

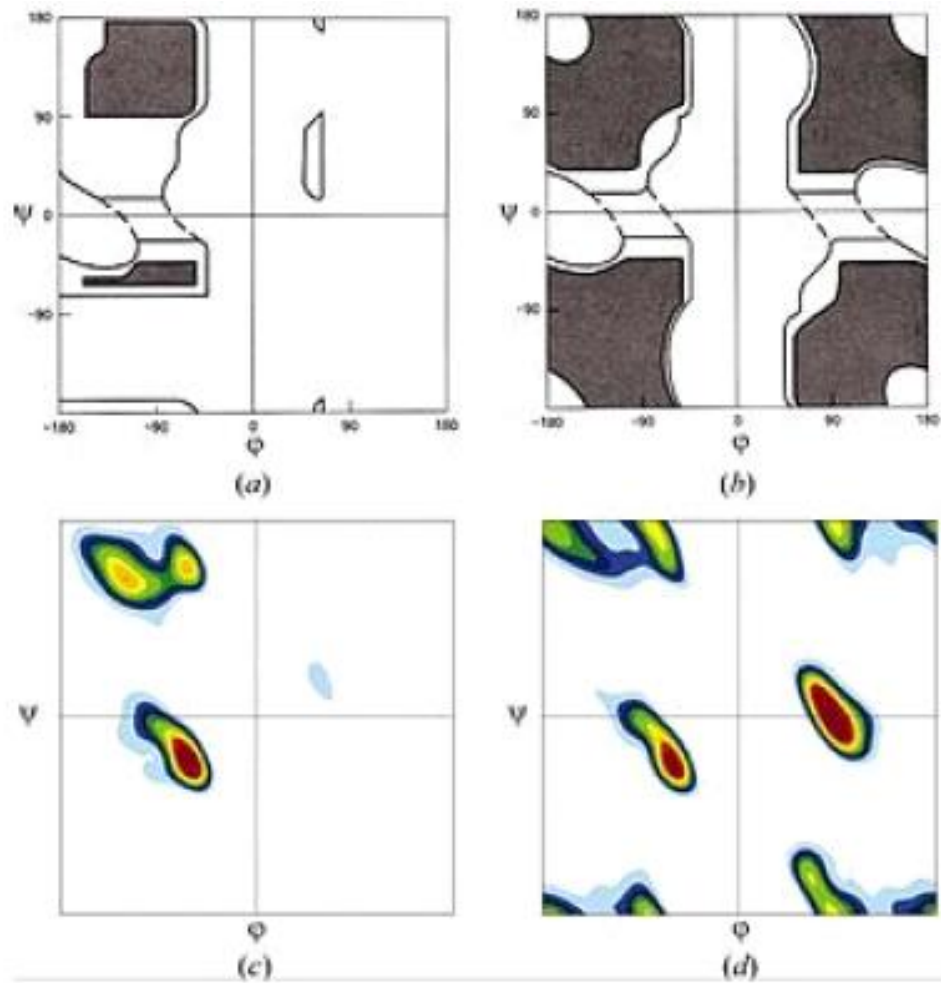
XRD

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-Ramachandran plot

A Ramachandran plot (also known as a Ramachandran map or a Ramachandran diagram or a $[\phi, \psi]$ plot), developed by Gopalasamudram Narayana Ramachandran and Viswanathan Sasisekharan is a way to visualize dihedral angles of amino acid residues in protein structure. It shows the possible conformations of ψ and ϕ angles for a polypeptide.

Also known as a Ramachandran map or a Ramachandran diagram or a $[\phi, \psi]$ plot). A peptidic bond has two degrees of freedom, the dihedral angles named ϕ and ψ by Ramachandran.



The classical version of the Ramachandran plot for (a) [alanine](#) and (b) glycine

According to Ramachandran & Sasisekharan (1968). The fully allowed regions are shaded; the partially allowed region is bounded by a solid line. The connecting regions enclosed by the dashed lines are permissible with slight flexibility of bond angles. These regions were arrived at by stereo-chemical modelling. Although some overall features of these plots are correct, the details differ from experimentally observed Ramachandran plots for (c) all 19 non-glycines and (d) glycine.

The most remarkable differences are that most regions show a 45 degree slope rather than being parallel to any of the axes. The region is split into two distinct maxima and the two most populated regions (red) for glycine seen in (d) were predicted to be permissible as shown in (b). There are five areas in the glycine plot; two with $\psi = 0$ and three with $\psi = 180$.