

USE OF THE HAUDSMITH - SAUNDERSON THEORY  
IN THE ENERGY RANGE 1-20 keV

G. E. Gorelik and S. G. Rozin

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It is shown by means of numerical calculation that the series representing the Haudsmith -Saunderson function is asymptotic.

In calculations related with an investigation of the effect of electrons on matter one usually uses the Haudsmith -Saunderson distribution function [1-3]

$$f_{H-S} = \sum_{l=0}^{\infty} \left( l + \frac{1}{2} \right) \exp \{ -G_l t \} P_l(\cos \theta), \quad (1)$$

where the  $P_l(\cos \theta)$  are Lagrange polynomials;  $\theta$  is the scattering angle;  $t$  is the electron mean free path; the coefficients  $G_l$  are functions of energy and are calculated from recursion relations [1].

TABLE 1. Values of Coefficients  $G_l$  at Energy 6.3 keV for Al

$l$	$2,87 \cdot 10^{-4} G_l$	$l$	$2,87 \cdot 10^{-4} G_l$	$l$	$2,87 \cdot 10^{-4} G_l$
1	3,66	40	51,60	80	-8,79
4	17,66	44	51,61	81	-7,96
8	32,61	48	51,59	82	-12,53
12	41,61	52	51,52	83	-26,89
16	46,53	56	51,34	84	-44,45
20	49,10	60	50,93	92	-429,86
24	50,40	64	50,00	96	-1024,29
28	51,05	68	47,90	99	-1913,60
32	51,37	72	43,18	100	-2350,84
36	51,52	76	32,58		

It is usually considered [2] that up to an energy of order 1 keV series (1) is weakly convergent and attempts are made to sum as few terms of this series as possible. However, as calculations showed, for energies of 1-20 keV series (1) diverges, since  $G_l$  increases with increase of  $l$  ( $G_0 = 0$ ), reaches a maximum at  $l = l_{max}$ , and then decreases, passing into the region of negative values (Table 1).

A comparison of the calculations with the experimental data, which is given in [2, 3], permits us to consider series (1) to be asymptotic and to

TABLE 2. Dependence of  $l_{max}$  on Energy for Al

$n$	$E, \text{ keV}$	$l_{max}$	$n$	$E, \text{ keV}$	$l_{max}$
1	21,3	95	11	10,6	63
2	20,1	—	12	9,5	65
3	19,0	—	13	8,5	55
4	17,9	—	14	6,7	48
5	16,9	—	15	6,3	44
6	15,9	—	16	5,6	46
7	15,0	79	17	5,0	—
8	14,2	75	18	4,7	38
9	13,4	70	19	4,5	—
10	11,9	—	20	4,2	34
			21	4,0	32
			22	3,8	30

Note. The dash indicates that a maximum value was not observed at the given energy in  $G_l$  ( $0 \leq l \leq 200$ ).

Institute of Heat and Mass Transfer, Academy of Sciences of the Belorussian SSR. Institute of Nuclear Power Engineering, Academy of Sciences of the Belorussian SSR, Minsk. Translated from *Inzhenerno-Fizicheskii Zhurnal*, Vol. 21, No. 2, pp. 359-360, August, 1971. Original article submitted September 3, 1970.

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approximate the distribution function with an accuracy to the first discarded term by a part of this series.

To achieve maximum accuracy series (1) should be cut off at a term containing  $(G_l)_{\max}$ . The values of  $l_{\max}$  for some energies are given in Table 2.

The program was compiled for the Minsk-22 computer for further use in Monte Carlo calculations. The calculation time of  $f_{H-S}$  for a fixed value of the angle and energy is 30 sec. The input data indicated in [2] were used.

#### LITERATURE CITED

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3. H. E. Bishop, Brit. J. Appl. Phys., 18, 703 (1967).