



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 04:54 AM UTC

PDB ID : 7UV8 / pdb_00007uv8
Title : Rad6-Bre1 Complex
Authors : Shukla, P.K.; Chandrasekharan, M.B.
Deposited on : 2022-04-29
Resolution : 2.70 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

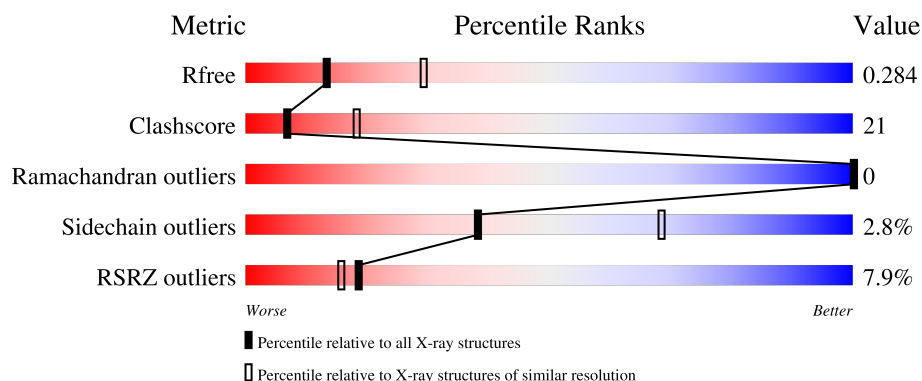
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 $\leq 5\%$. 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	150	9% 47% 51% ..
2	U	212	8% 41% 33% • 24%
2	V	212	3% 47% 27% • 25%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3861 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.

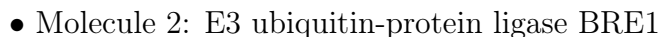
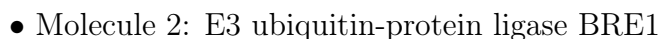
- Molecule 1 is a protein called Ubiquitin-conjugating enzyme E2 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	0	0
			1201	765	203	227	6			

- Molecule 2 is a protein called E3 ubiquitin-protein ligase BRE1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	V	160	Total	C	N	O	S	0	0	0
			1319	823	231	259	6			
2	U	161	Total	C	N	O	S	0	0	0
			1341	845	236	254	6			

- Molecule 1: Ubiquitin-conjugating enzyme E2 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	241.58Å 49.95Å 57.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.49 – 2.70 38.49 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.49-2.70) 99.9 (38.49-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.240 , 0.280 0.270 , 0.284	Depositor DCC
R_{free} test set	982 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	75.6	Xtriage
Anisotropy	0.419	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 74.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3861	wwPDB-VP
Average B, all atoms (Å ²)	105.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.85% 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

5.1 Standard geometry

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	0.58	0/1236	0.80	0/1683
2	U	0.75	2/1356 (0.1%)	1.00	1/1814 (0.1%)
2	V	0.67	0/1334	0.89	0/1789
All	All	0.67	2/3926 (0.1%)	0.90	1/5286 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	U	105	VAL	C-O	7.84	1.32	1.24
2	U	79	PHE	C-O	5.78	1.30	1.24

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	U	40	ARG	N-CA-C	-5.21	105.49	111.07

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	63	TYR	Peptide

5.2 Too-close contacts ⓘ

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	1201	0	1168	66	0
2	U	1341	0	1367	73	0
2	V	1319	0	1323	58	0
All	All	3861	0	3858	164	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 21.

All (164) 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)
2:V:35:PHE:HB2	2:U:26:VAL:HG21	1.43	0.99
1:A:80:ASN:HD21	1:A:117:ASN:HB3	1.30	0.96
2:U:92:LEU:HD11	2:U:107:ARG:HB3	1.56	0.84
2:U:80:ILE:HG21	2:U:111:PHE:HE1	1.48	0.79
2:U:60:ILE:O	2:U:64:ARG:HG3	1.83	0.78
2:V:112:MET:HE2	2:U:72:LEU:HD23	1.67	0.74
2:U:80:ILE:O	2:U:84:CYS:SG	2.47	0.73
2:V:35:PHE:HB2	2:U:26:VAL:CG2	2.21	0.69
2:U:100:ASP:O	2:U:104:ILE:HG12	1.92	0.68
2:U:92:LEU:CD1	2:U:107:ARG:HB3	2.25	0.67
1:A:43:PRO:HD3	1:A:110:GLN:HG3	1.76	0.66
2:U:80:ILE:HG21	2:U:111:PHE:CE1	2.31	0.66
2:U:189:VAL:HG12	2:U:190:PHE:CD1	2.30	0.65
1:A:129:LEU:HB3	1:A:137:TYR:HB2	1.78	0.65
2:V:58:GLU:OE2	2:V:159:GLN:HG2	1.96	0.64
2:V:173:ASN:OD1	2:V:176:ARG:NH2	2.26	0.64
1:A:110:GLN:HA	1:A:113:PHE:HD2	1.65	0.62
1:A:129:LEU:HD22	1:A:133:HIS:HB2	1.82	0.62
2:V:47:ALA:O	2:V:51:GLN:HG3	2.00	0.62
2:V:112:MET:HE3	2:V:144:LYS:HG3	1.82	0.61
2:V:58:GLU:O	2:V:62:VAL:HG23	2.00	0.61
1:A:3:THR:HG21	1:A:61:GLU:OE1	2.00	0.61
2:U:178:TYR:O	2:U:182:GLU:HG2	2.01	0.61
2:U:26:VAL:HA	2:U:29:PHE:CD2	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:PRO:HB3	1:A:36:TRP:CD1	2.36	0.61
2:V:161:THR:HG22	2:U:160:LEU:HD11	1.83	0.60
2:U:136:ILE:O	2:U:140:THR:HG23	2.01	0.60
2:U:23:GLN:O	2:U:26:VAL:HG12	2.02	0.60
2:U:156:GLU:O	2:U:160:LEU:HD23	2.02	0.60
1:A:143:GLU:O	1:A:147:LYS:HG3	2.02	0.59
1:A:105:ILE:O	1:A:109:ILE:HG13	2.03	0.59
2:U:170:TYR:O	2:U:174:ILE:HD12	2.02	0.59
2:U:189:VAL:HG12	2:U:190:PHE:HD1	1.68	0.59
2:U:86:ASP:HB3	2:U:88:ASN:OD1	2.01	0.59
2:V:168:LYS:HE2	2:U:167:LEU:HD11	1.86	0.58
2:V:147:ARG:NH1	2:U:78:ARG:HD2	2.19	0.58
1:A:43:PRO:HG2	1:A:46:THR:HG21	1.86	0.58
2:V:51:GLN:HE22	2:U:168:LYS:NZ	2.01	0.57
1:A:21:PRO:HG2	1:A:24:VAL:HG21	1.84	0.57
2:V:120:LYS:HB3	2:U:79:PHE:CZ	2.40	0.57
2:V:157:ASN:OD1	2:U:157:ASN:HB2	2.04	0.57
1:A:32:ASN:OD1	1:A:34:MET:HB3	2.05	0.57
2:V:54:LEU:O	2:V:58:GLU:HG3	2.05	0.57
2:U:30:GLN:OE1	2:U:31:LYS:N	2.38	0.57
1:A:63:TYR:CG	1:A:64:PRO:HA	2.39	0.56
1:A:124:VAL:O	1:A:128:THR:HG23	2.06	0.55
1:A:63:TYR:CD1	1:A:64:PRO:HA	2.41	0.55
1:A:130:PHE:O	1:A:134:LYS:HE2	2.06	0.55
1:A:28:PRO:HB3	1:A:36:TRP:NE1	2.22	0.55
2:V:32:GLU:HG2	2:U:23:GLN:HG2	1.89	0.55
2:V:148:LYS:HD3	2:U:68:ASN:HB2	1.89	0.55
1:A:78:HIS:HE1	1:A:116:PRO:HA	1.72	0.54
2:U:86:ASP:HB2	2:U:89:GLU:CD	2.33	0.54
1:A:79:PRO:HD3	1:A:126:ALA:HB1	1.90	0.54
2:U:109:ASP:OD1	2:U:113:LYS:HE3	2.07	0.53
2:V:80:ILE:HG13	2:V:115:LEU:HD23	1.89	0.53
1:A:41:ILE:HD12	2:U:189:VAL:HG11	1.91	0.52
1:A:74:SER:HB2	1:A:76:MET:HE1	1.91	0.52
1:A:78:HIS:ND1	1:A:79:PRO:HD2	2.24	0.52
1:A:3:THR:HG23	1:A:6:ARG:H	1.73	0.52
2:V:146:LEU:HD13	2:U:146:LEU:HB3	1.92	0.52
1:A:90:ASP:HA	1:A:93:GLN:HG2	1.91	0.52
2:V:176:ARG:O	2:V:180:ARG:HG3	2.10	0.52
2:V:78:ARG:O	2:V:82:THR:HG23	2.10	0.51
2:V:143:LEU:HB3	2:V:147:ARG:NH2	2.24	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:U:56:ARG:O	2:U:60:ILE:HG13	2.10	0.51
1:A:101:ASP:O	1:A:105:ILE:HG13	2.11	0.51
2:U:78:ARG:O	2:U:81:GLU:HB3	2.11	0.51
1:A:3:THR:CG2	1:A:6:ARG:H	2.24	0.50
1:A:3:THR:HG23	1:A:5:ALA:N	2.27	0.50
1:A:46:THR:O	1:A:49:GLU:HB2	2.12	0.50
2:V:48:LEU:HD23	2:U:48:LEU:HG	1.93	0.50
2:U:69:ILE:HA	2:U:72:LEU:HD12	1.92	0.50
2:V:27:ILE:O	2:V:31:LYS:HG3	2.11	0.50
1:A:40:ILE:CD1	1:A:109:ILE:HB	2.42	0.49
2:V:81:GLU:HB2	2:V:93:CYS:HB3	1.94	0.49
2:V:35:PHE:CB	2:U:26:VAL:HG21	2.29	0.49
1:A:13:PHE:CE2	1:A:28:PRO:HD2	2.48	0.49
1:A:72:PHE:HB3	1:A:76:MET:HE3	1.93	0.49
1:A:69:HIS:CE1	1:A:86:GLU:HG2	2.48	0.49
2:V:42:ARG:HA	2:U:41:ARG:HD2	1.94	0.49
2:V:146:LEU:HB3	2:U:146:LEU:HD23	1.95	0.49
1:A:7:ARG:O	1:A:7:ARG:HD2	2.12	0.49
2:V:95:GLU:HB3	2:V:103:LEU:CD2	2.42	0.49
2:V:36:ARG:HG2	2:U:179:ASP:HB3	1.95	0.48
2:V:33:ALA:HA	2:V:36:ARG:NH1	2.28	0.48
2:V:78:ARG:O	2:V:81:GLU:HB3	2.13	0.48
1:A:141:VAL:O	1:A:144:THR:HB	2.12	0.48
2:V:153:LEU:HB3	2:U:153:LEU:HB3	1.95	0.48
2:V:23:GLN:O	2:V:26:VAL:HG12	2.14	0.48
1:A:43:PRO:HD3	1:A:110:GLN:CG	2.42	0.47
2:U:167:LEU:HD22	2:U:171:TYR:CE2	2.49	0.47
2:V:55:SER:OG	2:U:56:ARG:HD2	2.15	0.47
2:V:120:LYS:HB3	2:U:79:PHE:CE1	2.50	0.47
1:A:56:LEU:HD23	1:A:71:LYS:NZ	2.29	0.46
2:V:94:ARG:O	2:V:98:GLN:HG3	2.15	0.46
2:U:23:GLN:HE22	2:U:183:SER:HB2	1.81	0.46
1:A:8:ARG:NH2	1:A:98:PRO:O	2.49	0.46
2:V:101:GLU:O	2:V:104:ILE:HG22	2.16	0.45
1:A:103:ALA:O	1:A:107:THR:HG23	2.16	0.45
2:V:74:VAL:O	2:V:78:ARG:HG3	2.16	0.45
1:A:42:GLY:HA3	1:A:48:TYR:O	2.16	0.45
2:V:143:LEU:HB3	2:V:147:ARG:HH22	1.80	0.45
2:U:147:ARG:O	2:U:151:GLU:HG3	2.16	0.45
1:A:89:LEU:HD21	1:A:91:ILE:HG12	1.98	0.45
2:U:96:ILE:HG23	2:U:104:ILE:HD12	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:ASP:O	1:A:15:ARG:HB3	2.17	0.44
1:A:13:PHE:O	1:A:14:LYS:C	2.60	0.44
2:U:171:TYR:O	2:U:175:ILE:HG13	2.17	0.44
2:U:112:MET:O	2:U:116:THR:HG23	2.18	0.44
1:A:49:GLU:O	1:A:50:ASP:HB2	2.17	0.44
1:A:56:LEU:HD23	1:A:71:LYS:HZ3	1.81	0.44
1:A:44:ALA:HA	1:A:49:GLU:HG3	1.99	0.44
2:V:136:ILE:HG23	2:U:139:LEU:HD11	2.00	0.44
2:V:150:LYS:HG2	2:U:153:LEU:HD12	1.99	0.44
1:A:75:GLU:OE1	1:A:147:LYS:NZ	2.43	0.44
2:V:178:TYR:HE2	2:U:175:ILE:HG23	1.82	0.44
1:A:84:ASN:OD1	1:A:84:ASN:N	2.50	0.44
2:V:54:LEU:HA	2:V:57:ARG:NH1	2.33	0.44
1:A:39:MET:HE2	1:A:39:MET:HB3	1.67	0.43
2:U:152:GLU:O	2:U:156:GLU:HG3	2.18	0.43
2:U:161:THR:HA	2:U:164:ILE:HD12	2.00	0.43
2:V:32:GLU:OE1	2:U:185:THR:OG1	2.37	0.43
2:U:39:ASN:O	2:U:43:VAL:HG13	2.18	0.43
1:A:41:ILE:HA	1:A:51:GLY:O	2.18	0.43
2:U:110:SER:HA	2:U:113:LYS:HD2	2.01	0.43
2:U:176:ARG:O	2:U:180:ARG:HG3	2.19	0.43
2:V:109:ASP:O	2:V:113:LYS:HG3	2.18	0.43
1:A:55:LEU:HD23	1:A:72:PHE:HA	2.01	0.42
1:A:39:MET:HG2	1:A:40:ILE:N	2.34	0.42
2:U:81:GLU:HB2	2:U:93:CYS:HB3	2.01	0.42
2:U:184:PHE:O	2:U:188:ARG:HG3	2.19	0.42
1:A:13:PHE:CD1	1:A:17:LYS:HE3	2.55	0.42
2:V:174:ILE:H	2:V:174:ILE:HG13	1.52	0.42
2:U:103:LEU:O	2:U:107:ARG:HG2	2.19	0.42
1:A:50:ASP:CG	2:U:180:ARG:HH12	2.26	0.42
1:A:14:LYS:HE3	1:A:18:GLU:OE1	2.20	0.42
1:A:137:TYR:CD1	1:A:137:TYR:C	2.97	0.42
2:V:111:PHE:O	2:V:115:LEU:HG	2.20	0.42
2:V:29:PHE:HA	2:U:185:THR:CB	2.50	0.41
2:V:38:ILE:HG12	2:U:38:ILE:HG13	2.01	0.41
1:A:138:VAL:O	1:A:142:LYS:HB2	2.20	0.41
2:V:23:GLN:O	2:V:26:VAL:CG1	2.69	0.41
2:V:29:PHE:HA	2:U:185:THR:HB	2.02	0.41
2:V:170:TYR:CZ	2:V:174:ILE:HD11	2.55	0.41
2:U:184:PHE:N	2:U:184:PHE:CD1	2.87	0.41
1:A:133:HIS:ND1	1:A:136:GLN:OE1	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:PRO:HG2	1:A:24:VAL:CG2	2.50	0.41
2:V:59:CYS:SG	2:U:59:CYS:HB2	2.60	0.41
1:A:89:LEU:O	1:A:93:GLN:HG2	2.21	0.41
1:A:34:MET:HE2	1:A:34:MET:HB2	1.72	0.41
1:A:71:LYS:HB2	1:A:85:GLY:HA3	2.02	0.41
1:A:134:LYS:O	1:A:138:VAL:HG12	2.20	0.41
2:V:27:ILE:H	2:V:27:ILE:HG13	1.60	0.41
2:U:96:ILE:HG23	2:U:104:ILE:CD1	2.50	0.41
2:V:146:LEU:HD12	2:U:143:LEU:CD2	2.51	0.41
2:V:167:LEU:HD13	2:U:168:LYS:HG3	2.02	0.41
1:A:140:ARG:O	1:A:143:GLU:HG2	2.21	0.40
1:A:28:PRO:HA	1:A:36:TRP:HA	2.02	0.40
1:A:57:LEU:HG	1:A:70:VAL:HG12	2.03	0.40
1:A:55:LEU:HG	1:A:72:PHE:CD1	2.56	0.40
2:U:90:LYS:O	2:U:94:ARG:HG3	2.22	0.40
2:V:120:LYS:HB3	2:U:79:PHE:HZ	1.86	0.40
2:U:48:LEU:HD12	2:U:48:LEU:HA	1.83	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	147/150 (98%)	137 (93%)	10 (7%)	0	100	100
2	U	157/212 (74%)	156 (99%)	1 (1%)	0	100	100
2	V	156/212 (74%)	153 (98%)	3 (2%)	0	100	100
All	All	460/574 (80%)	446 (97%)	14 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	134/135 (99%)	131 (98%)	3 (2%)	45	74
2	U	149/197 (76%)	145 (97%)	4 (3%)	39	69
2	V	148/197 (75%)	143 (97%)	5 (3%)	32	62
All	All	431/529 (82%)	419 (97%)	12 (3%)	38	68

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	VAL
1	A	91	ILE
1	A	125	GLU
2	V	21	LEU
2	V	24	SER
2	V	27	ILE
2	V	86	ASP
2	V	180	ARG
2	U	26	VAL
2	U	27	ILE
2	U	174	ILE
2	U	182	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	ASN
1	A	110	GLN
2	V	51	GLN
2	V	145	ASN
2	U	23	GLN
2	U	98	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	149/150 (99%)	0.86	13 (8%) 16 13	90, 130, 162, 202	0
2	U	161/212 (75%)	0.72	18 (11%) 10 8	59, 90, 146, 171	0
2	V	160/212 (75%)	0.57	6 (3%) 44 40	61, 85, 139, 180	0
All	All	470/574 (81%)	0.71	37 (7%) 18 16	59, 103, 156, 202	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	U	80	ILE	6.3
1	A	113	PHE	6.0
2	U	83	PHE	5.0
2	U	79	PHE	4.8
2	U	29	PHE	4.7
2	U	21	LEU	3.9
2	U	22	THR	3.8
2	U	26	VAL	3.8
2	U	132	ALA	3.5
2	U	188	ARG	3.4
2	U	25	ASP	3.1
2	V	80	ILE	3.1
2	V	79	PHE	3.1
1	A	110	GLN	3.0
2	V	81	GLU	2.9
2	U	101	GLU	2.9
2	V	114	LEU	2.8
2	U	34	LEU	2.8
2	U	24	SER	2.8
1	A	87	ILE	2.7
1	A	72	PHE	2.7
1	A	5	ALA	2.6
1	A	57	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	109	ILE	2.5
2	U	84	CYS	2.5
1	A	91	ILE	2.4
1	A	92	LEU	2.4
2	U	23	GLN	2.4
2	V	83	PHE	2.4
1	A	70	VAL	2.3
2	U	27	ILE	2.3
2	V	116	THR	2.2
1	A	23	GLY	2.2
1	A	68	PRO	2.1
1	A	102	VAL	2.1
2	U	38	ILE	2.1
2	U	121	PRO	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.