

PARAMETERIZATION OF SIMPLE BRANCHING SOLUTIONS
OF COMPLETE RANK AND ITERATIONS
IN THE NONLINEAR ANALYSIS

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Dedicated to 70-th anniversary of Professor V.A. Trenogin's birth

Introduction

The role of the geometry of planes, diagrams, and Newton's polyhedrons is well-known (see [1], [2]) in the asymptotical analysis of the theory of branching (see [1]–[5] and other papers). In the construction of branching solutions by iteration methods, both explicit and implicit parameterizations were used (see [6]–[12]), and in complex cases (see [10], [13]–[15]) — the geometrical constructions indicated above. In [6], [7], and [9]–[12], the rigid constraints ensuring access to the principal part of the equation of branching (e. b.) already at the first iteration were of a decisive importance for the convergence of methods. In particular cases a method for construction of the initial approximation and the choice of parameterization of the sought-for branch were indicated. In [6], [7], [9], [11], and [12] the uniformization parameter was chosen directly from the form of the Newton diagram of the coefficients of the projection QF . Therefore on each iteration it was necessary to solve one linear equation. Iteration methods for more complex situations in [6]–[12] were not developed, because the ways for the parameter of uniformization of branches to be chosen and iteration process to be organized in general case were not clear.

In [13], [14] it was shown that both the essential expansion of the class of equations and creation of more flexible algorithms are possible by means of application of N -step methods, in which on each iteration N linear equations are solved and different parameters of uniformization are applied in accordance with [2]. In [13] an N -step iteration method was suggested which converges in a neighborhood of the branching point under constraints which are weaker than those of other methods. The method possesses the advantages of both explicit and implicit parameterization (see [12]), since it gives us a certain freedom in the choice of the uniformization parameter. However, the iteration scheme suggested on the basis of this method in [13] (see § 2 there), requires, in the general case, a preliminary construction of an element \hat{x}^0 from the subspace $E_1^{\infty-n}$, which is used at the initial approximation. A method for construction of the element \hat{x}^0 was given in [13] (see lemma 2 there). Nevertheless, the preliminary computation of \hat{x}^0 makes both the computations and the proper iteration scheme more complicated.

The objective of this article is not a simple modification and expansion of the possibilities of iteration methods in [11], [12], [16], but also deduction of iteration formulas convenient for the theory of approximate methods and its applications. A more convenient iteration scheme with simplified choice of the initial approximation is suggested.

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