Benzene on Cu(111): I. Application of van der Waals-Density Functional Formalism to Determine Binding Sites and Energy Contour Map

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1Supported by (KB & PH) Swedish Vetenskapsrådet VR #621-2008-4346 and by (TLE) NSF CHE 07-50334, also secondarily by UMD MRSEC DMR 05-20471.
3KB, TLE, and PH, Phys. Rev. B 80 (2009) 155431