广州市材料信息学重点实验室 Guangzhou Municipal Key Laboratory of Materials Informatics

X-Ray photoelectron Spectrograph WPEM fitting

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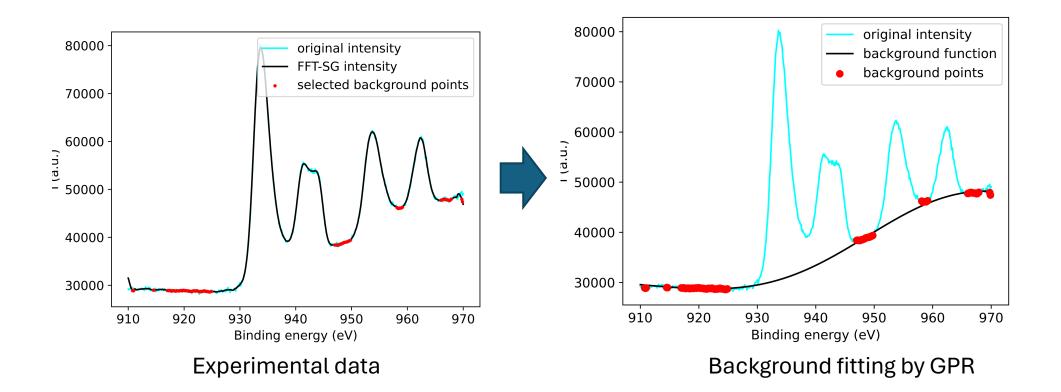
https://github.com/Bin-Cao (binjacobcao@gmail.com)



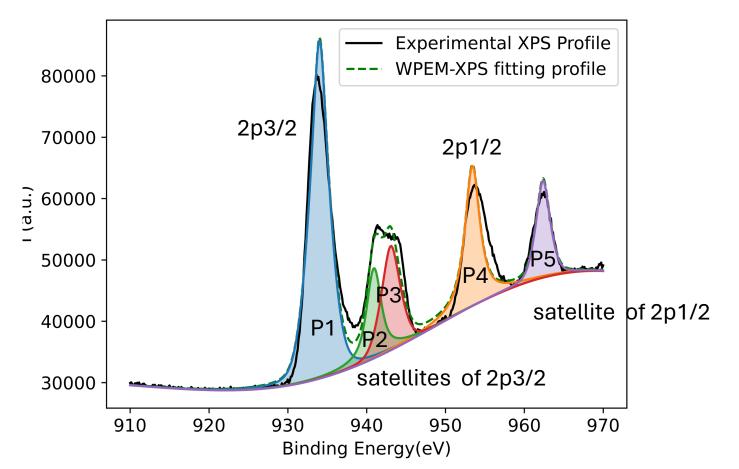
Cu (I): no satellites

Cu (II): one / two satellites

BIOINORGANIC CHEMISTRY 6, 45-59 (1976)



Peak decomposition

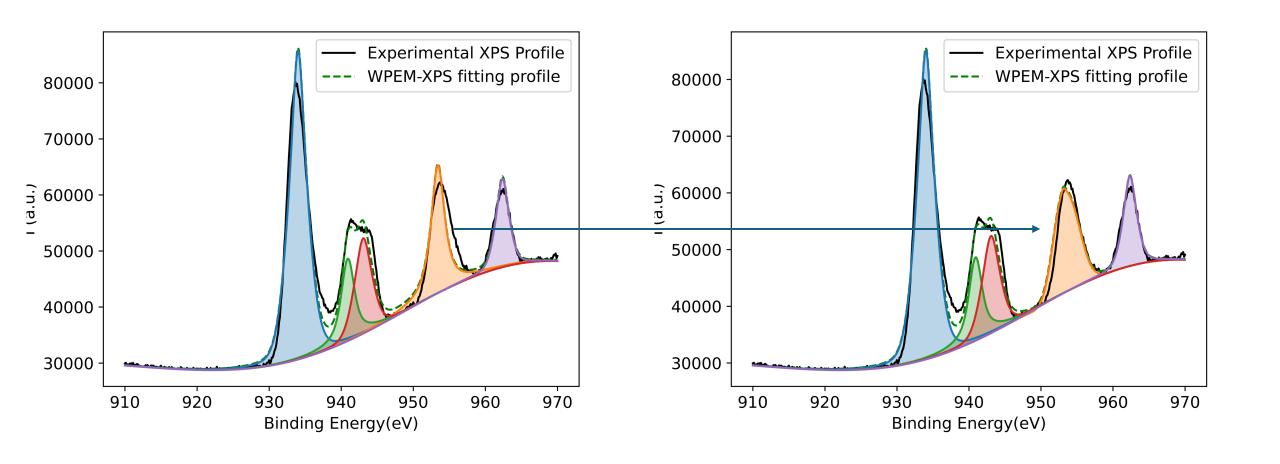


Integral area (P2+P3) / Integral area (P1) = Integral area (P5) / Integral area (P4)

Intensity P1 : P4 = 2 : 1

Energy gap follws central field approximation Zeff = 25.576

Asymetary peak shap



Asymetary factor = $w * (\Delta x)$



Thank you!

If you use the PPT, please quote it as follows:

Bin CAO. (2024). Bgolearn: A Bayesian global optimization package. Retrieved from https://github.com/Bin-Cao/Bgolearn



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