

Accepted Manuscript

Molecular dynamics simulation of functionalized graphene surface for high efficient loading of doxorubicin

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PII: S0022-2860(17)30438-6

DOI: [10.1016/j.molstruc.2017.04.007](https://doi.org/10.1016/j.molstruc.2017.04.007)

Reference: MOLSTR 23624

To appear in: *Journal of Molecular Structure*

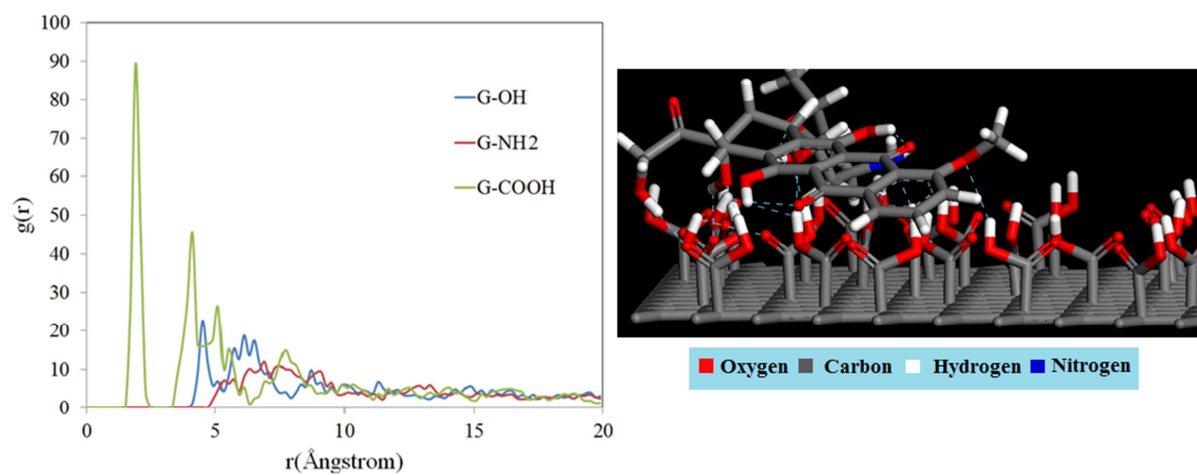
Received Date: 7 March 2017

Revised Date: 31 March 2017

Accepted Date: 3 April 2017

Please cite this article as: M.M. Mirhosseini, M. Rahmati, S.S. Zargarian, R. Khordad, Molecular dynamics simulation of functionalized graphene surface for high efficient loading of doxorubicin, *Journal of Molecular Structure* (2017), doi: 10.1016/j.molstruc.2017.04.007.

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