

<b>Modulbezeichnung:</b> Theory (MSM-ME2) (Theory)	<b>15 ECTS</b>	
Modulverantwortliche/r:	Andreas Görling	
Lehrende:	Andreas Heßelmann, Wolfgang Hieringer, Bernd Meyer, Andreas Görling	
Startsemester: WS 2019/2020	Dauer: 2 Semester	Turnus: halbjährlich (WS+SS)
Präsenzzeit: 210 Std.	Eigenstudium: 240 Std.	Sprache: Englisch

#### Lehrveranstaltungen:

##### A. Quantum Chemistry-I (2L/1S)

Quantum Chemistry I / Quantenchemie I (WS 2019/2020, Vorlesung, 2 SWS, Andreas Görling)

Quantum Chemistry I - Exercises / Übung zur Quantenchemie I (WS 2019/2020, Übung, 1 SWS, Jannis Erhard et al.)

Quantum Chemistry II (SS 2020, Vorlesung, 2 SWS, Andreas Görling)

Quantum Chemistry II (Seminar) (SS 2020, Übung, 1 SWS, Andreas Görling et al.)

##### B. Modeling of catalytic processes (2L/1S)

Modeling of Catalytic Processes (SS 2020, Vorlesung, 2 SWS, Bernd Meyer)

Modeling of Catalytic Processes (Praktikum) (SS 2020, Praktikum, 2 SWS, Bernd Meyer et al.)

##### C. Scientific programming (2LAB/1S)

Attendance in lab courses is compulsory!

Scientific Programming / Wissenschaftliches Programmieren (WS 2019/2020, Praktikum, 2 SWS, Andreas Heßelmann et al.)

##### D. Handling of computer systems in science (2LAB/1S)

Attendance in lab courses is compulsory!

Handling of computer systems in science (SS 2020, Praktikum, 2 SWS, Wolfgang Hieringer et al.)

##### E. Practical Training in Computer Chemistry (4LAB)

Attendance in lab courses is compulsory!

Practical Training in Computer Chemistry / Praktikum Computerchemie (WS 2019/2020, Praktikum, Andreas Görling et al.)

Practical training in computer chemistry (SS 2020, Praktikum, 4 SWS, Andreas Görling et al.)

#### Inhalt:

##### A: Quantum Chemistry I:

Consolidation of the mathematical backgrounds in quantum chemistry, Hartree-Fock method, configuration interactions; density functional theory and its application to molecular systems

##### B: Modeling of catalytic processes

Introduction to the theoretical concepts and methods to study catalytic processes: energetic, kinetics and dynamics of adsorbates, reactivity of surfaces; transition state theory, microkinetic modeling, kinetic Monte-Carlo techniques, molecular dynamics

C: **Lab course:** Scientific programming using FORTRAN

D: **Lab course:** Introduction to Linux systems;

E: **Lab course:** application of modern modeling techniques to investigate molecular systems

#### Lernziele und Kompetenzen:

The students

- get experience with advanced knowledge and techniques in theoretical chemistry
- are able to utilize advanced computer-based techniques to model research related problems in the field of chemistry, biochemistry, catalysis and material science
- learn to operate Linux-based and large-scale computing systems
- are able to summarize and to interpret theoretical calculations in written form (lab report).

#### Verwendbarkeit des Moduls / Einpassung in den Musterstudienplan:

Das Modul ist im Kontext der folgenden Studienfächer/Vertiefungsrichtungen verwendbar:

##### [1] Molecular Science (Master of Science)

(Po-Vers. 2013 | NatFak | Molecular Science (Master of Science) | Wahlpflichtmodul Molecular Science)

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**Studien-/Prüfungsleistungen:**

Theorie und Modellierung - Theory (Prüfungsnummer: 30802)

(englische Bezeichnung: Theory and Modelling - Theory)

Prüfungsleistung, schriftlich oder mündlich

Anteil an der Berechnung der Modulnote: 10%

weitere Erläuterungen:

Assessment and examinations: Oral examination (45 min) or alternative examination according to FAU Corona statutes!

Prüfungssprache: Englisch

Erstablingung: SS 2020, 1. Wdh.: WS 2020/2021

1. Prüfer: Andreas Göring

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**Organisatorisches:**

**Intended stage in the degree course:** Mandatory Elective Module (Wahlpflichtmodul) or Elective Module (Wahlmodul), semester 1-3

**Frequency of offer:** Annually, **A:** winter term **B:** summer term **C:** winter term **D:** summer term **E:** winter and summer term