



INTERNATIONAL ATOMIC ENERGY AGENCY
UNITED NATIONS EDUCATIONAL, SCIENTIFIC AND CULTURAL ORGANIZATION



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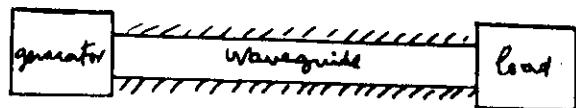
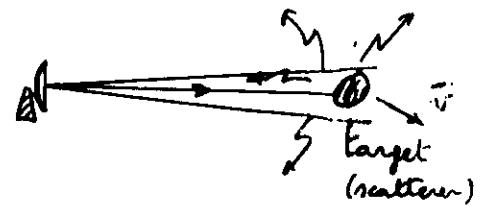
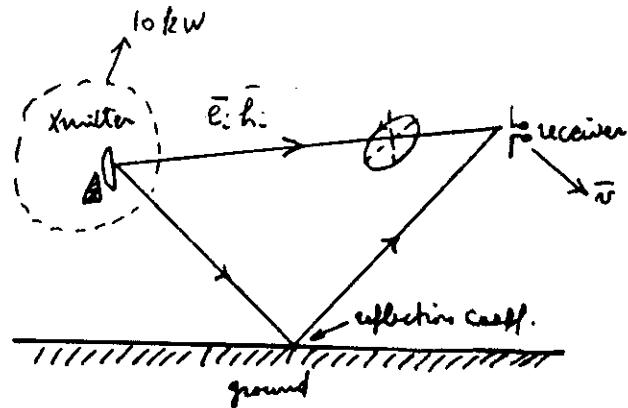
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COURSE ON BASIC TELECOMMUNICATIONS SCIENCE
9 January - 3 February 1989

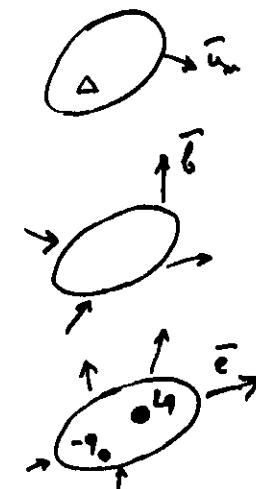
REFRESHER COURSE

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These notes are intended for internal distribution only.



0'



Divergence

$$\operatorname{div} \vec{v} = \frac{\partial v_x}{\partial x} + \frac{\partial v_y}{\partial y} + \frac{\partial v_z}{\partial z}$$

$$\iint_S \vec{v} \cdot \vec{n} dS = \iiint_V \operatorname{div} \vec{v} dV \approx \operatorname{div} \vec{v} \Delta V$$

$\operatorname{div} \vec{v} = 0$ "as much goes in as goes out"

$\operatorname{div} \vec{v} > 0$ "more goes out than goes in" the flux diverges

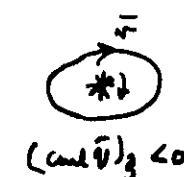
$\operatorname{div} \vec{v} < 0$ net convergence

Curl

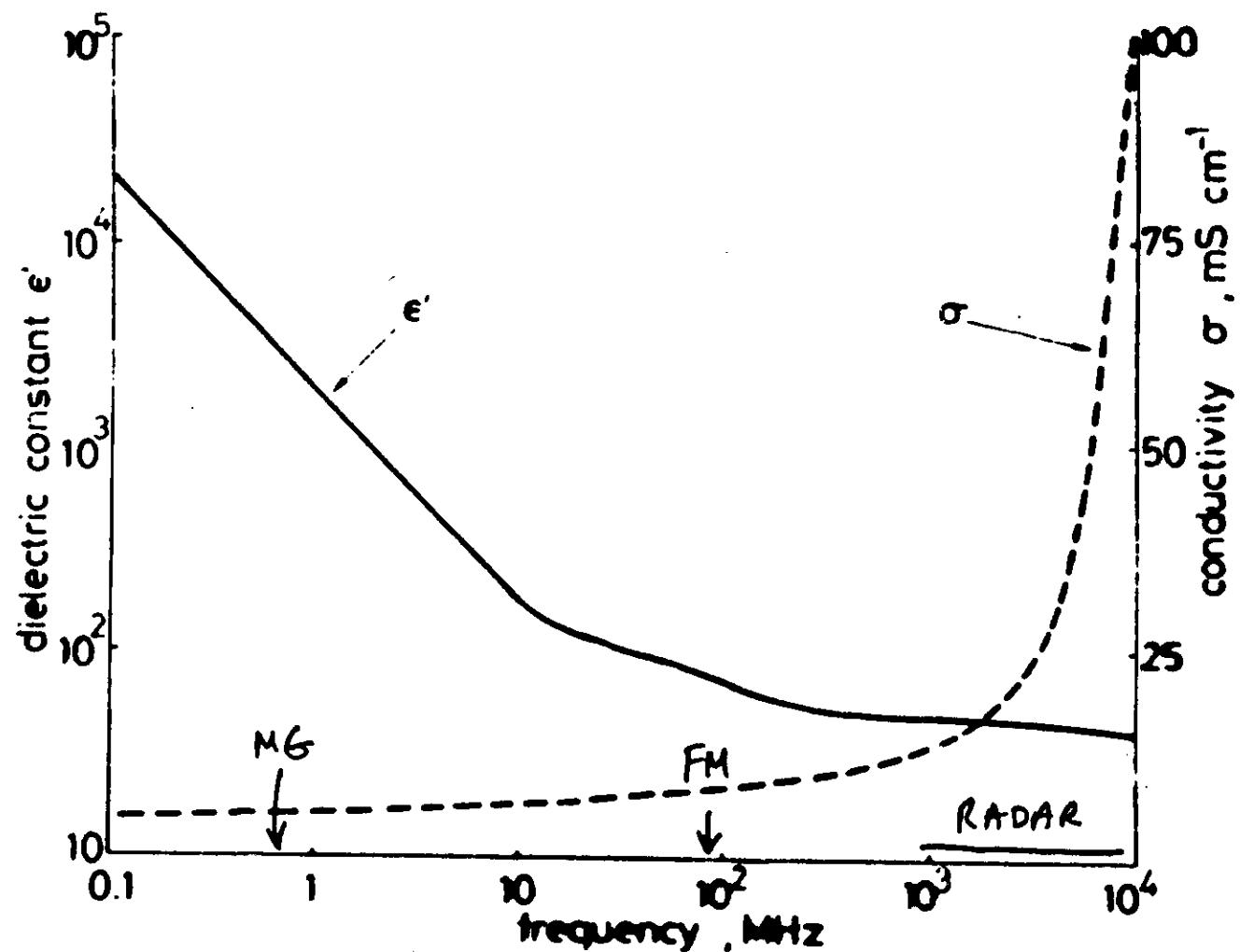
$$\operatorname{curl} \vec{v} = \left(\frac{\partial v_y}{\partial z} - \frac{\partial v_z}{\partial y} \right) \vec{u}_x + \left(\frac{\partial v_z}{\partial x} - \frac{\partial v_x}{\partial z} \right) \vec{u}_y + \left(\frac{\partial v_x}{\partial y} - \frac{\partial v_y}{\partial x} \right) \vec{u}_z$$

$$\iint_S \vec{v} \cdot \vec{n} dS = \iiint_V \vec{u}_z \cdot \operatorname{curl} \vec{v} dV$$

$\operatorname{curl} =$ tendency to rotate (in hydrodynamics)



$\operatorname{curl} = 0$ irrotational flow
paddle-wheel stationary



Dielectric constant ϵ' and conductivity $\sigma(\text{mS cm}^{-1})$ of muscle tissue as functions of frequency

Properties of Alternatives in biological medium

Frequency (MHz)	Wavelength In Air (cm)	Dielectric Constant (ϵ_r)	Conductivity σ (S/meter)	Wavelength λ_x (cm)	Depth of Penetration (cm)	Muscle, Skin, and Tissues with High Water Content		Reflection Coefficient Air-Muscle Interface	Reflection Coefficient Muscle-fat Interface	Θ
						r	s			
1	30,000	2.000	0.400	436	91.3	0.982	+179	--	--	0.28
10	3,000	110	0.625	118	21.6	0.956	+178	--	--	0.14
27.12	1,100	113	0.602	68.1	14.3	0.925	+177	0.651	-11.13	0.28
43.00	738	97.3	0.580	51.3	11.2	0.913	+176	0.652	-10.21	0.32
100	300	71.7	0.495	27	6.66	0.881	+175	0.650	-7.96	0.45
200	150	56.5	1.00	16.6	4.79	0.844	+175	0.612	-8.06	0.63
300	100	54	1.15	11.9	3.89	0.825	+175	0.592	-8.14	0.78
433	68.3	53	1.18	8.76	3.57	0.803	+175	0.562	-7.86	1.06
700	49	52	1.25	5.34	3.18	0.779	+176	0.532	-5.60	1.73
915	39.8	51	1.28	4.46	3.04	0.772	+177	0.519	-4.32	2.02
1200	30	50	1.35	3.66	2.42	0.761	+177	0.506	-3.05	2.62
2400	15.2	47	2.17	1.76	1.70	0.754	+177	0.500	-3.03	2.96
3000	11.1	45	2.27	1.45	1.61	0.751	+178	0.495	-3.20	3.38
3500	9	44	4.95	0.99	0.708	0.749	+177	0.492	-4.06	2.69
3600	8.67	43.3	4.93	0.775	0.720	0.746	+177	0.492	-4.29	2.83
3700	8.38	43	0.33	0.570	0.413	0.744	+176	0.513	-6.66	2.13
3800	8.10	43	10.00	0.464	0.343	0.743	+176	0.510	-5.96	2.22

Values are from Sato and others (1970) for a medium of 37°C .

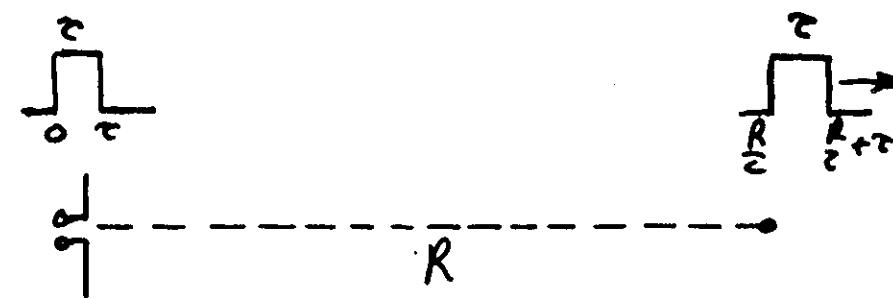
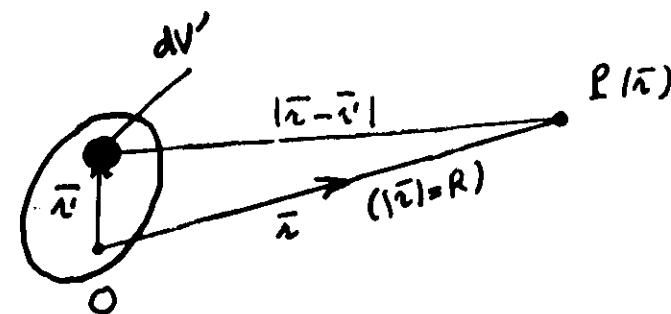
Properties of microwaves in biological media*

Fat, Bone, and Tissues with Low Water Content										
Frequency (MHz)	Wavelength In Air (cm)	Dielectric Constant (ϵ_L)	Conductivity σ_L (mS/m)	Wavelength λ_L (cm)	Depth of Penetration (cm)	Reflection Coefficient		Reflection Coefficient Fat-Muscle Interface		Q_{av}
						Air-Fat Interface		r	s	
1	30,000	--	--	--	--	--	--	--	--	--
10	3,000	--	--	--	--	--	--	--	--	--
27.12	1.106	20	10.9 - 43.2	241	159	0.660	+174	0.651	+169	1.11
40.68	.738	14.6	12.6 - 52.8	187	118	0.617	+173	0.652	+170	1.00
100	.300	7.45	19.1 - 75.9	106	60.4	0.511	+168	0.650	+172	0.87
200	.150	5.95	25.8 - 94.2	59.7	39.2	0.458	+168	0.612	+172	1.10
300	.100	5.7	31.6 - 107	41	32.1	0.438	+169	0.592	+172	1.38
433	.64.3	5.6	37.9 - 118	28.8	26.2	0.427	+170	0.562	+173	1.73
750	.40	5.6	49.8 - 130	16.8	23	0.415	+173	0.532	+174	2.48
915	.32.8	5.6	55.6 - 147	13.7	17.7	0.417	+173	0.519	+176	2.81
1500	.20	5.6	70.8 - 171	8.41	13.9	0.412	+174	0.506	+176	3.86
2450	.12.2	5.5	96.4 - 213	5.21	11.2	0.406	+176	0.500	+176	4.84
3000	.10	5.5	110 - 234	4.25	9.74	0.406	+176	0.495	+177	5.33
5000	.6	5.5	162 - 309	2.63	6.67	0.393	+176	0.502	+175	6.49
5900	.5.17	5.05	186 - 338	2.29	5.24	0.388	+176	0.502	+176	6.21
8000	.3.75	4.7	256 - 431	1.73	4.61	0.371	+176	0.513	+173	6.09
10000	.3	4.5	324 - 549	1.41	3.30	0.363	+175	0.518	+174	5.73

*Based on data from Schwan and Phamont (24) for temperature of 30° C.

6'

$$\phi(\vec{r}, t) = \frac{1}{4\pi\epsilon_0} \iiint \frac{\rho(\vec{r}', t - \frac{|\vec{r}-\vec{r}'|}{c})}{|\vec{r}-\vec{r}'|} dV'$$



Dielectric constants and loss factors of common materials

Material	Dielectric constant (ϵ')		Loss factor ($\epsilon''/\epsilon' \times 10^{-4}$)	
	at 10 MHz	at 3000 MHz	at 10 MHz	at 3000 MHz
Metals				
copper	39	46	1300	12
gold	4.5	2.5	4.2	0.18
Insulators and glasses				
fused quartz	3.78	3.78	0.00004	0.0002
pyrex (soda-lime-silicate)	4.04	4.02	0.0015	0.026
radio glass (SiO ₂ , Na ₂ O, CaO, MgO, Al ₂ O ₃)	6.3	-	0.14	-
Minerals				
calcite, dry	9.0	9.0	0.33	0.22
feldspar (K-feldspar)	5.4	5.4	0.0016	0.0016
mica, dry mica	2.55	2.05	0.04	0.016
Mineral Fibers				
cotton, 200 kg/m ² , 7% moisture	1.5	at 27	0.03	at 27
wool, 150 kg/m ² , 20% moisture	1.2	MHz	.01	MHz
Organic fibers				
hemp, yellow	2.05	2.39	0.020	0.018
white oil	2.2	2.2	0.009	0.004
rayon, viscose	2.25	2.25	0.00045	0.00045
Paper and board				
newsprint	1.5	at 27	0.4	0.4
cardboard	1.5	at 27	0.5	0.4
Plastics				
cellulose acetate (acetate rayon)	5.5	4.2	0.07	0.00
melamine-formaldehyde (formaldehyde molding compound)	4.3	3.7	0.23	0.22
phenol-formaldehyde (Bakelite)	4.3	3.0	0.18	0.15
polyacrylic (Plexiglas)	3.2	3.0	0.09	0.04
polyester	4.0	4.0	0.04	0.04
polyethylene (polythene)	2.25	2.25	0.00004	0.001
polymethyl methacrylate (Perspex, plexiglas)	2.7	2.6	0.027	0.015
polycarbonate	2.25	2.25	0.00005	0.00005
polytetrafluoroethylene (PTFE)	2.1	2.1	0.00005	0.00003
polycarbonate (PCV), clear	2.9	2.8	0.03	0.02
polycarbonate (PCV), colored	2.7	2.9	0.4	0.1
Glass				
borosilicate	2.4	2.4	0.000	0.007
borosilicate	2.5	2.5	0.08	0.03
Resins				
epoxy, room temp.	3.7	3.2	0.07	0.003
epoxy, room temp 25°C	70.0	70.0	0.30	12.0
epoxy, 100-meter 25°C	60.0	70.0	70.00	18.00
Other				
silicones, room temp	2.6	2.5	0.1	0.07
silicones, liquid	2.5	1.0	0.00	0.05

Source: Handbook of Chemistry and Physics, 1971-72

BASIC TELECOM - third week



```
C-----
C This is a simple ray tracing program
C The refractive index is a function of the height only
C common delta
C write(6,20)
20 format(' Exponential decay factor')
read(5,*) delta
write(6,1)
1 format(' Starting angle in degrees')
read(5,*) stangle
stangle=stangle*3.141592654/180.
write(6,2)
2 format(' Unit step in the horizontal direction')
read(5,*) delx
write(6,10)
10 format(' Number of steps')
read(5,*) nmax
x=0.
y=0.
yex=0.
write(6,3)
3 format(' Horizontal distance vers vs height')
write(6,*) x,y
C
C
ta=tan(stangle)
100 dely=delx*ta
y=y+dely
yex=yex+delx*tan(stangle-x/delta)
x=x+delx
write(6,*) x,y,yex
call refindex(y,res)
ta=(ta+res*delx)/(1.-ta*res*delx)
itot=itot+1
if(itot.eq.nmax) stop
goto 100
C
end
C-----
subroutine refindex(y,res)
common delta
C
C This subroutine provides (dN/dy)/N
C
example: exponential profile
N=Aexp(-y/delta)
C
res = -1./delta
return
end
C-----
C This is a simple ray tracing program
C The refractive index is a function of the height only
C common an0,an1,delta
C write(6,20)
20 format(' Exponential decay factor')
```

```

read(5,*)
write(6,30)
format(' Refractive index profile constants N0 and N1:')
1 read(5,*)
an0,an1
write(6,*)
an0,an1
write(6,1)
format(' Starting angle in degrees')
read(5,*)
stangle
stangle=stangle*3.141592654/180.
write(6,2)
format(' Unit step in the horizontal direction')
read(5,*)
delx
write(6,10)
format(' Number of steps')
read(5,*)
nmax
x=0.
y=0.
yex=0.
write(6,3)
format(' Horizontal distance versus height')
write(6,*)
x,y

ta=tan(stangle)
dely=delx*ta
y=y+dely
x=x+delx
write(6,*)
x,y
call refindex(y,res)
ta=(ta+res*delx)/(1.-ta*res*delx)
itot=itot+1
if(itot.eq.nmax) stop
goto 100

end
-----
subroutine refindex(y,res)
common an0,an1,delta
This subroutine provides (dN/dy)/N
example: exponential profile on top of constant
N=N0 + N1exp(-y/delta)
dummy=an1*exp(-y/delta)
res=(-dummy/delta)/(an0+dummy)
return
end
=====
This program implements the method of moments
for a straight wire on potential 1
The charge density is taken to be piecewise constant
and a point-matching technique is used
program wire
real*4 a(100,100),x(100),r(100)
a: matrix of the problem
x: unknown charge density
(normalised on a factor 4*pi*epsilon0)
r: right-hand member (=1 in this particular case)
Data input

```

```

210
1 write(6,1)
format(' Number of divisions on the wire')
read(5,*)
ndiv
if(ndiv.gt.100) goto 200
C
2 write(6,2)
format(' Total length of the wire')
read(5,*)
tl
C
3 write(6,3)
format(' Radius of the wire')
read(5,*)
rad
C
Elementary division
C
del=tl/ndiv
C
diagonal elements i=j
C
do 10 i=1,ndiv
a(i,i)=2.*alog(del/rad)
continue
C
non-diagonal elements
C
do 20 i=1,ndiv
do 20 j=1,ndiv
if(i.eq.j) goto 20
a(i,j)=1./iabs(i-j)
continue
C
right-hand member
C
do 30 i=1,ndiv
r(i)=1.
continue
C
Solution of the system of equations
C
call eqn(a,r,x,ndiv)
C
Write the final solution to a file result.dat
C
We now introduce the factor 4*pi*epsilon0 = (10**-9)/9.
The results in RESULT.DAT are expressed in Coulomb*(10**-13),
hence the multiplication factor becomes (10**4)/9.
fak=1.e4/9.
open(unit=1,file='result.dat')
write(1,*)
(fak*(i),i=1,ndiv)
close(unit=1)
C
Calculate the total charge on the wire
totchar=0.
do 500 i=1,ndiv
totchar=totchar+x(i)
continue
totchar=totchar*fak*del
write(6,501)
totchar
format(' The total charge is: ',620.6,' 10**-13 Coulomb')
stop
C
To many unknowns
C

```

```

200 write(6,66)
66 format(' Number of unknowns is limited to 100! ')
      goto 210
C
end
C=====
subroutine eqn(a,r,x,n)
C
Simple routine to solve linear system of equations
r(i)=Aij*xj
n:number of equations
maximum number is now 100
C
real*4 a(100,100),x(100),r(100)
C
do 60 k=1,n-1
do 10 i=k+1,n
  r(i)=a(k,k)*r(i)-r(k)*a(i,k)
do 10 j=k+1,n
  a(i,j)=a(k,k)*a(i,j)-a(i,k)*a(k,j)
10 continue
top=0.
do 70 i=k+1,n
do 70 j=k+1,n
  if(abs(a(i,j)).gt.top) top=abs(a(i,j))
70 continue
do 80 i=k+1,n
  r(i)=r(i)/top
do 80 j=k+1,n
  a(i,j)=a(i,j)/top
80 continue
60 continue
C
elimination finished - calcute unknowns
C
x(n)=r(n)/a(n,n)
do 30 k=1,n-1
do 20 j=1,k
  r(n-k)=r(n-k)-a(n-k,n-j+1)*x(n-j+1)
20 continue
x(n-k)=r(n-k)/a(n-k,n-k)
30 continue
C
return
end
C=====
program stubs
C
This program solves the double stub matching problem
It starts a given real characteristic impedance and a
given load impedance.
The distance between this load and the first stub and the
distance between the stubs must also be supplied to
the program
A proper warning is displayed if no solution turns out to be
possible
C
write(6,1)
1 format(' Characteristic impedance (real):')
read(5,*) r0
write(6,2)
2 format(' Real and imaginary part of load: ')
read(5,*) rr,xr
write(6,3)
3 format(' Distance from load to first stub: ')
read(5,*) an
C
100
4 write(6,4)
format(' Distance between the stubs: ')
read(5,*) anb
sig=1.
call double(r0,rr,xr,an,anb,wav11,wav12,sig)
write(6,5)
5 format(' First solution: ')
if(sig.ne.0.) write(6,102) wav11,wav12
if(sig.eq.0.) write(6,101)
format(' No solution possible! ')
sig=-1.
call double(r0,rr,xr,an,anb,wav11,wav12,sig)
write(6,6)
6 format(' Second solution: ')
if(sig.ne.0.) write(6,102) wav11,wav12
if(sig.eq.0.) write(6,101)
format(' First stub: ',G20.6,' Second stub: ',G20.6)
stop
end
C=====
subroutine double(r0,rr,xr,an,anb,wav11,wav12,sig)
wav11=0.
wav12=0.
pi=3.141592654
theta=anb*720.
call rangle(theta)
call reflect(r0,0.,rr,xr,gm,gad)
gad=gad+180.
call rangle(gad)
call spiral(gm,gad,an,0.,1.,gm,gad)
call imped(1.,0.,gm,gad,r1,x1)
thetar=theta*pi/180.
d=r1*(1.-cos(thetar))
if(d.gt.2..or.d.eq.0.) goto 100
a=2.*r1/(r1+1.)-cos(thetar)
b=sin(thetar)
d=r1/(1.+r1)
e=1./(1.+r1)
c=d-e
aa=a**2+b**2
bb=2.* (b**2*d+a*d-a*e)
cc=b**2*d**2+d**2-2.*d*k*+e**2-(b*e)**2
disc=bb**2-4*a**2*cc
if(disc.le.0..or.b.eq.0.) disc=0.
x=(bb+sig*sqrt(disc))/(2.*aa)
if(disc.ne.0.) y=(a*x-c)/b
if(disc.eq.0.) y=sig*sqrt(abs(e**2-(x-d)**2))
phir=atan2(y,x)
phiphi=phir*180./pi
gm=x/cos(phiphi)
call imped(1.,0.,gm,phi,r1,x1)
bl=x12-x1
call reflect(1.,0.,0.,bl,gml,ang)
if(ang.lt.0.) ang=360.+ang
wav11=(360.-ang)/720.
call spiral(gm,phi,anb,0.,1.,gm,gad)
if(gad.lt.-90.) gad=-180.-gad
call imped(1.,0.,gm,gad,r1,x1)
bl=-x1
call reflect(1.,0.,0.,bl,gm2,ang)
if(ang.lt.0.) ang=360.+ang
wav12=(360.-ang)/720.
return
sig=0.
return
end

```

```

C----- subroutine rectan(ra,thetad,x,y)
C
C pi=3.141592654
C thetad=thetad*pi/180.
C x=ra*cos(thetad)
C y=ra*sin(thetad)
C return
C end
C----- subroutine imped(r0,x0,gm,gad,rs,xs)
C
C call rectan(gm,gad,gr,gi)
C call polar(1.+gr,gi,an,anad)
C call polar(r0,x0,z0m,z0ad)
C call polar(1.-gr,-gi,ad,adad)
C call rectan(z0m*an/ad,z0ad+anad-acad,rs, xs)
C return
C end
C----- subroutine reflect(r0,x0,rr,xr,gm,gad)
C
C call polar(rr-r0,xr-x0,an,anad)
C call polar(rr+r0,xr+x0,ad,adad)
C gm=an/ad
C gad=anad-adad
C call rangle(gad)
C return
C end
C----- subroutine spiral(gm1,gad1,an,alpha,beta,gm2,gad2)
C
C dgad=720.*an
C gad2=gad1-dgad
C call rangle(gad2)
C pi=3.141592654
C attn=-4.*pi*alpha*an/beta
C gm2=gm1*exp(attn)
C return
C end
C----- subroutine rangle(thetad)
C
C 1 if(thetad.lt.360.) goto 2
C thetad=thetad-360.
C goto 1
C 2 if(thetad.ge.0.) goto 3
C thetad=thetad+360.
C goto 2
C 3 if(thetad.gt.180.) thetad=thetad-360.
C return
C end
C----- subroutine polar(x,y,ra,thetad)
C
C ra=sqrt(x*x+y*y)
C thetad=atan2(y,x)*180./3.141592654
C return
C end
C----- program findif
C
C This program solves Laplaces equation for the potential
C by means of finite differences
C
C It is written here to solve the symmetric stripline problem i.e.

```

a perfectly conducting strip of width bs (in arbitrary units) between two perfectly conducting ground planes. The distance between these planes and the strip is hsg (in arbitrary units)

In order to solve this problem the configuration is put into a box of potential zero. The boundaries to the left and to the right of the strip are specified by a distance bz (arbitrary units). This distance must be large enough to yield good results. The number of divisions on the strip and between the strip and the planes must also be large enough (hsg and bs must be > 15). We also determine the impedance for that configuration. The medium between the planes is air.

Typical example: strip width $bs=15$
 distance between strip and ground $hsg=15$
 distance between strip and box $bz=60$

This is a strip where the width is half the distance between the planes. Characteristic impedance: 100 Ohm (see literature)
 numerical result after about 500 iterations: 96.6 Ohm

```

real*4 v(202,202)
integer*4 bs,hsg,hs,bz

Data input

write(6,1)
format(' Strip width in units:')
read(5,*), bs
write(6,2)
format(' Distance above the ground:')
read(5,*), hsg
hs=hsg-1
write(6,4)
format(' Distance left and right of strip:')
read(5,*), bz

ista=0
600 write(6,5)
format(' Number of cycles [0=stop],number of iter./cycle:')
read(5,*), ncy,niter
if(ncy.eq.0) goto 888

icy=0
iter=0
if(ista.ne.0) goto 200

initialise

nx=bs+2*bz+2+1
ny=hsg+hs+1+1
do 20 i=1,nx
do 20 j=1,ny
v(i,j)=0.

20 continue
iy=hsg+1
do 21 i=bz+2,bz+bs+2
v(i,iy)=1.
21 continue

Start iterations

200 do 22 j=2,ny-1
do 22 i=2,nx-1
if(j.ne.iy) goto 50
if((i.lt.bz+2).or.(i.gt.bz+bs+2)) goto 50
v(i,j)=1.
50 continue
22 continue

```

```

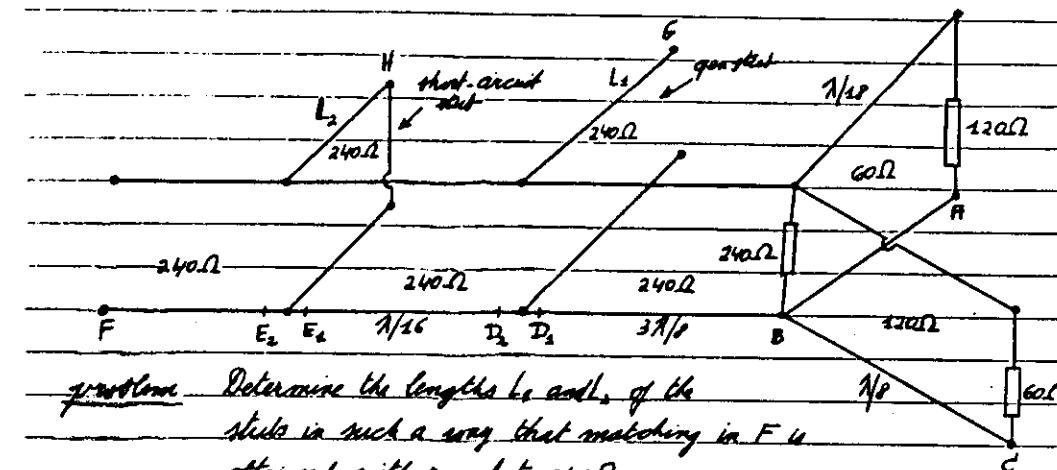
50      v(i,j)=(v(i-1,j)+v(i+1,j)+v(i,j+1)+v(i,j-1))/4.
22      continue
iter=iter+1
if(iter.ne.niter) goto 200
C
100    if(ista.ne.0) goto 100

Calculation of the total capacitance with respect to the
plane below the strip
C
301    write(6,301)
format(' Capacitance calculation')
C
Numerical calculation of the capacitance
C
100    ista=1
idis=2
sum=0.
100    iy=hsg+1+idis
ibegin=bz+2-dis
iend=bz+bs+2+idis
do 303 i=ibegin,iend
fak=1.
if((i.eq.ibegin).or.(i.eq.iend)) fak=0.5
sum=sum+fak*(v(i,iy+1)-v(i,iy-1))
303    continue
iy=hsg+1-dis
do 304 i=ibegin,iend
fak=1.
if((i.eq.ibegin).or.(i.eq.iend)) fak=0.5
sum=sum-fak*(v(i,iy+1)-v(i,iy-1))
304    continue
ix=bz+2-dis
ibegin=hsg+1-dis
iend=hsg+1+idis
do 305 i=ibegin,iend
fak=1.
if((i.eq.ibegin).or.(i.eq.iend)) fak=0.5
sum=sum+fak*(v(ix-1,i)-v(ix+1,i))
305    continue
ix=bz+bs+2+idis
do 306 i=ibegin,iend
fak=1.
if((i.eq.ibegin).or.(i.eq.iend)) fak=0.5
sum=sum-fak*(v(ix-1,i)-v(ix+1,i))
306    continue

sum=0.5*sum
sum=(1000./(36.*3.141592654))+sum
write(6,600) -sum
format(' The numerical calculation yields: ',620.6)
C
600    icy=icy+1
if(icy.eq.ncy) goto 800
iter=0
goto 200
C
800    write(6,809)
format(' Impedance of the stripline configuration:')
write(6,*)
stop
end
C=====

```

Typical double stub matching problem



problem Determine the lengths l_1 and l_2 of the stubs in such a way that matching in F is obtained with respect to 240Ω .

solution (see also on Smith chart)

$$1. \quad Z_A = 120\Omega \Rightarrow Z'_A = Z_A/R_0 = 2 \Rightarrow Y'_A = 1/2$$

$$2. \quad B \rightarrow A: \quad R_B = 1/10 \text{ or } 40^\circ; \text{ consequently } Y_{BA} = 0.55 + j0.275$$

$$Y_{BA} = (0.55 + j0.275)/60\Omega$$

$$3. \quad Z_B = 60\Omega \Rightarrow Z'_B = 0.5 \Rightarrow Y'_B = 2$$

$$4. \quad C \rightarrow B: \quad R_C = 1/10 \text{ or } 90^\circ; \quad Y_{CB} = 0.1 - j0.6$$

$$Y_{CB} = (0.1 - j0.6)/120\Omega$$

5. situation in B:

$$\begin{aligned} Y_{\text{total}} &= Y_0 + Y_{AB} + Y_{CB} \\ &= 240\Omega \quad | \quad Y_0 \quad | \quad Y_{AB} \quad | \quad Y_{CB} \quad | \quad Y_{\text{total}} = Y_{\text{total}} * 240\Omega \\ &= 4 \times (0.55 + j0.275) \\ &\quad + 2 \times (0.1 - j0.6) \\ &= \frac{1}{Z_0} = \frac{1}{240\Omega} \\ &= 1 \\ &= 4 \Omega = 0.1j \end{aligned}$$

$$6. B \rightarrow D_2 : BD_2 = 3\sqrt{2} \text{ or } 270^\circ ; Y'_{D_2} = 0.4 + j0.925$$

We now apply the matching procedure

step 1 : the circle with real part = 1 must be rotated in the towards load direction over $E_2 D_2 = \pi/16$ or 45° degrees (see Smith chart)

step 2 : going from D_2 to D_1 only changes the imaginary part of the admittance

whatever the length L_2 , going from D_2 to D_1 we stay on a circle with real part = 0.4

step 3 : we determine the intersection of the circles from step 1 and step 2; we have 2 solutions:

$$Y'_{D_2}^{(1)} = 0.4 + 0.8j \Rightarrow \text{admittance } B_2^{(1)N} \text{ of stat 1} = -0.125j$$

$$Y'_{D_2}^{(2)N} = 0.4 + 4j \Rightarrow \text{admittance } B_2^{(2)N} \text{ of stat 1} = 3.075j$$

$$\text{step 4 : } D_2 E_2 = \pi/16 \text{ or } 45^\circ \Rightarrow Y'_{E_2}^{(1)N} = 1 + 1.575j$$

$$Y'_{E_2}^{(2)N} = 1 - 6.4j$$

$$\text{step 5 : } Y'_{E_2} = 1 \text{ (matching)} \Rightarrow \text{admittance } B_2^{(1)N} \text{ of stat 2} = -1.575j$$

$$\text{admittance } B_2^{(2)N} \text{ of stat 2} = 6.4j$$

step 6 : length of the stubs

stub 1 starts in E ; open stub $\Rightarrow Y' = 0$ (left hand side of Smith chart)

$B_2^{(1)N} = -0.125j$ conjugate with $0.4j1$ (from E to $L_2^{(1)N}$ on Smith chart)

$B_2^{(2)N} = 3.075j$ conjugate with $0.2j1$ ($E \rightarrow L_2^{(2)N}$)

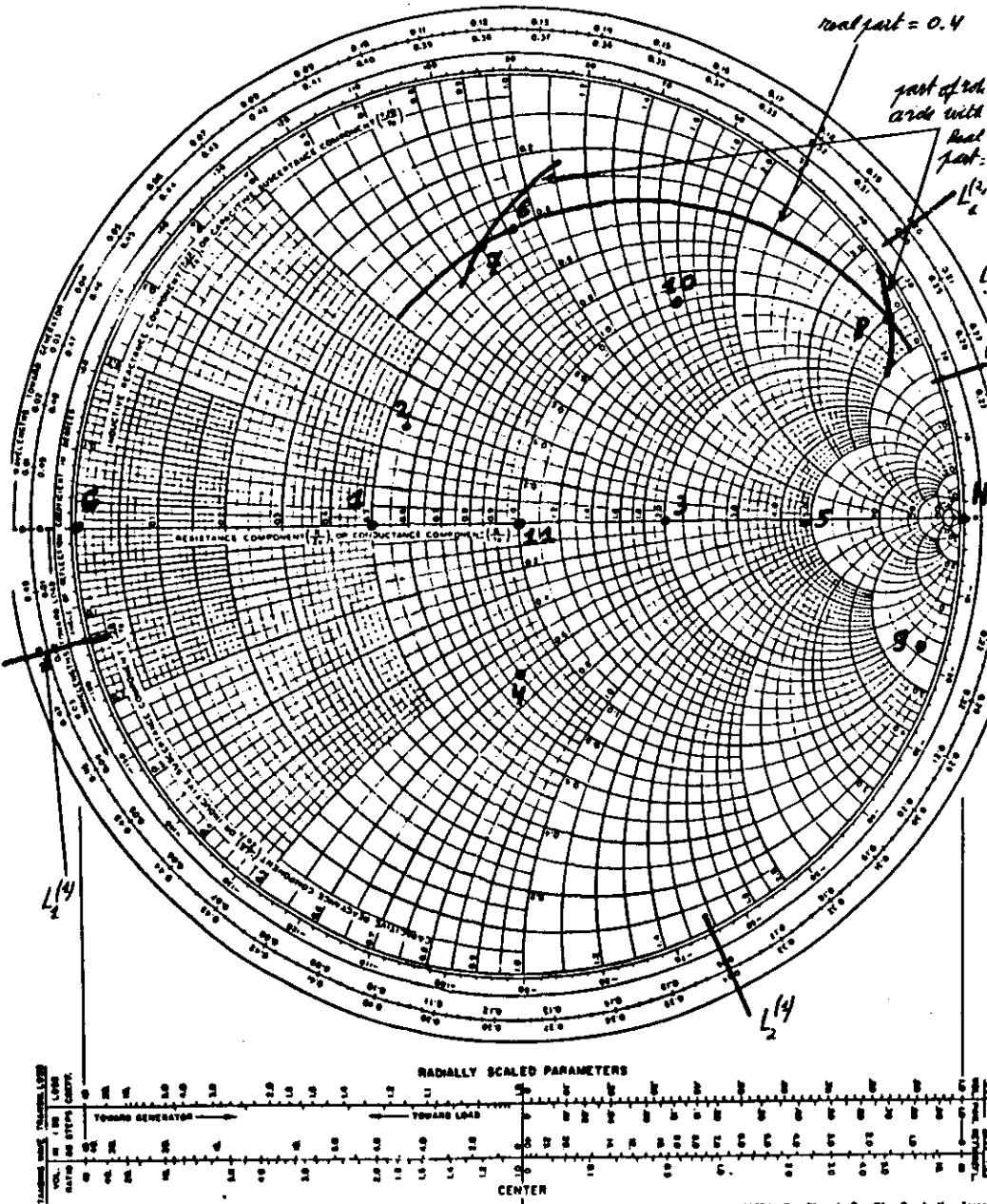
stub 2 starts in H ; short-circuited stub $\Rightarrow Y' = \infty$ (right hand side of Smith chart)

$B_2^{(1)H} = -1.575j$ conjugate with $0.09j1$ ($H \rightarrow L_2^{(1)H}$)

$B_2^{(2)H} = 6.4j$ conjugate with $0.476j1$ ($H \rightarrow L_2^{(2)H}$)

NAME	TITLE	DOUBLE STUB MATCHING	DWG. NO.
SMITH CHART Form 5301-7560-N	GENERAL RADIO COMPANY, WEST CONCORD, MASSACHUSETTS		DATE

IMPEDANCE OR ADMITTANCE COORDINATES



Electronics Vol. 17, No. 1, pp. 150-153, 310-323, Jan. 1944

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- 1: Y'_H 2: Y'_{AB} 3: Y'_C 4: Y'_{CB} 5: Y'_{total} 6: Y'_D
- 7: Y'_{D_2} 8: $Y'_{D_2}^{(1)}$ 9: $Y'_{E_2}^{(1)N}$ 10: $Y'_{E_2}^{(2)N}$ 11: $Y'_{E_2} = 1$

NAME

TITLE

DATE

SMITH CHART FORM S301-7560-N GENERAL RADIO COMPANY, WEST CONCORD, MASSACHUSETTS

IMPEDANCE OR ADMITTANCE COORDINATES

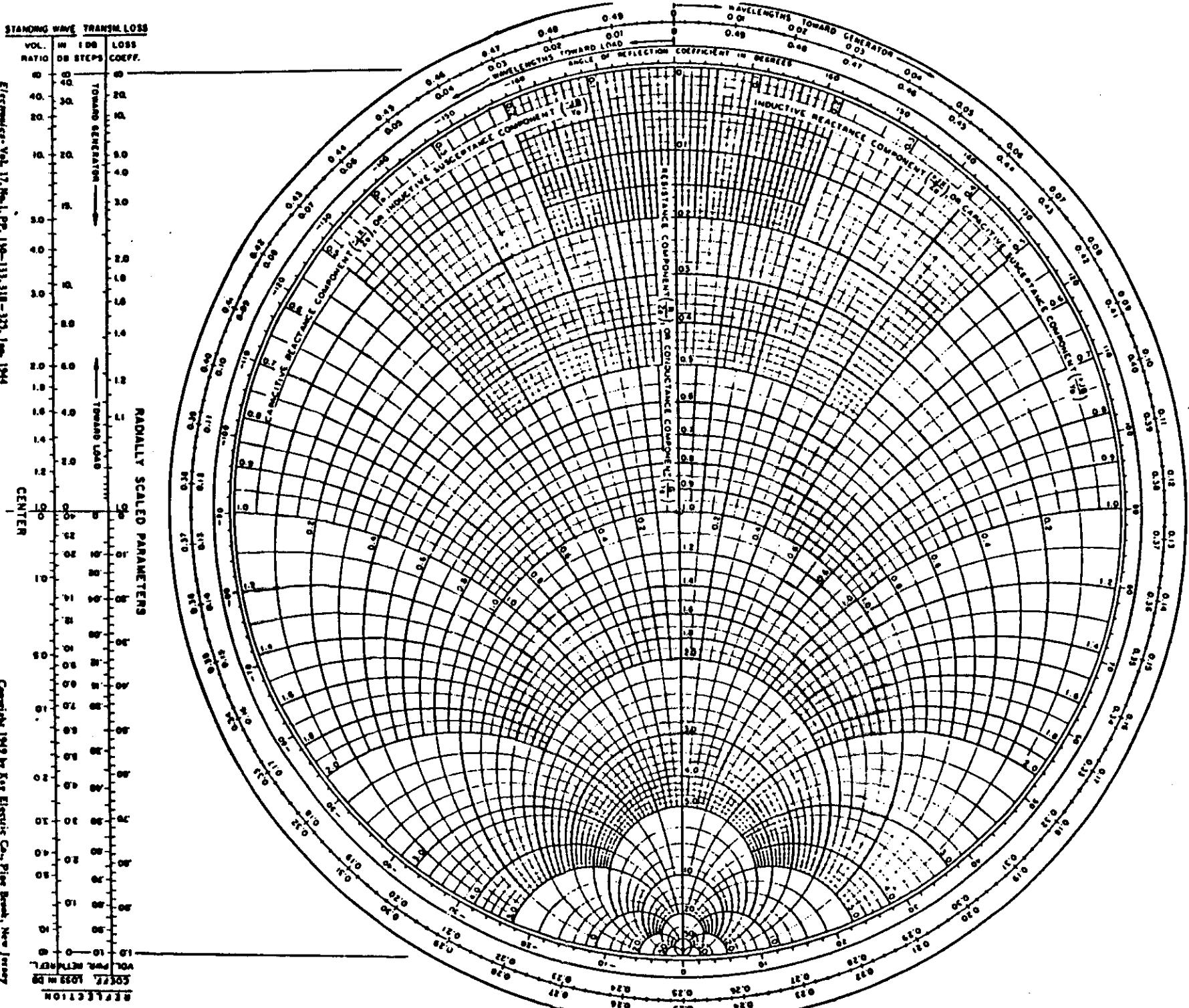


Table 1. Basic program for solution of double-stub tuners

```

1 SUBROUTINE DOUBLE (R0, PR, XPLA, GAM, WVL1, WVL2, SIG)
2   C DETERMINES THE LENGTHS (IN WAVELENGTHS) OF
3   C THE TWO SHORT CIRCUIT STUBS FOR A DOUBLE-STUB TUNER.
4   C
5   C INPUTS:
6   C   R0 - CHARACTERISTIC IMPEDANCE (REAL),
7   C   PR - REAL PART OF LOAD IMPEDANCE,
8   C   XPLA - IMAGINARY PART OF LOAD IMPEDANCE,
9   C   GAM - DISTANCE (IN WAVELENGTHS) FROM THE LOAD
10  C TO THE FIRST STUB.
11  C
12  C AN = DISTANCE (IN WAVELENGTHS) FROM THE FIRST
13  C STUB TO THE SECOND STUB.
14  C SIG = 1... GIVES ONE VALID SOLUTION,
15  C SIG = -1... GIVES ANOTHER VALID SOLUTION.
16  C
17  C OUTPUTS:
18  C   WVL1= LENGTH OF FIRST STUB (IN WAVELENGTHS),
19  C   WVL2= LENGTH OF SECOND STUB (IN WAVELENGTHS),
20  C   WVL1= WVL2=0.
21  C
22  C PI=4.ATAN(1)
23  C CALL RANGLE(THETA)
24  C ENTER SMITH CHART AT LOAD
25  C CALL REFLECT (R0, PR, GAM, GAM)
26  C ROTATE TO LOAD ADmittance.
27  C GAM=GAM+180.
28  C ROTATE ANWAVELENGTHS CLOCKWISE ON CONSTANT SWR
29  C CIRCLE.
30  C CALL SPIRAL(GAM,GAM,0.1,0,0,1,0,0)
31  C OBTAIN NORMALIZED ADMITTANCE AT FIRST STUB.
32  C CALL IMPED(1,0,0,1,0,0,0,1,0,0)
33  C TEST INPUT DATA TO SEE IF A SOLUTION IS POSSIBLE.
34  C D=R0*1.0-COS(THETA)
35  C FIDGE 2.0,0 DECIMAL PLACES
36  C FIDGE 2.0,0 DECIMAL PLACES
37  C
38  C SOLVE SIMULTANEOUS EQUATIONS FOR TWO CIRCLES.
39  C
40  C PI=2*ATAN(1)-COS(THETA)
41  C PI=PI-ATAN(1)
42  C D=PI*1.0-COS(THETA)
43  C FIDGE 2.0,0 DECIMAL PLACES
44  C CD=PI-PI
45  C CD=E
46  C CD=A-E*B^2
47  C
48  C
49  C
50  C
51  C SUBROUTINE DOUBLE 737CC OPT=0 TRACE DEBUG FTM4,64-629
52  C GAM=XCOSSOPH1
53  C CALL IMPED(1,0,0,0,GAM,PHI,RL,XL2)
54  C IF(DSC=.0-.4*ACCC)
55  C IF(DSC=.0-.4*ACCC)
56  C IF(DSC=.0-.4*ACCC)
57  C IF(DSC=.0-.4*ACCC)
58  C IF(DSC=.0-.4*ACCC)
59  C
60  C OBTAIN LENGTH OF STUB NEAREST LOAD --- WVL1.
61  C BL=XL-AL
62  C CALL REFLECT(1,0,0,BL,GAM1,ANG)
63  C FIDSC(NORM)*(X-A)*COS(PI*BL/WAVL1)
64  C FIDSC(NORM)*(X-A)*COS(PI*BL/WAVL1)
65  C
66  C PH1=ATAN(Y1*BL/WAVL1)
67  C
68  C ROTATE ANWAVELENGTHS CLOCKWISE ON CONSTANT SWR
69  C CIRCLE.
70  C CALL SPIRAL (GAM, PH1, AMB, 0, 1, 0, 0, GAM)
71  C FIDGE 1.7, 0, 0, 0, 0, 0, 0, 0, 0, 0
72  C OBTAIN LENGTH OF STUB FARTHEST FROM LOAD --- WVL2.
73  C CALL IMPED(1,0,0,0,GAM,PH1,RL,XL)
74  C
75  C BL=XL
76  C CALL REFLECT(1,0,0,BL,GAM2,ANG)
77  C FIDMOL(T,DUMG1,-300+ANG)
78  C WVL1=WVL1+ANG1/720.
79  C
80  C RETURN

```

Table 2: Subroutines called for in DOUBLE program

```

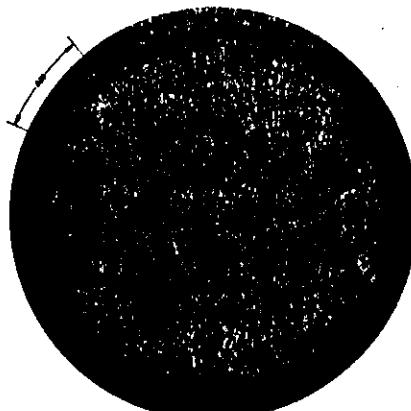
SUBROUTINE RECT2POLA(GAM,AD,AD,X,Y)
1  C CHANGE FROM POLAR TO RECTANGULAR COORDINATES
2  C
3  C INPUT PARAMETERS:
4  C   GAM=MAGNITUDE
5  C   AD=ANGLE
6  C
7  C OUTPUT PARAMETERS:
8  C   X=ABSCISSA
9  C   Y=ORDINATE
10  C
11  C X=AD*COS(AD)
12  C Y=AD*SIN(AD)
13  C
14  C RETURN
15  C
16  C
17  C SUBROUTINE IMPERPOL(GAM,GAM,PR,XS)
18  C LEAVE SMITH CHART BY CHANGING FROM REFLECTION
19  C COEFFICIENT TO IMPEDANCE
20  C
21  C INPUT PARAMETERS:
22  C   R0=REAL PART OF CHARACTERISTIC IMPEDANCE
23  C   GAM=IMAGINARY PART OF CHARACTERISTIC IMPEDANCE
24  C   GAM=ANGLE OF REFLECTION COEFFICIENT IN DEGREES
25  C
26  C OUTPUT PARAMETERS:
27  C   R0=REAL PART OF SENDING-END IMPEDANCE
28  C   GAM=IMAGINARY PART OF SENDING-END IMPEDANCE
29  C
30  C CALL RECT2POLA(GAM,GR,GR)
31  C CALL POLAR2P(XS,Z0,Z0,AD)
32  C CALL POLAR2P(0,Z0,Z0,AD)
33  C CALL RECT2P(Z0,AD,AD,AD,AD,PR,XS)
34  C
35  C RETURN
36  C
37  C
38  C SUBROUTINE REFLECT(AD,PR,30,GR,GR)
39  C ENTER SMITH CHART BY OBTAINING REFLECTION COEFFICIENT
40  C
41  C INPUT PARAMETERS:
42  C   R0=REAL PART OF CHARACTERISTIC IMPEDANCE
43  C   GAM=IMAGINARY PART OF CHARACTERISTIC IMPEDANCE
44  C   PR=REAL PART OF RECEIVING-END IMPEDANCE
45  C   X0=IMAGINARY PART OF RECEIVING-END IMPEDANCE
46  C
47  C OUTPUT PARAMETERS:
48  C   GAM=MAGNITUDE OF REFLECTION COEFFICIENT
49  C   AD=ANGLE OF REFLECTION COEFFICIENT IN DEGREES
50  C
51  C BETWEEN -180 and 180
52  C
53  C CALL POLAR2P(-PR,30,X0,AN,AD)
54  C CALL POLAR2P(+PR,X0+AD,AD,AD)
55  C CALL POLAR2P(GM,AN,AD,AD)
56  C CALL RANGLE(GAM)
57  C
58  C RETURN
59  C
60  C
61  C SUBROUTINE RANGLE(THETA)
62  C
63  C INPUT PARAMETERS:
64  C   GAM=ANGLE OF REFLECTION COEFFICIENT IN DEGREES
65  C   GAM=ANGLE OF REFLECTION COEFFICIENT IN DEGREES
66  C
67  C BETWEEN -180 and 180
68  C
69  C CALL POLAR2P(-PR,30,X0,AN,AD)
70  C CALL POLAR2P(+PR,X0+AD,AD,AD)
71  C CALL POLAR2P(GM,AN,AD,AD)
72  C CALL RANGLE(GAM)
73  C
74  C RETURN
75  C

```

```

SUBROUTINE SPIRAL(GAM1,AD1,AN,ALPHA,BETA,GR2,GR2)
1  C MOVES OPERATING POINT ALONG SPIRAL IN SMITH CHART FROM
2  C STARTING REFLECTION COEFFICIENT TO FINISHING COEFFICIENT
3  C
4  C INPUT PARAMETERS:
5  C   GAM1= MAGNITUDE OF STARTING REFLECTION COEFFICIENT
6  C   AD1= ANGLE IN DEGREES OF STARTING REFLECTION COEFFICIENT
7  C   GR1= FLOATER-1
8  C
9  C   AN= LENGTH OF LINE IN WAVELENGTHS
10  C   --POSITIVE FOR MOVEMENT TOWARD GENERATOR
11  C   --NEGATIVE FOR MOVEMENT TOWARD LOAD
12  C   ALL= ALPHA ATTENUTATION CONSTANT
13  C   BETA= PHASE CONSTANT
14  C   GAM2= MAGNITUDE OF FINISHING REFLECTION COEFFICIENT
15  C   GAM2= ANGLE (DEGREES) OF FINISHING REFLECTION COEFFICIENT
16  C   BETWEEN -180 AND 180
17  C
18  C DOAD=-720,100
19  C GR2=GAM1-DOAD
20  C CALL RANGLE(40,DOAD)
21  C CALL POLAR2P(X,GR2,AD2)
22  C PR=1.0/SQRT(X*X+GR2*GR2)
23  C ATTN=1.0/(ALPHA*AM(BETA))
24  C GAM2=GR2*EXP(ATTN)
25  C
26  C RETURN
27  C
28  C SUBROUTINE RANGLE(THETA)
29  C
30  C INPUT:
31  C   THETA= ANY ANGLE IN DEGREES
32  C
33  C OUTPUT:
34  C   THETAD= SAME ANGLE CONVERTED TO RANGE
35  C   -180LTTHETADLT+180
36  C
37  C SUBROUTINE 1,30,100 TO 2
38  C THETAD=THETA-360
39  C GO TO 1
40  C ATTHETAD GE 0, GO TO 3
41  C
42  C GO TO 2
43  C
44  C SUBROUTINE 0,180,100,THETAD=THETAD-360
45  C
46  C RETURN
47  C
48  C
49  C SUBROUTINE POLAR2P(X,Y,THETA)
50  C
51  C CHANGE FROM RECTANGULAR TO POLAR COORDINATES
52  C
53  C INPUT PARAMETERS:
54  C   X=ABSCISSA
55  C   Y=ORDINATE
56  C
57  C OUTPUT PARAMETERS:
58  C   GAM=MAGNITUDE
59  C   THETAD=ANGLE IN DEGREES BETWEEN -180 AND 180
60  C
61  C PR=SQRT(X*X+Y*Y)
62  C GAM=SIGN(X)*Y/PR
63  C THETAD=ATAN(Y/X)
64  C
65  C RETURN

```



2. If the circle of constant susceptance passing through Y_s/Y_0 doesn't intercept circle A, the computer reports that a solution is impossible.



3. The circle A intercepts Y_s/Y_0 and Y_s'/Y_0 , which are the two possible solutions. Subtracting Y_s/Y_0 from Y_s'/Y_0 provides WAVL1.

$Y_s/Y_0 = 1$. (line 62). Finally, also find the required wavelength WAVL2 of this stub by the single-stub method (lines 63 to 66).

How to use subroutine DOUBLE

The input and output arguments for subroutine DOUBLE are explained in the comment cards of the listing in Table 1. When you use SIG = 1.0 for the last argument of the subroutine, you will obtain solution #1 for WAVL1 and WAVL2, and when you use SIG = -1.0 for this argument, you will obtain the other possible solution. When first using subroutine DOUBLE, it might be wise to add print statements after statements 25, 29, 61, and 61. In this way, you can verify the entire analysis by plotting these points.

As an example of the use of DOUBLE, consider the problem where $Z_0 = 100 \Omega$, $Z_s = 200 - j100 \Omega$, the distance from load to the first stub is 0.05 wavelengths, and the separation of the stubs is 0.80 wavelengths. In the main program, use the statement:

CALL DOUBLE (100, 200, -100, 0.05, 0.80, WAVL1,
WAVL2, 1.0)

and then print out the values of WAVL1 and WAVL2. This constitutes one of two possible solutions. If you call DOUBLE again with -1.0 instead of 1.0 for the last argument, you obtain the other possible solution.

1. WAVL1 = 0.1005, WAVL2 = 0.0970
2. WAVL1 = 0.2084, WAVL2 = 0.8557

For another design problem, consider the case where the frequency varies over a given range. In this case, the arguments XR, AN, and ANB take on values that depend on the frequency. The CALL DOUBLE statement and the print statements can be put in the same Do-loop that defines XR, AN, and ANB. In this way, you obtain hundreds of solutions in a matter of seconds.

For a final example, consider a case that is difficult and inaccurately solved by the usual Smith-chart method. Suppose $Z_0 = 600 \Omega$, $Z_s = 1000 + j500 \Omega$, AN = 0.125, and



4. Moving Y_s/Y_0 clockwise along a circle of constant VSWR for ANB wavelength provides Y_s'/Y_0 . The admittance of the second stub is found by $(Y_s'/Y_0 - 1)$.

ANB = 0.256. Carrying out the Smith-chart solution, we find that the two crossings of the A circle with the b circle are close together and near the center of the chart. The computer program always gives the same results accurate to as many places as desired.

1. WAVL1 = 0.1202, WAVL2 = 0.2400
2. WAVL1 = 0.1250, WAVL2 = 0.2500

The author thanks his students who patiently tested subroutine for all possible variations of inputs.

References
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 3. D. W. Darmatoff and W. S. McCallum, *Electromagnetic Wave Propagation*, McGraw-Hill Book Company, New York, (1969).

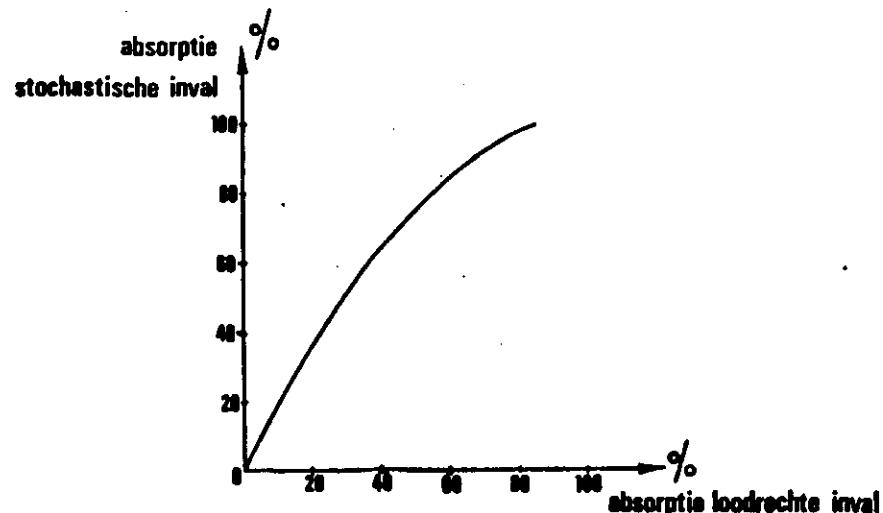


Fig 1

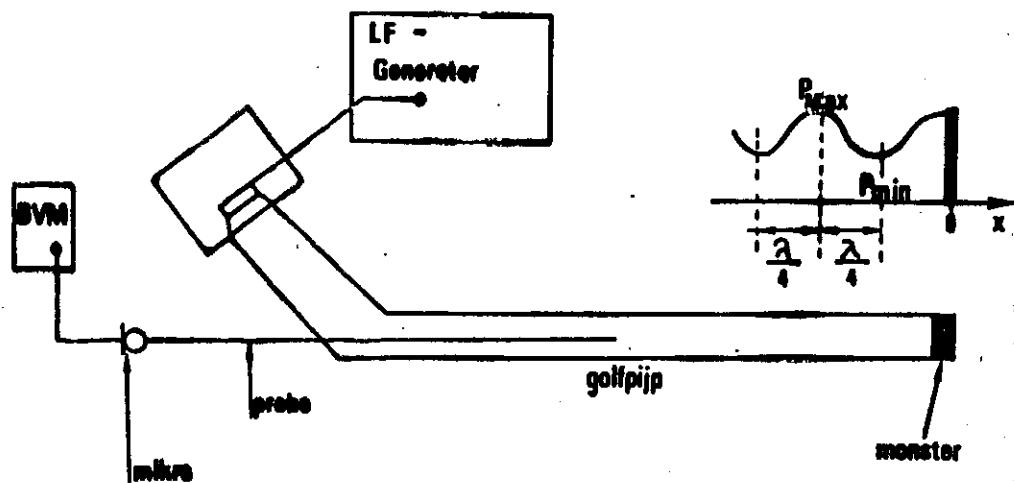


Fig 2