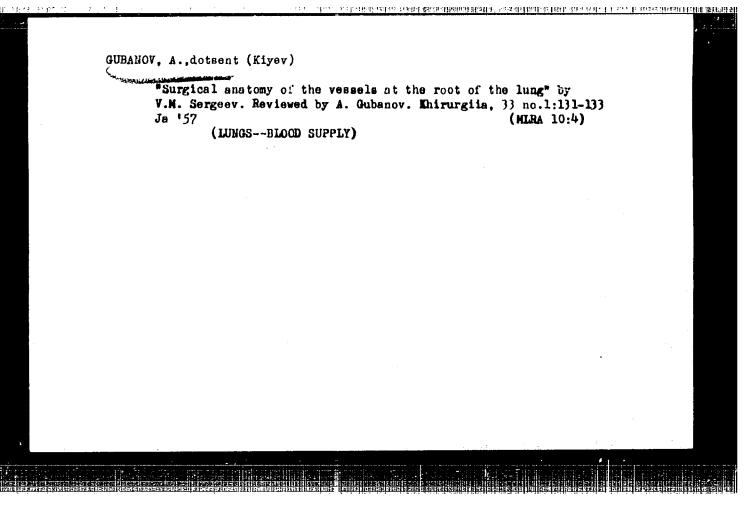
GUBANOV, A.C., dotsent

Development of segmental pulmonary resection in tuberculosis [with summary in French] Probletub. 34 no.6:31-37 N-D '56. (MIRA 10:2)

1. Iz Ukrainskogo nauchno-isaledovatel'skogo instituta tuberkuleza imeni F.G.Yanovskogo (dir. A.S.Mamolat)
(PNEUMONECTOMY, in various diseases, tuberc., segmental technic (Rus))



Intersegmental connective tissue borders of the lungs. Khirurgiia
33 no.8:48-53 Ag '57. (MIRA 11:4)

1. Is Urrainskogo nauchno-issledovatel skogo instituta tuberkulesa
in. akad. F.G. Ianovekogo (dir. A.S. Mamolat).

(LUNGS, anat. and histol.
intersegmental connective tissue borders, surg. anat.
aspects)

(COMBECTIVE TISSUE, anat. & hictol.
intersegmental of lungs, surg. aspects)

Characteristics of bethological changes of pulmonary braches of the vagus nerve in tuberculesis [with summary in French]. Probl.tub. 35 no.3:82-86 '57. (MIZA 10:10)

1. Iz Ukrainskogo nauchno-issledovatel'skogo instituts tuberkuleza imeni F.G.Yanovskogo (dir. A.S.Mamolat) (TUBERCULOSIS, PULMOMARY, pathology, vagus nerve (Rus)) (NERVSS, VAGUS, pathology, in tuberc., pulm. (Rus))

GUBANOV, A.G., dotsent (Kiyev, 54, ul.Chkalova, d.74, kv.7)

So-called new method of bronchial closure in pulmonary re-

So-called new method of bronchial closure in pulmonary resection for tuberculosis. Nov.khir.arkh. no.1:41-45 Ja-F 159. (MIRA 12:6)

1. Ukrainskiy nauchno-issledovatel skiy institut tuberkuleza.
(BRONCHI--SURGERY) (SUTURES)

GUBANOV, A.G., dotsent

Simplified methods for resection of lung tiesue. Pat., klin.i
terap.tub. no.8:333-338 '58. (MIRA 13:7)

1. Ix Ukrainskogo mauchno-issledovatel'skogo instituta tuberkuleza im. akad. F.G. Tanovskogo.
(LUNGS--SURGERY)

GUBAROV, A.G., dotsent (Kiyov, ul. Chkalova, d.74/7); GOMOVERKO, G.G.;

BEREZOVSKIY, K.K., starshiy nauchnyy sotrudnik

First experience in using porolon for plombage of the chest cavity in an experiment and in the clinic. Nov. khir. arkh. no.3:65-72 My. Je '60. (MIKA 15:2)

1. Pervoye khirurgicheskoye otdeleniye (zav. - dotsent G.G.Gorovenko) i 2-ye khirurgicheskoyo otdeleniye (zav. - prof. N.K.Amosov) Ukrainskogo nauchno-issledo vatel'skogo instituta tuberkuleza imeni akademika F.G.Yanovskogo. (PLASTICS IN NEDICIHE) (PLOMBAGE (SURGERY)) (CHEST._SURGERY)

sensia in nomen yan ngamin is ni hinin ngamig ina ing a tini ing pilang ang mang mahababababan.

GUBANOV, A. G.; LITVINOV, V. V.; SMIRNOV, A. A.; KHMELEVSKAYA, G. A.

Experimental data on the use of porolon for alloplasty. Grud. khir. no.4:66-71 '61. (MIRA 14:12)

1. Iz Kiyevskogo nauchno-issledovatel'skogo instituta tuberkuleza imeni akademika F. G. Yanovskogo i Nauchno-issledovatel'skogo instituta meditsinskoy klimatologii i klimatoterapii imeni I. M. Sechenova (Yalta). Adres avtorov: Krym, Yalta, ul. Dzerzhinskogo, d. 48. Institut imeni I. M. Sechnova, korp. 12

(PLASTICS—THERAPEUTIC USE) (LUNGS—SURGERY)

GUBANOV, A.G., dotsent; FEDOTOV, A.F.

Intrapleural plombage with porolon in accociation with pulmonary resection. Probl. tub. 39 no.3:44-49 '61. (MIRA 14:5)

1. Iz Ukrainskogo nauchno-issledovatel'skogo instituta tuberkulesa imeni F.G. Ianovskogo (dir. - kandamed.nauk A.S. Mamolat).

(LUNGS—SUNDERY) (LUNGS—COLLAPSE)

to an actual title of the contraction of the city of

GUBANOV, Aleksey Gavrilovich; UMOVIST, M.N., red.; FOTOTSKAYA, L.A., tekhn. red.

[Partial pulmonary resections in tuberculosis; anatomical and experimental materials] Chastichnye rezektsii legkikh pri tuberkuleze; anatomicheskie i eksperimental nye materialy. Kiev, Gosmedizdat USSR, 1961. 303 p. (MIRA 15:7) (TUBERCULOSIS) (LUNGS-SURGERY)

ा सम्भावता संभवता स्थान स्थान स्थान स्थान । स

GUBANOV, A.G.; SEVEROV, V.S.; OSINTSEVA, V.P.; FEDOTOV, A.F.

Use of porolon plombage in partial resections of the lungs in tuberculosis. Vest.khir. no.5:46-51 '61. (MIRA 15:1)

1. Iz Instituta tuberkuleza (dir. - prof. N.A. Shmelev) AMN SSSR
i Ukrainskogo nauchno-issledovatel'skogo instituta tuberkuleza
(dir. - kand.med.nauk A.S. Mamolat).
(LUNCS-SURGERY) (TUBERCULOSIS) (PLASTICS IN MEDICINE)

GOROVENKO, Grigoriy Gavrilovich, doktor med. nauk; GUBANOV, A.G., red.;
CHUCHUPAK, V.D., tekhn. red.

[Pulmonary resections following ineffective collapse therapy]Rszektsii legkikh posle neeffektivnoi kollapsoterapii. Kiev, Gosmediadat USSR, 1962. 277 p. (MIRA 16:1)

(TUBERCULOSIS) (PNEUMOTHORAX) (LUNGS—SURGERY)

GUBAROV, A.G., dotsent; NOVITSKIY, A.B.

Technic of extra-musculo-periosteal plombage of the thoracic cavity. Khirurgiia no.1:83-88 '62. (MIRA 15:11)

1. Iz Ukrainskogo nauchmo-issledovatel'skogo instituta tuberku-leza imeni akad. F.G. Yanovskogo (dir. - dotsent A.S. Mamolat)
i Simeizskogo khirurgicheskogo sanatoriya "Primor'ye" (glavnyy vrach I.T. Sokolova [decessed]).

(IUNGS—GOLLAPSE)

BAYANDIN, P.A. (Murmansk); SHVETSOV, I.M.; TIMOFEYEVA, M.V.; KOVAL', V.P.; KOZLOVA, E.Z.; TRET'YAKOV, N.I. (Kaliningrad); MAMEDOV, E.Sh. (Poselok Martuni, AzerSSR); BOROVYY, Ye.M.; DULAYEV, S.G. (Grodno); GERASIMOV, B.A. (Lugansk); MEL'NIK, L.A. (Chernovtsy); MIGAL', L.A.; GUBANOV, A.G.; GOROVENKO, G.G. (Kiyev); SHAROV, B.K. (Chelyabinsk); SHUVALOVA, Z.A. (Sverdlovsk) NEYMARK, I.I.; ARYAYEV, L.N. (Odessa); KABANOV, A.N.; KONOVALOV, Yu.S.; ZAK, V.I. (Orenburg); MIKHAYLOV, M.M.; SEZ'KO, A.D. (Voronezh); SHALAYEV, M.I.; DONIN, V.I. (Saratow).

Abstracts. Grudn. khir. 5 no.3:110-126 My-Je 63 (MIRA 17:1)

1. Iz kafedry normal'noy anatomii Ryazanskogo meditsinskogo instituta imeni akademika I.P.Pavlova (for Shevtsov). 2. Iz Sochinskogo nauchmo-issledovatel'skogo instituta kurortologii i fizioterapii Ministerstva zdravookhraneniya RSFSR (for Timofeyeva).
3. Iz khirurgicheskogo otdeleniya Ternopol'skoy klinicheskoy gorodskoy bol'nitsy (for Koval'). 4. Iz kafedry topograficheskoy anatomii i operativnoy khirurgii (zav. - prof. A.P. Sokolov). Permskogo meditsinskogo instituta (for Kozlova). 5. Iz khirurgicheskogo otdeleniya (zav. - Ye. M. Porovyy) Rovenskoy oblastnoy bol'nitsy (glavnyy vrach - UkrSSR V.ii. Vel'skiy) (for Borovyy).

(Continued on next card)

BAYANDIN, P.A. (continued) Card 2.

6.Iz fakul'tetskoy khirurgicheskoy kliniki (dir. - prof. I.M. Popov'yan) i gospital noy terapevticheskoy kliniki (dir. - prof. L.S.Shvarts) lechebnogo fakul'teta Saratovskogo meditsinskogo instituta (for Migal'). 7. Iz kafedry fakul'tetskoy khirurgii (zav. - prof. I.I.Neymark) Altayskogo meditsinskogo instituta (for Neymark). 8. Iz Novosibinskogo gorodskogo protivetuberkuleznogo dispansera (for Kabanov). 9. Iz kafedry fakul'tetskoy khirurgii (zav. - prof. I.A.Ivanov) Permskogo meditsinskogo instituta (for Shalayev).

GUBANOV, A.G., dotsent (Kiyev, ul. Chkalova, d.74, kv.7); FURMANOV, Yu.A.; MARULIN, B.A.

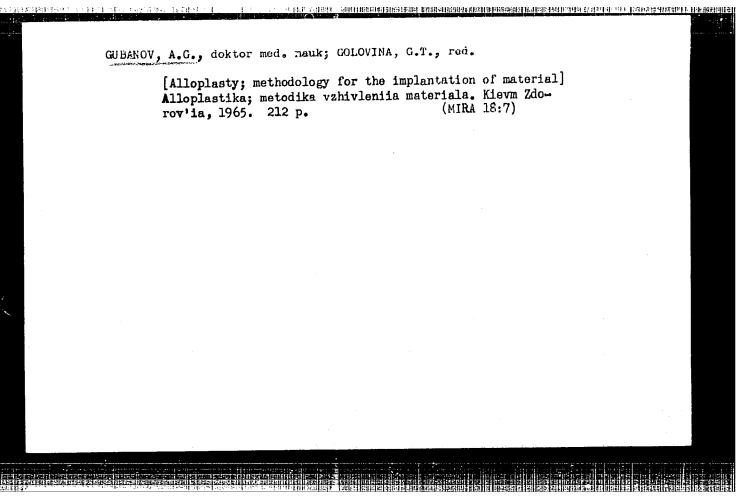
Soft elastic porous polymers as plastic material in surgery. Vest. khir. 89 no.10:65-72 0 '62. (MIRA 17:10)

1. Iz Ukrainskogo nauchno-issledovatel skogo instituta tuberkuleza i grudnoy khirurgii daeni akademika F.G. Yanovskogo (dir. - dotsent A.S. Mamolat).

GUBANOV, A.G., dotsent tKiyev, ul. Fuchtka, d..i, kv.n?;

Prevention of infection of the pleural cavity in alicelasty of the esophagus. Vest. khir. 91 no.11:21-25 N 163. (MIRA 17:12)

1. Lz otdela polimerov (rukcvoditel' - dotsent A.G.Gubanov) Ukrainskogo instituta tuberkuleza i grudnov khirurgii (direktor - dotsent A.S. Mamolat).

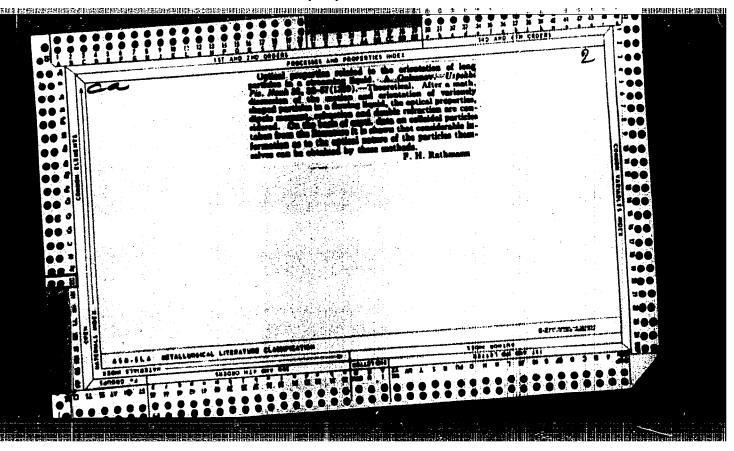


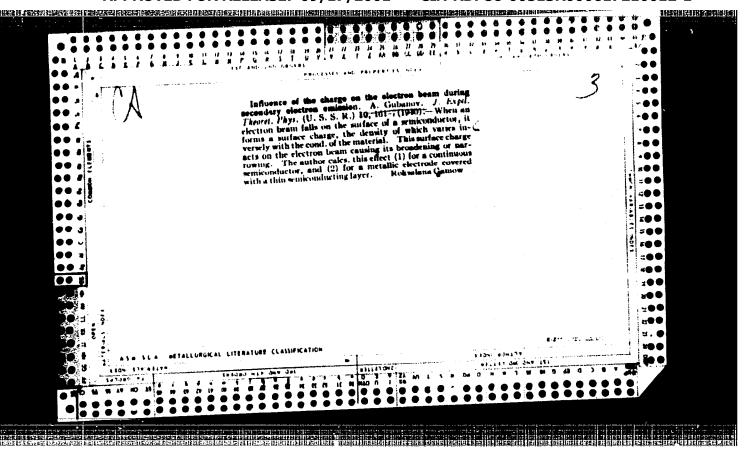
KUVAYEV, V.B.; VLASOV, M.I.; GUBANOV, I.A.

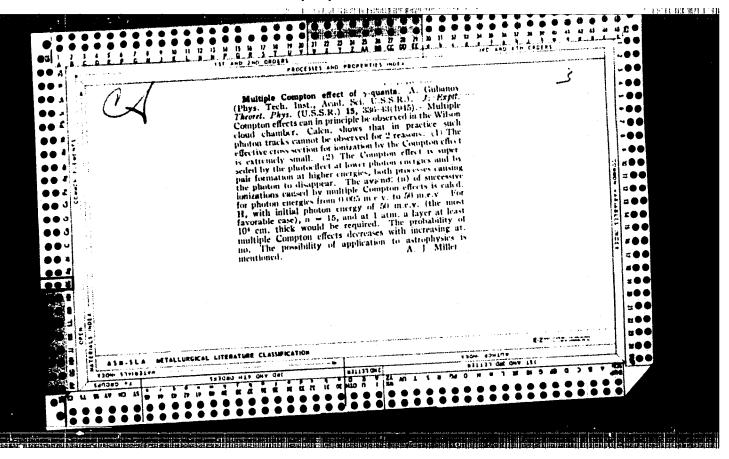
Larkspur Delphinium confusum M. Pop., a new medicinal plant.
Bot. zhur. 49 nc.78997-1002 Jl 164 (MIRA 1788)

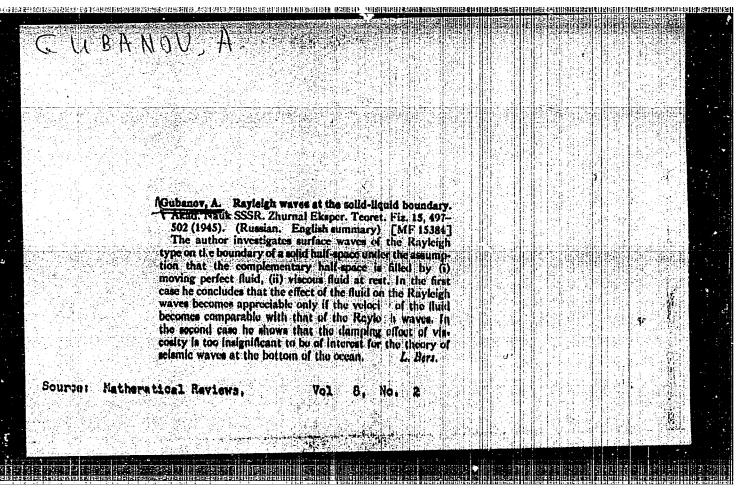
1. Vsesoyumnyy naucimo-isaledovatel skiy institut lekarstvennykh i aromaticheskikh rasteniy, Moskovskaya oblast.

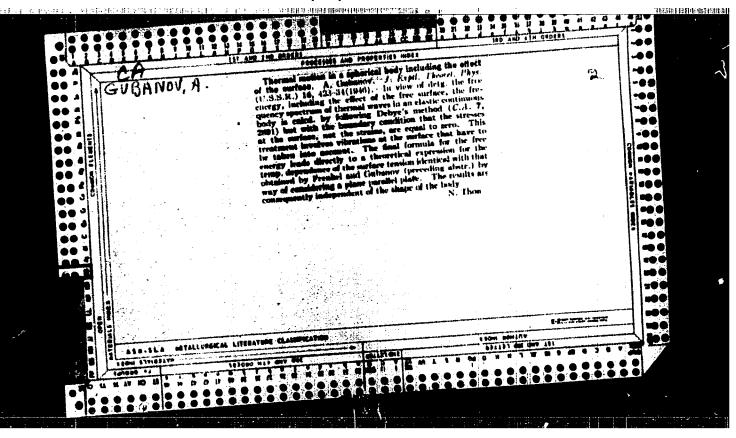
"APPROVED FOR RELEASE: 09/17/2001 CIA-RDP86-00513R000617210012-1

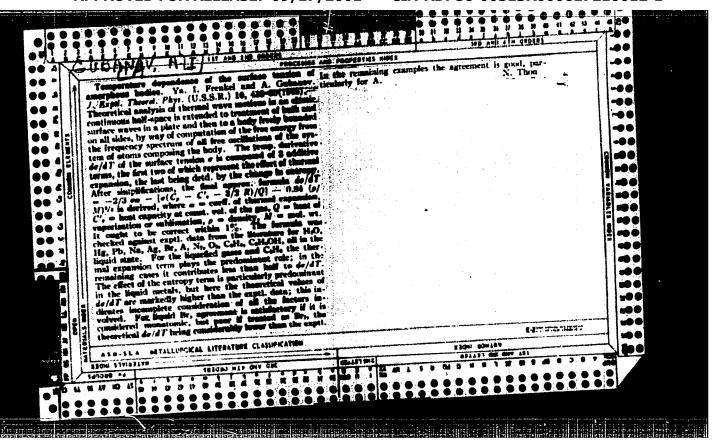


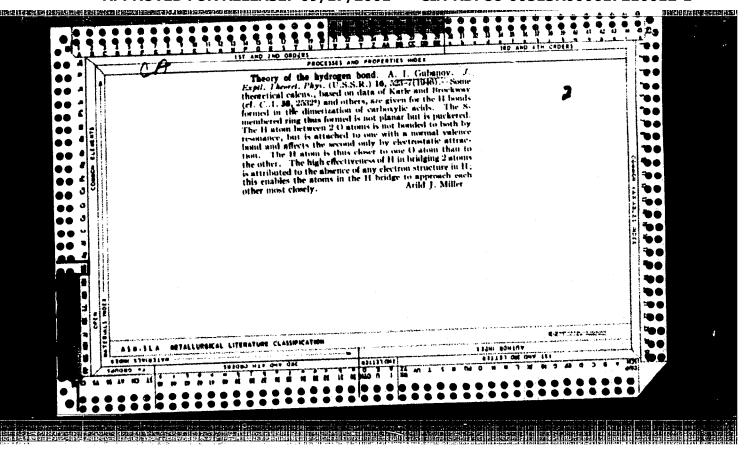


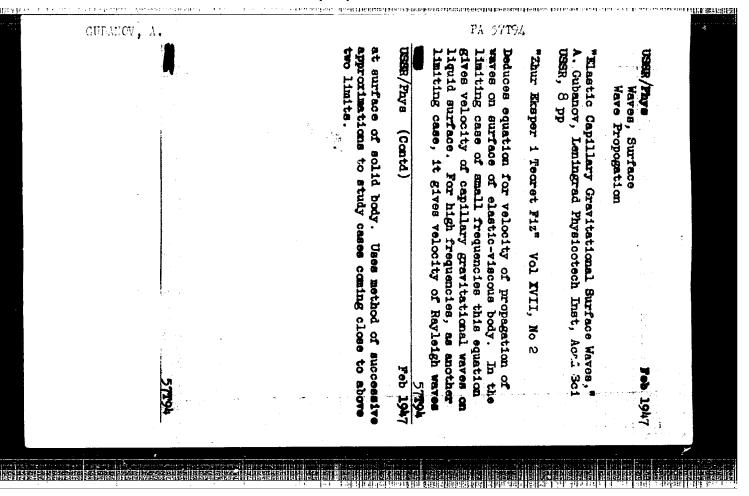












"APPROVED FOR RELEASE: 09/17/2001 CIA-RDP86

CIA-RDP86-00513R000617210012-1

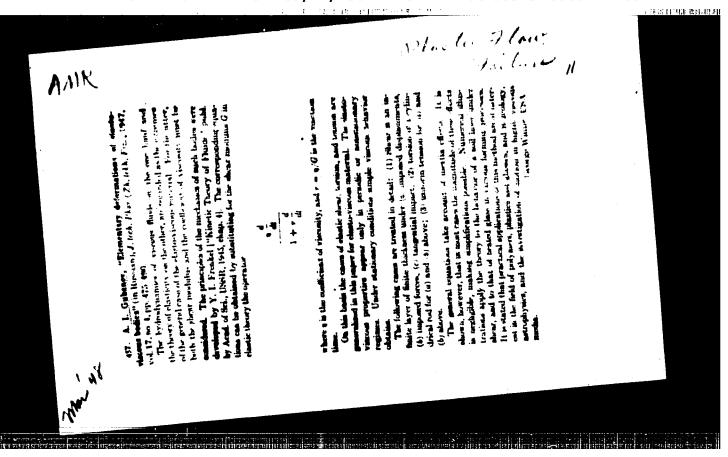
GUBANOV, A.

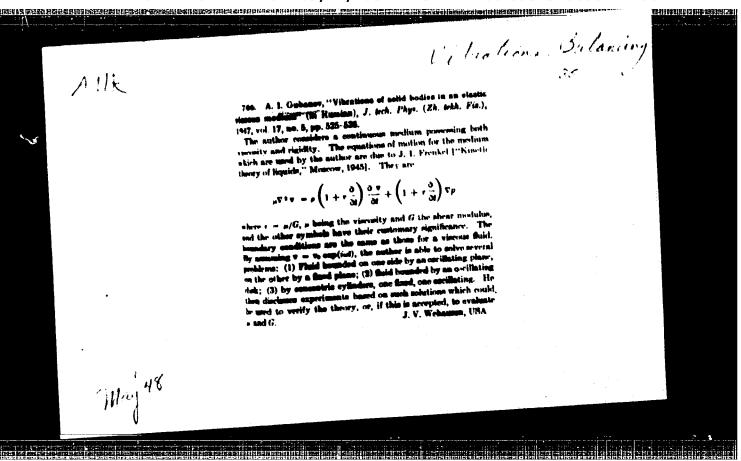
"BER/Mathematics, Applied Feb 1947
Wave propagation

"Dipping of a Cylinder into an Elastic-viscous
Medium," A. Gubanov, 10 pp

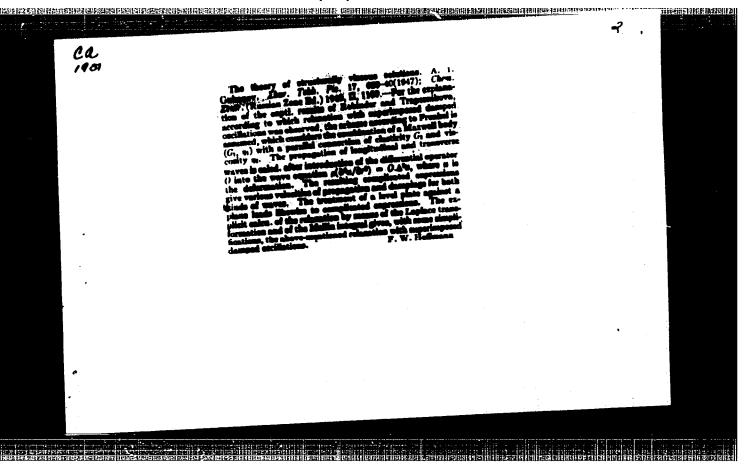
"Zhur Tekh Fiz" Vol XVII, No 2

Selution of Frenkel's equation of propagation of a
disturbance by the method of operators (Iaplacian
transformations).

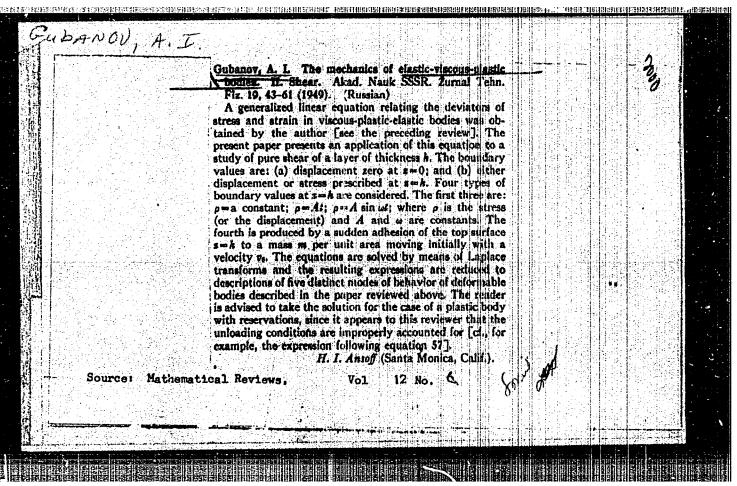




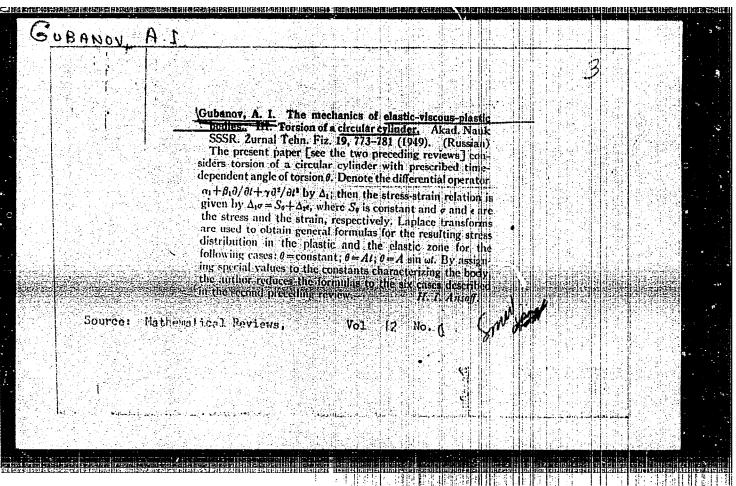
"APPROVED FOR RELEASE: 09/17/2001 CIA-RDP86-00513R000617210012-1

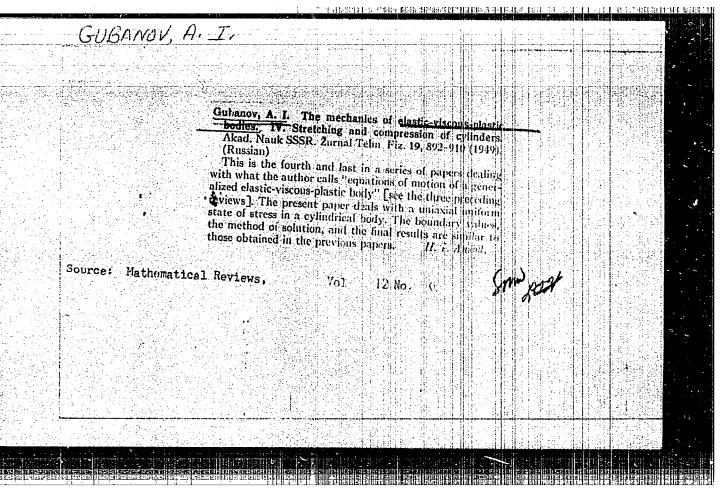


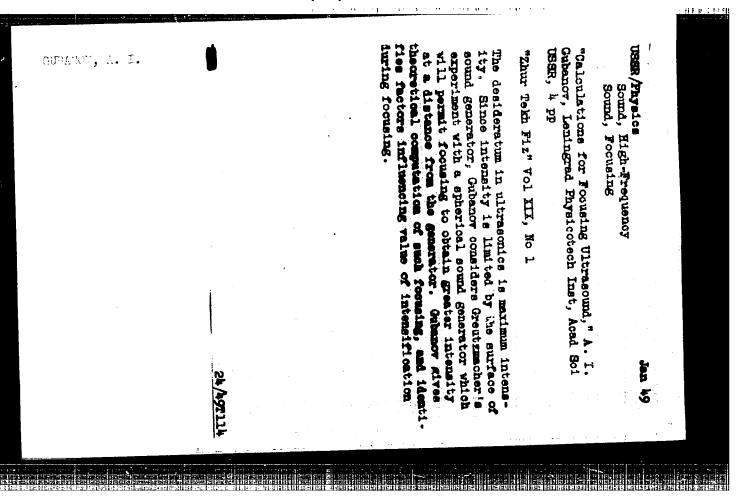
Gubanov, A. I. The mechanics of elastic-viscous-plastic horizon. A second representation of matter. A seneralized aquations of matter. A second representations of matter. A second representation of the second representation of series and parallel combinations of springs, dashpots, and "dry friction" elements which are introduced to account for the yield point in plastic behavior. Linear relations between the stress and strain deviators are assumed for all of these elementary deformation mechanisms. It is shown that under this assumption the most general model leads to a linear relation of the type (1) $\Delta_1 S = S_0 + \Delta_1 E$, where $\Delta_1 = \alpha_1 + \tau_1 d/dt + \nu_1 d^2/dt^2 + \cdots$		Akad. Akad. Akad. 1949). to describe six visclul to d	nown that, by issigning appropriate 2), and S, equation (1) can be made posity a and stiffness G in series) osity and stiffness in parallel); (e, a pair of a and G in parallel with (the previous combination in series only with hardening (no viscosity) out with hardening (no viscosity) out with hardening the friction G). els exhibiting plastic behavior with n is made to distinguish between onditions. Equation (1) is next use a viar-Stokes equation which incre-	
S and E are the and So is the valuable the present discussions.	stress and strain deviators, respecte of S.at.the yield point. For purposion only terms through the second Mathematical Reviews.	oses of terms. The Hencky-Mi	sea yield condition is assumed for	



"APPROVED FOR RELEASE: 09/17/2001 CIA-RDP86-00513R000617210012-1







5-10A+10-1, 4.1.

CA

Theory of the centact of two combonductors with different types of condentity. A.1. Gubanov (Leniegrad Phys. Tech. 18st.). Zhar. Irsh. "Fix." 20," 1287–1301(1980).—The contact of a pair of an electronic (I) and a hole semi-conductor (II), e.g. CiyO-TiGh, has rectifying properties, and the resistance R_0 of the pair at a low applied voltage V is considerably greater than the sum ZR of the resistances of the individual semi-conductors even in the transmitting direction. With V tending to zero, R_0 in either direction tends to the same limit; with V increasing in the transmitting direction, R_0 falls rapidly to ZR, whereas in the blocking direction R_0 falls rapidly to ZR, whereas in the blocking direction it mostly also decreases but much more slowly, if at all. These phenomena may be general in solid rectifiers where rectification possibly takes place actually at the boundary of Z semiconductors of different types, rather—than on a metal/semiconductor boundary. The usual some scheme of the contact shows that passage of electrons or holes across the boundary of Z semiconductors is difficult. Only a very small fraction z abs/Z (where z = electronic charge, Z_0 = distance between the bottoms of the condition of the Z semiconductors of the cortons can pass from II to II, and only an insignificant fraction of holes can pass from II to II. Recombination of electrons and holes at the boundary, producing space charges z which cause at the boundary, producing space charges z which cause at the boundary, producing space charges z which cause

and the sones curve downwards, whereas in II, $\rho > 0$, and the sones curve upwards, also in the absence of a contact potential difference V_{ν} . With the pos. voltage applied to I, the effects are reversed, and the sones curve in the opposite direction. Impoverishment of the boundary region in the direction. Impoverishment of the boundary region in the approan. (i.e. in the absence of current), the chempotential μ_{ν} of II is above μ_{ν} in the blocking direction (V > 0), and below μ_{ν} in the blocking direction (V > 0). Calcus, based on these schemes are made for a model involving a gap (adaorbed layer) between I and II. Zeroth approans, gives the relation between V and the concas, u_{ν} and u_{ν} of electrons and holes in I and II, resp. First approam, then gives the relation between the current V and V and V and V and hence, with the aid of the let approam, the relation between V and V and V and V and V are contact, passage of electrons into the conduction some of II and of holes into the filled sone of I gave no entisfactory quant, agreement with the data for CapOurlo. Use was made of recombination at the boundary, with inclusion of capture of electrons from I on local levels of II with nubequent recombination with holes of II, and capture of holes from II on local levels of I and subsequent recombination with holes of II, and capture of holes from II on local levels of I and subsequent recombination is derived, including all these contributions. Although the contact potential difference V is not a primary factor, it does play a definite role. At $V_{\nu} < 0$, direct recombination is predominant, and V is not direct recombination is predominant, and V is not direct recombination is predominant, and V is not direct recombination in predominant, and V is not

APPROVED FOR RELEASE: 09/17/2001 CIA-RDP86-00513R000617210012-1"

Theory of comboundances with mixed conductivity. A.

1. Combounce (Lembourd Phys. Tech. Inst.). Ann. Abril.

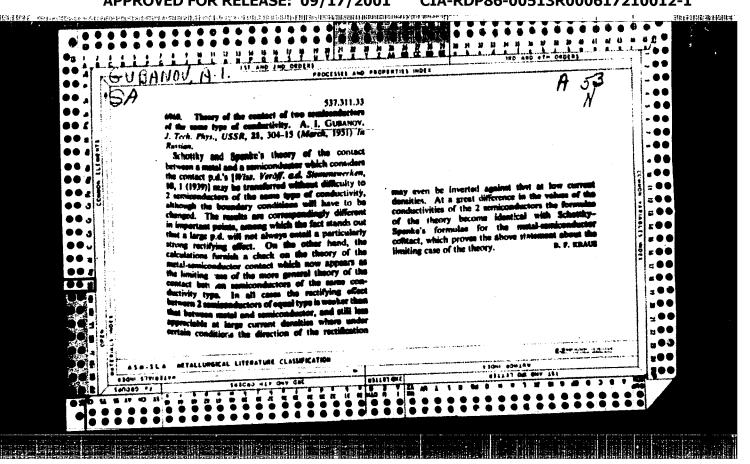
Front. Fiz. 21, 79-74 [Mill.).—The general system of equa
tions is art up for a syndrombuctor in which both types of

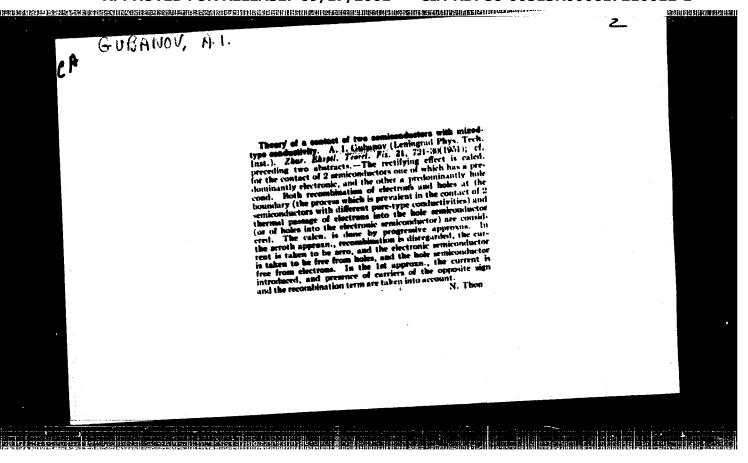
cartiers of current, electrons and holes, are present simul
taneously. The treatment differs front that of Davytov

(Zhu. 18th. Fiz. 7, 2121 [Mill.). Ch. M. Mah. Nanh

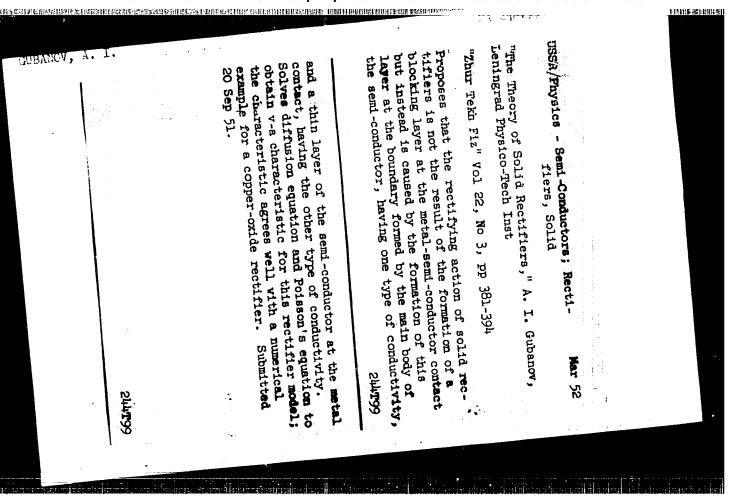
(Ch. 18th. Fiz. 7, 2121 [Mill.). The Mill. Mah. Nanh

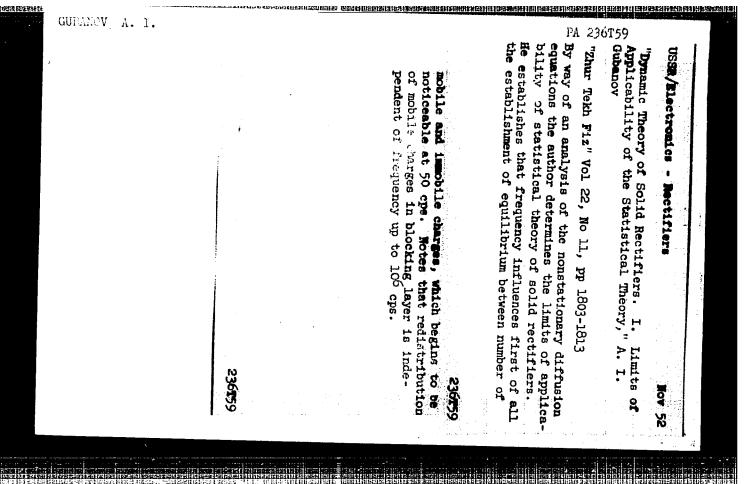
(Ch. 18th. Fiz. 18th. Mill. Mill.





GUBANOV, A. I. Feb 52 USSR/Physics - Semiconductors "Theory of Contact of Metal With Semiconductors in the Case of Great Contact Difference of Potential; A.I. Gubanov, Leningrad Phys Tech Inst, Acad Sci USSR "Zhur Eksper i Teoret Fiz" Vol XXII, No 2, pp 204-213 Calculates the theoretical v-amp characteristic of contact of metal with semiconductor with a potential difference on contact great enough to modify close to the contact the cond type of the semiconductor (electron to hole). Computation is performed for flat, knife, and point contacts. Received 3 May 52.

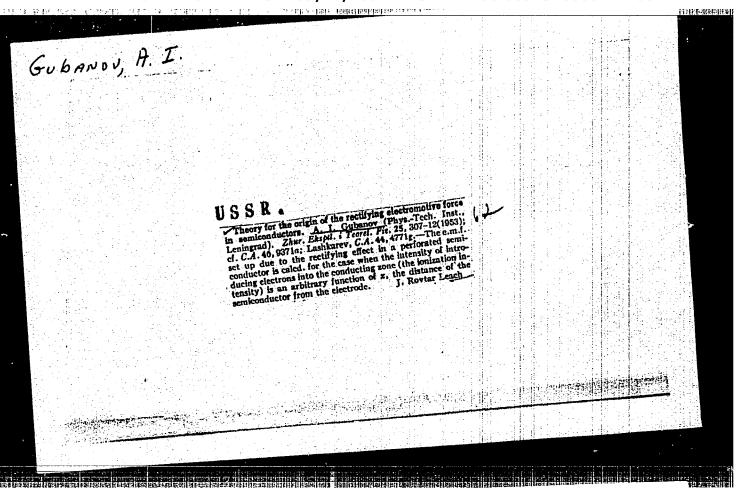




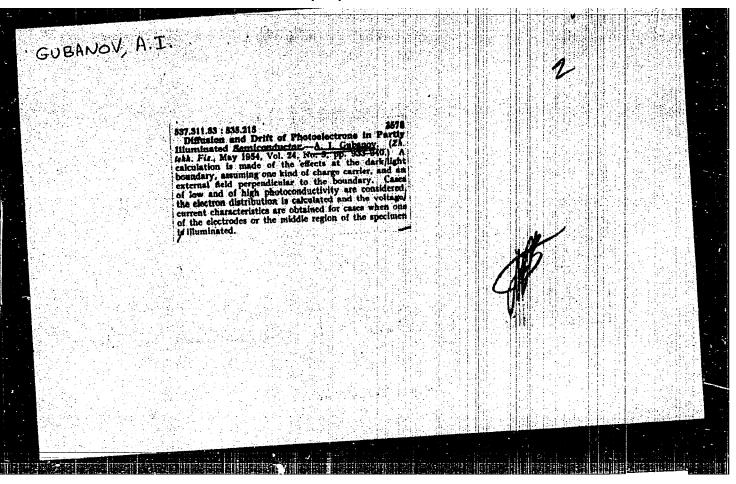
"Dynamic Theory of Solid Rectifiers. II. Frequency Characteristics," A. I. Gubanov
"Zhur Tekh Fiz" Vol 22, No 11, pp 1814-1826

Studies the influence of frequency of alternating voltage with amplitude of several volts, imposed on a solid rectifier, on the distribution of the immobile charges in the blocking layer. Calculates the distribution of potential, and also the conduction and displacement currents as functions of the frequency. Gives numerical examples for case of Cu₂O.

"APPROVED FOR RELEASE: 09/17/2001 CIA-RDP86-00513R000617210012-1



GUBANOV, A. I. FD-605 USSR/Physics - Semiconductor strong-field effect : Pub 153-17/22 Card 1/1 Gubanov, A. I. Author painted the transfer and security and described. : Theory of the strong-field effect in semiconductors Title : Zhur. tekh. fiz. 24, 308-319, Feb 1954 Periodical : Reviews the general theory. Discusses ionization by thermal electrons and by collisions. Derives corresponding formulas taking under son-Abstract sideration the image forces of transistor theory. Indebted to A. I. Anselm. 28 references, including 9 foreign. : July 20, 1953 Submitted



GUBANOV, A. I.

USSR/Physics - Fluid Dynamics

FD-612

Card 1/1

: Pub. 146-2/18

Author

Gubanov, A. I.

Title

alkalikaritati kiliki aliminta Zonal theory of a one-dimensional model of a fluid

Periodical

Zhur. eksp. i teor. fiz. 26, 139-149, February 1954

Abstract

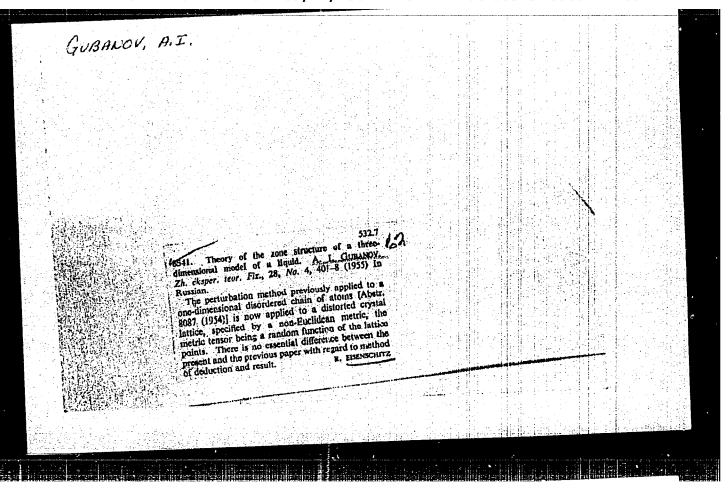
: Examines the behavior of an electron in the field of a one-dimensional chain of atoms by means of the methods of quantum mechanics. Using the solution of the Schroedinger equation in a deformed coordinate scale the author shows that when the distant order in the disposition of the atoms disappears, i.e. when the chain "melts," the energy spectrum of the electron preserves its zonal structure, and only the regions of the zones are slightly displaced. This is in accord with experimental results on conductivity. The author thanks A. I. Ansel'm and O. B. Firsov for having assisted him in the preparation of the

article.

Institution : Leningrad Physicotechnical Institute, Acad Sci USSR

Submitted

: July 18, 1953



GUBANOV, Aleksandr Ivanovich; NOVOZHILOV, Yu.V., redaktor; VOLCHOK, K.M.,

(Skhnicheskiy redaktor

[A theory of the rectifying activity of semiconductors] Teoria

vyprismliminshchege delutria poluprovodnikov. Moskva, Gos. ind-vo

tekhniko-teoret. lit-ry, 1956. 348 p.

(Semiconductors)

(Semiconductors)

GUBANOV, A1

USSR / PHYSICS

CARD 1 / 2

PA - 1264

SUBJECT

AUTHOR TITLE

The Length of the Free Path of an Electron in Liquid and in

Žurn. techn. fis, 26, fasc.8, 1651-1656 (1956)

reviewed 9 / 1956

Apart from the reasons enumerated in some of the author's previous works for PERIODICAL the diffusion of electrons in liquids, the following may still be mentioned: 1.) On the occasion of the displacement of atoms also the maximum and the minimum of the retential may change. 2.) In all cases even greater deformations of structure than those already found exist. All this tends to show that the potential of the field is not a strictly periodic function of the deformed as was previously believed. The character of these deviations remains the same as in the case of admixtures and other defects in crystal. The existence of defects causes an additional diffusion of the electrons. In the course of the present work diffusion was computed with the help of the method applied when computing alloy resistances. Next, the total reciprocal of the free path of the electron in amorphous bodies was computed and its dependence on the electron energy and on temperature is explained. Like in the case of the theory of alloy resistances the potential was divided into a periodic part and into a part deviating from it. Formulae taken from the author's previous works were used in order finally to obtain the free path:

APPROVED FOR RELEASE: 09/17/2001₆) GIABRDP\$6-00513R000617210012-1" Zurn.techn.fis, 26, fasc.8, 1651-1656

Afis, 26, fasc. 8, 1651-1656 (1956) Salar (1956)
$$\frac{\pi}{a^3 k^2} \left(\frac{dE}{dk} \right)^2 = \frac{1}{\left| \overline{V'_n} \right|} \left(1 + \frac{2}{9} a^2 k^2 \right)$$

The lower cross bar above the Vi denotes the average value after the n-th elementary cell, the upper one that after all elementary cells. In order to be able to take account of all different kinds of diffusion of the electron in a liquid, the reciprocal free paths must be added. According to the formulae developed in previous works the corresponding free paths for metals, atomic semiconductors, and ion semiconductors were computed. For these cases also the dependence of the length of the free path on the energy of the electron and on temperature was computed. In the case of temperature dependence it is necessary to distinguish between two cases. In the case of real liquids the exponent depends on the temperature and decreases without interruption during the process of heating the liquid from melting- to critical temperature. At present it is not possible to interpret this dependence theoretically, but it may be obtained, in the most favorable of cases, as the result of radiographical or electro-

Glasslike or amorphous semiconductors present a more simple problem. Here the exponent is "frozen" and no longer dependent on temperature. Temperature dependence is clearly expressed by the formulae.

INSTITUTION:

GUBANOV, A. 1.

USSR / PHYSICS SUBJECT

CARD 1 / 2

PA - 1384

GUBANOW, A.I., MAKOWSKIJ, L.L.

AUTHOR TITLE

On the Article by K.B.TOLPYGO and I.G.SASLAWSKAJA on "Bipolar Diffusion in Semiconductors in the Case of Strong Currents".

Zurn.techn.fis, 26, fasc. 9, 2126-2127 (1956) PERIODICAL

reviewed: 10 / 1956 Issued: 10 / 1956

On the occasion of the solution of the system of differential equations in the work by K.B.TOLPYGO and I.G.SASLAWSKAJA(Zurn.techn.fis, 25, 955 (1955) the method of successive approximation was incorrectly used. The authors investigated the system of equations (11a), (11b), (11w') and (11g'). They attempt to find the solution of (16) by exploiting the smallness of the parameter A which takes the recombination of the electrons and holes into account. The function (0), however, must necessarily appear in the following form according to

the equation (11g'): $\begin{cases} (\Theta) = \frac{F-1}{A} & (\Theta) \\ (\Theta) & (\Theta) \end{cases} + \begin{cases} (\Theta) + A \\ (\Theta) & (\Theta) \end{cases}$ (a)

Let us now investigate the equation (11a). By inserting the unknown functions which are expressed in series according to A-powers, we obtain:

$$\frac{1}{\lambda} \left[\frac{dN_0(\xi)}{d\xi} + A \frac{dN_1(\xi)}{d\xi} + \dots \right] = \Theta + N_0 y_0 + A(N_0 y_1 + y_0 N_1) + \dots$$

It should be noted that on the occasion of approximation towards zero, the

 $m z_{urn.techn.fis}, \, 26$, fasc.9, 2126-2127 (1956) CARD 2 / 2

equation obtained in the special case coincides with the initial equation. By dividing the equation (11a) by the equation (11g'), as was done by the authors (the parameter b is neglected in this equation), we obtain the

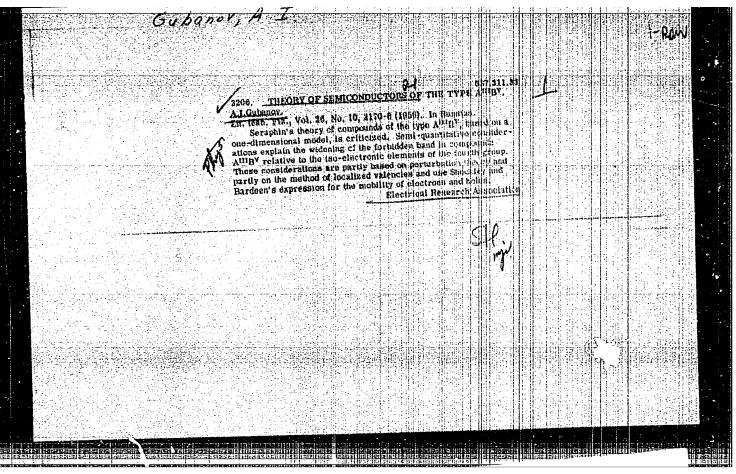
equation (15a).

In order to prove that such an operation contributes absolutely nothing towards solving the task we represent both parts of the equation (11g') in form of a series with respect to the powers A:

$$\frac{1}{A} \frac{d\xi_{-1}}{d\theta} + \frac{d\xi_{0}}{d\theta} + A \frac{d\xi}{d\theta} + \dots = \frac{\lambda}{A} \left[\frac{1}{N_{0}z_{0}} + A (\dots) + \dots \right], (b)$$
where
$$\frac{d\xi_{-1}}{d\theta} = \frac{\lambda}{N_{0}z_{0}}$$
 etc.

It is clear that, by multiplying both parts of the equation (11a) by one and the

same series, the equation (11a) is again obtained by approximation towards zero. The error committed by the authors consists in having failed to consider the fact that the quantity on the left side of the equation (15a) is proportional to the parameter A as follows from the equation (11g'), and from physical deliberations. Actually, 8 is the relative density of the electron current which, in the flat case and with lacking recombination (with A=0) is constant - Therefore Θ is proportinal to A in first approximation. In view of the fact that the results obtained by K.B.TOLPYGO and I.G.SASLAWSKAJA are based upon an incorrect solution of equations they are of doubtful value. INSTITUTION:



GUBANOY, ESSR/Physical Chemistry.

Liquids and Amorphous Bodies.

B--6

Gases.

Abs Jour: Ref Zhur-Khimiya, No 5, 1957, 14560

Author

A. I. Gubanov

Inst Title Electron Scattering in Liquid in Consequence of Distur-

bance of Long Range Order

Orig Pub:

Zh. eksperim. i teor. fiziki, 1956, 30, No 5, 862-872

Abstract:

With a view to evolve the zonal theory of liquids proposed by the author (RZhKhim, 1956, 35216, 38904), the motion of electrons in a liquid was considered using motion of electrons in a liquid was considered using Cartesian co-ordinates. Wave functions of zero approximation were proposed; these functions are analogous to imation were proposed; these functions are analogous to enter functions for a cartesian and entirely assertions. Bloch's functions for a crystal and satisfy equations differing from Schrödinger's equation by minor terms. The electron scattering in liquid in consequence of the disturbance of the long range order of atom arrangement was computed and the dependence of the free path length

Card 1/2

Upon rnysical Chemistry. Liquids and Amorphous Bodies. Gases.

B-6

Abs Japprovedቸውሮ የድርድልያቴ፣ 09/17/2001 ¹⁴ሮያል-RDP86-00513R000617210012-1"

of an electron, corresponding to this scattering, on Abstract: its wave number was obtained in the region of small figures.

Card 2/2

STRIECT

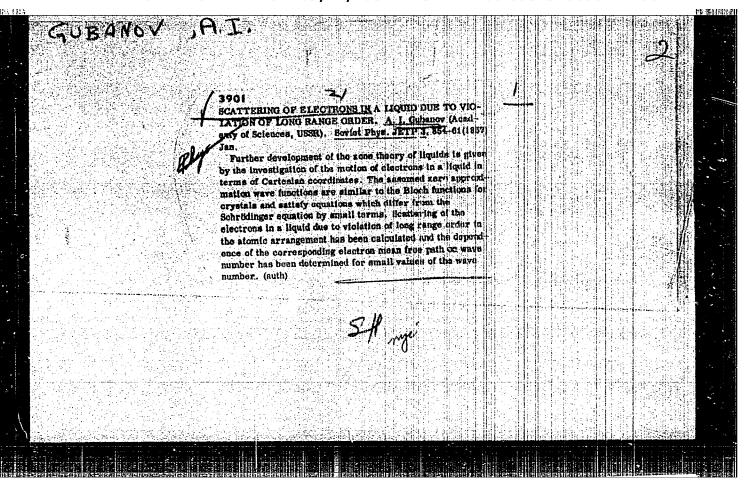
Zurn.eksp.i teor.fis, 31, fasc. 3, 462-472(1956) CARD 2 / 2 PA - 1662

tron by phonons which does not satisfy the momentum conservation theorem within the electron-phonon system, is possible, for, because of the destruction of the distant order on the occasion of the emission or absorption of a phonon by an electron, the lattice is able to absorb a certain additional momentum. This additional scattering of electrons is here called "phonon liquid scattering".

There follows the computation of the free length of path which is due to this phonon-liquid scattering. Also generalization for the case that, besides the electric field, there exists a temperature gradient or a concentration gradient, is not difficult. In most cases it is possible to disregard phonon liquid scattering in semiconductors; only at high temperatures does it make a slight correction to ordinary thermal scattering.

In conclusion the case of optic oscillations (and in what it differs from acoustic oscillations) is discussed. The usual thermal scattering of electrons in liquids can, with the exception of slight corrections, be computed by means of the same formulae that are employed in the case of solids. However, in liquids there is an additional phonon-liquid scattering, which is able to play an important part in liquid metals but not in semiconductors.

INSTITUTION: Leningrad Physical Technical Institute of the Academy of Science in the USSR



"APPROVED FOR RELEASE: 09/17/2001 CIA-RDP86-00513R000617210012-1 的数据表示的问题是还是我们是这些表现的记录是完全还有要组化的主义是这个文明和自己,但是是一个人,但是是一个人,但是是一个人,也是是一个人,也是一个人,也是一个人

Gusanov, HI

PA - 2353

AUTHOR: Title

Free Path of an Electron in Liquid and Amorphous Conductors (Dlina svobodnogo probega v zhidkikh i amorfnyk provodnikakh, Russian).

Izvestiia Akad. Nauk SSSR, Ser. Fiz., 1957, Vol 21, Nr 1, pp 104 -

PERIODICAL:

ABSTRACT:

104 (U.S.S.R.)

Reviewed: 5 / 1957

Received: 4 / 1957 In previous works the authors were able to show the following by

solving the Schrödinger equation in the deformed system of the coordinates & (in which the self-consisting-field for the electron is approximately periodic): he energy-spectrum has, as in the case of crystals, a zone-structure. The electrons are described by Bloch-

functions in the system of the coordinates & i.e. the electrons are

quasi-free.

The destruction of the distant order in the liquid, however, leads to a specific "liquid scattering" of electrons. On the occasion of the computation of this scattering, the following was shown: The quasi-Bloch-functions of the form $\psi = u(\xi)e^{ikr}$ (where u is a periodic function of the coordinates ξ) satisfy an equation which differs from the Schrödinger equation by small terms of the order of magnitude &. Here & denotes the order of magnitude of the small deformation of elementary cells in the case of the melting of the

crystal. The author considered the aformentioned function as zero-th approximation and the aforementioned small terms as per-

Card 1/3

PA - 2353

Free Path of an Electron in Liquid and Amorphous Conductors. turbation; by this method he computed the free path l of the electron on the occasion of "liquid-scattering". If the aforementioned quasi-Bloch-functions are used, the scattering of electrons in the liquid by thermal phonons is computed as in the case of crystals; this kind of computation has, however, also certain peculiarities. For the maintainance-law of momentum is not bound to apply in the case of the electron-phonon system because of the lack of a distant order in the liquid. Therefore, also electrons of an additional phonon-liquid scattering occur besides ordinary phonon-scattering. On the occasion of the computation of the free length of path Ph-fl in the case of this scattering for metals as well as for atomic and ion semiconductors it was found that in the case of metals 1 Ph-fl is comparable to 1 Ph. 1 Ph corresponds to the ordinary scattering by phonons. In the case of semiconductors, $l_{Ph-fl} \gg l_{Ph}$ applies. Finally also the free length of path l_d for scattering by the defects and deviations from periodicity was computed. For the resulting free path $(1/1) = (1/1_{fl}) + (1/1_{Ph}) +$ + $(1/l_{Ph-fl})$ + $(1/l_d)$ applies. In the case of atomic semiconductors

Card 2/3

PA - 2353

Free Path of an Electron in Liquid and Amorphous Conductors.

e.g. the dependence (1/1) = A + Bu applies. Here u denotes the energy of the electron, A and B are functions of temperature. (No illustrations).

ASSOCIATION: Leningrad Physical-Technical Institute of the Academy of Science of the U.S.S.R.

PRESENTED BY:

SUBMITTED:

AVAILABLE: Library of Congress.

and the state of t

Card 3/3

SUBJECT

PERIODICAL

USSR / PHYSICS

CARD 1 / 2

PA - 1990

AUTHOR TITLE

The Electric Conductivity, the Heat Conductivity, the Thermoelectromotoric Force, HALL'S Constant, and the NERNST Constant

for Amorphous Bodies with Electronic Conductivity

Zurn.techn.fis. 27, fasc.1, 3-11 (1957)

Issued: 2 / 1957

र वर्षा विद्यालया विद्यालया है।

In the course of this work the temperature dependence of various kinetic coefficients for amorphous conductors is computed in consideration of the scattering of electrons which is characteristic of this case and is connected with the lack of a remote order. If we investigate the scattering of electrons by thermal oscillations, we find that in semiconductors a universal free length of path of the electron exists at any temperature T, which holds good in the case of all kinetic processes. The approximation of the effective mass which is amply sufficient for semiconductors is here considered as being all that is necessary. On this assumption it is possible to compute the kinetic coefficients by means of the formulae of the theory of free electrons. Formulae are given for the following kinetic coefficients: electric conductivity, heat conductivity, THOMSON'S coefficient, PELTRIER'S effect, differential thermoelectromotoric force, isothermal electric conductivity in the magnetic field, isothermal HALL'S constant, isothermal NERNST'S constant. The author carried out computations separately for metals at high temperatures of atmit semiconductors and for ion semiconductors with weak binding between the oscillations of the lattice and the motion of the electrons. The case with strong binding

CARD 2 / 2 Zurn.techn.fis, 27, fasc.1, 3-11 (1957)

PA - 1990

(polarones) in amorphous ion semiconductors can be investigated separately.

At first liquid metals are computed at T >> 0 in consideration of the complete formulae by L. BRILLOUIN, Quantum Statistics (here the author quotes the

The author then computes the kinetic coefficients (more accurately expressed the explicit temperature dependence of these kinetic coefficients) for

amorphous atomic semiconductors. In conclusion amorphous semiconductors with ion bindings are studied. In this case various expressions hold for the free length of path of the electron at high as well as at low temperatures. Because of the very complicated nature of the expressions for this free length of path computation must be carried out in a rougher manner than in the case of atomic semiconductors. Computations are carried out separately for high and for low temperatures. At high temperatures the temperature dependence of the kinetic coefficients has a simple form. In the case of the approximation investigated here this is the same form as in the case of crystalline ion semiconductors. However, in the case of genuine liquids the temperature dependence of the kinetic coefficients has a very complicated character and can be represented only graphically for the individual concrete substances.

INSTITUTION: Physical-Technical Institute, LENINGRAD

and the state of t

CUBANOV, A. I.

AUTHOR:

57-11-13/33

TITLE:

Gubanov, A. I. On the Alteration of Semiconductor Properties with Fusion (Ob

izmenenii svoystv poluprovodnikov pri plavlenii)

PERIODICAL:

Zhurnal Tekhn. Fiz., 1957, Vol. 27, Nr 11, pp. 2510-2516 (USSR)

ABSTRACT :

With regard to the fact that in the earlier work of the author (ZhETF, 26, 139, 1954) in the case of the classification of the width of the forbidden zone not even the sign of the alteration of width with fusion could be given the whole investigation is carried out again here. The relatively small alterations of the energy level, of effective masses etc. with the fusion of crystals or with their transition to amorphous state are calculated the coordination number being maintained. The alteration of the zone width as well as of the effective masses of current carriers with fusion is brought into connection with the change of crystals with deformation. In the end the remarks on the alteration with the fusion of the free length of path of the "movability-currentcarriers" as well as of electric conductivity are given. The author shows that the free length of path of the current carriers with fusion must decrease for two reasons. In the case of an admixture-conductivity the matter is more complicated, and the problem of the admixture-level in liquid or amorphous bodies can, at

Card 1/2

On the Alteration of Semiconductor Properties with Fusion.

57-11-13/33

present, not be regarded solved. There are 8 Slavic references.

ASSOCIATION:

Leningrad Physical-Technical Institute AN USSR (Leningradskiy

fiziko-tekhnicheskiy institut AN SSSR)

SUBMITTED:

May 27, 1957

AVAILABLE:

Library of Congress

Card 2/2

Gubanov, A. I., Lun'kin, Yu. P.

57-11-25/33

TTTLE:

Kinetics Equations of Gas Dissociation with Account of Diffusion (Uravneniya kinetiki dissotsiatsii gaza s uchetom diffuzii)

Zhur_nal Tekh. Fiz., 1957, Vol. 27, Nr 11, pp. 2631-2639, (USSR)

ABSTRACT:

PERIODICAL:

A system of equations is deduced which describe the behavior of the gas in non-equilibrium diffusion and dissociation. The cases for a diatomic gas and for air as a five-components-mixture are investigated. The temperature dependence of the kinetic coefficients in the equations obtained is evaluated. However, as it is based on a series of approximations it requires an additional examination by experiments as well as a precision. It is shown that the specific velocities of the dissociation reaction in the case of air will essentially differ only by the exponential multiplicands that depend on the dissociation. It is shown that the thermo-diffusion coefficient is a very comprehensive expression and not a function of the binary thermo-diffusion coefficient, but that it is determined by the conduct of the whole multi-components-system. If a turbulent diffusion occurs in a multi-components-mixture it is difficult to put down the expression for the flow of particles in a general form. In this case it is difficult to approach from the point of view of the statistical theory of turbulency and it is more reasonable to base on the semi-empirical phemenological theory There are 2 Slavic references.

Card 1/2

CIA-RDP86-00513R000617210012-1" APPROVED FOR RELEASE: 09/17/2001

Kinetics Equations of Gas Dissociation with Account of Diffusion. 57-11-25/33

ASSOCIATION: Leningrad Physical-Technical Institute AN USSR (Leningradskiy fiziko-tekhnicheskiy institut AN SSSR)

SUBMITTED: May 3, 1957

AVAILABLE: Library of Congress

Card 2/2

APPROVED FOR RELEASE: 09/17/2001 CIA-RDP86-00513R000617210012-1"

es erseás elem cuels hesipe

AUTHOR:

Gubanov, A. I.

sov/57-23-9-26/33

TITLE:

Reflection and Refraction of Shock Waves at the Boundary of Two Media (Otrazheniye i prelomleniye udarnykh voln na granitse dvukh sred) I. Normal Incidence (I. Sluchay normal'nogo padeniya)

PERIGDICAL:

Zhurnal tekhnicheskoy fiziki, 1958, Nr 9, pp. 2035-2040 (USSR)

ABSTRACT:

The reflection and the refraction of shock waves of high intensity are essentially different from the reflection and the refraction of acoustic waves. This is a study of the reflection and the refraction of plane shock waves at the plane boundary of two media. This is done in a manner similar to that found in references 1 - 5 for sonic waves without a linear approximation. This investigation is limited to the most simple case, that of normal incidence. The next paper will deal with the more complicated case of a sloped incidence. It is assumed that the laws of the propagation of the shock waves in each of the homogeneous and contacting media are known and that the basic conditions prevailing at the front of the incident, of the reflected, and of the transmitted wave can be written down. A general equation is derived. This equation makes it possible to

Card 1/2

SUV/57-28-9-26/33

Reflection and Refraction of Shock Waves at the Boundary of Two Media. I. Normal Incidence

and the same

find the equation describing the pressure in the reflected shock wave which is generated by the normal incidence of a plane shock wave upon the plane boundary of two arbitrary media. Approximate solutions are derived for this equation: 1) For two perfect gases. 2) For a shock wave of small intensity. 3) For two slightly differing media. 4) For two greatly differing media. The subject was suggested by N. S. Levchen'. There are 1 figure and 7 references, 7 of which are Soviet.

ASSOCIATION:

Fiziko-tekhnicheskiy institut AN SSSR, Leningrad (Physical and Technical Institute, AS USSR, Leningrad)

Card 2/2

AUTHORS:

Gordeyev, G. V., Gubanov, A. I.

的行動的支持性的形式的影響的大型性的用的分類和表現的大學的影響性質的影響的影響的影響的一般的影響的計畫的影響的計畫的影響的影響的影響的影響的影響的影響的影響的影響的

SOV/57-28-9-28/33

TITLE:

On the Problem of Planta Acceleration in a Magnetic Field

(K voprosu uskoreniya plazmy v magnitnom pole)

PERIODICAL:

Zhurnal tekhnicheskoy fiziki, 1958/Val 28, pp. 2046-2054 (USSR)

ABSTRACT:

This report covers the investigation of a partly ionized plasma contained between two co-axial cylindrical electrodes to which a potential V is applied. It is assumed that a constant and uniform magnetic field is imposed on the plasma. The field strength H is at a direction parallel to the axis of the cyl-

inder (z-axis). Let the current I through the plasma be

considered constant independently of whether the magnetic field is applied or not (this is achieved by a suitable choice of the voltage U). The electric and the magnetic field arranged cross-wise effect a motion of the charged particles in the same direction at right angles to the fields. They carry away the neutral gas atoms effecting a circular motion of the gas about the inner cylinder. The acceleration of the plasma by the external magnetic field is examined. The steady plasma flow is computed taking into account the friction between the plasma

Card 1/2

On the Problem of Plasma Acceleration in a Magnetic Field

and the electrodes. The dependence of the stream velocity upon the magnetic flux and upon the electrode radii and the energy necessary for the maintenance of the current is computed. The computations showed that at realizable dimensions of the apparatus supersonic velocities can be attained in the plasma flow. There are 2 references, 1 of which is Soviet.

ASSOCIATION: Leningradskiy fiziko-tekhnicheskiy institut AN SSSR (Leningrad Physical and Technical Institute, AS USSR)

Card 2/2

是一个人,我们就是一个人,我们就是一个人,我们们我们的人,我们们就是一个人,我们们我们们的一个人,我们们我们的人,我们们我们的人,我们们我们的人,我们们我们的人 py: -7-28-10-2740 24(6) Gubanov. a. 1. CUTIFIED: bependency of the Zonal Structure of an eller of an AB, Type TATES: Unon Regularity (2007) impost sommer structury splays tips AB, ot uporyadochennosti) Znurmi: tekonicheskoy fiziki. Vol. 28, 62 70. up 2109-2114 (40 8) 1958 Partionic L: tarking to wine method of rigidly bound electrons and attriceing the averaged node potentials smirnov (set) carried out ar SBTBUS": approximate computation of the electronspectrum in a partial ordered solution. The general equation for the computation of the energy of an electron in the alley, which is a function of the wave vector K, takes the shape of equation (1). In the case of a binary allow with an stopic concentration of the first component c, with a node concentration c. Tiret kind v, and with the degree of ordering η the mean potentials produced by the nodes of first and second kind are, sucording to imirnov, specified by equations (3). In combination with the experimental investigations by Romer (Ref 2) these formulae are applied to an allow of an iP, type (which cryptailizes in the cubic face-Card 1/3

() 15 可控制 计图形性图像图像操作中操作的基础的基础的形式和操作中中等19 1 16 mm 1 1 1 1 JOV 1-7-28-10-9/40 Dependency of the Zonal Otructure of an Floy of an EB, Type $\operatorname{Up}_{\mathbb{Z}^n}$ n Regularity centered lattice with the edge length a). (aCu, is the alloy the investigation of which has been driven to the farthest point. In this case the elementary cell possesses 4 sites (one of first and three of second kind), $\mathbf{v} = \frac{1}{4}$ and equation (1) is transferred into equation (6). This equation (6) can callity be solved for three cases: 1) The alloy is completely disordered $(\eta=0)$ as all sites are, on the average, equivalent, 2) k (wave vector) coincides with the cube diagonal. () k coincides with the warface of the Brillouin-(Brillywon)-zone with a cube shape. From the formulae (7), (10), and (11) the electron energy can immediately be determined for a stoic dometrically (c - $v = \frac{1}{3}$) absolutely ordered $(\eta = 1)$ alloy. It is snown that the zonat structure is dependent upon a parameter : Formila (16). This parameter is distinctive of the incluence of the degree of ordering in the alloy. In the paper by Komar (Ref 2) the hypothesis as advanced, that when the alloy 100, 10 breezes & corresponding re-arrangement of the zones (about places The com-

Dependency of the Zone: Structure of an Alloy of an AB, Type Upon Regularity

putations presented in this paper offer a definite substantiation of this hypothesis and a theoretical explanation for the change in sign of the carriers during ordering of the alloy. 1. P. Komar, Member of the Academy of Potences Ukrainskaya SUR. suggested the subject and discussed it afth the author. There are I figure and 3 references, 5 of which are Loviet.

SUBMITTED: February 28, 1958

Card 3/3

s/169/61/000/010/037/053 D228/D304

AUTHORS:

Gordeyev, G. V., and Gubanov, A. I.

TITLE:

The question of plasma acceleration in a magnetic field

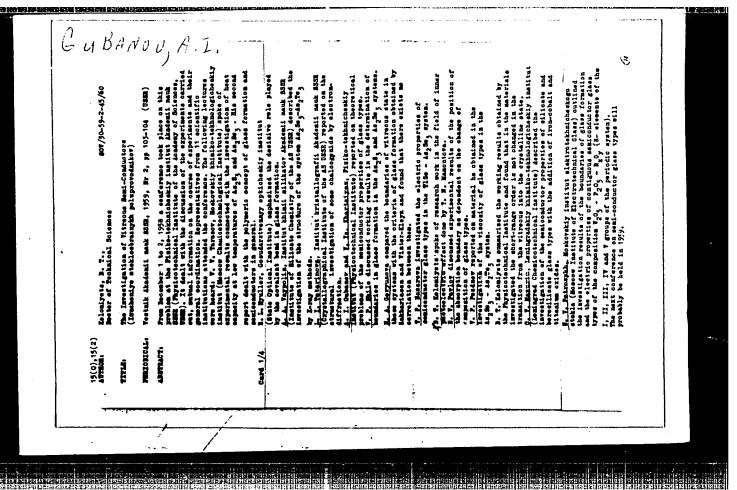
PERIODICAL:

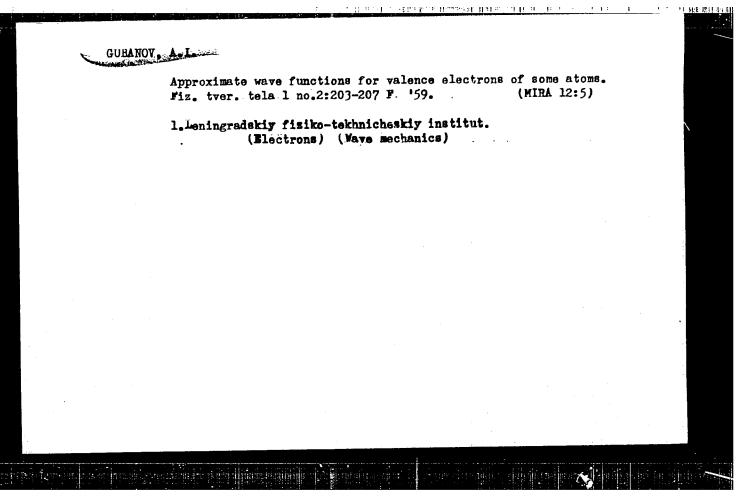
Referativnyy zhurnal, Geofizika, no. 10, 1961, 12, abstract 10G66 (V sb. Vopr. magn. gidrodinamiki i dinamiki plazmy, Riga, AN LatvSSR, 1959, 73, diskus.,

74-75)

TEXT: The movement of plasma between two infinite, coaxial, cylindrical electrodes in an external, axial magnetic field is considered in the stationary case. The calculation was made in the hydrodynamic approximation—disregarding the Hall current, but with due allowance for the plasma's viscosity. [Abstracter's note: Complete translation.]

Card 1/1





24(6) 24,7700

66250

AUTHORS:

Gubanov, A.I., Nran'yan, A.A.

SOV/181-1-7-6/21

TITLE:

Application of the Method of Equivalent Orbits to the Investigation of Zone Structure in Compounds of Type AIIIB

PERIODICAL:

Fizika tverdogo tela, 1959, Vol 1, Nr 7, pp 1044-1052 (USSR)

ABSTRACT:

The present paper compares the method of strong coupling with the method of equivalent orbits. After mentioning that in the method of strong coupling the electron is described by Bloch sums, the author proceeds to explain the essence of the other method. It is based on the fact that the orbits of the molecular type (crystal or molecule) are formed by orbits which are localized in various parts of the system. If there are symmetries in the system considered, the localized orbits separate into groups in such a way that all orbits of one group convey the same character of charge distribution in the space, and differ only in position and orientation. These orbits are said to be equivalent. Molecular and equivalent orbits have a unitary transformation (1) in common. This unitary matrix is set up according to G.G. Hall. The authors do not agree with the opinion expressed by Slater and Koster (Ref 10) that the method of equivalent orbits be of no advantage. The authors claim, instead, that it permits a linear combination,

Card 1/3

66250

Application of the Method of Equivalent Orbits to the SOV/181-1-7-6/21 Investigation of Zone Structure in Compounds of Type AIIIBV

经证券公司,我们的国际的工程,我们是一个人的工程,我们是一个人的工程,我们是一个人的工程,我们是一个人的工程,我们也是一个人的工程,我们也是一个人的工程,我们也是一个人的工程,我们也是一个人的工程,我们也是一个人的工程,我们也是一个人的工程,我们就是一个人的工程,我们也是一个人的工程,我们也是一个人的工程,我们也是一个人的工程,我们也是一个人的工程,我们也是一个人

of the functions of free atoms, that the order of the secular equation becomes lower, etc. They agree with the above-mentioned authors, however, in defining the method of equivalent orbits as an approximation to the general method of strong coupling. In opposition to Hall, the authors believe that the method of equivalent orbits offers the possibility of computing conduction bands. An analytical formula for the dependence of energy on the wave vector is then derived by the method described, and finally, an analytical survey of the results obtained is given. The purpose of the computations made by the authors was that of determining the rules governing the change in the zone structure in transitions from one substance to another in the group of compounds AIIIBV and in transitions from elements of the 4th group to AIIIBV. Such a compound was found to exhibit four valency bands and four conduction bands, and it is further stated

Card 2/3

66250

Application of the Method of Equivalent Orbits to the III $_{\rm BV}^{\rm SOV}/181-1-7-6/21$ Investigation of Zone Structure in Compounds of Type AIII $_{\rm BV}^{\rm SOV}/181-1-7-6/21$

that in the transition from 4th group elements to AIIIBIV

compounds within 4-valency bands as well as within the conduction bands an additional forbidden zone appears which does not occur in elements of the 4th group. There are 17 references, 3 of which are Soviet.

ASSOCIATION:

Fiziko-tekhnicheskiy institut AN SSSR, Leningrad

(Physical and Technical Institute of the AS USSR, Leningrad)

SUBMITTED: June 26, 1958

Card 3/3

24(2), 24(3)

AUTHORS: Gubánov, A. I., Gashimzade, F. M.

TITLE: Investigation of the Symmetry of the Energy Bands of

Electrons in the Type Crystals CdIn2Se4

PERIODICAL: Fizika tverdogo tela, 1959, Vol 1, Nr 9, pp 1411 - 1416 (USSR)

ABSTRACT: The purpose of the present paper is that of investigating the energy spectrum of the electrons in semiconductors of the CdIn₂Se₄ type by means of a group-theoretical method; this compound crystallizes in tetragonal syngony in the D_{2d} space

group, while most other such compounds exhibit a S_4^2 structure. First, the symmetric properties of these structural groups investigated here $(D_{2d}^1 - P\bar{4}2m)$ are carefully analyzed in order

to obtain provisional data of the type of energy bands. For the symmetric points of the Brillouin zone given in a figure, table 1 show the characteristic values for single space

groups, and table 2 for double ones. Table 3 likewise offers representations of single and double groups. The conditions

Card 1/2

Investigation of the Symmetry of the Energy Bands of SOV/181-1-9-15/31 Electrons in the Type Crystals CdIn₂Se₄

of consistency for single and double groups are compiled in table 4 and 5, respectively. The band structure is investigated by means of these data and those given in tables 6 and 7. It is shown that without considering the spin, the energy limit is in the center of the Brillouin zone. On principle, the limits can be situated in the points (000), $(\frac{\pi}{a}, \frac{\pi}{a}, 0)$, $(00, \frac{\pi}{b})$, $(\frac{\pi}{a} \frac{\pi}{a} \frac{\pi}{b})$, $(0 \frac{\pi}{a} 0)$, and $(0 \frac{\pi}{a} \frac{\pi}{b})$. Theoretical considerations (Ref 7) and experiments with the cyclotron resonance showed the edge valence band to be situated in the point K = (000) and to be triply degenerated. All this holds without consideration of the spin. It is shown that the group theory may not be employed to determine, which of the bands (T2) lie higher than others. Yu. Firsov is mentioned in the text. There are 1 figure, 7 tables, and 7 references, 2 of which are Soviet. Leningradskiy fiziko-tekhnicheskiy institut AN SSSR (Leningrad Institute of Physics and Technology of the AS USSR)

ASSOCIATION:

SUBMITTED:

January 19, 1959

Card 2/2

的表面感染。中国是全部的1918年,中国的1920年的中国主义员员的1918年,在1918年,在1918年,1918年

S/181/60/002/02/11/033 B006/B067

AUTHORS:

Gubanov, A. I., Gashimzade, F. M.

TITLE:

The Structure of the Energy Bands in Semiconductors of the

Cain₂Se₄-Type

PERIODICAL: Fizika tverdogo tela, 1960, Vol. 2, No. 2, pp. 255-260

TEXT: In continuation of a previous, paper (Ref. 1) the authors investigate the energy spectrum of CdIn_Se Crystals by means of the method of

localized valences. E(k) in the range of the energy extremes at k=0 was obtained by a perturbation-theoretical method developed by Shockley and Dresselhaus; spin-orbit interaction was taken into account in first approximation. The possible forms of the equipotential surfaces near the extremes were investigated by taking the spin into account. The possible positions of the energy extremes of the electrons in $CdIn_2Se_4$ had been

investigated in the paper of Ref. 1. In the method of localized valences molecular functions composed of hybridized atomic functions were used as zero approximation. It can be well applied to CdIn₂Se₄ which shows mainly

Card 1/3

The Structure of the Energy Bands in Semiconductors of the CdIn₂Se₄-Type

S/181/60/002/02/11/033 B006/B067

a covalent bond, and has 8 atoms and 16 valence lines per unit cell. The 16th order secular equation for the energy was obtained in the approximation of the second neighbors. The equations were treated by the group theory, and the 16 solutions were classified into four subgroups each of which is transformed according to one of the irreducible representations Γ_1 , Γ_4 , and Γ_5 of the point-symmetry groups of the crystal. It was found that, as in the case of diamond and sphalerite, roots exist in the approximation of the first neighbors which are independent of k and are transformed according to the irreducible representation of Γ_5 with k=0. In this approximation, Γ_4 and Γ_5 appear degenerate. The band edge is found at k=0, and is triply degenerate as is the case with sphalerite. Fig. 1 shows the approximate structure of the valence band. The effective hole mass in this doubly degenerate band is determined by the interaction integral of the second neighbors. There are 4 figures and 13 references; 6 Soviet, 3 American, 2 British, 1 German, and 1 Japanese.

Card 2/3

原基[5]

VC

The Structure of the Energy Bands in Semiconductors of the CdIn₂Se₄-Type

S/181/60/002/02/11/033 B006/B067

ASSOCIATION: Fiziko-tekhnicheskiy institut AN SSSR Leningrad (Physicotechnical Institute of the AS USSR, Leningrad)

SUBMITTED: May 11, 1959

K

Card 3/3

81362 S/181/60/002/03/20/028 B006/B017

24.7900 AUTHOR:

Gubanov, A. I.

TITLE:

Quasiclassical Theory of Amorphous Ferromagnetics

PERIODICAL:

Fizika tverdogo tela, 1960, Vol. 2, No. 3, pp. 502-505

TEXT: Up to now, theoretical and experimental investigations of the ferromagnetism of crystalline bodies only have been known from publications, and a possible existence of amorphous ferromagnetics has not been investigated. Liquid ferromagnetics have not been investigated as was said in the introduction - because their melting point is above the Curie point; hence, amorphous or vitreous ferromagnetics exist if these substances contain the corresponding elements, and if they have sufficiently low temperatures. Since amorphous vitreous ferromagnetics should be preferred to crystalline ones for many purposes, the author theoretically investigates in this paper their possible existence. It is supposed that the vitreous substance investigated contains only one type of atom which causes ferromagnetism (hence, the exchange integrals of

Card 1/3

Quasiclassical Theory of Amorphous Ferromagnetics

\$/181/60/002/03/20,'028 B006/B017

these atoms must be positive, and their amount must be high compared to that of the exchange integrals between the other atoms and between other atoms and the "ferromagnetic" ones). The system may then be regarded as a one-component system in first approximation; it contains only one type of atom, the others play the role of a neutral medium whose concentration determines the distances between the "ferromagnetic atoms". Furthermore, it is assumed that the distances between neighboring "ferromagnetic atoms" obey statistical laws. The computations were made by the so-called quasichemical method (Fouler and Gugenheim, Ref. 1). L. S. Stil'bans, S. V. Vonsovskiy, and Ya. S. Shur (Refs. 2 and 3) used this method for investigating alloys and ferromagnetism. The results show that the formulas obtained for an amorphous body are, in principle, the same as those obtained for a crystal. A quantitative difference is only found in the dependence of the exchange integral A on the distance of the "ferromagnetic atoms" r; however, A remains positive with increasing r. Hence, it was found that certain bodies can be ferromagnetic also in the amorphous state. In the amorphous state, however, the Curie point will be much lower. In conclusion, the author thanks A. P. Komar, Academician

Card 2/3

Quasiclassical Theory of Amorphous Ferromagnetics

81362 S/181/60/002/03/20/028 B006/B017

of the Ukrainskaya Akademiya nauk (Ukrainian Academy of Sciences) for having suggested the subject and for critical remarks. There are 4 references; 3 Soviet and 1 British.

ASSOCIATION: Fiziko-tekhnicheskiy institut AN SSSR Leningrad (Institute

of Physics and Technology, AS USSR, Leningrad)

SUBMITTED:

May 26, 1959

Card 3/3

Gicheria, A.D.

81956 \$/181/60/002/04/16/034 B002/B063

24,2100

AUTHOR:

Gubanov, A. I.

TITLE:

The Theory of Amorphous Conductors

PERIODICAL: Fizika tverdogo

Fizika tverdogo tela, 1960, Vol. 2, No. 4, pp. 651-655

TEXT: In contrast with the preceding papers of the author (Refs. 1 and 2), a rigorous proof is given in the present paper for the band structure of the energy spectrum of electrons in liquid or amorphous bodies. A new wave function describing the motion of an electron in an amorphous body is suggested in zero approximation:

$$\psi_{k} = \frac{1}{\sqrt{G}} u_{k} (\xi) \exp \left[i \sum_{\alpha=1}^{3} k_{\alpha} \lambda_{\alpha} \right]; \quad \lambda_{\alpha} \text{ are the coordinates of the}$$

electron in the system concerned, and \mathbf{k}_α are constants. The exponential factor yields the progressive motion of the electron, and the

Card 1/2

X

The Theory of Amorphous Conductors

81956 S/181/60/002/04/16/034 B002/B063

factor $u_{k}(\xi)$ which is periodic with respect to ξ gives the effect of the potential barrier. G. S. Gantsevich is mentioned. There are 3 Soviet references.

ASSOCIATION: Fiziko-tekhnicheskiy institut AN SSSR, Leningrad (Physicotechnical Institute of the AS USSR, Leningrad)

SUBMITTED:

July 14, 1959



Card 2/2

GubANOU, A.T.

82531

24.7100

S/181/60/002/007/003/042 B006/B070

AUTHORS:

Gubanov, A. I., Chevychelov, A. D.

TITLE:

Calculation of the Energy Spectrum of Strongly Anisotropic

Crystals 7

PERIODICAL:

Fizika tverdogo tela, 1960, Vol. 2, No. 7, pp. 1379-1389

TEXT: The purpose of the present work is to make a quantum-mechanical calculation of the energy spectrum of the electrons for two different models of an anisotropic crystal. In their theory of galvanomagnetic phenomena in metals I. M. Lifshits, M. Ye. Azbel', and M. I. Kaganov (Ref. 1) have assumed the existence of open isoenergetic surfaces in the k-space. The authors of the present paper have succeeded in establishing similar surfaces theoretically by investigating two models of a strongly anisotropic crystal: layer model and chain model. In a crystal with layer-model structure, for example, zinc, such planes appear since the valence forces are mainly acting between the atoms that lie in these planes; the binding between the planes is considerably weaker, and the

Card 1/4

82531

4.1.18.2.3.18.6. 主题投稿和如即往往的看法的的基础的重要的对比较级的基础的可能的工作的 1.4.1.1.1.1.1.1.1.1.1.1.1.1.

Calculation of the Energy Spectrum of Strongly Anisotropic Crystals

S/181/60/002/007/003/042 B006/B070

interatomic distance in the directions perpendicular to the layers are correspondingly large. In the direction of this crystallographic axis (z-direction), the motion of the electrons (in the single-electron crystal model) is assumed to be almost free, and in the other two directions strongly inhibited. Analogous assumptions are made for the chain model: A number of crystals have chain structure, that is, they have long spiral chains coiled round the hexagonal axis (selenium, tellurium), and the valence forces act along these chains; the binding between these chains is relatively weak, and the interatomic distances are large. The expressions for the potential of the electron is the same in both cases, with the difference only that in the first case the component $V(\overrightarrow{r})$ oc urs, which is a two-dimensional periodic function, and in the second case the function V(z) occurs which is periodic only in the z-direction. For both models, the spectrum is investigated by starting from the Schrödinger equations for these potentials. Later, the authors consider the shapes of the isoenergetic surfaces, first for the hexagonal lattice both for the layer and the chain models (Fig. 1). hexagonal lattice both for the layer and $\xi_2(p)$; $\vec{q} = (k_x, k_y)$ being the two-Fig. 2 shows the functions $\xi_1(\vec{q})$ and $\xi_2(p)$; $\vec{q} = (k_x, k_y)$

Card 2/4

8 2531

Calculation of the Energy Spectrum of Strongly Anisotropic Crystals

是一个人,我们就是我们的人,我们就是我们的人,我们们的人,我们们的人,我们们是不是一个人,我们就是我们的人,我们就是我们的人,我们就是我们的人,我们就是我们的人

S/181/60/002/007/003/042 B006/B070

dimensional k-vector and p its z-component. $\mathcal{E}_1(\vec{q})$ has a band shape, and for the first, third, and generally for odd bands $\epsilon_1(\vec{q})$ has a minimum in $\vec{q} = 0$ and a maximum in $\vec{q} = \vec{Q}$ (\vec{Q} is a vector that lies in the boundary for the first Brillouin zone). For even bands $\mathcal{E}_1(\vec{q})$ has a minimum in $\vec{q} = \vec{Q}$ and a maximum in $\vec{q} = 0$. Later, the isoenergetic curves for the planes formed by p and \vec{Q} are investigated. Fig. 3 shows some possible odd hands. It is seen that the topology of the isoenergetic curves. odd bands. It is seen that the topology of the isoenergetic surfaces for the two models are opposite to each other. For sufficiently large depths of potential both models have open surfaces of only one kind: corrugated cylinders in the case of layer structure, and corrugated planes in the case of chain structure. The theoretical results agree with the experimental ones. It is further shown that for lattices with the same crystallographic symmetry, but different chemical binding characteristics different laws hold for the dispersion of electrons. There are 3 figures and 4 Soviet references.

ASSOCIATION:

Fiziko-tekhnicheskiy institut AN SSSR Leningrad

(Institute of Physics and Technology of the AS USSR,

Leningrad)

Card 3/4

Calculation of the Energy Spectrum of Strongly Anisotropic Crystals
SUBMITTED: November 5, 1959

Card 4/4

84813

S/181/60/002/008/048/052/XX B006/B070

24.4500 AUTHORS:

Gubanov, A. I., Pushkarev, O. Ye.

TITLE:

Wave Functions of Valence Binding in Some Crystals

PERIODICAL: Fizika tverdogo tela, 1960, Vol. 2, No. 8, pp. 1776-1782

TEXT: The aim of the authors was to derive a more accurate (approximate) wave function of the valence electrons (localized along the valence lines) than was possible in a previous paper (Ref. 2). They use a variational method for this purpose. They consider the valence line of a diatomic molecule and choose a trial function which is analogous to that used in Ref. 3 for the hydrogen ion molecule. When the interatomic distance is increased, the trial function breaks into two atomic functions which, in the present case, are two sp functions of the neighboring atoms. To determine the trial function more conveniently, the atomic functions must be obtained first. The following approximate atomic function is assumed in §1: $R = ar[exp(-2r/r_k)-bexp(-4r/r_i)]$, where r_k is the covalent radius, and r_i is the ion radius, a and b are determined from the orthogonality

Card 1/3

84813

Wave Functions of Valence Binding in Some Crystals

S/181/60/002/008/048/052/XX B006/B070

and normalization conditions, respectively, $\int_0^{r_1} r^2 R dr = 0$, $\int_0^{\infty} r^2 R^2 dr = 1$.

b and a are explicitly given by (5) and (6). For a number of elements, numerical values of r_k , r_i , b, and a are given in a table. The atomic

functions for carbon, gallium, germanium, and arsenic are represented in Figs. 1-4. In the following, the authors determine the single-electron wave functions of the valence lines of diamond-type crystals by the method of variation. A numerical computation is made for germanium, and its results are given. The numerical example shows that in the approximation given here the wave function obtained by the variational method considered here is not much different from a linear combination of the atomic functions. In the present case, the maximum is shifted somewhat away from the atom, i.e., the electron density forms a cluster between the nearest atoms. There are 4 figures, 1 table, and 5 references: 2 Soviet, 1 US, 1 German, and 1 British.

ASSOCIATION: Fiziko-tekhnicheskiy institut AN SSSR, Leningrad (Institute of Physics and Technology AS USSR, Leningrad)

Card 2/3

Wave Functions of Valence Binding in S/181/60/002/008/048/052/XX B006/B070

SUBMITTED: December 22, 1959

84442

S/057/60/030/009/007/021 B019/B054

26.1410

AUTHORS:

Gubanov, A. I. and Lun'kin, Yu. P.

TITLE:

The Equations of Magnetoplasmadynamics

PERIODICAL:

Zhurnal tekhnicheskoy fiziki, 1960, Vol. 30, No. 9,

pp. 1046-1052

TEXT: It is usual in investigations of magnetic hydrodynamics to assume $\omega \tau \ll 1$ (ω is the Larmor frequency, τ the mean free time of charged particles). These assumptions are fulfilled in dense media and with weak magnetic fields in the plasma. Calculations in one- and two-liquid approximations were carried out for any $\omega \tau$. The introduction deals with a combination of the one-liquid approximation developed by Chapman et al. (Ref. 1) and the two-liquid theory developed by S. I. Braginskiy (Ref. 2). The relations (1) for pressure, temperature, the tensor of viscous tensions, and the heat flow are given. It is shown in the second part of the paper that expressions for the tensor of viscous tensions and the heat flow in one-liquid approximations can be easily obtained with the aid

Card 1/2

The Equations of Magnetoplasmadynamics

\$/057/60/030/009/007/021 B019/B054

of relations (1) from the formulas found by Braginskiy. When the resulting expressions are introduced into the motion- and energy equations of the one-liquid approximation, the equations of magnetoplasmadynamics are obtained after allowing for some transformations. The equations found are similar to those used in magnetohydrodynamics ($\omega\tau\ll$ 1), and the same methods as in magnetohydrodynamics can be used for their solution. The existence of additional terms in the equations leads, however, to new physical effects which do not follow from magnetohydrodynamics. There are 3 references: 2 Soviet and 1 British.

ASSOCIATION:

Fiziko-tekhnicheskiy institut AN SSSR, Leningrad

(Institute of Physics and Technology of the AS USSR,

Leningrad)

SUBMITTED:

July 2, 1959

Card 2/2

: .

84443

S/057/60/030/009/008/021 B019/B054

26. 1410

Gubanov, A. I. and Lun'kin, Yu. P.

TITLE:

AUTHORS:

The Cuettov Flow in Magnetoplasmadynamics ?

PERIODICAL:

Zhurnal tekhnicheskoy fiziki, 1960, Vol. 30, No. 9,

pp. 1053-1060

THE EXECUTION OF THE PROPERTY OF THE PROPERTY

TEXT: The authors investigated the flow between two parallel infinite plates, one of which is at rest while the other moves in its plane. It is assumed that a magnetic field $\overline{H_0}$ exists in various directions with respect to the plates and the motion \overline{u} of one plate. First, the case is studied where $\overline{H_0}$ is perpendicular to the plane of the plates. The authors show in a very long expansion that a flow originates here which is perpendicular to \overline{u} . This is called a specific effect of magnetoplasma-dynamics. Further, the case is studied where $\overline{H_0}$ lies in the plane of the plates. Here, the authors show that the magnetic field generates not only currents in the plasma but also currents running in the plates. The distribution of currents can only be given if the shape and dimension of

Card 1/2

84443

The Cuettov Flow in Magnetoplasmadynamics

\$/057/60/030/009/008/021 B019/B054

the plates and the position of the current-carrying conductors are given. Finally, the case is investigated where H_{O} has any direction. It appears that in all cases investigated, properties are found which do not follow from magnetohydrodynamics ($\omega \tau \ll 1$). $\overline{H_0}$ always shows an influence on flow and viscosity. There are 3 references: 1 Soviet and 2 US.

ASSOCIATION: Fiziko-tekhnicheskiy institut AN SSSR, Leningrad

(Institute of Physics and Technology of the AS USSR,

Leningrad)

SUBMITTED:

April 4, 1960

Card 2/2