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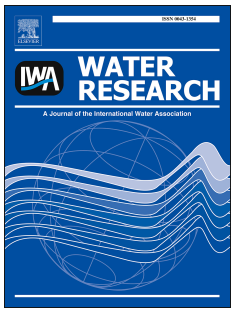

Correlation and prediction of adsorption capacity and affinity of aromatic compounds on carbon nanotubes

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Graphical abstract

Prediction model:

$\log q_e = \log \boldsymbol{Q^0} - (\varepsilon/\boldsymbol{E})^{\boldsymbol{b}}$

$\boldsymbol{Q^0} = f(\boldsymbol{A}, \boldsymbol{R}_{meso}, \mathbf{MP}, \boldsymbol{N})$

$\boldsymbol{E} = f(\alpha_m, \boldsymbol{R}_{meso}, \pi^*)$

$\boldsymbol{b} = f(\alpha_m, \boldsymbol{R}_{meso}, \pi^*)$

$\varepsilon = -\mathrm{RTln}(C_e/C_s)$