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**Title:** Investigations of Biomolecular Conformation and Dynamics using <sup>19</sup>F NMR

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**Abstract:** Fluorine NMR is a useful probe of protein structure, conformation and folding in fluorinated proteins, owing to the high sensitivity, large chemical shift of the <sup>19</sup>F nucleus and the lack of high background signal to be suppressed. Protein NMR studies using fluorinated amino acids are a valuable addition to the existing suite of experiments with isotopically enriched nuclei. Recently, there have been several efforts to develop novel methods of biosynthetically tagging proteins to contain <sup>19</sup>F labels using fluorinated amino acids. Each fluorine signal in the resultant protein then becomes a reporter of the protein's structure and conformational dynamics. <sup>19</sup>F NMR relaxation times (including T<sub>1</sub> relaxation and homonuclear and heteronuclear NOE relaxation) can give detailed structural information. The <sup>19</sup>F CSA mechanism is an important relaxation mechanism at high magnetic fields and several novel experiments have been designed, to determine the complete <sup>19</sup>F CSA tensor. Several groups are also trying to apply <sup>19</sup>F NMR techniques to study membrane proteins and intrinsically disordered proteins which have a fluorinated tag. This chapter will focus on novel NMR techniques using fluorine as a probe and their application to structure determination and conformational dynamics of fluorinated biomolecules.

**URI:** <https://www.semanticscholar.org/paper/Investigations-of-Biomolecular-Conformation-and-19-Dorai/be553c6dd5460b05f472f33450615218894e9c99>  
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