Table A. Atomic Coordinates and Thermal Displacement Parameters (Å²) for α-FOX-7 at 200 K obtained from a Single Crystal Investigation.

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<th>atom</th>
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<th>z/c</th>
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<th>U22</th>
<th>U33</th>
<th>U12</th>
<th>U13</th>
<th>U23</th>
<th>Ueq</th>
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<td>0.75066</td>
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<th>U12</th>
<th>U13</th>
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<th>Ueq</th>
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Table C. Atomic Coordinates and Thermal Displacement Parameters (Å)\(^2\)
for α-FOX-7 at 333 K obtained from a Single Crystal Investigation.

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**Table D.** Donor (D)-Acceptor (A) Geometry (Å, °) for Hydrogen Bonds in α-FOX-7 at 200 K Obtained from a Single Crystal Investigation. Intramolecular Bonds are marked *.

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<th>H···A (Å)</th>
<th>D···A (Å)</th>
<th>D – H···A</th>
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D – H: ± 0.04 Å, H···A: ± 0.04 Å, D···A: ± 0.004 Å, ∠D – H···A: ± 2°
<table>
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<th>D – H ··· A</th>
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D – H: ± 0.04 Å, H···A: ± 0.04 Å, D···A: ± 0.004 Å, ∠D – H···A: ± 2°
Table F. Donor (D)-Acceptor (A) Geometry (Å, °) for Hydrogen Bonds in α-FOX-7 at 333 K obtained from a Single Crystal Investigation. Intramolecular Bonds are marked *.

<table>
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<td>150</td>
</tr>
<tr>
<td>N22-H22⋯O21*</td>
<td>0.86</td>
<td>2.02</td>
<td>2.622</td>
<td>127</td>
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<tr>
<td>N22-H22⋯O12</td>
<td>0.86</td>
<td>2.36</td>
<td>3.034</td>
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<tr>
<td>averaged</td>
<td>0.92</td>
<td>2.24</td>
<td>3.04</td>
<td>137</td>
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</tbody>
</table>

D – H: ± 0.04 Å, H⋯A: ± 0.04 Å, D⋯A: ± 0.004 Å, ∠D–H⋯A: ± 2°