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**An abstract method for extending two-level
preconditioners to multilevel preconditioners of
comparable quality**

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We present an abstract method for designing a multilevel preconditioner given a two-level preconditioner for an operator with positive definite symmetric part.

If we denote by A_{fine} and A_{coarse} two discrete versions of a continuous operator, then a two-level preconditioner T_{fine} for A_{fine} can be described in general by a function $T_{\text{fine}} = \mathcal{F}(A_{\text{coarse}}^{-1}, A_{\text{fine}})$, where it is assumed that the evaluation of \mathcal{F} requires a level-independent number of applications of A_{fine} and k applications of A_{coarse}^{-1} ($k = 1$ or 2). The natural extension to a multilevel preconditioner, consisting in replacing in T_{fine} the call to A_{coarse}^{-1} with a recursive call to \mathcal{F} , is known to sometimes produce multilevel preconditioners of lower quality (e.g., for certain types of inverse problems). Based on the idea that inverting A_{fine} essentially means to solve the nonlinear equation $X^{-1} - A_{\text{fine}} = 0$, we define our multigrid preconditioner to be the first Newton iterate of the map $X \mapsto X^{-1} - A_{\text{fine}}$ starting at the “natural” multilevel preconditioner. For $k = 1$, the resulting algorithm has a W-cycle structure, and differs only slightly from the textbook version of the W-cycle. Moreover, the method guarantees that the resulting preconditioner maintains the approximation quality of the initial two-level preconditioner. The quality of approximation is measured using a certain distance function, which determines the degree to which two operators with positive definite symmetric parts are spectrally equivalent. We apply this method to designing and analyzing a multigrid preconditioner for a linear advection-diffusion-reaction equation.

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