



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:06 AM GMT

PDB ID : 1ZQY  
Title : DNA POLYMERASE BETA (POL B) (E.C.2.7.7.7), 31-KD DOMAIN;  
SOAKED IN THE PRESENCE OF MGCL2 (50 MILLIMOLAR)  
Authors : Pelletier, H.; Sawaya, M.R.  
Deposited on : 1996-04-19  
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

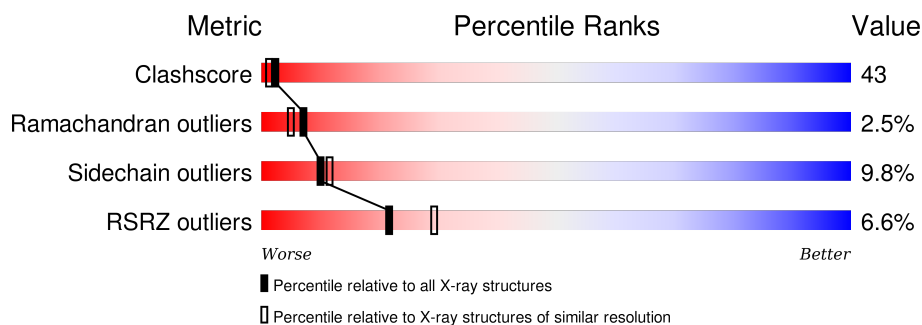
# 1 Overall quality at a glance

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))
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	248	<div> <div>6%</div> <div>33%</div> <div>46%</div> <div>16%</div> <div>••</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2037 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 POLYMERASE BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	1	0
			1935	1218	342	367	8			

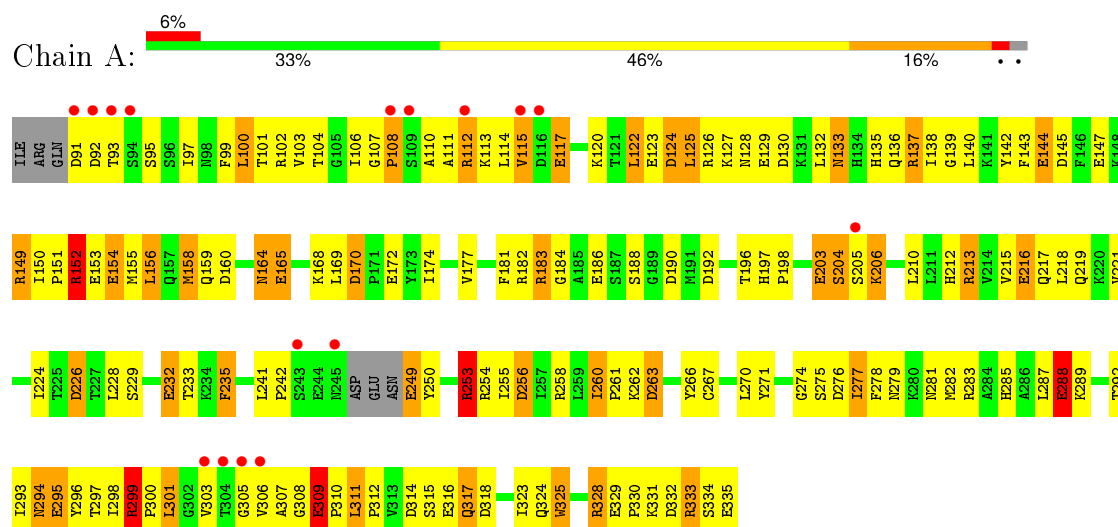
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	102	Total	O	0	0
			102	102		

### 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 ( $\text{RSRZ} > 2$ ). 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 POLYMERASE BETA



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	120.56Å 63.24Å 38.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 23.75 – 2.11	Depositor EDS
% Data completeness (in resolution range)	92.0 (20.00-2.30) 86.2 (23.75-2.11)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.37 (at 2.11Å)	Xtriage
Refinement program	TNT V. 5-D	Depositor
R, $R_{free}$	(Not available) , (Not available) 0.192 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	19.8	Xtriage
Anisotropy	0.786	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 106.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 15197 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	2037	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.92% 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 [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	1.47	17/1973 (0.9%)	1.91	54/2662 (2.0%)

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	1	0

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	295	GLU	CD-OE1	8.35	1.34	1.25
1	A	288	GLU	CD-OE1	8.21	1.34	1.25
1	A	216	GLU	CD-OE1	8.11	1.34	1.25
1	A	147	GLU	CD-OE1	8.03	1.34	1.25
1	A	309	GLU	CD-OE1	7.93	1.34	1.25
1	A	165	GLU	CD-OE1	7.85	1.34	1.25
1	A	172	GLU	CD-OE1	7.79	1.34	1.25
1	A	117	GLU	CD-OE1	7.62	1.34	1.25
1	A	154	GLU	CD-OE1	7.51	1.33	1.25
1	A	203	GLU	CD-OE1	7.43	1.33	1.25
1	A	335	GLU	CD-OE1	7.17	1.33	1.25
1	A	232	GLU	CD-OE1	7.16	1.33	1.25
1	A	153	GLU	CD-OE2	7.15	1.33	1.25
1	A	144	GLU	CD-OE1	6.62	1.32	1.25
1	A	129	GLU	CD-OE2	-6.50	1.18	1.25
1	A	249	GLU	CD-OE1	6.04	1.32	1.25
1	A	186	GLU	CD-OE1	5.79	1.32	1.25

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	182	ARG	NE-CZ-NH1	16.23	128.41	120.30
1	A	299[A]	ARG	NE-CZ-NH1	11.67	126.13	120.30
1	A	299[B]	ARG	NE-CZ-NH1	11.67	126.13	120.30
1	A	182	ARG	NE-CZ-NH2	-10.41	115.09	120.30
1	A	333	ARG	NE-CZ-NH1	9.98	125.29	120.30
1	A	314	ASP	CB-CG-OD1	-9.94	109.36	118.30
1	A	258	ARG	NE-CZ-NH2	-9.66	115.47	120.30
1	A	190	ASP	CB-CG-OD2	-9.13	110.08	118.30
1	A	258	ARG	NE-CZ-NH1	8.40	124.50	120.30
1	A	299[A]	ARG	NE-CZ-NH2	-8.02	116.29	120.30
1	A	299[B]	ARG	NE-CZ-NH2	-8.02	116.29	120.30
1	A	253	ARG	NE-CZ-NH1	7.90	124.25	120.30
1	A	145	ASP	CB-CG-OD2	7.84	125.36	118.30
1	A	333	ARG	NE-CZ-NH2	-7.74	116.43	120.30
1	A	213	ARG	NE-CZ-NH2	-7.57	116.52	120.30
1	A	170	ASP	CB-CG-OD2	7.54	125.09	118.30
1	A	170	ASP	CB-CG-OD1	-7.49	111.56	118.30
1	A	263	ASP	CB-CG-OD2	7.29	124.86	118.30
1	A	253	ARG	NE-CZ-NH2	-7.28	116.66	120.30
1	A	263	ASP	CB-CG-OD1	-7.12	111.89	118.30
1	A	192	ASP	CB-CG-OD1	-6.99	112.01	118.30
1	A	235	PHE	CB-CA-C	-6.98	96.44	110.40
1	A	160	ASP	CB-CG-OD1	-6.89	112.10	118.30
1	A	142	TYR	CB-CG-CD1	-6.81	116.91	121.00
1	A	183	ARG	NE-CZ-NH2	-6.80	116.90	120.30
1	A	190	ASP	CB-CG-OD1	6.78	124.40	118.30
1	A	254	ARG	NE-CZ-NH2	-6.77	116.91	120.30
1	A	152	ARG	N-CA-CB	6.76	122.78	110.60
1	A	124	ASP	CB-CG-OD1	-6.62	112.34	118.30
1	A	213	ARG	NE-CZ-NH1	6.61	123.60	120.30
1	A	328	ARG	NE-CZ-NH2	-6.48	117.06	120.30
1	A	256	ASP	CB-CG-OD1	-6.46	112.48	118.30
1	A	314	ASP	CB-CG-OD2	6.45	124.11	118.30
1	A	260	ILE	CB-CA-C	-6.35	98.89	111.60
1	A	226	ASP	CB-CG-OD1	6.33	124.00	118.30
1	A	276	ASP	CB-CG-OD1	-6.24	112.69	118.30
1	A	137	ARG	NE-CZ-NH1	6.19	123.39	120.30
1	A	102	ARG	CD-NE-CZ	-6.18	114.94	123.60
1	A	112	ARG	NE-CZ-NH1	6.18	123.39	120.30
1	A	318	ASP	CB-CG-OD2	6.04	123.73	118.30
1	A	132	LEU	CB-CA-C	-5.99	98.81	110.20
1	A	137	ARG	NE-CZ-NH2	-5.95	117.33	120.30
1	A	318	ASP	CB-CG-OD1	-5.94	112.95	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	183	ARG	NE-CZ-NH1	5.93	123.26	120.30
1	A	149	ARG	NE-CZ-NH1	5.87	123.24	120.30
1	A	229	SER	N-CA-CB	5.81	119.22	110.50
1	A	147	GLU	N-CA-CB	5.80	121.04	110.60
1	A	226	ASP	CB-CG-OD2	-5.57	113.29	118.30
1	A	256	ASP	CB-CG-OD2	5.47	123.22	118.30
1	A	332	ASP	CB-CG-OD2	5.31	123.08	118.30
1	A	130	ASP	C-N-CA	-5.28	108.51	121.70
1	A	125	LEU	CB-CG-CD1	-5.23	102.11	111.00
1	A	266	TYR	CB-CG-CD1	-5.14	117.92	121.00
1	A	267	CYS	N-CA-CB	-5.11	101.41	110.60

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	152	ARG	CA

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	1935	0	1868	164	0
2	A	102	0	0	12	2
All	All	2037	0	1868	164	2

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 43.

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)
1:A:317:GLN:NE2	1:A:317:GLN:H	1.43	1.16
1:A:277:ILE:H	1:A:277:ILE:HD13	1.26	0.97
1:A:317:GLN:HE21	1:A:317:GLN:N	1.63	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:GLN:HE21	1:A:317:GLN:H	0.97	0.91
1:A:294:ASN:ND2	1:A:296:TYR:H	1.75	0.84
1:A:294:ASN:HD22	1:A:296:TYR:H	1.25	0.84
1:A:270:LEU:HD21	1:A:282:MET:HE3	1.60	0.82
1:A:279:ASN:O	1:A:283:ARG:HG2	1.79	0.80
1:A:103:VAL:HB	1:A:106:ILE:HD13	1.61	0.80
1:A:133:ASN:ND2	1:A:136:GLN:H	1.79	0.80
1:A:270:LEU:HD21	1:A:282:MET:CE	2.13	0.78
1:A:154:GLU:O	1:A:158:MET:HG2	1.84	0.78
1:A:97:ILE:HD11	1:A:112:ARG:HG2	1.70	0.72
1:A:110:ALA:O	1:A:114:LEU:HD13	1.90	0.71
1:A:103:VAL:HB	1:A:106:ILE:CD1	2.21	0.71
1:A:277:ILE:N	1:A:277:ILE:HD13	2.05	0.70
1:A:100:LEU:O	1:A:103:VAL:HG23	1.91	0.70
1:A:92:ASP:O	1:A:95:SER:HB3	1.93	0.68
1:A:144:GLU:OE1	1:A:144:GLU:N	2.25	0.68
1:A:330:PRO:HD2	2:A:429:HOH:O	1.95	0.67
1:A:97:ILE:CD1	1:A:112:ARG:HG2	2.25	0.66
1:A:270:LEU:HD12	1:A:333:ARG:NH1	2.11	0.65
1:A:278:PHE:CE2	1:A:282:MET:HE2	2.31	0.65
1:A:107:GLY:O	1:A:111:ALA:N	2.29	0.65
1:A:317:GLN:NE2	1:A:317:GLN:N	2.26	0.65
1:A:120:LYS:N	1:A:124:ASP:OD2	2.27	0.64
1:A:263:ASP:N	1:A:263:ASP:OD1	2.26	0.64
1:A:122:LEU:HD12	1:A:123:GLU:N	2.12	0.63
1:A:123:GLU:O	1:A:127:LYS:HG2	1.99	0.63
1:A:295:GLU:H	1:A:295:GLU:CD	2.02	0.62
1:A:133:ASN:HD22	1:A:133:ASN:C	2.03	0.62
1:A:183:ARG:NE	2:A:499:HOH:O	2.31	0.62
1:A:315:SER:HB2	1:A:317:GLN:HE22	1.65	0.62
1:A:133:ASN:HD21	1:A:136:GLN:H	1.45	0.62
1:A:305:GLY:O	1:A:307:ALA:N	2.34	0.61
1:A:232:GLU:HG3	1:A:233:THR:HG23	1.82	0.61
1:A:133:ASN:H	1:A:136:GLN:NE2	2.00	0.59
1:A:103:VAL:CB	1:A:106:ILE:HD13	2.32	0.59
1:A:294:ASN:HD22	1:A:295:GLU:N	2.01	0.58
1:A:150:ILE:HG22	1:A:151:PRO:HD2	1.84	0.58
1:A:164:ASN:OD1	1:A:165:GLU:HG2	2.02	0.58
1:A:301:LEU:HD12	1:A:301:LEU:N	2.19	0.58
1:A:277:ILE:HG12	1:A:278:PHE:H	1.69	0.57
1:A:204:SER:CA	1:A:206:LYS:HE2	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:ILE:HB	1:A:228:LEU:CD2	2.34	0.57
1:A:274:GLY:HA2	1:A:279:ASN:OD1	2.05	0.57
1:A:150:ILE:CG2	1:A:151:PRO:HD2	2.35	0.57
1:A:152:ARG:NH2	1:A:181:PHE:O	2.39	0.56
1:A:204:SER:C	1:A:206:LYS:HE2	2.26	0.56
1:A:299[A]:ARG:HG3	1:A:299[A]:ARG:HH11	1.70	0.55
1:A:323:ILE:C	1:A:324:GLN:HG2	2.27	0.55
1:A:285:HIS:HD2	1:A:323:ILE:HD12	1.71	0.55
1:A:294:ASN:C	1:A:294:ASN:HD22	2.10	0.55
1:A:156:LEU:HD13	1:A:181:PHE:HE1	1.71	0.54
1:A:249:GLU:HG3	2:A:455:HOH:O	2.07	0.54
1:A:143:PHE:HB3	1:A:144:GLU:OE1	2.07	0.54
1:A:262:LYS:HG3	1:A:262:LYS:O	2.07	0.54
1:A:277:ILE:H	1:A:277:ILE:CD1	2.07	0.53
1:A:287:LEU:HD12	1:A:287:LEU:O	2.08	0.53
1:A:114:LEU:HD12	1:A:114:LEU:N	2.21	0.53
1:A:288:GLU:N	1:A:288:GLU:OE1	2.41	0.53
1:A:137:ARG:NH1	2:A:468:HOH:O	2.41	0.53
1:A:315:SER:HB2	1:A:317:GLN:NE2	2.24	0.52
1:A:285:HIS:O	1:A:288:GLU:N	2.42	0.52
1:A:150:ILE:O	1:A:188:SER:N	2.33	0.52
1:A:205:SER:O	1:A:206:LYS:HD3	2.10	0.52
1:A:299[B]:ARG:HD2	1:A:308:GLY:HA3	1.92	0.51
1:A:289:LYS:NZ	1:A:324:GLN:HG3	2.26	0.51
1:A:150:ILE:HD13	1:A:253:ARG:HG2	1.92	0.51
1:A:215:VAL:O	1:A:219:GLN:HG3	2.10	0.51
1:A:271:TYR:CG	1:A:295:GLU:HB3	2.46	0.51
1:A:329:GLU:HB3	1:A:330:PRO:HD2	1.92	0.51
1:A:106:ILE:CG2	1:A:111:ALA:HB2	2.41	0.51
1:A:197:HIS:CG	1:A:198:PRO:HD2	2.46	0.51
1:A:285:HIS:NE2	1:A:325:TRP:CD1	2.79	0.51
1:A:169:LEU:O	1:A:170:ASP:HB2	2.10	0.51
1:A:334:SER:HA	2:A:499:HOH:O	2.10	0.50
1:A:152:ARG:NH1	2:A:463:HOH:O	2.28	0.50
1:A:108:PRO:O	1:A:111:ALA:N	2.44	0.50
1:A:241:LEU:CD2	1:A:242:PRO:HD2	2.42	0.49
1:A:297:THR:O	1:A:299[A]:ARG:NH1	2.45	0.49
1:A:112:ARG:O	1:A:115:VAL:HG23	2.13	0.49
1:A:122:LEU:HD12	1:A:122:LEU:C	2.32	0.49
1:A:156:LEU:CD1	1:A:181:PHE:CE1	2.96	0.49
1:A:224:ILE:HD13	1:A:235:PHE:CE2	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:275:SER:OG	1:A:277:ILE:HG12	2.12	0.48
1:A:323:ILE:O	1:A:324:GLN:HG2	2.12	0.48
1:A:138:ILE:HB	1:A:228:LEU:HD23	1.95	0.48
1:A:150:ILE:HA	1:A:151:PRO:HD3	1.66	0.48
1:A:255:ILE:HG12	1:A:256:ASP:N	2.29	0.48
1:A:270:LEU:HD12	1:A:333:ARG:HH11	1.78	0.48
1:A:299[B]:ARG:HD2	1:A:308:GLY:CA	2.44	0.48
1:A:110:ALA:HA	1:A:113:LYS:HE3	1.96	0.48
1:A:204:SER:HB3	1:A:206:LYS:HE2	1.95	0.47
1:A:241:LEU:HD23	1:A:241:LEU:HA	1.73	0.47
1:A:212:HIS:O	1:A:216:GLU:HG3	2.13	0.47
1:A:156:LEU:CD1	1:A:181:PHE:HE1	2.26	0.47
1:A:217:GLN:O	1:A:221:VAL:HG22	2.14	0.47
1:A:299[B]:ARG:CD	1:A:309:GLU:H	2.28	0.47
1:A:260:ILE:HG22	1:A:261:PRO:N	2.28	0.47
1:A:126:ARG:H	1:A:126:ARG:HG3	1.47	0.47
1:A:155:MET:HA	1:A:158:MET:HG3	1.96	0.47
1:A:114:LEU:CD1	1:A:114:LEU:N	2.78	0.47
1:A:122:LEU:O	1:A:125:LEU:HB2	2.14	0.47
1:A:126:ARG:HG2	1:A:140:LEU:HD21	1.97	0.47
1:A:104:THR:HG23	1:A:139:GLY:HA3	1.96	0.46
1:A:270:LEU:CD1	1:A:333:ARG:NH1	2.77	0.46
1:A:104:THR:HG22	2:A:435:HOH:O	2.15	0.46
1:A:150:ILE:N	1:A:188:SER:O	2.30	0.46
1:A:103:VAL:HA	2:A:435:HOH:O	2.14	0.46
1:A:174:ILE:HB	1:A:196:THR:HG22	1.98	0.46
1:A:308:GLY:O	1:A:309:GLU:HB2	2.16	0.46
1:A:156:LEU:HA	1:A:156:LEU:HD12	1.82	0.46
1:A:315:SER:CB	1:A:317:GLN:NE2	2.79	0.46
1:A:328:ARG:NH2	2:A:470:HOH:O	2.49	0.46
1:A:278:PHE:CE2	1:A:282:MET:CE	2.98	0.45
1:A:113:LYS:HG3	1:A:114:LEU:HD12	1.98	0.45
1:A:113:LYS:O	1:A:117:GLU:HG2	2.17	0.45
1:A:277:ILE:HG12	1:A:278:PHE:N	2.30	0.45
1:A:283:ARG:NH2	1:A:294:ASN:HA	2.31	0.45
1:A:255:ILE:CG1	1:A:256:ASP:N	2.80	0.45
1:A:250:TYR:HB3	2:A:404:HOH:O	2.16	0.45
1:A:316:GLU:N	1:A:317:GLN:NE2	2.65	0.45
1:A:213:ARG:HD3	2:A:444:HOH:O	2.16	0.45
1:A:301:LEU:N	1:A:301:LEU:CD1	2.80	0.45
1:A:92:ASP:C	1:A:95:SER:HB3	2.37	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:MET:HA	1:A:158:MET:CG	2.47	0.44
1:A:159:GLN:HB2	1:A:177:VAL:HG21	2.00	0.44
1:A:317:GLN:CD	1:A:317:GLN:H	2.09	0.44
1:A:281:ASN:HA	1:A:281:ASN:HD22	1.65	0.44
1:A:299[B]:ARG:HG2	1:A:310:PRO:N	2.33	0.44
1:A:210:LEU:HA	1:A:210:LEU:HD23	1.83	0.43
1:A:152:ARG:NH2	1:A:184:GLY:CA	2.82	0.43
1:A:311:LEU:HA	1:A:311:LEU:HD12	1.70	0.43
1:A:149:ARG:HG3	2:A:447:HOH:O	2.18	0.43
1:A:205:SER:O	1:A:206:LYS:HB3	2.19	0.43
1:A:168:LYS:HB2	1:A:168:LYS:HZ2	1.82	0.43
1:A:197:HIS:ND1	1:A:198:PRO:HD2	2.34	0.43
1:A:158:MET:H	1:A:158:MET:HG2	1.71	0.43
1:A:138:ILE:HG23	1:A:138:ILE:HD12	1.61	0.43
1:A:285:HIS:CD2	1:A:323:ILE:HD12	2.52	0.43
1:A:285:HIS:CD2	1:A:325:TRP:CE2	3.07	0.43
1:A:293:ILE:HG21	1:A:293:ILE:HD13	1.62	0.43
1:A:299[B]:ARG:HB3	1:A:300:PRO:HD2	2.01	0.43
1:A:168:LYS:NZ	1:A:168:LYS:CB	2.81	0.43
1:A:133:ASN:HD21	1:A:135:HIS:HB3	1.83	0.42
1:A:127:LYS:HA	1:A:127:LYS:HD3	1.84	0.42
1:A:126:ARG:C	1:A:128:ASN:H	2.23	0.42
1:A:174:ILE:HD12	1:A:196:THR:CG2	2.50	0.42
1:A:303:VAL:C	1:A:305:GLY:H	2.23	0.42
1:A:150:ILE:CD1	1:A:253:ARG:HG2	2.50	0.42
1:A:152:ARG:NH2	1:A:184:GLY:HA2	2.35	0.42
1:A:218:LEU:HD23	1:A:218:LEU:HA	1.88	0.41
1:A:114:LEU:H	1:A:114:LEU:CD1	2.33	0.41
1:A:241:LEU:HD23	1:A:242:PRO:HD2	2.02	0.41
1:A:294:ASN:ND2	1:A:294:ASN:C	2.74	0.41
1:A:91:ASP:C	1:A:93:THR:H	2.24	0.41
1:A:168:LYS:HB2	1:A:168:LYS:NZ	2.35	0.41
1:A:289:LYS:HZ3	1:A:324:GLN:HG3	1.86	0.41
1:A:99:PHE:CD1	1:A:99:PHE:C	2.95	0.41
1:A:311:LEU:HA	1:A:312:PRO:HD3	1.77	0.41
1:A:298:ILE:C	1:A:299[A]:ARG:HG3	2.41	0.40
1:A:292:THR:N	1:A:301:LEU:HD11	2.37	0.40
1:A:299[B]:ARG:HD2	1:A:309:GLU:H	1.86	0.40

All (2) 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)
2:A:477:HOH:O	2:A:477:HOH:O[2_555]	0.72	1.48
2:A:502:HOH:O	2:A:502:HOH:O[2_555]	1.85	0.35

## 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	239/248 (96%)	213 (89%)	20 (8%)	6 (2%)	<b>7</b> <b>5</b>

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	206	LYS
1	A	306	VAL
1	A	309	GLU
1	A	204	SER
1	A	301	LEU
1	A	108	PRO

### 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	206/226 (91%)	185 (90%)	21 (10%)	<b>9</b> <b>10</b>

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	LEU
1	A	101	THR
1	A	115	VAL
1	A	122	LEU
1	A	133	ASN
1	A	152	ARG
1	A	156	LEU
1	A	158	MET
1	A	164	ASN
1	A	203	GLU
1	A	226	ASP
1	A	253	ARG
1	A	277	ILE
1	A	288	GLU
1	A	294	ASN
1	A	299[A]	ARG
1	A	299[B]	ARG
1	A	311	LEU
1	A	317	GLN
1	A	325	TRP
1	A	331	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	ASN
1	A	133	ASN
1	A	136	GLN
1	A	157	GLN
1	A	217	GLN
1	A	252	HIS
1	A	281	ASN
1	A	285	HIS
1	A	294	ASN
1	A	317	GLN

### 5.3.3 RNA ⓘ

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 [i](#)

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	242/248 (97%)	-0.09	16 (6%)	22 29	16, 36, 89, 100	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	205	SER	6.8
1	A	304	THR	6.8
1	A	303	VAL	5.1
1	A	93	THR	4.7
1	A	306	VAL	4.5
1	A	94	SER	4.1
1	A	91	ASP	4.0
1	A	115	VAL	3.9
1	A	112	ARG	3.8
1	A	116	ASP	3.4
1	A	108	PRO	3.2
1	A	109	SER	2.5
1	A	245	ASN	2.4
1	A	92	ASP	2.3
1	A	243	SER	2.2
1	A	305	GLY	2.2

### 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.