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Density functional calculations of molecular g-tensors in the zero-order regular approximation for relativistic effects

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| 274 | Calculation of Zero-Field Splittings, g-Values, and the Relativistic Nephelauxetic Effect in Transition Metal Complexes. Application to High-Spin Ferric Complexes. <i>Inorganic Chemistry</i> , <b>1998</b> , 37, 6568-6582                                    | 5.1  | 268       |
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