

# *X-Ray Crystal Structure Request Form*

Name \_\_\_\_\_ Date \_\_\_\_\_  
Advisor \_\_\_\_\_ Adv's Approval \_\_\_\_\_  
Room \_\_\_\_\_ Phone # \_\_\_\_\_ Sample ID \_\_\_\_\_

Structure determination to be preformed by:

\_\_\_\_\_ User \_\_\_\_\_  
\_\_\_\_\_ Dr. Eichhorn \_\_\_\_\_

Expected Molecular Formula \_\_\_\_\_  
Solvents Used for Crystallization \_\_\_\_\_  
Solvents Used in Synthesis \_\_\_\_\_

Known Properties \_\_\_\_\_ Air Sensitive \_\_\_\_\_ Water Sensitive \_\_\_\_\_ Light Sensitive  
\_\_\_\_\_ Solvent Loss Sensitive  
\_\_\_\_\_ Other (Describe) \_\_\_\_\_

Predicted Structure	Operator's Use Only
	Operators Name _____ Structure ID _____ Cell Parameters: a _____ $\alpha$ _____ b _____ $\beta$ _____ c _____ $\gamma$ _____ Crystal Dimensions _____ Space Group _____ Date Mounted _____ Date Collection Completed _____ Date Structure Completed _____ Temperature _____ Number of Data _____ Data $>4\sigma$ _____ $>3\sigma$ _____ $>2\sigma$ _____ Number of Parameters _____ Types of Absorption Correction: _____ _____ Min/Max Transmission _____ % Decay _____ R _____ $R_w$ _____ R(all) _____ $R_w$ (all) _____ g.o.f. _____ Min/Max Peak on Final Map _____ Comments: _____ _____ _____
Actual Structure	