

DESY 65/8

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Experimente

INTENSITY AND POLARISATION  
OF COHERENT BREMSSTRAHLUNG SPECTRA  
FROM A DIAMOND CRYSTAL  
IN UNIVERSAL REPRESENTATION

by

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Summary

Plots of intensity and polarisation functions are represented and their use is explained in obtaining the main features of coherent spectra produced from a diamond radiator for any desired electron energy and crystal orientation. The polarisation to be expected may be obtained from the curves for the first peak of the spectrum, which gives the only significant contribution. In addition the contribution of a single lattice point is represented.

Experimental situation

An electron beam with momentum  $\vec{p}_0$ , energy  $E_0$ , strikes the diamond radiator at the small angle  $\theta_0$  with respect to the  $[110]$ -axis of the crystal.  $\alpha$  is the angle between the planes  $\vec{p}_0$ ,  $[110]$  and  $[001]$ ,  $[110]$  in the reciprocal lattice space, see Fig.1.

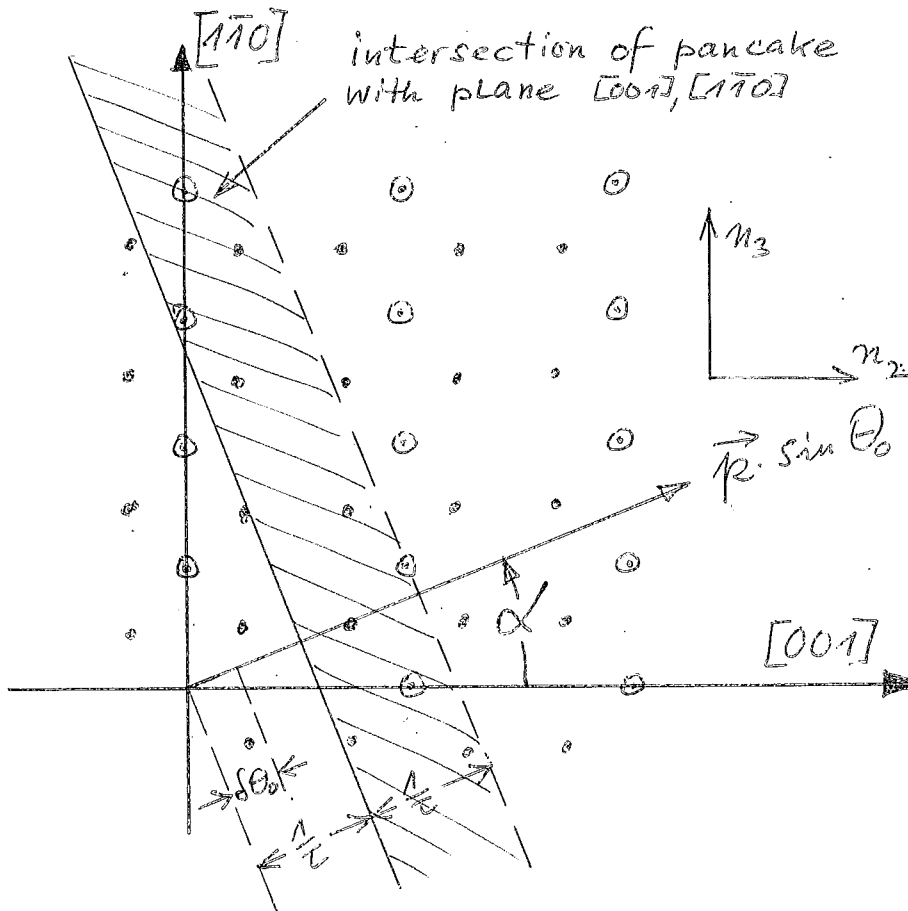


Fig.1 Reciprocal lattice plane, view  $\perp$   $[110]$ .

Coherent spectra for  $\alpha = 0, 90^\circ$

The resulting coherent spectrum, i.e. the dependence of the normalized intensity

$$(1) \quad I = \frac{x \, d\sigma/dx}{N\sigma} \quad (\text{compare AESY 64/9, equ. 39})$$

on the relative quantum energy

$$(2) \quad x = k/E_0$$

may be represented as a function of the product  $\theta_0 \cdot E_0$  and is in its coherent part proportional to  $E_0$ . Also, the position of the discontinuities  $x_d$  as well as the polarisation  $P$  are functions of  $\theta_0 \cdot E_0$ , which have the following form:

$$(3) \quad x_d = 1 / \left( 1 + \frac{37,6}{\theta_0 E_0 n_2} \right), \quad \alpha = 0^\circ, \quad n_2 \text{ integer, but } n_2 \neq 2, 6, 10, \dots$$

$$(4) \quad x_d = 1 / \left( 1 + \frac{26,7}{\theta_0 E_0 n_3} \right), \quad \alpha = 90^\circ, \quad n_3 \text{ integer}$$

$$(5) \quad I = E_0 \cdot I_k(x, \theta_0 E_0) + I_i(x), \quad E_0 \text{ in GeV}$$

$$(6) \quad P = E_0 \cdot p_0(x, \theta_0 E_0) / I$$

$I_k$  and  $I_i$  are the coherent and incoherent parts of the intensity. (5) and (6) are valid for any value of  $\alpha$ . The reciprocal lattice points in the  $[001], [110]$  plane are numbered by indices  $n_2, n_3$  as shown in Fig.1. The polarisation is defined such that the electrical vector is parallel to the plane  $\vec{p}_0, [110]$ .

Given  $\theta_0, E_0$  and  $\alpha = 0^\circ$  or  $90^\circ$ , one can construct a spectrum by the aid of the plots 1 A to 6 A, 6 AB ( $\alpha = 0$ ), or 1 B to 5 B, 6 AB ( $\alpha = 90^\circ$ ), e.g.:

1 A gives the position of discontinuities  $x_d$  as function of  $\theta_0 \cdot E_0$  for a number of lattice rows  $n_2$ . The dashed curve in addition gives the x-value at which the maximum intensity of the first peak

has its half value. With  $x_d$  fixed,  $I_k$  becomes a function of  $\theta_o E_o$  only. 2 A to 6 A give the upper and lower intensities,  $\hat{I}_k$  and  $\check{I}_k$  for the value  $\theta_o E_o$  chosen. These have to be multiplied according to (5) and then  $I_i(x_d)$  must be added (plot 6AB). In this way the spectrum is defined by 2 intensity points for each step and by a half-intensity point on the slope of the first peak. At the high energy end the spectrum approaches the incoherent part 6AB, which is not dependent from  $E_o$  or  $\theta_o$ .

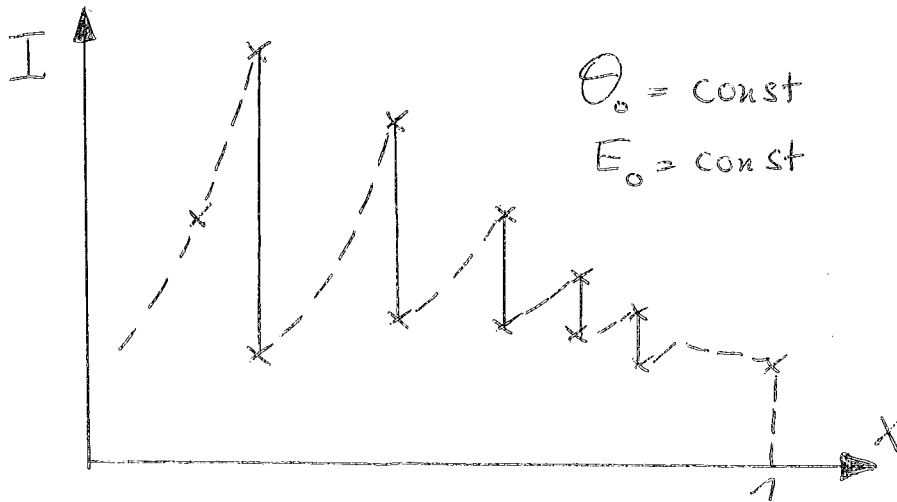


Fig.2 Spectrum constructed from the plots 1A,B-6A,B

The polarisation function  $p(\theta_o E_o)$  is given in 7 AB. With the intensity of the first peak now being known, it serves to calculate the polarisation of the photons in the first peak.

$\theta_o$ -dependence of intensity

For the alignment of the crystal, it is useful to know the intensity as a function of  $\theta_o$ , with  $x = x_o$  and  $E_o$  fixed. Again the plots 1 A, 1 B give the values  $\theta_d = \frac{\theta_o \cdot E_o}{E_o}$  for  $x = x_o$  at which the discontinuities appear. The intensity steps are taken from curves 2 A to 6 A, 2 B to 5 A, whichever case applies, as in the case of an energy spectrum. In addition one now has to use the plot 8 AB, which represents the total intensity, integrated over  $x$ , as function of  $\theta_o E_o$ :

$$(7) \quad I_T = \int_0^1 I(x, \theta_0 E_0) dx = E_0 \int_0^1 I_k(x, \theta_0 E_0) dx + \int_0^1 I_i(x) dx$$

$$= E_0 \cdot I_{kT}(\theta_0 E_0) + I_{iT}$$

This has the following reason: Experimentally, the intensity distribution of the coherent photon beam is measured with a pair spectrometer, taking the total intensity of the beam as monitor. Taking an energy spectrum,  $\theta_0 E_0 = \text{const}$ , the total intensity is constant. However it changes with  $\theta_0$  if a  $\theta$ -dependence is measured, such that the intensity distribution actually observed is:

$$(8) \quad \frac{I(\theta_0)}{I_T} = \frac{E_0 I_k(x_0, \theta_0 E_0) + I_i(x_0)}{E_0 I_{kT}(\theta_0 E_0) + I_{iT}}$$

The  $\theta$ -distribution is symmetrical to  $\theta_0 = 0$ .

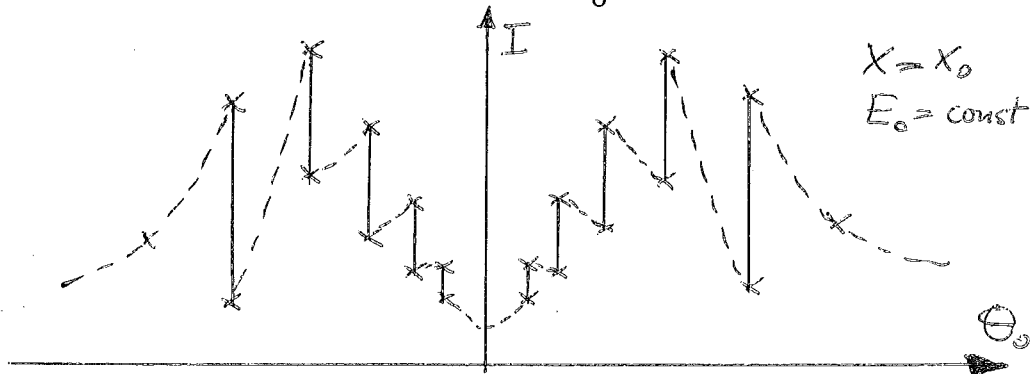


Fig.3  $\theta$ -Dependence of intensity from plots 1A,B-8A,B.

Contribution of the single lattice point (0,2)

In the general case  $0 \leq \alpha \leq 90^\circ$  the following relation gives the position of all possible discontinuities in the spectrum:

$$(9) \quad X_d = 1 / \left[ 1 + \frac{10^3}{\theta_0 E_0 (266 n_2 \cos \alpha + 377 n_3 \sin \alpha)} \right], \quad 0 \leq \alpha \leq 90^\circ$$

The intensity now does not allow a simple representation as in the cases  $\alpha = 0, 90^\circ$ , but it has been shown experimentally by

Bologna, Lutz, Schulz, Timm, Zimmermann (to be published) that orientations with small  $\alpha$  and large  $\theta_0$  are possible to allow the separation of certain single lattice points such that the intensity from this point dominates the spectrum, contributions from all other points being small.

The most favorable lattice point with respect to intensity and polarisation is  $(n_2, n_3) = (0, 2)$ ; this contribution is therefore also represented here. The position of this, single, discontinuity, the intensity and polarisation are given by the following formulas:

$$(10) \quad x_d = 1 / \left[ 1 + \frac{13,27}{\theta_0 E_0 \sin \alpha} \right], \quad 0 \leq \alpha \leq 90^\circ$$

$$n_2 = 0; \quad n_3 = 2$$

$$(11) \quad I = E_0 \cdot I_k(x_d) + I_i$$

$$(11') \quad I_k = \frac{1-x_d}{x_d} [1 + (1-x_d)^2] \cdot 7,36 \quad [GeV^{-1}]$$

$$n_2 = 0; \quad n_3 = 2$$

$$(12) \quad P = E_0 \cdot p(x_d) \cdot \cos \alpha / I$$

$$(12') \quad p = \frac{(1-x_d)^2}{x_d} \cdot 14,72 \quad [GeV^{-1}], \quad n_2 = 0; \quad n_3 = 2$$

The functions  $x_d$  ( $\theta_0 E_0 \sin \alpha$ ),  $I_k(x_d)$ , and  $p(x_d)$  for the single point (0,2) are plotted on pages 1 C, 2 C.  $I_k(x_d) = 0$  in this case because contributions from other lattice points are not included.

Given  $\theta_0$ ,  $E_0$  and  $\alpha$  one can find the position of the discontinuity  $x_d$  from 1 C. The curves 2 C then give the necessary information to calculate intensity and polarisation according to (11) and (12). The polarisation obtained from this single lattice point is higher than in the case  $\alpha = 0, 90^\circ$ . For small  $x_d$  the polarisation approaches

$$(13) \quad P \rightarrow \frac{1}{1 + \frac{I_i}{E_0} \frac{x_d}{(1-x_d)^2}} \rightarrow 1; \quad x_d \ll 1$$

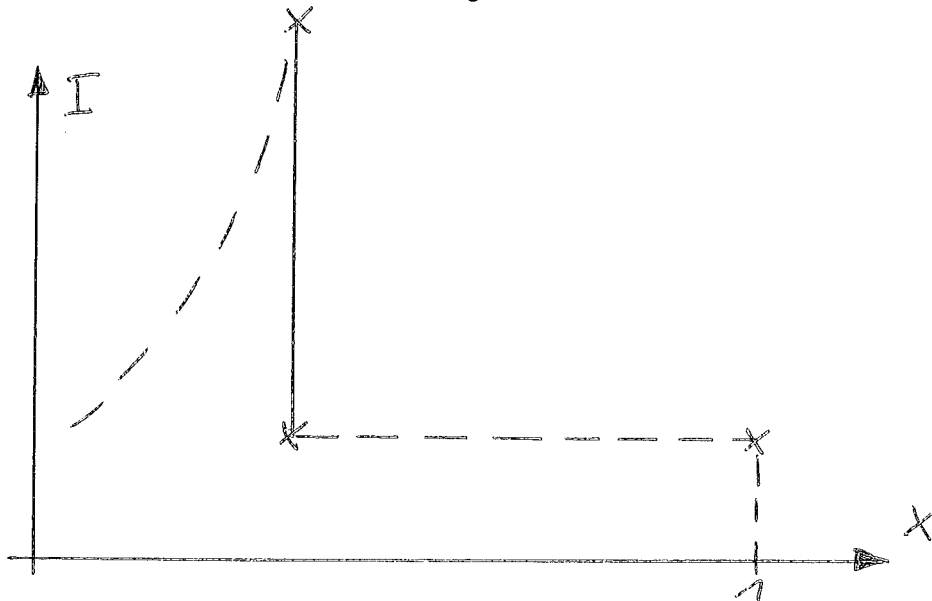


Fig.4 Spectrum from a single lattice point as constructed from 1 C, 2 C, 6 AB.

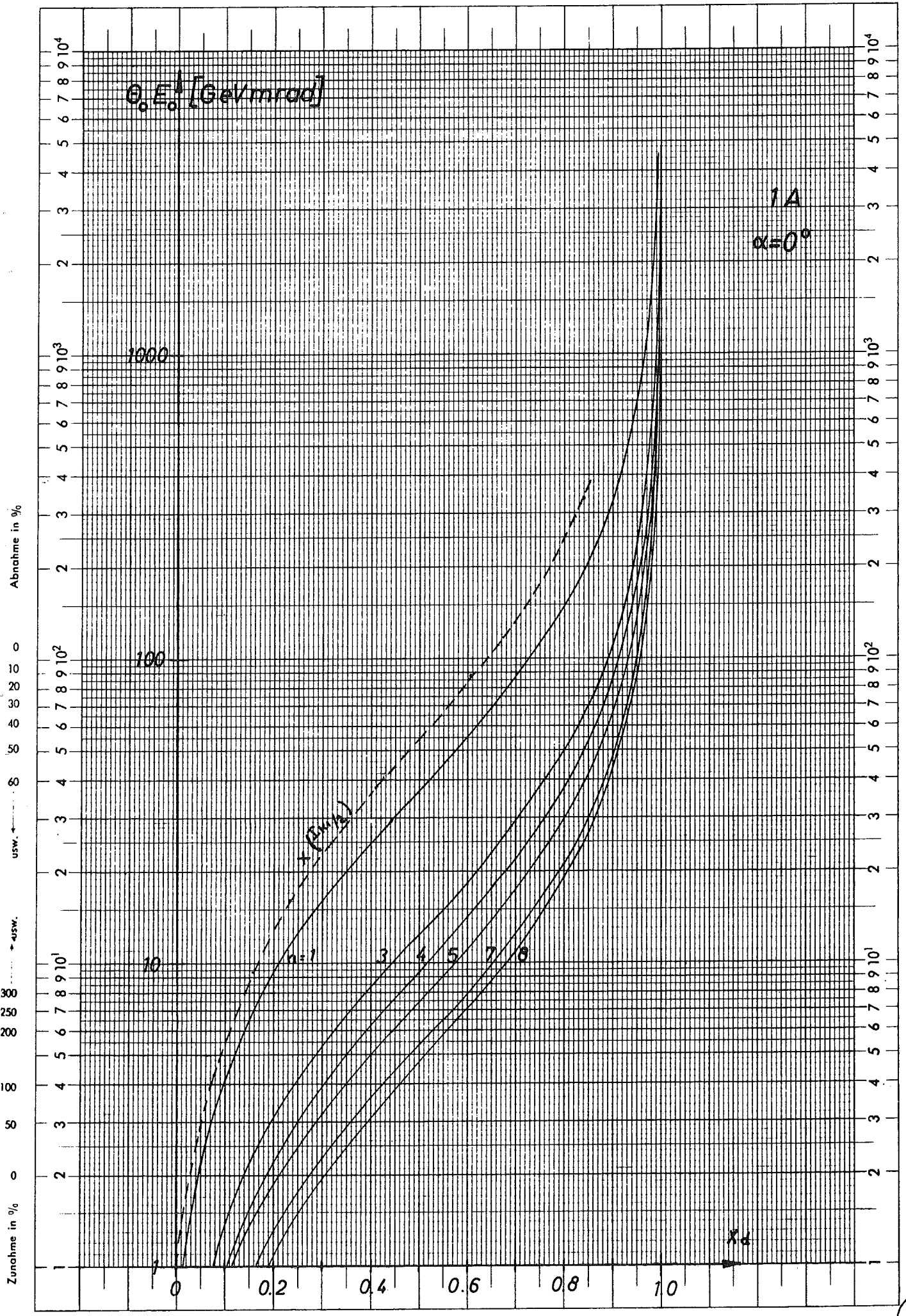
Final remarks

The spectral intensities obtained in this way do not, of course, take into account all the experimental effects which tend to smooth out the discontinuities such as primary divergence, multiple scattering, mosaic spread of the crystal and apparatus resolution.

The universal character of the intensity- and polarisation functions which are given here can be recognised in the formulas for the cross section of coherent Bremsstrahlung production which have been given by Barbiellini, Bologna and Murtas<sup>\*)</sup>.

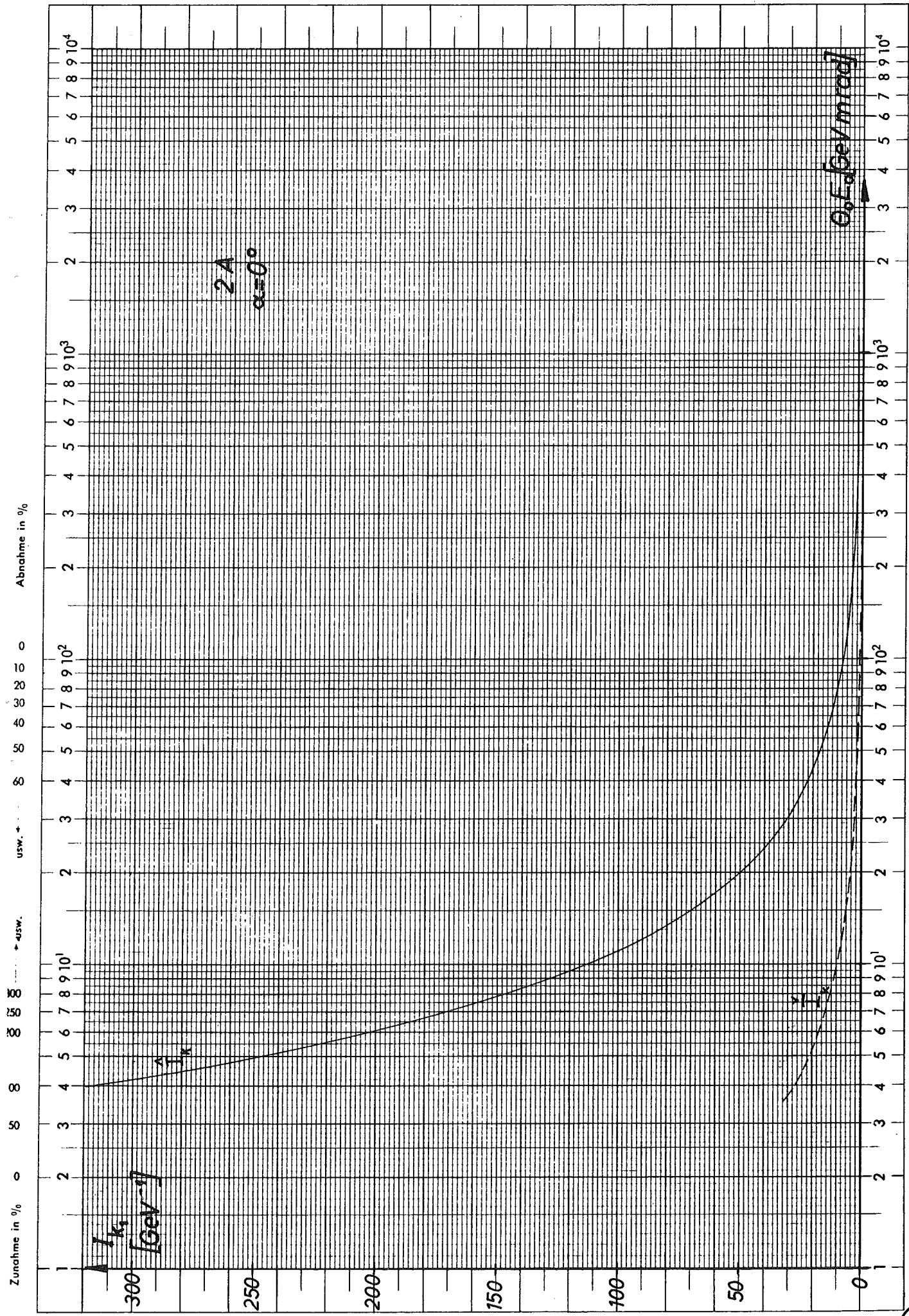
I thank Dr.G.Bologna und Mr.G.Lutz for discussions and Miss W.Kuffner for help in evaluating the data and for preparing the drawings.

<sup>\*)</sup> Phys.Rev.Lett. 8, 454, (1962), eqns. 3, 4 for  $\alpha = 0, 90^\circ$ .  
Nuovo Cim. 28, 435, (1963), eqn.3 for any  $\alpha$ .

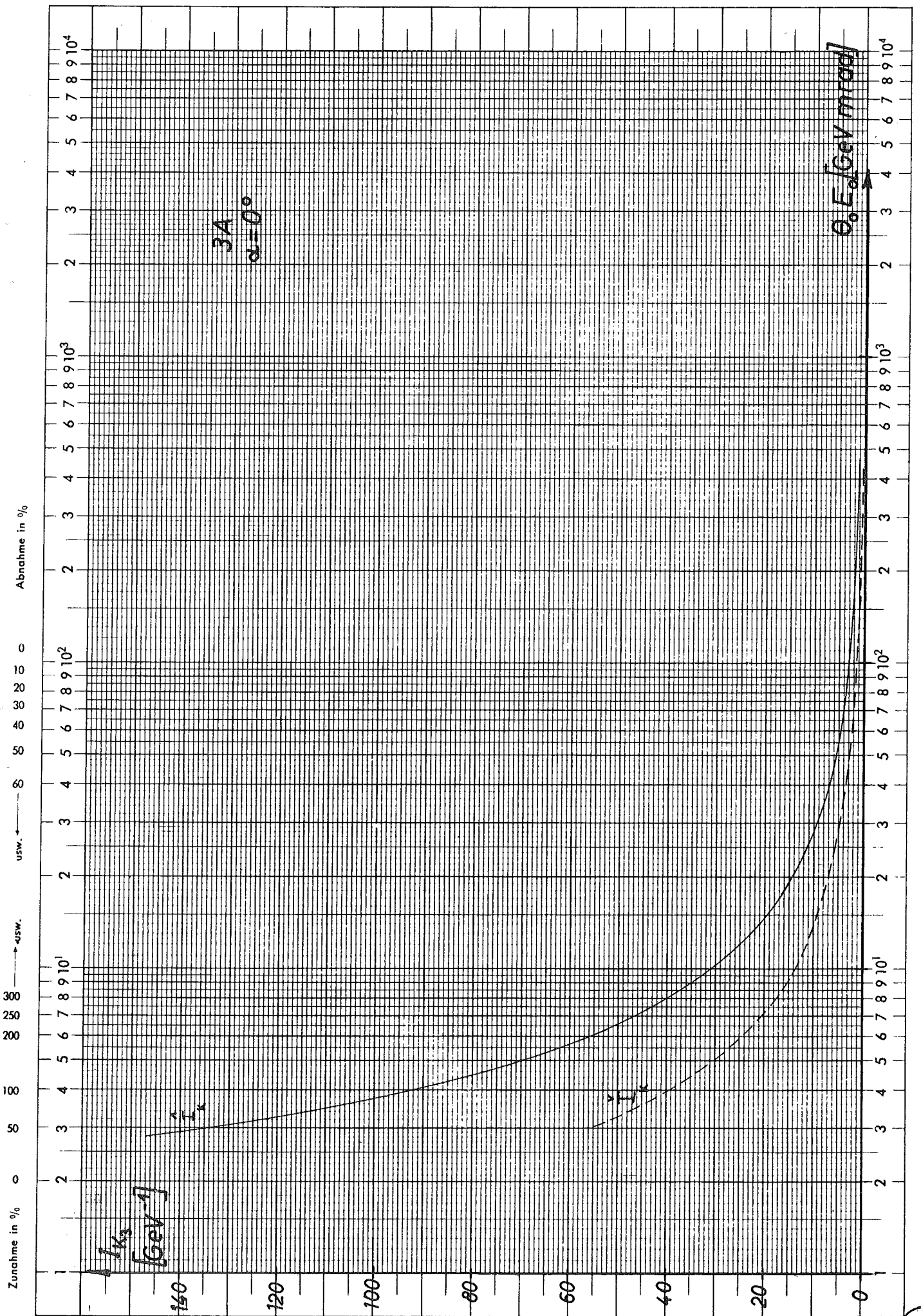


Eine Achse logar. geteilt von 1 bis 10000, Einheit 62,5 mm, die andere in mm mit Prozentmaßstab

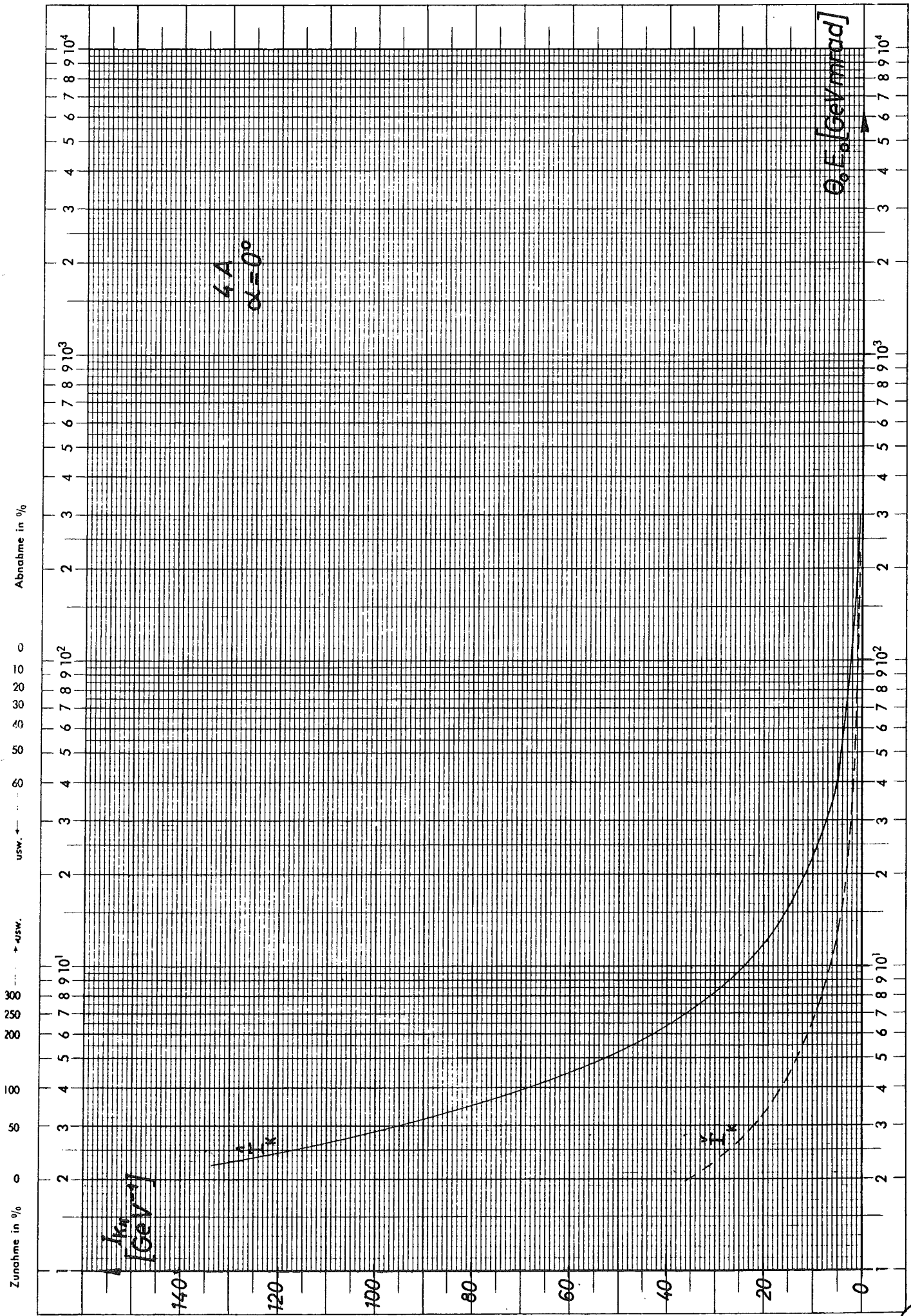




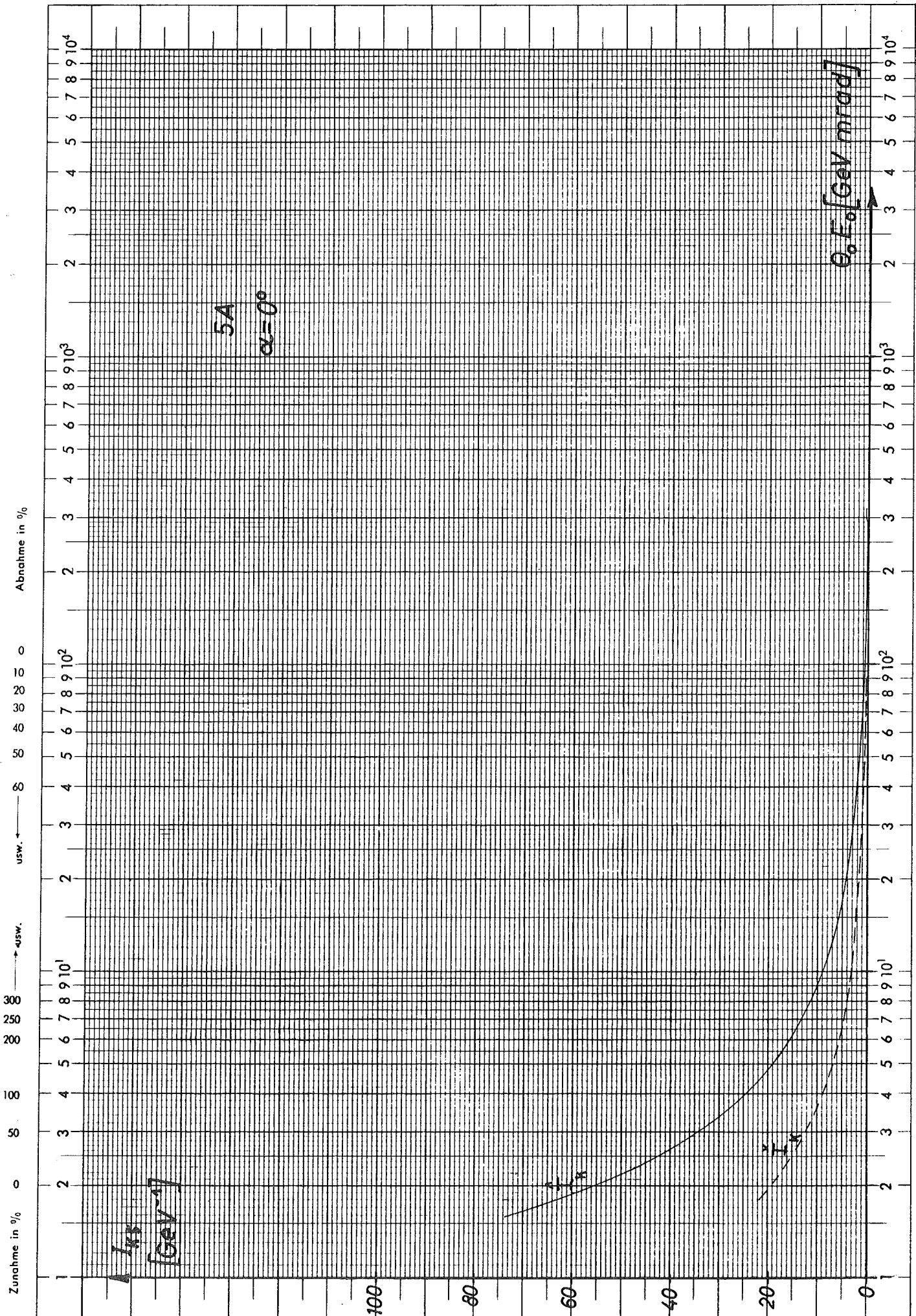
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Abnahme in %

0

10

20

30

40

50

60

usw. ←

→ usw.

200

250

300

4

5

6

7

8

9

10<sup>1</sup>

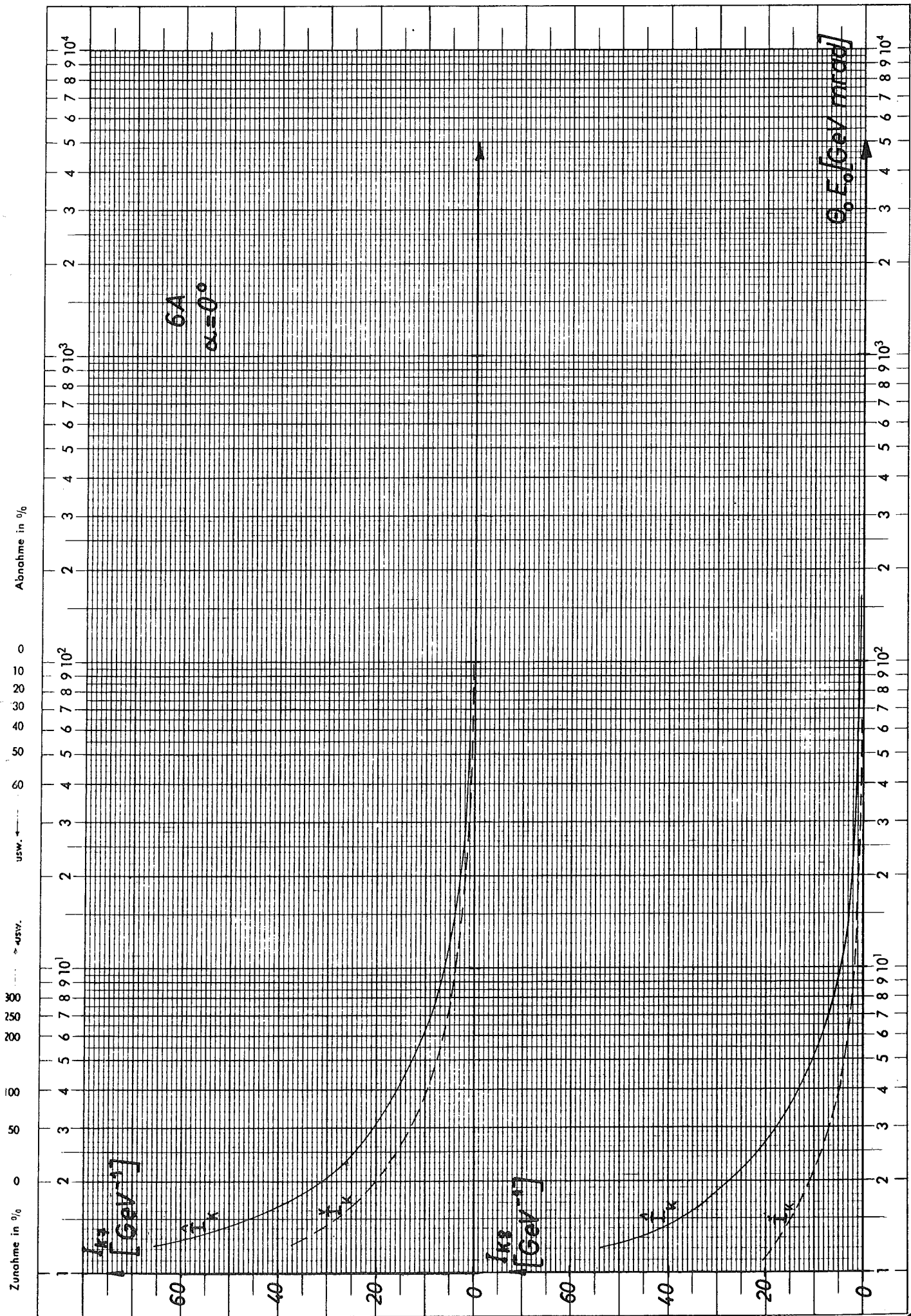
$I_k$

[GeV<sup>-1</sup>]

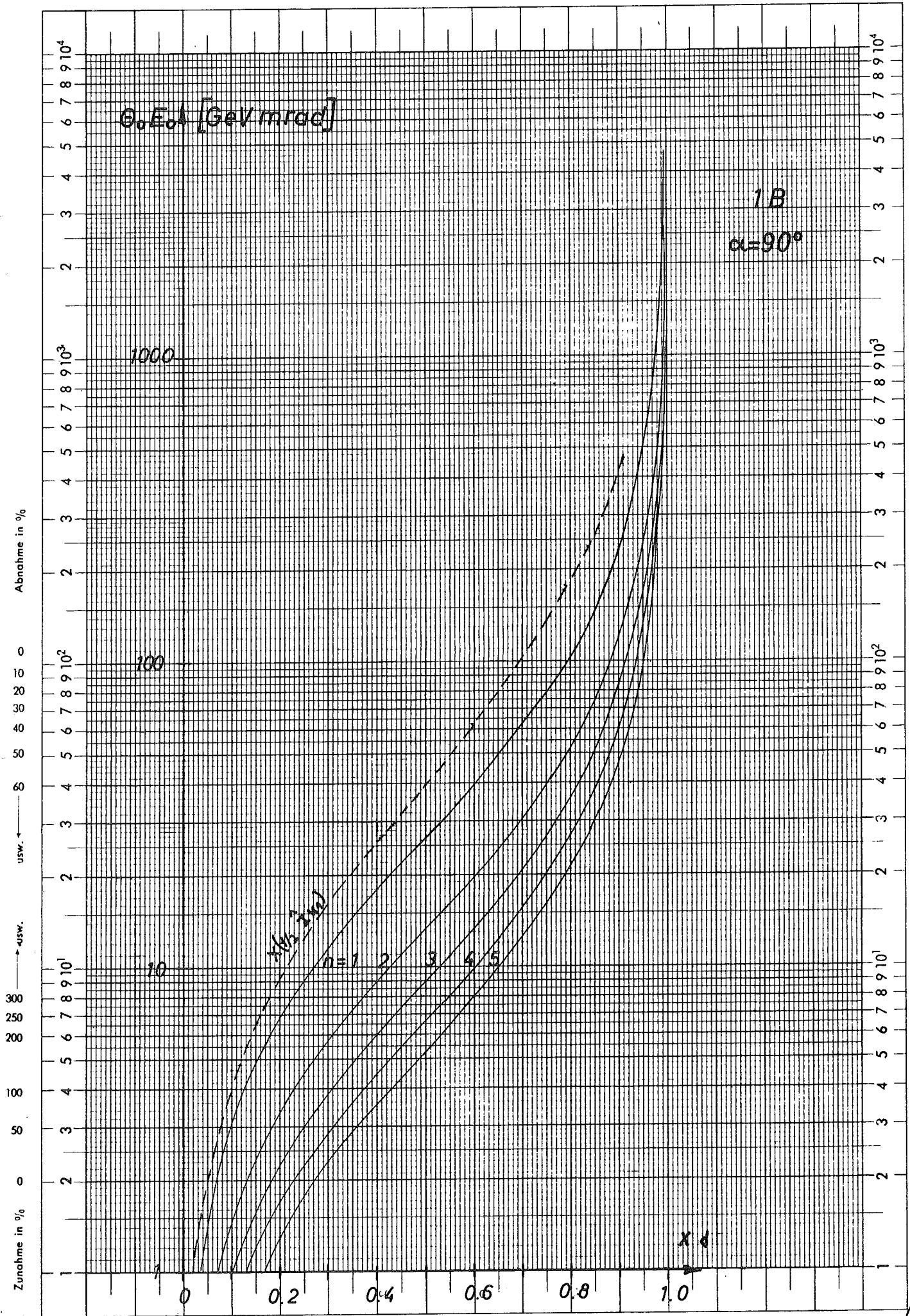
5A

$\alpha = 0^\circ$

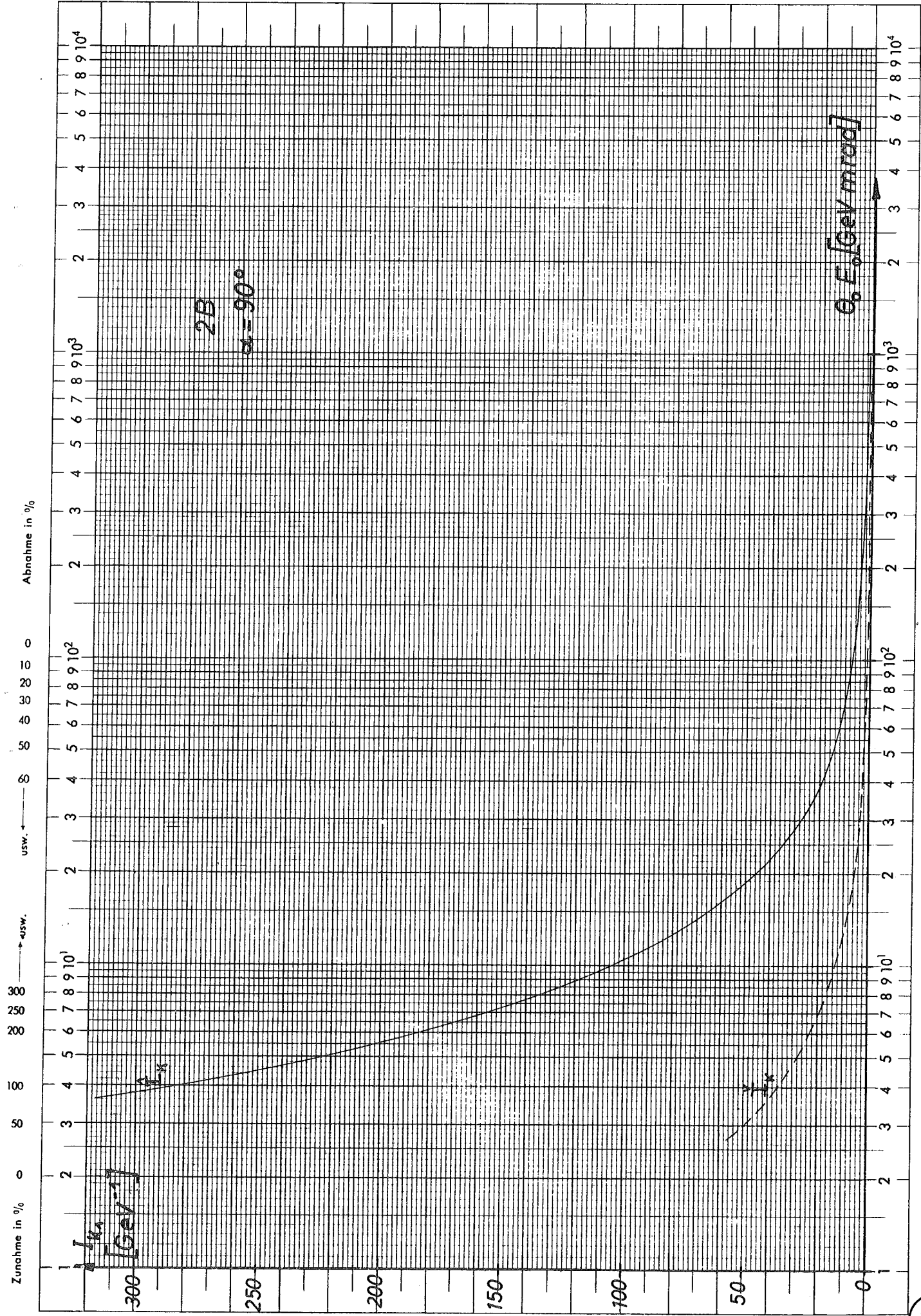
$\theta_0 F_0$  [GeV mrad]



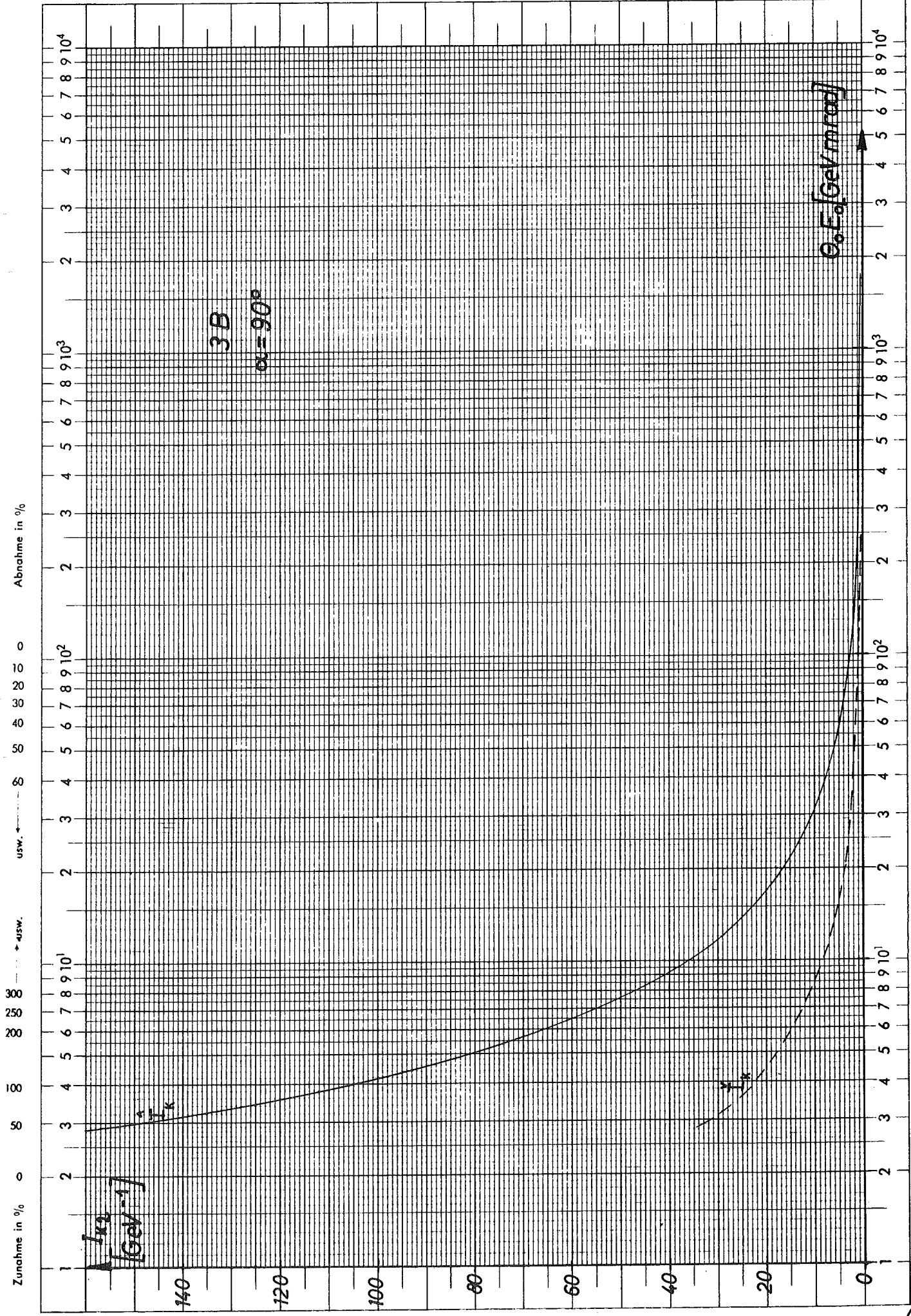
Eine Achse logar. geteilt von 1 bis 10000, Einheit 62,5 mm, die andere in mm mit Prozentmaßstab



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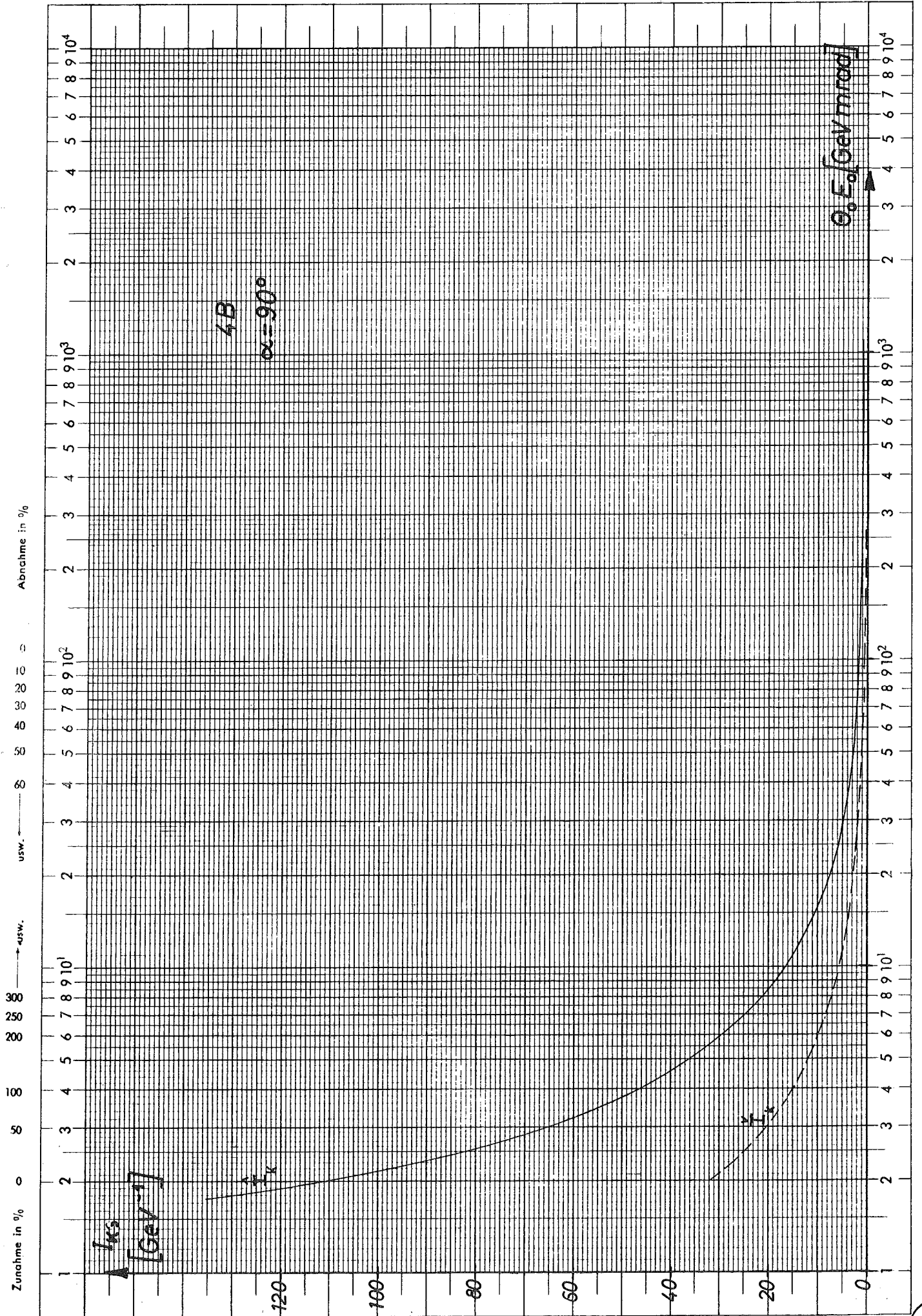


Eine Achse logar. geteilt von 1 bis 10000, Einheit 62,5 mm, die andere in mm mit Prozentmaßstab

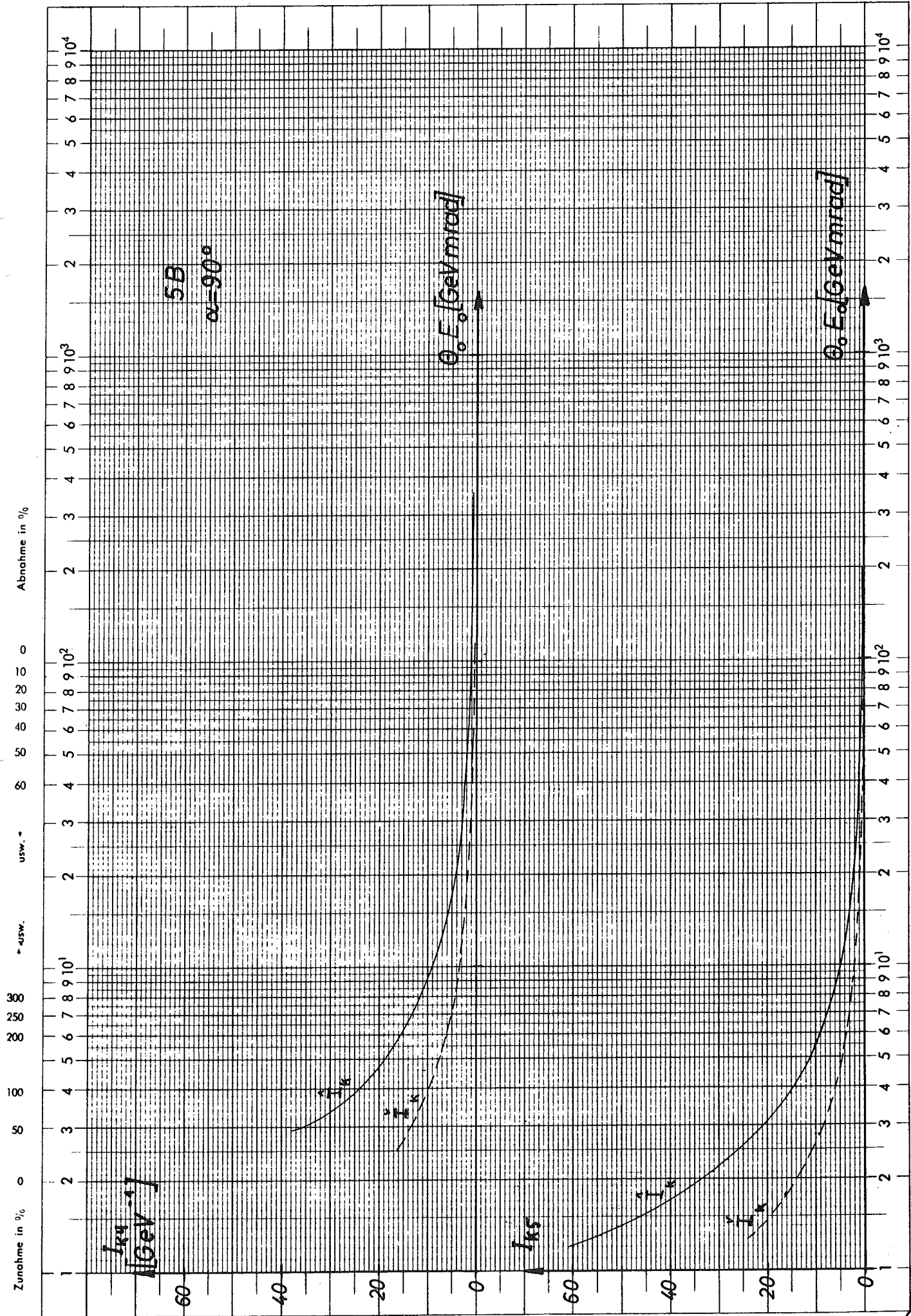


Eine Achse logar. geteilt von 1 bis 10000, Einheit 62,5 mm, die andere in mm mit Prozentmaßstab

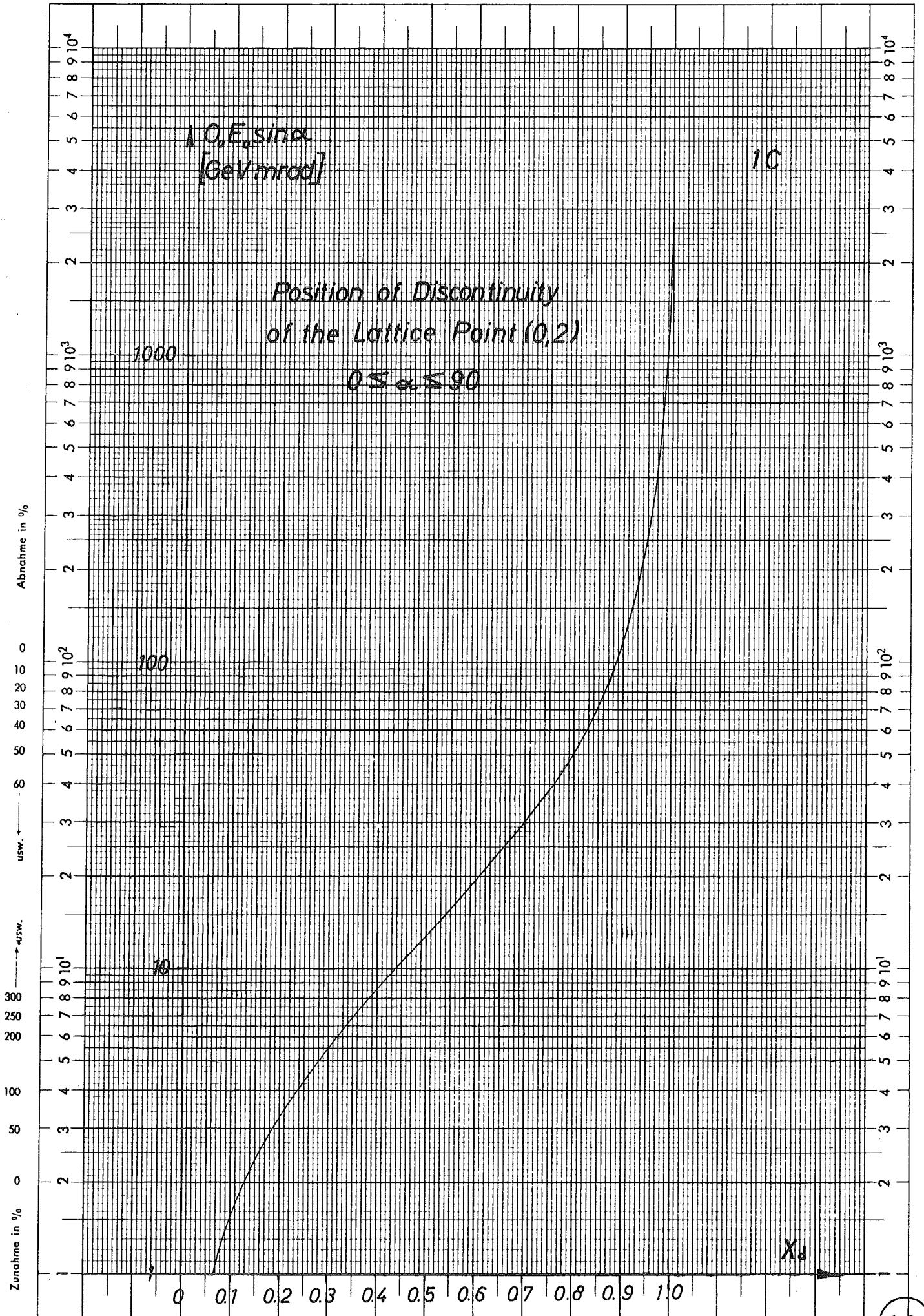




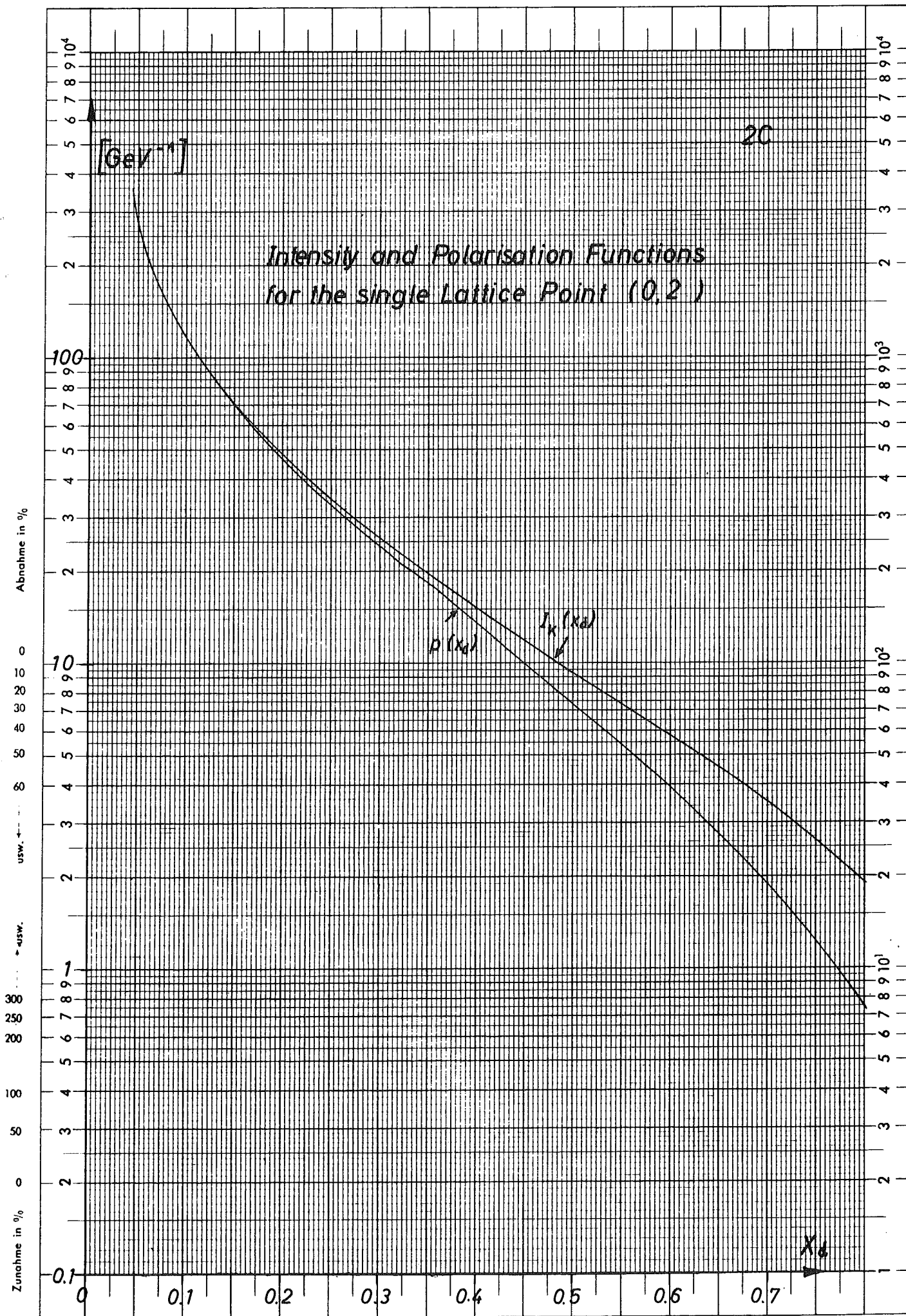
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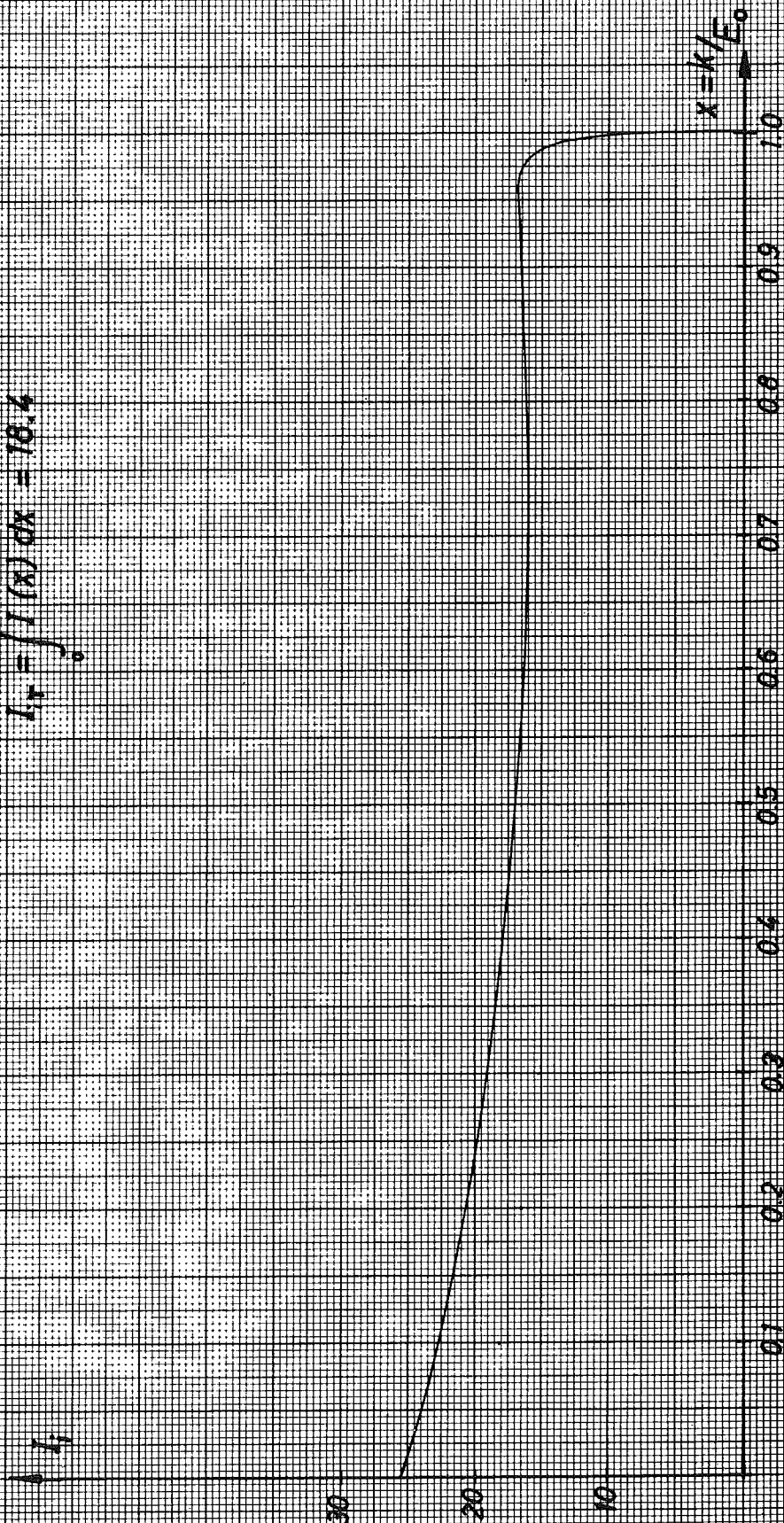


Eine Achse logar. geteilt von 1 bis 10000, Einheit 62,5 mm, die andere in mm mit Prozentmaßstab

6AB

Incoherent Contribution

$$I_T = \int_0^1 I(x) dx = 18.4$$



A O.E.  
[GeV mrad]

7AB

Polarisation Function

130  
120  
110  
100  
90  
80  
70  
60  
50  
40  
30  
20  
10  
0

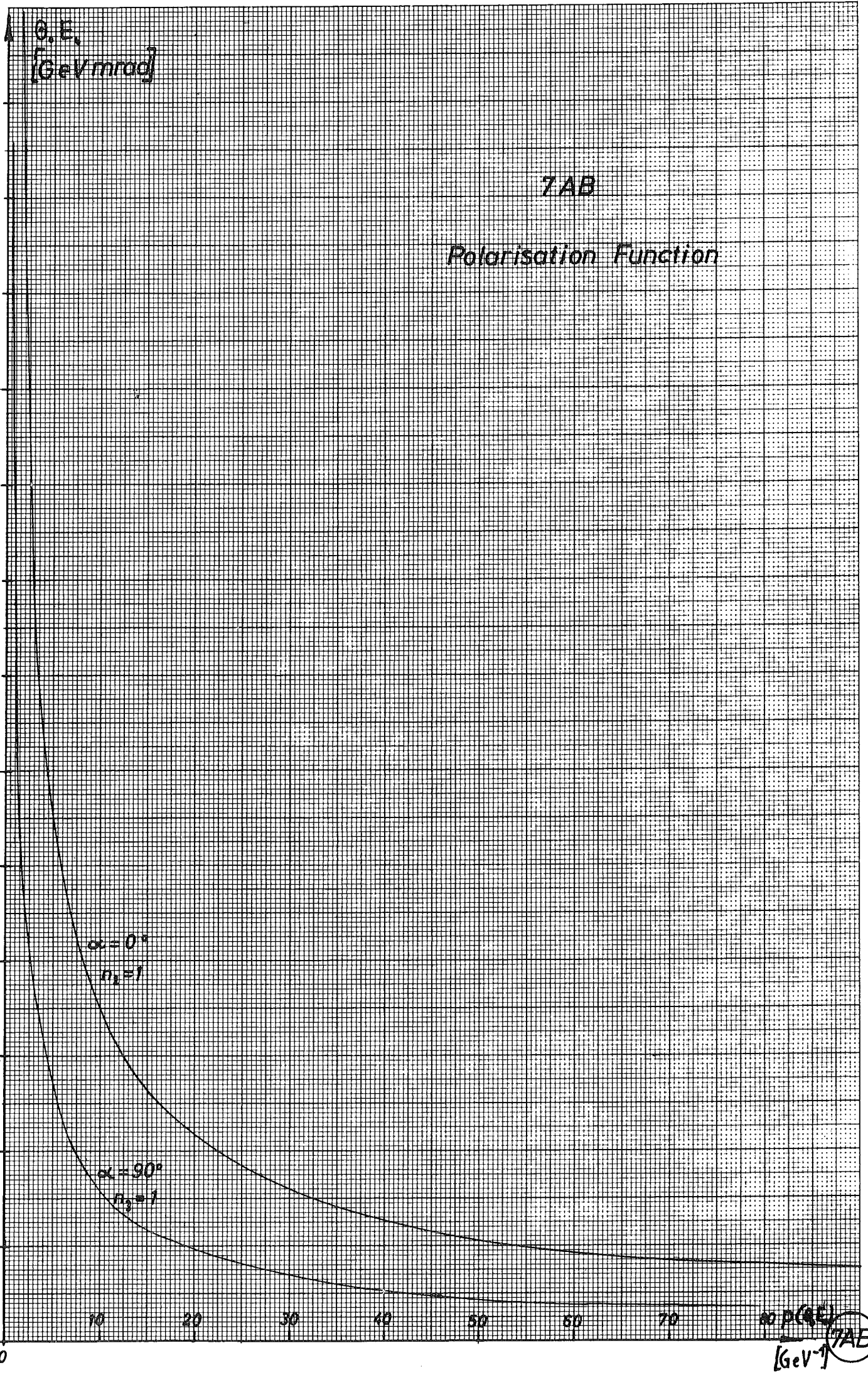
$\alpha = 0^\circ$   
 $n_2 = 1$

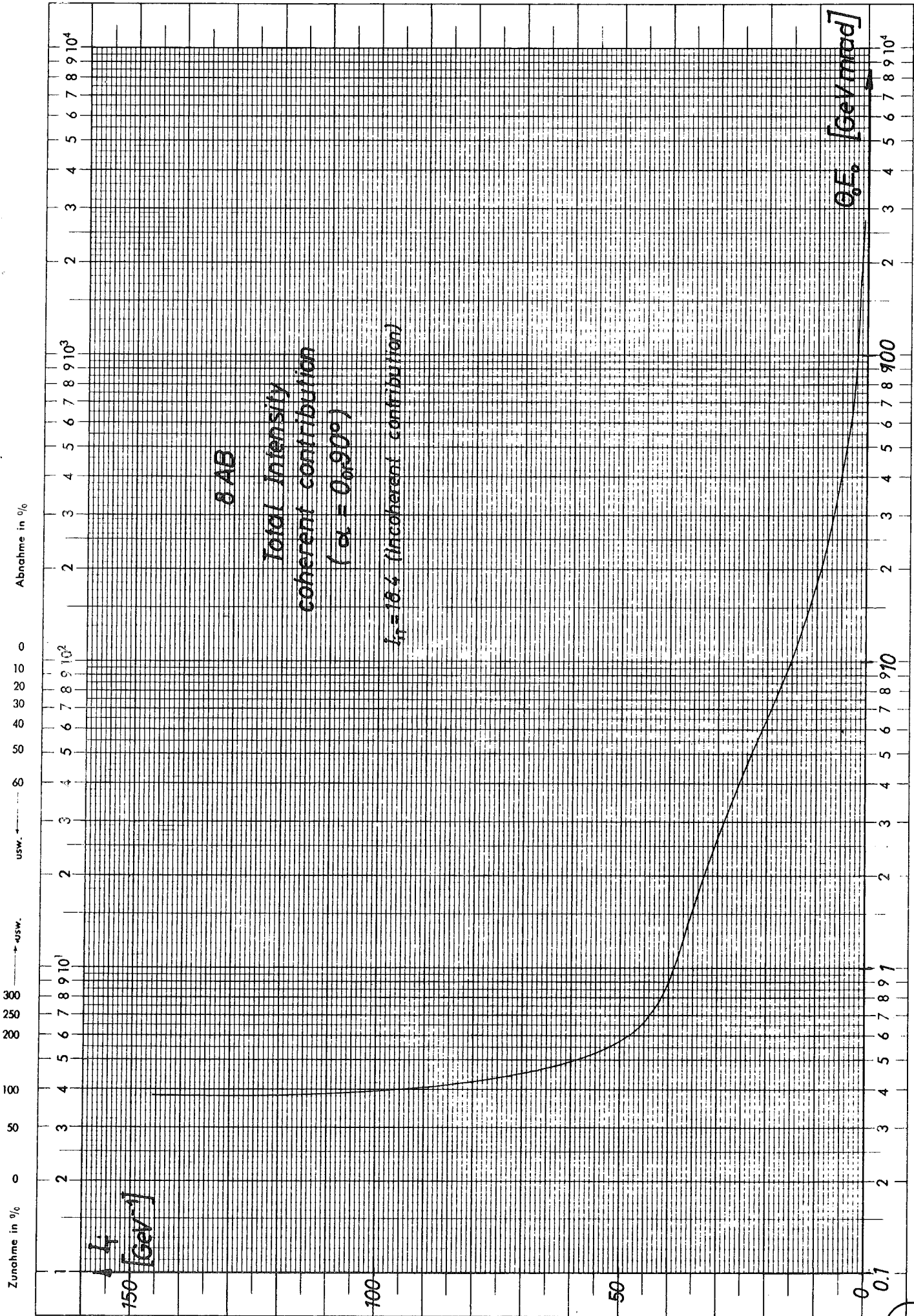
$\alpha = 90^\circ$   
 $n_3 = 1$

10 20 30 40 50 60 70 80

80 p (GeV)  
[GeV<sup>-1</sup>]

7AB





Eine Achse logar. geteilt von 1 bis 10000, Einheit 62,5 mm, die andere in mm mit Prozentmaßstab