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| Author Inst Title | : Blokhin. M.A. : Rostov-on-Don University, USSR : Calculation of the Distortion of X-ray Spectra Due to the Apparatus. |
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| Orig Pul | o: Dokl. AN SSSR, 1956, 107, No 2, 229-232 |
| Abstract | t : Description of a method permitting the distortion of an x-ray spectrum of any shape and to use this method further on to restore the true shape of the spectrum. The method is based on the use of trigonometric series and can also serve to obtain the shape of the density distri- bution curve of the free and filled electronic states. In the latter |
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| | case it is necessary to carry out additional correction of the spec- trum for "distortion" introduced by the internal level of the atom. |
| | case it is necessary to carry out additional correction of the spec- trum for "distortion" introduced by the internal level of the atom. |

PHASE I BOOK EXPLOITATION 1073

Blokhin, Mikhail Arnol'dovich

Fizika rentgenovskikh luchey (Physics of X-rays) 2d. Ed. Rev. Moscow, Gostekhizdat, 1957. 518 p. 7,000 copies printed.

Ed.: Kuznetsova, Ye.B.; Tech. Ed.: Gavrilov, S.S.

FURPOSE: This is a textbook on the physics of X-rays and X-ray spectral analysis. It presumes a knowledge of electrodynamics, basic quantum mechanics, atomic theory and differential and integral calculus.

COVERAGE: The author presents the general principles of the physics of X-rays necessary for a further study of specialized fields involving X-rays. The author states that the approach of this book is especially suited for the experimental physicist. He deals in detail with the fine structure of X-ray spectra and the energy levels of atoms, paying particular attention to the physical aspects. The author states that X-ray analysis of metals, alloys, minerals, etc. is widely used in the USSR and lists the following scientists as working in this field: Academician, A.F. Ioffe, G.V. Kurdyumov, N.V. Belov, S.T. Konobeyevskiy, N.Ye. Uspenskiy, G.S. Zhdanov, A.I. Kitaygorodskiy, N.V. Ageyev, Ya.S. Umanskiy, B.M. Rovinskiy, V.I. Iveronova, Yu.S. Terminasov. The fol-

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Physics of X-rays

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lowing scientists have developed new methods and equipment used in chemical X-ray spectral analysis: I.B. Borovskiy, V.N. Protopop, E.Ye. Vaynshteyn, N.D. Borisov, Ya.M. Fogel', and M.A. Blokhin. A.I. Kostarev and E.Ye. Vaynshteyn are mentioned as contributors to the theory of absorption spectra. I. B. Borovskiy, E.Ye. Vaynshteyn, K.I. Narbutt, R.L. Barinskiy, Ya.M. Fogel', A.I. Krasnikov, and M.A. Blokhin are mentioned for their research in the fine structure of emission spectra and absorption spectra. P.I. Lakirskiy, A.I. Alikhanov, G.V. Kurdyumov, and A.P. Komar are mentioned in conjunction with the Institute of Physics and Technology of the Academy of Sciences of the USSR as having contributed to the field of X-ray spectral analysis. There is an extensive bibliography of both Soviet and non-Soviet sources at the end of the book. The author states that the first edition of this book was used as a textbook at Rostov University.

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BLOITH IN, MA. BLOKHIN, M. A., NESTERENKO, P.S., and SHUVAYEV, A.T (RGU)

"X-ray Spectral Investigatian of Sulphur-Containing Samplesd"

Materials of the 2nd All-Union Conference on X-ray Spectroscopy; Moscow, January, 31 February 4, 1957 (Materialy II Vsesoyuznogo soveshchaniya po rentgenovskoy spektroskopii; Moskva, 31 yanvarya - 4 fevralya g.)

Izvestiya Akademii nauk SSSR, Seriya fizicheskaya 1957, Vol 2, Nr 10, pp 1341n - 1342 (USSR)

Assoc: RGU

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48-10-2/20

Materials of the 2nd All-Union Conference on X-ray Spect oscopy; Moscow, January 31, to February 4, 1957

A. I. Kozlenkov (Fizfak MGU); Interrelationship of Some X-ray Spectral and Magnetic Characteristics of Iron-Base Alloys by S. A. Nemnonov and K. M. Kolabova (UFAN SSSR); Investigation of Binding Forces in Solid Iron-Molybdenum Solutions According to the Fine Structures of X-ray Absorption Spectra by V. A. Trapeznikov and S. A. Nemnonov (UFAN SSSR); On the Theory of Solid Solutions Based on Transitional Metals by I. B. Borovskiy and K. P. Gurov (IMET AN SSSR); Relationship of Temperature an Concentration of Fine Structure of X-ray Absorption Spectra of Solids and an Investigation of Binding Forces by V. A. Trapeznikov; Investigation of X-ray L-Spectra of Some Rare-earth Element Compounds by N. V. Troneva, I. D. Marchukova and I. B. Borovskiy (Fizfak MGU); Investigation of X-ray Emission K Lines of B-Group Titanium in Carbides and Some Other Compounds by E. Ye. Vaynshteyn and Yu. N. Vasil'yev (GEOKhI AN SSSR); X-ray Spectral Investigation of Molybdenum L Spectra in Some Alloys and Compounds by V. A. Batyrev, I. B. Borovskiy and S. A. Ditsman (IMET AN SSSR); Some Satelites of Spectral Lines by T. I. Kakushadze (Georgian Teacher's Institute); X-ray Spectral Investigation of Sulphur-containing Samples by M. A. Blokhin, P. S. Nesterenko and A. T. Shuvayev (RGU).

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CIA-RDP86-00513R000205530008-8

Blockhin, M.A. 48-10-3/20 Blokhin, M.A., Sachenko, V.P. AUTHOR: The Breadths of the Internal Levels and the Distribution of Electron TITLE: States According to the Energies of the Elements of the Iron Transition Group (Shiriny wutrennikh urovney i raspredeleniye plotnosti elektronnykh sostoyaniy po energiyam elementov perekhodnoy gruppy zheleza) Izvestiya Akad.Nauk SSSR, Ser.Fiz., 1957, Vol. 21, Nr 10, PERIODICAL: pp. 1343-1350 (USSR) The attempt is made here to determine the breadth of the K-levels ABSTRACT: of some elements by means of the interpolation on these elements of the experimental values for the K-level breadths of intransitive elements (Mg, Al, Ar, K, Fe). When detecting the changes of the breadth of K-levels by means of the atomic number two essentially different types of transition, the radiation- and the radiationless transitions, were taken into account. The entire probability for the emergence of the atom from the given state is equal to the sum of the probability of the radiation transition P_p and the radiationless P_1 . The level breadth $\gamma = A(P_p+P_1)$. A - coefficient, the value of which is determined by the shape of the level. The ratio of these probabilities can be computed also experimentally. It is shown that the com-Card 1/2

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48-10-3/20

The Breadths of the Internal Levels and the Distribution of Electron States According to the Energies of the Elements of the Iron Transition Group

> puted data agree with the experimental ones. It is further shown that the dependence $P_p(Z)$ may be approximatively well expressed by $P_p=DZ^{4,71}$. In order to obtain the probability values for radiationless transitions, computed in approximation, for various elements, these probabilities were calculated for K-transitions. It is shown that the coefficient $D = 1,00.10^{\circ}$ transitions per second. According to the curve given here, the γ -values for the various elements are found. According to the breadths of the $K_{cl_{1,2}}$ lines, it is possible, if the breadth of the K-levels is known, to compute the breadth of the L_{II} - and L_{III} levels. The curves obtained for the densities of the electron states cannot be interpreted uniquely. The most acceptable interpretation is that by Beeman, W.W., and H. Friedman (Phys.Rev. 56, 392, 1939). There are 4 figures, 5 tables and 44 references, 6 of which are Slavic.

ASSOCIATION: Rostov State University (Rostovskiy gosudarstvennyy universitet)

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| • | 24(2),24(7) AUTHORS: | Blokhin, M. A., Shuvayev, A. T. SOV/48-22-12-10/33 |
|---|-------------------------|---|
| | TITLE: | Investigation of Compounds With the Structure of Perovskite by Ecentgen Spectra (Issledovaniye soyedineniy so strukturoy perovskita po rentgenovskim spektram) |
| | PERIODICAL: | Izvestiya Akademii nauk SSSR. Seriya fizicheskaya, 1958; Vol 22; Nr 12, pp 1453-1455 (USSR) |
| | ABSTRACT: | Emission and absorption K spectra of Ti, emission and absorption L spectra of Zr, the absorption K spectrum of Fe and absorption L spectra of Sr and Ba as well as of BaTiO ₃ , SrTiO ₃ , SrFeO ₃ , |
| | | BaZrO, and PbZrO, were investigated in the present paper. It |
| | Card 1/2 | was ascertained that the band of valence electrons of compounds having a percessite structure, is a hybrid band with a strong admixture of p-states. The formation of a generalized conduction band was confirmed experimentally. On principle, the conduction band has a d-character with a small admixture of p-states. The following band of free states has, on principle, a p-character with an irregular distribution of the state densities. The Ti-ions charge in BaTiO ₃ amounts by no means to more than 2.7. |
| | Jaru 172 | |

| | Investigation of Perovskite | of Compounds With the Structure by Rosatgen Spectra | SOV/48-22-12-10/33 |
|-------|------------------------------------|---|---|
| : | | In the proximity of the Curie (Kyuri) po of the electron states of the conduction There are 3 figures and 7 references, 1 | oint an energy decrease n band was ascertained. of which is Soviet. |
| | ASSOCIATION: | Rostovskiy-na-Donu gos. universitet (Rostov-na-Donu State University) | |
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24(4)

PHASE I BOOK EXPLOITATION

SOV/3040

Blokhin, Mikhail Arnol'dovich

Metody rentgeno-spektral'nykh issledovaniy (Methods of X-ray Spectrum Analysis) Moscow, Fizmatgiz, 1959. 386 p. 5,000 copies printed.

Ed.: V. I. Rydnik; Tech. Ed.: S. S. Gavrilov.

PURPOSE: This book is intended for scientific and laboratory workers, engineers, and technicians in x-ray spectra research and analysis. It will be of interest to advanced students of the field at vuzes.

doverAGE: This book treats the field of x-ray spectroscopy and discusses the use of apparatus and equipment in x-ray analysis, experimental studies of the fine structure of x-ray emission and absorption spectra, and processing of data. It describes the theory and practice of quantitative and qualitative x-ray analyses of alloys, ores, and minerals. It lists a number of Soviet institutes at which experimental work in x-ray spectroscopy is

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| Methods of X-ray (Contd) | SOV/3040 | |
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| carried out and mentions N. D. Bor Borovskiy, V. N. Protopopov, Ye. J K. I. Narbutt, and A. B. Gil'varg field of study. There are 433 refe Soviet, 38 German, 24 French, 5 Swe | L. Kostrov, E. Ye. Vayn as having contributed to | shteyn, |
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|-----|--------------|--|--|--|
| | AUTHORS | Blokhin, M. A., Busler, I. P. | , I. V., Kramarov, O | . P., Chernyavskaya, |
| | TITLE: | The Use of a Monitor in monitora pri rentgeno-s | | |
| | PERIODI | CAL: Pribory i tekhnika (USSR) | eksperimenta, 1959, | Nr 1, pp 106-111 |
| · · | ABSTRAC | T: In the continuous re by means of ionisation ility source of the rad it is difficult to ensu age at the relatively h | Or scintillation cou liation is necessary, ure a high stability high power used by th | Inters, a high stab- At the same time in the anode volt- ne tube. This prob- |
| | | lem is particularly con varied within wide limit the intensity ratio of this and other reasons ods for measuring line sity is directly stabil | its, for example, in a very weak and a ve the present authors intensity ratios eit | the measurement of ery bright line. For have developed meth- ther when the inten- |
| × 1 | | ion is not stabilized a counters are used for t | at all. Ionisation of this purpose as monit | chambers or geiger cors. The device is |
| • | Card 1/ | shown diagrammatically anode of the X-ray tube jincident on the specime | in Fig 1. In this f | figure 1 is the |

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The Use of a Monitor in X-Ray Spectral Analysis

ional specimen 3 . Fluorescence radiation leaving 2 is analyzed in a spectrometer which uses a geiger counter as the detector. The radiation from the additional specimen 3 enters the monitor 7 through a collimator 4 . The monitor is in the form of a geiger counter. The additional specimen is made from a pure element (or its oxide). The stabilization is ensured by using the output signal of the monitor to stabilize the cathode supply of the X-ray tube. The system is completely automatic, the control circuit being shown in Fig 2. It is shown that the use of a monitor in conjunction with good collimation of the direct fluorescence radiation from the additional specimen enables one to carry out accurate measurements of X-ray intensities without any stabilization of the supplies. Fig 4 shows a typical spectrum obtained with this instrument. Fig 3 shows the root mean square error in the intensity of the K line as a function of the atomic num-

ber Z of the specimen under investigation, the additional specimen being Ni . It follows from this figure that if a Card 2/3

SOV/120-59-1-26/50

The Use of a Monitor in X-Ray Spectral Analysis

relative error of 3% is sufficient (the number of counts taken being sufficiently high, i.e. the statistical error being low) then the atomic number of the specimen under investigation may differ from the corresponding number of the additional specimen by 4 . Hence altogether nine neighbouring elements may be investigated whose atomic numbers are symmetrically placed on either side of the atomic number of the additional specimen. If the relative statistical counting error does not exceed 4%, then for the above 3% the final relative error would be 4%, then for the above 2% the final relative error would be less than 5%. Thus almost the entire spectral region normally used in analysis by long wave spectrometers may be covered, using a single additional specimen, for example, a chromium specimen. Typical results are shown in Fig 4. There are 4 figures, 2 tables and 12 references, of which 8 are English, 1 is Japanese in English and the rest are Soviet.

ASSOCIATION: Rostovskiy-na-Donu gosudarstvennyy universitet (Rostovna-Domu State University)

SUBMITTED: January 18, 1958.

Card 3/3

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CIA-RDP86-00513R000205530008-8

80896

S/048/60/024/04/06/009 B006/B017

246300 AUTHORS:

Sec. The

On the Problem of the Shape of the Energy Bands of a Solid

Blokhin, M. A.,

TITLE:

PERIODICAL: Izvestiya Akademii nauk SSSR. Seriya fizicheskaya, 1960, Vol. 24, No. 4, pp. 397-406

Sachanko, V. P.

TEXT: The present article is a reproduction of a lecture delivered at the <u>Ath All-Union Conference on X-Ray Spectroscopy</u> (Rostov-na-Donu, June 29 - July 6, 1959). Since the experimentally observed shape of X-ray emission bands is only in approximate agreement with the theoretically computed one, the authors endeavor to improve the theory by taking into account the distortions due to the apparatus and the line broadening. In the introduction, the results of some investigations undertaken by other authors are discussed. Among others, Landsberg (Ref. 4) computed the form of the L_{III} sodium emission band by taking into account

the broadening of the conduction band levels, i.e., in free-electron approximation by means of a perturbation of the form $\exp(-ar_{12})/r_{12}$.

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On the Problem of the Shape of the Energy Bands of a Solid

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These computations show, however, some deficiencies which are discussed here. Above all, the level broadening in the conduction band was investigated only in connection with the emission of an X-ray photon; however, it occurs in all experiments on energy bands of solids. Hence, it must be taken into account not only in X-ray spectroscopy but also in investigating optical spectra, the photoeffect, and electrical conductivity. The authors of the present paper give an exact computation of the problem investigated by Landsberg (in free-electron approximation). They study again the influence exercised by level broadening on the shape of X-ray emission bands and the energy distribution of electrons in the bands of the solid. Two conduction electrons with the wave vectors k_2 and k_3 are

examined. One is to fill a vacancy (with the wave vector \vec{k}_1) after a collision, and the other is to pass over into a state with \vec{k}_4 . For $\vec{k}_2 + \vec{k}_3 = \vec{k}_1 + \vec{k}_4$ an explicit expression is obtained for the probability $W_A(k_1)$ after some operations. $2W_A(k_1)$ is the total transition probability to the \vec{k}_1 level when the exchange effects are

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On the Problem of the Shape of the Energy Bands of a Solid S/048/60/024/04/06/009 B006/B017

neglected. The dependence of the amount of level broadening on the position of the level in the band is illustrated in Fig. 1. The two curves $\Delta E(E)$ are given for $E_{\mathbf{F}} = 7$ ev (position of the Fermi surface in

copper) and $E_{\rm p} = 3.2$ ev (Fermi surface in sodium). As may be seen from the diagram, the level broadening in the conduction band on the band edge is very small and increases rapidly for lower levels. The following considerations deal with the energy distribution of the conduction electrons. By taking into account the transition probability one obtains

 $N_{1}(E) = \frac{1}{\pi} \int_{-\frac{1}{\pi}}^{E_{F}} \frac{N(t)P(t)}{a(t)\left[1 + \left(\frac{t-E}{a(t)}\right)^{2}\right]} dt$ In Fig. 2 the following transition

curves are plotteds $N_1(E)$, $N(E) = \sqrt{E}$, N(E) = E/E, and $N_1(E)$. Fig. 3 shows that the theoretical curve $N_1(E)$ agrees very well with the experimental shape of the L_{III} band of sodium. The following considerations concern the influence exercised by level broadening on the shape of the 3d bands of Cu, Ni, and Fe (Fig. 4). A great number of details concerning this

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| | On the Problem of the Shape of the Energy Bands of a Solid | S/048/60/024/04/06/009 B006/B017 | |
| | subject are discussed. In conclusion, the aut broadening in the energy bands explains some X-ray spectra, and that it is of universal im physics. A. F. Ioffe and Samoylovich are ment and 11 references: 3 Soviet, 4 American, and | portance in solid-state tioned. There are 4 figures | |
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NIKIFOROV, I.Ya.; SACHENKO, V.P.; BLOKHIN, M.A. Comparison of different methods for improving the form of spectra. Izv. AN SSSR. Ser. fiz. 25 no.8:1054-1059 Ag '61. (MIRA 14:8) 1. Rostovskiy-na-Donu gosudarstvennyy universitet. (X-ray spectroscopy)

APPROVED FOR RELEASE: 08/22/2000



BLOKHIN, M.A.; LOSEV, N.F.

Modern methods of X-ray spectroscopic fluorescent analysis. (MIRA 14:9) Zav.lab. 27 ho.9:1091-1099 '61. (X-ray spectroscopy) . .

APPROVED FOR RELEASE: 08/22/2000

CIA-RDP86-00513R000205530008-8

BLOKHIN, M.A.; VOLKOV, V.F.

Determination of the thickness of a deposited layer by means of the KRFS-2 X-ray spectromater. Zav.lab. 27 no.9:1110-1111 '61. (MIRA 14:9) 1. Rostovskiy gosudarstvennyy universitet. (Thickness measurement) (Spectrometry)

APPROVED FOR RELEASE: 08/22/2000

S/048/62/026/003/010/015 B142/B104

AUTHORS:

Blokhin, M. A., Giltvarg, A. B., Nikiforov, I. Ya., and Sachenko, V. P.

TITLE:

tenii:

Two-crystal X-ray spectrometer

PERIODICAL:

Akademiya nauk SSSR. Izvestiya. Seriya fizicheskaya, v. 26, no. 3, 1962, 397 - 404

TEXT: The adjustment of the new spectrometer is comparatively simple and takes only a few hours. The crystals can be taken out of the apparatus without disturbing the adjustment. The distance between the rotating axes of the crystals is 100 mm. The focus of the X-ray tube is 300 mm distant from the rotating axis of the first crystal. The distance of the rotating axis of the second crystal from the window of the Geiger counter is 100 mm. The second crystal can be rotated by $\pm 1.5^{\circ}$ from the middle position reading accuracy 0.01°). The spectrometer is not adjusted by means of the crystals but by glass plates. After adjustment, the crystals are inserted to determine the CuKo₁ - line and the angle between crystal surface and lattice

planes. Eight horizontal plates were built into the collimator to reduce Card 1/3

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Two-crystal X-ray spectrometer

s/048/62/026/003/010/015 B142/B104

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the vertical scattering of the beam to a minimum and yet to obtain high radiation intensities. A beryllium plate inserted between the collimator and the first crystal is to eliminate the focus drift and the effect of feeding-voltage fluctuations. It was difficult to choose the suitable crystals since extreme optical uniformity is required, and the angle between crystal surface and lattice planes shall be as small as possible. Its maximum was 105". Plates parallel to (1010) and (1120) were cut from various quartz crystals and investigated after etching. The purity of the two crystals is determined by the width of the reflection curves. The quality of the plates is estimated from the shadows produced by deviations of the refractive indices. A final examination carried out by means of a polarization system indicates optical inequality of the plates by bright a c spots. There are 6 figures and 6 references: 1 Soviet and 5 non-Soviet. Thee two English-language references are: L. G. Parrat, Rev. Scient. Instrum. 5, no. 11, 113 (1934); Rev. Scient. Instrum., 6, no. 5, 113 (1935).

Card 2/3

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CIA-RDP86-00513R000205530008-8

e.e.... s/048/62/026/003/014/015 B102/B104 Blokhin, M. A., Demekhin, V. F., and Shveytser, I. G. AUTHORS: Correction of the X-ray emission spectrum for self- absorption TITLE: Akademiya nauk SSSR. Izvestiya. Seriya fizicheskaya, v. 26, PERIODICAL: no. 3, 1962, 419 - 422 TEXT: Corrections for self-absorption are considered for the continuous and the characteristic spectrum separately. In the first case, it is not necessary to know the absolute values of the absorption coefficients. In the second case, the intensity of the characteristic spectrum can be given by I = $Ae^{-C_1 c} \left[c_2 + c_3 c \right]$ with $C_{1} = \frac{x}{\sin\psi} \frac{V^{3}}{V^{2} - V_{1}^{2}}, \quad C_{2} = \frac{V - V_{1}}{V_{1}} - \lg \frac{V}{V_{1}}, \quad C_{3} = \frac{x}{\sin\psi} \frac{2V^{3} - 3V^{3}V_{1} + V_{1}^{3}}{6V_{1}(V^{2} - V_{1}^{2})}$ (5), where x is the maximum penetration depth of electrons, and τ is the absorption coefficient. A practical correction for self-absorption is demonstrated for the L_{β} band and the L_{III} spectrum of metallic Mo. The intensities Card 1/

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Correction of the ... S/048/62/026/003/014/015 B102/B104 with and without absorber (I and I) were measured. The scattered background itself was also measured with and without absorber $(I_b and I_{bo})$, allowing for the cosmic background intensity I_c . Then $\tau = 2.3\frac{e}{m} \log \left[(I_o - I_c) / (I - I_b - I_c) \right]$. For a particular case x = 0.04 was obtained; I = $I \cdot 0.53/e^{-2.6 \cdot 10^{-5\tau}}$ (1+7.8.10⁻⁶ τ). The effect of voltage unself-absorption was studied at 3.5, 5.5, and 12 kv. The self-absorption of the characteristic spectrum increases with increasing voltage while that of the continuous spectrum decreases. Since, however, the intensity of the former spectrum rises with increasing voltage more rapidly than that of the latter, it depends on geometry if the self-absorption of the continuous spectrum increases or decreases. There are 2 figures and 6 references: 4 Soviet and ASSOCIATION: Rostovskiy gos. universitet (Rostov State University) Card 2/2

APPROVED FOR RELEASE: 08/22/2000

ale leste BLOKHIN, M.A.; SHUVAYEV, A.T. Effect of chemical bonds on the X-ray emission spectrum of titanium. Izv. AN SSSR. Ser. fiz. 26 no.3:429-432 Mr 162. (MIRA 15:2) 1. Rostovskiy gosudarstvennyy universitet. (Chemical bonds) (X-ray spectroscopy)) (Titanium)

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CIA-RDP86-00513R000205530008-8

S/048/63/027/003/002/025 B108/B114 Nikiforov, I. Ya., and Blokhin, M. A. AUTHORS: About the form of the K_{β_5} emission band of iron. II. TITLE: The transition probability as a function of energy Akademiya nauk SSSR. Izvestiya. Seriya fizicheskaya, v. 27, no. 3, 1963, 314-318 PERIODICAL: TEXT: This paper was presented at the 6th Conference on X-ray Spectroscopy, Odessa, July 2 - 10, 1962. The transition probability, was calculated in the single-electron theory. The method of cells (M. F. Manning, Phys. Rev., 63, 190, 1943) was used to calculate the phase of the K emission band N(F). For this purpose, the dispersion shape of the K_{β_5} emission band, N(E). For this purpose, the dispersion curves were determined from the coefficients a₁(E) accounting for the spherical harmonics of 1-symmetry in the total wave function of the valency electrons. The calculated form of the band agrees well with the Card 1/2

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APPROVED FOR RELEASE: 08/22/2000

BLOKHIN, M.A.; DUYMAKAYEV, Sh.I.

Optimal X-ray spectral analysis of solutions. Zav. lab. 29 no.9:1061-1064 '63. (MIRA 17:1)

1. Rostovskiy gosudarstvennyy universitet.

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BLOKHIN, M.A.; DRUZ', V.V.
X-ray spectral analysis of multicomponent mixtures. Zav. lab. 29 no.9:1070-1074 '63. (MIRA 17:1)
1. Rostovskiy gosudarstvennyy universitet.

APPROVED FOR RELEASE: 08/22/2000

BLOKHIN, M.A.; DUYMAKAYEV, Sh.I.

X-ray spectral analysis of multicomponent mixtures. Zav.lab. 30 no.4:425-426 '64. (MIRA 17:4)

1. Rostovskiy-na-Donu gosudarstvennyy universitet.

APPROVED FOR RELEASE: 08/22/2000

S/0048/64/028/005/0780/0785 ACCESSION NR: AP4038761 AUTHOR: Blokhin, M.A.; Nikiforov, I.Ya. TITLE: Shape of the KC1,2 lines of the iron group elements /Report, Seventh Conference on X-Ray Spectroscopy held in Yerevan 23 58p-1 Oct 19637 SOURCE: AN SSSR. Izvestiya. Seriya fizicheskaya, v.28, no.5, 1964, 780-785 TOPIC TAGS: x-ray spectrum, chromium, mangangee, iron, cobalt, copper, nickel, zinc ABSTRACT: The KX doublets of the elements of atomic number from 24 (Cr) through 30 (Zn) were recorded with the high resolution two-crystal spectrometer of the Rostov State University. The instrument and the experimental procedure are discussed elsewhere (M.A.Blokhin, A.B.Gil'varg, I.Ya.Nikiforov, V.P.Sachenko, Izv.AN SSSR, Ser.fiz 26,397,1962). The resolving power was approximately 38 000, the dispersion was 0.01 X per second of arc, and the angle could be measured to ±0.5". The Cr and Ni spectra were obtained with Cu anodes on which Cr or Ni had been electroplated. The other spectra were obtained with Cu anodes into which powders of the corresponding metals had been pressed. The double reflection curves with parallel crystals were quite narrow (0.15 to 0.21 eV). The widths of the XX1 lines were corrected for in-Card 1/3 الجاني يعكمه الاستان فيعمد الفاقة متصافية المتربح متركبة المحتوجين والمراجع الحاري

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strumental broadening by simply subtracting the width of the corresponding parallel crystal double reflection curve. One spectrum (Cu) was corrected by a more rigorous method (V.P.Sachenko and I.Ya.Nikiforov, Optika i spektro. 13,447, 1962) with practically identical results; the corrections are therefore believed to be adequate. Of the published measurements of the $K \alpha_1$ widths of the iron group metals, the present measurements, those of G.Brogren (Arkiv.fyz.23,219,1963), and those of A.Meisel and W. Nefedow (Z.phys.Chem.(DDR),219,397,1962) were obtained under the most advantageous conditions with regard to instrumental broadening. Although there is considerable agreement among the three groups of data, there is also some disagreement among them. The KO1 asymmetry indices were calculated, and they are compared with data of other workers and with the magnetic moments. Although the correlation between KX1 asymmetry and magnetic moment is strong, it is not perfect. Notable deviants are Cr and Cu, both of which are much too asymmetric for their small (or vanishing) magnetic moments. The copper spectrum was corrected for the width of the K level, and it is concluded from the shape of the corrected curve that the asymmetry of the Cu spectrum is due to complex structure of the LIII level. The KX1 line of metallic chromium was found to be complex. This fine structure is much more prominent in the spectrum of the oxide, where it appears also in the KC2 line and has previously been reported by others. In order to make the MX doublet shapes conveniently available for

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| | cally, but als | o in tabular placements is | ucture analysis, the data are prese form. The accuracy claimed for the s 1% of the peak value, and 0.05 eV ables. | a tabulated intensities | |
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ACCESSION NR: AP4038762

5/0048/64/028/005/0786/0789

AUTHOR: Nikiforov, I.Ya.; Blokhin, M.A.

TITLE: Concerning the shape of the x-ray emission bands of transition metals of the iron group /Report, Seventh Conference on X-Ray Spectroscopy held in Yerevan 23 Sep-1 Oct 19637

SCURCE: AN SSSR. Izvestiya. Seriya fizicheskaya, v.28, no.5, 1964, 786-789

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TOPIC TAGS: x-ray spectrum, band spectrum, line spectrum, iron, nickel, copper, x-ray structure analysis

ABSTRACT: The K\$5 bands of Fe, Ni and Cu, and the K\$1 lines of Fe and Cu were recorded with the two-crystal spectrometer (resolution 38 000) of Rostov State University (M.A.Blokhin, A.B.Gil'varg, I.Ya.Nikiforov and V.P.Sachenko, Izv.AN SSSR, Ser. fiz.26,397,1962). Quartz crystals cut parallel to the (1120) planes were employed. The Fe and Cu spectra were obtained with anodes of the respective metals. The Ni spectrum was obtained with a Cu anode on which Ni had been electroplated. The x-rzy tube wes operated at 35 kV and 20 mA. An accuracy of 2% is claimed for the ordinates of the published spectral intensity curves. The shapes of the Cu and Fe K\$1 lines

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were recorded because of their technical importance for precision x-ray structure analysis. They are presented graphically but are not discussed. It is not possible to obtain reliable widths of the $K\beta_5$ bands without correcting for the width of the K level and removing the K3"' satellite. This analysis was not performed. The widths at half maximum increased with increasing atomic number, corresponding to the increasing density of free electrons between the 1s and 3p shells. The principal peak of the Cu band was found to be double. The two peaks are ascribed to the $k\beta_5$ and $K\beta_2$ lines, although the assignment is regarded as arbitrary because of the strong hybridization of the conduction band. The Cu $K\beta^{\prime\prime\prime}$ satellite was clearly re-solved into two satellites, which are designated by $K\beta^{\prime\prime\prime}$ and $K\beta^{\prime\prime\prime}$. Weak structure was found on the long wavelength side of the Ni and Ro (2). Weak structure was found on the long wavelength side of the Ni and Fe KB5 bands. This could be due to a long wavelength satellite, or to zonal structure of the electron states in the lattice. The fe KB5 peak was broad and nearly flat. Previous calculations of the Fe K\$5 band shape (I.Ya.Nikiforov, and M.A.Blokhin, Izv.AN SSSR, Ser.fiz.27, 314, 1963) are compared with the present measurements, and considerably better agreement is found than was previously obtained with the measurements of J.A.Bearden and C.H. Shaw (Phys.Rev. 48, 18, 1935). The calculations do not reproduce the structure on the long wavelength sids. The paper closes with a short essay on the role of X-ray spectroscopy in the development of solid state physics. Although it is not possible

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to obtain electron densities by simply dividing x-ray band intensities by transition probabilities as envisaged 30 years ago by H.Jones, N.F.Mott and H.W.B.Skinner (Phys. Rev.45,379,1934), one can nevertheless employ different approximate methods for dealing with the many body problem to calculate x-ray band shapes, and by comparing the calculated shapes with experimental data one can select the most promising mathematical methods for further development in connection with solids of particular types. Orig.art.has: 3 formulas and 3 figures.

ASSOCIATION: Rostovskiy-na-Donu gosudarstvenny*y universitet (Rostov-on-the-Don-State University)

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| • | ACCESSION NR: AP4038765 | S/0048/64/028/005/0801/0804 |
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| | AUTHOR: Blokhin, M.A.; Shuvaye | ev, A.T.; Gorskiy, V.V. |
| · · · · | | nvestigations of chemical bonds in sulfur compounds n. X-Ray Spectroscopy held in Yerevan 23 Sep-1 Oct 1963 |
| | SOURCE: AN SSSR. Izvestiya. S | eriya fizicheskaya, v.28, no.5, 1964, 801-804 |
| | TOPIC TAGS: x-ray spectrum, 1 | ine shift, sulfur, sulfur compound, chemical bond |
| | AP40387587) the shift of the l chemical compounds is due to the neighboring atoms. This pl sulfur, and a short table is the shifts produced by various cal bond shifts of the S XX 1 KX fluorescence spectra of S excited by 20 kV Cu bremsstral | huvayev (Izv.AN SSSR,Ser.fiz.28,758,1964 /see Abstract KX lines of sulfur (and other Period 2 elements) in the charge on the atom arising from the influence of henomenon is discussed at some length for the case of presented, based on data in the literature, showing is chemical bonds and bond configurations. These chemi- tines are believed to be approximately additive. The in several compounds were recorded. The spectra were hlung and formed by reflection from the (1010) planes systal. The temperature of the samples did not exceed |
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50°C. The shifts of the KX lines were measured with respect to their position in rhombic sulfur. The shifts obtained for the compounds H_7C_7 -S-S-C7H7 and S = = $C(N(CH_3)_2)_2$ were 0 and -0.13 eV, respectively, and are in accord with the structures as written. A shift of -0.02 eV was observed for methylene blue; of the two structures proposed for this compound, the x-ray data favor that in wich the chlorine is attached to one of the nitrogen atoms and not to the sulfur. The polymer $(CuS_2N_4C_24H_{18})_n$ showed a shift of -0.06 eV; from this it is concluded that the S-C bond is single. Three compounds containing the SCN group were investigated: KSCN, CuSCN and NH2C6H4SCN, for which the KX aline shifts were -0.10, 0 and 0.07 eV, respectively. These data favor the structure $-S-C \equiv N$ for the SCN group in the aniline derivative and the copper salt, and a structure between this and -S=C=N in KSCN. It is concluded that x-ray spectroscopy can be a useful tool for investigating che-"The authors are grateful to Z.V. Zvonkova (Physical-chemical Institute mical bonds. im.L.YA.Karpov) and I.G.Mochalina (Special Organic Synthesis Laboratory of the Moscow State University) for preparing the samples." Orig.art.has: 3 tables.

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CIA-RDP86-00513R000205530008-8

ACCESSION NR: AP4038771

\$/0048/64/028/005/0825/0829

AUTHOR: Demokhin, V.F.; Blokhin, M.A.

TITLE: Fluorescence spectra of silivon in some compounds /Report, Seventh Conference on X-Ray Spectroscopy held in Yerevan 23 Sep to 1 Oct 19637

SOURCE: AN SSSR. Izvestiya. Seriya fizicheskaya, v.28, no.5, 1964, 825-829

TOPIC TAGS: x-ray spectrum, siligon, silicon dioxide, silicon carbide, carbon chemical bond

ABSTRACT: The $K\alpha_1,\alpha_2$, $K\alpha',\alpha_3,\alpha'_3,\alpha'_4$ and $k\beta_{1,x}$ lines of silicon were observed in SiO₂, SiC and crystalline silicon, both pure (resistivity 0.5 ohm cm). The investigation was undertaken to determine the true shape of the $k\beta_x$ band, for which widely differing shapes have been reported (J.Farineau,Ann.Phys.10,20,1938; Ya.M.Fogel', Zhur.eksp.i teor.fiz.9,1217,1935), and to determine the effect of chemical bonding on the lines. The spectra were obtained by reflection from the (1010) planes of a bent (50 cm radius) quartz crystal. The resolving power was 16 000. The x-ray tube was operated at 25 kV and 30 mA and the samples were at temperatures between 50 and 70°C. The continuous spectrum was observed with 3.8 kV on the x-ray tube in order to

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determine whether the K absorption of the silicon in the diffracting crystal would distort the spectrum in the $K\beta_X$ region; no such distortion was observed. A dry run was made with an aluminum sample held with the same cement as was used in the investigation proper; no lines were found in the region of interest. The presence of phosphorus in the silicon crystal produced no observable change in the spectrum. The experimental error in these measurements was ±0.07 eV for position, ±0.03 eV for line widths, and 2% for relative intensities. The width of both the XX1 and the KG_2 line, corrected for instrumental broadening, was 0.45 ± 0.05 eV, and their separation was 0.56 eV. The shifts due to chemical bonding of the KO1 line, as well as those of the α_4 , α_3 , α'_3 and α' satellites, agreed with the values obtained by N.G. Johnson (Diss.Lund, 1939) and H.Karlsson-Flemberg (Z.Phys. 96, 167, 1935). When the charge on the silicon ion increased, the satellites shifted approximately twice as far as the KC1 line, and the intensities relative to KC1 of the C4 and C' satellites increased and those of α_3 and α'_3 decreased. This behavior is discussed in terms of the charge between the K and L shells due to the valence electrons. The shape of the 131,x band agreed well with that reported by Fogel' (loc.cit.). Fine structure was perceptibel in the $K\beta_X$ line. The $K\beta_1$ line was observed in pure silicon under conditions that are said to proclude its being due to the presence of SiO2. The KB Eands of Si and C in SiC and of Si in the pure crystal were found to be very si

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| bands of these subst mond structure in bo | is concluded that the distribution of el ances is determined mainly by the lattic th cases. Orig. art. has: 4 figures and | l 1 table. |
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CIA-RDP86-00513R000205530008-8

S/0048/64/028/005/0830/0831 ACCESSION NR: AP4038772 AUTHOR: Blokhin, M.A.; Demekhin, V.F. TITLE: Emission spectra of scandium in Sc203 /Report, Seventh Conference on X-Ray Spectroscopy held in Yerevan 23 Sep to 1 Oct 19637 SOURCE: AN SSSR. Izvestiya. Seriya fizicheskaya, v.28, no.5, 1964, 830-831 TOPIC TAGS: x-ray spectrum, x-ray emission, scandium, scandium compound ABSTRACT: The Sc K fluorescence spectrum was excited by 30 kV bremsstrahlung and -by-Ti KX radiation, and the spectra were recorded photographically with a quartz crystal spectrometer (resolution 16 000). The Ti KX1,2 lines were chosen to excite the "tertiary" spectrum because they lie in the long wavelength region of the Sc K -absorption band and therefore, according to the exciton model (L.G. Parratt, Revs. Mod. Phys.31,616,1959), they cannot ionize the Sc atom in the K shell, but only excite it. The tertiary spectrum was excited in a special apparatus, the design of which is shown schematically in Fig.1 of the Enclosure. Exposures up to 80 hours were required to record some portions of this spectrum. Intensity measurements in the KX3,4 region were performed by simultaneously recording the spectrum on two films and com-Card 1/3 -

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paring the image of the KX₃,4 region on the foremost film with the weakened image of the KX₂ peak on the other. The secondary and tertiary spectra differed in only one respect: the KX₃ and KX₄ lines were entirely absent from the tertiary spectrum. All the other lines observed (α_1 , α_2 , β_1 , β_5 , β^* and $\beta^{"}$) were identical in shape and relative intensity in the two spectra. The absence of KX₃,4 from the tertiary spectrum suggests that the double ionization theory of the origin of these lines is correct; it is in sharp contradiction with the interpretation of the KX satellites given by T.I.Kakushadze (Izv.vy*sshikh uchebn.zaved.Fizika, No.3,142,1963). The identity of the K spectra from ionized and from merely excited atoms is regarded as confirmation of the suggestion of J.Friedel (Philos.Mag.(7),43,153,1952) that after K ionization the valence band releases an electron which shields the hole in the K shell. Orig.art.has: 3 figures.

ASSOCIATION: Rostovskiy-ma-donu gosudarstvenny*y universitet (Rostov-on-the-Don State University)

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ACCESSION NR: AP4038774

AUTHOR: Blokhin, M.A.; Demekhin, V.F.; Shveytser, I.G.

TITLE: L Spectra of some molybdenum compounds /Report, Seventh Conference on X-Ray Spectroscopy held in Yerevan 23 Sep to 1 Oct 19637

SOURCE: AN SSSR. Izvestiya. Seriya fizicheskaya, v.28, no.5, 1964, 834-835

TOPIC TAGS: x-ray spectrum, x-ray absorption, molybdenum, molybdenum compound, chemical bond

ABSTRACT: In order to obtain information concerning the extent to which electrons in the incomplete 4d shell of transition metals participate in chemical bonding, the LII and LIII absorption spectra and the I β_2 emission bands of metallic molybdenum, and Mo in MoO3, CaNoO4 and MoS2 were recorded. Although some of these spectra have been previously reported, the results of different workers are not all in agner level. The spectra were recorded photographically with a spectrometer having a resolution of 12 000, and the LIII edge was observed with a second instrument having half this resolving power and employing an ionization chamber for redording. The ob-

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served spectra were corrected for the width of the inner level, the width of the No LIII level being assumed to be 1.76 eV. The corrected absorption curve for metallic Mo did not break sharply; this indicates that the LIII level is in fact somewhat wider than assumed. A gap between the emission and absorption was perceptible in the insulators MoO3 and CaMoO4. The LIII spectra were in good agreement with those obtained by I.V.Borovskiy, K.P.Gurov, et al (Izv.AN SSSR, Ser.fiz.21,1401,1957). As the valence increased, the absorption edge shifted toward shorter wavelengths. This shift, which attained 3.4 eV for the LIII edge of CaMoO4, is ascribed to decreased shielding of the inner portion of the atom by the valence electrons that become inshielding of the inner portion after a basorption line was observed in the LIII spectrum of CaMoO4 at 13 eV from the primary line. Such lines have been previously observed in molybdenum compounds and are ascribed to transitions of 2p electrons to the incomplete 4d shell. Orig.art.has: 3 figures and 1 table.

ASSOCIATION: Rostovskiy-na-Donu gosudarstvenny*y universitet (Rostov-on-the-Don State University)

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| be as follows: La - 0. | asitivity of this method for all the ratio gm/ltr , Ge - 0.1, Pr - 0.15, Nd - 0. | 15. Sm - 0.2. Eu - 0.3. |
| Gd - 0.3, Tb - 0.35, D Orig. art. has: 1 table | y - 0.45, Ho - 0.5, Er - 0.6, Tu - 0.9, | Yb - 0.9, Lu - 0.95. |
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DUYMAKAYEV, Sh.I.; BLOKHIN, M.A.

Degeneracy reading in X-ray spectral analysis by means of the addition of diluents or the element being analyzed. Zav. lab. 31 no.9:1072-1076 (MIRA 18:10) *65.

1. Rostovskiy-na-Donu gosudarstvennyy universitet.

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BLOKHIN, M.A.; OVCHABENKO, Ye.Ya.; MYAGKOV, P.I.; SOTNIKOV, V.A.; MAMONOV, Yu.M.; BELKINA, G.L.

> Improving the accuracy of X-ray spectral analysis by a dual channel method. Zav.lab. 31 no.4:423-426 ^{165.} (MIRA 18:12)

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PA 4/47103 BLOKHIN, N. N. PROF Jan 48 UBSR/Medicine - Metabolism Medicine - Carbohydrates, Metabolism "The School of Professor Yefim Semenovich London in Teningrad University," Prof N. N. Blokhin, 8 pp West Leningrad U" No 1 Sol Res Inst of Physiol of Leningrad U was founded 18 1932. From then on Prof London has been in obarge of the Metabologic Lab. Describes his achievements. He was first to observe changes in block flow through organs in vivo. Reviews laboratory work on carbohydrates, albumin, fats, gases, and ter metabolism. 1/49163

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