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USER'S MANUAL FOR LINEAR, A COMPUTER  
PROGRAM THAT CALCULATES THE LINEAR  
CHARACTERISTICS OF A GYROTRON

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## I. GENERAL DESCRIPTION

### A. Purpose

Calculates the linear characteristics of a gyrotron. This program is capable of:

- 1) calculating the starting current or frequency detuning for each gyrotron mode.
- 2) generating mode spectra
- 3) plotting these linear characteristics as a function of device parameters (e.g., beam voltage)
- 4) doing the above for any axial RF field profile

### B. Compatibility

ANSI Standard FORTRAN IV, 1966, with some features of FORTRAN 77.

This program was written using the MULTICS operating system on the Honeywell 68/DPS computer.

### C. Limitations

- 1) The electron beam parameters are calculated assuming that adiabatic theory is valid in the gun region [1].
- 2) The electron beam is assumed to be annular, azimuthally symmetric, and to have no radial thickness. It is also monoenergetic and has no velocity spread.
- 3) The cavity is assumed to be a cylindrical open resonator.

### D. Method

The starting current of a gyrotron [2] can be written as:

$$I_{ST} = - \frac{\epsilon_0}{2} \frac{\omega}{Q_T} \left| p_0 \right|^2 \frac{m_e}{e} \frac{(K_{II} V_{II})^2}{G(r_e)} \left[ J'_n(K_L r_L) \right]^{-2} \left[ F_c - \frac{1}{2} \frac{S V_1^2}{c^2} \frac{dF_c}{dx} \right]^{-1} \quad (1)$$

where  $Q_T$  is the total Q of the cavity,  $\omega$  is the resonator frequency,  $m_e$  is the relativistic electron mass,  $r_L$  is the electron Larmor radius,  $S = \omega/K_{II} V_{II}$ , and  $V_{II}$  and  $V_L$  are the electron beam's parallel and perpendicular velocity respectively. The parameter  $|p_0|^2$  represents the stored energy in the cavity and contains the factor:

$$E_S = \frac{1}{K_{II} L} \int_{Z_{IN}}^{Z_{OUT}} d\bar{Z} \left[ g(\bar{Z}) \right]^2 \quad (2)$$

where  $g(\bar{Z})$  = axial RF field profile

$$\bar{Z} = K_{II} Z$$

$Z_{IN}$ ,  $Z_{OUT}$  = limits of the beam-field interaction

$K_{II} = qK/L$  = axial wavenumber (K is a constant)

The variable  $G(r_e)$  is the beam-field coupling term, and can be written as follows depending on the structure of the field in the azimuthal direction:

$$(Standing) \quad G(r_e) = 0.25 \left[ J_{m+n}^2 (K_L r_e) + J_{m-n}^2 (K_L r_e) \right] \quad (3)$$

$$(Rotating) \quad G(r_e) = 0.5 \left[ J_{m+n}^2 (K_L r_e) \right]$$

where  $n$  is the harmonic number and  $r_e$  is beam radius. The gain function  $F_c$  is expressed in terms of the detuning parameter  $x = (n\omega_c - \omega)/(K_{II} V_{II})$  and has the following form:

$$F_c(x) = 0.5 \left[ \int_{\bar{Z}_{IN}}^{\bar{Z}_{OUT}} g(\bar{Z}) e^{ix\bar{Z}} d\bar{Z} \right]^2 \quad (4)$$

The frequency detuning that a gyrotron experiences can be written as:

$$\frac{2Q_T(\omega - \omega')}{\omega} = \frac{F_s - \frac{1}{2} \left( \frac{sv_1^2}{c^2} \right) \frac{dF_s}{dx}}{F_c - \frac{1}{2} \left( \frac{sv_1^2}{c^2} \right) \frac{dF_c}{dx}} \quad (5)$$

where  $\omega$  = cavity resonant frequency with beam

$\omega'$  = cavity resonant frequency without beam

$F_s$  = detuning gain function

In Eq.(5), the assumption  $s \ll Q_{ohmic}$  is made.

#### E. Calculational procedure

In this code, the starting current or frequency detuning is considered the y parameter. It is calculated as a function of two other parameters x and z. A range of x must be supplied, as well as discrete values of z. As a typical example, for a mode spectra the starting current would be calculated as a function of resonator magnetic field (x) for each gyrotron mode (z).

#### F. Notes

- 1) All parameters are in MKS units.
- 2) All data is inputted as real numbers unless otherwise noted.

#### REFERENCES

1. A.L. Gol'denberg and M.I. Petelin, Radiophysics and Quantum Electronics 16, No. 1 (1973) 106-111.

2. K.E. Kreischer and R.J. Temkin, Int. J. Infrared and Millimeter Waves 1, No. 2 (1980) 195-223. Also, K.E. Kreischer, M.I.T. Plasma Fusion Center Research Report PFC/RR-81-1, to be published in Infrared and Millimeter Waves, edited by K.J. Button.
3. K.E. Kreischer and R.J. Temkin, Int.J. Infrared and Millimeter Waves 2, No. 2 (1981) 175.

## II. INPUT AND OUTPUT

### A. Variables

Eleven device parameters are required to calculate linear characteristics. Each is referenced by a two letter abbreviation. The following is a table of these parameters with their abbreviations and dimensions:

<u>PARAMETER</u>	<u>ABBREVIATION</u>	<u>DIMENSIONS</u>
Beam voltage	VB	volts
Anode voltage	VA	volts
Cavity magnetic field	BR	tesla
Cathode magnetic field	BC	tesla
Mode	MO	none
Cavity radius	RR	meter
Cavity length	LE	meter
Cathode/anode distance	DC	meter
Ohmic Q	QO	none
Diffractive Q	QD	none
Cathode radius	RC	meter

### B. File INPUT.DAT

This file, which consists of three lines, provides information needed by the program

1. PAR(1), PAR(2),..., PAR(11)/

PAR(N) = default value of n<sup>th</sup> variable listed in A.

2. MODE(1), ZERO(1), MODE(2), ZERO(2),..., MODE(N), ZERO(N)/

MODE(N) = index of n<sup>th</sup> mode. For TE<sub>mpq</sub>, MODE = mp.

ZERO(N) = transverse index of n<sup>th</sup> mode. For TE<sub>mpq</sub>, the p<sup>th</sup> zero of

$$J_m^l(x) = 0$$

3.  $K, E_s, QTOT, Q(1), X(1), Q(2), X(2), \dots, Q(QTOT), X(QTOT)$ /  
 $K$  = axial wavenumber constant (see Section I.D.).  
 $E_s$  = stored energy factor (see Section I.D.).  
 $QTOT$  = number of q branches calculated for each  $TE_{mp}$  mode.  
 $Q(N)$  = value of q for  $n^{th}$  branch.  
 $X(N)$  = approximate value of detuning parameter X corresponding to  
 $n^{th}$  branch [3].

#### C. Other Information required during program execution

##### 1. VAR, VALUE

An opportunity to change default value is given.

##### 2. IMODE

IMODE = parameter indicating whether starting current or detuning is desired.

##### 3. PARX, XMIN, XMAX

PARX = two letter representation of x parameter.

XMIN = lower limit of x parameter.

XMAX = upper limit of x parameter.

##### 4. YMIN, YMAX

YMIN = minimum value of starting current or frequency detuning.

YMAX = maximum value of starting current or frequency detuning.

##### 5. ZPAR, ZNO, Z(1), Z(2), ..., Z(ZNO), /

ZPAR = two letter representation of z parameter.

ZNO = number of z values to be inputted.

Z(N) = value of  $n^{th}$  z parameter.

##### 6. ICROS

ICROS = parameter indicating cross-sectional structure of RF field.

7. IPROF

IPROF = parameter specifying type of axial RF profile.

(see Section II.D if you are specifying profile)

8. NHAR

NHAR = harmonic number.

At this point, desired data will be calculated, followed by a prompt for type of output.

9. IOUT

IOUT = parameter specifying type of output. IOUT = 3 terminates program.

D. Files and information needed if profile is specified

1. LINPROF. FORTRAN

A Fortran subroutine must be included with the program that contains the axial profile of the RF field. It should have the following form:

SUBROUTINE LINPROF (Q,Z)

Z = G (Q,Z)

RETURN

END

In this routine, you would provide the axial profile as the function G. Note that Q is the axial index of  $TE_{mpq}$ .

2. If the profile is specified, the parameters K and  $E_s$  of INPUT.DAT are ignored. The constant K is asked for specifically, and  $E_s$  is calculated within the program.

E. Tabular output

1. Heading title.
2. Date and time of program execution.
3. List of variables and their values. The x and z parameters are set to zero.
4. Tables of detuning or starting current data for each case requested.  
Each table is headed by q and z parameter values.

### III. SAMPLE RUN

#### A. The file INPUT.DAT

```
6.5e4,2.e4,5.57,.219,3.,3.47e-3,1.5e-2,4.e-3,2.e4,1.16e3,9.18e-3,30./  
11.,1.8412,21.,3.0542,1.,3.8317,31.,4.2012,41.,5.3176,12.,5.3314,  
51.,6.4156,22.,6.7061,2.,7.0156,61.,7.5013,32.,8.0152,13.,8.5363,  
71.,8.5778,42.,9.2824,81.,9.6474,23.,9.9695,3.,10.1735,52.,10.5199,  
91.,10.7114,33.,11.3459,14.,11.7060,62.,11.7349,101.,11.7709,  
43.,12.6819,111.,12.8265,72.,12.9324,24.,13.1704,4.,13.3237,121.,  
13.8788,53.,13.9872,82.,14.115,34.,14.5858,15.,14.8636,131.,14.9284,  
63.,15.2682,92.,15.2867,44.,15.9641,141.,15.9754,25.,16.3475,102.,  
16.4479,5.,16.4706,73.,16.5294,151.,17.0203,54.,17.3128,112.,17.6003/  
3.14159,0.5,3.,1.,-1.,2.,-4,2.,-1.4/
```

## B. Terminal input

```
vb= 0.650E+05      va= 0.200E+05      br= 0.557E+01
bc= 0.219E+00      mo= 0.300E+01      rr= 0.347E-02
le= 0.150E-01      dc= 0.400E-02      go= 0.200E+05
qd= 0.116E+04      rc= 0.918E-02
```

Any changes? Print 'no,/' or 'parameter,value'

```
no,/
Choose plot type:(inteser)
(1)Starting current vs. two parameters
(2)Retuning vs. two parameters(sine profile)
```

```
1
Input x parameter:aa,minimum,maximum
```

```
br,5.2,5.8
Input y parameter minimum,maximum
```

```
0.,10.
Input z parameter:aa,no. of values,value1,value2,...,/
mo,2.,3.,23.,/

```

```
Choose cross-sectional structure:(inteser)
(1)Rotatins
(2)Standing
```

```
2
Choose axial profile:(inteser)
(1)Sine
(2)Gaussian
(3)Profile you have specified
```

```
1
Input harmonic number(inteser)
```

```
1
```

```
Select output format:(inteser)
(1)Table
(2)Graphics
(3)Neither
```

```
1
Input table heading in apostrophes
```

```
Starting current for TE(031) and TE(231) modes'
```

### C. Terminal Output

Starting current for TE(031) and TE(231) modes  
01/21/82 1131.5 est Thu

vb= 0.650E+05      va= 0.200E+05      br= 0.000E+00  
bc= 0.219E+00      mo= 0.000E+00      rr= 0.347E-02  
le= 0.150E-01      dc= 0.400E-02      eo= 0.200E+05  
qd= 0.116E+04      rc= 0.918E-02

Standing cross sectional structure.

All data is in MKS units.

mo= 0.300E+01  
a = 0.100E+01

mo= 0.300E+01  
a = 0.200E+01

mo= 0.300E+01  
a = 0.200E+01

br	ISTART	br	ISTART	br	ISTART
--	-----	--	-----	--	-----
0.5625E+01	0.1290E+02	0.5800E+01	0.3155E+01	0.5504E+01	0.1465E+03
0.5610E+01	0.2891E+01	0.5795E+01	0.2984E+01	0.5492E+01	0.1730E+02
0.5596E+01	0.1739E+01	0.5790E+01	0.2862E+01	0.5480E+01	0.9847E+01
0.5581E+01	0.1320E+01	0.5785E+01	0.2781E+01	0.5468E+01	0.7300E+01
0.5567E+01	0.1123E+01	0.5780E+01	0.2735E+01	0.5456E+01	0.6113E+01
0.5552E+01	0.1028E+01	0.5774E+01	0.2721E+01	0.5444E+01	0.5515E+01
0.5538E+01	0.9934E+00	0.5769E+01	0.2739E+01	0.5432E+01	0.5249E+01
0.5523E+01	0.1003E+01	0.5764E+01	0.2789E+01	0.5420E+01	0.5209E+01
0.5509E+01	0.1051E+01	0.5759E+01	0.2873E+01	0.5408E+01	0.5357E+01
0.5494E+01	0.1140E+01	0.5754E+01	0.2997E+01	0.5396E+01	0.5688E+01
0.5480E+01	0.1279E+01	0.5749E+01	0.3169E+01	0.5384E+01	0.6224E+01
0.5465E+01	0.1481E+01	0.5744E+01	0.3401E+01	0.5372E+01	0.7015E+01
0.5451E+01	0.1776E+01	0.5739E+01	0.3714E+01	0.5360E+01	0.8150E+01
0.5436E+01	0.2206E+01	0.5733E+01	0.4140E+01	0.5348E+01	0.9778E+01
0.5422E+01	0.2854E+01	0.5728E+01	0.4733E+01	0.5336E+01	0.1216E+02
0.5408E+01	0.3869E+01	0.5723E+01	0.5594E+01	0.5324E+01	0.1577E+02
0.5393E+01	0.5562E+01	0.5718E+01	0.6928E+01	0.5313E+01	0.2155E+02
0.5379E+01	0.8651E+01	0.5713E+01	0.9223E+01	0.5301E+01	0.3161E+02
0.5364E+01	0.1520E+02	0.5708E+01	0.1401E+02	0.5289E+01	0.5175E+02
0.5350E+01	0.3395E+02	0.5703E+01	0.2974E+02	0.5277E+01	0.1050E+03

C. Terminal Output (continued)

$m_0 = 0.230E+02$   
 $a = 0.100E+01$

$m_0 = 0.230E+02$   
 $a = 0.200E+01$

$m_0 = 0.230E+02$   
 $a = 0.200E+01$

br	ISTART	br	ISTART	br	ISTART
--	-----	--	-----	--	-----
0.5512E+01	0.3672E+02	0.5730E+01	0.1652E+02	0.5387E+01	0.4064E+03
0.5497E+01	0.4108E+01	0.5723E+01	0.9316E+01	0.5377E+01	0.2895E+02
0.5482E+01	0.2340E+01	0.5715E+01	0.6624E+01	0.5367E+01	0.1588E+02
0.5467E+01	0.1744E+01	0.5708E+01	0.5250E+01	0.5358E+01	0.1146E+02
0.5453E+01	0.1472E+01	0.5701E+01	0.4443E+01	0.5348E+01	0.9329E+01
0.5438E+01	0.1343E+01	0.5693E+01	0.3935E+01	0.5338E+01	0.8156E+01
0.5423E+01	0.1296E+01	0.5686E+01	0.3611E+01	0.5328E+01	0.7487E+01
0.5408E+01	0.1308E+01	0.5679E+01	0.3410E+01	0.5318E+01	0.7130E+01
0.5393E+01	0.1372E+01	0.5671E+01	0.3303E+01	0.5308E+01	0.6996E+01
0.5379E+01	0.1490E+01	0.5664E+01	0.3276E+01	0.5298E+01	0.7042E+01
0.5364E+01	0.1674E+01	0.5657E+01	0.3324E+01	0.5289E+01	0.7249E+01
0.5349E+01	0.1943E+01	0.5649E+01	0.3450E+01	0.5279E+01	0.7621E+01
0.5334E+01	0.2334E+01	0.5642E+01	0.3671E+01	0.5269E+01	0.8171E+01
0.5319E+01	0.2906E+01	0.5634E+01	0.4015E+01	0.5259E+01	0.8933E+01
0.5304E+01	0.3767E+01	0.5627E+01	0.4536E+01	0.5249E+01	0.9958E+01
0.5290E+01	0.5118E+01	0.5620E+01	0.5341E+01	0.5239E+01	0.1133E+02
0.5275E+01	0.7368E+01	0.5612E+01	0.6659E+01	0.5230E+01	0.1316E+02
0.5260E+01	0.1147E+02	0.5605E+01	0.9082E+01	0.5220E+01	0.1564E+02
0.5245E+01	0.2015E+02	0.5598E+01	0.1473E+02	0.5210E+01	0.1909E+02
0.5230E+01	0.4474E+02	0.5590E+01	0.4092E+02	0.5200E+01	0.2402E+02

Select output format:(integer)

- (1)Table
- (2)Graphics
- (3)Neither

3  
r 11:32 12.899 90

IV. PROGRAM LISTING

```

c Program Linear
c Written by K. E. Kreischer, M.I.T.
c Latest Revision: 12/1/81
c Purpose: Calculates linear characteristics of gyrotron.

c external references for Multics graphics
c external plot $setup(descriptors)
c external plot $scale(descriptors)
c external plot_ (descriptors)
c external reference for data in table
c external date time
c common par,xdat
c common/prof/zmin,ph,istep
c common/gun/wnoz,wnp,u,w,freq,freqb,x,gfac,se,q,x1,xm,xfun(4)
c dimension ch1(60,2).par(12).data(2400,2).zparam(20).xdat(50)
c &.xpts(2).x1im(2).ica(3).ija(3).ta(3)
c character parx*2,parz*2,ytitle*50,xttitle*50
c dimension xdat1(50),ydat1(50)
c data data,zparam/4820*0./

c file input.dat opened for reading
c open(7,form='formatted',file='input.dat')
c first line of data from input.dat is read and displayed
c read(7,)(par(1),i=1,12)
c write(6,5)(par(1),i=1,11)
c format(' vb='',e10.3,5x, va='',e10.3,5x, br='',e10.3/, bc='',e10.3,
c 85x,'mo='',e10.3,5x, rr='',e10.3/, 1e='',e10.3,5x, dc='',e10.3,5x,
c 8,'qo='',e10.3/, qd='',e10.3,5x, rc='',e10.3/)

c write(6,6)
c format(' Any changes? print "no." or "parameter,value"/')
c changes to default values accepted
c read(5,40)parx,parval
c if(parx .eq. 'no') goto 15
c call iden(parx,1)
c if(1.eq.99) goto 10
c par(1)=parval
c goto 7
c write(6,11)
c format(' Error in last entry/')
c goto 7
c rest of input.dat is read
c 15 read(7,)(ch1(1,1),j=1,2),i=1,60)
c read(7,)(xdat(1),i=1,50)
c type of calculation is chosen
c write(6,25)
c format(' Choose plot type:(integer) /' (1)Starting current vs. two
c 8parameters /, (2)Detuning vs. two parameters(sine profile)./')
c read(5,) imode
c info on x,y parameters and their ranges is obtained
c write(6,35)
c format(' Input x parameter:aa,minimum,maximum /')
c read(5,40)parx,xmin,xmax
c format(v)
c call iden(parx,1par)
c write(6,43)
c format(' Input y parameter minimum,maximum /')
c read(5,) ymin,ymax
c z parameter and its discrete values are inputted
c write(6,50)

```

```

50      format(' Input z parameter:aa,no. of values,value1,value2
8...//')
     read(5,40) parz,zno, (zparam(i), i=1,20)
     call iden(parz,izpar)
c   Info on beam and rf field are inputted
     write(6,55)
     format(' Choose cross-sectional structure:(integer) '/',
&(1)Rotating', '(2)Standing')
     read(5,) icros
     write(6,60)
     format(' (3)profile you have specified/')
     read(5,) iprof
     if(iprof.ne.3) goto 80
c   info on specified profile is obtained
     write(6,61)
     format(' Input k(parallel) constant/')
     read(5,) xdat(1)
     write(6,62)
     format(' Input z limit of axial profile:lower,upper')
     read(5,) zmin1,zmax1
     write(6,63)
     format(' Input number of integration steps(integer)')
     read(5,) istep
c   Info on harmonic is inputted
     write(6,81)
     format(' Input harmonic number (integer)')
     read(5,) nhar
c   determine if detuning parameter is a function of x parameter
     1xvrx=0
     if(ixpar .ne. 8) 1xvrx=1
     11=2
     if(icros .ne. 1) 11=1
c   do loop varying q of TE(mpq)
     do 500 j=1,xdat(3)
        q=xdat(2*j+2)
        if(ixvrx.eq.0) goto 95
        if(iprof.ne.3) goto 87
        zmin=q*zmin1*xdat(1)/par(7)
        zmax=q*zmax1*xdat(1)/par(7)
        ph=(zmax-zmin)/istep
        x=xdat(2*j+3)
        call xcal(iprof,imode)
        df=xfun(2)
        call gun(nhar,iprof)
        fac=0.5*nhar*freb*w**2/wnoz/u/9.e16
        do 90 k1=1,2
           x=xdat(2*j+3)
           do 85k2=1,2
              step=(0.2/q)/10***(k2-1)
              do 82k3=1,10
                 x=x+(-1.)*(k1+k2)*step
                 call xcal(iprof,imode)
                 test=-xfun(1)+fac*xfun(2)
                 if(k2.eq.1.and.test.lt.df/10) goto 85
                 if(k2.eq.2.and.test.gt.df/10) goto 85
                 continue
               x11m(k1)=x
85

```

```

90 continue
c branch to separate algorithm if profile is specified
if (iprof.eq.3) goto 401
c do loop varying z parameter
do 500k=1,zno
  par(izpar)=zparam(k)
c select correct mode index
  do 70 j1=1,60
    if (int(par(5)).ne. int(chi(j1,1))) goto 70
    x1=chi(j1,2)
    xm=int(chi(j1,1))/10
  goto 75
70   continue
c branches depending on whether x parameter is a function of
c detuning parameter
  75  if (ixvrx .eq. 1) goto 130
c direct calculation of data/ixvrx=0
  call gun(nhar,iprof)
  call xcal(iprof,imode)
  do 120 i=1,11
    rot=(-1.)**i
    do 120i=1,50
      par(ixpar)=(i-1)*(xmax-xmin)/49.+xmin
      call istart(nhar,imode,iprof,icros,rot,val,1,0,1)
      ic=(1-1)*zno*xdat(3)+(k-1)*xdat(3)+j
c input of final data into file
      data((ic-1)*51+i,1)=par(ixpar)
      data((ic-1)*51+i,2)=val
      igap=51
      goto 500
c direct calculation of data/ixvrx=1
130  del=.01/q**2
c vary x parameter until detuning parameter is matched
c determine range for each mode in terms of x parameter
  do 160 i=1,2
    ex=x1*lm(i)
    icnt=0
    call match(xmin,xmax,par(ixpar),del)
    call gun(nhar,iprof)
    test=abs(x-ex)
    call match1(test,par(ixpar),icnt)
    goto (155,157,158,159) icnt
    xpts(i)=par(ixpar)
    goto 160
    xpts(i)=xmax
    goto 160
    xpts(i)=xmin
159   continue
  do 170 i=1,11
    rot=(-1.)**i
    do 170i=1,20
      par(ixpar)=(i-1)*(xpts(2)-xpts(1))/19.+xpts(1)
      call istart(nhar,imode,iprof,icros,rot,val,1,1,1)
      ic=(1-1)*zno*xdat(3)+(k-1)*xdat(3)+j
c. input of final data into file
      data((ic-1)*21+i,1)=par(ixpar)
      data((ic-1)*21+i,2)=val
      igap=21
      goto 500
c calculate data for case of iprof=3 and ixvrv=1.

```

```

c minimize number of integrations in calculation of x functions
c in subroutine xcal.
c energy constant is calculated
401 icnt=0
val=0
call integ(zmin,ph,istep+1,val,icnt)
if (icnt.eq.1) goto 425
call lmpoff(val,q)
val=val**2
goto 420
xdat(2)=val/xdat(1)/q
do 409 i=1,20
ex=x1im(1)+(i-1)*(x1im(2)-x1im(1))/19.
x=ex
call xcal(iprof,imode)
do 409kd=1,zno
par(lzpar)=zparam(kl)
do 406 j=1,60
if (int(par(5)).ne.int(chi(j1,1))) goto 406
xi=chi(j1,2)
xm=int(chi(j1,1))/10..
goto 407
406 continue
c find x parameter corresponding to each detuning parameter
407 icnt=0
del=.01/q**2
call match(xmin,xmax,par(ixpar),del)
call gun(nhar,iprof)
test=abs(x-ex)
call match1(test,par(ixpar),icnt)
goto(404,403,405,405) icnt
val=-999.
do 409 i=1,11
rot=(-1.)**i
if (icnt.ge.3) goto 408
call startnhar,imode,iprof,lcros,rot,val,0,0,1)
mm=i+21*((i-1)*zno+xdat(3)+(k-1)*xdat(3)+j-1)
403
405
408
409
500 continue
c selection of output format
501 write(6,502)
502 format(' Select output format: (integer) /' '(1)Table /'
& '(2)Graphics /' '(3)Neither /')
read(5, ) tout
goto(504,600,900) tout
c output of data in tabular form
504 write(6,506)
506 format(' Input table heading in apostrophes /')
c read(5,40) mttitle
write(6,508) mttitle
format(10(/),1x,a50)
call date_time
do 510 i = 1,12
if (j.eq.ipar.or.j.eq.lzpar) par(j) = 0
510 write(6,511)
format(//)
write(6,5)(par(j),j=1,11)
511

```

```

if (icros.eq.1) goto 513
write(6,512)
format(,' Standing cross sectional structure.')
512
      goto 517
      write(6,516)
      format(,' Rotating cross sectional structure.')
513
      516
      write(6,511)
517
      write(6,515)
      format(,' All data is in MKS units.'//)
514
      n=0
      do 550 k=1,xdat(3)*zno+11
n=n+1
      i1a(n)=aint(1.01+(k-1)/(xdat(3)*11))
      i1ja(n)=aint((1.01+(k-1)/11)-xdat(3)*(i1a(n)-1))
      i1ca(n)=k-1
      if (n.eq.3.or.k.eq.(xdat(3)*zno+11)) goto 520
      goto 530
      write(6,522)parz,zparam(i1a(1)),parz,zparam
      8,(i1a(2)),parz,zparam(i1a(3))
      format(8x,2(a2," ",e10.3,13x),a2,'=',e10.3)
520
      write(6,524)xdat(2*i1ja(kk)+2),kk=1,3
      format(8x,2('q ',e10.3,13x),'q ',e10.3/)
522
      format(8x,2('q ',e10.3,13x),'q ',e10.3/)
524
      ytitle=" I START "
      if (imode.eq.2)ytitle="DETUNING"
      write(6,526)parx,ytitle,parx,ytitle
      format(7x,2,a2,8x,a8,8x,a8)
526
      write(6,528)
      format(7x,2(' ',8x,8(' '_'),8x),'_',8x,8(' '_'))/
528
      do 540 i=1,igap-1
      write(6,530)((data(i1ca(m)*igap+j,kk),kk=1,2),m=1,n)
530
      format(6(2x,e11.4))
540
      continue
      write(6,511)
      n=0
      continue
      goto 501
c output of data in graphic form
550
      write(6,605)
      format(,' Input heading title in apostrophes')
      read(5,40)mtitle
      write(6,610)
      format(,' Input x-axis title in apostrophes')
      read(5,40)xtitle
      write(6,615)
      format(,' Input y-axis title in apostrophes')
      read(5,40)yttitle
      write(6,620)
      format(,' Input plot type:(integer) /' ,(1)Llinear-Llinear'/' ,
      & (2)Log x-Linear y /', (3)Llinear x-Log y /', (4)Log-Log /')
      read(5,40)itype
      base=10.
      if (itype.eq.1) base=0.
      call plot$setup(mttitle,xttitle,ytitle,itype,base,0.0)
      call plot$scale(xmin,xmax,ymin,ymax)
      do 640 i=0,xdat(3)*zno+11-1
      1xyd1m=0
      do 630 j=1,igap-1
      if (data(i*igap+j).gt.ymin.or.data(i*igap+j,2).gt.ymax)
      & goto 630
      1xyd1m=1xyd1m+1

```

- 19 -

```

xdat1(1xydim)=data(l*lgap+j,1)
ydat1(1xydim)=data(l*lgap+j,2)
continue
if(1xydim.eq.0) goto 640
call plot_(xdat1,ydat1,1xydim,1,'.')
continue
goto 501
continue
end

subroutine iden(a,ib)
c determines numerical code for device variables
character*2 a

ib=99
if(a .eq. 'vb') ib=1
if(a .eq. 'va') ib=2
if(a .eq. 'br') ib=3
if(a .eq. 'bc') ib=4
if(a .eq. 'mo') ib=5
if(a .eq. 'rr') ib=6
if(a .eq. 'le') ib=7
if(a .eq. 'dc') ib=8
if(a .eq. 'qo') ib=9
if(a .eq. 'qd') ib=10
if(a .eq. 'rc') ib=11
if(a .eq. 'cb') ib=12
return
end

subroutine xcail(iprof,imode)
c calculates gain functions and their derivatives
common par(12),xdat(50)
common/prof/zmin,ph,lstep
common/gun/wnoz,wnop,u,w,frer,freb,x,gfac,se,q,xm,xfun(4)
do 2 j=1,4
  xfun(j)=0.
  p1=3.141592654
  do 50k=1,2
    x1=x-.001*(-1)**k
    if (iprof .ne. 1) goto 5
    fc=2*sin((x1+1)*p1*q/2)**2/(1-x1**2)**2
    if (iprof .ne. 1 .or. imode .ne. 2) goto 10
    fs=-1/(1-x1**2)**2*(sin(q*p1*(x1-1))+x1*(1-x1**2)*q*p1/2)
    5   if (iprof .ne. 2) goto 15
    fc=p1/2*exp(-.5*x1**2)
    10  if (iprof .ne. 3) goto 20
    15  c gain function calculated by integration
      fc=0
      do 19 j=1,2
        icnt=0
        val=0
        call integ(zmin,ph,lstep+1,val,1cnt)
        val1=val
        if (icnt .eq. 1) goto 19
        call linprof(val,q)
        if (j .eq. 1) val=val*cos(x1*val1)
        if (j .eq. 2) val=val*sin(x1*val1)
        goto 17
        17  fc=0.5*val*val+fc
        xfun(1)=xfun(1)+.5*fc
        xfun(2)=xfun(2)-500*(-1)**k*fc
        xfun(3)=xfun(3)+.5*fc
      19

```

```

50 xfun(4)=xfun(4)-500*(-1)**k*fs
      return
end
c integration done by Simpson approximation
if (icnt .eq. 0) accum=0
if (icnt .eq. 0) index=0
if (index .eq. 1 .or. index .eq. 1step) val=val/2
accum=accum+val*(3+(-1)**index)*h/3
icnt=2
index=index+1
val=sminth*(index-1)
if (index .eq. (1+istep)) val=accum
if (index .eq. (1+istep)) icnt=1
return
c subroutine bessel(xm,arg,func)
c function J(arg) of order xm is calculated
c result returned as func.
icnt=0
val=0
istep=10*max1(abs(xm),arg)+1
istep=max0(istep,30)
h=3.141592654/(istep-1)
call integ(0.,h,istep,val,icnt)
if (icnt .eq. 1) goto 10
val=cos(arg*sin(val)-xm*val)
goto 5
func=val/3.14159
return
end
c remaining device variables and detuning parameter are calculated
c assuming adiabatic theory
common par(12),xdat(50)
common/gun/wnoz,wnop,u,w,frer,freb,x,gfac,se,q,x1,xm,xfun(4)
ec=par(2)/par(8)
ec=1+par(1)/par(8)
gam=1+par(1)/5.1e5
al=par(3)/par(4)
beta=sqrt(1-1/gam**2)
w=sqrt(al)*ec/par(4)
wnop=x1/par(6)
wnoz=xdat(1)/par(7)*q
frer=3.e8*sqrt(wnoz**2+wnop**2)
freb=1.75758e11*par(3)/gam
arg=9.e16*beta**2-w**2
if (arg .lt. 0.) goto 10
u=sqrt(arg)
x=(nhar*freb-frer)/wnoz/u
return
x=0.
u=1.e10
return
end
c beam-field coupling and stored energy factors are calculated
common par(12),xdat(50)
common/gun/wnoz,wnop,u,w,frer,freb,x,gfac,se,q,x1,xm,xfun(4)
call bessel(xm,x1,func)
se=1.57*par(7)/wnop**2*(x1**2-xm**2)*

```

```

&func**2*xdat(2)
  rf(rbeam-sqrt(par(4)/par(3))*par(11)
  if (rbeam.ge.par(6)) gfac=-1.e-10
  if (rbeam.ge.par(6)) goto 50
  arg=wnop*rbeam
  if (icros.ne.1) goto 15
  call bessel(xm+rot*nhar,arg,func)
  gfac=.5*func**2
  goto 50
  har=nhar
  end

15   call bessel(xm-har,arg,func)
  call bessel((xm+har,arg,func)
  gfac=.25*(func1**2+func**2)
  return

  subroutine match(amin1,amax1,var,del1)
c 'var' is varied until 'val' is within limit set by 'del1'.
c also determines if match is outside parameter limits 'amin1,amax1'
  amin=amin1
  amax=amax1
  del1=del11
  del2=(amax-amin)/400.
  fmax=amax
  fmin=amin
  val=amin+0.618034*(amax-amin)
  return
  entry match1(val,var,icnt)
  if (icnt.ne.0) goto 10
  var1=val
  var1=var
  icnt=1
  goto 20
  if (val.gt.var1) goto 15
  if (var1.le.var) amin=var1
  if (var1.gt.var) amax=var1
  var1=var
  val=var1
  goto 20
  if (var.le.var1) amin=var
  if (var.gt.var1) amax=var
  if (abs(fmax-var1).lt.del1) icnt=3
  if (abs(fmin-var1).lt.del1) icnt=4
  if (val1.lt.del1) icnt=2
  if (icnt.eq.2) var=var1
  if (icnt.eq.2) val=val1
  if (icnt.gt.1) goto 30
  varf=amin+0.381966*(amax-amin)
  var=amin+0.618034*(amax-amin)
  if (abs(var-var1).lt.abs(varf-var1)) var=varf
  return
  end

30   subroutine istart(nhar,imode,iprof,icros,rot,val,11,12,13)
c starting current or frequency detuning is calculated
  common par(12),xdat(50)
  common/gun/wnoz,wnop,u,w,freq,freq,x,prof,xfun(4)
  if (11.eq.1) call gun(nhar,iprof)
  if (12.eq.1) call xcaltiprof(imode)
  if (13.eq.1) call crosec(nhar,icros,rot,iprof)
  fac=.5*nhar*freq*w**2/wnoz/u/9.e16
  if (imode.eq.2) goto 5

```

```
nr=q*whop*w/freb
har=nhar
call bessel(har-1.,arg,func1)
call bessel(har+1.,arg,func2)
harfac=func1-func2
qt=par(10)/q**2*(x1/10.1735)**2
c starting current is calculated
val=-1.e-22*frer/qt*se*wnoz**2*u**2/gfac/harfacc**2/
8(xfun(1)-fac*xfun(2))*(1+par(1)/5.11e5)
goto 10
c frequency detuning is calculated
5      val=(xfun(3)-fac*xfun(4))/(xfun(1)-fac*xfun(2))
10    return
end
```

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