



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 01:33 am BST

PDB ID : 5GN2  
Title : Crystal structure of Uracil DNA glycosylase (BdiUNG) from Bradyrhizobium diazoefficiens  
Authors : Patil, V.V.; Chembazhi, U.V.; Varshney, U.; Woo, E.  
Deposited on : 2016-07-19  
Resolution : 1.95 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

<https://www.wwpdb.org/validation/2017/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  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

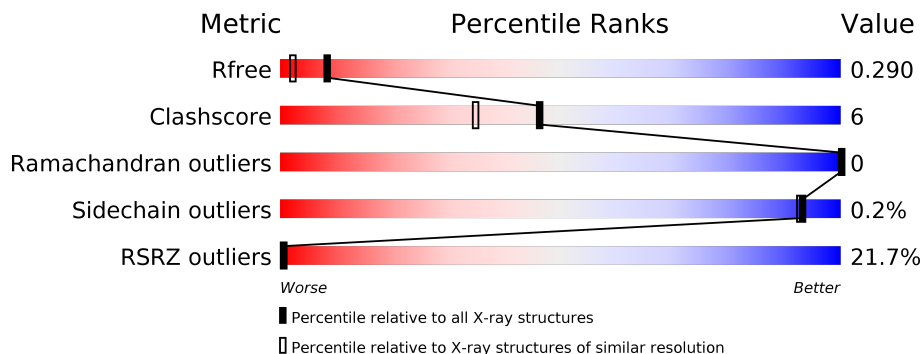
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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))
$R_{free}$	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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 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	272	
1	B	272	
1	C	272	
1	D	272	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8963 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 Blr0248 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	272	2099	1344	368	381	6	0	0	0
1	B	272	2099	1344	368	381	6	0	0	0
1	C	272	2099	1344	368	381	6	0	0	0
1	D	272	2099	1344	368	381	6	0	0	0

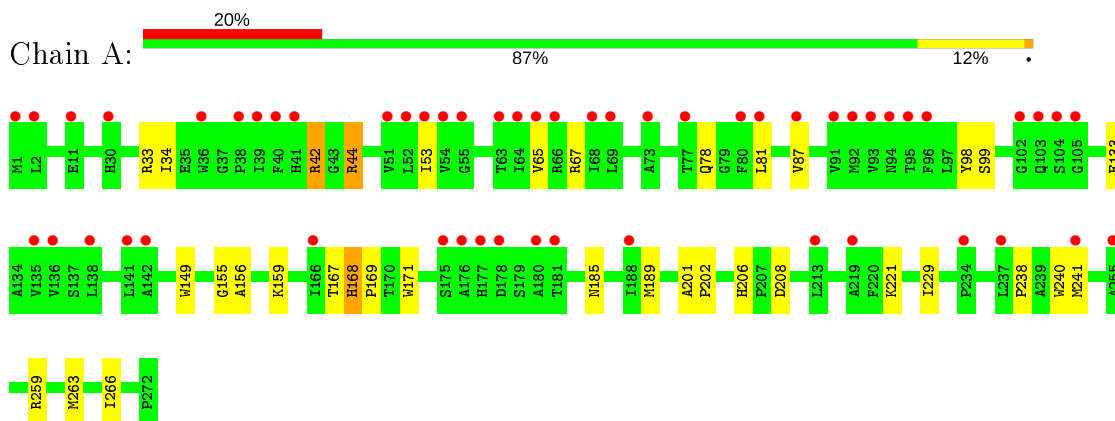
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	198	198	198	0	0
2	B	176	176	176	0	0
2	C	48	48	48	0	0
2	D	145	145	145	0	0

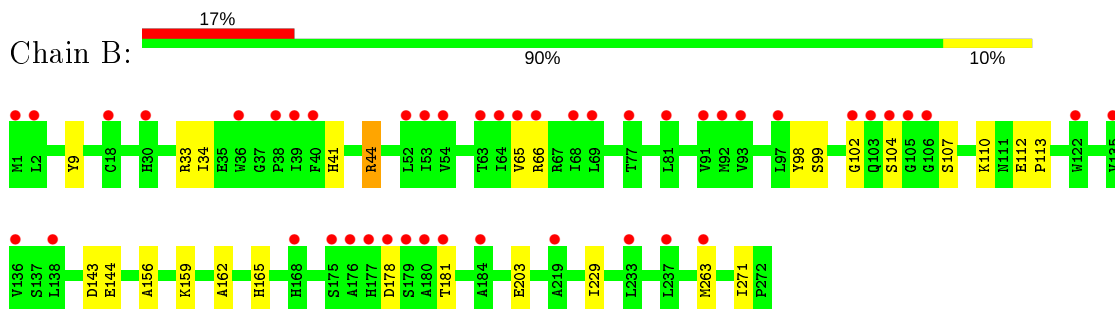
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

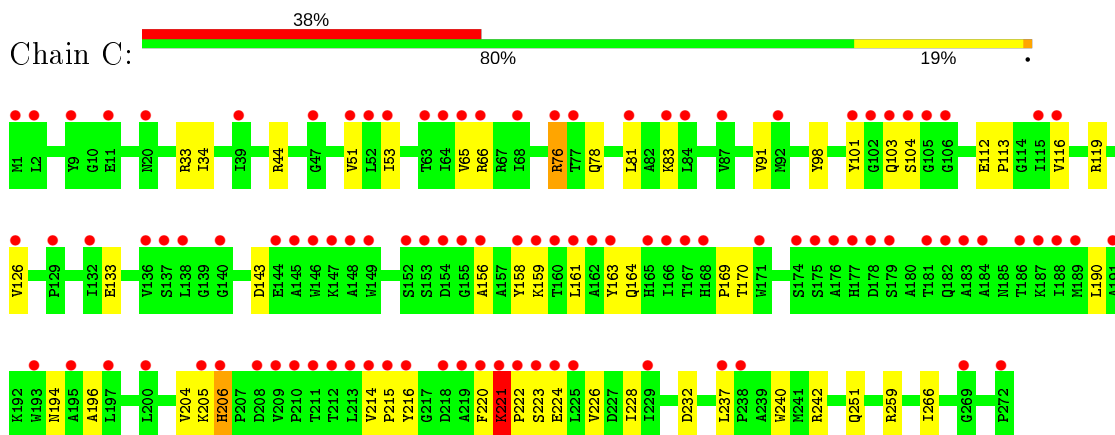
- Molecule 1: Blr0248 protein



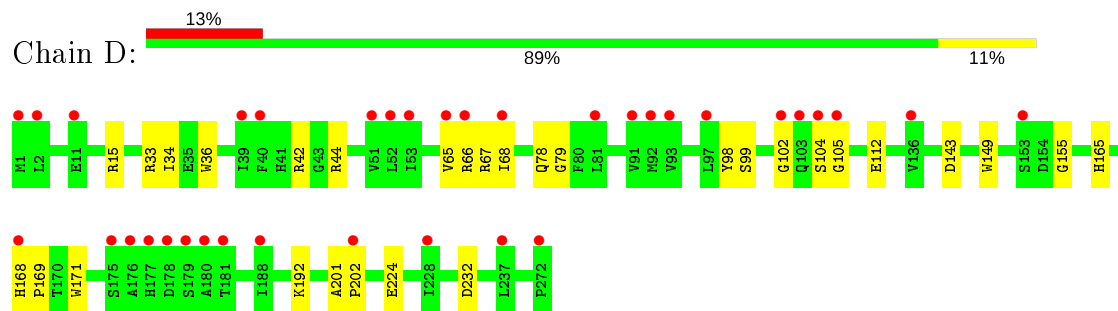
- Molecule 1: Blr0248 protein



- Molecule 1: Blr0248 protein



## ● Molecule 1: Blr0248 protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.65Å 90.03Å 255.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.52 – 1.95 30.52 – 1.95	Depositor EDS
% Data completeness (in resolution range)	98.7 (30.52-1.95) 98.7 (30.52-1.95)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.99 (at 1.95Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.245 , 0.290 0.251 , 0.290	Depositor DCC
$R_{free}$ test set	1994 reflections (1.69%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	33.2	Xtrriage
Anisotropy	0.748	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 49.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8963	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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, 0.333 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	0.47	1/2159 (0.0%)	0.63	3/2944 (0.1%)
1	B	0.51	0/2159	0.62	1/2944 (0.0%)
1	C	0.50	1/2159 (0.0%)	0.67	5/2944 (0.2%)
1	D	0.45	0/2159	0.62	1/2944 (0.0%)
All	All	0.48	2/8636 (0.0%)	0.64	10/11776 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	222	PRO	N-CD	5.25	1.55	1.47
1	A	169	PRO	N-CD	5.12	1.55	1.47

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	44	ARG	NE-CZ-NH2	-8.31	116.15	120.30
1	C	206	HIS	C-N-CD	6.06	141.13	128.40
1	A	42	ARG	NE-CZ-NH1	-5.93	117.33	120.30
1	A	168	HIS	C-N-CD	5.53	140.02	128.40
1	A	44	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	C	221	LYS	C-N-CD	5.43	139.81	128.40
1	C	237	LEU	C-N-CD	5.41	139.76	128.40
1	D	102	GLY	N-CA-C	-5.29	99.87	113.10
1	C	76	ARG	CG-CD-NE	5.13	122.57	111.80
1	C	76	ARG	CB-CG-CD	5.01	124.62	111.60

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	2099	0	2072	21	0
1	B	2099	0	2072	18	0
1	C	2099	0	2069	48	0
1	D	2099	0	2072	21	0
2	A	198	0	0	1	0
2	B	176	0	0	1	0
2	C	48	0	0	2	0
2	D	145	0	0	3	0
All	All	8963	0	8285	105	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 6.

All (105) 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:C:216:TYR:CE1	1:C:224:GLU:OE2	1.67	1.44
1:C:216:TYR:HE1	1:C:224:GLU:OE2	0.92	1.23
1:C:216:TYR:HE1	1:C:224:GLU:CD	1.46	1.18
1:C:220:PHE:HD2	1:C:224:GLU:HG2	1.30	0.95
1:C:216:TYR:CE1	1:C:224:GLU:CD	2.28	0.92
1:C:216:TYR:CZ	1:C:224:GLU:OE2	2.25	0.89
1:D:42:ARG:NH2	1:D:78:GLN:OE1	2.08	0.87
1:C:220:PHE:CD2	1:C:224:GLU:HG2	2.16	0.79
1:D:44:ARG:NH2	1:D:232:ASP:OD1	2.15	0.79
1:D:104:SER:H	1:D:105:GLY:HA2	1.46	0.77
1:B:110:LYS:NZ	1:B:144:GLU:OE1	2.18	0.75
1:C:51:VAL:HG21	1:C:81:LEU:HD22	1.68	0.73
1:D:104:SER:N	1:D:105:GLY:HA2	2.07	0.69
1:D:143:ASP:HB2	1:D:165:HIS:HD2	1.60	0.66
1:D:42:ARG:HG3	1:D:67:ARG:HB3	1.80	0.64
1:C:220:PHE:HD2	1:C:224:GLU:CG	2.09	0.63
1:C:44:ARG:NH2	1:C:232:ASP:OD1	2.33	0.62
1:B:143:ASP:HB2	1:B:165:HIS:HD2	1.64	0.61
1:C:53:ILE:HD11	1:C:81:LEU:HD21	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:76:ARG:NH2	1:C:169:PRO:O	2.37	0.58
1:B:178:ASP:OD2	1:B:181:THR:OG1	2.12	0.58
1:A:42:ARG:NH1	1:A:78:GLN:OE1	2.25	0.57
1:C:216:TYR:OH	1:C:224:GLU:OE2	2.22	0.57
1:A:53:ILE:HD11	1:A:81:LEU:HD21	1.87	0.56
1:D:104:SER:N	1:D:105:GLY:CA	2.68	0.56
1:C:251:GLN:NE2	2:C:305:HOH:O	2.38	0.56
1:C:158:TYR:HD1	1:C:161:LEU:HB2	1.71	0.56
1:C:204:VAL:HG12	1:C:204:VAL:O	2.05	0.55
1:D:149:TRP:O	1:D:155:GLY:HA3	2.07	0.55
1:D:33:ARG:HD2	1:D:99:SER:HB3	1.89	0.54
1:A:34:ILE:HA	1:A:98:TYR:CD1	2.42	0.54
1:B:143:ASP:HB2	1:B:165:HIS:CD2	2.43	0.54
1:D:34:ILE:HA	1:D:98:TYR:CD1	2.43	0.54
1:C:156:ALA:O	1:C:159:LYS:HG3	2.08	0.54
1:C:143:ASP:OD1	1:C:163:TYR:OH	2.21	0.53
1:C:78:GLN:HB3	1:C:226:VAL:HG13	1.91	0.52
1:A:238:PRO:HD2	1:A:241:MET:SD	2.49	0.52
1:C:158:TYR:CD1	1:C:161:LEU:HD22	2.44	0.51
1:C:116:VAL:HG12	1:C:119:ARG:NH2	2.26	0.51
1:B:102:GLY:C	1:B:104:SER:H	2.15	0.51
1:D:42:ARG:HD3	1:D:68:ILE:O	2.11	0.50
1:A:168:HIS:O	1:A:189:MET:HE1	2.11	0.50
1:C:78:GLN:CB	1:C:226:VAL:HG13	2.42	0.50
1:B:33:ARG:HD2	1:B:99:SER:HB3	1.94	0.50
1:C:116:VAL:HG12	1:C:119:ARG:HH22	1.77	0.49
1:A:259:ARG:HG2	1:B:271:ILE:HD13	1.95	0.49
1:A:44:ARG:NH2	1:A:229:ILE:HD11	2.28	0.49
1:C:83:LYS:HE3	1:C:214:VAL:O	2.12	0.49
1:C:34:ILE:HA	1:C:98:TYR:CD1	2.49	0.48
1:C:103:GLN:O	1:C:104:SER:OG	2.23	0.48
1:D:168:HIS:ND1	2:D:301:HOH:O	2.11	0.48
1:A:263:MET:SD	1:B:263:MET:HG3	2.54	0.47
1:B:162:ALA:CB	1:B:203:GLU:HG2	2.44	0.47
1:A:149:TRP:O	1:A:155:GLY:HA3	2.15	0.47
1:C:216:TYR:CE1	1:C:224:GLU:OE1	2.65	0.47
1:B:34:ILE:HA	1:B:98:TYR:CD1	2.49	0.47
1:C:101:TYR:CD2	1:C:259:ARG:NH1	2.83	0.46
1:B:44:ARG:NH2	1:B:229:ILE:HD11	2.30	0.46
1:B:65:VAL:O	1:B:66:ARG:HB2	2.16	0.46
1:C:44:ARG:HA	1:C:44:ARG:HD3	1.83	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:221:LYS:NZ	2:A:309:HOH:O	2.45	0.45
1:D:168:HIS:CE1	1:D:171:TRP:HB2	2.52	0.45
1:A:156:ALA:O	1:A:159:LYS:HG3	2.17	0.45
1:C:91:VAL:HG11	1:C:126:VAL:HG11	1.99	0.45
1:C:91:VAL:HG11	1:C:126:VAL:CG1	2.46	0.45
1:C:76:ARG:NH2	1:C:170:THR:HA	2.32	0.45
1:A:167:THR:HG22	1:A:189:MET:HE2	1.98	0.45
1:B:44:ARG:HD3	1:B:44:ARG:HA	1.81	0.44
1:C:76:ARG:HD3	2:C:334:HOH:O	2.17	0.44
1:C:205:LYS:O	1:C:206:HIS:CG	2.70	0.44
1:C:240:TRP:CZ2	1:C:266:ILE:HG12	2.53	0.44
1:C:33:ARG:HG3	1:C:101:TYR:HB2	2.00	0.44
1:C:190:LEU:HD21	1:C:220:PHE:CE2	2.52	0.44
1:B:156:ALA:HA	1:B:159:LYS:HE3	1.99	0.44
1:C:112:GLU:HA	1:C:113:PRO:HD3	1.90	0.43
1:A:201:ALA:HB3	1:A:202:PRO:HD3	2.00	0.43
1:C:133:GLU:OE1	1:C:206:HIS:ND1	2.37	0.43
1:D:65:VAL:O	1:D:66:ARG:HB2	2.18	0.43
1:D:112:GLU:OE1	2:D:302:HOH:O	2.22	0.43
1:C:65:VAL:O	1:C:66:ARG:HB2	2.18	0.42
1:D:15:ARG:NH2	2:D:305:HOH:O	2.30	0.42
1:A:171:TRP:CZ3	1:A:185:ASN:HB3	2.54	0.42
1:B:112:GLU:HA	1:B:113:PRO:HD3	1.81	0.42
1:C:83:LYS:HE2	1:C:83:LYS:HB2	1.79	0.42
1:B:9:TYR:CE1	1:B:41:HIS:CE1	3.07	0.42
1:A:33:ARG:HD2	1:A:99:SER:HB3	2.01	0.42
1:D:79:GLY:HA3	1:D:224:GLU:O	2.19	0.42
1:A:65:VAL:HG22	1:A:241:MET:HE2	2.01	0.42
1:D:192:LYS:HB2	1:D:192:LYS:HE3	1.64	0.42
1:C:164:GLN:HG2	1:C:196:ALA:HA	2.02	0.42
1:B:98:TYR:HA	2:B:333:HOH:O	2.19	0.41
1:C:194:ASN:OD1	1:C:216:TYR:HB2	2.20	0.41
1:D:201:ALA:HB3	1:D:202:PRO:HD3	2.02	0.41
1:C:101:TYR:CZ	1:C:259:ARG:HD2	2.56	0.41
1:A:42:ARG:HG3	1:A:67:ARG:HB3	2.02	0.41
1:C:228:ILE:O	1:C:242:ARG:HD2	2.20	0.41
1:D:168:HIS:HA	1:D:169:PRO:HD3	1.89	0.41
1:A:87:VAL:HG22	1:A:208:ASP:CB	2.51	0.41
1:B:107:SER:HB2	1:B:110:LYS:HD3	2.03	0.41
1:C:266:ILE:HD11	1:D:36:TRP:CE2	2.56	0.40
1:A:133:GLU:OE2	1:A:206:HIS:ND1	2.42	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:THR:HG22	1:A:189:MET:CE	2.51	0.40
1:A:240:TRP:CZ2	1:A:266:ILE:HG12	2.56	0.40
1:C:214:VAL:HA	1:C:215:PRO:HD2	1.91	0.40
1:C:221:LYS:O	1:C:224:GLU:HB3	2.21	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	270/272 (99%)	263 (97%)	7 (3%)	0	100	100
1	B	270/272 (99%)	260 (96%)	10 (4%)	0	100	100
1	C	270/272 (99%)	255 (94%)	15 (6%)	0	100	100
1	D	270/272 (99%)	263 (97%)	7 (3%)	0	100	100
All	All	1080/1088 (99%)	1041 (96%)	39 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	A	213/213 (100%)	213 (100%)	0	100	100
1	B	213/213 (100%)	213 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	213/213 (100%)	211 (99%)	2 (1%)	78	77
1	D	213/213 (100%)	213 (100%)	0	100	100
All	All	852/852 (100%)	850 (100%)	2 (0%)	93	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	221	LYS
1	C	223	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	165	HIS

### 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 carbohydrates 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

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	272/272 (100%)	1.24	54 (19%) <b>1</b> <b>1</b>	28, 40, 60, 87	0
1	B	272/272 (100%)	1.12	45 (16%) <b>1</b> <b>2</b>	30, 40, 57, 94	0
1	C	272/272 (100%)	1.91	102 (37%) <b>0</b> <b>0</b>	36, 62, 86, 94	0
1	D	272/272 (100%)	1.14	35 (12%) <b>3</b> <b>5</b>	33, 43, 59, 95	0
All	All	1088/1088 (100%)	1.35	236 (21%) <b>0</b> <b>1</b>	28, 43, 78, 95	0

All (236) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	176	ALA	9.7
1	C	104	SER	8.9
1	D	1	MET	8.7
1	A	104	SER	8.3
1	C	105	GLY	8.2
1	A	103	GLN	7.6
1	D	103	GLN	6.9
1	C	103	GLN	6.9
1	C	209	VAL	6.8
1	D	102	GLY	6.7
1	B	103	GLN	6.4
1	B	1	MET	6.3
1	C	175	SER	6.1
1	C	188	ILE	5.8
1	C	219	ALA	5.7
1	A	177	HIS	5.6
1	C	184	ALA	5.6
1	C	171	TRP	5.6
1	D	104	SER	5.6
1	C	177	HIS	5.5
1	D	2	LEU	5.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	B	104	SER	5.2
1	B	176	ALA	5.1
1	C	210	PRO	5.0
1	B	53	ILE	5.0
1	A	53	ILE	4.9
1	C	102	GLY	4.8
1	B	177	HIS	4.7
1	C	1	MET	4.7
1	C	81	LEU	4.7
1	C	205	LYS	4.7
1	A	54	VAL	4.7
1	C	181	THR	4.6
1	D	177	HIS	4.6
1	A	1	MET	4.6
1	C	224	GLU	4.5
1	A	69	LEU	4.3
1	C	65	VAL	4.2
1	A	68	ILE	4.2
1	C	2	LEU	4.2
1	C	222	PRO	4.1
1	B	175	SER	4.1
1	A	176	ALA	4.1
1	D	176	ALA	4.0
1	A	138	LEU	4.0
1	D	65	VAL	4.0
1	C	206	HIS	4.0
1	B	68	ILE	3.9
1	C	167	THR	3.9
1	B	65	VAL	3.9
1	C	106	GLY	3.9
1	C	53	ILE	3.9
1	A	237	LEU	3.9
1	C	211	THR	3.9
1	D	181	THR	3.7
1	D	175	SER	3.7
1	C	83	LYS	3.7
1	C	147	LYS	3.7
1	C	159	LYS	3.7
1	A	81	LEU	3.7
1	A	93	VAL	3.6
1	C	214	VAL	3.6
1	B	39	ILE	3.6

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	D	39	ILE	3.6
1	C	158	TYR	3.6
1	C	51	VAL	3.5
1	B	178	ASP	3.5
1	C	191	ALA	3.5
1	A	77	THR	3.5
1	A	178	ASP	3.5
1	C	152	SER	3.4
1	B	52	LEU	3.4
1	B	237	LEU	3.4
1	C	39	ILE	3.4
1	A	66	ARG	3.4
1	C	178	ASP	3.4
1	C	20	ASN	3.3
1	D	40	PHE	3.3
1	C	52	LEU	3.3
1	D	11	GLU	3.3
1	B	93	VAL	3.3
1	C	237	LEU	3.3
1	C	182	GLN	3.3
1	A	55	GLY	3.3
1	A	141	LEU	3.3
1	A	64	ILE	3.2
1	A	39	ILE	3.2
1	C	116	VAL	3.2
1	C	146	TRP	3.2
1	C	149	TRP	3.2
1	B	64	ILE	3.1
1	C	208	ASP	3.1
1	C	225	LEU	3.1
1	B	180	ALA	3.1
1	D	272	PRO	3.1
1	D	68	ILE	3.1
1	C	165	HIS	3.1
1	C	155	GLY	3.1
1	A	11	GLU	3.0
1	C	166	ILE	3.0
1	A	52	LEU	3.0
1	C	168	HIS	3.0
1	B	2	LEU	2.9
1	A	102	GLY	2.9
1	C	189	MET	2.9

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	B	168	HIS	2.9
1	A	65	VAL	2.9
1	B	105	GLY	2.9
1	C	195	ALA	2.9
1	A	95	THR	2.9
1	B	219	ALA	2.9
1	B	122	TRP	2.9
1	A	2	LEU	2.9
1	D	237	LEU	2.9
1	C	221	LYS	2.9
1	B	92	MET	2.9
1	C	156	ALA	2.9
1	D	180	ALA	2.9
1	D	51	VAL	2.8
1	C	64	ILE	2.8
1	B	38	PRO	2.8
1	C	193	TRP	2.8
1	B	54	VAL	2.8
1	C	126	VAL	2.8
1	C	136	VAL	2.8
1	B	81	LEU	2.8
1	C	92	MET	2.8
1	B	69	LEU	2.7
1	A	92	MET	2.7
1	B	102	GLY	2.7
1	C	87	VAL	2.7
1	C	161	LEU	2.7
1	D	92	MET	2.7
1	A	166	ILE	2.7
1	C	68	ILE	2.7
1	C	162	ALA	2.7
1	B	40	PHE	2.7
1	C	84	LEU	2.7
1	C	229	ILE	2.6
1	C	216	TYR	2.6
1	D	179	SER	2.6
1	C	101	TYR	2.6
1	A	30	HIS	2.6
1	C	11	GLU	2.6
1	A	40	PHE	2.6
1	A	136	VAL	2.6
1	C	186	THR	2.6

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	80	PHE	2.5
1	C	215	PRO	2.5
1	A	87	VAL	2.5
1	A	63	THR	2.5
1	C	163	TYR	2.5
1	C	197	LEU	2.5
1	C	66	ARG	2.5
1	C	77	THR	2.5
1	B	30	HIS	2.5
1	B	97	LEU	2.5
1	C	153	SER	2.5
1	D	153	SER	2.5
1	A	51	VAL	2.5
1	D	178	ASP	2.5
1	C	137	SER	2.5
1	B	184	ALA	2.4
1	A	188	ILE	2.4
1	D	53	ILE	2.4
1	D	168	HIS	2.4
1	B	136	VAL	2.4
1	A	94	ASN	2.4
1	C	212	THR	2.4
1	B	138	LEU	2.4
1	B	233	LEU	2.4
1	A	135	VAL	2.4
1	C	63	THR	2.4
1	C	183	ALA	2.4
1	D	93	VAL	2.4
1	A	41	HIS	2.4
1	C	138	LEU	2.4
1	C	144	GLU	2.4
1	C	148	ALA	2.4
1	B	91	VAL	2.4
1	D	52	LEU	2.4
1	B	179	SER	2.3
1	C	179	SER	2.3
1	C	223	SER	2.3
1	C	200	LEU	2.3
1	A	105	GLY	2.3
1	A	175	SER	2.3
1	D	136	VAL	2.3
1	C	220	PHE	2.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	C	154	ASP	2.3
1	C	132	ILE	2.3
1	B	63	THR	2.3
1	B	181	THR	2.3
1	A	38	PRO	2.3
1	A	219	ALA	2.3
1	A	181	THR	2.3
1	C	160	THR	2.3
1	C	218	ASP	2.3
1	A	73	ALA	2.2
1	C	213	LEU	2.2
1	C	145	ALA	2.2
1	C	47	GLY	2.2
1	D	105	GLY	2.2
1	A	234	PRO	2.2
1	C	174	SER	2.2
1	A	255	ALA	2.2
1	C	9	TYR	2.2
1	C	269	GLY	2.2
1	C	115	ILE	2.1
1	A	241	MET	2.1
1	B	263	MET	2.1
1	A	36	TRP	2.1
1	B	36	TRP	2.1
1	C	187	LYS	2.1
1	B	135	VAL	2.1
1	C	272	PRO	2.1
1	A	213	LEU	2.1
1	D	81	LEU	2.1
1	A	96	PHE	2.1
1	A	142	ALA	2.1
1	B	18	CYS	2.1
1	D	188	ILE	2.1
1	B	106	GLY	2.1
1	C	140	GLY	2.1
1	A	180	ALA	2.1
1	C	129	PRO	2.1
1	C	238	PRO	2.1
1	D	228	ILE	2.1
1	C	76	ARG	2.0
1	A	91	VAL	2.0
1	D	91	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	202	PRO	2.0
1	D	66	ARG	2.0
1	D	97	LEU	2.0
1	B	66	ARG	2.0
1	B	77	THR	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.