THE VARIANCES AND COVARIANCES OF 17.5 - 4MEASURED INTENSITIES IN PRECISE LATTICE-CON-STANT DETERMINATION BY THE BOND METHOD. By E. Gałdecka, Institute for Low Temperature and Structure Research, Polish Academy of Sciences, pl. Katedralny 1, 50-950 Wrocław, Poland.

Precision of measurements on the Bond diffractometer depends on several factors as discussed in my papers (Urbanowicz E., (1981) . Acta Cryst. A<u>37</u>, 364-368 and 369-373). In theoretical considerations it is often assumed that the var-iance of measured intensity is equal to number of counts. This result from Poisson statistics and from the assumption that the only source of attistical errors is counting statistics and all the other factors are negligible. In present paper beside counting statistics the following parameters which influence the statistical errors are taken into account: a)  $\mathcal{G}(V)/V$  - fluctuation of the power voltage, b)  $\mathcal{G}(\omega)_{M}$  - statistical angle-reading error,

c)  $G(\Delta \omega)$  the error in positioning of the angle by stepping motor. As a result, the following formula for the var-

iance of measured intensities was derived:  $G^{2}(h_{i}) = e_{1}h_{i} + e_{2}h_{i}^{2} + G^{2}(\omega_{i})(h_{i})^{2}$ where:  $h_{i}$  - the number of counts in i-th posi-

tion of the sample (i = -p,...,p and 2p + 1 - the number of the scanning points),  $\omega_1$  - angle position of the sample,

 $h_1^*$  - the first derivative  $\partial h/\partial \omega$  in the i-th position.

$$\tilde{G}^{2}(\omega_{1}) = e_{3} + i^{2}e_{4}, e_{1} = 1, e_{2} = \tilde{G}^{2}(V)/V^{2},$$

$$\mathbf{e}_{3} = \left[\mathbf{G}^{2}(\boldsymbol{\omega})_{\mathsf{M}} + \mathbf{p}\mathbf{G}^{2}(\Delta\boldsymbol{\omega})\right]/2,$$

 $\mathbf{e}_{4} = [G^{2}(\omega)_{\Pi} - pG^{2}(\Delta \omega)]/(2p^{2}).$ 

An analysis of the covariances of angle posi-tions  $\omega_{\rm i}$  of the sample leads to the conclusion that in the case of automatic scanning the covariances between the values  $\omega_{1}$  and  $\omega_{1}$ 

 $(i, j = -p, \dots, p \text{ and } i \neq j)$  cannot be neglected, as their values given by the formula

$$\operatorname{cov}(\omega_{i}, \omega_{j}) = a_{3} + i j a_{4} - |i - j| a_{5},$$

where  $e_5 = 6^2 (\Delta \omega)/2$ ,

are comparable with the values of the variances.

Covariances between  $\omega_i, \omega_i$  (angle positions) result in correlations between measured intensities h<sub>i</sub>, h<sub>j</sub>:

$$cov(h_{i},h_{j}) = h_{i}h_{j}cov(\omega_{i},\omega_{j})$$

The model of variances and covariances of the measured intensities has been verified on the basis of several series of experimental data. Experimental conditions were discussed when: 1) we can take  $G^2(h_i) = h_i$ ,

- 2) fluctuations of the power voltage are
- essential, dispersion of position angles of the sample 3) cannot be neglected.

17.5-5 PSEUDOTRANSLATIONS AND STATISTICS. By G. Cascarano, C. Giacovazzo & M. Luić, Dipartimento Geomineralogico, University, 70121 Bari, Italy.

Pseudotranslation symmetry frequently occurs in crystal structures. Their effects on the structure factor statistics has been studied by different authors in some aspects and in few space groups. It seems interesting to describe a gene ral theory which provides the distribution of the structure factors in any space group for any combination of the pseudotranslations. An important complementary problem is that of deriving, from the observed statistical effects, information about the pseudotranslations : for example the nature of the pseudotranslations, the percentage of atoms which are related by the pseudotranslations, .... Such an information is of primary importance in the field of direct methods, where the particular features of the structure have to be taken into account.

EFFECT OF HEAVY ATOMS ON INTENSITY 17.5–6 STATISTICS OF X-RAY DIFFRACTION FROM CRYSTALS. By D. Pradhan, S. Ghosh and <u>G. D. Nigam</u>, Depart-ment of Physics, Indian Institute of Techno-logy, Kharagpur-721 302, India.

It is well known that if the unit cell con-tains one or few heavy atoms in addition to a large number of light atoms, the distri-bution of x-ray intensities usually deviate considerably from those inferred by Wilson statistics. In such cases, strong correlation between atomic positions results from the finite radii of atoms, which prevents complete randomness in the distribution of atoms in A modified cumulative distribution function

of normalised intensities is derived for a triclinic cell when heavy atom occupies special position in the unit cell. The distribution being considered, the remaining light atoms would no longer be randomly distributed in the midst of heavy atom in the <u>whole</u> unit cell. It corresponds, in fact, to an inaccessible volume, which is assumed spherical, around the heavy atom with radius equal to the sum of the radius of heavy atom and the aver-age radius of light atoms. This inaccessible volume would modulate the average intensity and the distribution parameter as a function of s. Numerical study is made to investigate the effect of the inaccessible volume on the cumulative distribution function. The results are compared with experimental distributions for rubidium di-o-nitrobenzoate and potassium hydrogen di-o-nitrobenzoate.