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## Effect of Boron Additive on the Formation Ferrous Phases in Alloys Al - Fe - Si

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Abstract. The article presents an analysis of the diagram of the state of the AI - Fe - Si system using the Thermo-Calc program: polythermal sections were constructed and analyzed in order to determine the concentration limits of the appearance of primary crystals of Fe-containing phases, the mass fractions of phases at certain temperatures were calculated. A search was performed for optimal concentrations and temperatures at which the formation of a highly symmetric  $AI_8Fe_2Si$  phase capable of perceiving plastic deformation takes place. The Al-30Fe-9Si-1B system was also investigated. Based on the calculation of the parameters of the primary crystallization of iron in AI - Fe - Si alloys, the possibility of evaluating the effectiveness of modifying this alloy with boron is shown. The boron additive was considered as a modifier contributing to the grinding of aluminium grains and the formation of a wide area of existence of the  $AI_8Fe_2Si$  phase. The concentration and temperature ranges at which the maximum number of phases  $AI_8Fe_2Si$  and  $AIB_2$  can be achieved have been determined. The minimum amount of the  $AI_9Fe_2Si_2$  phase has been calculated, which is undesirable because it leads to a decrease in the mechanical properties of the alloy. From a practical point of view, the complex of proposed studies will allow to scientifically substantiate and develop new technologies for the production and processing of semi-finished products and finished products from aluminium-based alloys and functional materials with a given structure.

Keywords: aluminium, crystallization, boron, phase, polythermic section, modification.

#### Introduction

Aluminum is one of the key export products of Kazakhstan. The country holds a leading position in Central Asia for its production. However, there are problems associated with this industry (high world market value, high energy costs) [1]. Nevertheless, the prospects for the development of the aluminum industry in Kazakhstan remain quite high. Kazakh companies are actively working to modernize their production facilities and introduce new technologies. This increases production efficiency and reduces energy costs [1].

In the conditions of saving material and energy resources, the issue of recycling aluminum alloys becomes especially relevant, which inevitably leads to an increase in the content of such poorly removed impurities as iron and silicon [2-5]. The increased content of the latter limits the possibility of obtaining such grades of aluminum alloys that are capable of deep drawing, and some operational properties are also reduced. At the same time, recycled aluminum alloys can be very effectively used as raw materials for the production of metal-ceramic materials based on aluminum with an increased content of both iron and silicon. Such metal-ceramic materials have high hardness values and can be used to manufacture products operating under high wear conditions. A distinctive feature of the alloy considered in this paper is the possibility of forming the  $\alpha$ -phase, which is characterized by high symmetry of the crystal lattice, which predicts its susceptibility to plastic deformation to a greater extent compared to the  $\beta$  and  $\theta$  phases, which have monoclinic crystal lattices [6-10].

Considering the increased content of alloying elements in the alloy, significantly exceeding the solubility limits in aluminum, it is quite labor-intensive to obtain such materials by traditional melting. For this purpose, the use of additive manufacturing methods can solve the problem of obtaining metal-ceramic materials.

Often, to ensure the formation of a favorable microstructure, small additives of modifiers are used in aluminum alloys, among which the most common are Ti, Zr, V, TiC, TiB<sub>2</sub>, Na, Sr, Sb, Ba [11-15].

Boron is also a well-known modifier of aluminum alloys, which helps to grind the primary crystallizing phases in aluminum alloys, improving the quality of mechanical processing. Aluminum-matrix composite materials alloyed with boron have low specific density, strength, corrosion resistance, good thermal conductivity, and the ability to absorb thermal neutrons. At the same time, there is insufficient literary data showing the effect of boron on the phase composition of aluminum alloys rich in both iron and silicon, and they are mainly fragmentary.

One of the promising areas for producing high-quality aluminum alloys is the additive method of composite material synthesis. The aim of the work was a fundamental study of the Al-Fe-Si system. Search for optimal concentrations and temperatures at which a highly symmetric phase of  $Al_8Fe_2Si$  is formed, capable of receiving plastic deformation. Al-30Fe-9Si-1B was also investigated. The boron additive was seen as a modifier, contributing to the grinding of aluminum grains and the formation of a wide range of phase  $Al_8Fe_2Si$ . Aluminum alloys, in which boron is used as a modifier (up to 0.04% at melt temperature 690-710°C immediately before casting) and grain shredder are used in the electrical industry. Boron is also introduced to increase electrical conductivity and as a refining additive in aluminum alloys, which are used in the nuclear power industry in the amount of 0095 - 0.1% [16].

### 1. Research methodology

At the initial stage, the Al-Fe-Si system from the side of the iron corner was considered. This analysis was carried out using the modern computer program Thermo-Calc.

The Thermo-Calc computer product is a program for calculating phase equilibrium, the basic principle of operation: the algorithm of global minimization of Gibbs energy of multicomponent systems. It is also possible to calculate thermodynamic properties of phases (Gibbs energy, enthalpy, etc.), metastable equilibrium. The main characteristic of this program is its modularity, extensibility, replenishment of the databank of various elements for different systems (metal, salt, oxide, aqueous solutions, etc.). This program also allows the calculation of phase diagrams of multicomponent systems (construction of polythermic and isothermal sections, phase composition calculation, calculation of cooling curves), which was done in this work.

Using the Thermo-Calc program (version TCW8, database TTAL8.2) the aluminum-based system was analyzed to determine the concentration limits of the appearance of primary Fe-containing phase.

The Al-Fe-Si aluminum corner state diagram (Figure 1) is considered as an example. In equilibrium with (Al), in addition to the phases from double systems, there can be two triple connections:  $Al_8Fe2Si (\alpha)$  and  $Al_9Fe_2Si_2 (\beta)$ . In this system there are three three-phase reactions involving (Al): (Al):  $L + Al_9Fe_2Si_2 \rightarrow L + Al_9Fe_2Si_2 + (Al)$ ;  $L + Al_9Fe_2Si_2 + (Al) \rightarrow Al_8Fe_2Si + Al_9Fe_2Si_2 + (Al)$ ;  $Al_9Fe_2Si_2 + (Al) \rightarrow Al_9Fe_2Si_2 + (Si) + (Al)$ , three three-phase reactions involving liquid phase L and phase  $\theta$ :  $L + \theta + \tau_2 \rightarrow L + \theta + Al_8Fe_2Si$ ;  $L + \theta + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si$ ;  $L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si$ ;  $Al_9Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si$ ;  $Al_9Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si$ ;  $Al_9Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si$ ;  $Al_9Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si$ ;  $Al_9Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si$ ;  $Al_9Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si$ ;  $Al_9Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si$ ;  $Al_9Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + Al_8Fe_2Si$ ;  $Al_9Fe_2Si \rightarrow L + Al_8Fe_2Si \rightarrow L + A$ 



Region number	Phase region
1	L
2	$L + \theta$
3	$L + \theta + \tau_{11}$
4	$L+\theta+\tau_2$
5	$L + \theta + Al_8Fe_2Si$
6	$L + Al_8Fe_2Si$
11	$L + Al_8Fe_2Si + Al_9Fe_2Si_2$
12	$L + Al_9Fe_2Si_2$
13	$L + Al_9Fe_2Si_2 + (Al)$
16	$Al_8Fe_2Si + Al_9Fe_2Si_2 + (Al)$
15	$Al_9Fe_2Si_2 + (Al)$
14	$Al_9Fe_2Si_2 + (Si) + (Al)$
17	$Al_8Fe_2Si + Al_9Fe_2Si_2$
18	$\theta + Al_9Fe_2Si_2$
7	$\theta + Al_8Fe_2Si$
8	$\theta + Al_8Fe_2Si + \tau_2$
9	$\theta$ + $\tau_{11}$ + $\tau_2$
10	$\theta + \tau_2$
19	$\theta + Al_9Fe_2Si_2 + \tau_2$
20	$\theta$ + Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub> + (Si)
21	$\theta + A \overline{l_9 F e_2 S i_2 + \tau_8}$
22	$\theta + Al_9Fe_2Si_2$

Fig. 1. - Polythermic section of the system AI - 32Fe - 9Si on the side of the iron corner

Primary crystals  $\theta$  - phase, tend to grow along the orientation [010], forming coarse needles, scales and plate shapes in Al-Fe hypereutectic alloys, which seriously impairs the mechanical properties of alloys [17], [18].

The Al<sub>8</sub>Fe<sub>2</sub>Si phase has a rather narrow crystallization interval and is suppressed by Al<sub>9</sub>Fe<sub>2</sub>Si<sub>2</sub> phase formation, which is a very undesirable process, as the presence of Al<sub>9</sub>Fe<sub>2</sub>Si<sub>2</sub> phase in the structure causes deterioration of the mechanical properties of the material [19], [20].

Phase	%	Content					
	mass	of the components, %					
		Al	Si				
t = 1072  °C (1  region)							
L	100	81,20	9,75	9,00			
	t = 925 °	°C (2 regio	on)				
L	84,5	69,71	19,52	10,63			
θ	15,5	59,51	39,60	0,82			
	t = 928 °	°C (3 regio	on)				
L	31,27	53,21	29,33	17,38			
θ	28,63	58,77	39,71	1,51			
τ <sub>11</sub>	40,09	51,11	40,69	8,18			
	t = 831 °	°C (4 regio	on)				
L	24,93	71,37	14,99	13,63			
θ	40,25	59,17	39,39	1,42			
τ2	34,80	51,05	33,89	15,05			
	t = 730 °	C (5 regio	on)	•			
L	8,88	84,45	6,63	8,90			
θ	12,79	59,60	39,21	1,17			
Al <sub>8</sub> Fe <sub>2</sub> Si	78,23	57,11	32,55	10,33			
	t = 671  °C (6  region)						
L	70,61	87,71	3,75	8,53			
Al <sub>8</sub> Fe <sub>2</sub> Si	29,38	57,50	32,55	9,94			
t = 665  °C (7 region)							
θ	6,08	59,64	39,15	1,19			
Al <sub>8</sub> Fe <sub>2</sub> Si	93,91	57,91	32,56	9,52			
t = 648  °C (8  region)							
θ	23,96	59,34	39,16	1,49			
Al <sub>8</sub> Fe <sub>2</sub> Si	43,25	57,67	32,55	9,76			
$\tau_2$	32,77	52,44	33,91	13,64			
	t = 828 °	°C (9 regio	on)				
θ	41,76	58,53	39,57	1,89			
$\tau_{11}$	16,63	51,11	40,69	8,18			
$\tau_2$	41,60	49,44	33,87	16,68			
t = 571 °C (10 region)							
θ	51,97	58,00	39,23	2,75			
τ2	48,02	50,16	33,88	15,94			
	t = 618 °	C (11 regi	on)				
L	61,99	90,20	1,97	7,81			
Al <sub>8</sub> Fe <sub>2</sub> Si	28,91	58,07	32,56	9,36			
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	9,08	58,28	26,90	14,81			

## Table 1. Phase composition of the system AI-32Fe-9Si

Phase	%	Content				
	mass	of the components, %				
		Al	Fe	Si		
	t = 624 °C	C (12 regio	n)			
L	97,37	89,03	2,16	8,79		
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	2,62	57,64	26,89	15,45		
	t = 589 °C	C (13 regio	n)			
L	25,09	88,17	1,12	10,7		
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	34,68	57,46	26,89	15,64		
(Al)	40,22	98,76	0,00	1,22		
	t = 379 °C	C (14 regio	n)			
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	52,15	63,49	26,95	9,54		
(Si)	3,74	0,00	0,00	100		
(Al)	44,10	99.79	0,00	0,20		
	$t = 491 \circ 0$	C (15 regio	n)	,		
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	81.11	62.32	26.94	10.73		
(Al)	18,88	99,68	0,00	0.31		
	$t = 550 \circ 0$	C (16 regio	n)	,		
Al <sub>8</sub> Fe <sub>2</sub> Si	30,58	60,35	32,59	7,05		
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	59.49	61.78	26.94	11.27		
(Al)	9.91	99.67	0.31	0.00		
t = 494  °C (17  region)						
Al <sub>8</sub> Fe <sub>2</sub> Si	49.72	60.32	32.59	7.08		
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	50.27	62.09	26.94	10.96		
t = 362  °C (18  region)						
θ	6,38	57,98	38,99	3,02		
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	93.61	63.60	26.96	9.43		
	$t = 367 \circ 0$	C (19 regio	n)	,,		
θ	33.68	56.72	38.96	4.31		
AleFeeSie	46.53	63.16	26.95	9.88		
T1191 02012	19.78	50.88	33.89	12 55		
62	$t = 229 \circ 0$	7 (20 regio	n)	12,55		
0	(-22)	53 03	38.82	8 1/		
AlaFasSia	50,00 60,01	64.00	26.06	0.02		
A191 62512	09,01	04,00	20,90	9,05		
(31)	0,51	$\frac{0,00}{1}$	0,00	100		
0	1 - 120	51.64	n) 20.75	0.50		
H Ea Si	0,70	51,04	30,13	9,39		
A19F e2S12	93,18	04,11	20,90	8,92		
τ8	0,11	20,10	50,26	29,62		
	t = 93  °C	(22 region	1)	0.0 <b>-</b>		
θ	49,58	52,15	38,77	9,07		
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	50,41	64,11	26,96	8,92		

To evaluate the morphology of the obtained phases, an alloy of the Al-Fe-Si system with a basic composition was synthesized. The technological conditions of the synthesis are presented in detail in [21, 22]. For metallographic analysis, microsections were prepared using a standard technique, using cutting, grinding, and polishing operations. The microstructure was studied immediately after synthesis, without additional processing in the "as prepared" state. The synthesized ingot has two main regions: the intermetallic region, which occupies the bulk of the material, and the eutectic region (Figure 2).



Fig. 2. - Microstructure of the alloy system AI-Fe-Si

Metallographic analysis shows that the structure of the material predominantly contains a combination of  $\beta/\theta$  phases, with the growth of the  $\beta$  phase occurring on the primary crystals of the  $\theta$  phase, which is in good agreement with the results of computer modeling for region 22 (Figure 1). At the same time, thin sections of the FCC solid solution of silicon and iron in aluminum were detected, formed as products of decomposition and phase transformation as a result of nonequilibrium cooling conditions during synthesis. The obtained results demonstrate the validity of using this approach to modeling for the purpose of predicting the phase composition of the alloy during its modification.

In this connection, the effect of boron on the formation and formation of the  $Al_8Fe_2Si$  phase was further considered. As an example, a diagram of the state of Al-Fe-Si-B (Figure 3) was constructed.



Fig. 3. – Polythermic section of the system AI-30Fe-9Si-1B

In this system, with the participation of the  $\theta$  phase, there are four three-phase reactions with the formation of Al\_8Fe\_2Si and AlB\_2: L +  $\theta$  + AlB<sub>2</sub> +  $\tau_2 \rightarrow$  L +  $\theta$  + Al\_8Fe\_2Si + AlB<sub>2</sub>; L +  $\theta$  + Al\_8Fe\_2Si + AlB<sub>2</sub>  $\rightarrow$  L + Al\_8Fe\_2Si + AlB<sub>2</sub>; L + Al\_8Fe\_2Si + AlB<sub>2</sub>  $\rightarrow$   $\theta$  + Al\_8Fe\_2Si + AlB<sub>2</sub>. There is one four-phase reaction involving (Al): L + Al\_8Fe\_2Si + Al\_9Fe\_2Si\_2 + AlB<sub>2</sub>  $\rightarrow$   $\theta$  + Al\_8Fe\_2Si + AlB<sub>2</sub> + (Al). The Al\_8Fe\_2Si phase has three three-phase reactions: Al\_8Fe\_2Si + Al\_9Fe\_2Si\_2 + AlB<sub>2</sub>  $\rightarrow$   $\theta$  + Al\_8Fe\_2Si + AlB<sub>2</sub> +  $\tau_2$ ;  $\theta$  + Al\_8Fe\_2Si + AlB<sub>2</sub> +  $\tau_2 \rightarrow$  Al\_8Fe\_2Si + AlB<sub>2</sub>. The formulations of each phase at a certain temperature are given in table 2. As seen from the reactions, the separation of phase Al\_8Fe\_2Si and AlB<sub>2</sub> goes through the entire temperature range.

Table 2. Phase co	mposition of the syster	n Al-30Fe-9Si-1B
	1 1	

Phase	%	Content				
	mass	0	of the components, %			
		Al	Fe	Si	В	
	t =	1236 °C (	1 region)			
L	100	64,21	25,78	9,00	1,00	
	t =	1236 °C (	2 region)			
L	100	64,21	25,78	9,00	1,00	
	t =	818 °C (.	3 region)			
L	44,65	72,57	14,10	13,12	1,89	
θ	33,80	59,21	39,37	14,1	0,00	
AlB <sub>2</sub>	2,85	17,21	0,00	0,00	82,78	
τ2	18,68	51,20	33,9	14,89	0,00	
t = 730  °C (4  region)						
L	23,54	84,25	6,73	8,93	0,07	
θ	8,13	59,6	39,21	1,18	0,00	
Al <sub>8</sub> Fe <sub>2</sub> Si	64,09	57,1	32,55	10,34	0,00	
AlB <sub>2</sub>	4,22	55,51	0,00	0,00	4,48	
t = 685  °C (5  region)						

Phase	%	Content				
	mass	0	f the com	ponents, '	%	
		Al	Fe	Si	В	
	t =	568 °C (7	7 region)			
Al <sub>8</sub> Fe <sub>2</sub> Si	30,53	59,72	32,58	7,69	0,00	
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	51,75	60,97	26,93	12,08	0,00	
AlB <sub>2</sub>	4,24	55,51	0,00	0,00	44,48	
(Al)	13,46	99,56	0,01	0,41	0,00	
	t =	568 °C (8	8 region)			
Al <sub>8</sub> Fe <sub>2</sub> Si	79,22	58,98	32,57	8,44	0,00	
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	16,38	60,04	26,92	13,02	0,00	
AlB <sub>2</sub>	4,39	55,51	0,00	0,00	44,48	
t = 677  °C (9  region)						
θ	4,64	59,66	39,16	1,17	0,00	
Al <sub>8</sub> Fe <sub>2</sub> Si	90,92	57,8	32,55	9,63	0,00	
AlB <sub>2</sub>	4,42	55,51	0,00	0,00	44,48	
	$t = 628 \ ^{\circ}C \ (10 \ region)$					
θ	22,45	59,31	39,15	1,54	0,00	

L	61,31	86,94	4,37	8,63	0,04		
Al <sub>8</sub> Fe <sub>2</sub> Si	34,76	57,39	32,55	10,05	0,00		
AlB <sub>2</sub>	3,91	55,51	0,00	0,00	44,48		
t = 625  °C (6  region)							
L	60,90	89,26	2,23	8,47	0,02		
Al <sub>8</sub> Fe <sub>2</sub> Si	31,08	57,80	32,55	9,63	0,00		
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	4,06	57,79	26,89	15,31	0,00		
AlB <sub>2</sub>	3,93	55,51	0,00	0,00	44,48		

Al <sub>8</sub> Fe <sub>2</sub> Si	38,90	57,9	32,56	9,53	0,00		
AlB <sub>12</sub>	3,23	17,21	0,00	0,00	82,78		
$\tau_2$	35,40	52,49	33,91	13,58	0,00		
t = 520 °C (11 region)							
Al <sub>8</sub> Fe <sub>2</sub> Si	68,86	59,46	32,58	7,95	0,00		
Al <sub>9</sub> Fe <sub>2</sub> Si <sub>2</sub>	27,97	61,02	26,93	12,04	0,00		
AlB <sub>12</sub>	3,16	17,21	0,00	0,00	82,78		

### 2. Results and Discussing

One of the main criteria for producing alloys based on aluminium is the process of modification (grinding of grains of cast aluminium with the introduction of special additives). The process of grinding aluminum grain at crystallization is the result of special interaction of transition metals with aluminium [23], [24]. The ideal modifier is a particle that satisfies the following requirements: it should be effectively crushed grain at a minimum concentration; in the melt is thermally stable and dispersed; it has a minimum structural difference with the lattice of the modifying alloy; do not lose their modifying properties when melting. None of the modifiers currently known have a complete set of these properties [16], [25], [26]. One such modifier in aluminum alloys is boron, which is a transition metal. It has good refining effect on Al-Si alloys [27], [28].

Analyzing the polythermic section of the system Al - 32Fe - 9Si on the side of the iron angle sections, we can conclude that the phase of interest to us Al<sub>8</sub>Fe<sub>2</sub>Si has a fairly wide crystallization interval at a temperature from 730 C to 490 °C, but its formation is suppressed by the formation of phase Al<sub>9</sub>Fe<sub>2</sub>Si<sub>2</sub>. For example, at 494 °C (table 2, 17 region), the phase region consists of Al<sub>8</sub>Fe<sub>2</sub>Si (49.72%) and Al<sub>9</sub>Fe<sub>2</sub>Si<sub>2</sub> (50.27%). However, with the drop in temperature, a phase transition is underway with the disappearance of the Al<sub>8</sub>Fe<sub>2</sub>Si phase, which is completely transitioned to the Al<sub>9</sub>Fe<sub>2</sub>Si<sub>2</sub> phase.

The Al<sub>9</sub>Fe<sub>2</sub>Si<sub>2</sub> phase is known to be the most harmful, reducing the mechanical properties of aluminum alloys. In particular, the plate or needle phase of Al<sub>9</sub>Fe<sub>2</sub>Si<sub>2</sub> not only cleaves the matrix, but also contributes to the formation of casting defects such as pores and shrink pores. These phases easily cause stress concentration and crack propagation, thus becoming the weakest sites in the alloys Al-Si [29], [30], [31]. Therefore, a section of the system Al-30Fe-9Si-1B was built and the main optimal phase compositions were determined.

Boron is an effective grain shredder in Al-Si alloys, but the mechanism responsible for the effect of boron on the formation and growth of phases rich in iron is still under discussion [31]. The mechanism of primary grinding of boron grains in the Al matrix in the alloy is based on the formation of particles AlB<sub>2</sub> as a result of eutectic reaction (table 3, region 3). And the formation of this phase has a fairly wide temperature crystallization interval (818 - 520 °C). It can also be concluded that this phase contributes to the formation of the Al<sub>8</sub>Fe<sub>2</sub>Si phase with a fairly good phase composition (min 31.08%, max 90.92%). Phase AlB<sub>2</sub> is crystallized by eutectic reaction at 818 C (table 4, area 3). Since the AlB<sub>2</sub> and Al<sub>8</sub>Fe<sub>2</sub>Si phases have different crystallographic lattices, it can be assumed that the AlB<sub>2</sub> phase is the catalyst for the onset of phase Al<sub>8</sub>Fe<sub>2</sub>Si [32], [33], [34].

#### Conclusions

Quantitative analysis of the phase diagram Al - Fe - Si as the basis of the high-strength alloy was carried out using the program Thermo-Calc (version TCW8, database TTAL8.2). The standard alloy Al - 32Fe - 9Si was considered, built a polythermic section, calculated the temperature of phase transformations. The phase composition obtained during modeling was compared with the real phase composition, and the microstructure of the alloy was studied. Optimum concentrations of iron and silicon were determined for  $Al_8Fe_2Si$  phase crystallization. The mass fractions of each phase are calculated at certain temperatures.

A polythermic cut Al-30Fe-9Si-1B was also calculated. The concentration and temperature regions at which the maximum number of  $Al_8Fe_2Si$  and  $AlB_2$  phases can be achieved have been determined. The minimum number of phase  $Al_9Fe_2Si_2$  is calculated, which is undesirable because it leads to a decrease in the mechanical properties of the alloy.

The data presented in this article are the initial stage in carrying out research on the development of technology for the synthesis of composite ceramic materials of the AlxFeySi system using an additive method to improve the quality of the structure of materials based on them. In the future, a complete thermodynamic analysis will be carried out to select the optimal concentrations of alloying elements in order to develop a rational mode of heat treatment of the alloy under study.

In the future, using the above described thermodynamic calculations to identify and modify phases rich in iron in Al-Fe-Si alloys, it is planned to experiment with the addition of boron by the method of additive synthesis.

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