atoms N1 from two independent heterocyclic cations participate in the bifurcatic hydrogen bonds N–H...O, O'. Three dimensional spatial package of crystals of the compound I is more dense (ρ_c =1.547 g.cm 3) than lamellar package of 8-aminoquinolinium bis(citrato)borate terahydrate crystals (ρ_c =1.453 g.cm 3) [1]. Crystals I are triclinic, space group P $\bar{1}$: a=10.6678(3) Å, b=14.2920(3) Å, c=16.4891(4) Å, α =78.106(1)°, β =75.368(1)°, γ =85.919(1)°, V=2379.8(1) Å 3 , Z=4, R=0.054, wR2=0.133 for 10562 independent reflections with R(int)=0.0266.

[1] Zviedre I., Belyakov S., Latvian Journ. Chem., 2006, 3, 233

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Isothermal section at 800°C of the Gd-Fe-Ge ternary system R. Ben Hassen^a, M. Jemmali^{a,b}, S. Walha^a, O. Tougait^b, H. Noël^b, ^aUnité de chimie des matériaux, Université Tunis El Manar, Tunis, Tunisia. ^bSciences chimiques de Rennes UMR 6226, CNRSuniversité de Rennes 1, Avenue du Général Leclerc, F-35042 Rennes, France.

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The investigation of the Gd-Fe-Ge system is part of an ongoing research project with the aim to clarify the phase equilibrium in ternary system of Gd and iron with a p-

block element. This ternary system has been partially investigated; the phase diagrams established at 800°C show two compounds. A: $Gd_1Fe_2Ge_2$ (a = 3.995, c = 10.46 Å, I4/mmm, $\hat{ThCr_2Si_2}$ structure type) [1], **B**: $Gd_1Fe_6Ge_6$ (a = 5.128, c = 4.076 Å, P6/m 2/m2/m, YCo₆Ge₆ structure type) [2]. We present here our experimental results on the Gd-Fe-Ge ternary system, studied in the whole concentration range at an isotherm of 800° C. All the samples were prepared by arc-melting the elemental components, followed by heat-treatment of one week. The phases in alloys were determined by electron-probe microanalysis and examined by X-ray powder diffraction analysis (X-ray diffractometer with $CoK\alpha$ radiation with iron filters) in order to determine the phase compositions and the equilibrium lines within the ternary system. In addition to the known phases, a new ternary phase, Gd₃Fe₁Ge₆, has been found and its crystal structure was refined from powder. Gd₃Fe₁Ge₆, this new ternary phase crystallizes in the orthorhombic space group Cmcm (n° 63) with the lattice parameters a = 4.151, b = 16.062 and c= 4.0239 Å, structure type CeNiSi₂. This phase shows a significant homogeneity range which extends between the compositions $GdFe_{1-x}Ge_2$ (0.55 $\leq x \leq 0.75$).

[1] Venurini G., Welter R., Malaman B., Journal of alloys and compounds, 1992, 185, 99.

[2] Duong N. P., Bruck E., Brommer P. E., Klaasse J. C. P., De Boer F. R., Buschow K. H. J., Journal of Magnetism and Magnetic Materials, 2002, 242-245, 813

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