



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:53 PM GMT

PDB ID : 1HQO  
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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

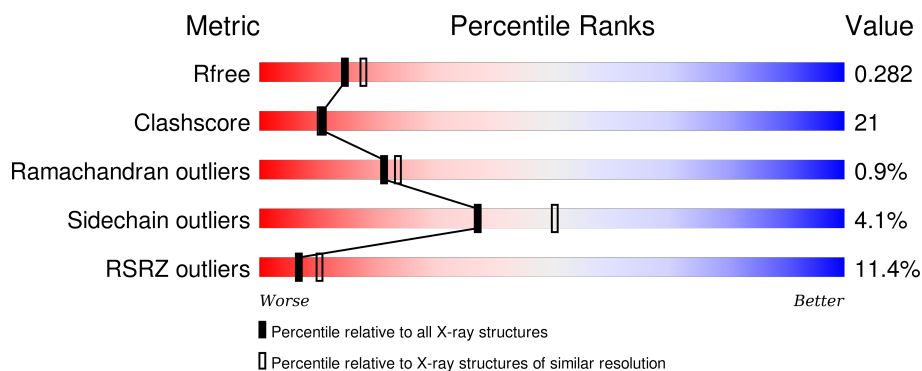
# 1 Overall quality at a glance

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))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	
1	B	258	

## 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

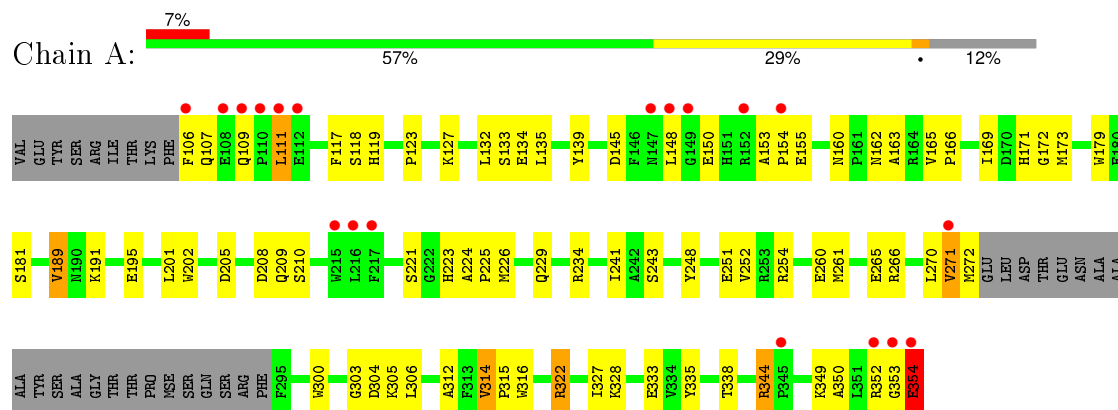
- Molecule 2 is water.

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

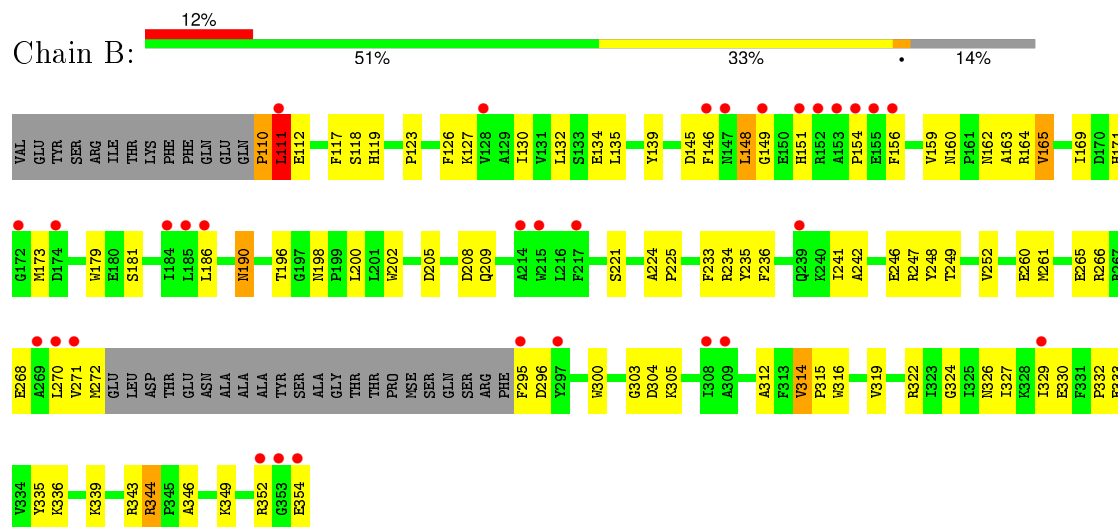
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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, $\alpha$ , $\beta$ , $\gamma$	64.31Å 69.19Å 149.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 29.34 – 2.25	Depositor EDS
% Data completeness (in resolution range)	88.8 (20.00-2.30) 94.1 (29.34-2.25)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	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	1427 reflections (4.95%)	DCC
Wilson B-factor (Å <sup>2</sup> )	47.0	Xtriage
Anisotropy	0.401	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 74.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 31560 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3868	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	1/1898 (0.1%)	0.74	1/2567 (0.0%)
1	B	0.57	0/1859	0.75	2/2514 (0.1%)
All	All	0.58	1/3757 (0.0%)	0.75	3/5081 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	354	GLU	CB-CG	5.40	1.62	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	111	LEU	N-CA-C	6.04	127.32	111.00
1	B	148	LEU	CA-CB-CG	5.87	128.80	115.30
1	A	354	GLU	N-CA-C	-5.63	95.80	111.00

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	75	0
1	B	1813	0	1782	82	0
2	A	98	0	0	5	0
2	B	106	0	0	17	0
All	All	3868	0	3594	153	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 (153) 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:HE	1:A:322:ARG:HA	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:OE1	1:B:354:GLU:HA	1.89	0.73
1:B:271:VAL:HG13	1:B:271:VAL:O	1.88	0.72
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:CA	1:A:322:ARG:HE	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:HB3	1:A:344:ARG:HH11	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:B:202:TRP:NE1	2:B:390:HOH:O	2.31	0.61
1:A:189:VAL:HG22	1:A:202:TRP:HB2	1.81	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:A:353:GLY:O	1:A:354:GLU:C	2.40	0.59
1:B:165:VAL:HG22	2:B:458:HOH:O	2.01	0.58
1:B:268:GLU:O	1:B:271:VAL:HG12	2.03	0.58
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:HG	1:A:111:LEU:O	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:111:LEU:O	1:B:111:LEU:HG	2.07	0.54
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: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:295:PHE:CG	1:B:296:ASP:N	2.77	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:B:118:SER:OG	1:B:119:HIS:N	2.46	0.49
1:B:314:VAL:CG1	1:B:315:PRO:HD3	2.37	0.49
1:A:133:SER:HB3	1:A:349:LYS:HE2	1.93	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: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
1:B:179:TRP:CD1	1:B:179:TRP:N	2.81	0.48
1:B:135:LEU:CD2	1:B:200:LEU:HD23	2.43	0.48
1:B:354:GLU:HB3	2:B:409:HOH:O	2.14	0.47
1:A:145:ASP:H	1:A:150:GLU:CD	2.17	0.47
1:B:127:LYS:HE2	1:B:181:SER:HB2	1.95	0.47
1:A:314:VAL:HG23	1:A:338:THR:HG23	1.97	0.47
1:A:322:ARG:NE	1:A:322:ARG:HA	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:145:ASP:CG	1:A:148:LEU:HG	2.35	0.46
1:A:271:VAL:HG13	1:A:271:VAL:O	2.15	0.46
1:A:328:LYS:HG3	1:A:335:TYR:CE2	2.51	0.46
1:A:225:PRO:O	1:A:229:GLN:HG2	2.16	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:B:339:LYS:O	1:B:343:ARG:HG3	2.16	0.45
1:B:117:PHE:CE1	1:B:169:ILE:HD11	2.50	0.45
1:A:171:HIS:C	1:A:173:MSE:H	2.20	0.45
1:A:172:GLY:HA3	2:A:421:HOH:O	2.16	0.45
1:A:352:ARG:HA	1:A:352:ARG:HD3	1.79	0.45
1:B:179:TRP:O	2:B:360:HOH:O	2.21	0.44
1:B:260:GLU:CD	1:B:333:GLU:HB2	2.36	0.44
1:A:349:LYS:HG3	1:A:350:ALA:N	2.32	0.44
1:B:349:LYS:O	1:B:352:ARG:HB3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:354:GLU:HB2	2:B:452:HOH:O	2.16	0.44
1:A:179:TRP:N	1:A:179:TRP:CD1	2.85	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:260:GLU:CD	1:A:333:GLU:HB2	2.37	0.44
1:A:132:LEU:HD13	1:A:139:TYR:CG	2.53	0.44
1:B:234:ARG:HG3	1:B:234:ARG:O	2.18	0.44
1:A:171:HIS:C	1:A:173:MSE:N	2.71	0.43
1:A:344:ARG:HH11	1:A:344:ARG:CB	2.30	0.43
1:A:118:SER:HA	1:A:165:VAL:HG11	2.00	0.43
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:N	2.71	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:H	2.22	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:B:271:VAL:O	1:B:271:VAL:CG1	2.60	0.42
1:B:303:GLY:O	1:B:304:ASP:C	2.58	0.42
1:A:305:LYS:HG2	1:A:306:LEU:N	2.33	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:314:VAL:CG1	1:A:315:PRO:CD	2.99	0.41
1:A:171:HIS:O	1:A:173:MSE:N	2.54	0.41
1:A:261:MSE:O	1:A:265:GLU:HG3	2.20	0.41
1:B:134:GLU:OE1	1:B:346:ALA:HB3	2.20	0.40
1:A:261:MSE:HE3	1:A:261:MSE:HB2	1.92	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
1:A:303:GLY:O	1:A:304:ASP:C	2.60	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	223/258 (86%)	207 (93%)	14 (6%)	2 (1%)	21	24
1	B	219/258 (85%)	203 (93%)	14 (6%)	2 (1%)	21	24
All	All	442/516 (86%)	410 (93%)	28 (6%)	4 (1%)	21	24

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%)	186 (95%)	9 (5%)	33	44
1	B	191/214 (89%)	184 (96%)	7 (4%)	41	55
All	All	386/428 (90%)	370 (96%)	16 (4%)	37	50

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

Mol	Chain	Res	Type
1	A	189	VAL
1	A	210	SER

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Mol	Chain	Res	Type
1	A	234	ARG
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
1	B	272	MSE
1	B	314	VAL
1	B	344	ARG

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

Mol	Chain	Res	Type
1	A	109	GLN
1	A	160	ASN
1	B	160	ASN
1	B	219	GLN

### 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 carbohydrates 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	221/258 (85%)	0.46	19 (8%) 13 18	20, 52, 81, 93	0
1	B	217/258 (84%)	0.64	31 (14%) 4 6	26, 51, 81, 90	0
All	All	438/516 (84%)	0.55	50 (11%) 7 10	20, 52, 81, 93	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	106	PHE	6.3
1	A	354	GLU	6.1
1	B	152	ARG	5.4
1	B	146	PHE	5.2
1	B	271	VAL	4.7
1	A	110	PRO	4.4
1	A	111	LEU	4.4
1	B	154	PRO	4.3
1	A	149	GLY	4.3
1	A	148	LEU	4.1
1	B	111	LEU	4.1
1	B	155	GLU	3.9
1	B	354	GLU	3.8
1	B	147	ASN	3.8
1	A	353	GLY	3.7
1	B	153	ALA	3.7
1	B	295	PHE	3.4
1	B	270	LEU	3.3
1	A	109	GLN	3.3
1	B	185	LEU	3.2
1	B	239	GLN	3.1
1	B	215	TRP	3.0
1	A	154	PRO	3.0
1	B	269	ALA	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	108	GLU	2.7
1	B	149	GLY	2.7
1	B	297	TYR	2.7
1	B	172	GLY	2.6
1	B	217	PHE	2.6
1	B	308	ILE	2.6
1	A	112	GLU	2.5
1	B	214	ALA	2.5
1	A	152	ARG	2.5
1	B	352	ARG	2.5
1	A	147	ASN	2.4
1	B	329	ILE	2.4
1	A	215	TRP	2.3
1	B	309	ALA	2.3
1	B	174	ASP	2.3
1	B	353	GLY	2.2
1	B	156	PHE	2.2
1	A	217	PHE	2.2
1	B	184	ILE	2.1
1	B	186	LEU	2.1
1	A	271	VAL	2.1
1	A	216	LEU	2.1
1	A	352	ARG	2.1
1	A	345	PRO	2.1
1	B	151	HIS	2.0
1	B	128	VAL	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 carbohydrates 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.