

Supporting Information

The Nature of Interactions and UV Induced Response within α -Zirconium Phosphate Intercalation Compounds with Azobenzenes

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Table S1. Selected interatomic distances in compound $Zr_{0.5}(HPO_4)C_{12}N_3H_{11}$

Atom1 (sym. code)	Atom2 (sym. code)	distance (Å)
Octahedral coordination		
Zr	O1	2.106(14)
Zr	O2 (-x,y+1/2,-z+1/2)	2.10(3)
Zr	O3 (-x,y-1/2,-z+1/2)	1.90(3)
Zr	O1(-x,-y+2,-z)	2.106(14)
Zr	O2 (x,-y+3/2,z-1/2)	2.10(3)
Zr	O3 (x,-y+5/2,z-1/2)	1.90(3)
Tetrahedral coordination		
P	O1	1.506(17)
P	O2	1.50(3)
P	O3	1.50(3)
P	O4	1.509(13)
Inorganic part		
C2	C3	1.39(6)
C3	C4	1.39(5)
C4	C5	1.39(3)
C5	C6	1.39(6)
C6	C7	1.39(4)
C7	C2	1.39(3)
C5	N8	1.45(8)
N8	N9	1.28(4)
N9	C10	1.46(4)
C10	C11	1.39(5)
C11	C12	1.39(5)
C12	C13	1.39(4)
C13	C14	1.39(6)
C14	C15	1.39(5)
C15	C10	1.39(4)
C13	N16	1.45(5)

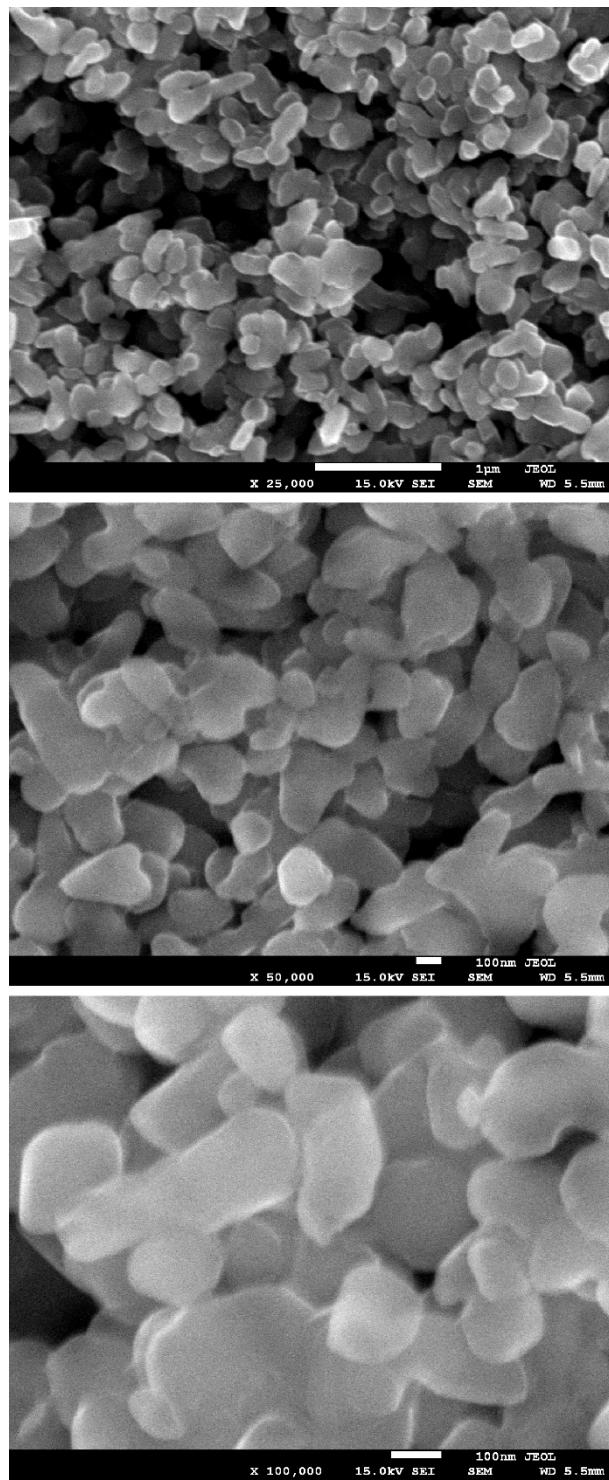


Figure S1. The SEM images of pure ZrP sample.

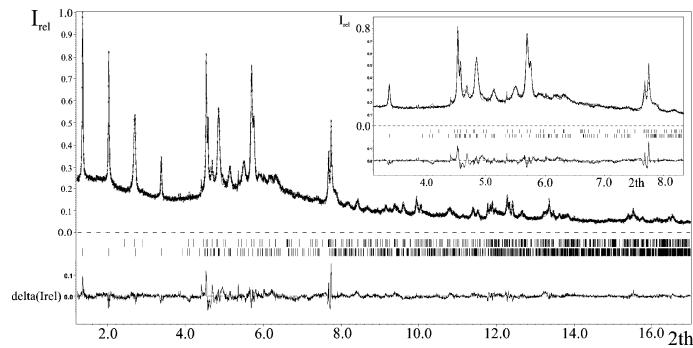


Figure S2. Synchrotron generated XRD pattern and the Rietveld plot for the ZpA sample.

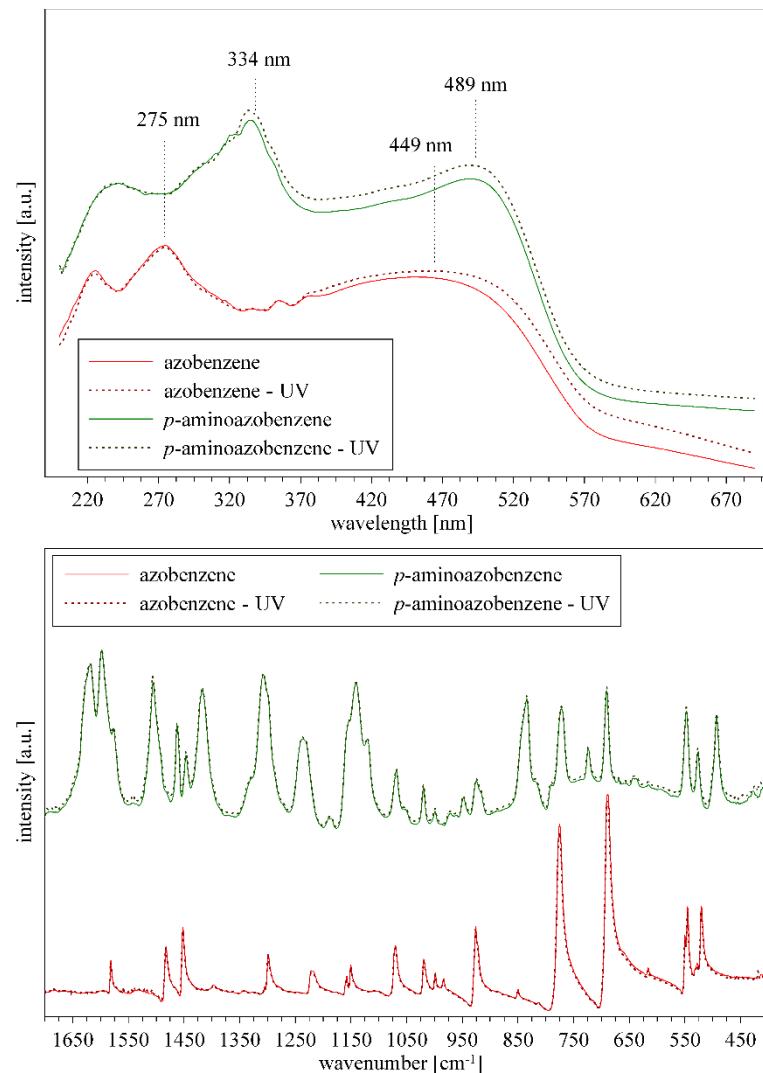


Figure S3. The UV-Vis (upper graph) and FTIR (lower graph) spectra of crystalline Az and pAz before and after UV irradiation.

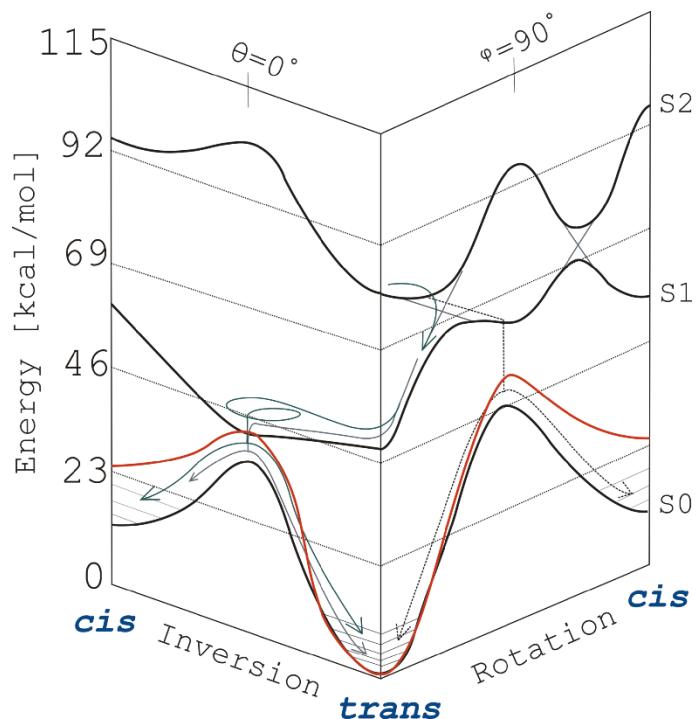


Figure S4. Scheme of the trans–cis isomerization process after $\pi\pi^*$ excitation.



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