DARR - Solid-state Correlation NMR Spectroscopy for Large Molecules -

Structural analysis of large molecules such as proteins using NMR is getting popular in the field of solid-state NMR. DARR (Dipolar Assisted Rotational Resonance) is one of techniques for analyzing large molecules in solids. The method can yield ¹³C-¹³C correlations for ¹³C-labeled samples, leading to intermolecular and intramolecular structures.

Figure shows DARR spectra of 13 C-labeled bacteriochlorophyll *c* (BChl *c*). Different rates of labelling distinguish intermolecular and intramolecular correlations, while variation of mixing times gives information on 13 C- 13 C distances.





Sample: 100% labelled BChl c (upper left) 50% labelled BChl c (lower right) Spectrometer: JNM-ECA600 Probe: 4mm CPMAS

Mixing time: 200 ms

Reference

Y. Kakitani, K. Harada, T. Mizoguchi, and Y. Koyama, Biochemistry, 46, 6513-6524 (2007): "Isotopic Replacement of Pigments and a Lipid in Chlorosomes from *Chlorobium Limicola*: Characterization of the Resultant Chlorosome."

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