

Polarized ATR-FTIR studies of DPPC/DPPG lipid bilayers doped with PLL

Autorzy

Paulina Trombik

Katarzyna Cieślik-Boczula

Rok wydania

2025

Czasopismo

Journal of Molecular
Structure

Numer woluminu

1321

Strony

140051/1-140051/9

DOI

10.1016/j.molstruc.2024.140051

Kolekcja

Naukowa

Język

Angielski

Typ publikacji

Artykuł

Streszczenie

The structural characterization of lipid bilayers is crucial for understanding various biological processes and designing effective systems of drugs. Attenuated Total Reflection Fourier Transformed Infrared (ATR-FTIR) spectroscopy is a powerful technique for probing lipid membrane properties, but conventional methods face challenges in discerning molecular orientation and intermolecular interactions. The polarized ATR-FTIR spectroscopy combined with Principal Component Analysis (PCA) was used to investigate the interactions between positively charged peptide (poly-L-lysine (PLL)) and anionic lipid bilayers (dipalmitoylphosphatidylcholine/dipalmitoylphosphatidylglycerol, DPPC/DPPG). This integrated approach offers deeper insight into mutual interactions between proteins and lipids of membrane, providing a comprehensive understanding of structure of both protein and lipid bilayers components. The findings of this study shed light on the orientational order of lipid molecules in DPPC/DPPG membrane structure as well as secondary structures of PLL peptides. It was shown that the presence of PLL peptide disrupts lipid molecule ordering within the lipid bilayer, primarily through electrostatic and H-bond interactions. The utility of polarized ATR-FTIR spectroscopy supported by PCA calculations in elucidating molecular organizations within complex membrane systems is emphasized.

Słowa kluczowe

DPPC/DPPG membranes, Poly-L-lysine, Polarized light, Principal Component Analysis (PCA), ATR-FTIR, Degree of orientational order

Adres publiczny

<http://dx.doi.org/10.1016/j.molstruc.2024.140051>

Strona internetowa wydawcy

<http://www.elsevier.com>

Plik został wygenerowany dnia 2026-06-22 11:15:13

Adres w repozytorium <https://old.chem.uni.wroc.pl/pl/repozytorium/XjqpVEU>.