**Figure 2\_source data 1.** Crystallographic data collection and refinement statistics.

|  |  |  |
| --- | --- | --- |
|  | **WT-1** | **Pat-1** |
|  |  |  |
| **Crystal parameters** |  |  |
| Space group | P6322 | P6122 |
| Cell constants | a = b = 104.5 Åc = 55.2 Å | a = b = 127.3 Åc= 81.8 Å |
| Subunits / AUa | 1 | 2 |
|  |  |  |
| **Data collection** |  |  |
| Beam line | ID-30, ESRF | ID-30, ESRF |
| Wavelength (Å) | 0.976 | 0.976 |
| Resolution range (Å)b | 50 – 1.55 (1.65 – 1.55) | 50 – 2.5 (2.6 – 2.5) |
| No. observed reflections | 162247 | 115993 |
| No. unique reflectionsc | 26220 | 13924 |
| Completeness (%)b | 99.7 (99.8) | 99.4 (100) |
| Rmerge (%)b, d | 4.3 (58.7) | 5.5 (49.6) |
| I/σ (I)b | 20.9 (3.2) | 20.6 (4.2) |
|  |  |  |
| **Refinement (REFMAC5)** |  |  |
| Resolution range (Å) | 15 – 1.55 | 15 – 2.5 |
| No. refl. working set | 24867 | 13148 |
| No. refl. test set | 1309 | 692 |
| No. non hydrogen | 1025 | 1653 |
| Solvent (H2O, ions, PEG) | 178 | 39 |
| Rwork / Rfree (%)e | 14.2 / 16.5 | 16.2 / 22.7 |
| r.m.s.d. bond (Å) / (°)f | 0.007 / 1.3 | 0.003 / 1.2 |
| Average B-factor (Å2) | 25.4 | 83.6 |
| Ramachandran Plot (%)g | 95.4 / 4.6 / 0 | 95.3 / 4.7 / 0 |
|  |  |  |
| PDB accession code | 6SM1 | 6SM2 |

[a] Asymmetric unit

[b] The values in parentheses for resolution range, completeness, Rmerge and I/σ (I) correspond to the highest resolution shell

[c] Data reduction was carried out with XDS and from a single crystal.

[d] Rmerge(I) = ΣhklΣj | I(hkl)j - <I(hkl)> | / Σhkl Σj I(hkl)j, where I(hkl)j is the jth measurement of the intensity of reflection hkl and <I(hkl)> is the average intensity

[e] R = Σhkl | |Fobs| - |Fcalc| |/Σhkl |Fobs|, where Rfree is calculated without a sigma cut off for a randomly chosen 5% of reflections, which were not used for structure refinement, and Rwork is calculated for the remaining reflections

[f] Deviations from ideal bond lengths/angles

[g] Percentage of residues in favored / allowed / outlier region