



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:44 AM UTC

PDB ID : 1HQO / pdb_00001hqo
Title : CRYSTAL STRUCTURE OF THE NITROGEN REGULATION FRAGMENT OF THE YEAST PRION PROTEIN URE2P
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Deposited on : 2000-12-18
Resolution : 2.30 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

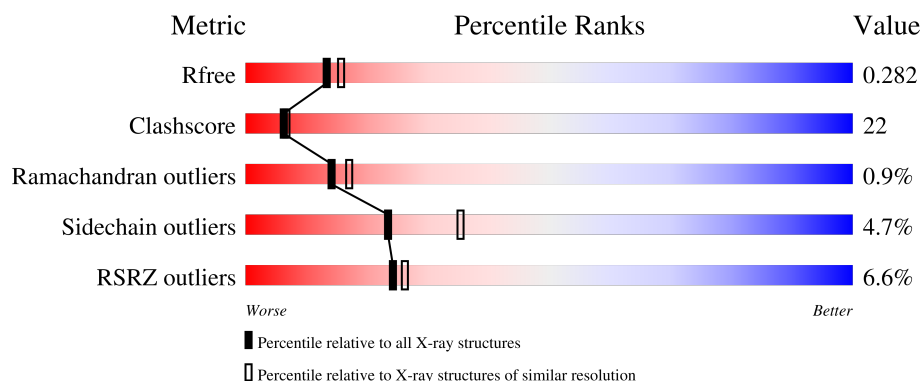
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	258	4% 53% 31% • • 12%
1	B	258	7% 49% 33% • • 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3868 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 URE2 PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	227	Total	C	N	O	Se	0	0	0
			1851	1201	318	326	6			
1	B	223	Total	C	N	O	Se	0	0	0
			1813	1177	312	318	6			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	173	MSE	MET	modified residue	UNP P23202
A	226	MSE	MET	modified residue	UNP P23202
A	261	MSE	MET	modified residue	UNP P23202
A	272	MSE	MET	modified residue	UNP P23202
A	289	MSE	MET	modified residue	UNP P23202
A	341	MSE	MET	modified residue	UNP P23202
A	342	MSE	MET	modified residue	UNP P23202
B	173	MSE	MET	modified residue	UNP P23202
B	226	MSE	MET	modified residue	UNP P23202
B	261	MSE	MET	modified residue	UNP P23202
B	272	MSE	MET	modified residue	UNP P23202
B	289	MSE	MET	modified residue	UNP P23202
B	341	MSE	MET	modified residue	UNP P23202
B	342	MSE	MET	modified residue	UNP P23202

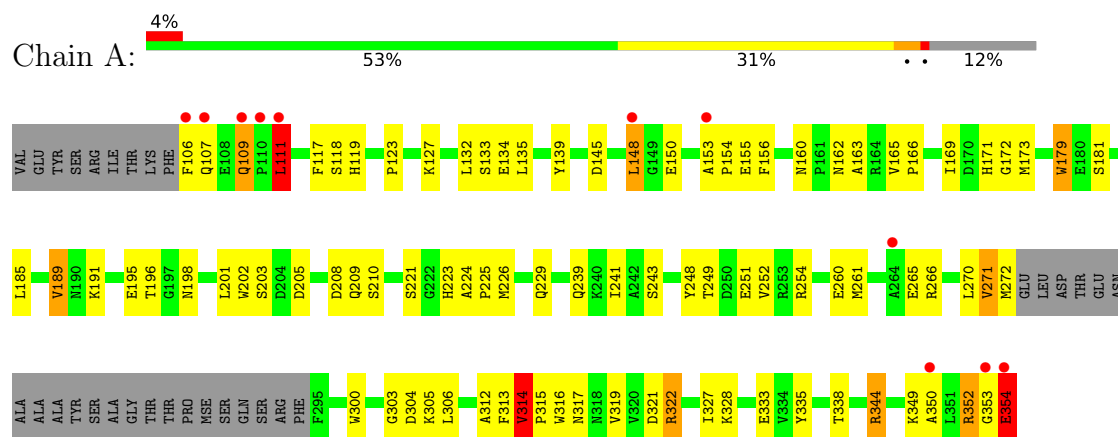
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	98	Total	O	0	0
			98	98		
2	B	106	Total	O	0	0
			106	106		

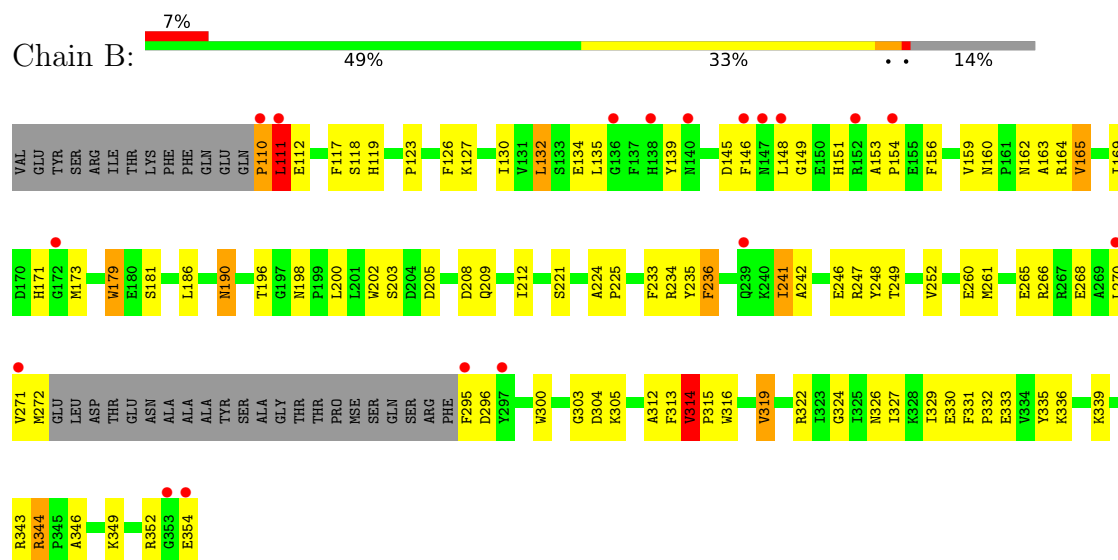
3 Residue-property plots [i](#)

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: URE2 PROTEIN



• Molecule 1: URE2 PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.31Å 69.19Å 149.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 20.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	88.8 (20.00-2.30) 94.6 (20.00-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 2.26Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.221 , 0.275 0.229 , 0.282	Depositor DCC
R_{free} test set	1568 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	47.0	Xtriage
Anisotropy	0.401	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 75.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3868	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 4.27% 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	0.71	0/1898	1.13	14/2567 (0.5%)
1	B	0.69	0/1859	1.13	12/2514 (0.5%)
All	All	0.70	0/3757	1.13	26/5081 (0.5%)

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	179	TRP	N-CA-C	8.68	124.07	112.88
1	B	111	LEU	N-CA-C	7.76	127.32	110.80
1	B	179	TRP	N-CA-C	7.29	121.89	112.92
1	A	109	GLN	CA-C-N	7.18	128.31	119.98
1	A	109	GLN	C-N-CA	7.18	128.31	119.98
1	B	331	PHE	CA-C-N	7.05	127.36	119.32
1	B	331	PHE	C-N-CA	7.05	127.36	119.32
1	A	352	ARG	N-CA-C	-6.83	103.84	111.28
1	B	236	PHE	N-CA-C	6.67	122.39	113.72
1	B	241	ILE	N-CA-C	6.65	118.10	107.73
1	A	319	VAL	N-CA-C	6.40	119.31	113.10
1	A	111	LEU	N-CA-C	6.32	124.27	110.80
1	B	132	LEU	N-CA-C	-6.15	104.27	110.97
1	B	153	ALA	CA-C-N	6.05	127.40	119.84
1	B	153	ALA	C-N-CA	6.05	127.40	119.84
1	A	314	VAL	N-CA-C	6.00	120.18	112.67
1	B	212	ILE	N-CA-C	-5.98	104.69	110.72
1	A	313	PHE	N-CA-C	5.91	119.67	112.23
1	B	313	PHE	N-CA-C	5.88	118.65	111.82
1	A	156	PHE	N-CA-C	-5.87	104.47	111.69
1	A	317	ASN	N-CA-C	-5.77	105.07	111.71
1	A	249	THR	N-CA-C	-5.77	105.13	111.82
1	A	148	LEU	N-CA-C	-5.68	106.20	113.02
1	B	314	VAL	N-CA-C	5.49	119.53	112.67
1	A	354	GLU	N-CA-C	-5.43	95.80	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	321	ASP	N-CA-C	-5.15	105.85	111.82

There are no chirality outliers.

There are no planarity outliers.

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	1851	0	1812	79	0
1	B	1813	0	1782	86	0
2	A	98	0	0	5	0
2	B	106	0	0	17	0
All	All	3868	0	3594	161	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 22.

All (161) 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:B:162:ASN:HB2	2:B:395:HOH:O	1.58	1.03
1:A:322:ARG:HA	1:A:322:ARG:HE	1.37	0.90
1:B:145:ASP:HB3	1:B:148:LEU:HG	1.54	0.87
1:B:314:VAL:HG12	1:B:315:PRO:HD3	1.56	0.85
1:A:226:MSE:HE3	2:A:402:HOH:O	1.81	0.79
1:A:314:VAL:HG12	1:A:315:PRO:HD3	1.65	0.78
1:B:117:PHE:HE1	1:B:169:ILE:CD1	1.99	0.76
1:A:205:ASP:HB3	1:A:208:ASP:HB2	1.69	0.75
1:A:106:PHE:HB3	1:A:107:GLN:OE1	1.88	0.74
1:B:151:HIS:ND1	2:B:420:HOH:O	2.20	0.74
1:A:109:GLN:HB2	1:A:171:HIS:CE1	2.23	0.73
1:B:344:ARG:HH11	1:B:344:ARG:HB3	1.52	0.73
1:B:205:ASP:O	1:B:209:GLN:HG3	1.87	0.73
1:B:354:GLU:HA	1:B:354:GLU:OE1	1.89	0.73
1:B:271:VAL:HG13	1:B:271:VAL:O	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:HIS:ND1	2:A:362:HOH:O	2.23	0.72
1:A:107:GLN:HB3	1:A:109:GLN:NE2	2.05	0.72
1:A:327:ILE:HG22	1:A:335:TYR:HB2	1.71	0.72
1:B:202:TRP:O	1:B:305:LYS:NZ	2.23	0.72
1:B:266:ARG:HD3	1:B:300:TRP:O	1.91	0.71
1:A:271:VAL:HG23	2:A:446:HOH:O	1.91	0.70
1:B:242:ALA:O	1:B:246:GLU:HG2	1.91	0.70
1:B:190:ASN:HA	2:B:390:HOH:O	1.92	0.70
1:B:110:PRO:O	1:B:112:GLU:N	2.23	0.69
1:A:191:LYS:O	1:A:195:GLU:HG3	1.93	0.69
1:A:221:SER:OG	1:B:221:SER:HB3	1.93	0.69
1:A:241:ILE:HG12	1:B:241:ILE:HD13	1.75	0.69
1:B:111:LEU:HD21	2:B:396:HOH:O	1.93	0.69
1:A:322:ARG:HE	1:A:322:ARG:CA	2.06	0.68
1:A:107:GLN:HB3	1:A:109:GLN:HE21	1.56	0.68
1:A:205:ASP:O	1:A:209:GLN:HG3	1.93	0.67
1:B:266:ARG:O	1:B:270:LEU:HG	1.97	0.65
1:A:145:ASP:N	1:A:150:GLU:OE1	2.29	0.65
1:A:117:PHE:HE1	1:A:169:ILE:HD12	1.62	0.65
1:A:344:ARG:HH11	1:A:344:ARG:HB3	1.62	0.65
1:B:260:GLU:OE2	1:B:333:GLU:HB2	1.97	0.64
1:B:205:ASP:HB3	1:B:208:ASP:HB2	1.79	0.64
1:B:248:TYR:O	1:B:252:VAL:HG23	1.98	0.64
1:A:111:LEU:HA	1:A:171:HIS:HB3	1.79	0.63
1:A:353:GLY:O	1:A:354:GLU:C	2.40	0.63
1:B:246:GLU:HB3	2:B:423:HOH:O	1.99	0.63
1:A:118:SER:HA	1:A:165:VAL:CG1	2.29	0.62
1:A:118:SER:HA	1:A:165:VAL:HG13	1.82	0.62
1:A:165:VAL:HG13	1:A:166:PRO:HA	1.81	0.61
1:A:189:VAL:HG22	1:A:202:TRP:HB2	1.81	0.61
1:B:202:TRP:NE1	2:B:390:HOH:O	2.31	0.61
1:A:160:ASN:ND2	1:A:179:TRP:HB2	2.16	0.61
1:B:146:PHE:HA	2:B:420:HOH:O	2.01	0.60
1:B:127:LYS:HD2	1:B:312:ALA:HA	1.84	0.60
1:A:266:ARG:HD3	1:A:300:TRP:O	2.01	0.60
1:A:145:ASP:H	1:A:150:GLU:CG	2.14	0.60
1:A:254:ARG:NH1	1:B:164:ARG:NH2	2.50	0.59
1:A:109:GLN:CB	1:A:171:HIS:CE1	2.86	0.59
1:B:165:VAL:HG22	2:B:458:HOH:O	2.01	0.58
1:B:179:TRP:H	1:B:179:TRP:CD1	2.22	0.58
1:B:268:GLU:O	1:B:271:VAL:HG12	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:261:MSE:O	1:B:265:GLU:HG3	2.04	0.58
1:A:162:ASN:O	1:A:163:ALA:HB3	2.04	0.57
1:A:248:TYR:O	1:A:252:VAL:HG23	2.03	0.57
1:A:111:LEU:O	1:A:111:LEU:HG	2.06	0.55
1:B:135:LEU:HD23	1:B:200:LEU:HD23	1.88	0.55
1:A:224:ALA:HB3	1:A:225:PRO:HD3	1.86	0.55
1:B:343:ARG:HD2	2:B:401:HOH:O	2.06	0.55
1:B:117:PHE:HE1	1:B:169:ILE:HD11	1.71	0.55
1:B:326:ASN:HB3	1:B:329:ILE:HG12	1.89	0.55
1:A:117:PHE:HE1	1:A:169:ILE:CD1	2.19	0.55
1:B:117:PHE:CE1	1:B:169:ILE:CD1	2.86	0.54
1:B:111:LEU:HD12	1:B:171:HIS:O	2.07	0.54
1:B:145:ASP:OD2	1:B:148:LEU:HD23	2.07	0.54
1:A:314:VAL:CG1	1:A:315:PRO:HD3	2.36	0.54
1:B:111:LEU:O	1:B:111:LEU:HG	2.07	0.54
1:B:149:GLY:HA2	2:B:420:HOH:O	2.07	0.53
1:B:327:ILE:HG22	1:B:335:TYR:HB2	1.89	0.53
1:A:123:PRO:HB2	1:A:316:TRP:CE2	2.43	0.53
1:B:156:PHE:O	1:B:159:VAL:HG22	2.08	0.53
1:A:127:LYS:HD2	1:A:312:ALA:HA	1.90	0.52
1:A:254:ARG:NE	2:A:394:HOH:O	2.41	0.52
1:B:111:LEU:HA	1:B:171:HIS:HB3	1.91	0.52
1:A:145:ASP:H	1:A:150:GLU:CD	2.17	0.52
1:A:266:ARG:O	1:A:270:LEU:HG	2.10	0.52
1:A:118:SER:OG	1:A:119:HIS:N	2.43	0.51
1:A:145:ASP:CG	1:A:148:LEU:HG	2.35	0.51
1:B:314:VAL:HG12	1:B:315:PRO:CD	2.35	0.50
1:B:123:PRO:HB2	1:B:316:TRP:CE2	2.47	0.50
1:B:162:ASN:O	1:B:163:ALA:HB3	2.11	0.50
1:A:171:HIS:C	1:A:173:MSE:H	2.20	0.49
1:B:118:SER:OG	1:B:119:HIS:N	2.46	0.49
1:B:260:GLU:CD	1:B:333:GLU:HB2	2.36	0.49
1:A:133:SER:HB3	1:A:349:LYS:HE2	1.93	0.49
1:B:314:VAL:CG1	1:B:315:PRO:HD3	2.37	0.49
1:B:233:PHE:HD2	2:B:358:HOH:O	1.96	0.49
1:B:186:LEU:O	1:B:190:ASN:HB2	2.13	0.49
1:B:295:PHE:CG	1:B:296:ASP:N	2.77	0.49
1:A:260:GLU:CD	1:A:333:GLU:HB2	2.37	0.48
1:B:132:LEU:HD13	1:B:139:TYR:CG	2.48	0.48
1:B:332:PRO:O	1:B:336:LYS:HG3	2.13	0.48
1:A:127:LYS:HE2	1:A:181:SER:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:TRP:CD1	1:B:179:TRP:N	2.81	0.48
1:B:171:HIS:C	1:B:173:MSE:N	2.71	0.48
1:B:135:LEU:CD2	1:B:200:LEU:HD23	2.43	0.48
1:B:196:THR:C	1:B:198:ASN:H	2.22	0.47
1:B:354:GLU:HB3	2:B:409:HOH:O	2.14	0.47
1:B:303:GLY:O	1:B:304:ASP:C	2.58	0.47
1:B:127:LYS:HE2	1:B:181:SER:HB2	1.95	0.47
1:B:202:TRP:CG	1:B:203:SER:H	2.33	0.47
1:A:314:VAL:HG23	1:A:338:THR:HG23	1.97	0.47
1:A:171:HIS:C	1:A:173:MSE:N	2.71	0.47
1:A:322:ARG:HA	1:A:322:ARG:NE	2.18	0.47
1:A:153:ALA:O	1:A:155:GLU:N	2.48	0.47
1:A:134:GLU:OE2	1:A:344:ARG:NH1	2.48	0.46
1:B:134:GLU:HG3	2:B:425:HOH:O	2.14	0.46
1:A:271:VAL:O	1:A:271:VAL:HG13	2.15	0.46
1:A:225:PRO:O	1:A:229:GLN:HG2	2.16	0.46
1:A:328:LYS:HG3	1:A:335:TYR:CE2	2.51	0.46
1:B:235:TYR:HB2	1:B:236:PHE:CE1	2.51	0.46
1:B:224:ALA:HB3	1:B:225:PRO:HD3	1.97	0.45
1:B:344:ARG:HH11	1:B:344:ARG:CB	2.25	0.45
1:B:247:ARG:HH11	1:B:247:ARG:HG3	1.82	0.45
1:B:171:HIS:O	1:B:173:MSE:N	2.49	0.45
1:A:202:TRP:CG	1:A:203:SER:H	2.35	0.45
1:B:339:LYS:O	1:B:343:ARG:HG3	2.16	0.45
1:A:172:GLY:HA3	2:A:421:HOH:O	2.16	0.45
1:A:303:GLY:O	1:A:304:ASP:C	2.60	0.45
1:B:117:PHE:CE1	1:B:169:ILE:HD11	2.50	0.45
1:A:352:ARG:HA	1:A:352:ARG:HD3	1.79	0.45
1:A:349:LYS:HG3	1:A:350:ALA:N	2.32	0.44
1:B:179:TRP:O	2:B:360:HOH:O	2.21	0.44
1:B:349:LYS:O	1:B:352:ARG:HB3	2.17	0.44
1:A:179:TRP:CD1	1:A:179:TRP:N	2.85	0.44
1:B:354:GLU:HB2	2:B:452:HOH:O	2.16	0.44
1:B:329:ILE:HG13	1:B:330:GLU:HG3	2.00	0.44
1:A:109:GLN:HB2	1:A:171:HIS:HE1	1.78	0.44
1:A:179:TRP:CD1	1:A:179:TRP:H	2.34	0.44
1:A:196:THR:C	1:A:198:ASN:H	2.24	0.44
1:A:132:LEU:HD13	1:A:139:TYR:CG	2.53	0.44
1:B:234:ARG:O	1:B:234:ARG:HG3	2.18	0.44
1:A:118:SER:HA	1:A:165:VAL:HG11	2.00	0.43
1:A:344:ARG:HH11	1:A:344:ARG:CB	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:GLU:OE2	1:A:254:ARG:NH1	2.52	0.43
1:A:241:ILE:HG21	1:B:241:ILE:CD1	2.48	0.43
1:B:171:HIS:C	1:B:173:MSE:H	2.26	0.43
1:A:241:ILE:HG22	1:A:243:SER:OG	2.19	0.43
1:B:196:THR:C	1:B:198:ASN:N	2.77	0.43
1:B:162:ASN:ND2	1:B:179:TRP:HE3	2.17	0.42
1:B:335:TYR:OH	1:B:339:LYS:HE3	2.19	0.42
1:B:160:ASN:HD22	1:B:179:TRP:HB2	1.84	0.42
1:B:319:VAL:O	1:B:322:ARG:HB2	2.20	0.42
1:A:305:LYS:HG2	1:A:306:LEU:N	2.33	0.42
1:B:271:VAL:O	1:B:271:VAL:CG1	2.60	0.42
1:B:126:PHE:O	1:B:130:ILE:HG13	2.19	0.42
1:A:135:LEU:HD21	1:A:201:LEU:HD12	2.02	0.41
1:B:110:PRO:N	2:B:432:HOH:O	2.54	0.41
1:B:324:GLY:HA3	2:B:373:HOH:O	2.20	0.41
1:A:171:HIS:O	1:A:173:MSE:N	2.54	0.41
1:A:314:VAL:CG1	1:A:315:PRO:CD	2.99	0.41
1:A:196:THR:C	1:A:198:ASN:N	2.78	0.41
1:A:261:MSE:O	1:A:265:GLU:HG3	2.20	0.41
1:A:261:MSE:HE3	1:A:261:MSE:HB2	1.92	0.40
1:B:134:GLU:OE1	1:B:346:ALA:HB3	2.20	0.40
1:A:145:ASP:OD2	1:A:148:LEU:HG	2.21	0.40
1:A:160:ASN:HD22	1:A:179:TRP:HB2	1.85	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	223/258 (86%)	207 (93%)	14 (6%)	2 (1%)	14	17
1	B	219/258 (85%)	203 (93%)	14 (6%)	2 (1%)	14	17

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	442/516 (86%)	410 (93%)	28 (6%)	4 (1%)	14 17

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	111	LEU
1	B	111	LEU
1	A	154	PRO
1	B	154	PRO

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	195/214 (91%)	185 (95%)	10 (5%)	21 32
1	B	191/214 (89%)	183 (96%)	8 (4%)	26 40
All	All	386/428 (90%)	368 (95%)	18 (5%)	23 35

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	185	LEU
1	A	189	VAL
1	A	210	SER
1	A	239	GLN
1	A	271	VAL
1	A	272	MSE
1	A	314	VAL
1	A	322	ARG
1	A	344	ARG
1	A	354	GLU
1	B	110	PRO
1	B	165	VAL
1	B	190	ASN
1	B	249	THR

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Mol	Chain	Res	Type
1	B	272	MSE
1	B	314	VAL
1	B	319	VAL
1	B	344	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	160	ASN
1	A	317	ASN
1	B	160	ASN
1	B	318	ASN

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	221/258 (85%)	0.43	11 (4%) 34 35	20, 52, 81, 93	0
1	B	217/258 (84%)	0.51	18 (8%) 17 19	26, 51, 81, 90	0
All	All	438/516 (84%)	0.47	29 (6%) 24 26	20, 52, 81, 93	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	146	PHE	4.3
1	A	353	GLY	4.2
1	B	354	GLU	3.7
1	B	271	VAL	3.3
1	B	136	GLY	3.1
1	A	354	GLU	3.0
1	A	110	PRO	3.0
1	B	138	HIS	2.9
1	B	110	PRO	2.9
1	B	152	ARG	2.9
1	B	239	GLN	2.7
1	B	270	LEU	2.7
1	B	148	LEU	2.7
1	A	153	ALA	2.6
1	B	154	PRO	2.6
1	B	295	PHE	2.5
1	B	297	TYR	2.5
1	A	106	PHE	2.4
1	B	111	LEU	2.4
1	A	264	ALA	2.3
1	B	140	ASN	2.3
1	B	353	GLY	2.3
1	B	147	ASN	2.2
1	A	111	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	109	GLN	2.1
1	B	172	GLY	2.1
1	A	350	ALA	2.0
1	A	107	GLN	2.0
1	A	148	LEU	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.