14-20$)

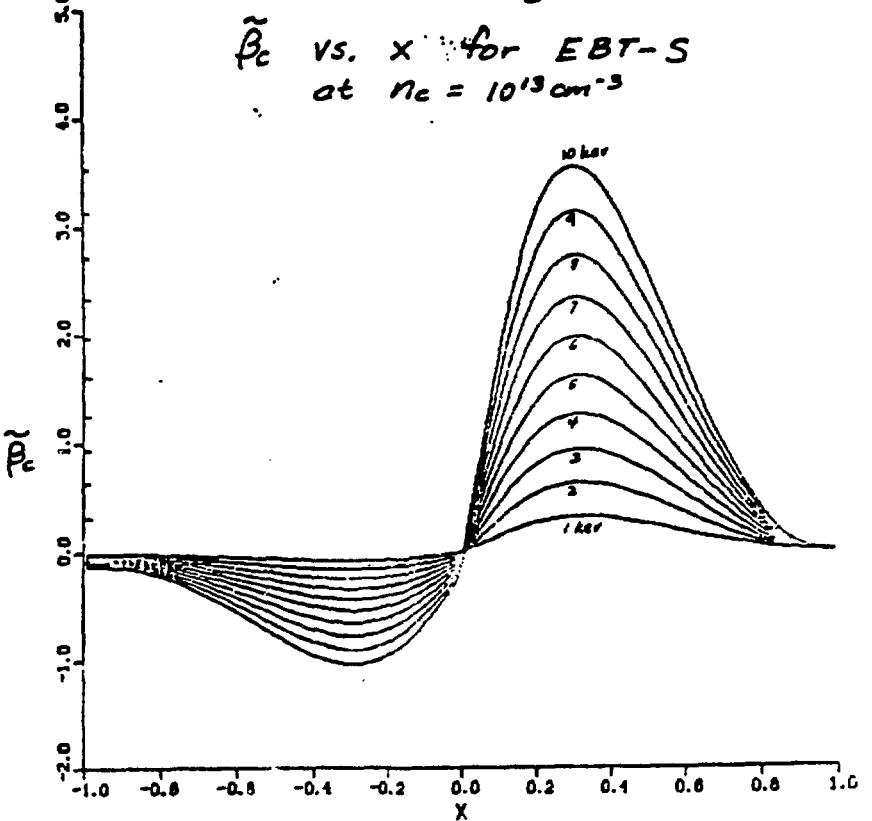
ORNL-DWG 82-3488

FED



TYPICAL $\tilde{\beta}_c$ PROFILES USED
IN RADIALLY - DEPENDENT
CALCULATIONS

$$\tilde{\beta}_c = -r \frac{d}{dr} \left(\frac{\beta_c}{B^2} \right) \approx \beta_c \left(\frac{r}{2\Delta} \right) (1 + \beta_H)$$



HOT PLASMA DECOUPLING CONDITIONS:

- WKB ANALYSIS:

$$1 - \tilde{\beta}_e \leq 3 \left[\frac{\pi}{\delta} \tilde{A} \left(1 + \frac{2 \tilde{\beta}_e}{\tilde{\beta}_H} - \frac{2}{\tilde{\beta}_H} \right) \right]^{1/3}$$

defining $\Omega_{cv} = \frac{\omega_{cv}}{\delta m \omega_0} = \sqrt{\frac{k \Delta g_0}{P}}$

$$\Omega_{cv} < \frac{3 (k \delta)^{1/4} \tilde{A}^{3/2} \sqrt{1 + 2 \tilde{\beta}_e / \tilde{\beta}_H - 2 / \tilde{\beta}_H}}{(k \delta) (1 - \tilde{\beta}_e)^{3/2}}$$

- LAYER ANALYSIS:

$$\tilde{\beta}_e \ll 1$$

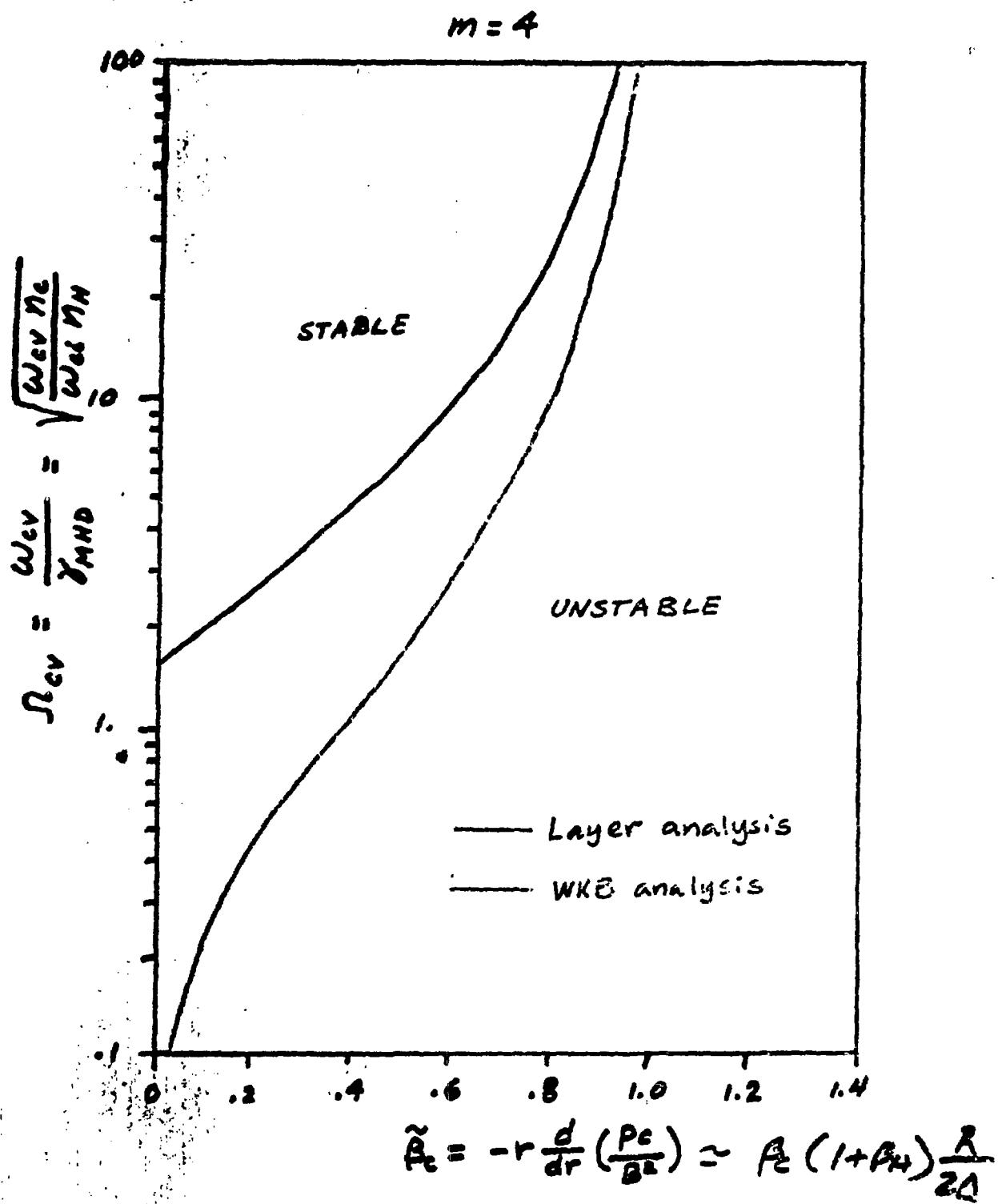
$$\Omega_{cv} < \text{Max} \left[2 \left(\frac{k}{R} \delta \right)^{1/4}, 4 \left(\frac{k}{R} \tilde{\beta}_e \right)^{1/2} \right]$$

$$\delta \approx |k| \Delta \quad K = \frac{1}{\delta(R-\Delta)} \frac{d\delta(R-\Delta)}{dr} \sim |k|$$

$$\tilde{\beta}_e \lesssim 1$$

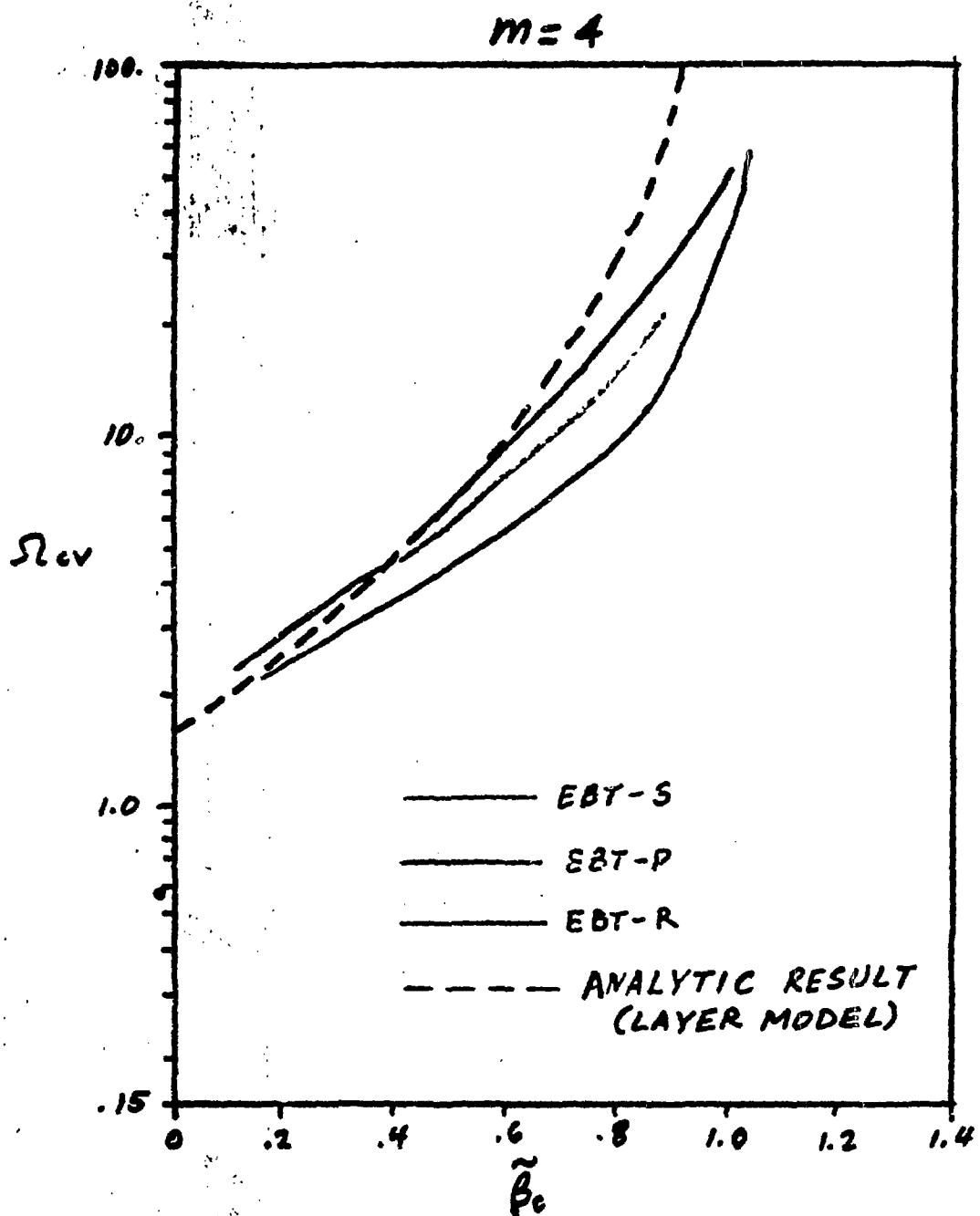
$$\Omega_{cv} > \frac{\delta^{3/4} |k|}{2 K^{1/4} (1 - \tilde{\beta}_e)^{1/2}}$$

HOT PLASMA DECOUPLING CONDITION FOR PRESSURE-DRIVEN INTERCHANGE

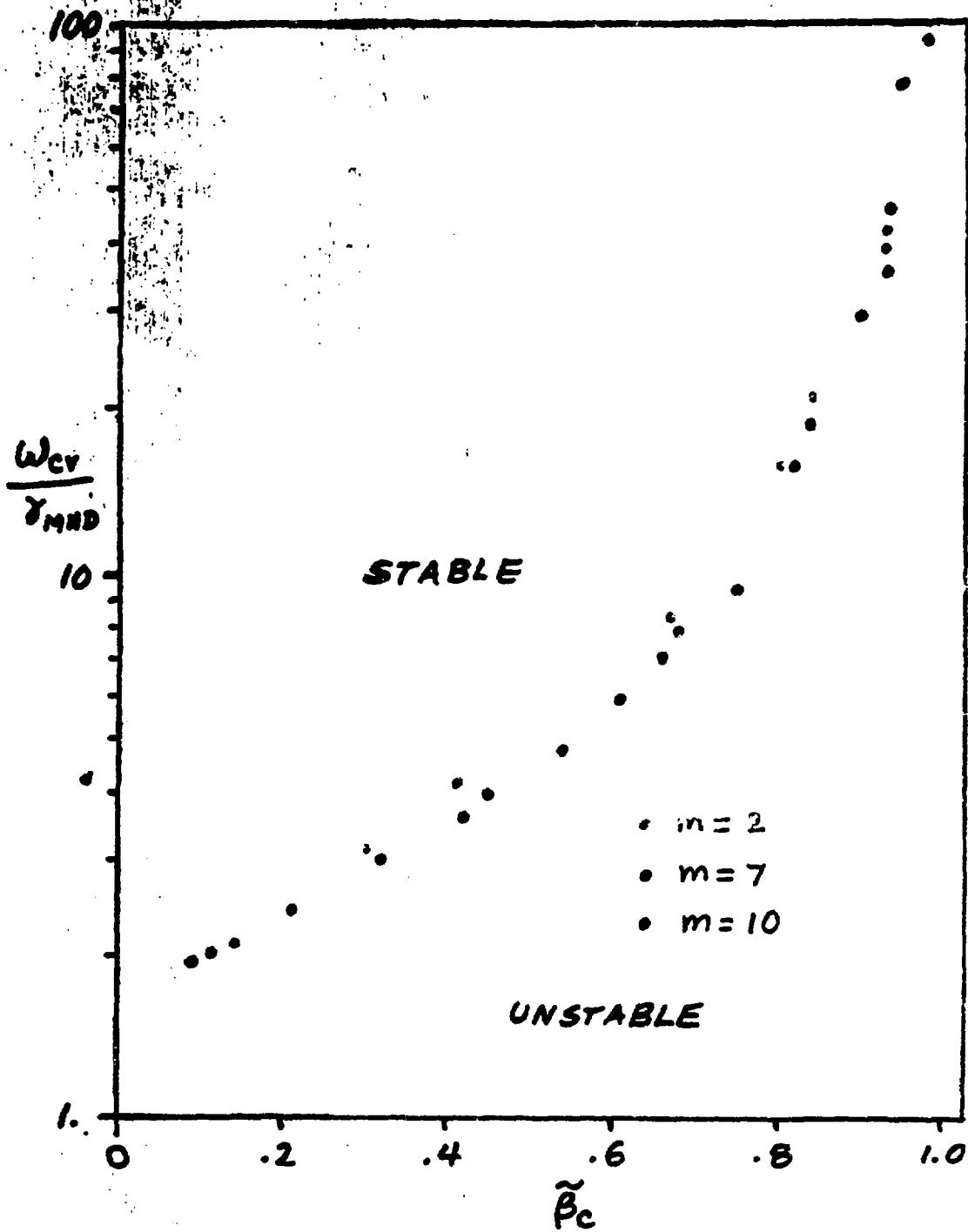


$$\tilde{\beta}_c = -r \frac{d}{dr} \left(\frac{P_c}{B^2} \right) \approx \beta_c (1 + \beta_H) \frac{R}{2Q}$$

DECOUPLING CONDITION FROM NONLOCAL CALCULATION

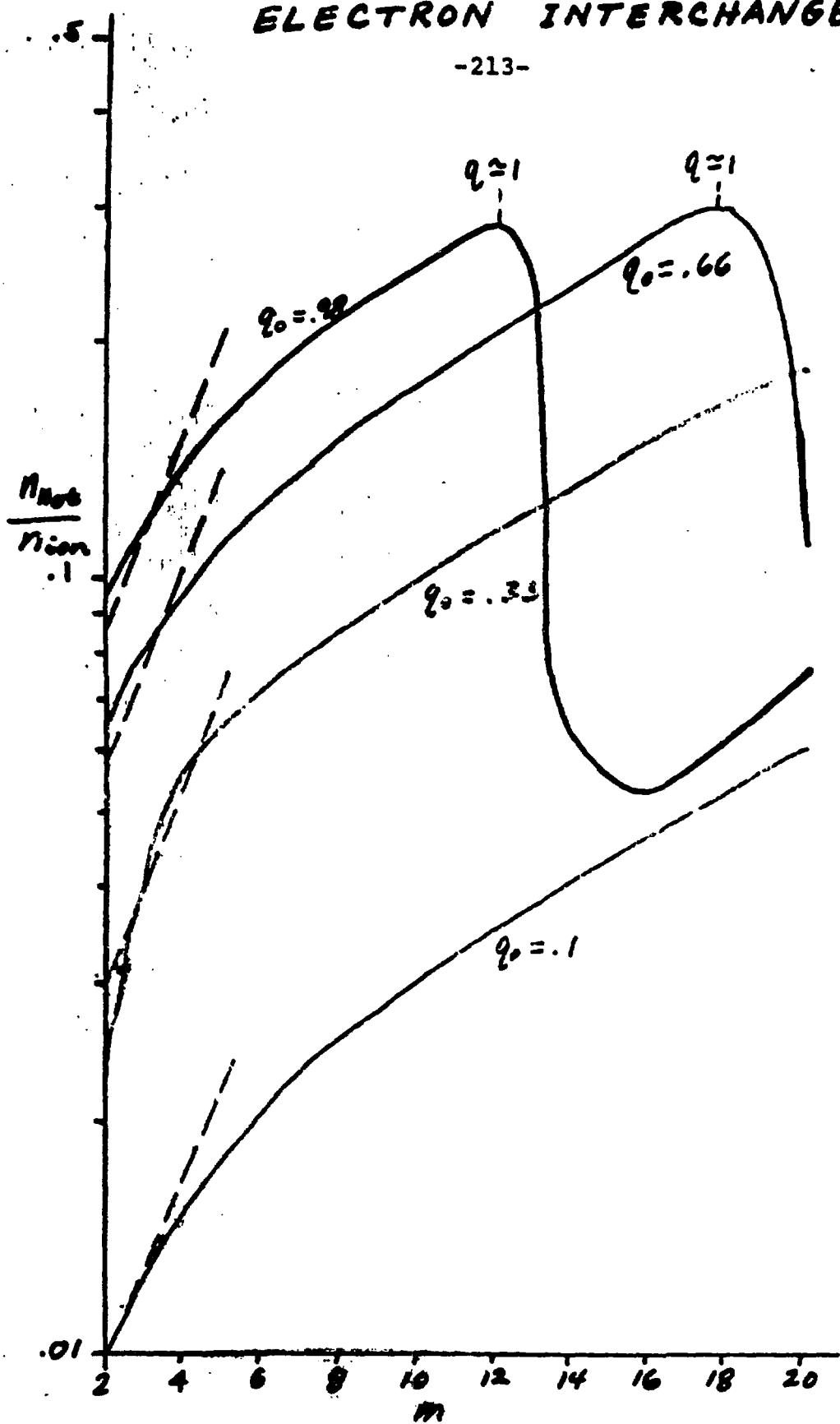


THE HOT PLASMA DECOUPLING
CONDITION IS NOT STRONGLY
DEPENDENT ON M (except through $\frac{\omega_c}{\delta_{MHD}}$)



ELECTRON INTERCHANGE

-213-



CONCLUSIONS:

• HOT ELECTRON INTERCHANGE

$m \geq 6$ somewhat more optimistic
than local theory due to $k_r \delta < 2$

$m \leq 6$ instability band present ($q < 1$)
can be moved to $m < 1$
by profile change (surface plasma)

frequency and $(n_H/n_c)_{\text{crit}}$ in
reasonable agreement with expt.

• COMPRESSATIONAL ALFVEN

stability { no node mode for $m \leq 9$
determined by { one node mode for $m \geq 9$

$(k_r)_{\text{eff.}}$ higher than local theory
→ larger density threshold

decoupling condition at different
 m than local theory due to
radial mode structure

one node unstable band at ~ 80 MHz
seen in expt. throughout T mode

CONCLUSIONS: (cont'd.)

- LOW FREQUENCY HOT ELECTRON INTERCHANGE

SIMILAR TO HIGH FREQUENCY BRANCH,
BUT WITH LOWER q_0

LAYER SCALING FITS LOW m RESULTS.

- INTERACTING RING-CORE INTERCHANGE

RELEVANT PARAMETER IS: $\tilde{\beta}_c = -r \frac{d}{dr} \left(\frac{P_c}{B^2} \right)_{MAX}$

EBT-S , EBT-P $\tilde{\beta}_c \lesssim .2$
(20 kev) (1-2 kev)

RING-CORE DECOUPLING LOST AT:

$$\frac{\omega_{cv}}{\delta_{MHD}} \simeq 1.5$$