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ES-MS spectrum of Inclusion of the complex [Pillar-2a G2-Br] + S13 Crystal Dimension/mm 022 X 014 X 011 021 X 012 X 009 Crystal Color, Habit Colorless, block Colorless, block Formula weight C116 H162 Br F3 O12 C116 H162 Br F3 O12

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S13 3) Details to the DFT Calculations of 1 and 2 General Considerations: DFT calculations of model compounds to cis/trans-1 and cis/trans-2 were carried out on the B3LYP/def2-SVP level of theory implementing the GAUSSIAN03 program[S4] pCartesian coordinates of the optimized structures are shown in Table S3, S4, S5 and S6, respectively