COMPUTING DERIVATIVES OF A DAE SOLUTION IN PARALLEL

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ABSTRACT

We use Pryce's Σ -method for structural analysis of differential-algebraic equation systems (DAEs). The equation and variable offsets it computes often give the correct index of a DAE. They also prescribe a stage by stage solution scheme for computing derivatives of a solution, strictly from lower order to higher. This solution scheme can naturally develop into a Taylor series method, such as the Nedialkov-Pryce daets code. We observe the parallelism of the derivative computations from a DAE's block triangular form, which is based on the sparsity pattern of a Jacobian matrix. We illustrate this parallel computing idea with the chemical Akzo Nobel DAE, and seek to develop efficient algorithms for solving large and sparse problems.

References

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