

In table 1 we give the values of E_w obtained from equation (9) for u_c equal to $1/\sqrt{3}$ as well as $1/\sqrt{6}$. We have computed for those metals which were considered by Smith

Workfunction Equation as Determined by Mehrotra & Mahanty

$$E_w = -(V + \Delta T)$$

$$= \frac{e^2 \omega_p}{2 \beta} \int_0^{u_c} \frac{du}{[u + (1 + u^2)^{1/2}]^2}$$

$$= \frac{3\mu}{(18\pi)^{1/3}} \frac{e^2}{a_0 \sqrt{r_s}}$$

- a_0 = Bohr radius;
- ω_p = plasmon frequency;
- μ = numerical value for integral;
- $\beta = (1/3)^{0.5} v_f$
- v_f = Fermi velocity;
- r_0 = radius of equilibrium density distribution of free electrons;
- r_s = electron density parameter defined by $no = \frac{3}{4} P i r_s^3 a_0$;
- ΔT = change in kinetic energy of electron added to surface of metal;
- V = electrostatic energy due to external electron at surface;

Table 1. Comparison of the workfunctions for u_c equal to $1/\sqrt{3}$ as well as $1/\sqrt{6}$. Values with superscripts a, b and c are taken from Lang and Kohn (1971), Rivière (1969) and Smith (1969), respectively. r_s values are computed from the electron densities given by Smith (1969).

Metal	r_s	Workfunctions (eV)		Previous theories	Experimental
		This work $u_c = 1/\sqrt{3}$	$u_c = 1/\sqrt{6}$		
Cs	5.64	3.10	2.51	2.49 ^a	2.14 ^a
Rb	5.23	3.22	2.61	2.63 ^a	2.21 ^a
K	4.96	3.30	2.68	2.74 ^a	2.39 ^a
Na	3.99	3.68	2.98	3.06 ^a	2.7 ^a
Li	3.25	4.08	3.31	3.37 ^a	3.1 ^a
Ag	3.01	4.24	3.43	3.49 ^a	4.0 ^a
Au	3.00	4.25	3.44	3.49 ^a	5.22 ^a
Cu	2.67	4.50	3.65	3.65 ^a	4.65 ^a
Ca	3.26	4.07	3.30	3.11 ^c	2.80 ^c
Mg	2.65	4.52	3.66	3.66 ^a	3.66 ^a
Cd	2.59	4.57	3.70	3.36 ^c	4.22 ^b
Zn	2.30	4.85	3.93	3.80 ^a	4.33 ^a
Be	1.88	5.37	4.35	3.75 ^c	3.905 ^b
La	2.71	4.47	3.62	3.30 ^c	3.3 ^c
Tl	2.49	4.66	3.78	3.40 ^c	3.7 ^c
In	2.41	4.74	3.84	3.44 ^c	3.8 ^c
Ga	2.20	4.96	4.02	3.56 ^c	3.96 ^c
Al	2.07	5.11	4.14	3.87 ^a	4.19 ^a
Sn	2.39	4.76	3.85	3.45 ^c	4.425 ^b
Pb	2.31	4.84	3.92	3.80 ^a	4.01 ^a
Ta	1.79	5.50	4.45	3.80 ^c	4.22 ^b
Nb	1.79	5.50	4.45	3.81 ^c	4.37 ^b
W	1.62	5.78	4.68	3.91 ^c	4.55 ^b
Mo	1.61	5.80	4.70	3.92 ^c	4.21 ^b
Re	1.50	6.01	4.87	3.98 ^c	5.0 ^c
Ir	1.41	6.20	5.02	4.02 ^c	5.3 ^c