

UNIVERSITY OF 8 MAI 1945 GUELMA
FACULTY OF MATHEMATICAL, INFORMATICS AND MATERIAL SCIENCES
LABORATORY OF COMPUTATIONAL CHEMISTRY AND NANOSTRUCTURES -LCCN-

Organize

3RD NATIONAL SEMINAR OF MOLECULAR MODELING

- APRIL 15TH, 2026



جامعة 8 ماي 1945 قالمة
UNIVERSITE 8 MAI 1945 GUELMA



✉ 3snmm2025@gmail.com

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Topics

- 1 AB-INITIO AND DFT
- 2 DOCKING ,MOLECULAR DYNAMICS AND MONTE CARLO
- 3 REACTIVITY AND MODELING INCLUSION COMPLEXES

Important Dates

Deadline for receipt : 28/02/2026

Acceptance : 18/03/2026

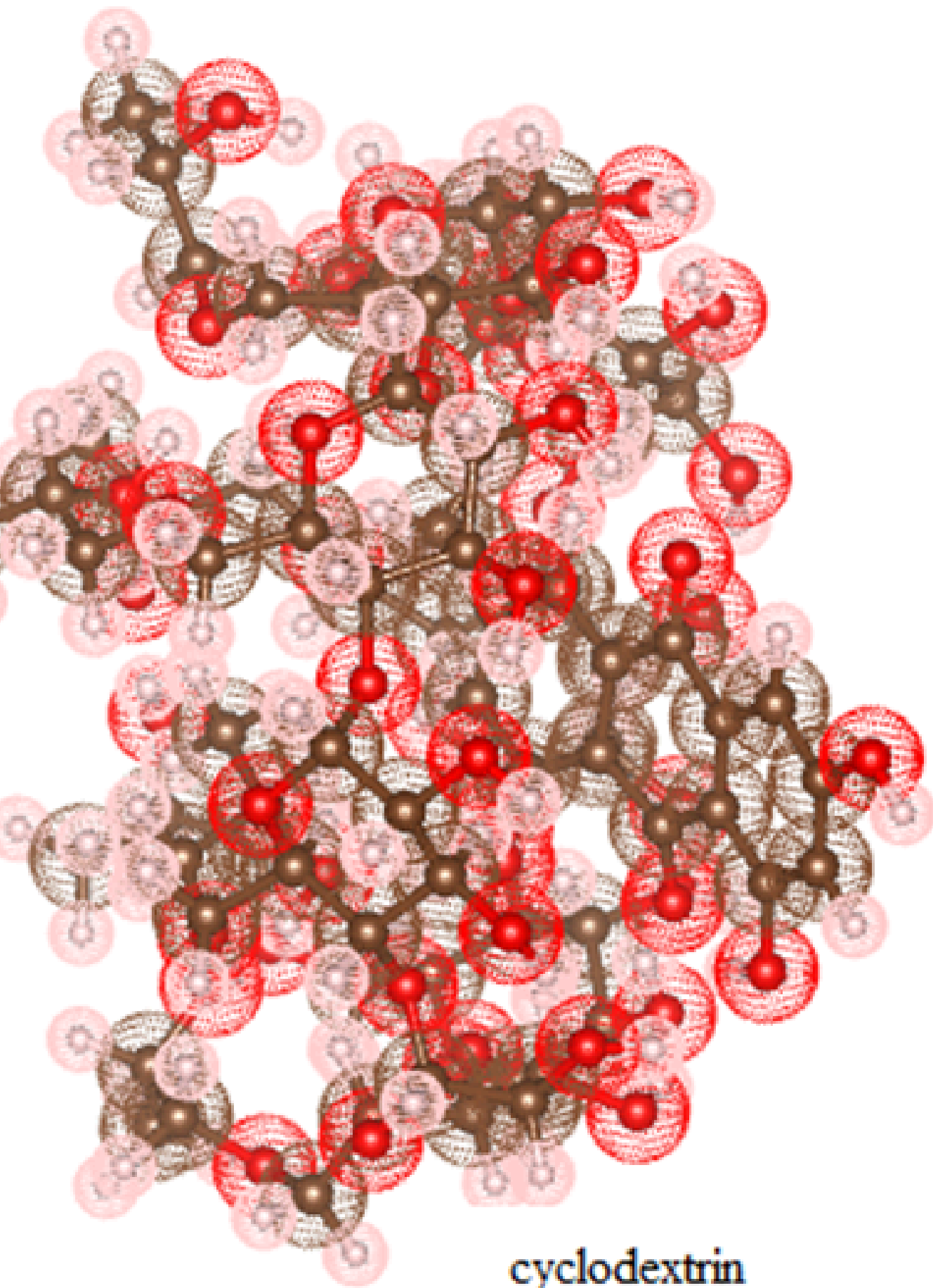
Seminar: 15/04/2026

Registration fees

• Research Teachers of the University	Amount : 8 000,00 DA
• Research teachers from other universities	Amount : 12 000,00 DA
• Research students	Amount : 5 000,00 DA

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cyclodextrin
inclusion complex

Objective

This is the third time the LCCN has organized a molecular modeling seminar, with the aim of bringing together doctoral students, research professors, and researchers to discuss advances in molecular modeling methods and techniques, structures, and molecular interactions in various chemical, physical, and biological systems.

PhD students and research professors are strongly encouraged to participate in informal discussions through oral presentations or posters.

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Languages of Presentation:

Instructions for Authors

Participants are invited to submit their abstracts for oral or poster presentations.

Abstracts should be submitted by email. They should be written on an A4 page with 2.5 cm margins on each side, in Times New Roman, size 12 font.

Titles should be in bold, capital letters. Author names should be in Times New Roman, size 12 font, with the corresponding author's name underlined. Author addresses should be in italics (Times New Roman, size 12).

The texts should be right- and left-justified with single-line spacing.

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- French
- English
- Arabic



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Guests of Honor:

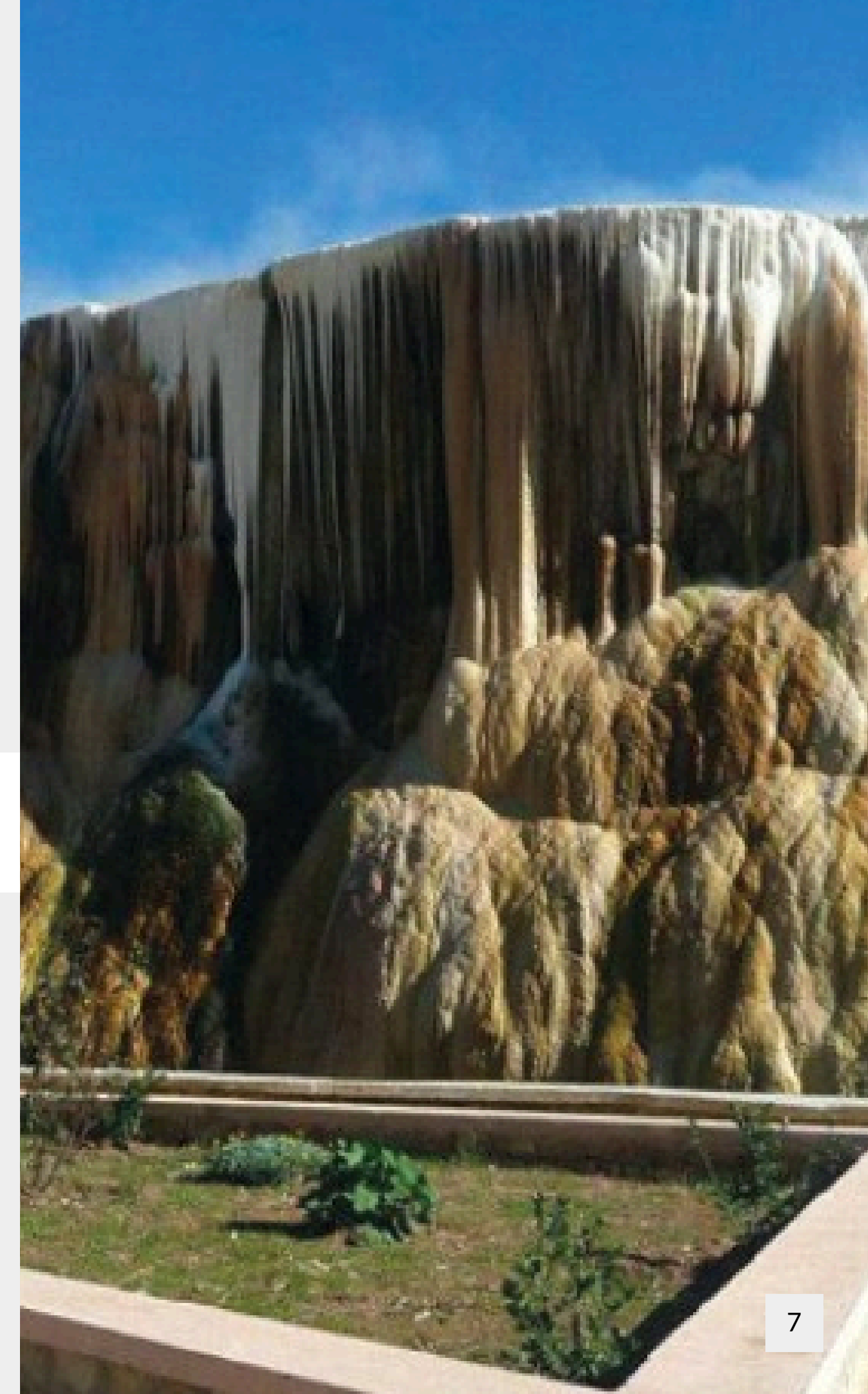
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Thank you very much for your participation

FOR MORE INFORMATION



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