Figure S1. Plots of the metal-metal bonding energy (ΔE_{mbe}) as a function of the square of the spin density (ρ_s^2) in the S=S_{max} state for same-valence dⁿdⁿ [M₂Cl₉]³⁻ and [M'M"Cl₉]³⁻ systems.

Figure S2. Plots of the spin polarization energy (ΔE_{spe}) as a function of the square of the spin density (ρ_s^2) in the S=S_{max} state for same-valence dⁿdⁿ [M₂Cl₉]³⁻ and [M'M"Cl₉]³⁻ systems.

Figure S3. Plots of the metal-metal bonding energy (ΔE_{mbe}) and spin polarization energy (ΔE_{spe}) as a function of the square of the spin density (ρ_s^2) in the S=S_{max} state for mixed-valence dⁿ⁻¹dⁿ [M₂Cl₉]²⁻ systems.

Figure S1





Figure S2





6.0

8.0

 ${\rho_{\text{s}}}^2 \, (\text{S=S}_{\text{max}})$

4.0

Ru

2.0

Os

Figure S3



