

## Supplementary Information

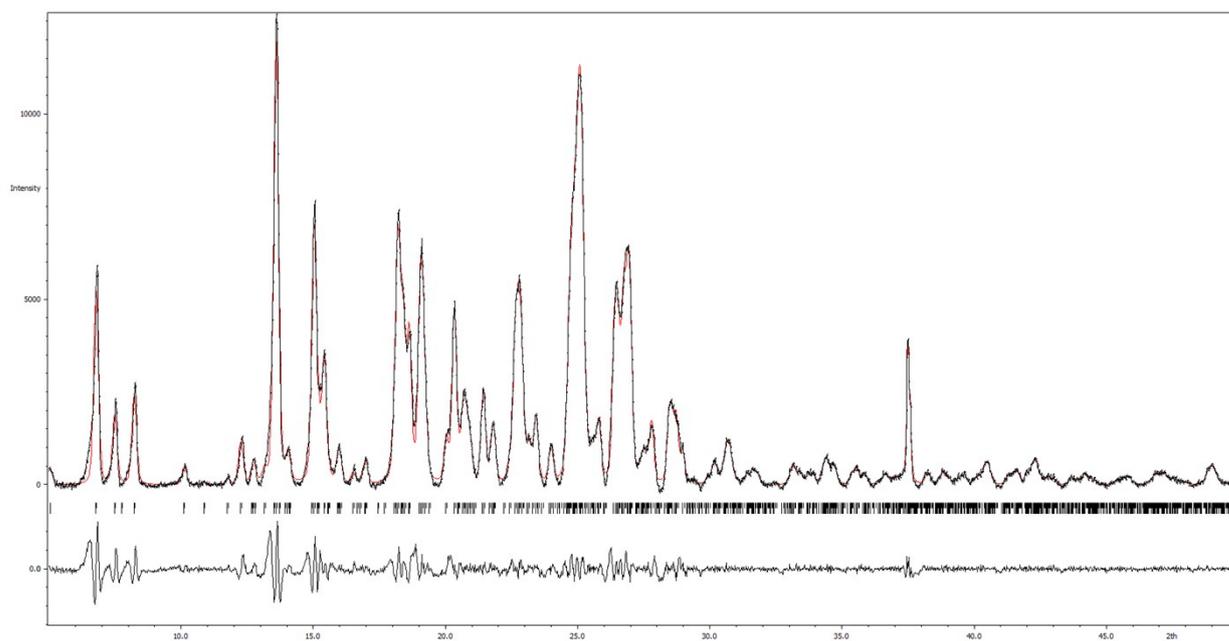
for the article “Study of influence of anion nature on crystal packing features and feasibility of [2 + 2] photocycloaddition reaction in protonated forms of dimethoxystyrylheterocycles”

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**Table S1.** Unit cell parameters in single crystal and powder of **14**.

Compound	<i>T</i> , K		<i>a</i> , Å	<i>b</i> , Å	<i>c</i> , Å	<i>α</i> , °	<i>β</i> , °	<i>γ</i> , °	<i>V</i> , Å <sup>3</sup>
<b>14</b>	298	Powder	7.2729(6)	25.9951(19)	23.7579(19)	90	97.486(6)	90	4454.2(7)
	150	Single-crystal	7.1809(5)	25.9269(18)	23.5772(16)	90	97.595(3)	90	4351.1(5)

PXRD experiment was conducted on a Haoyuan DX2700 diffractometer (CuK $\alpha$ ,  $\lambda = 1.5418$  Å, Ni-filter, reflection geometry,  $2\theta$  from 5 to 50°) equipped with a scintillator detector. The PXRD patterns were analyzed using JANA2006 [1]. For refinement, the SCXRD data were used as initial unit cell parameters.



**Figure S1.** PXRD pattern for compound **14** ( $T = 298$  K). Red line – the calculated fit superimposed with the experimental curve (black points). Difference curve is at the bottom of the figure.

[1] Petricek V., Dusek M., Palatinus L. Crystallographic Computing System JANA2006: General features // Zeitschrift für Kristallographie - Crystalline Materials. 2014. Vol. 229. No. 5. pp. 345-352.