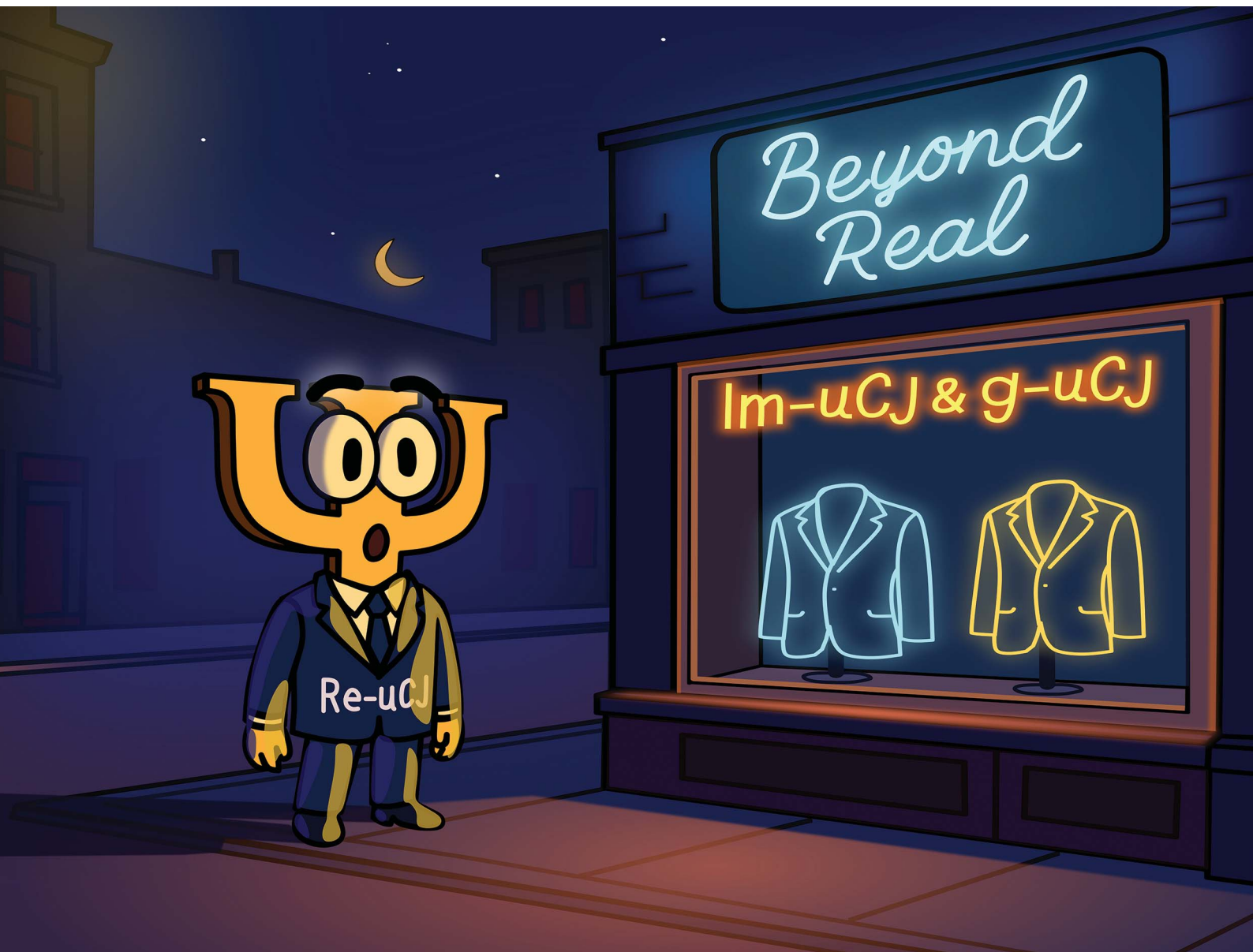


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EDGE ARTICLE

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molecular electronic structure calculations on near-term
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