Comparing inclusion techniques on chemical engineering problems

Ali Baharev, Endre Rév
Department of Chemical and Environmental Process Engineering
Budapest University of Technology and Economics
Budafoki ut 6-8.
1111 Budapest, HUNGARY
ali.baharev@gmail.com

Abstract

Solving general nonlinear systems of equations and/or finding the global optimum of nonconvex functions constitute an important part of the everyday practice in chemical engineering. Standard methods cannot provide theoretical guarantee for convergence to a solution, cannot find multiple solutions, and cannot prove non-existence of solutions. This is the main motive to apply interval methods. Interval analysis has been applied to a wide variety of problems in chemical engineering, e.g. [1]. Similarly impressive results can be achieved with αBB , a fairly general global optimization method [2, 3].

Computing steady states of multistage separation processes requires solving large-scale nonlinear systems of equations. Despite of the outstanding results referred above, computation of these problems with interval arithmetic have not yet been considered in the literature, according to the authors' best knowledge. The only globally convergent methods used in this area seem to be the homotopy-continuation methods [4]. Leading edge software packages, such as [5], are also available to find zeros of nonlinear systems. Unfortunately, finding all solutions can be guaranteed only in special cases e.g. polynomial systems with no constraints [6].

The authors aim to compute steady states of homogeneous and heterogeneous azeotropic distillation columns with interval methods, keeping the algorithm as problem independent as possible. The results achieved so far are presented here.

Numerical evidence published in the literature, e.g. [7, 8], seem to indicate superiority of the linear interval approximation (LIA, $\mathbf{L}(x) = Ax + \mathbf{b}$, A is a real matrix), proposed by Kolev in a number of publications e.g. [9], compared to the traditional interval linear approximation (ILA, $\mathbf{L}(x) = \mathbf{A}(x-z) + f(z)$, \mathbf{A} is an interval matrix) such as the interval Newton method. LIA has the following advantages over ILA when applied to root-finding. (i) The solution set of the LIA has a much simpler form, the hull solution is straightforward: $\mathbf{X} \cap -A^{-1}\mathbf{b}$. (ii) Linear programming is directly applicable to prune the current box. The automatic

computation of LIA is possible with affine arithmetic [10] which in turn (iii) automatically keeps track of correlation between the computed partial results yielding tighter enclosures. There is no significant difference in the computation time per iteration between LIA and ILA.

In [11] LIA and ILA are compared as linearization techniques applying them to chemical engineering problems of real complexity. The examples considered are highly structured and are full of dependency. LIA outperforms the traditional 'textbook' interval Newton algorithm (IN/GS) by an order of magnitude in the case of the studied examples. Note that state-of-the-art variants of the interval Newton methods, e.g. [12, 13], also outperform the IN/GS used for comparison. Linear programming may be preferable as pruning technique for LIA because of its robustness. Considering the conclusions of [11], the C++ class has been re-implemented, and the LP pruning method has been revised. The improvement is significant; real life medium-scale problems are successfully solved. Some of the problems used for comparison are suitable for benchmarks, they will be contributed soon.

References:

- [1] Y. Lin, C. R. Gwaltney, M. A. Stadtherr; Reliable Modeling and Optimization for Chemical Engineering Applications: Interval Analysis Approach, Reliable Computing, 2006; 12:427-450.
- [2] C. S. Adjiman, S. Dallwig, C. A. Floudas, A. Neumaier; A Global Optimization Method, αBB, for General Twice-Differentiable Constrained NLPs

 I. Theoretical Advances, Computers and Chemical Engineering, 1998;
 22:1137-1158.
- [3] C. S. Adjiman, I. P. Androulakis, C. A. Floudas, A Global Optimization Method, αBB, for General Twice-Differentiable Constrained NLPs - II. Implementation and Computational Results, Computers and Chemical Engineering, 1998; 22:1159-1179.
- [4] J. W. Kovach, W. D. Seider; Heterogeneous Azeotropic Distillation: Homotopy-Continuation Methods, Comput. Chem. Eng., 1987; 11(6):593-605.
- [5] L. T. Watson, M. Sosonkina, R. C. Melville, A. P. Morgan, H. F. Walker; Algorithm 777: HOMPACK90: A Suite of Fortran 90 Codes for Globally Convergent Homotopy Algorithms, ACM Transactions on Mathematical Software (TOMS), 1997; 23(4):514-549.
- [6] A. P. Morgan; Solving Polynomial Systems Using Continuation for Engineering and Scientific Problems, PenticeHall, Inc., 1987.
- [7] K. Yamamura, T. Kumakura, Y. Inoue; Finding All Solutions of Nonlinear Equations Using Inverses of Approximate Jacobian Matrices, IEICE Trans. Fundamentals, 2001; E84-A(11):2950-2952.
- [8] I. P. Nenov, D. H. Fylstra; Interval Methods for Accelerated Global Search in the Microsoft Excel Solver, Reliable Computing, 2003; 9:143-159.

- [9] L. V. Kolev; A New Method for Global Solution of Systems of Non-Linear Equations, Reliable Computing, 1998; 4:125-146.
- [10] J. Stolfi, L. H. Figueiredo; Self-Validated Numerical Methods and Applications, Monograph for the 21st Brazilian Mathematics Colloquium (CBM'97), IMPA. Rio de Janeiro, Brazil; 1997.
- [11] A. Baharev, E. Rév; Reliable Computation of Equilibrium Cascades with Affine Arithmetic, AIChE Journal, 2008, 54(7):1782-1797.
- [12] R. B. Kearfott; Decomposition of Arithmetic Expressions to Improve the Behavior of Interval Iteration for Nonlinear Systems, Computing, 1991; 47(2):169-191.
- [13] T. Beelitz, A. Frommer, B. Lang, P. Willems; Symbolic-Numeric Techniques for Solving Nonlinear Systems, PAMM, 2005; 5:705-708.

Keywords: affine arithmetic, linear programming, root finding, branch-and-prune, dependency problem