

TEMPERATURE DEPENDENCE OF FIELD EMISSION OF SINGLE CRYSTAL PLANES OF TUNGSTEN

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T-F emission (W. W. Dolan, W. P. Dyke, *Phys. Rev.* **95**, 327 (1954)) of single crystal planes of tungsten as an emitter was investigated in Müller's tube. A comparison of the experimental data with the theory of Murphy and Good is made. A new method of computing the geometrical factor β for any point of the tip is shown.

A field emission tube suitable for measurements of local currents has been constructed ([2]–[5]). The collector current $i(T)$ was measured as a function of temperature T at a fixed anode voltage U . Measurements were made for several values of U . According to Murphy and Good [6] the approximate formula [7] for the *T-F* current is

$$i(T) = i(0) (1 + K^2 T^2 / 6) \quad (1)$$

where $i(T)$ is field emission current at temperature T , $i(0)$ field emission current at temperature 0° K, and

$$K = 2.77 \times 10^4 \varphi^{1/2} t(\gamma) / F \quad (2)$$

where $t(\gamma) \ll 1$ stands for the slowly-variable Nordheim function, F is electric field determined by $F = \beta U$, where β is a geometrical factor, depending on geometry and shape of the tip and tube. Equation (2) was transformed into the equilibrate shape

$$\frac{\Delta i}{i(T_p)} = \frac{(T/T_p)^2 - 1}{1 + 6/K^2 T^2} \equiv f(x), \quad (3)$$

where $\Delta i = i(T) - i(T_p)$, $T_p = 295^\circ\text{K}$ is the temperature at which measurements began. In the coordinate system $\Delta i/i(295) = f[(T/295)^2 - 1]$, the relationship (3) should be a straight line whose slope is

$$(1 + 8.99 \times 10^{-14} F^2 / \varphi)^{-1}. \quad (4)$$

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Figure 1 shows that in the above-mentioned coordinate system the experimental plots for different crystallographic directions have different runs within the measured temperature range, *i. e.* from 295° to 1150°K; however, within the temperature range from 295° to (450–550)°K (depending on the crystallographic planes) all plots may be approximated by straight lines. It for a definite plane, *e. g.*, (221) in the above-mentioned coordinate system, the family of characteristics is drawn, with U treated as a parameter (Fig. 2), it is evident that the slopes of the plots approximating initial runs of the function $\Delta i/i(295) = f(x)$ are changed for different U [7]. Denoting by ϑ the angle of the slope of the straight line $f(x)$

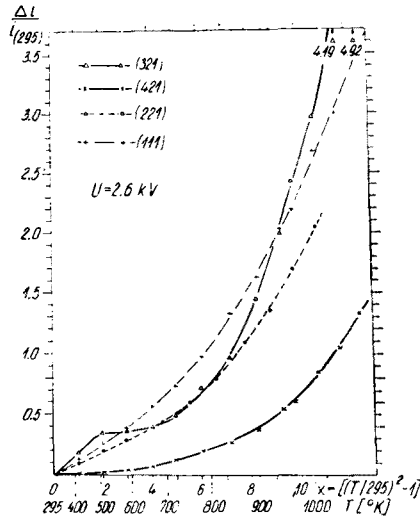


Fig. 1. Relative increase of collector current $\Delta i/i(295)$ as a function of temperature T or variable $x = [(T/295)^2 - 1]$ for different crystallographic planes

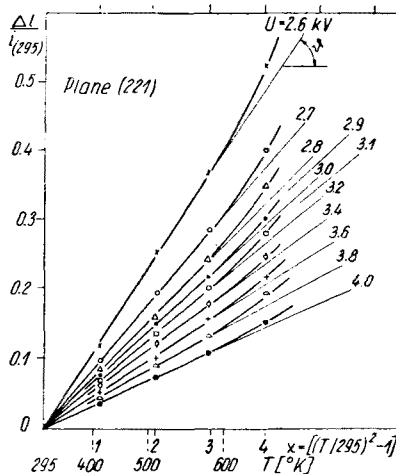


Fig. 2. Relative increase of collector current $\Delta i/i(295)$ as a function of temperature T or variable $x = [(T/295)^2 - 1]$ for (221) crystallographic plane, for several values of voltage U

with respect to the x -axis the relation F^2/φ may be determined with the help of Eq. (4) by $\operatorname{tg} \vartheta$:

$$F^2/\varphi = 1.11 \times 10^{14} (\operatorname{ctg} \vartheta - 1). \quad (5)$$

Now, when we plot the values calculated on the basis of Eq. (5) in the coordinate system $F^2/\varphi = g(U^2)$, we again obtain a straight line (Fig. 3), whose slope is

$$\operatorname{tg} \xi = \beta^2/\varphi. \quad (6)$$

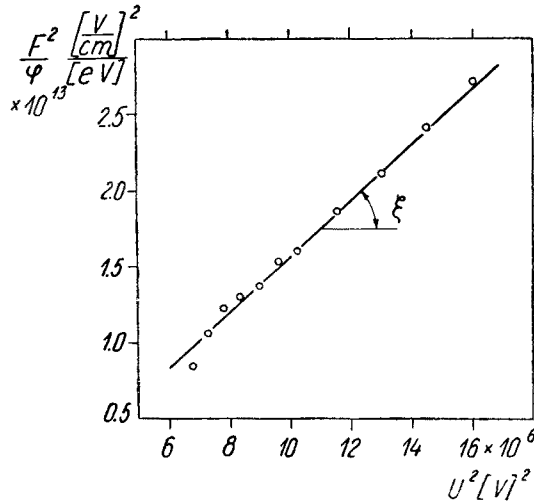


Fig. 3. Values of F^2/φ computed from Eq. (5) as a function of voltage squared

If then we have the F. N.¹ plot for a chosen plane, its slope is

$$m = -B\varphi^{3/2}/\beta, \quad (7)$$

where $B = 6.83 \times 10^{-7} v(y)$ ($v(y)$ is the Nordheim function near unity for the applied fields). Finally, φ and β were calculated from equations (6) and (7). For the (221) plane they have the values $\varphi_{(221)} = 4.4$ eV and $\beta = 9.2 \times 10^3$ cm⁻¹. We see that the value of $\varphi_{(221)}$ calculated in this way is in good agreement with values calculated by others using different methods (Schmidt: $\varphi_{(221)} = 4.37$ eV [4], and Müller: $\varphi_{(221)} = 4.35$ eV [8]). This confirms the correctness of the formula of Murphy and Good for temperatures up to (450–550)°K. This temperature depends on the structure of the crystal faces. The method described here is suitable for computing the distribution of β along a chosen “azimuth” on the tip. The value of β can also be computed by another method² based on the suggestion of Ovchinnikov and Carev [9], but taking into account the coefficient of the shape of the tip, α used by Drechsler and Henkel [10].

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¹ Fowler Nordheim.

² Publication in preparation.

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