betrug: $w^{-1} = \sigma^2(F) + 0.0002F^2$. Abschliessende Differenz-Fourier-Synthesen enthielten keine wesentlichen Maxima.

Die Verfeinerungen konvergierten bei R 0.083, R_w 0.065 für IIIk und R 0.080, R_w 0.059 für IIIl. Die Atomkoordinaten sowie die Bindungsabstände und -winkel sind in den Tab. 6–11 zusammengestellt. Tabellen der Strukturfaktoren, anisotropen Temperaturfaktoren und Wasserstoffparameter können bei den Autoren angefordert werden.

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