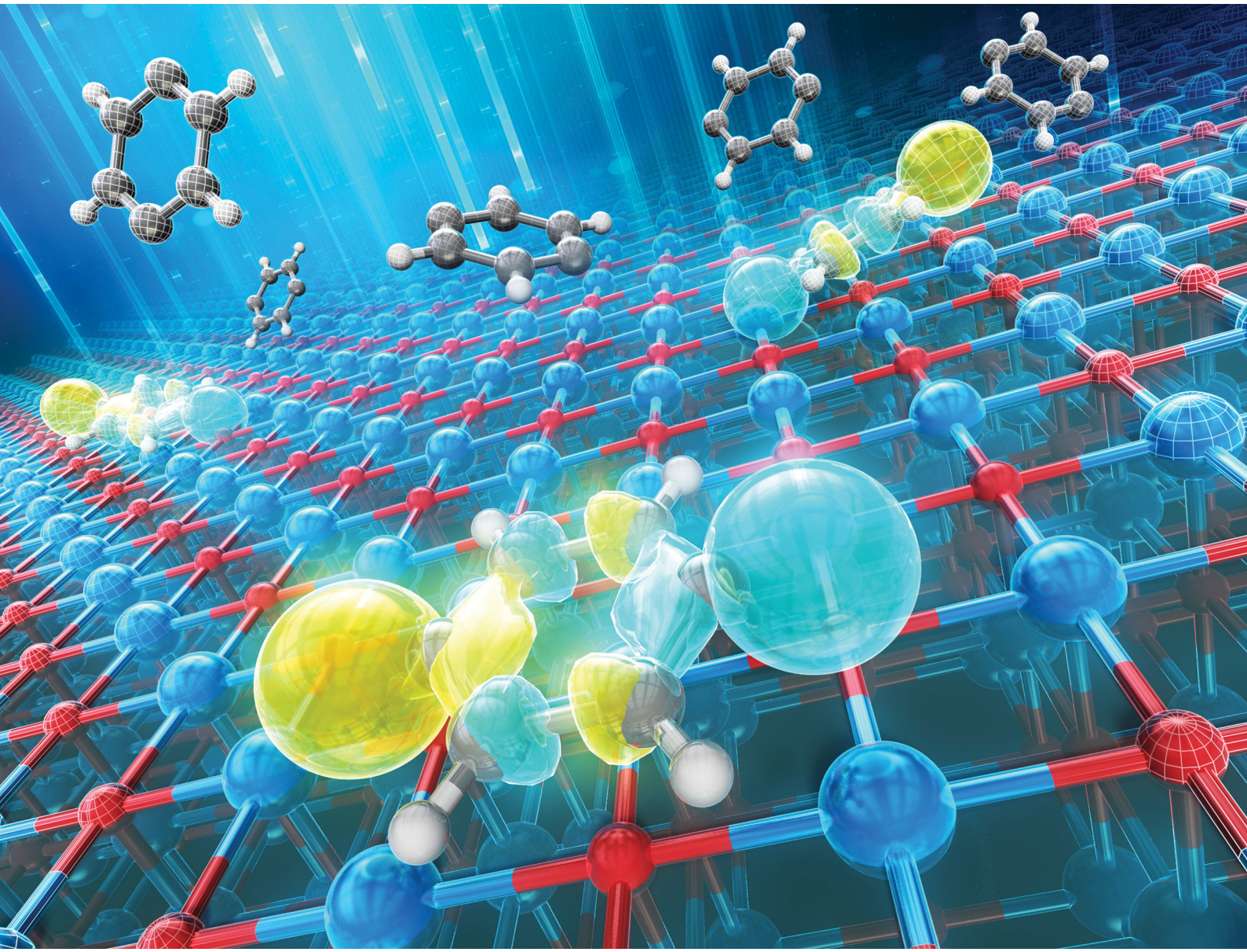
Volume 25
Number 43
21 November 2023
Pages 29367–30080

PCCP

Physical Chemistry Chemical Physics

rsc.li/pccp



ISSN 1463-9076

ROYAL SOCIETY OF CHEMISTRY

PAPER

Kohei Tada *et al.*

Model calculations for the prediction of the diradical character of physisorbed molecules: *p*-benzyne/MgO and *p*-benzyne/SrO

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