



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:56 am GMT

PDB ID : 1NUI  
Title : Crystal Structure of the primase fragment of Bacteriophage T7 primase-helicase protein  
Authors : Kato, M.; Ito, T.; Wagner, G.; Richardson, C.C.; Ellenberger, T.  
Deposited on : 2003-01-31  
Resolution : 2.90 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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

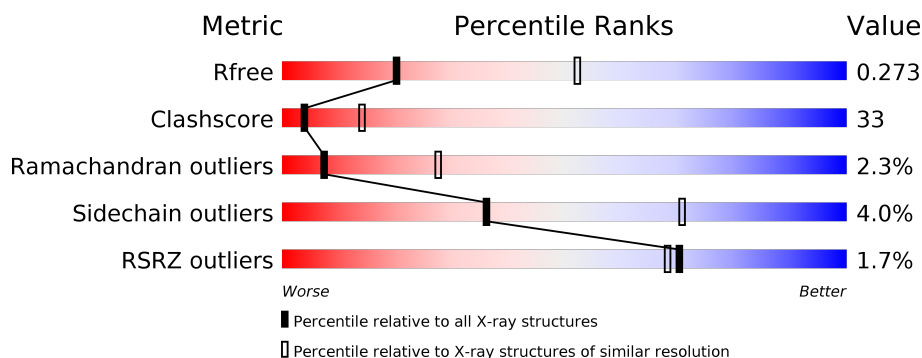
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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))
$R_{free}$	100719	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

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	255	<div> <div> <div>0%</div> <div>39%</div> <div>52%</div> <div>5%</div> </div> </div>
1	B	255	<div> <div>2%</div> <div>53%</div> <div>38%</div> <div>6%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3783 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 DNA primase/helicase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	0	0
			1891	1184	331	361	15			
1	B	239	Total	C	N	O	S	0	0	0
			1872	1173	328	356	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	64	GLY	MET	ENGINEERED	UNP P03692
B	64	GLY	MET	ENGINEERED	UNP P03692

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Mg	0	0
			2	2		
3	A	2	Total	Mg	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	O	0	0
			4	4		

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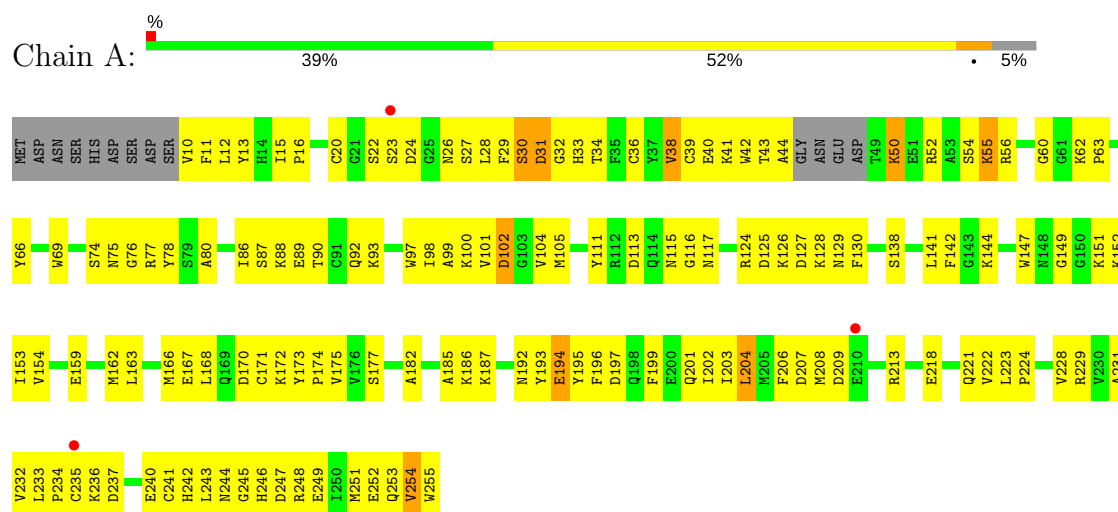
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	10	Total	O	0	0
			10	10		

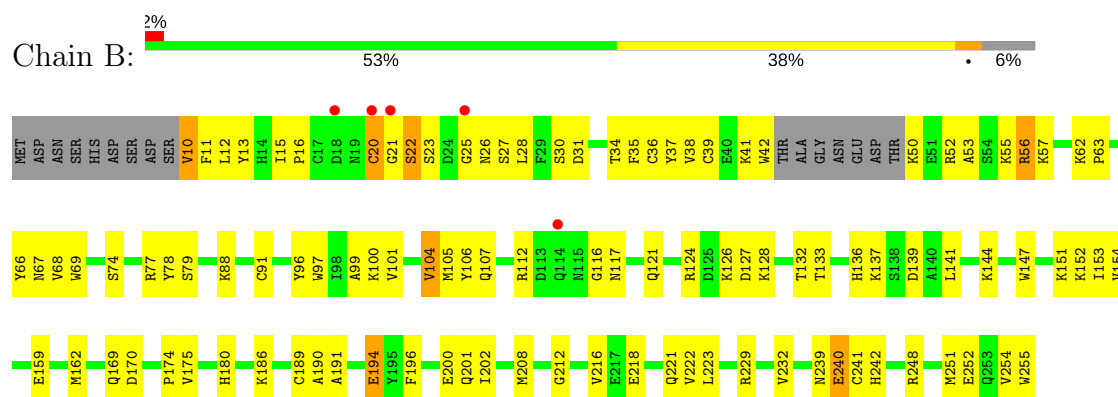
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: DNA primase/helicase



#### • Molecule 1: DNA primase/helicase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	136.46Å 136.46Å 85.44Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.90 53.32 – 2.58	Depositor EDS
% Data completeness (in resolution range)	99.8 (10.00-2.90) 98.4 (53.32-2.58)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.58 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, $R_{free}$	0.235 , 0.277 0.231 , 0.273	Depositor DCC
$R_{free}$ test set	1048 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	54.0	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 58.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	3783	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MG

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	0.40	0/1931	0.61	0/2599
1	B	0.44	0/1912	0.63	0/2572
All	All	0.42	0/3843	0.62	0/5171

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	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	66	TYR	Sidechain

### 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	1891	0	1821	159	0
1	B	1872	0	1802	103	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	4	0	0	0	0
4	B	10	0	0	0	0
All	All	3783	0	3623	246	1

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 33.

All (246) 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:56:ARG:HB2	1:B:56:ARG:HH11	1.13	1.08
1:A:20:CYS:HB3	1:A:39:CYS:HB3	1.45	0.95
1:A:10:VAL:N	1:A:30:SER:HG	1.68	0.92
1:B:56:ARG:HB2	1:B:56:ARG:NH1	1.83	0.91
1:B:11:PHE:HE1	1:B:27:SER:HG	1.25	0.84
1:A:55:LYS:HA	1:B:116:GLY:HA3	1.60	0.84
1:A:208:MET:HE3	1:A:233:LEU:H	1.42	0.82
1:B:41:LYS:NZ	1:B:41:LYS:HB3	1.94	0.82
1:A:13:TYR:HE1	1:A:15:ILE:HD11	1.48	0.77
1:A:233:LEU:HB3	1:A:241:CYS:SG	2.26	0.76
1:A:56:ARG:HG2	1:A:56:ARG:HH11	1.51	0.76
1:B:10:VAL:HA	1:B:30:SER:OG	1.87	0.74
1:A:208:MET:CE	1:A:233:LEU:H	2.00	0.73
1:A:93:LYS:HD3	1:B:52:ARG:HH22	1.54	0.72
1:B:248:ARG:O	1:B:252:GLU:HG3	1.90	0.72
1:A:105:MET:HB2	1:A:126:LYS:HE3	1.71	0.72
1:A:115:ASN:HB2	1:A:117:ASN:ND2	2.05	0.71
1:B:201:GLN:HE22	1:B:229:ARG:HH11	1.35	0.71
1:A:76:GLY:HA2	1:A:98:ILE:O	1.90	0.71
1:A:235:CYS:HB3	1:A:240:GLU:HG3	1.72	0.71
1:A:50:LYS:HE2	1:A:50:LYS:HA	1.73	0.71
1:A:69:TRP:CD2	1:B:63:PRO:HB3	2.26	0.70
1:A:236:LYS:H	1:A:240:GLU:HG2	1.55	0.70
1:A:20:CYS:CB	1:A:39:CYS:HB3	2.19	0.70
1:A:104:VAL:HG12	1:A:105:MET:N	2.09	0.67
1:A:228:VAL:HG12	1:A:229:ARG:N	2.09	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:ALA:O	1:A:185:ALA:HB2	1.94	0.66
1:B:41:LYS:HZ2	1:B:41:LYS:HB3	1.59	0.66
1:B:79:SER:HA	1:B:88:LYS:HD2	1.78	0.66
1:A:20:CYS:HB3	1:A:39:CYS:CB	2.20	0.65
1:B:77:ARG:HH21	1:B:105:MET:HE3	1.60	0.64
1:A:104:VAL:HG12	1:A:105:MET:H	1.63	0.64
1:B:77:ARG:HH21	1:B:105:MET:CE	2.11	0.64
1:B:12:LEU:HD21	1:B:30:SER:HB3	1.79	0.64
1:B:10:VAL:O	1:B:30:SER:N	2.23	0.64
1:A:251:MET:HE3	1:A:255:TRP:HE1	1.62	0.64
1:A:28:LEU:HA	1:A:34:THR:HG22	1.81	0.63
1:A:235:CYS:HB3	1:A:240:GLU:CG	2.29	0.63
1:A:13:TYR:CE1	1:A:15:ILE:HD11	2.32	0.63
1:A:204:LEU:HD22	1:A:206:PHE:CZ	2.33	0.62
1:A:234:PRO:HG3	1:A:249:GLU:OE1	1.99	0.62
1:B:16:PRO:CB	1:B:23:SER:HA	2.29	0.62
1:A:228:VAL:HG12	1:A:229:ARG:H	1.63	0.62
1:A:52:ARG:HG2	1:A:52:ARG:HH11	1.64	0.62
1:A:125:ASP:OD1	1:A:129:ASN:HB2	1.99	0.62
1:A:115:ASN:HB2	1:A:117:ASN:HD21	1.65	0.61
1:A:105:MET:HB2	1:A:126:LYS:CE	2.29	0.61
1:B:186:LYS:HB2	1:B:218:GLU:HB3	1.83	0.60
1:A:89:GLU:O	1:A:93:LYS:HG3	2.01	0.60
1:A:16:PRO:HA	1:A:23:SER:O	2.01	0.60
1:A:15:ILE:HG23	1:A:16:PRO:HD2	1.84	0.60
1:A:56:ARG:HG2	1:A:56:ARG:NH1	2.16	0.60
1:A:194:GLU:H	1:A:194:GLU:CD	2.05	0.59
1:A:41:LYS:HE3	1:A:43:THR:HG21	1.84	0.59
1:B:16:PRO:HG3	1:B:23:SER:HA	1.82	0.59
1:B:189:CYS:HB3	1:B:223:LEU:HD21	1.83	0.59
1:A:54:SER:C	1:A:56:ARG:H	2.06	0.59
1:A:13:TYR:OH	1:A:52:ARG:NH1	2.37	0.58
1:A:54:SER:O	1:A:56:ARG:N	2.35	0.58
1:A:33:HIS:HB2	1:A:44:ALA:HA	1.86	0.58
1:B:97:TRP:CZ3	1:B:99:ALA:HB2	2.39	0.58
1:A:207:ASP:OD2	1:A:237:ASP:HB3	2.04	0.58
1:A:243:LEU:O	1:A:244:ASN:ND2	2.35	0.58
1:A:142:PHE:HB3	1:A:177:SER:HB3	1.86	0.58
1:A:56:ARG:HG3	1:A:56:ARG:O	2.02	0.58
1:A:88:LYS:O	1:A:92:GLN:HG3	2.04	0.57
1:B:56:ARG:CB	1:B:56:ARG:HH11	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:TYR:O	1:A:175:VAL:HG23	2.03	0.57
1:A:209:ASP:O	1:A:213:ARG:HG2	2.04	0.57
1:A:33:HIS:CB	1:A:44:ALA:HA	2.35	0.56
1:B:194:GLU:H	1:B:194:GLU:CD	2.09	0.56
1:B:190:ALA:HB2	1:B:222:VAL:HG11	1.86	0.56
1:A:90:THR:HG22	1:A:163:LEU:HD22	1.88	0.56
1:A:60:GLY:HA3	1:B:68:VAL:HG12	1.87	0.56
1:A:78:TYR:CE2	1:A:97:TRP:HB3	2.40	0.56
1:B:154:VAL:HG21	1:B:254:VAL:HG22	1.87	0.56
1:B:201:GLN:HE22	1:B:229:ARG:NH1	2.04	0.56
1:B:15:ILE:HG13	1:B:26:ASN:HD22	1.71	0.56
1:A:41:LYS:HE3	1:A:43:THR:CG2	2.36	0.55
1:B:194:GLU:CD	1:B:194:GLU:N	2.60	0.55
1:A:74:SER:O	1:A:75:ASN:HB2	2.07	0.55
1:A:78:TYR:HE1	1:B:13:TYR:HD2	1.55	0.55
1:B:196:PHE:HB3	1:B:202:ILE:CD1	2.36	0.55
1:B:96:TYR:OH	1:B:107:GLN:NE2	2.37	0.54
1:B:21:GLY:O	1:B:22:SER:HB3	2.08	0.54
1:A:206:PHE:N	1:A:206:PHE:CD2	2.76	0.54
1:A:36:CYS:O	1:A:40:GLU:HA	2.07	0.54
1:B:152:LYS:HB3	1:B:201:GLN:HG3	1.90	0.54
1:A:153:ILE:HB	1:A:199:PHE:CD2	2.42	0.54
1:B:153:ILE:HA	1:B:174:PRO:HB2	1.89	0.54
1:A:208:MET:HE1	1:A:232:VAL:HA	1.90	0.53
1:B:20:CYS:SG	1:B:21:GLY:N	2.82	0.53
1:B:189:CYS:HB3	1:B:223:LEU:CD2	2.39	0.53
1:A:182:ALA:O	1:A:185:ALA:CB	2.57	0.53
1:A:78:TYR:CE1	1:B:13:TYR:HD2	2.26	0.53
1:B:16:PRO:CG	1:B:23:SER:HA	2.39	0.53
1:A:235:CYS:SG	1:A:241:CYS:HA	2.49	0.53
1:A:202:ILE:HD13	1:A:223:LEU:HD22	1.90	0.52
1:B:147:TRP:CD2	1:B:174:PRO:HA	2.44	0.52
1:A:39:CYS:O	1:A:40:GLU:HB2	2.09	0.52
1:A:162:MET:HG2	1:A:175:VAL:HG12	1.92	0.52
1:A:218:GLU:O	1:A:221:GLN:HB2	2.09	0.52
1:B:201:GLN:NE2	1:B:229:ARG:HD3	2.25	0.52
1:A:228:VAL:O	1:A:229:ARG:HG3	2.09	0.52
1:A:33:HIS:HB2	1:A:43:THR:O	2.10	0.52
1:A:93:LYS:HD3	1:B:52:ARG:NH2	2.23	0.52
1:B:139:ASP:HB3	1:B:191:ALA:HB1	1.92	0.52
1:B:34:THR:HG23	1:B:42:TRP:O	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:ILE:HD12	1:A:86:ILE:N	2.25	0.51
1:A:116:GLY:HA3	1:B:55:LYS:HA	1.92	0.51
1:A:223:LEU:CD2	1:A:224:PRO:HD2	2.40	0.51
1:A:204:LEU:HD21	1:A:223:LEU:HD11	1.91	0.51
1:A:170:ASP:O	1:A:172:LYS:HG3	2.11	0.51
1:A:203:ILE:N	1:A:203:ILE:HD13	2.25	0.51
1:A:193:TYR:CZ	1:A:224:PRO:HG3	2.45	0.51
1:B:96:TYR:CE2	1:B:107:GLN:HB3	2.46	0.51
1:A:78:TYR:HD1	1:B:13:TYR:HB3	1.75	0.51
1:A:87:SER:N	1:A:167:GLU:OE2	2.36	0.50
1:A:228:VAL:CG1	1:A:229:ARG:H	2.23	0.50
1:A:78:TYR:O	1:A:88:LYS:HG2	2.12	0.50
1:A:102:ASP:C	1:A:104:VAL:H	2.15	0.50
1:B:162:MET:HG2	1:B:175:VAL:HG12	1.92	0.50
1:A:127:ASP:O	1:A:128:LYS:HB2	2.12	0.50
1:A:228:VAL:CG1	1:A:229:ARG:N	2.75	0.50
1:B:26:ASN:HB2	1:B:36:CYS:SG	2.52	0.50
1:B:77:ARG:NH2	1:B:105:MET:HE3	2.27	0.50
1:B:35:PHE:HA	1:B:41:LYS:O	2.11	0.50
1:A:206:PHE:N	1:A:206:PHE:HD2	2.10	0.49
1:B:37:TYR:CD1	1:B:37:TYR:C	2.86	0.49
1:B:77:ARG:NH2	1:B:100:LYS:HE3	2.27	0.49
1:A:235:CYS:HB2	1:A:241:CYS:SG	2.52	0.49
1:A:248:ARG:O	1:A:252:GLU:HG3	2.12	0.49
1:B:121:GLN:HB2	1:B:133:THR:OG1	2.13	0.49
1:A:15:ILE:CG2	1:A:16:PRO:HD2	2.42	0.49
1:A:20:CYS:HB3	1:A:39:CYS:SG	2.53	0.49
1:A:186:LYS:O	1:A:222:VAL:HG21	2.13	0.48
1:B:41:LYS:NZ	1:B:41:LYS:CB	2.71	0.48
1:A:138:SER:OG	1:A:187:LYS:HE3	2.14	0.48
1:A:15:ILE:O	1:A:26:ASN:HB3	2.13	0.48
1:B:126:LYS:C	1:B:128:LYS:H	2.17	0.48
1:B:196:PHE:HB3	1:B:202:ILE:HD11	1.95	0.48
1:B:11:PHE:HE1	1:B:27:SER:OG	1.92	0.47
1:A:149:GLY:HA2	1:A:199:PHE:CE1	2.49	0.47
1:B:41:LYS:HZ3	1:B:41:LYS:HB3	1.74	0.47
1:A:42:TRP:O	1:A:42:TRP:HE3	1.98	0.47
1:A:154:VAL:HG21	1:A:254:VAL:HG22	1.97	0.47
1:A:101:VAL:HG21	1:B:63:PRO:HG3	1.96	0.47
1:B:104:VAL:HG23	1:B:106:TYR:CE1	2.49	0.47
1:A:207:ASP:O	1:A:213:ARG:NH2	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:CYS:O	1:A:40:GLU:N	2.48	0.47
1:B:26:ASN:HA	1:B:36:CYS:HA	1.97	0.47
1:B:39:CYS:SG	1:B:41:LYS:HB2	2.55	0.47
1:A:26:ASN:ND2	1:A:34:THR:HB	2.30	0.46
1:B:77:ARG:HH22	1:B:100:LYS:HE3	1.80	0.46
1:A:151:LYS:HD2	1:A:152:LYS:HG3	1.96	0.46
1:A:104:VAL:CG1	1:A:105:MET:N	2.78	0.46
1:B:208:MET:CE	1:B:232:VAL:HA	2.45	0.46
1:A:223:LEU:HD22	1:A:224:PRO:HD2	1.97	0.46
1:A:30:SER:O	1:A:31:ASP:C	2.54	0.46
1:B:152:LYS:CB	1:B:201:GLN:HG3	2.46	0.46
1:B:16:PRO:CA	1:B:23:SER:HA	2.46	0.46
1:B:91:CYS:HB3	1:B:96:TYR:O	2.15	0.46
1:A:201:GLN:C	1:A:202:ILE:HG13	2.36	0.46
1:A:55:LYS:HA	1:B:116:GLY:CA	2.39	0.46
1:B:16:PRO:HB3	1:B:23:SER:HA	1.98	0.46
1:A:241:CYS:O	1:A:244:ASN:N	2.40	0.45
1:A:10:VAL:HG12	1:A:11:PHE:N	2.31	0.45
1:B:251:MET:HE1	1:B:255:TRP:HE1	1.81	0.45
1:A:27:SER:O	1:A:34:THR:HA	2.17	0.45
1:B:151:LYS:O	1:B:200:GLU:HB2	2.16	0.45
1:A:30:SER:O	1:A:32:GLY:N	2.50	0.45
1:B:251:MET:CE	1:B:255:TRP:HE1	2.29	0.45
1:B:13:TYR:CE1	1:B:28:LEU:HD23	2.52	0.45
1:A:101:VAL:O	1:A:102:ASP:HB2	2.15	0.45
1:A:223:LEU:HD23	1:A:224:PRO:CD	2.47	0.45
1:A:231:ALA:HA	1:A:253:GLN:OE1	2.16	0.45
1:A:78:TYR:CD2	1:A:97:TRP:HB3	2.52	0.45
1:B:239:ASN:O	1:B:242:HIS:HB3	2.17	0.44
1:B:15:ILE:O	1:B:26:ASN:HB3	2.18	0.44
1:A:113:ASP:OD1	1:A:113:ASP:C	2.55	0.44
1:B:137:LYS:O	1:B:180:HIS:HE1	2.00	0.44
1:A:130:PHE:N	1:A:130:PHE:CD1	2.84	0.44
1:A:166:MET:HE2	1:A:171:CYS:HB3	1.98	0.44
1:A:52:ARG:HG2	1:A:52:ARG:NH1	2.31	0.44
1:A:80:ALA:HB1	1:A:87:SER:HA	1.98	0.44
1:B:10:VAL:CA	1:B:30:SER:OG	2.62	0.44
1:B:104:VAL:HG23	1:B:106:TYR:HE1	1.82	0.43
1:B:16:PRO:HA	1:B:22:SER:O	2.17	0.43
1:A:97:TRP:CZ3	1:A:99:ALA:HB2	2.53	0.43
1:B:196:PHE:HB3	1:B:202:ILE:HD13	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:TYR:CD1	1:B:13:TYR:HB3	2.53	0.43
1:B:41:LYS:HZ2	1:B:41:LYS:CB	2.26	0.43
1:B:50:LYS:HA	1:B:53:ALA:HB3	2.00	0.43
1:A:50:LYS:CA	1:A:50:LYS:HE2	2.44	0.43
1:A:54:SER:C	1:A:56:ARG:N	2.70	0.43
1:B:132:THR:HG21	1:B:136:HIS:HB2	2.00	0.43
1:B:141:LEU:HB2	1:B:144:LYS:HB2	1.99	0.43
1:B:240:GLU:O	1:B:241:CYS:C	2.56	0.43
1:A:104:VAL:CG1	1:A:105:MET:H	2.29	0.43
1:B:56:ARG:O	1:B:56:ARG:HG3	2.18	0.43
1:A:29:PHE:N	1:A:29:PHE:CD1	2.85	0.43
1:A:242:HIS:HA	1:A:247:ASP:OD2	2.19	0.43
1:A:141:LEU:HG	1:A:192:ASN:ND2	2.33	0.43
1:A:130:PHE:N	1:A:130:PHE:HD1	2.16	0.42
1:B:78:TYR:O	1:B:88:LYS:HD2	2.18	0.42
1:B:16:PRO:HA	1:B:23:SER:HA	1.99	0.42
1:A:75:ASN:ND2	1:A:100:LYS:O	2.44	0.42
1:A:201:GLN:NE2	1:A:229:ARG:HD3	2.34	0.42
1:A:63:PRO:HB3	1:B:69:TRP:CD2	2.54	0.42
1:A:111:TYR:CZ	1:A:142:PHE:HB2	2.55	0.42
1:A:147:TRP:CD2	1:A:174:PRO:HA	2.54	0.42
1:A:204:LEU:HD21	1:A:223:LEU:CD1	2.50	0.42
1:A:246:HIS:HB3	1:A:249:GLU:OE1	2.19	0.42
1:A:12:LEU:HA	1:A:56:ARG:HH12	1.84	0.42
1:A:69:TRP:CE3	1:B:63:PRO:HB3	2.53	0.42
1:A:102:ASP:O	1:A:104:VAL:HG23	2.20	0.42
1:A:141:LEU:HB3	1:A:144:LYS:HB2	2.02	0.42
1:B:212:GLY:O	1:B:216:VAL:HG23	2.19	0.42
1:B:169:GLN:NE2	1:B:254:VAL:HG11	2.35	0.42
1:A:66:TYR:C	1:B:67:ASN:HB2	2.40	0.42
1:B:100:LYS:HB3	1:B:105:MET:CE	2.50	0.42
1:B:112:ARG:HA	1:B:117:ASN:O	2.20	0.42
1:A:100:LYS:HA	1:A:104:VAL:O	2.20	0.42
1:A:254:VAL:HG13	1:A:254:VAL:O	2.20	0.42
1:A:12:LEU:HD22	1:A:56:ARG:HG2	2.02	0.42
1:B:10:VAL:HG12	1:B:31:ASP:OD1	2.20	0.41
1:A:22:SER:HB3	1:A:38:VAL:HG21	2.02	0.41
1:A:127:ASP:HB2	1:A:129:ASN:ND2	2.35	0.41
1:A:22:SER:CB	1:A:38:VAL:HG21	2.50	0.41
1:A:36:CYS:O	1:A:40:GLU:CA	2.67	0.41
1:A:204:LEU:HD22	1:A:206:PHE:CE2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:ARG:NH2	1:B:159:GLU:HB2	2.36	0.41
1:B:221:GLN:HB2	1:B:221:GLN:HE21	1.67	0.41
1:A:56:ARG:CG	1:A:56:ARG:NH1	2.81	0.41
1:B:190:ALA:CA	1:B:222:VAL:HG11	2.51	0.41
1:A:105:MET:HB2	1:A:126:LYS:NZ	2.35	0.41
1:A:80:ALA:HB2	1:A:88:LYS:N	2.36	0.41
1:A:254:VAL:CG1	1:A:254:VAL:O	2.69	0.41
1:A:10:VAL:O	1:A:29:PHE:HA	2.20	0.41
1:A:168:LEU:HD21	1:A:242:HIS:CE1	2.56	0.40
1:A:241:CYS:HB3	1:A:246:HIS:HB2	2.03	0.40
1:A:63:PRO:HD3	1:B:74:SER:OG	2.22	0.40
1:A:124:ARG:HH22	1:A:159:GLU:HB2	1.86	0.40
1:A:195:TYR:O	1:A:197:ASP:N	2.54	0.40
1:B:101:VAL:O	1:B:104:VAL:HG22	2.21	0.40
1:A:89:GLU:HG2	1:A:93:LYS:HE3	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:SER:CB	1:B:22:SER:CB[4_556]	2.11	0.09

## 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	238/255 (93%)	191 (80%)	41 (17%)	6 (2%)	6	25
1	B	235/255 (92%)	205 (87%)	25 (11%)	5 (2%)	8	30
All	All	473/510 (93%)	396 (84%)	66 (14%)	11 (2%)	7	27

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	55	LYS
1	B	22	SER
1	A	31	ASP
1	A	196	PHE
1	A	204	LEU
1	A	38	VAL
1	A	245	GLY
1	B	25	GLY
1	B	104	VAL
1	B	127	ASP
1	B	38	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	200/212 (94%)	192 (96%)	8 (4%)	36	71
1	B	198/212 (93%)	190 (96%)	8 (4%)	36	71
All	All	398/424 (94%)	382 (96%)	16 (4%)	36	71

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

Mol	Chain	Res	Type
1	A	24	ASP
1	A	30	SER
1	A	50	LYS
1	A	62	LYS
1	A	77	ARG
1	A	102	ASP
1	A	194	GLU
1	A	254	VAL
1	B	10	VAL
1	B	20	CYS
1	B	56	ARG
1	B	57	LYS
1	B	62	LYS

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Mol	Chain	Res	Type
1	B	170	ASP
1	B	194	GLU
1	B	240	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	ASN
1	A	117	ASN
1	A	129	ASN
1	A	198	GLN
1	A	201	GLN
1	A	221	GLN
1	B	107	GLN
1	B	145	HIS
1	B	169	GLN
1	B	180	HIS
1	B	201	GLN
1	B	221	GLN
1	B	244	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.



There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	242/255 (94%)	0.08	3 (1%) 79 77	4, 19, 29, 41	0
1	B	239/255 (93%)	0.00	5 (2%) 64 60	4, 18, 32, 45	0
All	All	481/510 (94%)	0.04	8 (1%) 70 68	4, 19, 31, 45	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	21	GLY	5.7
1	B	18	ASP	2.9
1	B	114	GLN	2.8
1	B	25	GLY	2.7
1	B	20	CYS	2.5
1	A	210	GLU	2.5
1	A	23	SER	2.4
1	A	235	CYS	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	MG	B	603	1/1	0.96	0.18	0.95	38,38,38,38	0
2	ZN	B	502	1/1	0.95	0.08	-1.46	83,83,83,83	0
2	ZN	A	501	1/1	0.96	0.06	-2.15	84,84,84,84	0
3	MG	B	604	1/1	0.90	0.31	-	45,45,45,45	0
3	MG	A	602	1/1	0.52	0.35	-	88,88,88,88	0
3	MG	A	601	1/1	0.47	0.32	-	87,87,87,87	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.