

Full wwPDB X-ray Structure Validation Report (i)

Oct 5, 2023 – 05:23 PM EDT

| PDB ID : | 6DEC |
|----------------|---|
| Title : | Crystal structure of Bos taurus $Arp2/3$ complex binding with C-terminus of |
| | Homo sapiens SPIN90 |
| Authors : | Nolen, B.J.; Luan, Q. |
| Deposited on : | 2018-05-11 |
| Resolution : | 4.60 Å(reported) |

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

| MolProbity | : | 4.02b-467 |
|--------------------------------|---|--|
| Mogul | : | 1.8.5 (274361), CSD as541be (2020) |
| Xtriage (Phenix) | : | 1.13 |
| EDS | : | 2.35.1 |
| buster-report | : | 1.1.7(2018) |
| Percentile statistics | : | 20191225.v01 (using entries in the PDB archive December 25th 2019) |
| Refmac | : | 5.8.0158 |
| CCP4 | : | 7.0.044 (Gargrove) |
| Ideal geometry (proteins) | : | Engh & Huber (2001) |
| Ideal geometry (DNA, RNA) | : | Parkinson et al. (1996) |
| Validation Pipeline (wwPDB-VP) | : | 2.35.1 |

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 4.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



| Motria | Whole archive | Similar resolution |
|-----------------------|---------------------|---|
| wietric | $(\# { m Entries})$ | $(\# { m Entries}, { m resolution} { m range}({ m \AA}))$ |
| R _{free} | 130704 | 1062 (5.40-3.80) |
| Clashscore | 141614 | 1130 (5.40-3.80) |
| Ramachandran outliers | 138981 | 1074 (5.40-3.80) |
| Sidechain outliers | 138945 | 1055 (5.40-3.80) |
| RSRZ outliers | 127900 | 1113 (5.50-3.70) |

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

| Mol | Chain | Length | Quality of chain | | |
|-----|-------|--------|------------------|------|------|
| 1 | А | 418 | 88% | 6% | • 5% |
| 1 | Н | 418 | 86% | 6% • | 8% |
| 2 | В | 394 | 87% | 7% | 6% |
| 2 | Ι | 394 | 50% •• 47% | | |
| 3 | С | 372 | 86% | ••• | 9% |



| Mol | Chain | Length | Quality of chain | | |
|-----|-------|--------|------------------|----|---------|
| 3 | J | 372 | 64% 5% • | | 31% |
| 4 | D | 300 | 90% | | • 7% |
| 4 | K | 300 | 90% | | • 6% |
| 5 | Е | 178 | 84% | | 6% • 8% |
| 5 | L | 178 | 75% | • | 22% |
| 6 | F | 168 | 96% | | |
| 6 | Ν | 168 | 93% | | |
| 7 | G | 151 | 74% | •• | 23% |
| 7 | 0 | 151 | % • 85% | | •• 13% |
| 8 | М | 455 | 69% | • | 28% |
| 8 | Р | 455 | 68% | • | 28% |
| 9 | Q | 6 | 100% | | |
| 10 | R | 9 | 100% | | |



2 Entry composition (i)

There are 12 unique types of molecules in this entry. The entry contains 20563 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

• Molecule 1 is a protein called Actin-related protein 3.

| Mol | Chain | Residues | | Ator | ns | | ZeroOcc | AltConf | Trace |
|-----|---------|----------|------|------|-----|-----|---------|---------|-------|
| 1 | 1 A 308 | Total | С | Ν | Ο | 0 | 0 | 0 | |
| | 390 | 2010 | 1202 | 404 | 404 | 0 | 0 | 0 | |
| 1 | Ц 206 | Total | С | Ν | Ο | 0 | 0 | 0 | |
| | 11 | 380 | 1948 | 1169 | 389 | 390 | 0 | | 0 |

• Molecule 2 is a protein called Actin-related protein 2.

| Mol | Chain | Residues | | Ator | ns | | ZeroOcc | AltConf | Trace |
|-----|---------|----------|-------|------|-----|-----|---------|---------|-------|
| 2 | 2 B 360 | 360 | Total | С | Ν | Ο | 0 | 0 | 0 |
| | D | 505 | 1850 | 1103 | 372 | 375 | 0 | 0 | 0 |
| 9 | т | I 209 | Total | С | Ν | 0 | 0 | 0 | 0 |
| Z | 1 | | 1030 | 612 | 209 | 209 | 0 | 0 | 0 |

• Molecule 3 is a protein called Actin-related protein 2/3 complex subunit 1B.

| Mol | Chain | Residues | | Ato | \mathbf{ms} | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|---------------|---|---------|---------|-------|
| 3 | С | 339 | Total | C | N | 0 | 0 | 0 | 0 |
| | | 1757 | 1070 | 347 | 340 | | | | |
| 3 | Т | I 959 | Total | С | Ν | Ο | 0 | 0 | 0 |
| 0 1 | 230 | 1274 | 758 | 258 | 258 | 0 | 0 | 0 | |

• Molecule 4 is a protein called Actin-related protein 2/3 complex subunit 2.

| Mol | Chain | Residues | | Ato | ms | | ZeroOcc | AltConf | Trace |
|-----|---------|----------|-------|-----|-----|-----|---------|---------|-------|
| 4 | 4 D 270 | Total | С | Ν | Ο | 0 | 0 | 0 | |
| 4 D | 219 | 1389 | 830 | 279 | 280 | 0 | 0 | | |
| 4 | 4 V 99 | 263 | Total | С | Ν | Ο | 0 | 0 | 0 |
| 4 N | Γ | 200 | 1429 | 855 | 288 | 286 | 0 | 0 | 0 |

• Molecule 5 is a protein called Actin-related protein 2/3 complex subunit 3.



| Mol | Chain | Residues | | Ato | ms | | ZeroOcc | AltConf | Trace |
|-----|---------|----------|-------|-----|-----|-----|---------|---------|-------|
| 5 | 5 F 162 | Total | С | Ν | Ο | 0 | 0 | 0 | |
| D E | 105 | 815 | 488 | 163 | 164 | | | | |
| 5 | E I 120 | 128 | Total | С | Ν | Ο | 0 | 0 | 0 |
| 5 Г | | 199 | 685 | 409 | 138 | 138 | 0 | 0 | 0 |

• Molecule 6 is a protein called Actin-related protein 2/3 complex subunit 4.

| Mol | Chain | Residues | | Ato | \mathbf{ms} | | ZeroOcc | AltConf | Trace |
|-----|---------|----------|-------|-----|---------------|-----|---------|---------|-------|
| 6 | 6 F 165 | Total | С | Ν | Ο | 0 | 0 | 0 | |
| | Ľ | 100 | 829 | 496 | 168 | 165 | 0 | 0 | 0 |
| 6 | 6 N | 163 | Total | С | Ν | Ο | 0 | 0 | 0 |
| 0 | 1 | | 825 | 496 | 166 | 163 | 0 | 0 | 0 |

• Molecule 7 is a protein called Actin-related protein 2/3 complex subunit 5.

| Mol | Chain | Residues | | Ato | ms | | ZeroOcc | AltConf | Trace |
|-----|---------|----------|-------|-----|-----|-----|---------|---------|-------|
| 7 | 7 C 116 | Total | С | Ν | Ο | 0 | 0 | 0 | |
| I G | 110 | 577 | 345 | 116 | 116 | 0 | 0 | 0 | |
| 7 | | 139 | Total | С | Ν | Ο | 0 | 0 | 0 |
| 1 | 0 | 152 | 648 | 384 | 132 | 132 | 0 | 0 | 0 |

• Molecule 8 is a protein called NCK-interacting protein with SH3 domain.

| Mol | Chain | Residues | Atoms | ZeroOcc | AltConf | Trace |
|-----|-------|----------|--|---------|---------|-------|
| 8 | М | 329 | Total C N O S 1638 979 329 329 1 | 0 | 0 | 0 |
| 8 | Р | 328 | Total C N O 1656 995 332 329 | 0 | 0 | 0 |

There are 2 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------------|------------|
| М | 268 | SER | - | expression tag | UNP Q9NZQ3 |
| Р | 268 | SER | - | expression tag | UNP Q9NZQ3 |

• Molecule 9 is a protein called unidentified.

| Mol | Chain | Residues | Atoms | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------------|---------|--------|--------|---------|---------|-------|
| 9 | Q | 6 | Total 30 | C 18 | N 6 | O 6 | 0 | 0 | 0 |

• Molecule 10 is a protein called unidentified.



| Mol | Chain | Residues | Atoms | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------------|---------|--------|--------|---------|---------|-------|
| 10 | R | 9 | Total 45 | С 27 | N 9 | O 9 | 0 | 0 | 0 |

• Molecule 11 is CALCIUM ION (three-letter code: CA) (formula: Ca).

| Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|-----------------|---------|---------|
| 11 | А | 1 | Total Ca 1 1 | 0 | 0 |
| 11 | В | 1 | Total Ca 1 1 | 0 | 0 |
| 11 | Н | 1 | Total Ca 1 1 | 0 | 0 |
| 11 | Ι | 1 | Total Ca 1 1 | 0 | 0 |

• Molecule 12 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



| Mol | Chain | Residues | Atoms | | | | ZeroOcc | AltConf | | | |
|-----|-------|----------|-------|----|-------|----|---------|---------|---|---|---|
| 19 | 19 A | 1 | Total | С | Ν | Ο | Р | 0 | 0 | | |
| 12 | Л | L | 31 | 10 | 5 | 13 | 3 | 0 | 0 | | |
| 19 | В | 1 | Total | С | Ν | Ο | Р | 0 | 0 | | |
| 12 | D | | 31 | 10 | 5 | 13 | 3 | 0 | 0 | | |
| 19 | п | Ц | Ц | 1 | Total | С | Ν | Ο | Р | 0 | 0 |
| 12 | 11 | 1 | 31 | 10 | 5 | 13 | 3 | 0 | 0 | | |
| 12 | Ι | I 1 | Total | С | Ν | Ο | Р | 0 | 0 | | |
| | | | 31 | 10 | 5 | 13 | 3 | 0 | 0 | | |



3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.



• Molecule 1: Actin-related protein 3





• Molecule 5: Actin-related protein 2/3 complex subunit 3 Chain E: 84% 6% 8% • Molecule 5: Actin-related protein 2/3 complex subunit 3 Chain L: 75% 22% MET PRO ALA ALA HIS SER SER SER SER SER ASP PRO ASP PRO ASP PRO LEU LEU • Molecule 6: Actin-related protein 2/3 complex subunit 4 Chain F: 96% • Molecule 6: Actin-related protein 2/3 complex subunit 4 Chain N: 93% • Molecule 7: Actin-related protein 2/3 complex subunit 5 Chain G: 74% 23% • Molecule 7: Actin-related protein 2/3 complex subunit 5 Chain O: 85% 13% • • MET SER LYS ASN ASN THR VAL SER SER • Molecule 8: NCK-interacting protein with SH3 domain Chain M: 69% 28%



| HIS GLU CTS CTS CTS CTS CTS CTS ALL ALL CTLE CTLE CTLE CTLE CTLE CTLE CTLE CT | SER SER SER VAL VAL VAL CALU VAL LEU VAL LEU VAL LEU VAL LEU VAL LEU VAL LEU VAL | SER MET ALLA LLEU PRO SER CGLY CGLN CGLN CGLN ASP ASP ASP ASP ASP ASP ASP ASP ASP ASP |
|--|---|---|
| P426 P426 P426 P519 P519 P543 P543 P546 P551 P551 P560 P600 | 0000 0000 0000 0000 0000 0000 0000 0000 0000 | VAL LEU GLY ALA PRO SER |
| • Molecule 8: NCK-interact | ing protein with SH3 dor | nain |
| Chain P: | 68% | •• 28% |
| SER PRO PRO PRO GLU GLU GLU GLU CLU GLU THR THR THR THR THR SER ALA SER ALA SER ALA SER CLU CLU CLU CLU CLU CLU CLU CLU CLU CLU | ALA ALA CLEU CLEU CLEU CLEU CLEU CLU CLU CLU CLU CLU CLU CLU CLU CLU CL | ALA VAL VAL THR THR THR TLE GLY VAL CLU VAL CLU VAL ARG ARG CLV VAL CLU VAL CLU VAL CLU VAL CLU VAL CLU VAL CLU VAC TEU VAC THR THR CLU VAC THR THR THR THR THR THR THR THR THR THR |
| HIS GLU LEU LEU LEU VAL VAL TILE TILE TILE TILE TILE TILE TILE TIL | SER SER PRO PRO MET MET CLU CLU CLU CLU CLU CLU CLU CLU CLU CLU | SER MET MET ALLA LEU PRO SER CILU VAL VAL ARG C376 AR3 AR3 AR3 AR3 AR3 AR40 GLN GLN GLN GLN GLN GLN AR5 AR5 AR5 AR6 AR6 AR6 AR6 AR6 AR6 AR6 AR6 AR6 AR6 |
| C412 TV412 LEU LEU C412 LEU C411 C411 C411 C411 C411 C411 C411 C41 | ASP 1546 1551 9551 9551 960 960 9611 9611 9611 9611 | P672 L5672 P661 P661 P661 GLY GLY GLY GLY SER |
| • Molecule 9: unidentified | | |
| Chain Q: | 100% | |
| There are no outlier residues | s recorded for this chain. | |
| • Molecule 10: unidentified | | |
| Chain R: | 100% | |

There are no outlier residues recorded for this chain.



4 Data and refinement statistics (i)

| Property | Value | Source |
|--|---|-----------|
| Space group | P 21 21 21 | Depositor |
| Cell constants | 179.84Å 197.38Å 202.07Å | Deperitor |
| a, b, c, α , β , γ | 90.00° 90.00° 90.00° | Depositor |
| $\mathbf{P}_{\text{assolution}}(\hat{\mathbf{A}})$ | 48.94 - 4.60 | Depositor |
| Resolution (A) | 48.94 - 4.47 | EDS |
| % Data completeness | 94.0 (48.94-4.60) | Depositor |
| (in resolution range) | 89.4(48.94-4.47) | EDS |
| R_{merge} | (Not available) | Depositor |
| R _{sym} | 0.12 | Depositor |
| $< I/\sigma(I) > 1$ | $1.67 (at 4.45 \text{\AA})$ | Xtriage |
| Refinement program | PHENIX 1.9_1692 | Depositor |
| D D. | 0.275 , 0.314 | Depositor |
| Π, Π_{free} | 0.295 , 0.320 | DCC |
| R_{free} test set | 1975 reflections (5.01%) | wwPDB-VP |
| Wilson B-factor $(Å^2)$ | 196.7 | Xtriage |
| Anisotropy | 0.329 | Xtriage |
| Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$ | 0.13, 311.7 | EDS |
| L-test for twinning ² | $< L >=0.46, < L^2>=0.29$ | Xtriage |
| Estimated twinning fraction | 0.025 for -h,l,k | Xtriage |
| F_o, F_c correlation | 0.86 | EDS |
| Total number of atoms | 20563 | wwPDB-VP |
| Average B, all atoms $(Å^2)$ | 179.0 | wwPDB-VP |

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 2.68% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: ATP, CA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with |Z| > 5 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

| Mal | Chain | Bond | lengths | E | Bond angles |
|-----|-------|------|----------|------|-------------------------------|
| | Unain | RMSZ | # Z > 5 | RMSZ | # Z > 5 |
| 1 | А | 0.35 | 0/2015 | 0.66 | 26/2799~(0.9%) |
| 1 | Н | 0.34 | 0/1947 | 0.66 | 26/2699~(1.0%) |
| 2 | В | 0.33 | 0/1851 | 0.57 | 17/2567~(0.7%) |
| 2 | Ι | 0.33 | 0/1020 | 0.59 | 11/1403~(0.8%) |
| 3 | С | 0.32 | 0/1772 | 0.54 | 14/2466~(0.6%) |
| 3 | J | 0.32 | 0/1250 | 0.58 | 12/1702~(0.7%) |
| 4 | D | 0.33 | 0/1387 | 0.50 | 9/1929~(0.5%) |
| 4 | К | 0.34 | 0/1432 | 0.50 | 9/1994~(0.5%) |
| 5 | Е | 0.32 | 0/810 | 0.64 | 10/1120~(0.9%) |
| 5 | L | 0.32 | 0/682 | 0.53 | 5/947~(0.5%) |
| 6 | F | 0.36 | 0/828 | 0.49 | 4/1155~(0.3%) |
| 6 | Ν | 0.35 | 0/824 | 0.48 | 4/1147~(0.3%) |
| 7 | G | 0.33 | 0/577 | 0.53 | 4/800~(0.5%) |
| 7 | 0 | 0.32 | 0/645 | 0.50 | 4/891~(0.4%) |
| 8 | М | 0.32 | 0/1635 | 0.58 | 16/2279~(0.7%) |
| 8 | Р | 0.33 | 0/1656 | 0.60 | $16\overline{/2305}\ (0.7\%)$ |
| All | All | 0.33 | 0/20331 | 0.57 | 187/28203~(0.7%) |

There are no bond length outliers.

All (187) bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | $\mathbf{Observed}(^{o})$ | $Ideal(^{o})$ |
|-----|-------|-----|------|---------|------|---------------------------|---------------|
| 8 | М | 566 | PRO | N-CA-CB | 6.33 | 110.89 | 103.30 |
| 8 | Р | 608 | PRO | N-CA-CB | 6.31 | 110.87 | 103.30 |
| 8 | Р | 566 | PRO | N-CA-CB | 6.26 | 110.81 | 103.30 |
| 8 | Р | 551 | PRO | N-CA-CB | 6.24 | 110.79 | 103.30 |
| 8 | Р | 681 | PRO | N-CA-CB | 6.20 | 110.74 | 103.30 |
| 3 | С | 106 | PRO | N-CA-CB | 6.20 | 110.74 | 103.30 |
| 1 | А | 295 | PRO | N-CA-CB | 6.18 | 110.72 | 103.30 |
| 1 | Н | 6 | PRO | N-CA-CB | 6.17 | 110.71 | 103.30 |



| Mol | Chain | Res | Type | Atoms | Z | Observed(^o) | Ideal(°) |
|-----|-------|-----|------|---------|------|--------------------------|----------|
| 1 | А | 279 | PRO | N-CA-CB | 6.15 | 110.68 | 103.30 |
| 5 | L | 31 | PRO | N-CA-CB | 6.14 | 110.67 | 103.30 |
| 1 | А | 120 | PRO | N-CA-CB | 6.13 | 110.66 | 103.30 |
| 1 | Н | 217 | PRO | N-CA-CB | 6.13 | 110.66 | 103.30 |
| 1 | А | 216 | PRO | N-CA-CB | 6.11 | 110.63 | 103.30 |
| 3 | J | 246 | PRO | N-CA-CB | 6.10 | 110.62 | 103.30 |
| 6 | F | 108 | PRO | N-CA-CB | 6.10 | 110.62 | 103.30 |
| 5 | L | 33 | PRO | N-CA-CB | 6.10 | 110.62 | 103.30 |
| 2 | В | 262 | PRO | N-CA-CB | 6.10 | 110.62 | 103.30 |
| 6 | N | 36 | PRO | N-CA-CB | 6.09 | 110.60 | 103.30 |
| 1 | А | 27 | PRO | N-CA-CB | 6.08 | 110.60 | 103.30 |
| 8 | М | 608 | PRO | N-CA-CB | 6.08 | 110.60 | 103.30 |
| 1 | Н | 32 | PRO | N-CA-CB | 6.08 | 110.59 | 103.30 |
| 4 | K | 206 | PRO | N-CA-CB | 6.07 | 110.59 | 103.30 |
| 1 | Н | 137 | PRO | N-CA-CB | 6.07 | 110.59 | 103.30 |
| 1 | Н | 295 | PRO | N-CA-CB | 6.07 | 110.58 | 103.30 |
| 1 | А | 290 | PRO | N-CA-CB | 6.06 | 110.57 | 103.30 |
| 1 | А | 179 | PRO | N-CA-CB | 6.05 | 110.56 | 103.30 |
| 1 | Н | 405 | PRO | N-CA-CB | 6.05 | 110.56 | 103.30 |
| 8 | Р | 529 | PRO | N-CA-CB | 6.04 | 110.55 | 103.30 |
| 3 | С | 184 | PRO | N-CA-CB | 6.04 | 110.54 | 103.30 |
| 7 | 0 | 62 | PRO | N-CA-CB | 6.03 | 110.54 | 103.30 |
| 5 | Е | 11 | PRO | N-CA-CB | 6.03 | 110.54 | 103.30 |
| 3 | С | 182 | PRO | N-CA-CB | 6.02 | 110.53 | 103.30 |
| 1 | Н | 179 | PRO | N-CA-CB | 6.02 | 110.53 | 103.30 |
| 1 | А | 115 | PRO | N-CA-CB | 6.02 | 110.53 | 103.30 |
| 8 | Р | 652 | PRO | N-CA-CB | 6.02 | 110.52 | 103.30 |
| 1 | А | 388 | PRO | N-CA-CB | 6.01 | 110.51 | 103.30 |
| 7 | G | 117 | PRO | N-CA-CB | 6.00 | 110.50 | 103.30 |
| 2 | В | 348 | PRO | N-CA-CB | 6.00 | 110.50 | 103.30 |
| 2 | В | 33 | PRO | N-CA-CB | 6.00 | 110.50 | 103.30 |
| 8 | М | 529 | PRO | N-CA-CB | 5.99 | 110.49 | 103.30 |
| 2 | Ι | 116 | PRO | N-CA-CB | 5.99 | 110.49 | 103.30 |
| 8 | Р | 484 | PRO | N-CA-CB | 5.99 | 110.49 | 103.30 |
| 3 | J | 10 | PRO | N-CA-CB | 5.99 | 110.48 | 103.30 |
| 1 | Н | 115 | PRO | N-CA-CB | 5.99 | 110.48 | 103.30 |
| 1 | А | 362 | PRO | N-CA-CB | 5.98 | 110.47 | 103.30 |
| 1 | Н | 290 | PRO | N-CA-CB | 5.98 | 110.47 | 103.30 |
| 6 | F | 50 | PRO | N-CA-CB | 5.97 | 110.47 | 103.30 |
| 7 | G | 63 | PRO | N-CA-CB | 5.97 | 110.47 | 103.30 |
| 1 | А | 308 | PRO | N-CA-CB | 5.97 | 110.47 | 103.30 |
| 5 | Е | 33 | PRO | N-CA-CB | 5.97 | 110.46 | 103.30 |



| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | $Ideal(^{o})$ |
|-----|-------|-----|------|---------|------|-------------|---------------|
| 4 | Κ | 103 | PRO | N-CA-CB | 5.96 | 110.46 | 103.30 |
| 2 | В | 28 | PRO | N-CA-CB | 5.96 | 110.45 | 103.30 |
| 3 | С | 288 | PRO | N-CA-CB | 5.96 | 110.45 | 103.30 |
| 7 | 0 | 37 | PRO | N-CA-CB | 5.96 | 110.45 | 103.30 |
| 7 | G | 62 | PRO | N-CA-CB | 5.96 | 110.45 | 103.30 |
| 5 | Е | 156 | PRO | N-CA-CB | 5.96 | 110.45 | 103.30 |
| 8 | М | 652 | PRO | N-CA-CB | 5.96 | 110.45 | 103.30 |
| 1 | А | 116 | PRO | N-CA-CB | 5.96 | 110.45 | 103.30 |
| 5 | Е | 2 | PRO | N-CA-CB | 5.94 | 110.43 | 103.30 |
| 8 | Р | 610 | PRO | N-CA-CB | 5.94 | 110.43 | 103.30 |
| 4 | D | 87 | PRO | N-CA-CB | 5.94 | 110.43 | 103.30 |
| 8 | М | 519 | PRO | N-CA-CB | 5.94 | 110.43 | 103.30 |
| 2 | В | 116 | PRO | N-CA-CB | 5.93 | 110.42 | 103.30 |
| 2 | В | 168 | PRO | N-CA-CB | 5.93 | 110.42 | 103.30 |
| 8 | М | 551 | PRO | N-CA-CB | 5.93 | 110.42 | 103.30 |
| 2 | Ι | 347 | PRO | N-CA-CB | 5.93 | 110.42 | 103.30 |
| 3 | J | 190 | PRO | N-CA-CB | 5.93 | 110.42 | 103.30 |
| 1 | Н | 412 | PRO | N-CA-CB | 5.93 | 110.41 | 103.30 |
| 8 | Р | 672 | PRO | N-CA-CB | 5.93 | 110.41 | 103.30 |
| 2 | Ι | 314 | PRO | N-CA-CB | 5.92 | 110.41 | 103.30 |
| 3 | J | 27 | PRO | N-CA-CB | 5.92 | 110.41 | 103.30 |
| 5 | Е | 112 | PRO | N-CA-CB | 5.92 | 110.40 | 103.30 |
| 3 | J | 88 | PRO | N-CA-CB | 5.91 | 110.40 | 103.30 |
| 3 | J | 184 | PRO | N-CA-CB | 5.91 | 110.39 | 103.30 |
| 6 | F | 7 | PRO | N-CA-CB | 5.91 | 110.39 | 103.30 |
| 3 | J | 106 | PRO | N-CA-CB | 5.91 | 110.39 | 103.30 |
| 2 | В | 247 | PRO | N-CA-CB | 5.90 | 110.38 | 103.30 |
| 7 | 0 | 117 | PRO | N-CA-CB | 5.90 | 110.38 | 103.30 |
| 3 | J | 182 | PRO | N-CA-CB | 5.90 | 110.38 | 103.30 |
| 8 | Р | 698 | PRO | N-CA-CB | 5.90 | 110.38 | 103.30 |
| 4 | D | 205 | PRO | N-CA-CB | 5.89 | 110.38 | 103.30 |
| 2 | В | 112 | PRO | N-CA-CB | 5.89 | 110.36 | 103.30 |
| 2 | В | 176 | PRO | N-CA-CB | 5.88 | 110.35 | 103.30 |
| 5 | L | 156 | PRO | N-CA-CB | 5.88 | 110.35 | 103.30 |
| 2 | В | 113 | PRO | N-CA-CB | 5.88 | 110.35 | 103.30 |
| 4 | D | 196 | PRO | N-CA-CB | 5.88 | 110.35 | 103.30 |
| 8 | P | 519 | PRO | N-CA-CB | 5.88 | 110.35 | 103.30 |
| 8 | Р | 611 | PRO | N-CA-CB | 5.88 | 110.35 | 103.30 |
| 5 | L | 23 | PRO | N-CA-CB | 5.87 | 110.35 | 103.30 |
| 1 | A | 412 | PRO | N-CA-CB | 5.87 | 110.34 | 103.30 |
| 2 | В | 73 | PRO | N-CA-CB | 5.87 | 110.35 | 103.30 |
| 3 | С | 88 | PRO | N-CA-CB | 5.87 | 110.34 | 103.30 |

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| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | $Ideal(^{o})$ |
|-----|-------|-----|------|---------|------|-------------|---------------|
| 1 | Н | 27 | PRO | N-CA-CB | 5.87 | 110.35 | 103.30 |
| 1 | А | 314 | PRO | N-CA-CB | 5.87 | 110.34 | 103.30 |
| 5 | Е | 109 | PRO | N-CA-CB | 5.87 | 110.34 | 103.30 |
| 3 | J | 180 | PRO | N-CA-CB | 5.86 | 110.33 | 103.30 |
| 1 | А | 32 | PRO | N-CA-CB | 5.86 | 110.33 | 103.30 |
| 1 | Н | 360 | PRO | N-CA-CB | 5.86 | 110.33 | 103.30 |
| 3 | С | 152 | PRO | N-CA-CB | 5.85 | 110.33 | 103.30 |
| 6 | Ν | 7 | PRO | N-CA-CB | 5.85 | 110.32 | 103.30 |
| 3 | J | 288 | PRO | N-CA-CB | 5.85 | 110.32 | 103.30 |
| 2 | В | 39 | PRO | N-CA-CB | 5.84 | 110.31 | 103.30 |
| 2 | В | 311 | PRO | N-CA-CB | 5.83 | 110.30 | 103.30 |
| 1 | Н | 314 | PRO | N-CA-CB | 5.83 | 110.30 | 103.30 |
| 6 | Ν | 108 | PRO | N-CA-CB | 5.83 | 110.30 | 103.30 |
| 4 | D | 206 | PRO | N-CA-CB | 5.83 | 110.29 | 103.30 |
| 1 | Н | 388 | PRO | N-CA-CB | 5.82 | 110.29 | 103.30 |
| 2 | Ι | 262 | PRO | N-CA-CB | 5.82 | 110.29 | 103.30 |
| 5 | L | 151 | PRO | N-CA-CB | 5.82 | 110.28 | 103.30 |
| 1 | А | 360 | PRO | N-CA-CB | 5.82 | 110.28 | 103.30 |
| 4 | D | 103 | PRO | N-CA-CB | 5.82 | 110.28 | 103.30 |
| 8 | М | 610 | PRO | N-CA-CB | 5.82 | 110.28 | 103.30 |
| 2 | Ι | 348 | PRO | N-CA-CB | 5.81 | 110.28 | 103.30 |
| 8 | Р | 543 | PRO | N-CA-CB | 5.81 | 110.27 | 103.30 |
| 4 | D | 28 | PRO | N-CA-CB | 5.81 | 110.27 | 103.30 |
| 1 | Н | 279 | PRO | N-CA-CB | 5.81 | 110.27 | 103.30 |
| 8 | Р | 600 | PRO | N-CA-CB | 5.81 | 110.27 | 103.30 |
| 6 | F | 36 | PRO | N-CA-CB | 5.81 | 110.27 | 103.30 |
| 4 | K | 196 | PRO | N-CA-CB | 5.80 | 110.27 | 103.30 |
| 5 | Е | 151 | PRO | N-CA-CB | 5.80 | 110.26 | 103.30 |
| 3 | С | 190 | PRO | N-CA-CB | 5.80 | 110.26 | 103.30 |
| 1 | Н | 285 | PRO | N-CA-CB | 5.80 | 110.26 | 103.30 |
| 3 | С | 180 | PRO | N-CA-CB | 5.79 | 110.25 | 103.30 |
| 1 | Н | 120 | PRO | N-CA-CB | 5.79 | 110.25 | 103.30 |
| 7 | 0 | 63 | PRO | N-CA-CB | 5.79 | 110.25 | 103.30 |
| 3 | J | 267 | PRO | N-CA-CB | 5.79 | 110.25 | 103.30 |
| 8 | М | 672 | PRO | N-CA-CB | 5.79 | 110.24 | 103.30 |
| 1 | Н | 216 | PRO | N-CA-CB | 5.79 | 110.24 | 103.30 |
| 1 | А | 6 | PRO | N-CA-CB | 5.78 | 110.24 | 103.30 |
| 2 | Ι | 247 | PRO | N-CA-CB | 5.78 | 110.24 | 103.30 |
| 7 | G | 37 | PRO | N-CA-CB | 5.78 | 110.23 | 103.30 |
| 2 | Ι | 311 | PRO | N-CA-CB | 5.78 | 110.23 | 103.30 |
| 4 | D | 282 | PRO | N-CA-CB | 5.77 | 110.23 | 103.30 |
| 3 | J | 62 | PRO | N-CA-CB | 5.77 | 110.23 | 103.30 |

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| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | $Ideal(^{o})$ |
|-----|-------|-----|------|---------|------|-------------|---------------|
| 4 | Κ | 28 | PRO | N-CA-CB | 5.77 | 110.22 | 103.30 |
| 4 | K | 48 | PRO | N-CA-CB | 5.77 | 110.22 | 103.30 |
| 4 | K | 205 | PRO | N-CA-CB | 5.77 | 110.22 | 103.30 |
| 1 | А | 217 | PRO | N-CA-CB | 5.76 | 110.22 | 103.30 |
| 5 | Е | 31 | PRO | N-CA-CB | 5.76 | 110.22 | 103.30 |
| 1 | А | 137 | PRO | N-CA-CB | 5.76 | 110.22 | 103.30 |
| 1 | Н | 362 | PRO | N-CA-CB | 5.76 | 110.21 | 103.30 |
| 2 | Ι | 112 | PRO | N-CA-CB | 5.76 | 110.21 | 103.30 |
| 5 | Е | 104 | PRO | N-CA-CB | 5.76 | 110.21 | 103.30 |
| 5 | Е | 106 | PRO | N-CA-CB | 5.75 | 110.21 | 103.30 |
| 1 | А | 236 | PRO | N-CA-CB | 5.75 | 110.20 | 103.30 |
| 6 | Ν | 50 | PRO | N-CA-CB | 5.75 | 110.20 | 103.30 |
| 2 | Ι | 176 | PRO | N-CA-CB | 5.74 | 110.19 | 103.30 |
| 8 | М | 426 | PRO | N-CA-CB | 5.74 | 110.19 | 103.30 |
| 2 | Ι | 268 | PRO | N-CA-CB | 5.74 | 110.18 | 103.30 |
| 1 | А | 285 | PRO | N-CA-CB | 5.73 | 110.18 | 103.30 |
| 8 | М | 484 | PRO | N-CA-CB | 5.73 | 110.18 | 103.30 |
| 1 | Н | 236 | PRO | N-CA-CB | 5.73 | 110.17 | 103.30 |
| 3 | С | 27 | PRO | N-CA-CB | 5.72 | 110.17 | 103.30 |
| 1 | А | 70 | PRO | N-CA-CB | 5.72 | 110.17 | 103.30 |
| 3 | С | 140 | PRO | N-CA-CB | 5.72 | 110.16 | 103.30 |
| 4 | Κ | 229 | PRO | N-CA-CB | 5.72 | 110.16 | 103.30 |
| 2 | В | 347 | PRO | N-CA-CB | 5.71 | 110.16 | 103.30 |
| 3 | С | 246 | PRO | N-CA-CB | 5.71 | 110.16 | 103.30 |
| 3 | С | 267 | PRO | N-CA-CB | 5.70 | 110.14 | 103.30 |
| 8 | Р | 625 | PRO | N-CA-CB | 5.70 | 110.14 | 103.30 |
| 8 | М | 611 | PRO | N-CA-CB | 5.70 | 110.13 | 103.30 |
| 1 | А | 77 | PRO | N-CA-CB | 5.69 | 110.13 | 103.30 |
| 8 | М | 600 | PRO | N-CA-CB | 5.67 | 110.10 | 103.30 |
| 1 | Н | 105 | PRO | N-CA-CB | 5.66 | 110.09 | 103.30 |
| 8 | М | 625 | PRO | N-CA-CB | 5.65 | 110.08 | 103.30 |
| 1 | Н | 70 | PRO | N-CA-CB | 5.65 | 110.08 | 103.30 |
| 3 | С | 10 | PRO | N-CA-CB | 5.65 | 110.08 | 103.30 |
| 1 | Н | 116 | PRO | N-CA-CB | 5.64 | 110.07 | 103.30 |
| 1 | А | 105 | PRO | N-CA-CB | 5.64 | 110.07 | 103.30 |
| 2 | В | 95 | PRO | N-CA-CB | 5.63 | 110.06 | 103.30 |
| 2 | Ι | 113 | PRO | N-CA-CB | 5.62 | 110.05 | 103.30 |
| 2 | В | 314 | PRO | N-CA-CB | 5.61 | 110.03 | 103.30 |
| 8 | Р | 426 | PRO | N-CA-CB | 5.60 | 110.02 | 103.30 |
| 2 | В | 268 | PRO | N-CA-CB | 5.57 | 109.99 | 103.30 |
| 4 | D | 48 | PRO | N-CA-CB | 5.57 | 109.98 | 103.30 |
| 1 | Н | 308 | PRO | N-CA-CB | 5.57 | 109.98 | 103.30 |

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| Mol | Chain | Res | Type | Atoms | Z | $Observed(^{o})$ | $Ideal(^{o})$ |
|-----|-------|-----|------|---------|------|------------------|---------------|
| 8 | М | 543 | PRO | N-CA-CB | 5.57 | 109.98 | 103.30 |
| 4 | K | 87 | PRO | N-CA-CB | 5.54 | 109.95 | 103.30 |
| 8 | М | 698 | PRO | N-CA-CB | 5.54 | 109.94 | 103.30 |
| 4 | K | 282 | PRO | N-CA-CB | 5.53 | 109.94 | 103.30 |
| 4 | D | 229 | PRO | N-CA-CB | 5.53 | 109.94 | 103.30 |
| 1 | Н | 194 | PRO | N-CA-CB | 5.51 | 109.92 | 103.30 |
| 3 | С | 62 | PRO | N-CA-CB | 5.46 | 109.85 | 103.30 |
| 1 | Н | 77 | PRO | N-CA-CB | 5.45 | 109.84 | 103.30 |
| 1 | А | 405 | PRO | N-CA-CB | 5.43 | 109.82 | 103.30 |
| 1 | A | 194 | PRO | N-CA-CB | 5.36 | 109.73 | 103.30 |
| 8 | М | 681 | PRO | N-CA-CB | 5.36 | 109.73 | 103.30 |

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There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 1 | А | 2010 | 0 | 906 | 2 | 0 |
| 1 | Н | 1948 | 0 | 873 | 2 | 0 |
| 2 | В | 1850 | 0 | 839 | 4 | 0 |
| 2 | Ι | 1030 | 0 | 449 | 1 | 0 |
| 3 | С | 1757 | 0 | 843 | 3 | 0 |
| 3 | J | 1274 | 0 | 551 | 5 | 0 |
| 4 | D | 1389 | 0 | 632 | 0 | 0 |
| 4 | K | 1429 | 0 | 663 | 3 | 0 |
| 5 | Е | 815 | 0 | 349 | 0 | 0 |
| 5 | L | 685 | 0 | 291 | 0 | 0 |
| 6 | F | 829 | 0 | 371 | 0 | 0 |
| 6 | N | 825 | 0 | 370 | 1 | 0 |
| 7 | G | 577 | 0 | 282 | 0 | 0 |
| 7 | 0 | 648 | 0 | 311 | 0 | 0 |
| 8 | М | 1638 | 0 | 737 | 0 | 0 |
| 8 | Р | 1656 | 0 | 760 | 1 | 0 |
| 9 | Q | 30 | 0 | 8 | 0 | 0 |
| 10 | R | 45 | 0 | 11 | 0 | 0 |



| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 11 | А | 1 | 0 | 0 | 0 | 0 |
| 11 | В | 1 | 0 | 0 | 0 | 0 |
| 11 | Н | 1 | 0 | 0 | 0 | 0 |
| 11 | Ι | 1 | 0 | 0 | 0 | 0 |
| 12 | А | 31 | 0 | 12 | 1 | 0 |
| 12 | В | 31 | 0 | 12 | 0 | 0 |
| 12 | Н | 31 | 0 | 12 | 0 | 0 |
| 12 | Ι | 31 | 0 | 12 | 0 | 0 |
| All | All | 20563 | 0 | 9294 | 20 | 0 |

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 1.

All (20) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

| Atom 1 | Atom 2 | Interatomic | Clash | |
|-----------------|------------------|--------------|-------------|--|
| Atom-1 | Atom-2 | distance (Å) | overlap (Å) | |
| 3:J:224:SER:O | 3:J:242:SER:N | 2.30 | 0.65 | |
| 4:K:2:ILE:O | 4:K:254:HIS:NE2 | 2.26 | 0.60 | |
| 2:B:222:TYR:CZ | 2:B:259:PHE:HB3 | 2.42 | 0.54 | |
| 2:B:241:VAL:HA | 2:B:255:GLY:HA2 | 1.90 | 0.53 | |
| 4:K:250:TYR:OH | 6:N:160:VAL:O | 2.24 | 0.52 | |
| 3:C:247:LEU:HA | 3:C:262:GLY:HA3 | 1.92 | 0.51 | |
| 3:J:270:PHE:HA | 3:J:282:GLY:H | 1.75 | 0.51 | |
| 3:C:261:ALA:HB2 | 3:C:267:PRO:HA | 1.93 | 0.49 | |
| 3:J:125:PHE:HA | 3:J:132:TRP:HA | 1.95 | 0.49 | |
| 3:J:79:TRP:HA | 3:J:88:PRO:HA | 1.96 | 0.47 | |
| 3:C:270:PHE:HA | 3:C:282:GLY:H | 1.81 | 0.46 | |
| 8:P:527:GLY:O | 8:P:530:PHE:N | 2.48 | 0.45 | |
| 1:A:169:ASP:HA | 1:A:322:SER:O | 2.18 | 0.44 | |
| 1:H:21:TYR:CE1 | 4:K:36:ALA:HB3 | 2.52 | 0.44 | |
| 1:A:171:GLY:HA3 | 12:A:502:ATP:O1G | 2.19 | 0.43 | |
| 2:B:75:GLU:N | 2:B:78:ILE:O | 2.51 | 0.43 | |
| 3:J:69:THR:O | 3:J:77:TYR:N | 2.51 | 0.43 | |
| 2:B:222:TYR:CE1 | 2:B:259:PHE:HB3 | 2.56 | 0.41 | |
| 1:H:21:TYR:OH | 1:H:103:ALA:HB2 | 2.21 | 0.41 | |
| 2:I:185:ALA:O | 2:I:189:ILE:N | 2.42 | 0.40 | |

There are no symmetry-related clashes.



5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

| Mol | Chain | Analysed | Favoured | Allowed | Outliers | Perce | ntiles |
|-----|--------------|-----------------|------------|----------|----------|-------|--------|
| 1 | А | 392/418~(94%) | 374 (95%) | 13 (3%) | 5 (1%) | 12 | 48 |
| 1 | Η | 372/418~(89%) | 355~(95%) | 13~(4%) | 4 (1%) | 14 | 52 |
| 2 | В | 365/394~(93%) | 337~(92%) | 25~(7%) | 3~(1%) | 19 | 60 |
| 2 | Ι | 189/394~(48%) | 180 (95%) | 7 (4%) | 2(1%) | 14 | 52 |
| 3 | С | 333/372~(90%) | 305~(92%) | 24 (7%) | 4 (1%) | 13 | 50 |
| 3 | J | 210/372~(56%) | 202 (96%) | 7 (3%) | 1 (0%) | 29 | 68 |
| 4 | D | 273/300~(91%) | 268~(98%) | 5 (2%) | 0 | 100 | 100 |
| 4 | Κ | 281/300~(94%) | 275~(98%) | 6(2%) | 0 | 100 | 100 |
| 5 | Ε | 151/178~(85%) | 141 (93%) | 4(3%) | 6 (4%) | 3 | 26 |
| 5 | L | 132/178~(74%) | 131 (99%) | 1 (1%) | 0 | 100 | 100 |
| 6 | \mathbf{F} | 163/168~(97%) | 160 (98%) | 3~(2%) | 0 | 100 | 100 |
| 6 | Ν | 159/168~(95%) | 154 (97%) | 4 (2%) | 1 (1%) | 25 | 65 |
| 7 | G | 114/151~(76%) | 112 (98%) | 1 (1%) | 1 (1%) | 17 | 56 |
| 7 | Ο | 126/151~(83%) | 123~(98%) | 2(2%) | 1 (1%) | 19 | 60 |
| 8 | М | 323/455~(71%) | 311 (96%) | 10 (3%) | 2(1%) | 25 | 65 |
| 8 | Р | 320/455~(70%) | 307~(96%) | 8 (2%) | 5(2%) | 9 | 45 |
| All | All | 3903/4872~(80%) | 3735 (96%) | 133 (3%) | 35 (1%) | 17 | 56 |

All (35) Ramachandran outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | А | 77 | PRO |
| 1 | А | 157 | GLN |
| 3 | С | 140 | PRO |
| 5 | Ε | 108 | GLU |
| 5 | Е | 112 | PRO |
| 8 | М | 608 | PRO |



| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 8 | Р | 608 | PRO |
| 1 | А | 70 | PRO |
| 3 | С | 27 | PRO |
| 3 | С | 190 | PRO |
| 5 | Е | 104 | PRO |
| 8 | Р | 681 | PRO |
| 2 | В | 7 | LYS |
| 5 | Е | 109 | PRO |
| 1 | Н | 70 | PRO |
| 1 | Н | 77 | PRO |
| 3 | J | 190 | PRO |
| 2 | Ι | 348 | PRO |
| 8 | Р | 680 | LEU |
| 8 | Р | 698 | PRO |
| 1 | А | 236 | PRO |
| 5 | Е | 103 | PHE |
| 8 | М | 566 | PRO |
| 1 | Н | 236 | PRO |
| 8 | Р | 566 | PRO |
| 6 | Ν | 6 | ARG |
| 1 | А | 6 | PRO |
| 2 | В | 346 | ASP |
| 1 | Н | 6 | PRO |
| 2 | В | 385 | VAL |
| 5 | Е | 111 | PHE |
| 3 | С | 62 | PRO |
| 7 | G | 37 | PRO |
| 7 | 0 | 37 | PRO |
| 2 | Ι | 176 | PRO |

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles | |
|-----|-------|------------|-----------|----------|-------------|-----|
| 1 | А | 8/363~(2%) | 8 (100%) | 0 | 100 | 100 |
| 1 | Н | 7/363~(2%) | 7 (100%) | 0 | 100 | 100 |



| Mol | Chain | Analysed | Rotameric | Outliers | Perce | ntiles |
|-----|-------|--------------|-----------|----------|-------|--------|
| 2 | В | 6/345~(2%) | 6 (100%) | 0 | 100 | 100 |
| 3 | С | 10/313~(3%) | 10 (100%) | 0 | 100 | 100 |
| 4 | D | 1/264~(0%) | 1 (100%) | 0 | 100 | 100 |
| 4 | Κ | 5/264~(2%) | 5 (100%) | 0 | 100 | 100 |
| 5 | Ε | 1/159~(1%) | 1 (100%) | 0 | 100 | 100 |
| 6 | F | 1/155~(1%) | 1 (100%) | 0 | 100 | 100 |
| 6 | Ν | 2/155~(1%) | 2~(100%) | 0 | 100 | 100 |
| 7 | G | 1/123~(1%) | 1 (100%) | 0 | 100 | 100 |
| 8 | М | 1/392~(0%) | 1 (100%) | 0 | 100 | 100 |
| 8 | Р | 5/392~(1%) | 5 (100%) | 0 | 100 | 100 |
| All | All | 48/3288 (2%) | 48 (100%) | 0 | 100 | 100 |

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond



| Mol Tuno Chain | | Dec | Link | Bond lengths | | | Bond angles | | | |
|----------------|------|------|------|--------------|----------|------|-------------|----------|------|----------|
| | туре | Unam | nes | | Counts | RMSZ | # Z > 2 | Counts | RMSZ | # Z > 2 |
| 12 | ATP | Н | 502 | 11 | 26,33,33 | 0.94 | 1 (3%) | 31,52,52 | 1.55 | 5 (16%) |
| 12 | ATP | А | 502 | 11 | 26,33,33 | 0.96 | 1 (3%) | 31,52,52 | 1.55 | 5 (16%) |
| 12 | ATP | Ι | 502 | 11 | 26,33,33 | 0.93 | 1 (3%) | 31,52,52 | 1.52 | 5(16%) |
| 12 | ATP | В | 502 | 11 | 26,33,33 | 0.96 | 1 (3%) | 31,52,52 | 1.43 | 5 (16%) |

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with |Z| > 2 is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

| Mol | Type | Chain | Res | Link | Chirals | Torsions | Rings |
|-----|------|-------|-----|------|---------|------------|---------|
| 12 | ATP | Н | 502 | 11 | - | 3/18/38/38 | 0/3/3/3 |
| 12 | ATP | А | 502 | 11 | - | 0/18/38/38 | 0/3/3/3 |
| 12 | ATP | Ι | 502 | 11 | - | 1/18/38/38 | 0/3/3/3 |
| 12 | ATP | В | 502 | 11 | - | 6/18/38/38 | 0/3/3/3 |

All (4) bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(Å) | $\mathrm{Ideal}(\mathrm{\AA})$ |
|-----|-------|-----|------|-------|------|-------------|--------------------------------|
| 12 | В | 502 | ATP | C5-C4 | 2.56 | 1.47 | 1.40 |
| 12 | Н | 502 | ATP | C5-C4 | 2.54 | 1.47 | 1.40 |
| 12 | А | 502 | ATP | C5-C4 | 2.53 | 1.47 | 1.40 |
| 12 | Ι | 502 | ATP | C5-C4 | 2.48 | 1.47 | 1.40 |

All (20) bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | $Observed(^{o})$ | $Ideal(^{o})$ |
|-----|-------|-----|------|-----------|-------|------------------|---------------|
| 12 | Η | 502 | ATP | PB-O3B-PG | -3.63 | 120.36 | 132.83 |
| 12 | А | 502 | ATP | PB-O3B-PG | -3.52 | 120.74 | 132.83 |
| 12 | Ι | 502 | ATP | PA-O3A-PB | -3.44 | 121.03 | 132.83 |
| 12 | А | 502 | ATP | PA-O3A-PB | -3.40 | 121.17 | 132.83 |
| 12 | Н | 502 | ATP | PA-O3A-PB | -3.39 | 121.21 | 132.83 |
| 12 | В | 502 | ATP | N3-C2-N1 | -3.36 | 123.43 | 128.68 |
| 12 | Ι | 502 | ATP | N3-C2-N1 | -3.25 | 123.59 | 128.68 |
| 12 | В | 502 | ATP | PA-O3A-PB | -3.23 | 121.76 | 132.83 |
| 12 | А | 502 | ATP | N3-C2-N1 | -3.17 | 123.72 | 128.68 |
| 12 | Н | 502 | ATP | N3-C2-N1 | -3.17 | 123.73 | 128.68 |



| Mol | Chain | Res | Type | Atoms | Z | $Observed(^{o})$ | $Ideal(^{o})$ |
|-----|-------|-----|------|-------------|-------|------------------|---------------|
| 12 | А | 502 | ATP | C3'-C2'-C1' | 3.14 | 105.71 | 100.98 |
| 12 | Ι | 502 | ATP | PB-O3B-PG | -3.14 | 122.04 | 132.83 |
| 12 | Н | 502 | ATP | C3'-C2'-C1' | 3.08 | 105.61 | 100.98 |
| 12 | Ι | 502 | ATP | C3'-C2'-C1' | 3.06 | 105.58 | 100.98 |
| 12 | В | 502 | ATP | C4-C5-N7 | -2.82 | 106.45 | 109.40 |
| 12 | Н | 502 | ATP | C4-C5-N7 | -2.80 | 106.48 | 109.40 |
| 12 | А | 502 | ATP | C4-C5-N7 | -2.78 | 106.50 | 109.40 |
| 12 | Ι | 502 | ATP | C4-C5-N7 | -2.73 | 106.56 | 109.40 |
| 12 | В | 502 | ATP | C3'-C2'-C1' | 2.65 | 104.97 | 100.98 |
| 12 | В | 502 | ATP | PB-O3B-PG | -2.28 | 125.00 | 132.83 |

There are no chirality outliers.

| | , | | | |
|-----|-------|-----|------|------|
| Mol | Chain | Res | Type | Atom |
| | 1 | | 1 | |

All (10) torsion outliers are listed below:

| Mol | Chain | \mathbf{Res} | Type | Atoms |
|-----|-------|----------------|------|-----------------|
| 12 | В | 502 | ATP | O4'-C4'-C5'-O5' |
| 12 | В | 502 | ATP | C3'-C4'-C5'-O5' |
| 12 | В | 502 | ATP | PG-O3B-PB-O1B |
| 12 | В | 502 | ATP | PG-O3B-PB-O2B |
| 12 | В | 502 | ATP | PB-O3A-PA-O1A |
| 12 | Н | 502 | ATP | PB-O3A-PA-O1A |
| 12 | Н | 502 | ATP | PB-O3A-PA-O2A |
| 12 | В | 502 | ATP | PB-O3A-PA-O2A |
| 12 | Ι | 502 | ATP | PB-O3A-PA-O2A |
| 12 | Н | 502 | ATP | C3'-C4'-C5'-O5' |

There are no ring outliers.

1 monomer is involved in 1 short contact:

| Mol | Chain | Res | Type | Clashes | Symm-Clashes |
|-----|-------|-----|------|---------|--------------|
| 12 | А | 502 | ATP | 1 | 0 |

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier.



The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

















5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

| Mol | Chain | Number of breaks |
|-----|-------|------------------|
| 5 | Ε | 1 |

All chain breaks are listed below:

| Model | Chain | Residue-1 | Atom-1 | Residue-2 | Atom-2 | Distance (Å) |
|-------|-------|-----------|--------|-----------|--------|--------------|
| 1 | E | 19:MET | С | 21:ALA | Ν | 4.49 |



6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

| Mol | Chain | Analysed | $<$ RSRZ $>$ | #RSRZ>2 | $\mathbf{OWAB}(\mathbf{\AA}^2)$ | Q<0.9 |
|-----|-------|-----------------|---------------------|--------------|---------------------------------|-------|
| 1 | А | 398/418~(95%) | -0.82 | 0 100 100 | 76, 138, 200, 295 | 0 |
| 1 | Н | 386/418~(92%) | -0.81 | 0 100 100 | 82, 157, 246, 311 | 0 |
| 2 | В | 369/394~(93%) | -0.78 | 1 (0%) 94 90 | 66, 155, 253, 296 | 0 |
| 2 | Ι | 209/394~(53%) | -0.65 | 1 (0%) 91 85 | 121, 215, 330, 447 | 0 |
| 3 | С | 339/372~(91%) | -0.67 | 0 100 100 | 78, 140, 193, 229 | 0 |
| 3 | J | 258/372~(69%) | -0.46 | 1 (0%) 92 87 | 140, 245, 321, 379 | 0 |
| 4 | D | 279/300~(93%) | -0.73 | 0 100 100 | 71, 219, 342, 423 | 0 |
| 4 | К | 283/300 (94%) | -0.86 | 0 100 100 | 76, 146, 213, 256 | 0 |
| 5 | Е | 163/178~(91%) | -0.73 | 0 100 100 | 117, 178, 251, 281 | 0 |
| 5 | L | 138/178 (77%) | -0.52 | 0 100 100 | 145, 242, 330, 351 | 0 |
| 6 | F | 165/168~(98%) | -0.84 | 0 100 100 | 58, 118, 203, 233 | 0 |
| 6 | N | 163/168~(97%) | -0.81 | 0 100 100 | 78, 156, 224, 288 | 0 |
| 7 | G | 116/151 (76%) | -0.88 | 0 100 100 | 90, 179, 281, 305 | 0 |
| 7 | Ο | 132/151~(87%) | -0.55 | 1 (0%) 86 79 | 142, 248, 305, 333 | 0 |
| 8 | М | 329/455~(72%) | -0.70 | 1 (0%) 94 90 | 102, 202, 289, 315 | 0 |
| 8 | Р | 328/455~(72%) | -0.88 | 0 100 100 | 71, 156, 277, 312 | 0 |
| 9 | Q | 0/6 | - | _ | - | - |
| 10 | R | 0/9 | - | _ | - | - |
| All | All | 4055/4887 (82%) | -0.74 | 5 (0%) 95 94 | 58, 172, 284, 447 | 0 |

All (5) RSRZ outliers are listed below:

| Mol | Chain | Res | Type | RSRZ |
|-----|-------|-----|------|------|
| 8 | М | 623 | GLY | 2.7 |
| 3 | J | 95 | ILE | 2.7 |
| 2 | Ι | 225 | TYR | 2.5 |



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| Mol | Chain | \mathbf{Res} | Type | RSRZ |
|-----|-------|----------------|------|------|
| 7 | 0 | 151 | VAL | 2.4 |
| 2 | В | 383 | VAL | 2.2 |

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

| Mol | Type | Chain | Res | Atoms | RSCC | RSR | $B-factors(Å^2)$ | Q<0.9 |
|-----|------|-------|-----|-------|------|------|------------------|-------|
| 11 | CA | Ι | 501 | 1/1 | 0.82 | 0.32 | 301,301,301,301 | 0 |
| 12 | ATP | Ι | 502 | 31/31 | 0.85 | 0.18 | 177,240,301,313 | 0 |
| 12 | ATP | Н | 502 | 31/31 | 0.93 | 0.25 | 129,147,198,226 | 0 |
| 12 | ATP | В | 502 | 31/31 | 0.93 | 0.22 | 100,117,145,151 | 0 |
| 12 | ATP | А | 502 | 31/31 | 0.95 | 0.16 | 106,119,162,181 | 0 |
| 11 | CA | Н | 501 | 1/1 | 0.98 | 0.25 | 144,144,144,144 | 0 |
| 11 | CA | А | 501 | 1/1 | 0.98 | 0.23 | 87,87,87,87 | 0 |
| 11 | CA | В | 501 | 1/1 | 0.99 | 0.27 | 116,116,116,116 | 0 |

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.











6.5 Other polymers (i)

There are no such residues in this entry.

