



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:12 AM UTC

PDB ID : 1QZ7 / pdb_00001qz7
Title : Beta-catenin binding domain of Axin in complex with beta-catenin
Authors : Xing, Y.; Clements, W.K.; Kimelman, D.; Xu, W.
Deposited on : 2003-09-15
Resolution : 2.20 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

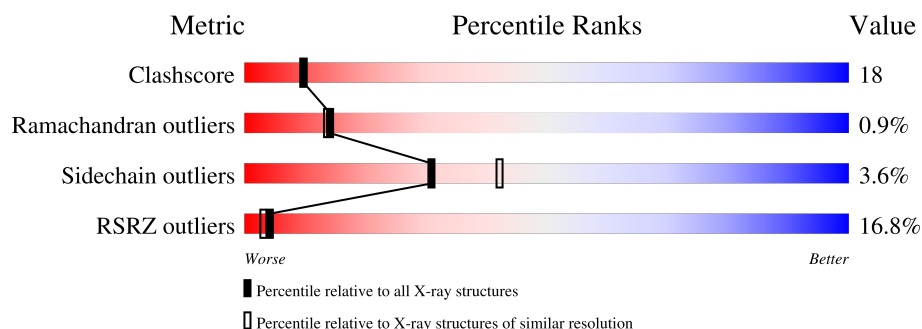
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 $\leq 5\%$. 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	A	533	16% 70% 27% ..
2	B	70	11% 11% 9% 76% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4148 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 Beta-catenin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	524	Total	C	N	O	S	0	0	0
			3979	2496	721	735	27			

- Molecule 2 is a protein called Axin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	17	Total	C	N	O	S	0	0	0
			138	83	24	30	1			

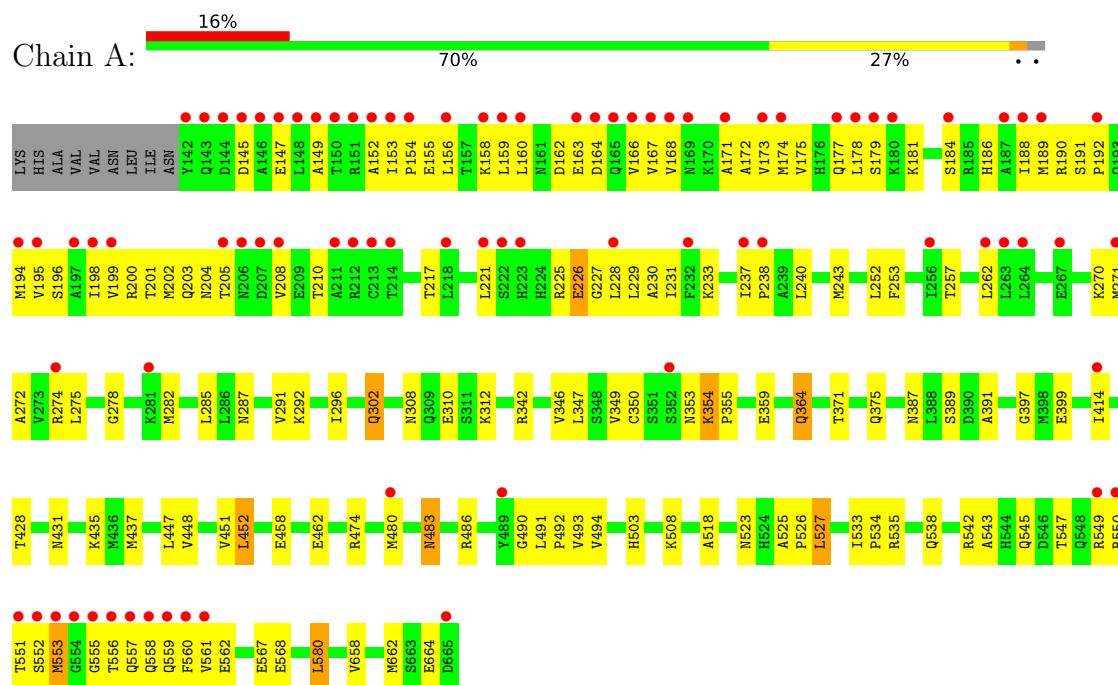
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	31	Total	O	0	0
			31	31		

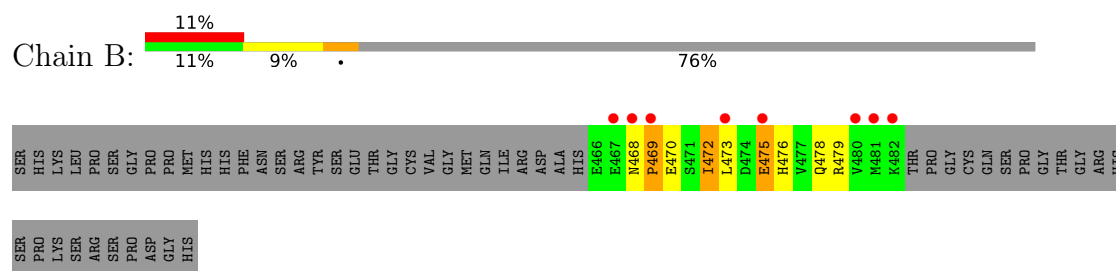
3 Residue-property plots

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: Beta-catenin



• Molecule 2: Axin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	85.14Å 75.04Å 101.40Å 90.00° 97.47° 90.00°	Depositor
Resolution (Å)	42.21 – 2.20 42.21 – 2.20	Depositor EDS
% Data completeness (in resolution range)	81.1 (42.21-2.20) 95.5 (42.21-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.256 0.223 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	35.2	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4148	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.76% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.41	0/4035	0.86	4/5476 (0.1%)
2	B	0.34	0/139	1.22	1/187 (0.5%)
All	All	0.41	0/4174	0.87	5/5663 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	475	GLU	CB-CA-C	12.33	130.38	110.90
1	A	551	THR	N-CA-C	-7.53	103.98	113.01
1	A	291	VAL	N-CA-C	6.91	117.70	110.72
1	A	474	ARG	N-CA-C	6.30	119.94	111.24
1	A	389	SER	N-CA-C	5.22	118.43	111.75

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	A	3979	0	4094	131	0
2	B	138	0	124	23	0
3	A	31	0	0	0	0
All	All	4148	0	4218	147	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 18.

All (147) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:LEU:HD23	1:A:171:ALA:HB2	1.48	0.96
1:A:658:VAL:HG12	1:A:662:MET:HE2	1.49	0.91
1:A:271:MET:O	1:A:275:LEU:HG	1.84	0.78
1:A:550:ARG:HH22	1:A:561:VAL:CG2	1.99	0.76
1:A:272:ALA:HA	1:A:275:LEU:HD12	1.68	0.75
1:A:364:GLN:HE21	1:A:364:GLN:H	1.35	0.74
1:A:364:GLN:H	1:A:364:GLN:NE2	1.85	0.74
1:A:355:PRO:O	1:A:359:GLU:HG2	1.87	0.74
1:A:270:LYS:HB3	1:A:274:ARG:HH12	1.55	0.70
1:A:154:PRO:O	1:A:158:LYS:HG3	1.91	0.70
2:B:468:ASN:HD22	2:B:469:PRO:CD	2.05	0.70
1:A:257:THR:HG23	2:B:476:HIS:ND1	2.07	0.69
1:A:153:ILE:HB	1:A:154:PRO:HD3	1.75	0.69
1:A:508:LYS:HE3	1:A:568:GLU:OE1	1.92	0.68
1:A:201:THR:O	1:A:205:THR:HG22	1.93	0.68
1:A:188:ILE:HG23	1:A:194:MET:HB3	1.77	0.67
1:A:371:THR:HG22	1:A:371:THR:O	1.95	0.66
2:B:468:ASN:HD22	2:B:469:PRO:HD2	1.60	0.66
1:A:200:ARG:HA	1:A:203:GLN:HE21	1.61	0.66
1:A:292:LYS:O	1:A:296:ILE:HG12	1.95	0.66
1:A:278:GLY:O	1:A:282:MET:HG3	1.97	0.65
1:A:189:MET:SD	1:A:226:GLU:HB2	2.38	0.63
1:A:483:ASN:ND2	1:A:486:ARG:NH2	2.48	0.62
1:A:189:MET:HB2	1:A:221:LEU:HD22	1.80	0.61
1:A:270:LYS:HB3	1:A:274:ARG:NH1	2.14	0.61
1:A:399:GLU:H	1:A:399:GLU:CD	2.08	0.61
1:A:533:ILE:HB	1:A:534:PRO:HD3	1.83	0.60
1:A:190:ARG:O	1:A:192:PRO:HD3	2.02	0.60
2:B:468:ASN:HD22	2:B:469:PRO:N	2.00	0.59
1:A:173:VAL:HG12	1:A:177:GLN:HE21	1.68	0.58
1:A:458:GLU:OE2	1:A:503:HIS:HD2	1.86	0.58
1:A:557:GLN:O	1:A:558:GLN:HG2	2.03	0.58
1:A:550:ARG:HD3	1:A:558:GLN:OE1	2.04	0.58
1:A:159:LEU:HD11	1:A:167:VAL:HG12	1.85	0.57
1:A:296:ILE:HD12	2:B:476:HIS:HD2	1.69	0.57
1:A:181:LYS:HB2	1:A:184:SER:OG	2.03	0.57
2:B:472:ILE:C	2:B:472:ILE:HD13	2.29	0.57
1:A:518:ALA:CB	1:A:580:LEU:HD13	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200:ARG:HA	1:A:203:GLN:NE2	2.22	0.55
1:A:159:LEU:CD2	1:A:171:ALA:HB2	2.29	0.54
2:B:468:ASN:C	2:B:470:GLU:H	2.14	0.54
1:A:371:THR:O	1:A:371:THR:CG2	2.55	0.54
1:A:525:ALA:HB3	1:A:526:PRO:HD3	1.90	0.53
1:A:171:ALA:HA	1:A:174:MET:HE3	1.90	0.53
1:A:347:LEU:O	1:A:350:CYS:HB3	2.09	0.53
1:A:174:MET:HA	1:A:177:GLN:NE2	2.23	0.53
1:A:188:ILE:HG23	1:A:194:MET:CB	2.38	0.53
1:A:559:GLN:HG2	1:A:560:PHE:N	2.23	0.53
1:A:171:ALA:O	1:A:175:VAL:HG23	2.09	0.53
1:A:547:THR:HA	1:A:558:GLN:NE2	2.23	0.52
1:A:164:ASP:OD2	1:A:167:VAL:HG23	2.09	0.52
1:A:558:GLN:OE1	1:A:558:GLN:HA	2.10	0.52
1:A:205:THR:O	1:A:205:THR:HG23	2.10	0.52
1:A:550:ARG:HH22	1:A:561:VAL:HG21	1.73	0.51
1:A:543:ALA:O	1:A:547:THR:HG23	2.10	0.51
1:A:542:ARG:HG2	1:A:542:ARG:HH11	1.75	0.51
2:B:468:ASN:HD22	2:B:468:ASN:C	2.18	0.51
1:A:535:ARG:HH11	1:A:538:GLN:NE2	2.09	0.51
1:A:200:ARG:O	1:A:204:ASN:ND2	2.44	0.51
1:A:177:GLN:C	1:A:179:SER:H	2.18	0.50
1:A:149:ALA:O	1:A:153:ILE:HG12	2.11	0.50
1:A:226:GLU:H	1:A:226:GLU:CD	2.19	0.50
1:A:202:MET:HE3	1:A:243:MET:HG3	1.92	0.50
1:A:535:ARG:NH1	1:A:538:GLN:NE2	2.59	0.50
1:A:230:ALA:HA	1:A:233:LYS:CE	2.41	0.50
1:A:227:GLY:O	1:A:231:ILE:HG13	2.12	0.49
1:A:257:THR:HG23	2:B:476:HIS:CE1	2.46	0.49
1:A:483:ASN:HD21	1:A:486:ARG:NH2	2.11	0.49
2:B:479:ARG:CB	2:B:479:ARG:HH11	2.26	0.48
1:A:567:GLU:H	1:A:567:GLU:CD	2.20	0.48
1:A:159:LEU:HD21	1:A:167:VAL:HG12	1.94	0.48
1:A:178:LEU:HB3	1:A:184:SER:HB2	1.95	0.48
1:A:160:LEU:HD12	1:A:201:THR:OG1	2.14	0.48
1:A:518:ALA:HB2	1:A:580:LEU:HD13	1.95	0.48
1:A:202:MET:HE1	1:A:240:LEU:HD23	1.94	0.48
1:A:243:MET:HA	1:A:243:MET:HE2	1.96	0.48
1:A:431:ASN:O	1:A:435:LYS:HG3	2.13	0.48
1:A:448:VAL:O	1:A:452:LEU:HD22	2.15	0.47
1:A:552:SER:O	1:A:553:MET:HB2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:PHE:HB3	2:B:472:ILE:HD12	1.97	0.47
1:A:159:LEU:HD11	1:A:167:VAL:CG1	2.45	0.47
1:A:354:LYS:HE3	1:A:387:ASN:O	2.15	0.47
1:A:354:LYS:HB2	1:A:354:LYS:NZ	2.30	0.47
1:A:547:THR:HA	1:A:558:GLN:HE22	1.80	0.47
1:A:145:ASP:O	1:A:149:ALA:HB3	2.15	0.47
1:A:253:PHE:HD2	2:B:469:PRO:O	1.98	0.47
1:A:490:GLY:O	1:A:493:VAL:HG12	2.15	0.47
1:A:149:ALA:HA	1:A:152:ALA:HB3	1.97	0.46
1:A:172:ALA:CB	1:A:210:THR:HG23	2.46	0.46
1:A:308:ASN:OD1	1:A:310:GLU:HB2	2.15	0.46
1:A:428:THR:O	1:A:435:LYS:HE2	2.15	0.46
2:B:475:GLU:O	2:B:478:GLN:HB2	2.15	0.46
1:A:179:SER:OG	1:A:217:THR:HG23	2.16	0.46
1:A:346:VAL:O	1:A:349:VAL:HG22	2.16	0.46
1:A:491:LEU:HB2	1:A:492:PRO:HD3	1.98	0.45
1:A:152:ALA:O	1:A:155:GLU:HB3	2.16	0.45
1:A:173:VAL:HG12	1:A:177:GLN:NE2	2.30	0.45
1:A:164:ASP:O	1:A:168:VAL:HG23	2.16	0.45
1:A:538:GLN:NE2	1:A:542:ARG:HH21	2.15	0.45
1:A:186:HIS:HA	1:A:189:MET:HE2	1.98	0.45
1:A:447:LEU:O	1:A:451:VAL:HG23	2.16	0.45
1:A:414:ILE:H	1:A:414:ILE:HG13	1.60	0.45
1:A:561:VAL:O	1:A:562:GLU:HB2	2.16	0.45
1:A:198:ILE:HD13	1:A:217:THR:HG21	1.99	0.45
1:A:518:ALA:HB1	1:A:580:LEU:HD13	1.99	0.45
2:B:468:ASN:O	2:B:470:GLU:N	2.50	0.44
1:A:173:VAL:O	1:A:177:GLN:HG3	2.16	0.44
1:A:308:ASN:O	1:A:312:LYS:HG3	2.16	0.44
1:A:523:ASN:O	1:A:527:LEU:HB2	2.17	0.44
1:A:195:VAL:O	1:A:199:VAL:HG23	2.17	0.44
1:A:202:MET:HE3	1:A:243:MET:CG	2.47	0.44
1:A:354:LYS:HD3	1:A:391:ALA:HB2	2.00	0.44
1:A:560:PHE:N	1:A:560:PHE:CD1	2.86	0.44
1:A:458:GLU:OE2	1:A:503:HIS:CD2	2.70	0.44
1:A:196:SER:O	1:A:200:ARG:HG3	2.18	0.43
1:A:350:CYS:HB3	1:A:353:ASN:HB2	1.99	0.43
1:A:556:THR:O	1:A:557:GLN:C	2.60	0.43
2:B:479:ARG:HH11	2:B:479:ARG:HB3	1.82	0.43
1:A:162:ASP:HB3	1:A:168:VAL:HG22	2.01	0.43
2:B:479:ARG:HB2	2:B:479:ARG:NH1	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:ARG:HH11	1:A:225:ARG:HG3	1.84	0.43
1:A:545:GLN:O	1:A:549:ARG:HG3	2.19	0.43
1:A:145:ASP:C	1:A:147:GLU:H	2.25	0.43
1:A:237:ILE:HB	1:A:238:PRO:HD3	2.01	0.43
1:A:257:THR:HG23	2:B:476:HIS:CG	2.53	0.43
1:A:658:VAL:O	1:A:662:MET:HG3	2.19	0.43
1:A:270:LYS:O	1:A:274:ARG:HG3	2.18	0.43
1:A:523:ASN:C	1:A:526:PRO:HD2	2.44	0.43
1:A:549:ARG:HH11	1:A:549:ARG:HG2	1.83	0.42
2:B:479:ARG:CB	2:B:479:ARG:NH1	2.82	0.42
1:A:397:GLY:HA2	1:A:399:GLU:OE1	2.19	0.42
2:B:472:ILE:HG23	2:B:473:LEU:N	2.35	0.42
1:A:296:ILE:HD12	2:B:476:HIS:CD2	2.52	0.42
1:A:230:ALA:HA	1:A:233:LYS:HE3	2.00	0.41
2:B:468:ASN:C	2:B:470:GLU:N	2.78	0.41
2:B:468:ASN:C	2:B:468:ASN:ND2	2.77	0.41
1:A:153:ILE:HG21	1:A:191:SER:OG	2.21	0.41
1:A:160:LEU:HD22	1:A:171:ALA:CB	2.51	0.41
1:A:228:LEU:HD22	1:A:262:LEU:HD23	2.03	0.41
1:A:483:ASN:HD21	1:A:486:ARG:HH22	1.68	0.41
1:A:156:LEU:HD12	1:A:178:LEU:HD12	2.02	0.41
1:A:302:GLN:NE2	1:A:342:ARG:HH21	2.19	0.40
1:A:354:LYS:HB3	1:A:355:PRO:CD	2.51	0.40
1:A:493:VAL:HG13	1:A:494:VAL:N	2.36	0.40
1:A:229:LEU:HG	1:A:233:LYS:HE2	2.02	0.40
1:A:285:LEU:C	1:A:287:ASN:H	2.30	0.40
2:B:468:ASN:ND2	2:B:469:PRO:HD2	2.33	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	Percentiles	
1	A	522/533 (98%)	488 (94%)	30 (6%)	4 (1%)	16	16
2	B	15/70 (21%)	13 (87%)	1 (7%)	1 (7%)	1	0
All	All	537/603 (89%)	501 (93%)	31 (6%)	5 (1%)	14	14

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	553	MET
1	A	208	VAL
1	A	555	GLY
2	B	469	PRO
1	A	166	VAL

5.3.2 Protein sidechains ⓘ

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	A	428/442 (97%)	413 (96%)	15 (4%)	32	43
2	B	16/62 (26%)	15 (94%)	1 (6%)	16	19
All	All	444/504 (88%)	428 (96%)	16 (4%)	31	42

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	163	GLU
1	A	226	GLU
1	A	252	LEU
1	A	302	GLN
1	A	354	LYS
1	A	364	GLN
1	A	375	GLN
1	A	437	MET
1	A	452	LEU
1	A	462	GLU
1	A	480	MET

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Mol	Chain	Res	Type
1	A	483	ASN
1	A	527	LEU
1	A	580	LEU
1	A	664	GLU
2	B	472	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	177	GLN
1	A	186	HIS
1	A	193	GLN
1	A	203	GLN
1	A	204	ASN
1	A	220	ASN
1	A	266	GLN
1	A	302	GLN
1	A	326	ASN
1	A	364	GLN
1	A	415	ASN
1	A	483	ASN
1	A	503	HIS
1	A	538	GLN
1	A	601	GLN
1	A	611	GLN
2	B	468	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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, 95th 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	OWAB(Å ²)	Q<0.9
1	A	524/533 (98%)	0.83	83 (15%) 5 3	22, 47, 99, 108	0
2	B	17/70 (24%)	1.84	8 (47%) 0 0	71, 81, 94, 94	0
All	All	541/603 (89%)	0.86	91 (16%) 4 3	22, 48, 99, 108	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	552	SER	8.0
1	A	557	GLN	5.6
1	A	148	LEU	5.3
1	A	160	LEU	5.2
1	A	551	THR	5.0
1	A	208	VAL	4.9
1	A	556	THR	4.9
1	A	665	ASP	4.9
1	A	168	VAL	4.3
1	A	177	GLN	4.1
1	A	489	TYR	3.9
1	A	166	VAL	3.9
1	A	167	VAL	3.8
2	B	482	LYS	3.8
1	A	142	TYR	3.6
2	B	481	MET	3.5
1	A	144	ASP	3.5
1	A	205	THR	3.4
1	A	195	VAL	3.4
1	A	212	ARG	3.4
1	A	159	LEU	3.4
1	A	146	ALA	3.3
1	A	237	ILE	3.3
1	A	221	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	151	ARG	3.2
1	A	152	ALA	3.1
1	A	554	GLY	3.1
1	A	154	PRO	3.1
1	A	550	ARG	3.1
2	B	480	VAL	3.1
1	A	149	ALA	3.0
1	A	271	MET	3.0
1	A	192	PRO	3.0
1	A	559	GLN	2.9
1	A	199	VAL	2.9
1	A	211	ALA	2.8
1	A	174	MET	2.8
1	A	178	LEU	2.8
1	A	171	ALA	2.8
1	A	206	ASN	2.8
1	A	558	GLN	2.8
1	A	232	PHE	2.8
1	A	165	GLN	2.7
1	A	223	HIS	2.7
1	A	555	GLY	2.7
1	A	553	MET	2.7
1	A	560	PHE	2.6
1	A	150	THR	2.6
1	A	480	MET	2.6
2	B	469	PRO	2.6
1	A	549	ARG	2.6
1	A	238	PRO	2.6
1	A	281	LYS	2.5
1	A	173	VAL	2.5
1	A	184	SER	2.5
1	A	143	GLN	2.5
1	A	188	ILE	2.5
1	A	256	ILE	2.5
1	A	197	ALA	2.4
1	A	198	ILE	2.4
1	A	228	LEU	2.4
1	A	352	SER	2.4
1	A	180	LYS	2.4
1	A	561	VAL	2.4
1	A	145	ASP	2.4
2	B	468	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	262	LEU	2.4
1	A	179	SER	2.3
1	A	267	GLU	2.3
2	B	467	GLU	2.3
1	A	264	LEU	2.3
1	A	169	ASN	2.2
1	A	207	ASP	2.2
1	A	189	MET	2.2
1	A	274	ARG	2.2
1	A	187	ALA	2.2
1	A	213	CYS	2.2
1	A	156	LEU	2.2
2	B	475	GLU	2.2
1	A	222	SER	2.2
1	A	194	MET	2.2
1	A	263	LEU	2.1
2	B	473	LEU	2.1
1	A	158	LYS	2.1
1	A	164	ASP	2.1
1	A	214	THR	2.1
1	A	147	GLU	2.1
1	A	218	LEU	2.1
1	A	414	ILE	2.1
1	A	153	ILE	2.0
1	A	163	GLU	2.0

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.