

Final schedule of AK-Ulrich Group Meeting in Freudenstadt 18-22 September 2011

Time	Sunday 18/9	Monday 19/9	Tuesday 20/9	Wednesday 21/9	Thursday 22/9
8.00		Breakfast	Breakfast	Breakfast	Breakfast
9.00	8.45 Meeting at Karlsruhe hbf 9.05 Leaving Karlsruhe Travel with RB + bus Ca 11.00 arrival at Hotel in Freudenstadt	Session 4. Peptides II (Parvesh) Ronald <i>Linking structure and dynamics of membrane-active peptides</i> Papia <i>Membrane-bound structure of a CPP derived from Kaposi fibroblast growth factor</i> Jonathan <i>Investigation of synergistic re-orientation of magainin 2 and PGLa in lipid membranes via ¹⁹F- and ¹⁵N-ssNMR</i>	Session 8. Hardware (Stephan) * Markus <i>Probe development</i> * Martin, Daniela, Julian, Tobias, Patrick, Andreas <i>Konstruktion eines Kondensatorenteststandes und Prototypenfertigung für die NMR #</i> * Hartmut & Johannes P <i>Protein Origami – a new way of protein synthesis</i>	Session 9. Simulations I (Chair: Erik) † Wolfgang Wenzel <i>Protein folding and structure prediction with free energy methods</i> † Moritz Wolf <i>Implicit hydrophobic pore potential simulations running in Gromacs</i>	Session 13. Mirror motifs I (Johannes R.) Sebastian <i>Mirror motifs in membrane proteins and peptides</i> Thorsten <i>The Tat dependent translocation: Experimental proof of the charge ZIPper hypothesis - Part 1</i> Christina <i>The Tat dependent translocation: Experimental proof of the charge ZIPper hypothesis - Part 2</i>
10.45	11.30 Coffee	Discussion and coffee	Discussion and coffee	Discussion and coffee	Discussion and coffee
11.15	12.00 Session 1. Introduction (Chair: Birgid) Anne <i>Welcome and introduction</i> Erik <i>Seminar information</i>	Session 5. Peptides and beyond (Stephan) Julia <i>Biomembranes for ssNMR</i> Sergii <i>Epsin 1-18 and its interactions with lipid membranes</i> Oleg <i>Diarylethylene-derived photoswitching peptides – toward the light-controlled antimicrobial peptides</i>	Free for activities	Session 10. Simulations II (Chair: Marcus Elstner) † Jakob Ulmschneider <i>In silico partitioning and transmembrane insertion of hydrophobic peptides under equilibrium conditions</i> † Tomas Kubar <i>Simulating peptides in biomembranes with coarse grained models</i>	Session 14. Session 5. Mirror motifs II / Gramicidin S (Torsten) ‡ Eva <i>Untersuchungen der Oligomerisation der TatA-Translocase</i> Marco Jan <i>What we can get from the head</i> Mareike <i>Establishment of the PhoD translocation assay #</i> Marina <i>Synthetic biology of gramicidin S producers</i> END
13.00	Lunch	Discussion and lunch		Discussion and lunch	13.15 Discussion and lunch
14.30	Session 2. Dynamics (Jochen) Erik <i>Dynamics of membrane-bound peptides studied by solid-state NMR</i> † Jesús Salgado <i>Non-classical hydrophobic mismatch: Effect of helix geometry and azimuthal rotation</i> Stephan <i>Fitting NMR data of peptides in oriented bilayers</i>	Session 6. Fluorine chemistry (Sergii) † Pavel Myhailiuk <i>The many roles of fluorine in biochemistry</i> Anton <i>Design and synthesis of novel fluorinated serine and phenylalanine analogues as potential ¹⁹F-NMR labels #</i> Vladimir <i>Fluorine-labelled gramicidin S analogs: orientational behavior #</i>		Session 11. Simulations III (Chair: Jesús Salgado) † Thomas Steinbrecher <i>A model of tIsB-biomembrane insertion from MD simulations</i> ‡ Benjamin <i>Struktur und Funktionsuntersuchung am TisB</i> Nico <i>Structure analysis of the AMP Temporin A</i>	Travel to Karlsruhe
16.25	Discussion and coffee	Discussion and coffee		Discussion and coffee	
16.55	Session 3. Peptides I (Erik) * Andrea <i>Festphasen-Peptidsynthese</i> * Kerstin <i>Peptidreinigung mit HPLC</i> Parvesh <i>Solid-state ¹⁹F-NMR structure analysis of the β-sheeted antimicrobial peptide (KIGAKI)₃ in aligned lipid bilayers</i> Susanne <i>Structure analysis and aggregation behavior of the cell-penetrating peptide TP10</i>	Session 7. Optical spectroscopy (Dirk) Jochen <i>Commissioning of the SRCD beamline at ANKA</i> * Siegmur <i>CD12: Status der Strahlrohr- und Messplatzsysteme</i> Johannes R <i>Some tools to investigate peptide-lipid interactions by fluorescence spectroscopy</i> Joana <i>Optimization of hemolysis</i>		Session 12. E5/PDGFR (Marina) Colin <i>Structural investigation of the orientation of E5 in different membrane-mimicking systems</i> Marcel <i>Structure and orientation of the transmembrane domain of PDGFR β</i> Dirk <i>Reconstitution of E5 and PDGFR-TM in lipid bilayers</i>	Symbols: † Guest speakers (45 min) *Technicians (15 min) ‡ Students short talk (15 min) Students, postdocs (35 min) # Full titles not shown due to lack of space
18.35	Discussion	Discussion		Discussion	Final version last updated 110913
19.30	Dinner	Dinner	Dinner	Dinner	