

**MICROELECTRONIC ENGINEERING
ROCHESTER INSTITUTE OF TECHNOLOGY**

A Review of IC Fabrication Technology

Dr. Lynn Fuller

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Microelectronic Engineering

Rochester Institute of Technology

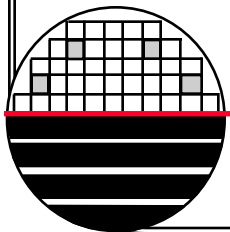
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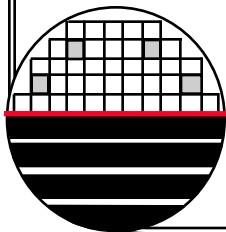
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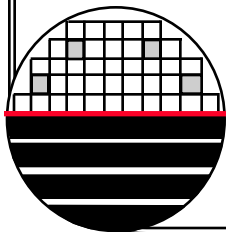
ADOBE PRESENTER

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OUTLINE

- § **Constants**
- § **Periodic Table**
- § **Material Properties**
- § **Oxide Growth**
- § **Diffusion**
- § **Resistivity, Sheet Resistance, Resistance**
- § **Mobility**
- § **pn Junction**
- § **MOSFET V_t**
- § **Ion Implantation**
- § **Conclusion**



CONSTANTS

Electronic charge	q	1.602 E -19 Coulomb
Speed of light in vacuum	c	2.998E8 m/s
Permittivity of vacuum	ϵ_0	8.854 E -14 F/cm
Free electron Mass	m_0	9.11E-31 Kg
Planck constant	h	6.625E-34 J s
Boltzmann constant	k	1.38 E-23 J /°K = 8.625E-5 eV/°K
Avogadro's number	A_0	6.022E23 molecules/gm- mole
Thermal voltage	kT/q	@ 300 °K = 0.02586

PLAY

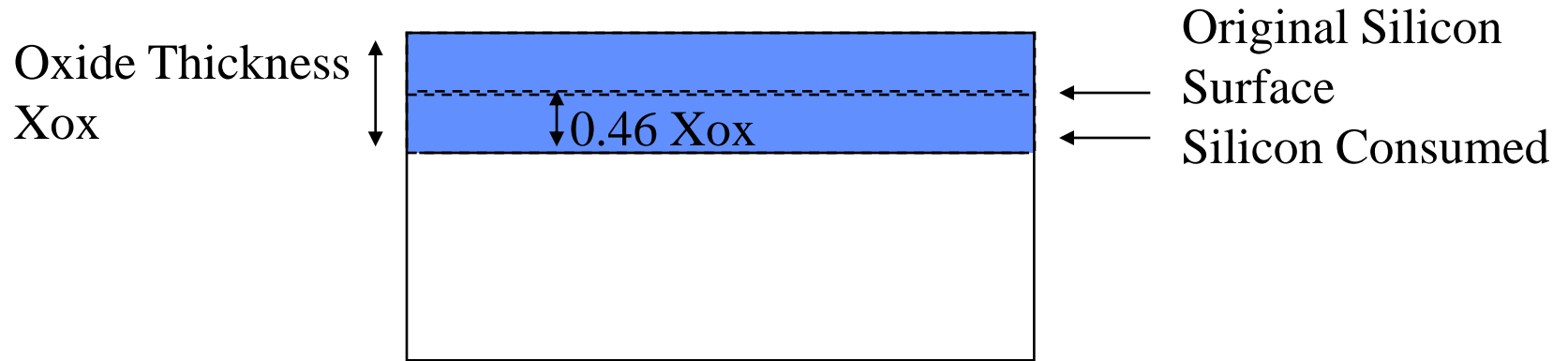
PERIODIC TABLE OF THE ELEMENTS

1 1.0079 H 0.0899 Hydrogen	<div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;"> <p>Atomic Number</p> <p>Density g/cm³</p> </div> <div style="border: 1px solid black; padding: 5px;"> <p>14 20.086</p> <p>Si</p> <p>2.33 Silicon</p> </div> <div style="text-align: center;"> <p>Atomic Weight</p> <p>Symbol</p> <p>Name</p> </div> </div>																2 4.0026 He 0.1787 Helium
3 6.941 Li 0.53 Lithium	4 9.01218 Be 1.85 Beryllium											5 10.81 B 2.34 Boron	6 12.011 C 2.62 Carbon	7 14.0067 N 1.251 Nitrogen	8 15.9994 O 1.429 Oxygen	9 18.9984 F 1.696 Fluorine	10 20.179 Ne 0.901 Neon
11 22.9898 Na 0.97 Sodium	12 24.305 Mg 1.74 Magnesium											13 26.9815 Al 2.70 Aluminum	14 20.086 Si 2.33 Silicon	15 30.97376 P 1.82 Phosphorus	16 32.06 S 2.07 Sulfur	17 35.453 Cl 3.17 Chlorine	18 39.948 Ar 1.784 Argon
19 39.0983 K 0.86 Potassium	20 40.08 Ca 1.55 Calcium	21 44.9559 Sc 3.0 Scandium	22 47.90 Ti 4.50 Titanium	23 50.941 V 5.8 Vanadium	24 51.996 Cr 7.19 Chromium	25 54.938 Mn 7.43 Manganese	26 55.847 Fe 7.86 Iron	27 58.9332 Co 8.90 Cobalt	28 58.70 Ni 8.90 Nickel	29 63.546 Cu 8.96 Copper	30 65.238 Zn 7.14 Zinc	31 69.72 Ga 5.98 Gallium	32 72.59 Ge 5.32 Germanium	33 74.9216 As 5.72 Arsenic	34 78.96 Se 4.80 Selenium	35 79.904 Br 3.12 Bromine	36 83.80 Kr 3.74 Krypton
37 85.468 Rb 1.53 Rubidium	38 87.62 Sr 2.8 Strontium	39 88.906 Y 4.5 Yttrium	40 91.22 Zr 6.49 Zirconium	41 92.906 Nb 8.55 Niobium	42 95.94 Mo 10.2 Molybdenum	43 98 Tc 11.5 Technetium	44 101.07 Ru 12.2 Rhodium	45 102.9055 Rh 12.4 Rhodium	46 106.4 Pd 12.0 Palladium	47 107.868 Ag 10.5 Silver	48 112.41 Cd 8.65 Cadmium	49 114.82 In 7.31 Indium	50 118.69 Sn 7.30 Tin	51 121.75 Sb 6.68 Antimony	52 127.60 Te 6.24 Tellurium	53 127.60 I 6.49 Iodine	54 131.30 Xe 5.89 Xenon
55 132.90 Cs 1.87 Cesium	56 137.33 Ba 3.5 Barium	57 138.906 La 6.7 Lanthanum	58 178.49 Hf 13.1 Hafnium	59 180.95 Ta 16.6 Tantalum	60 183.85 W 19.3 Tungsten	61 186.207 Re 21.0 Rhenium	76 190.2 Os 22.4 Osmium	77 192.22 Ir 27.16 Iridium	78 195.09 Pt 21.4 Platinum	79 196.9665 Au 19.3 Gold	80 200.59 Hg 13.53 Mercury	81 204.37 Tl 11.85 Thallium	82 207.2 Pb 11.4 Lead	83 206.980 Bi 9.8 Bismuth	84 209 Po 9.4 Polonium	85 210 At ???	86 222 Rn 9.91 Radon
87 223 Fr ???	88 226.02 Ra 5 Radium	89 227.02 Ac 10.07 Actinium	104 261 Unq ???? Unnilquadium	105 262 Unp ???? Unnilpentium	106 263 Unh ???? Unnilhexium	PLAY											
			58 140.12 Ce 6.78 Cerium	59 140.91 Pr 6.77 Praseodymium	60 144.24 Nd 7.00 Neodymium	61 145 Pm 6.475 Promethium	62 150.4 Sm 7.54 Samarium	63 151.96 Eu 5.26 Europium	64 157.25 Gd 7.89 Gadolinium	65 158.92 Tb 8.27 Terbium	66 162.5 Dy 8.54 Dysprosium	67 164.93 Ho 8.90 Holmium	68 167.26 Er 9.06 Erbium	69 169.93 Tm 9.33 Thulium	70 173.04 Yb 6.98 Ytterbium	71 174.97 Lu 9.84 Lutetium	
			90 232.0 Th 11.7 Thorium	91 231 Pa 15.4 Protactinium	92 238.02 U 18.90 Uranium	93 237 Np 20.4 Neptunium	94 237 Pu 19.8 Plutonium	95 243 Am 13.6 Americium	96 247 Cm 13.511 Curium	97 247 Bk ???? Berkelium	98 251 Cf ???? Californium	99 252 Es ???? Einsteinium	100 257 Fm ???? Fermium	101 258 Md ???? Mendelevium	102 259 No ???? Nobelium	103 260 Lr ???? Lawrencium	

MATERIAL PROPERTIES

	Symbol	Units	Si	Ge	GaAs	GaP	SiO ₂	Si ₃ N ₄
Atoms per unit cell			8	8	8	8		
Atomic Number	Z		14	32	31/33	31/15	14/8	14/7
Atomic weight	MW	g/g-mole	28.09	72.59	144.64	100.70	60.08	140.28
Lattice constant	ao	nm	0.54307	0.56575	0.56532	0.54505		0.775
Atomic density	No	cm ⁻³	5.00E22	4.42E22	2.21E22	2.47E22	2.20E22	1.48E22
Density	d	g cm ⁻³	2.328	5.323	5.316	4.13	2.19	3.44
Energy Gap 300°K	Eg	eV	1.124	0.67	1.42	2.24	8~9	4.7
Relative permittivity	εr		11.7	16.0	13.1	10.2	3.9	7.5
Index of refraction	n		3.44	3.97	3.3	3.3	1.46	2.0
Melting point	Tm	°C	1412	937	1237	1467	1700	
Specific heat	Cp	J (gK) ⁻¹	0.70	0.32	0.35		1.4	0.17
Thermal diffusivity	K	w(cmK) ⁻¹	0.87	0.36	0.44	0.004	0.32	
Coefficient expansion	Dth	K ⁻¹	2.5E-6	5.7E-6	5.9E-6	5.3E-6	5E-6	2.8E-6
Intrinsic carrier conc	ni	cm ⁻³	1.45E10	2.4E13	9.0E6			
Electron Mobility	μn	cm ² /Vs	1417	3900	8800	300	20	
Hole Mobility	μp	cm ² /Vs	471	1900	400	100	10E-8	
Density of States conduction	Nc	cm ⁻³	2.8E19	1.04E19	4.7E17			
Density of States valance	Nv	cm ⁻³	1.04E19	6.0E18	7.0E18			
Breakdown Electric Field	E	V/cm	3E5	8E4	3.5E5		6~9E6	
Effective mass electron	mn*/mo		1.08	0.55	0.068	0.5		
Effective mass hole	mp*/mo		0.81	0.3	0.5	0.5		
Electron affinity	qX	eV	4.05	4.00	4.07	4.3	1.0	

OXIDE GROWTH

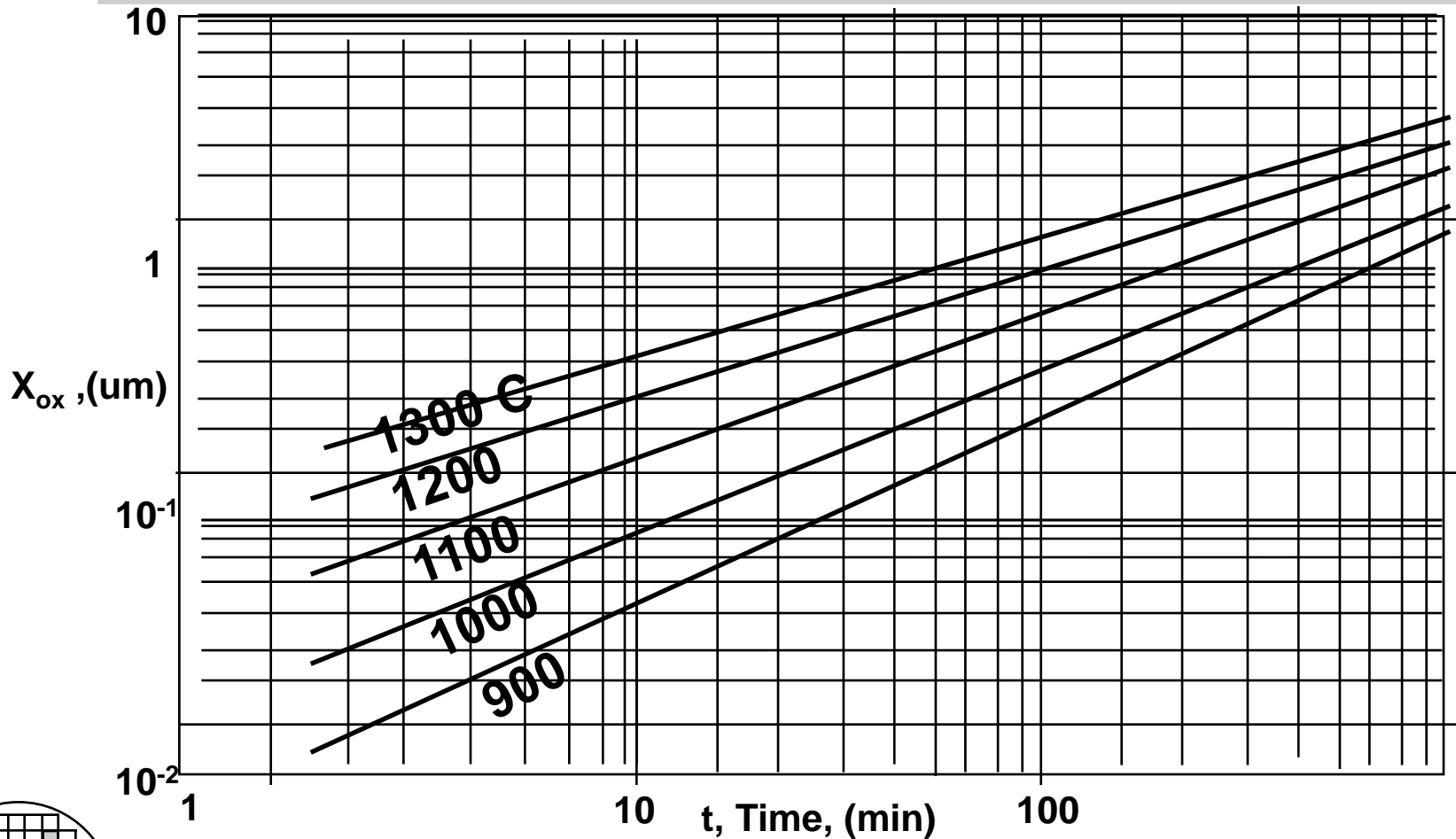


Dry oxide O_2 only

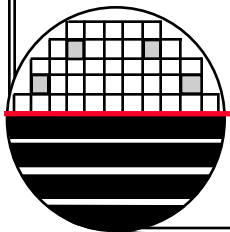
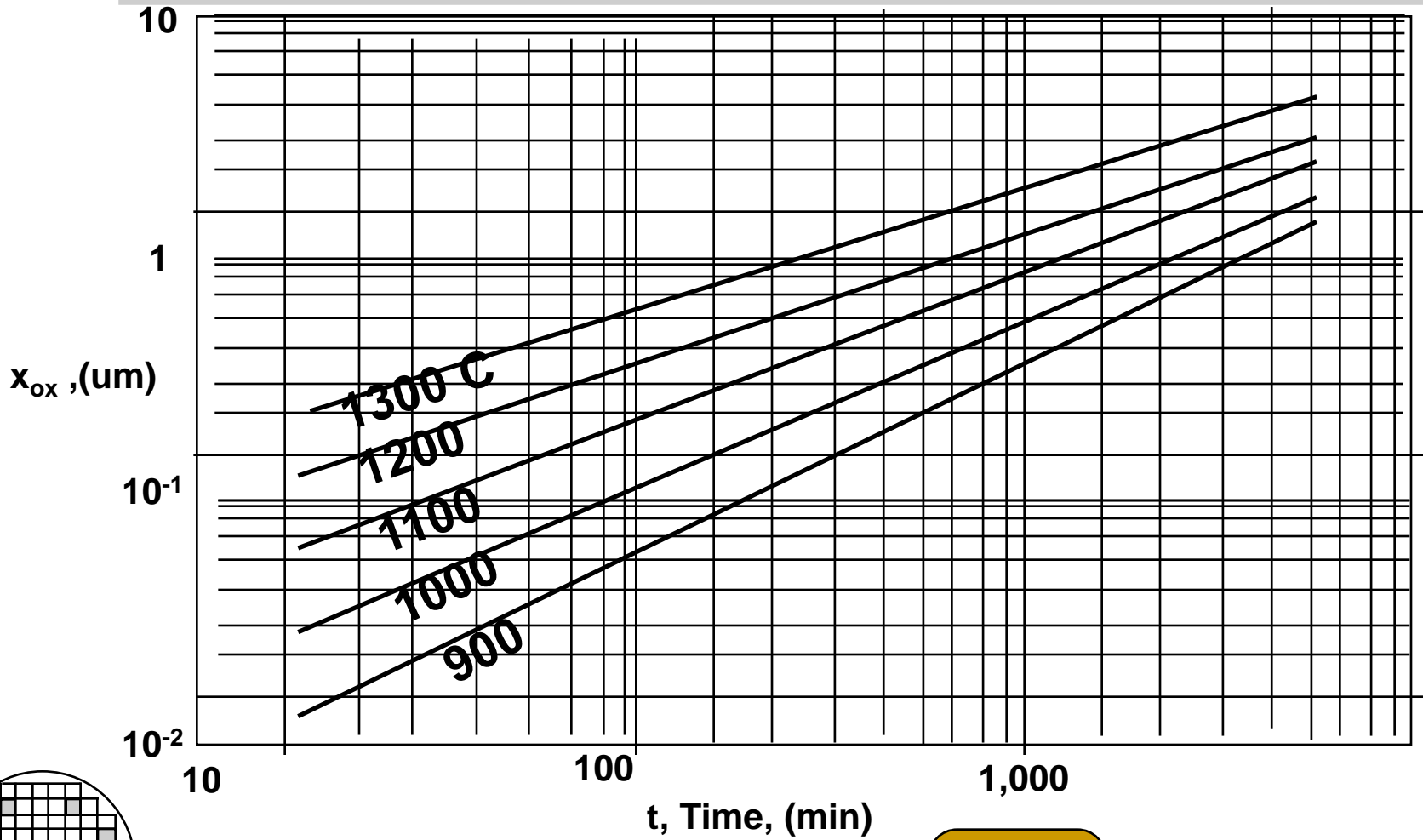
Wet oxide O_2 bubbled through water

Steam burn H_2 in O_2 to make H_2O (steam)

WET OXIDE GROWTH CHART



DRY OXIDE GROWTH CHART



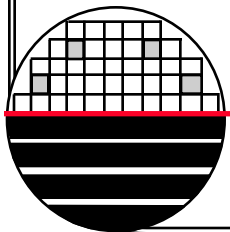
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OXIDE GROWTH CALCULATOR

	A	B	C	D	E	F	G	H	I	J	
1	ROCHESTER INSTITUTE OF TECHNOLOGY							OXIDE.XLS			
2	MICROELECTRONIC ENGINEERING							7/4/2014			
3											
4	CALCULATION OF OXIDE THICKNESS							Dr. Lynn Fuller / Jamie Wasiewicz			
5											
6	To use this spreadsheet change the values in the white boxes. The rest of the sheet is										
7	protected and should not be changed unless you are sure of the consequences. The										
8	calculated results are shown in the purple boxes. O2 bubbled through warm water is "wet",										
9	burning H2 with O2 is called "steam"										
10											
11	CONSTANTS			VARIABLES				CHOICES			
12	K	1.38E-23 J/K						1=yes, 0=no			
13	(Bo/Ao) dry	6230000 μm/hr		Temp =	900 °C		wet	0			
14	Ea (dry)	2 eV		time =	42 min		dry	0			
15	(Bo/Ao) wet	89500000 μm/hr		Partial Pressure, p =	1.00 Atm		steam	1			
16	Ea (wet)	2.05 eV						<100>	1		
17	(Bo/Ao) steam	1.63E+08 μm/hr						<111>	0		
18	Ea (steam)	2.05 eV									
19	Bo dry	7.72E+02 μm ² /hr		Xint=	0 Å						
20	Ea (dry)	1.23 eV									
21	Bo wet	2.14E+02 μm ² /hr		Silicon VLSI Technology, Plummer, Deal, Griffin							
22	Ea (wet)	0.71 eV		Prentice Hall, 2000, pg 319-369							
23	Bo steam	3.86E+02 μm ² /hr									
24	Ea (steam)	0.78 eV		(Bo/Ao)/1.68 for <100>							
25											
26	CALCULATIONS:										
27											
28	Xox (Oxide thickness)=(A/2){[1+(t+Tau)4B/A^2]^0.5 - 1} =							1005 Å			
29											
30	B = [Bo exp (-Ea/KTemp)]*p	0.1737431 μm ² /hr									
31	B/A = [(Bo/Ao) exp (-Ea/KTemp)]*p	1.57E-01 μm/hr									
32	A	1.1100164 μm									
33	Tau = (Xi2+AXi)/B	0 hr									
34											
35											
36	Xox	↑		Oxide-SiO2		←		Original Silicon Surface Prior to Oxide Growth			
37			Silicon				↑		0.46 Xox (silicon consumed)		
38											
39											
40											
41											
42											

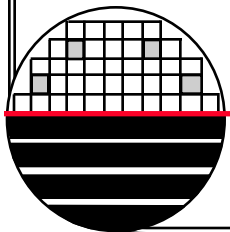
OXIDE.XLS



EXAMPLES

1. Estimate the oxide thickness resulting from 50 min. soak at 1100 °C in wet oxygen.

2. If 1000 Å of oxide exists to start with, what is resulting oxide thickness after an additional 50 min. soak at 1100 °C in dry oxygen.



OXIDE THICKNESS COLOR CHART

Thickness	Color
500Å	Tan
700	Brown
1000	Dark Violet - Red Violet
1200	Royal Blue Blue
1500	Light Blue - Metallic Blue
1700	Metallic - very light Yellow Green
2000	Light Gold or Yellow - Slightly Metallic
2200	Gold with slight Yellow Orange
2500	Orange - Melon
2700	Red Violet
3000	Blue - Violet Blue
3100	Blue Blue
3200	Blue - Blue Green
3400	Light Green
3500	Green - Yellow Green
3600	Yellow Green
3700	Yellow
3900	Light Orange
4100	Carnation Pink
4200	Violet Red
4400	Red Violet
4600	Violet
4700	Blue Violet

Thickness	Color
4900	Blue Blue
5000	Blue Green
5200	Green
5400	Yellow Green
5600	GreenYellow
5700	Yellow - "Yellowish"(at times appears to be Lt gray or metal)
5800	Light Orange or Yellow - Pink
6000	Carnation Pink
6300	Violet Red
6800	"B'ur'sh"(appears violet red, Blue Green, looks Blue)
7200	Blue Green - Green
7700	"Yellowish"
8000	Orange
8200	Salmon
8500	Dull, Light Red Violet
8600	Violet
8700	Blue Violet
8900	Blue Blue
9200	Blue Green
9500	Dull Yellow Green
9700	Yellow - "Yellowish"
9900	Orange
10000	Carnation Pink

Nitride Thickness = (Oxide Thickness)(Oxide Index/Nitride Index)
 Eg. Yellow Nitride Thickness = (2000)(1.46/2.00) = 1460

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PLAY

DIFFUSION FROM A CONSTANT SOURCE

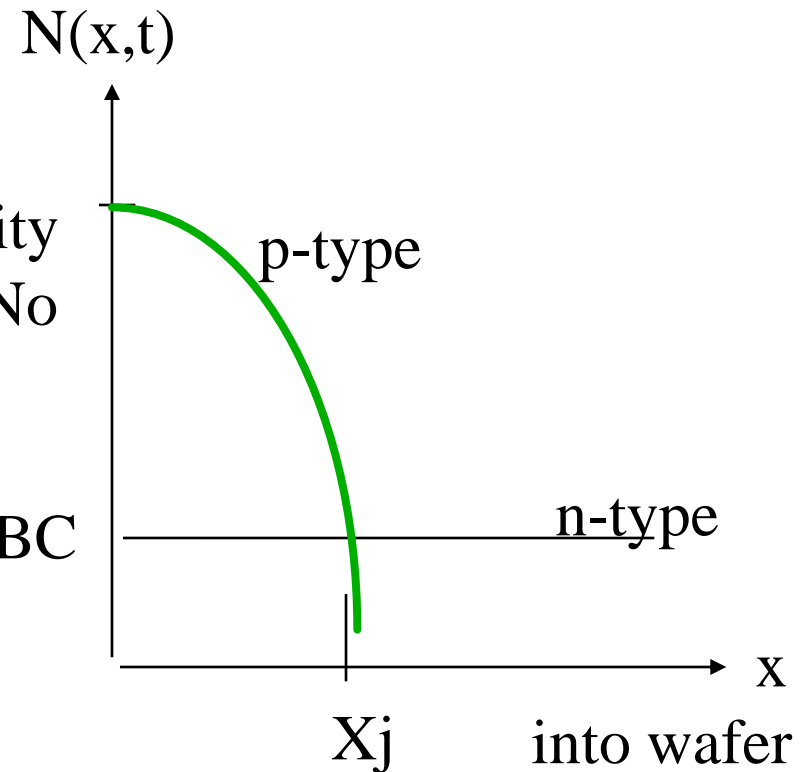
PLAY

STOP

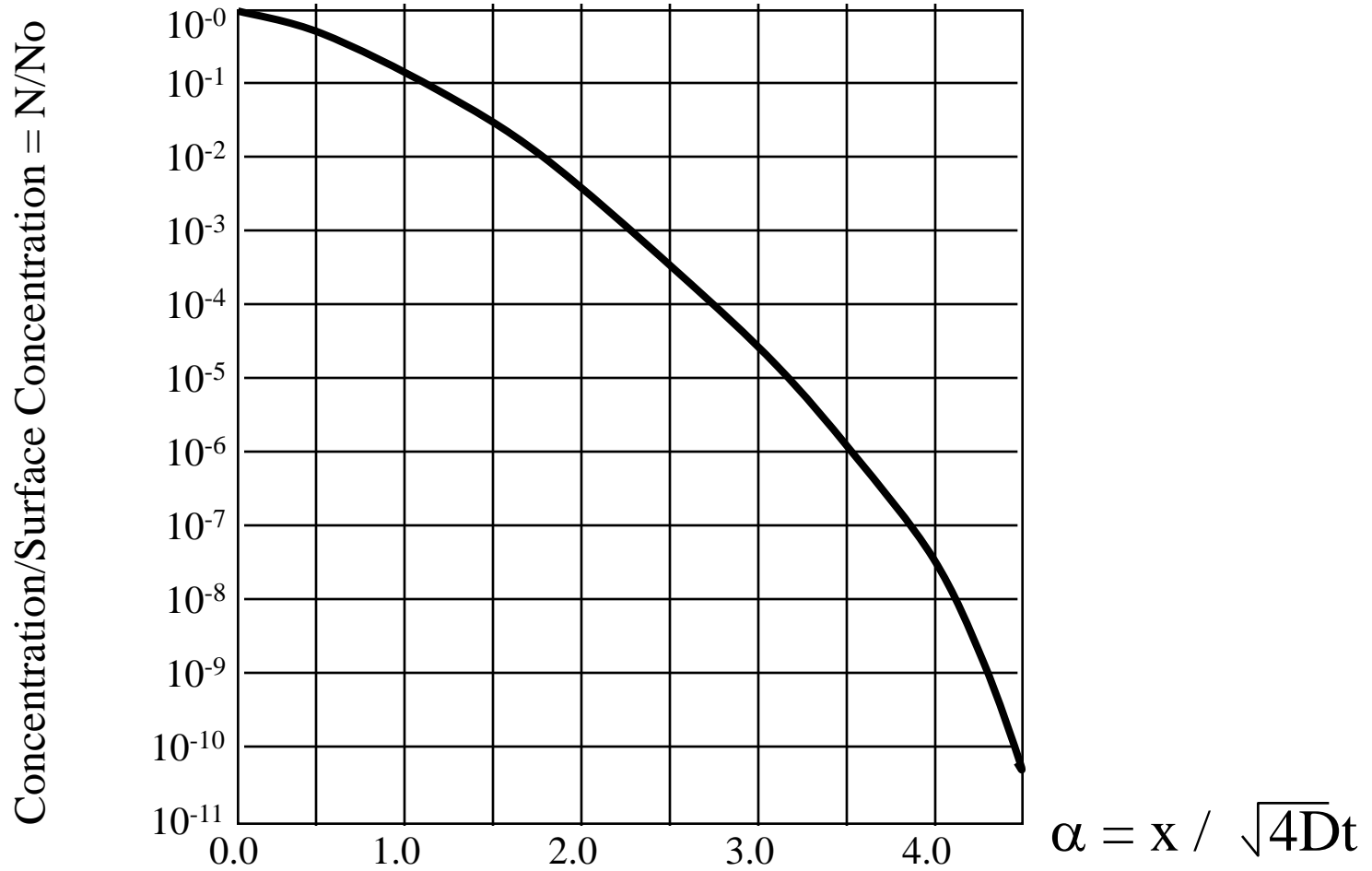
$$N(x,t) = N_0 \operatorname{erfc} \left(\frac{x}{2 \sqrt{Dt}} \right)$$

Solid Solubility Limit, N_0

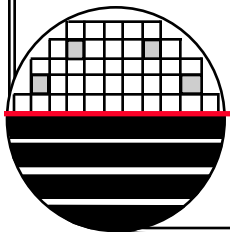
Wafer Background Concentration, NBC



ERFC FUNCTION



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DIFFUSION CONSTANTS AND SOLID SOLUBILITY

DIFFUSION CONSTANTS

TEMP	BORON DRIVE-IN	PHOSPHOROUS PRE	PHOSPHOROUS DRIVE-IN	BORON SOLID SOLUBILITY NOB	PHOSPHOROUS SOLID SOLUBILITY NOP
900 °C	1.07E-15 cm²/s	2.09e-14 cm ² /s	7.49E-16 cm ² /s	4.75E20 cm⁻³	6.75E20 cm ⁻³
950	4.32E-15	6.11E-14	3.29E-15	4.65E20	7.97E20
1000	1.57E-14	1.65E-13	1.28E-14	4.825E20	9.200E20
1050	5.15E-14	4.11E-13	4.52E-14	5.000E20	1.043E21
1100	1.55E-13	9.61E-13	1.46E-13	5.175E20	1.165E21
1150	4.34E-13	2.12E-12	4.31E-13	5.350E20	1.288E21
1200	1.13E-12	4.42E-12	1.19E-12	5.525E20	1.410E21
1250	2.76E-12	8.78E-12	3.65E-12	5.700E20	1.533E21

PLAY

TEMPERATURE DEPENDENCE OF DIFFUSION CONSTANTS

Temperature Dependence:

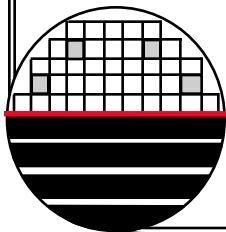


$$D = D_0 \text{Exp} (-E_A/kT) \text{ cm}^2/\text{sec} \quad k = 8.625\text{E-}5 \quad \text{eV}/^\circ\text{K} \quad T \text{ in Kelvins}$$

Boron	$D_0 = 0.76$	Phosphorous	$D_0 = 3.85$
	$E_A = 3.46$		$E_A = 3.66$

Temperature Dependence of the Solid Solubility of Boron and Phosphorous in Silicon

$\text{NOB} = 3.5\text{E}17T + 1.325\text{E}20 \text{ cm}^{-3}$	T in Celsius
$\text{NOP} = 2.45\text{E}18T - 1.53\text{E}21 \text{ cm}^{-3}$	T in Celsius



DIFFUSION FROM A LIMITED SOURCE

$$N(x,t) = \frac{Q'_A(tp) \text{ Exp } (-x^2/4Dt)}{\sqrt{\pi Dt}}$$

PLAY

for erfc predeposit

$$Q'_A(tp) = Q_A(tp)/\text{Area} = 2 N_0 \sqrt{(Dp)tp} / \pi = \text{Dose}$$

PLAY

for ion implant predeposit

$$Q'_A(tp) = \text{Dose}$$

Where D is the diffusion constant at the drive in temperature and t is the drive in diffusion time, Dp is the diffusion constant at the predeposit temperature and tp is the predeposit time

DIFFUSION MASKING CALCULATOR

Select
Boron or Phosphorous
Enter
Temperature and Time

	A	B	C	D	E	F	G	H
1	Rochester Institute of Technology						Raymond Krom	
2	Microelectronic Engineering						Dr. Lynn Fuller	
3	9.5.07						Raymond Krom	
4								
5	Diffusion Mask Calculator				Enter 1-Yes 0-No in white boxes			
6	Temperatures must be between 1000C and 1200C							
7	or result will be in error.							
8	Dopant			Diffusion				
9	Boron	<input type="checkbox"/>		Temp.	<input type="text" value="1100"/>	°C		
10	Phosphorous	<input type="checkbox"/>		Time	<input type="text" value="100"/>	minutes		
11						Boron	1867 Angstroms	
12	Oxide					Phosp	6399 Angstroms	
13	Fitted to data taken from Hamilton and Howard						<input type="text" value="6399"/>	Angstroms
14								

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From: Hamilton and Howard

DIFFUSION MASKING

Phosphorous Masking

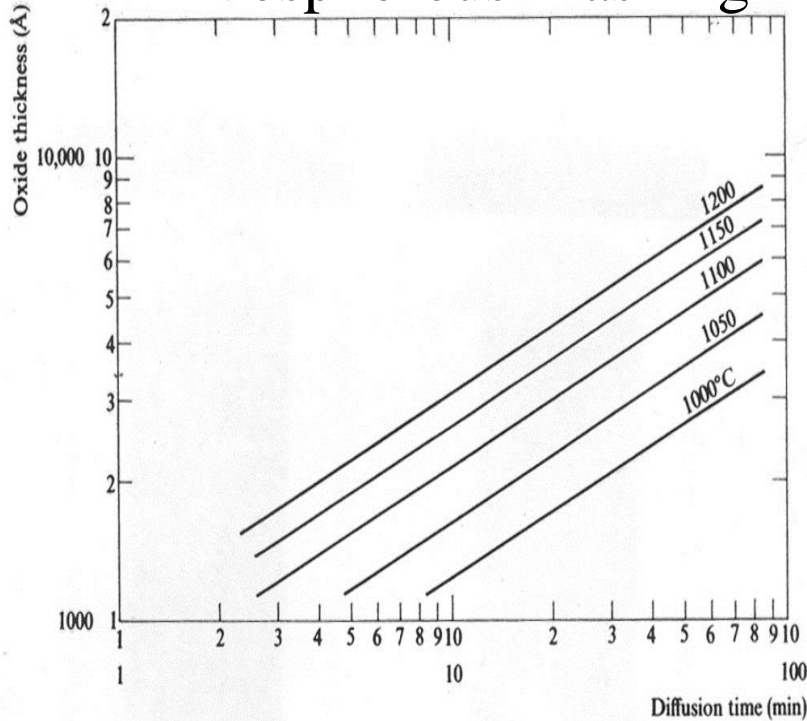


FIGURE 2-28 Minimum oxide thickness required for phosphorus diffusion masking as a function of time (after Burger and Donovan⁴).

Boron Masking

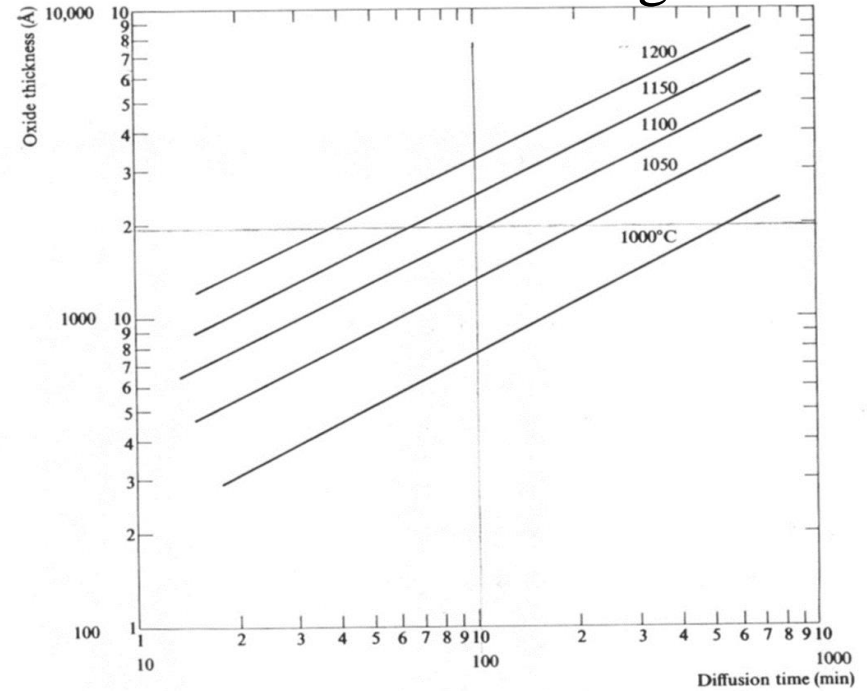


FIGURE 2-27 Minimum oxide thickness required to mask boron at various temperatures (after Burger and Donovan⁴).

From: Hamilton and Howard

DIFFUSION AND DRIVE IN CALCULATIONS

GIVEN	VALUE	UNITS
Starting Wafer Resistivity	10	ohm-cm
Starting Wafer Type	1	1 or 0
	0	1 or 0
Pre Deposition Temperature	950	°C
Pre Deposition Time	15	min
Drive-in Temperature	1100	°C
Drive-in Time	480	min

CALCULATE	VALUE	UNITS
Solid Solubility at Temperature of Pre Deposition	4.65E+20	cm-3
Diffusion Constant at Temperature of Pre Deposition	3.93E-15	cm/sec
Diffusion Constant at Temperature of Drive-in	1.43E-13	cm/sec

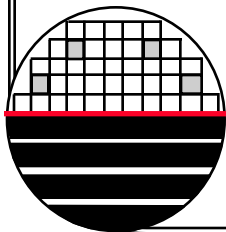
CALCULATION OF DIFFUSION CONSTANTS

	D0 (cm ² /s, EA (eV))	
Boron	0.76	3.46
Phosphorous	3.85	3.66
NOB = 3.5E17 (T) + 1.325E20		
NOP = 2.45E18(T) - 1.53E21		

DIFFUSION.XLS

CALCULATIONS	VALUE	UNITS
Substrate Doping = 1 / (q μ _{max} Rho)	4.42E+14	cm-3
Ratio of N _{sub} /N _s =	9.51E-07	
Approximate inverse erfc from erfc(u) ≈ (e ^{-u²})/(uπ) ^{0.5} inv_erfc =	3.47	

RESULTS	VALUE	UNITS
x _j after pre deposition = ((4D _p t _p) ^{0.5})*(inv_erfc(N _{sub} /N _s))	0.13	μm
Pre deposition Dose, Q _A = 2N _o (D _p t _p /π) ^{0.5}	9.87E+14	atoms/cm ²
x _j after drive-in = ((4 D _d t _d /Q _A) ln (N _{sub} (πD _d t _d) ^{0.5})) ^{0.5}	4.03	μm
average doping N _{ave} = Dose/x _j	2.45E+18	atoms/cm ³
mobility (μ) at Doping equal to N _{ave}	109	cm ² /V-s
Sheet Resistance = 1/(q (μ(N _{ave}))Dose)	58	ohms
Surface Concentration After Drive-in = Dose/ (pDt) ^{0.5}	8.68E+18	cm-3



DIFFUSION FROM A LIMITED SOURCE

GIVEN

Starting Wafer Resistivity
Starting Wafer Type

	VALUE	UNITS
Rho =	10	ohm-cm
n-type = 1	1	1 or 0
p-type = 1	0	1 or 0

Pre Deposition Ion Implant Dose

4.00E+15	ions/cm ²
----------	----------------------

Drive-in Temperature

1000	°C
------	----

Drive-in Time

360	min
-----	-----

CALCULATE

Diffusion Constant at Temperature of Drive-in

VALUE	UNITS
1.43E-14	cm/sec

CALCULATION OF DIFFUSION CONSTANTS

	D0 (cm ² /s)	EA (eV)
Boron	0.76	3.46
Phosphorous	3.85	3.66

IonImplt.xls

CALCULATIONS

Substrate Doping = $1 / (q \mu_{max} Rho)$

VALUE	UNITS
4.42E+14	cm ⁻³

RESULTS

Pre deposition Dose

x_j after drive-in = $((4 D_d t_d / QA) \ln (N_{sub} (\pi D_d t_d)^{0.5}))^{0.5}$

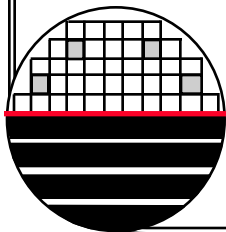
average doping $N_{ave} = Dose / x_j$

mobility (μ) at Doping equal to N_{ave}

Sheet Resistance = $1 / (q (\mu(N_{ave}))Dose)$

Surface Concentration = $Dose / (pDt)^{0.5}$

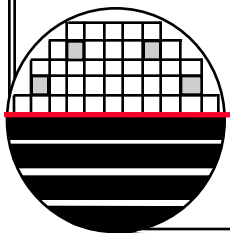
VALUE	UNITS
4.00E+15	atoms/cm ²
1.25	μm
3.21E+19	atoms/cm ³
57	cm ² /V-s
27.6	ohms
1.28E+20	cm ⁻³



EXAMPLE

1. A predeposit from a p-type spin-on dopant into a $1\text{E}15\text{ cm}^{-3}$ wafer is done at 1100°C for 10 min. Calculate the resulting junction depth and dose.

2. The spin-on dopant is removed and the Boron is driven in for 2 hours at 1100°C . What is the new junction depth?



RESISTANCE, RESISTIVITY, SHEET RESISTANCE

Resistance = $R = \rho L/\text{Area} = \rho_s L/w$ ohms

Resistivity = $\rho = 1/(q\mu_n n + q\mu_p p)$ ohm-cm

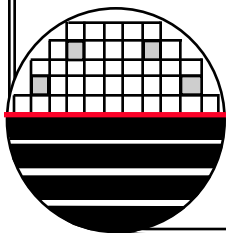
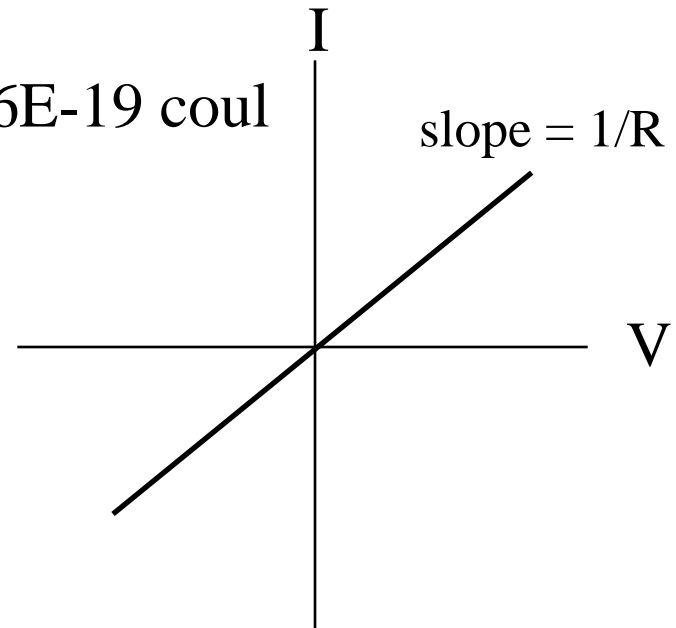
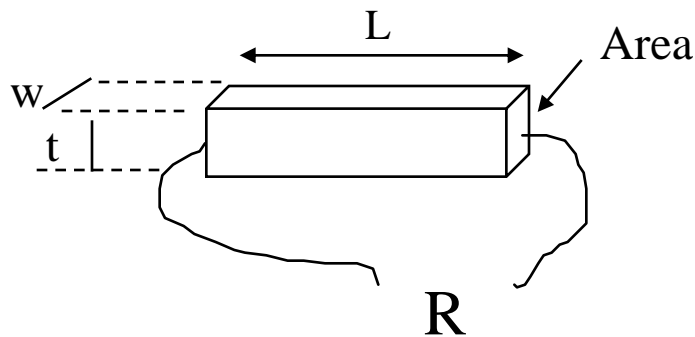
PLAY

Sheet Resistance = $\rho_s = 1/(\int q \mu(N) N(x) dx) \sim 1/(q\mu \text{Dose})$ ohms/square

PLAY

$$\rho_s = \rho / t$$

$$q = 1.6E-19 \text{ coul}$$



CALCULATION OF CARRIER CONCENTRATIONS

B	1.11E+03	Nd =	<input type="text" value="3.00E+16"/>	cm-3	Donor Concentrat
h	6.63E-34 Jsec	Ed=	<input type="text" value="0.049"/>	eV below Ec	
εo	8.85E-14 F/cm	Na =	<input type="text" value="8.00E+15"/>	cm-3	Acceptor Concen
εr	11.7	Ea=	<input type="text" value="0.045"/>	eV above Ev	
ni	1.45E+10 cm-3	Temp=	<input type="text" value="300"/>	°K	
Nc/T^3/2	5.43E+15				
Nv/T^3/2	2.02E+15				

Donor and Acceptor Levels (eV above or below Ev or Ec)

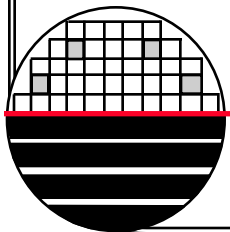
Boron	0.044
Phosphorous	0.045
Arsenic	0.049

carrier_conc.xls

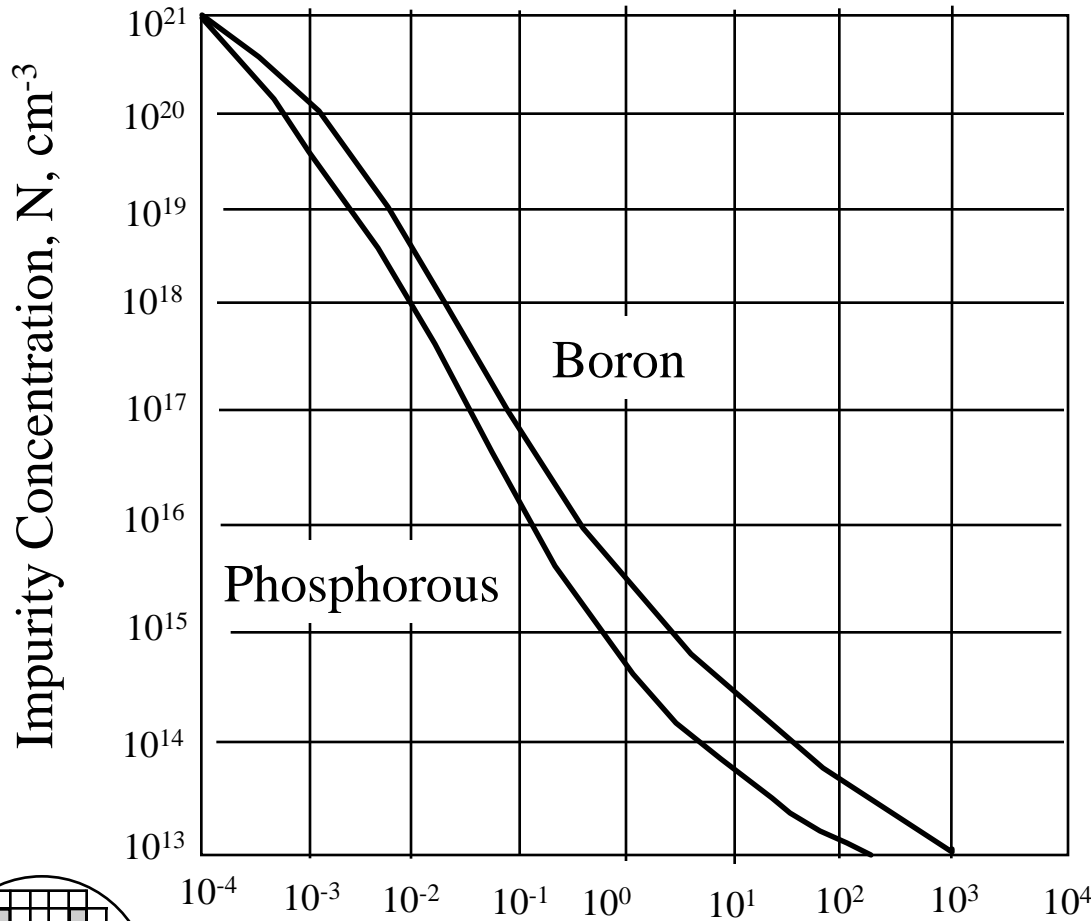
CALCULATIONS: (this program makes a guess at the value of the fermi level and trys to minimize the charge balance)

KT/q	<input type="text" value="0.026"/>	Volts
Eg=Ego-(aT^2/(T+B))	<input type="text" value="1.115"/>	eV
Nc	<input type="text" value="2.82E+19"/>	cm-3
Nv	<input type="text" value="1.34E+01"/>	cm-3
Fermi Level, Ef	<input type="text" value="0.9295"/>	eV
free electrons, n = Nc exp(-q(Ec-Ef)/KT)	<input type="text" value="2.17E+16"/>	cm-3
Ionized donors, Nd+ = Nd*(1+2*exp(q(Ef-Ed)/KT))^-1)	<input type="text" value="2.97E+16"/>	cm-3
holes, p = Nv exp(-q(Ef-Ev)/KT)	<input type="text" value="3.43E-15"/>	cm-3
Ionized acceptors, Na- = Na*(1+2*exp(q(Ea-Ef)/KT))^-1)	<input type="text" value="8.00E+15"/>	cm-3
Charge Balance = p + Nd+ - n - Na-	<input type="text" value="3.22E+12"/>	cm-3

Click on Button to do Calculation



RESISTIVITY OF SILICON VS DOPING

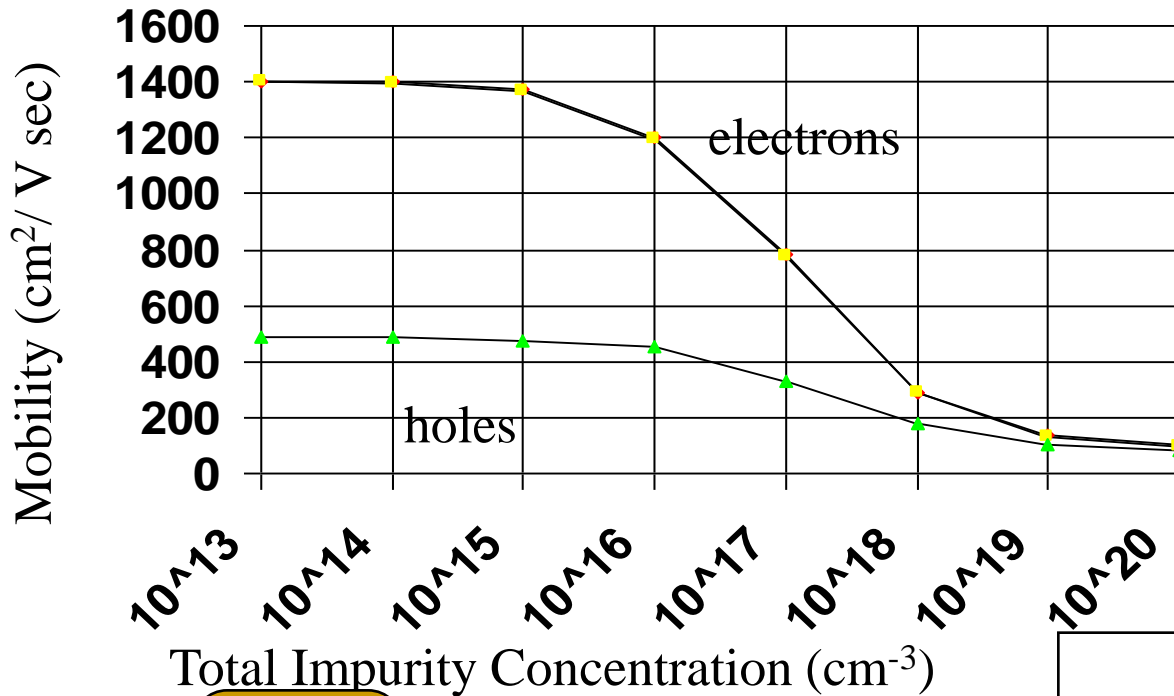


$$\rho = 1/(q\mu(N)N)$$

Because μ is a function of N and N is the doping, the relationship between resistivity ρ and N is given in the figure shown, or calculated from equations for $\mu(N)$



ELECTRON AND HOLE MOBILITY



PLAY

Electron and hole mobilities in silicon at 300 K as functions of the total dopant concentration (N). The values plotted are the results of the curve fitting measurements from several sources. The mobility curves can be generated using the equation below with the parameters shown:

PLAY

$$\mu(N) = \mu_{mi} + \frac{(\mu_{max} - \mu_{min})}{\{1 + (N/N_{ref})^\alpha\}}$$

Parameter	Arsenic	Phosphorous	Boron
μ_{min}	52.2	68.5	44.9
μ_{max}	1417	1414	470.5
N_{ref}	9.68X10 ¹⁶	9.20X10 ¹⁶	2.23X10 ¹⁷
α	0.680	0.711	0.719

From Muller and Kamins, 3rd Ed., pg 33

TEMPERATURE EFFECTS ON MOBILITY

Derived empirically for silicon for T in K between 250 and 500 °K and for N (total dopant concentration) up to 1 E20 cm-3

$$\mu_n(T,N) = 88 T_n^{-0.57} + \frac{1250 T_n^{-2.33}}{1 + [N / (1.26E17 T_n^{2.4})]^{0.88} T_n^{-0.146}}$$

PLAY

$$\mu_p(T,N) = 54.3 T_n^{-0.57} + \frac{407 T_n^{-2.33}}{1 + [N / (2.35E17 T_n^{2.4})]^{0.88} T_n^{-0.146}}$$

EXCELL WORKSHEET TO CALCULATE MOBILITY

MICROELECTRONIC ENGINEERING

3/13/2005

CALCULATION OF MOBILITY

Dr. Lynn Fuller

To use this spreadsheet change the values in the white boxes. The rest of the sheet is protected and should not be changed unless you are sure of the consequences. The calculated results are shown in the purple boxes.

CONSTANTS

VARIABLES

CHOICES

$T_n = T/300 = 1.22$

Temp= °K

N total cm-3

1=yes, 0=no

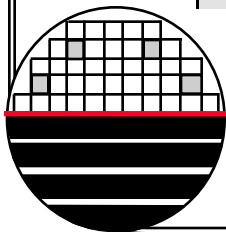
n-type

p-type

<100>

Kamins, Muller and Chan; 3rd Ed., 2003, pg 33

mobility= cm²/(V-sec)



EXCELL WORKSHEET TO CALCULATE RESISTANCE

CALCULATION OF RESISTANCE FROM LENGTH, WIDTH, THICKNESS AND IMPLANT DOSE

To use this spreadsheet change the values in the white boxes. The rest of the sheet is protected and should not be changed unless you are sure of the consequences. The calculated results are shown in the purple boxes.

Resistors_Poly.xls

Calculation of Mobility of Single Crystal Silicon

CONSTANTS	VARIABLES	CHOICES
$T_n = T/30$ 1.00	Temp = <input style="width: 50px;" type="text" value="300"/> °K	1=yes, 0=no
Concentration from Dose / thickness, $N = \text{Dose}/t$	<input style="width: 50px;" type="text" value="2.00E+20"/> cm ⁻³	n-type <input style="width: 50px;" type="text" value="1"/> p-type <input style="width: 50px;" type="text" value="0"/>
Kamins, Muller and Chan; 3rd Ed., 2003, pg 33		
		mobility = $\mu =$ <input style="width: 50px;" type="text" value="90"/> cm ² /(V-sec)

Calculation of Resistance

L length is the drawn length	Length, L = <input style="width: 50px;" type="text" value="600"/> μm
Width is the drawn width	Width, W = <input style="width: 50px;" type="text" value="100"/> μm
Thickness is known if poly, or Xj from Diffusion.xls	Thickness, t = <input style="width: 50px;" type="text" value="0.5"/> μm
Implanter setting if doped by ion implant or from Diffusion.xls if doped by c	Dose = <input style="width: 50px;" type="text" value="1.00E+16"/> /cm ²
	Poly ? <input style="width: 50px;" type="text" value="1"/> Yes=1, No=0
	resistance/poly grain boundary <input style="width: 50px;" type="text" value="0.3"/> ohm

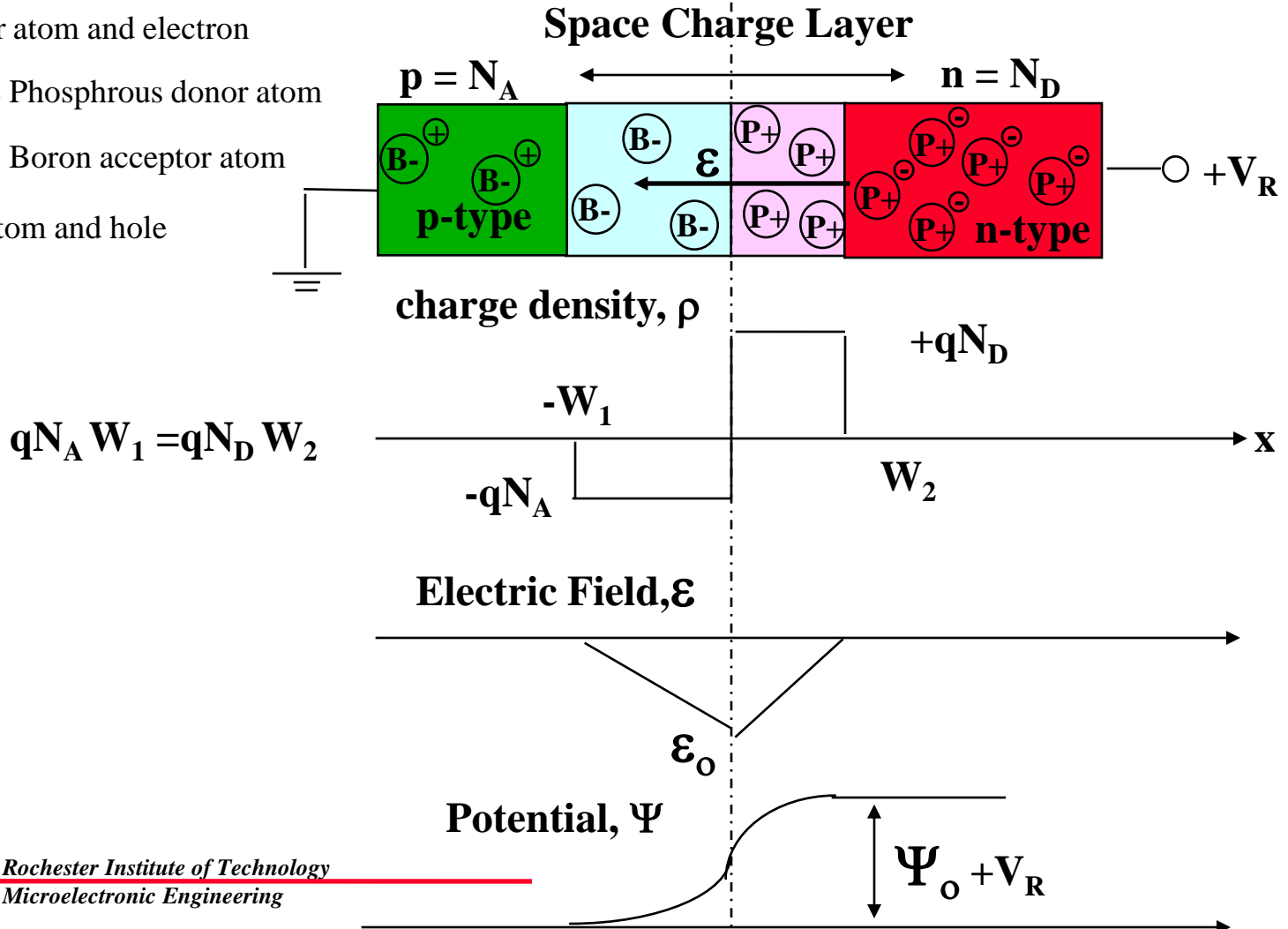
Calculation of Resistance

approximate number of grain boundaries in path = $L/t =$	<input style="width: 50px;" type="text" value="1200"/>
Average Doping = Dose/Thickness =	<input style="width: 50px;" type="text" value="2.00E+20"/> atoms/cm ³
Mobility, $\mu =$	<input style="width: 50px;" type="text" value="90"/> cm ² /v-sec
$q = 1.6\text{e-}19$ coulomb / ion	Rhos = sheet resistance = $1/(q \mu \text{Dose}) =$ <input style="width: 50px;" type="text" value="7"/> ohms/sq
$R = \text{Rho} L / W / t$	Rho = bulk resistivity = <input style="width: 50px;" type="text" value="14"/> ohm-cm
$R = \text{Rhos} L / W$	Resistance = <input style="width: 50px;" type="text" value="402"/> ohms
	If Poly the effective sheet resistance = <input style="width: 50px;" type="text" value="67"/> ohms/sq

We assume the grain size is equal to the poly film thickness. We calculate the number of grains from the length, L, divided by the grain size, t. We also assume the grain boundary adds a fixed resistance that is not a function of temperature or doping. The resistance of a grain boundary is found from resistance measurements of poly resistors.

UNIFORMLY DOPED PN JUNCTION

- $(P^+)^{\ominus}$ Phosphorous donor atom and electron
- (P^+) Ionized Immobile Phosphorous donor atom
- (B^-) Ionized Immobile Boron acceptor atom
- $(B^-)^{\oplus}$ Boron acceptor atom and hole



UNIFORMLY DOPED PN JUNCTION

Built in Voltage:

$$\Psi_0 = KT/q \ln (N_A N_D / n_i^2)$$

$$n_i = 1.45E10 \text{ cm}^{-3}$$

Width of Space Charge Layer, W: with reverse bias of V_R volts

$$W = (W_1 + W_2) = [(2\epsilon/q) (\Psi_0 + V_R) (1/N_A + 1/N_D)]^{1/2}$$

W_1 width on p-side

W_2 width on n-side

$$W_1 = W [N_D / (N_A + N_D)]$$

$$W_2 = W [N_A / (N_A + N_D)]$$

Maximum Electric Field:

$$E_0 = - [(2q/\epsilon) (\Psi_0 + V_R) (N_A N_D / (N_A + N_D))]^{1/2}$$

Junction Capacitance per unit area:

$$C_j' = \epsilon_0 \epsilon_r / W = \epsilon_0 \epsilon_r / [(2\epsilon/q) (\Psi_0 + V_R) (1/N_A + 1/N_D)]^{1/2}$$

$$\begin{aligned} \epsilon &= \epsilon_0 \epsilon_r = 8.85E^{-12} \text{ (11.7) F/m} \\ &= 8.85E^{-14} \text{ (11.7) F/cm} \end{aligned}$$

EXAMPLE

Example: If the doping concentrations are $N_A=1E15$ and $N_D=3E15$ cm^{-3} and the reverse bias voltage is 0, then find the built in voltage, width of the space charge layer, width on the n-side, width on the p-side, electric field maximum and junction capacitance. Repeat for reverse bias of 10, 40, and 100 volts.

$$\Psi_o = V_{bi} = \frac{KT}{q} \ln \left(\frac{N_A N_D}{n_i^2} \right) =$$

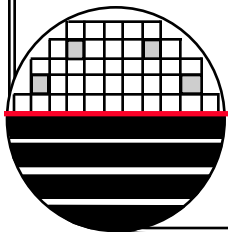
$$W = (W_1 + W_2) = \left[\frac{2\epsilon}{q} (\Psi_o + V_R) \left(\frac{1}{N_A} + \frac{1}{N_D} \right) \right]^{1/2} =$$

$$W_1 =$$

$$W_2 =$$

$$E_{max} =$$

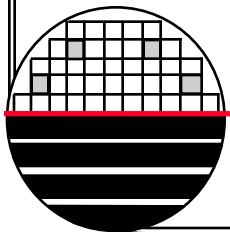
$$C_j =$$



EXAMPLE CALCULATIONS

	A	B	C	D	E	F	G	H	I	J	
1	ROCHESTER INSTITUTE OF TECHNOLOGY							PN.XLS			
2	MICROELECTRONIC ENGINEERING							3/4/2009			
3											
4	CALCULATIONS FOR PN JUNCTION (ELECTROSTATICS)							DR. LYNN FULLER			
5											
6	To use this spreadsheet change the values in the white boxes. The rest of the sheet is										
7	protected and should not be changed unless you are sure of the consequences. The										
8	calculated results are shown in the purple boxes.										
9											
10	CONSTANTS					VARIABLES					
11	K	1.38E-23 J/K									
12	q	1.60E-19 Coul			Temp=	<input type="text" value="300"/>		K			
13	Eg	1.12 eV									
14	so	8.85E-14 F/cm			Nd =	<input type="text" value="1.00E+19"/>		cm-3			
15	sr	11.7			Na =	<input type="text" value="5.00E+14"/>		cm-3			
16	ni	1.45E+10 cm-3			Vr =	<input type="text" value="0"/>		Volts Reverse Bias Voltage			
17											
18											
19											
20											
21	CALCULATIONS:										
22											
23	KT/q =							<input type="text" value="0.025887"/>	Volts		
24	Vbi = (KT/q) ln (NaNd/ni ²)							<input type="text" value="0.80"/>	Volts		
25	W = [(2s/q)(Vbi+Vr)(1/Na + 1/Nd)] ^{0.5}							<input type="text" value="1.44"/>	µm		
26	W1 = W[Nd/(Na+Nd)]							<input type="text" value="1.44"/>	µm		
27	W2 = W[Na/(Na+Nd)]							<input type="text" value="0.00"/>	µm		
28	Eo = -[(2q/eoer)(Vbi+Va)(NaNd/(Na+Nd))] ^{0.5}							<input type="text" value="-1.11E+04"/>	V/cm		
29	Cj' = eoer/W							<input type="text" value="7.21E-09"/>	F/cm ²		

PN.XLS



LONG CHANNEL THRESHOLD VOLTAGE, V_T

Flat-band Voltage $V_{FB} = \phi_{ms} - \frac{Q_{ss}}{C'_{ox}} - \frac{1}{C'_{ox}} \int_0^{X_{ox}} \frac{X}{X_{ox}} \rho(x) dx$

p-type substrate
(n-channel)
n-type substrate
(p-channel)
 $Q_{ss} = q N_{ss}$

Bulk Potential : $\phi_p = -KT/q \ln (N_A/n_i)$

$\phi_n = +KT/q \ln (N_D/n_i)$

Work Function Difference $\phi_{MS} = \phi_M - (X + Eg/2q + [\phi_p])$

$\phi_{MS} = \phi_M - (X + Eg/2q - [\phi_n])$

Maximum Depletion Width:
(W_{dmax}) $\sqrt{\frac{4 \epsilon_s [\phi_p]}{qNa}}$

$\sqrt{\frac{4 \epsilon_s [\phi_n]}{qNd}}$

Threshold Voltage: $V_T = V_{FB} + 2 [\phi_p] + \frac{1}{C'_{ox}} \sqrt{2 \epsilon_s q Na (2[\phi_p])}$

p-type substrate

Threshold Voltage: $V_T = V_{FB} - 2 [\phi_n] - \frac{1}{C'_{ox}} \sqrt{2 \epsilon_s q Nd (2[\phi_n])}$

n-type substrate

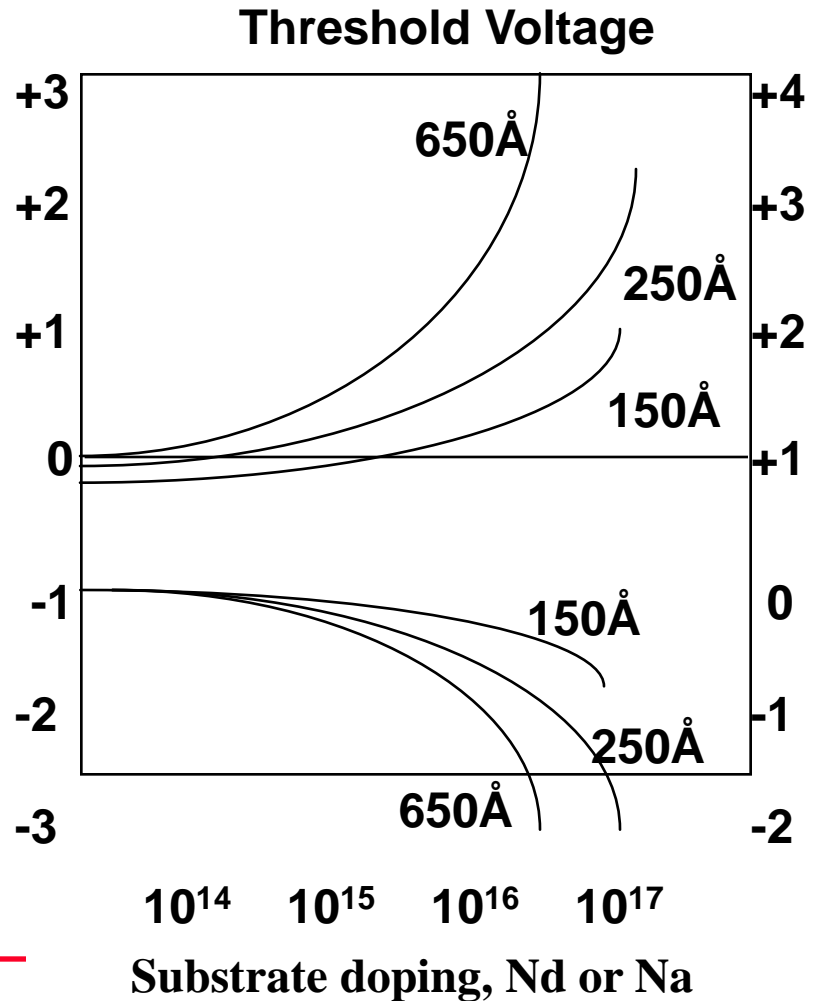


LONG CHANNEL V_t

Gate work function, n+, p+, aluminum
 Substrate doping, Nd or Na
 Oxide thickness, X_{ox}
 Surface State Density, N_{ss} or Q_{ss}
 also
 Substrate to Source voltage difference



n+ poly gate left scale
 p+ poly gate right scale
 $Q_f = 0$
 $V_{bs} = 0$
 implant dose = zero



VT ADJUST IMPLANT

Assume that the total implant is shallow (within W_{dmax})

$$\pm \Delta V_t = q \text{Dose}^* / C_{ox}'$$

where Dose^* is the dose that is added to the Si
 C_{ox}' is gate oxide capacitance/cm²
 $C_{ox}' = \epsilon_0 \epsilon_r / X_{ox}$

Boron gives + shift

Phosphorous gives - shift

Example: To shift +1.0 volts implant Boron through 1000 Å Kooi oxide at an energy to place the peak of the implant at the oxide/silicon interface. Use a $\text{Dose} = \Delta V_t C_{ox}' / q$
 $= (1.0)(3.9)(8.85E-14) / (1.6E-16) = 2.16E11$ ions/cm²
 but multiply by 2 since 1/2 goes into silicon

MOSFET THRESHOLD VOLTAGE CALCULATION

CALCULATION OF MOSFET THRESHOLD VOLTAGE LYNN FULLER

To use this spreadsheet change the values in the white boxes. The rest of the sheet is protected and should not be changed unless you are sure of the consequences. The calculated results are shown in the purple boxes.

CONSTANTS	VARIABLES	CHOICES
T = 300 K	Na = 4.00E+17 cm ⁻³	Aluminum gate
KT/q = 0.026 volts	Nd = 4.00E+17 cm ⁻³	n+ Poly gate
ni = 1.45E+10 cm ⁻³	Nss = 3.00E+11 cm ⁻²	p+ Poly gate
Eo = 8.85E-14 F/cm	Xox = 50 Å	N substrate
Er si = 11.7		P substrate
Er SiO2 = 3.9		
E affinity = 4.15 volts		Desired VT
q = 1.60E-19 coul		or
Eg = 1.124 volts		Delta VT
		Given Dose (Boron)

1=yes, 0=No	
<input type="checkbox"/>	Select one type of gate
<input type="checkbox"/>	Select one type of substrate

CALCULATIONS:

CALCULATIONS:	RESULTS
METAL WORK FUNCTION	= 4.122989 volts
SEMICONDUCTOR POTENTIAL	= +/- 0.445453 volts
OXIDE CAPACITANCE / CM2	= 6.9E-07 F/cm2
METAL SEMI WORK FUNCTION I	= -0.143558 volts
FLAT BAND VOLTAGE	= -0.213093 volts
THRESHOLD VOLTAGE	= -1.601792 volts
DELTA VT = VTdesired - VT	= 0.601792 volts
IMPLANT DOSE	= 2.6E+12 ions/cm2 x 2 = 5.19271E+12
	where + is Boron, - is Phosphorous
IMPLANT DOSE FOR GIVEN Delta	= 8.63E+13 ions/cm2 x 2 = 1.72575E+14
Vt WITH GIVEN DOSE	= -1.451133 volts assume 1/2 dose in Si

MOSFETVT.XLS

ION IMPLANT EQUATIONS

Gaussian Implant Profile

$$N(x) = \frac{N'}{\sqrt{2\pi} \Delta R_p} \exp \left[\frac{-(X-R_p)^2}{2\Delta R_p^2} \right]$$

$R_p = \text{Range}$
 $\Delta R_p = \text{Straggle}$

} From Curves

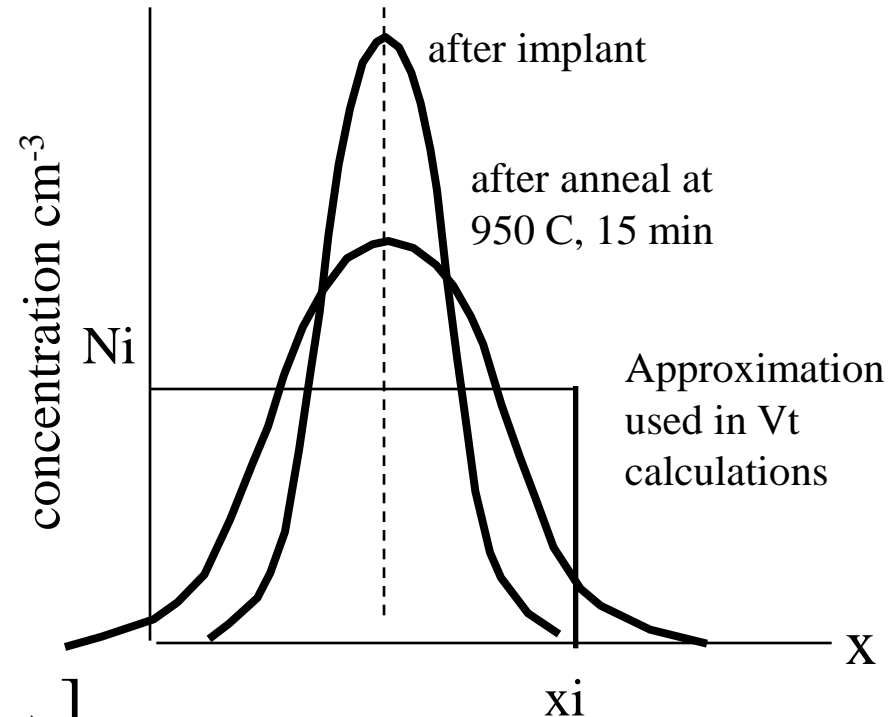
$$N' = \text{Dose} = \int \frac{I}{mqA} dt$$

After Anneal

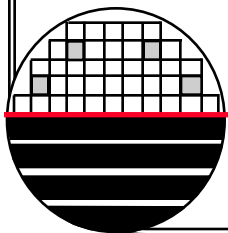
$$N(x) = \frac{N'}{\sqrt{2\pi} \sqrt{\Delta R_p^2 + 2Dt}} \exp \left[\frac{-(X-R_p)^2}{2(\Delta R_p^2 + Dt)} \right]$$

where D is diffusion constant at the anneal temperature
 t is time of anneal

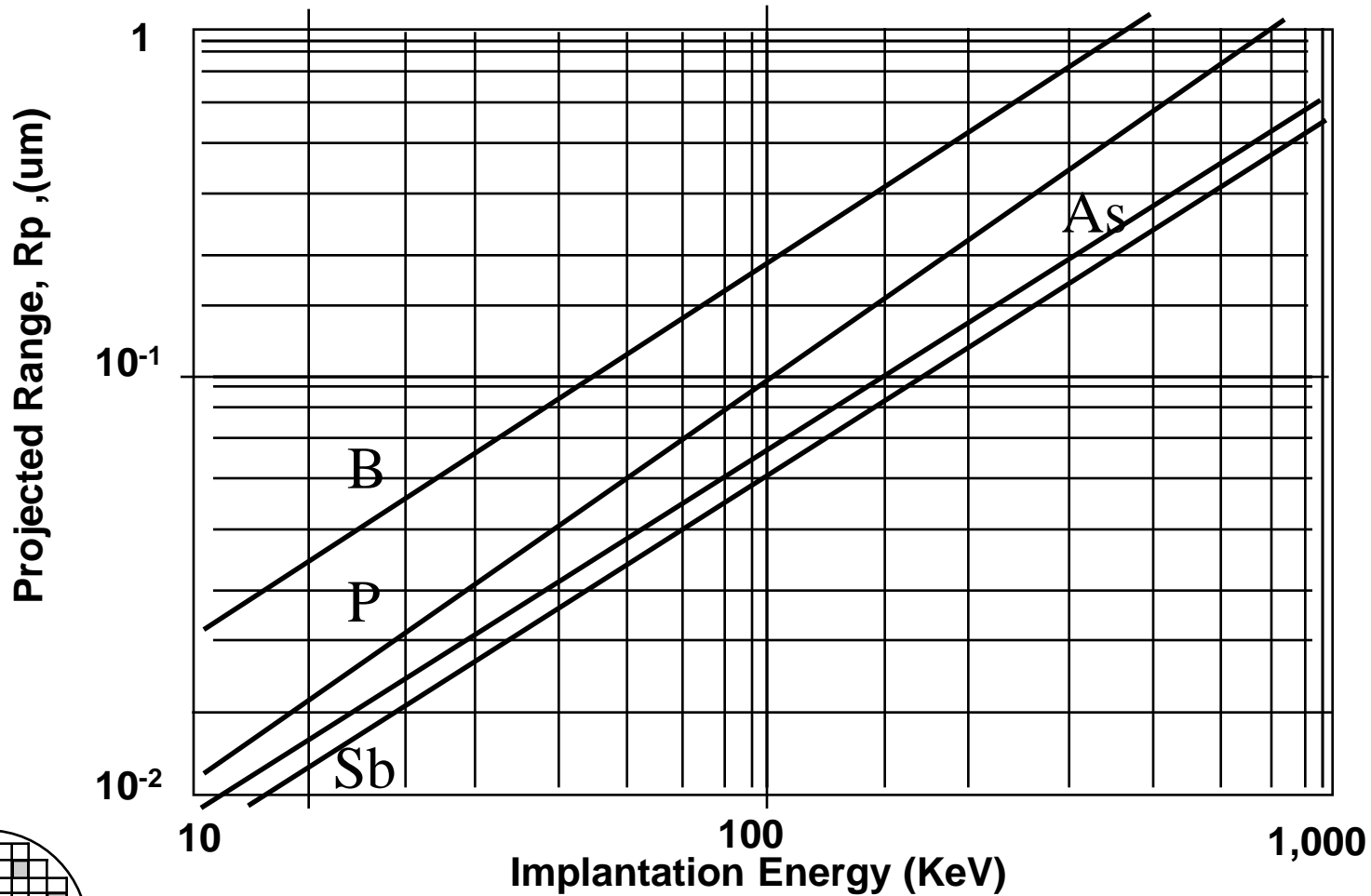
Rochester Institute of Technology
 Microelectronic Engineering



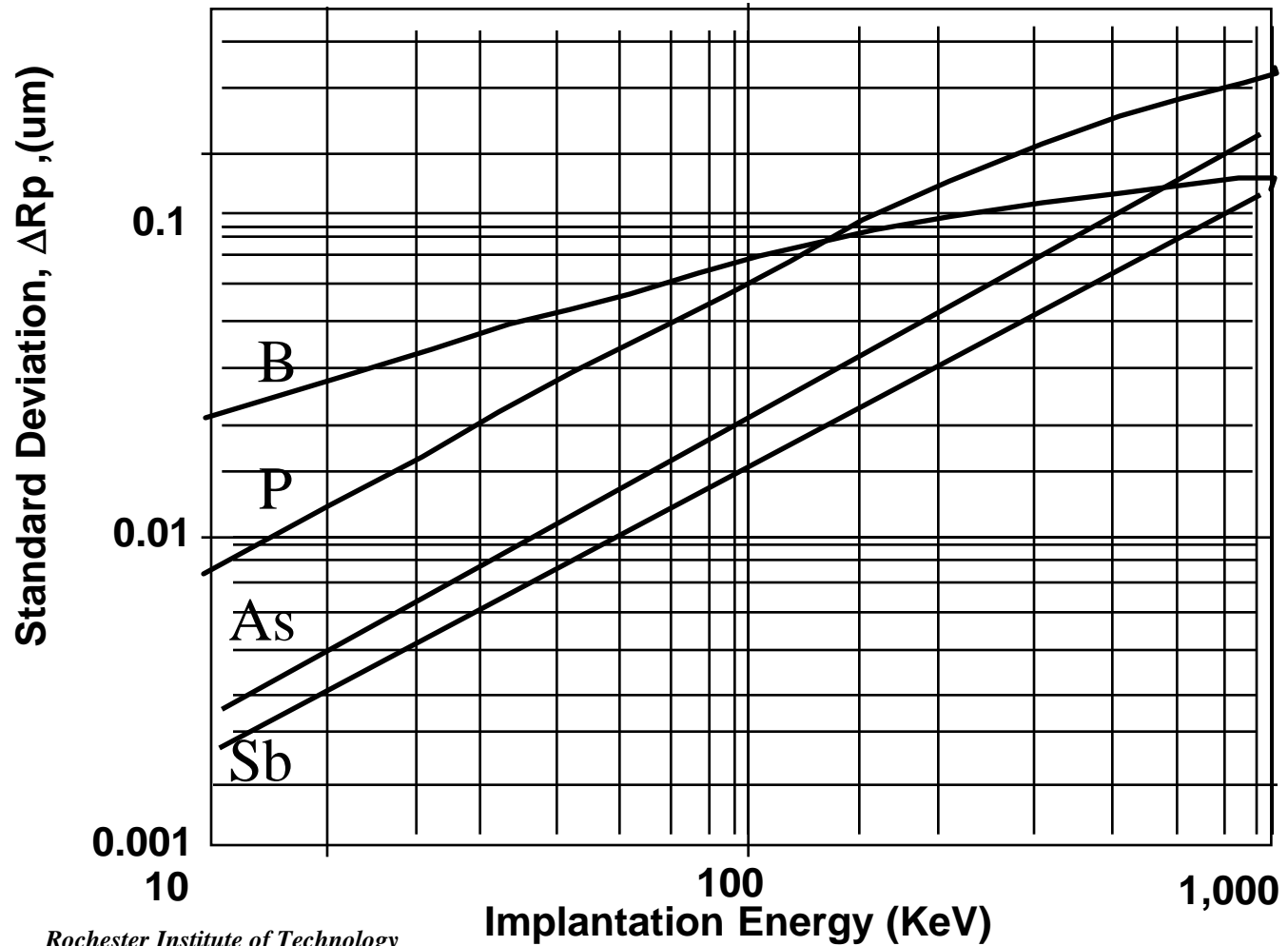
Approximation
 $N' = N_i x_i$



ION IMPLANT RANGE



ION IMPLANT STANDARD DEVIATION



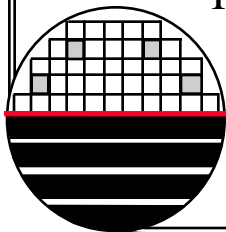
ION IMPLANT MASKING CALCULATOR

Rochester Institute of Technology			Lance Barron
Microelectronic Engineering			Dr. Lynn Fuller
11/20/2004			

IMPLANT MASK CALCULATOR		Enter 1 - Yes 0 - No in white boxes	
DOPANT SPECIES	MASK TYPE	ENERGY	
B11	Resist	60	KeV
BF2	Poly		
P31	Oxide		
	Nitride		
Thickness to Mask >1E15/cm3 Surface Concentration		4073.011	Angstroms

This calculator is based on Silvaco Suprem simulations using the Dual Pearson model.

In powerpoint click on spread sheet to change settings for a new calculation

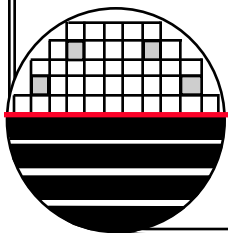


REFERENCES

1. Basic Integrated Circuit Engineering, Douglas J. Hamilton, William G. Howard, McGraw Hill Book Co., 1975.
2. Micro Electronics Processing and Device Design, Roy a. Colclaser, John Wiley & Sons., 1980.
3. Device Electronics for Integrated Circuits, Richard S. Muller, Theodore I. Kamins, Mansun Chan, John Wiley & Sons., 3rd Ed., 2003.
4. VLSI Technology, Edited by S.M. Sze, McGraw-Hill Book Company, 1983.
5. Silicon Processing for the VLSI Era, Vol. 1., Stanley Wolf, Richard Tauber, Lattice Press, 1986.
6. The Science and Engineering of Microelectronic Fabrication, Stephen A. Campbell, Oxford University Press, 1996.

HOMWORK - REVIEW OF IC TECHNOLOGY

1. If a window is etched in 5000 \AA of oxide and the wafer is oxidized again for 50 min in wet O_2 at $1050 \text{ }^\circ\text{C}$ what is the new thickness (where it was 5000 \AA), the thickness in the etch window, and the step height in the silicon if all the oxide is etched off the wafer. Draw a picture showing original Si surface.
2. A Boron diffusion is done into 5 ohm-cm n-type wafer involving two steps. First a short predeposit at 950 C for 30 min., followed by removal of the diffusion source and a drive in at 1100 C for 2 hours. Calculate the junction depth and the sheet resistance of the diffused layers. Estimate the oxide thickness needed to mask this diffusion.
3. For a pn junction with the p side doping of $1\text{E}17$ and the n side at $1\text{E}15$ calculate, width of space charge layer, width on p side, on n side, capacitance per unit area, max electric field.
4. Calculate the threshold voltage for an aluminum gate PMOSFET fabricated on an n-type wafer with doping of $5\text{E}15$, a surface state density of $7\text{E}10$, and gate oxide thickness of 150 \AA . What is the threshold voltage if the surface state density is $3\text{E}11$?
5. Calculate the ion implant dose needed to shift the threshold voltage found in the problem above to -1 Volts .



HOMWORK - EXACT CALCULATION OF SHEET RESISTANCE FOR A DIFFUSED LAYER

1. A Boron p-type layer is diffused into an n-type silicon wafer (1E15 cm⁻³) at 1100 °C for 1 hour. Calculate the exact value of the sheet resistance and compare to the approximate value.

Sheet Resistance = $\rho_s = 1 / \left(\int q \mu(N) N(x) dx \right) \sim 1 / (q \mu \text{ Dose})$ ohms/square

$$\mu(N) = \mu_{\min} + \frac{(\mu_{\max} - \mu_{\min})}{\{1 + (N/N_{\text{ref}})^\alpha\}}$$

for Boron

μ_{\min}	44.9
μ_{\max}	470.5
N_{ref}	2.23×10^{17}
α	0.719

Let $Q'_A(tp) = 5.633 \times 10^{15} \text{ cm}^{-2}$
 $D = 1.55 \times 10^{-13} \text{ cm}^2/\text{s}$
 $t = 1 \text{ hour}$

$$N(x,t) = \frac{Q'_A(tp) \text{ Exp} \left(-x^2/4Dt \right)}{\sqrt{\pi Dt}}$$

HW SOLUTION - EXACT CALCULATION OF SHEET RESISTANCE FOR A DIFFUSED LAYER

Divide the diffused layer up into 100 slices and for each slice find the doping and exact mobility. Calculate the sheet resistance from the reciprocal of the sum of the conductance of each slice.

