Albumin binding and ligand-exchange processes of the Ru(III) anticancer agent NAMI-A and its bis-DMSO analogue determined by ENDOR spectroscopy

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Supporting Information

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Table S1: g values and line widths, used in the simulation of all EPR spectra.

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<th>Compound</th>
<th>g values</th>
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Figure S1 EPR spectra of Ru-bis-DMSO after incubation in PBS for 0, 10, 20, 30, and 60 minutes at 37 °C. For spectral parameters see experimental section in main text.
Figure S2 (a-e) Deconvolution of EPR spectra from Ru-bis-DMSO following incubation in PBS for 0, 10, 20, 30, and 60 minutes at 37 °C. For g values and linewidths used in simulations see Table S1.
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