Self-interaction correction to density-functional approx systems

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Citation Report

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8862	O<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:mrow><mml:mn>1</mml:mn><mml:mi>s</mml:mi></mml:mrow></mml:math>core-level shifts at the anatase TiO<mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mrow /><mml:mn>2</mml:mn></mml:msub></mml:math>(101)/N3 photovoltaic interface: Signature of H-bonded supramolecular assembly. Physical Review B, 2011, 84, .	1.1	24
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14072	xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline" overflow="scroll" id="d1e967" altimg="si17.gif"><mml:msub><mml:mrow /><mml:mrow><mml:mi>x</mml:mi></mml:mrow></mml:msub> Band structures in coupled-cluster singles-and-doubles Greenâ€™s function (GFCCSD). Journal of Chemical Physics, 2018, 148, 204109.	1.2	20
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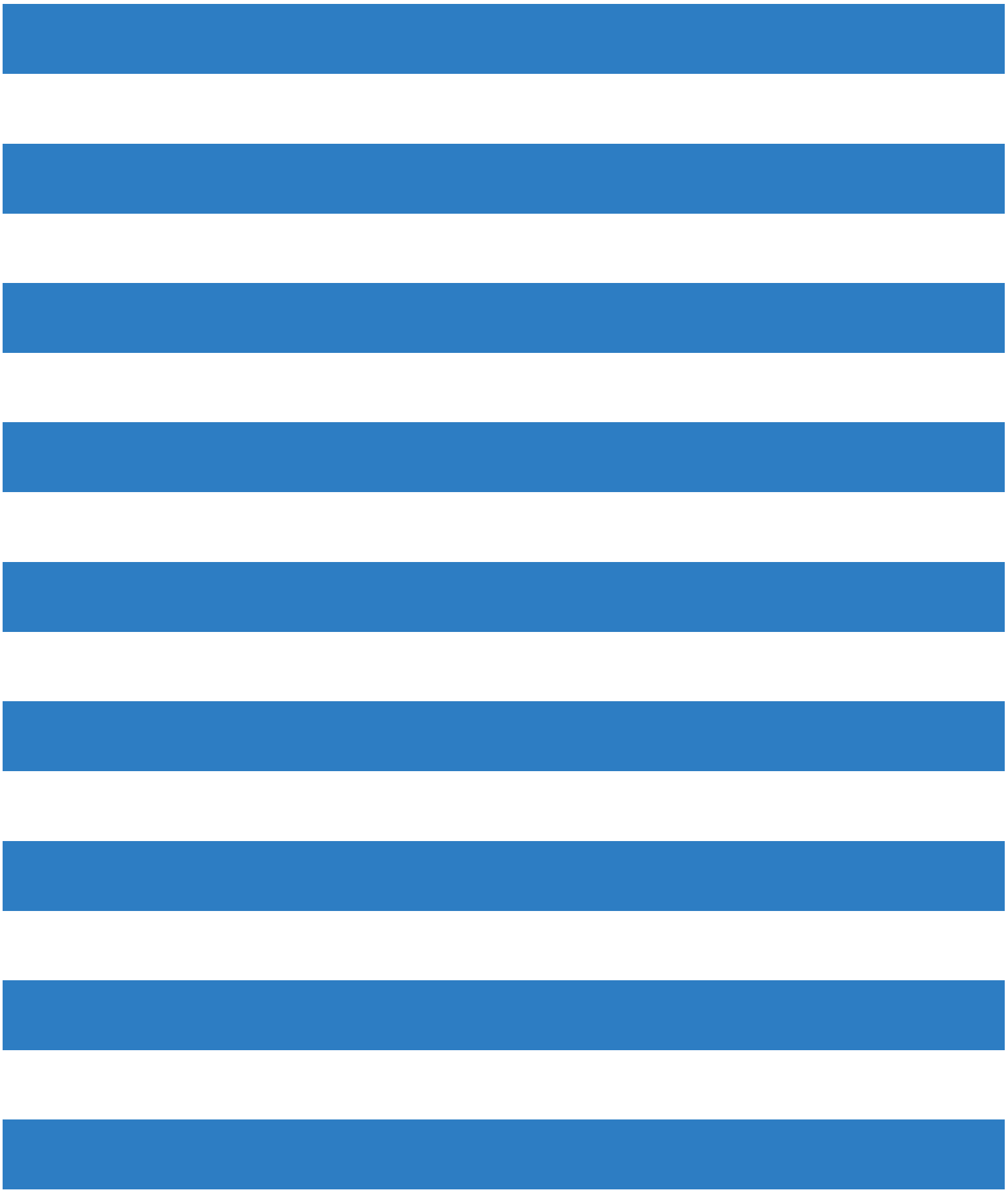

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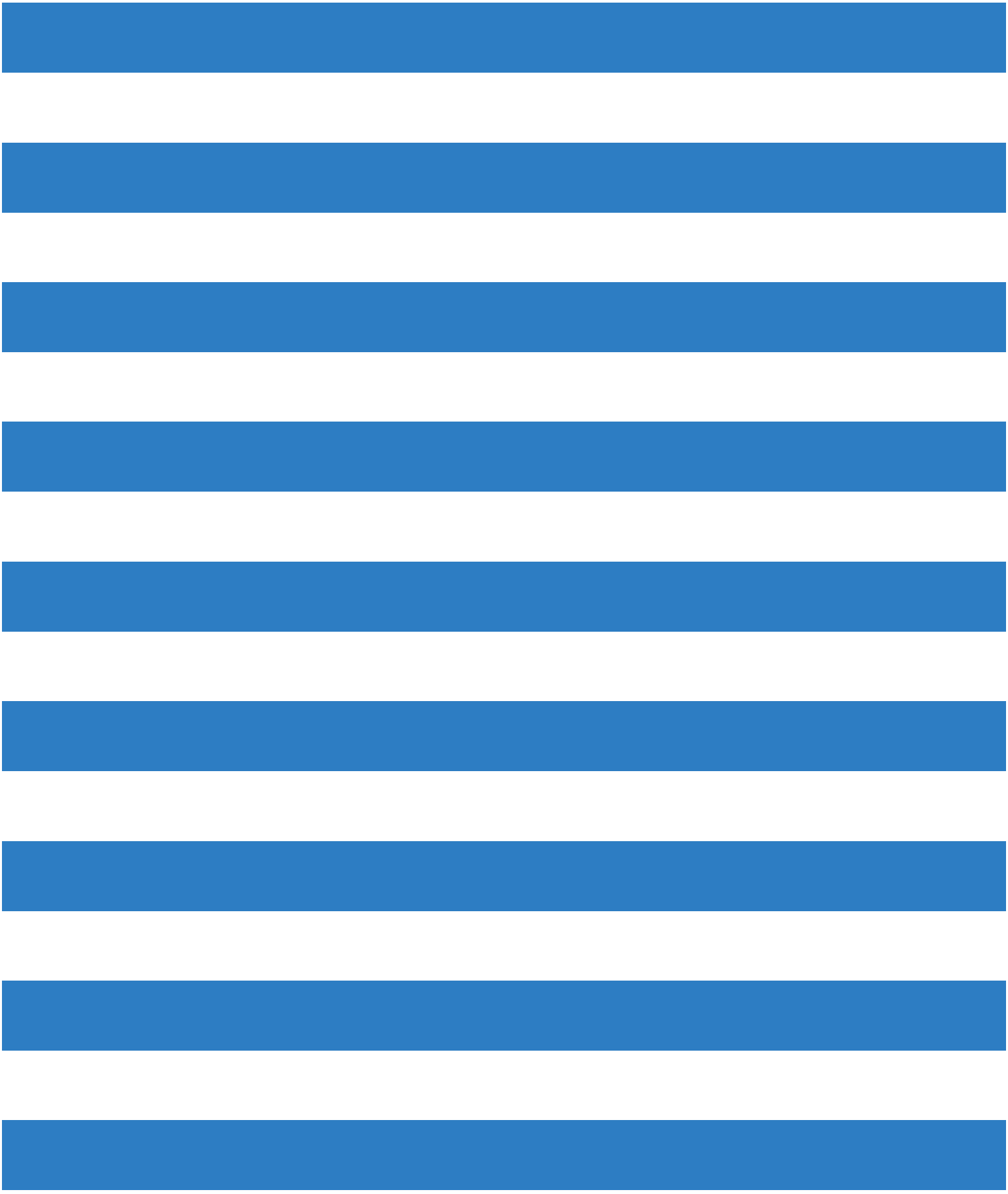

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