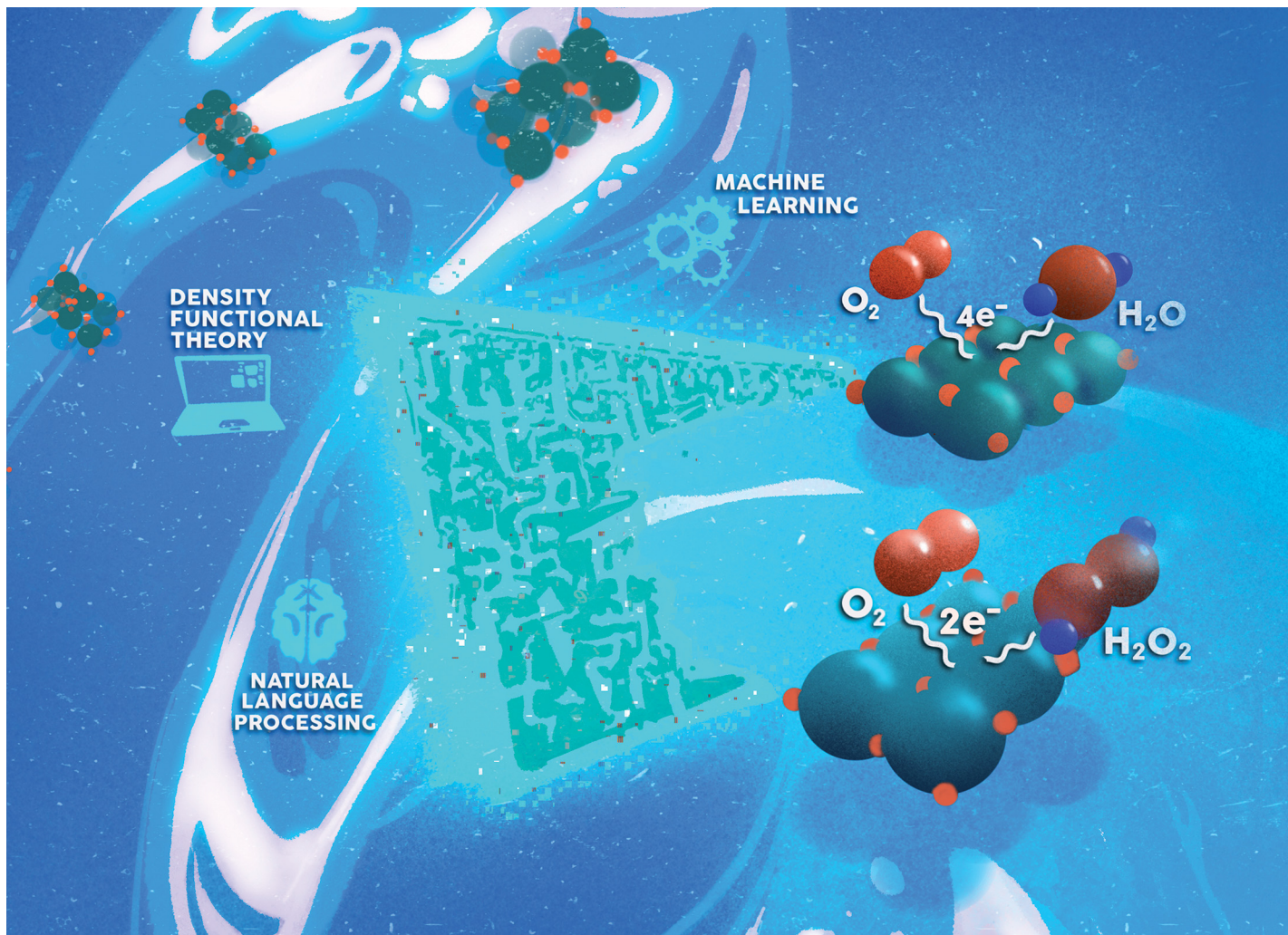


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Revolutionizing ORR catalyst design through computational methodologies and materials informatics

Recent advancements in computational methodologies, particularly density functional theory (DFT) and materials informatics, have revolutionized catalyst design for oxygen reduction reactions (ORR).

As featured in:



See Samira Siahrostami *et al.*, *EES. Catal.*, 2024, 2, 1037.