

Title :**Development of molecular pseudopotentials for the study of properties and chemical reactions in condensed phase.****Department :** Institut des Sciences Moléculaires de Marseille

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Team: Chimie Théorique et Modélisation**Supervisor :** Dr. Yannick Carissan**Co-supervisor :** Dr. Paola Nava**Contact :** Yannick Carissan

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Position : 3-years scholarship available from October 2016 (gross salary: about 1600 euros/month)**Deadline :** 20 may 2016 – deadline for sending the candidature to the ED250 (Doctoral School)corinne.esquiva@univ-amu.fr

The candidate must contact the supervisor or the co-supervisor before this date for a pre-selection.

31 may 2016 – Interview at the ED250 (Doctoral School)

Project :

This theoretical chemistry project aims to simplify the study of experiences made in rare-gas matrices for reproducing interstellar conditions.

From a computational point of view, the description of the environment (i.e. the rare-gas matrix) is resource-consuming. Within the CTOM group, pseudopotentials have been already developed to treat the environment of chemical systems by reducing the computational cost. We propose to further develop those methods and to adapt them to reproduce effects of more complex chemical means, such as the gas-rare matrices. This project will imply some programming (Fortran and/or python).

Context :

This study continues efforts on pseudopotential development of the CTOM group, as well as current and established joint work with experimentalists on systems in matrices. [1,2]

References:

[1] J. Drujon and Y. Carissan, *Journal of Computational Chemistry*, 2013, **34**, 49–59.

[2] T. Butscher, F. Duvernay, P. Theule, G. Danger, Y. Carissan, D. Hagebaum-Reignier, and T. Chiavassa *MNRAS*, 2015, **453**, 1587-1596.