

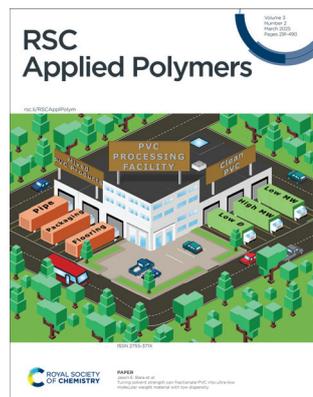
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eISSN 2755-371X CODEN RAPSBD 3(2) 291–490 (2025)



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See Jason E. Bara *et al.*,
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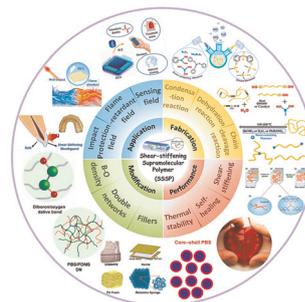
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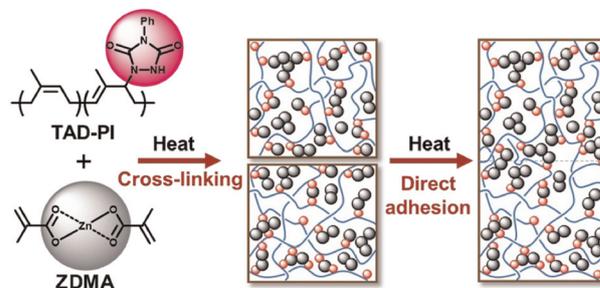
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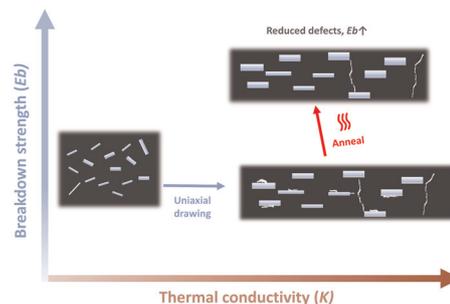
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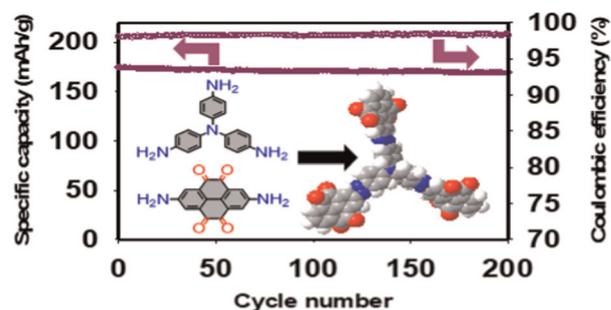
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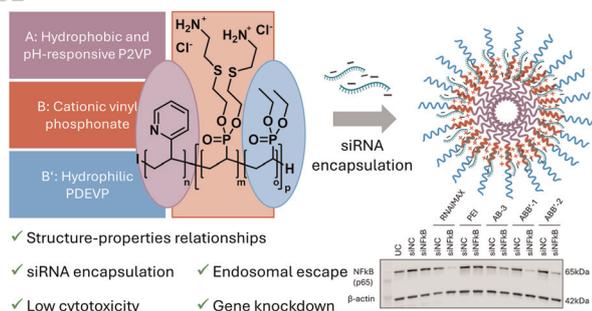
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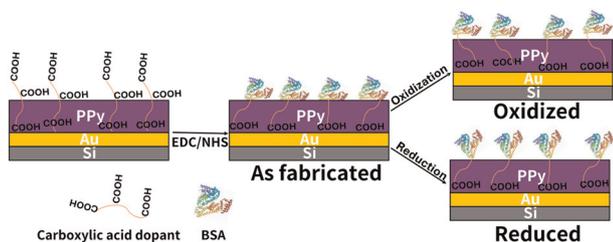
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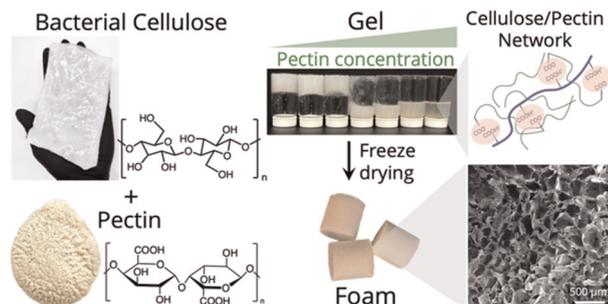
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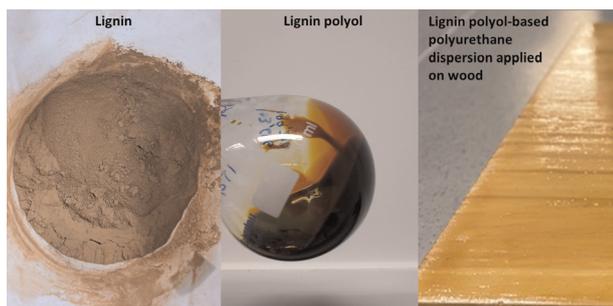
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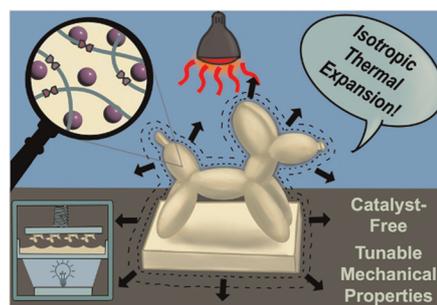
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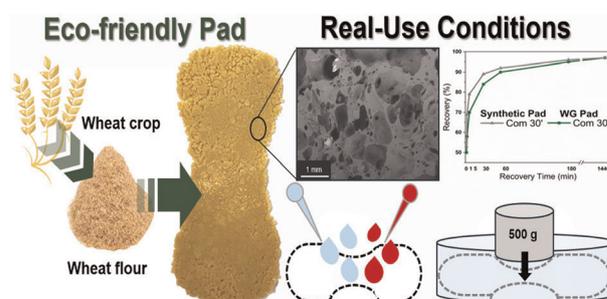
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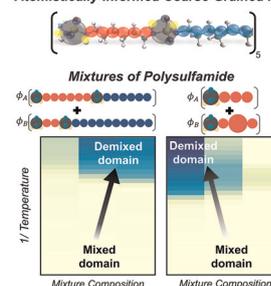


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Coarse-grained molecular dynamics simulations of mixtures of polysulfamides

Jay Shah and Arthi Jayaraman*

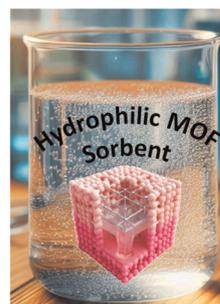
Atomistically-informed Coarse-Grained model

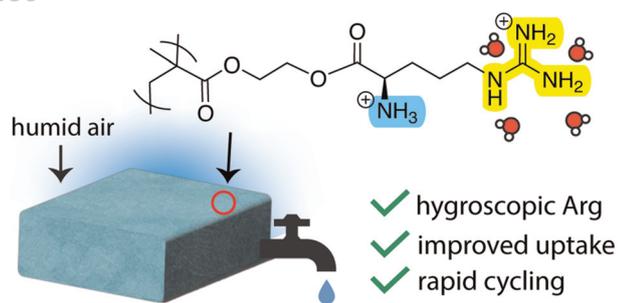


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Moki K. Thanusing, Brett L. Pollard and Luke A. Connal*

