

SCIENTIFIC AND METHOD MODULES

| Module name | From Molecules to Materials | | | |
|---|---|--|--|--|
| Number | 2016-T4 | | | |
| Aims | This module links molecular sciences and materials science, teaches how materials with optimized catalytic activity and adjustable magnetic, electronic, o optical properties are obtained from molecules, and provides an understanding o the properties and applications of these materials. | | | |
| Basics | covered in basic modules B1–B3 (molecular precursors, supramolecular chemistry, polymers, organic and inorganic nanostructures) and MOFs covered in the module "Smart Molecules" (T1) | | | |
| Contents | Materials from "hard" (synthetic molecules, crystalline nanostructures, quantum- electronic structures) building blocks and/or "soft" (polymers, biomolecules, MOFs, metal-binding peptides, thin films) building blocks, Supramolecular arrangements (proteins and peptides together with modifications to improve the material qualities [pegylation, lipidation, glycosylation]), Properties of these materials (magnetic, electronic, and optical properties, electronic, photonic, and magnetoresistive devices, superconductivity), Applications (catalysis [immobilised catalysts, MOFs], gas separation or gas storage [MOFs], sensors in electronics and photonics, quantum information technology [spintronics], energy conversion [including solar energy]), Theory (structure and interfaces within systems containing complex nano-aggregates), Atomistic particle transport through MOFs. | | | |
| Methods | Immobilisation techniques, Deposition (chemical vapour deposition [CVD, PECVD, MOVPE], physical deposition [PLD], atomic layer deposition [ALD]), Surface analysis (MIES), Device characterization (e.g. <i>I</i> - <i>V</i> , <i>C</i> - <i>V</i> , S-Parameter), Modelling (mesoscale simulations, classical molecular dynamics simulations, multi-scalar approach, QM/MM simulations, advanced Monte Carlo techniques). | | | |
| Туре | Two-day block course/ bi-yearly recurrence with modification | | | |
| Date (month/year) | 19/20 September 2016 | | | |
| Time | 10:00-19:00; 10:00-15:20 | | | |
| Work load | 15 hours presence/ 45 hours self-study | | | |
| Examination | Poster presentation about a self-chosen topic about "Artificial molecules and solid state crystals" (own research or from literature) and discussion (oral) in front of the poster with the organizer(s) | | | |
| Credit points | 2 | | | |
| Responsible scientists | Prof. Jan Meijer, Prof. Pablo Esquinazi | | | |
| International guest lecturers | J. M. van Ruitenbeek (Leiden University, Netherlands), P. Olivero (University of Torino, Italy) | | | |
| Industrial partners | | | | |
| Recommendations for literature, e- learning | JM. Spaeth, H. Overhof: "Point Defects in Semiconductors and Insulators"; Susan Shannon (Editor): "Trends in Quantum Computing Research"; M.A. Mielsen and I. L.Chuang: "Quantum Computation and Quantum Information" | | | |

SCHEDULE for Module 2016-T4

| Time | Lecturer | Programme | Location | |
|-------|----------|-----------|----------|--|
| Day 1 | | | | |
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| Day 2 | | | | |
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Didactic elements:

Lecture, discussions, practical training – lab demonstration, etc.

Expected performance: Active participation in discussions during lab demonstration etc.