

Ni-Pb

10

PbNiF₃

Syono Y. u. gp.

1968

Techn. Rept. ISSP

A, N337, 23pp., ill.

homomorphous.

(Ce. BaMnO₃)I

V 852

1897

Mosnier

Ann. chim. phys., 1897, 12, 374

ZnJ_2 , aq)

$2\text{ZnJ}_2 \cdot \text{PbJ}_2$) ΔHf

$2\text{CdJ}_2 \cdot \text{PbJ}_2$)

$2\text{FeJ}_2 \cdot \text{PbJ}_2$)

$2\text{NiJ} \cdot \text{PbJ}_2$)

$2\text{NiJ} \cdot \text{PbJ}_2 \cdot 3\text{H}_2\text{O}$) ΔHf

CoJ_2 , FeJ_2)

MnJ)

W, M

F

адм. н. / 05

ЕСТЬ в. н.

$\text{MnJ}_2 \cdot \text{PbJ}_2, \text{MnJ}_2, \text{PbJ}_2 \cdot 3\text{H}_2\text{O}, \text{CrJ}_2,$

$2\text{CrJ}_2 \cdot \text{PbJ}_2, 2\text{CrJ}_2 \cdot \text{PbJ}_2 \cdot 3\text{H}_2\text{O} \quad (\Delta \text{Hf})$

II Me PbF_6 , Me = Hg, Zn, Cd, Hg, Ni 1969
or. str.

VI 7023

IX
Homann R., Hoppe R.

Z. anorgan. und allgem. Chem., 1969, 368,
NS-6, 271-278

Neue Hexafluoroplumbate (IV). HgPbF_6 ,
 ZnPbF_6 , CdPbF_6 , HgPbF_6 und NiPbF_6 .

PZ, 65 551 (1970). ML 11

139-VI-6709

1969

Ni Pb Se
= 3 - 2 2

1-6969
T_m
Rep 8/9/69

48117g Lead selenide-nickel system. Myuller, N. N.;
Sotnikova, L. I. (Giprotvetmetobrabotka, Moscow, USSR).

Izv. Akad. Nauk SSSR, Neorg. Mater. 1969, 5(11), 1899-902
(Russ). The polythermal PbSe-Ni section of the ternary Pb-
Se-Ni system was constructed. The formation by a peritectic
reaction of the ternary compd. Ni₃Pb₂Se₂ m. 700° was estab-
lished. The unit cell parameters for this compd. were detd.
The formation of the eutectic between this compd. and PbSe
contg. 20 wt. % Ni and having a m.p. of 670° was established.
X-ray, microhardness, and DTA analyses were used for the
study. S. A. Mersol

C.A. 1970. 72. 10

30411.8755
TE, Ph

Ni - Pb 42530

1973

Candland C.T., Vanfleet H.B.

Effect of pressure on the interstitial
diffusion of nickel in lead to 50 kbar.
"Phys. Rev.B: Solid State", 1973, 7, N 2,
575-580

(англ.)

0851 БИК

839 841

ВИНИТИ

$Ni_3Pb_2S_2$

B9-XVI-3342

1976

85: 184994m Reaction of lead(II) sulfide with nickel. Zar-
garova, M. I.; Kakhramanov, K. Sh.; Roshal, R. M.; Guseinov,
G. G. (Spets. Konstr. Byuro, Inst. Fiz., Baku, USSR). *Izv.
Akad. Nauk SSSR, Neorg. Mater.* 1976, 12(9), 1557-9 (Russ).
The quasi-binary system $PbS-Ni$ was examd., and the phase
diagram constructed. A complex chem. interaction was obsd. in
this system with the formation according to a peritectic reaction
of $Ni_3Pb_2S_2$ (chem. analog of $Ni_3Pb_2Se_2$). Close phys. properties
of the $Ni_3Pb_2S_2$ and $Ni_3Pb_2Se_2$ are noted, and also a semimetallic
charcter of their elec. parameters. A thermal effect and a sharp
discontinuity in elec. properties were detected at $550 \pm 10^\circ$,
which is attributed to a structural transformation of $Ni_3Pb_2S_2$ in
the solid state.

(Tr)

C. A. 1976. 25 N24.

№3 Pb₂S₂

ВФ-ХVI-3342

1976

2 Б861. Взаимодействие PbS с Ni. Заргарова М. И., Кахраманов К. Ш., Рошаль Р. М., Гусейнов Г. Г. «Изв. АН СССР. Неорган. материалы», 1976, 12, № 9, 1557—1559

(Ттх) Показана квазибинарность системы PbS—Ni и построена диаграмма состояния. Выявлено сложное хим. взаимодействие в этой системе с образованием по перитектич. р-ции соединения $Ni_3Pb_2S_2$ — хим. аналога $Ni_3Pb_2Se_2$. Отмечена близость физ. св-в сульфоплюмбата и селеноплюмбата Ni, а также полуметаллич. характер их электрофиз. параметров. Обнаружены термич. эффект и резкий скачок электрич. св-в при $550 \pm 10^\circ$, объясняющиеся структурными превращениями $Ni_3Pb_2S_2$ в тв. состоянии.

Резюме

X1977N2

Ni-Pb

1986

22 Б3064. Термодинамические свойства свинца в растворах Ni—Pb. Thermodynamische Eigenschaften des Bleis in Ni—Pb-Lösungen. P o m l a n e k Т. «Z. Metallk.», 1986, 77, № 6, 388—392 (нем., рез. англ.)

Методом равновесного насыщения в вакууме с использованием в кач-ве станд. р-ра Cu—Pb измерены коэф. активности Pb в жидк. р-рах Ni—Pb для $x_{Pb} < 0,10$ при т-рах 1430, 1480 и 1510° С. Результаты описаны ур-нием $\ln \gamma_{Pb} = (7245 \pm 3,13)/T(1635 \pm 0,071) + [(42\,906 \pm 1853)/T + (18,54 \pm 0,80)] \cdot x_{Pb}$. С использованием лит. данных рассчитаны коэф. активности Pb и Ni в тв. р-рах Pb—Ni и вычислен состав монотектич. точки $x_{Pb} = 0,105$.

В. Ф. Байбуз

термод. св-ва
Pb в р-рах
Ni—Pb.

Х. 1986, 19, № 22

Ni-Pb

1987

108: 117032b The Ni-Pb (nickel-lead) system. Nash, P. (Ed.).
Coll. Eng., Illinois Inst. Technol., Chicago, IL 60616 USA). In
Alloy Phase Diagrams 1987, 8(3), 264-8 (Eng). The Ni-Pb
diagram was assessed. Crystal structures and lattice parameters
Ni, Pb, and metastable NiPb and activity coeffs. for Ni in liq. Pb
given. Extensive Ni-Pb liquidus data was included.

pass. group.

C.A. 1988, 108, N14

1999

F: Ni-Pb-Sn

P: 1

131:173327 Thermodynamic modeling of the nickel-lead-tin system. Ghosh, G (Department of Materials Science and Engineering, Robert R. McCormick Sch Engineering and Applied Science, Northwestern University, Evanston, IL 60 3108, USA). Metall. Mater. Trans. A, 30A(6), 1481-1494 (English) 1999 A set of self-consistent thermodyn. model parameters is presented to describe the phase equil. in the Ni-Pb and Ni-Sn systems. Sublattice descriptions are used for thermodyn.

modeling of the η -Ni₃Sn, λ - η -Ni₃Sn₂, λ -Ni₃Sn₂, and Ni₃Sn₄ phases. A three-sublattice and sublattice model are used to describe the molar Gibbs energies of η -Ni and λ -Ni₃Sn₂, resp., and also to describe the second-order phase transition from η -Ni₃Sn₂ to λ -Ni₃Sn₂. In a majority of the cases the agreement between the exptl. data and calcd. values was satisfactory. The exptl. Ni-Pb-Sn ternary phase diagram is not known, several isotherma sections are calcd. based on thermodyn. principles.

Ni-Pb

2001

134: 153192z Thermodynamic assessment of the Cu-Ni-Pb system. Wang, Cui Ping; Liu, Xing Jun; Ohnuma, Ikuo; Kainuma, Ryo-suke; Ishida, Kiyohito. (Department of materials Science, Graduate School of Engineering, Tohoku University, Sendai, Japan 980-8579). *CALPHAD: Comput. Coupling Phase Diagrams Thermochem.* 2000 (Pub. 2001), 24(2), 149-167 (Eng), Elsevier Science Ltd. A thermodyn. assessment of the binary systems Ni-Pb, Cu-Pb and the ternary system Cu-Ni-Pb was carried out by the CALPHAD method. The Gibbs free energies of both the liq. and solid soln. phases were described by sub-regular soln. models. A set of parameters describing the Gibbs energies of the different phases in this ternary system was optimized using exptl. phase diagram and thermodyn. information. Good agreement between the calcd. phase diagrams and exptl. data was obtained in the binary Ni-Pb, Cu-Pb and ternary Cu-Ni-Pb systems.

(al., ml. g-pr
DF, pay. gray)

(+2) 

C.A. 2001, 134, N11