



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:14 AM UTC

PDB ID : 2O0K / pdb_00002o0k
Title : T4 gp17 ATPase domain mutant
Authors : Sun, S.; Rossmann, M.G.
Deposited on : 2006-11-27
Resolution : 2.50 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
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

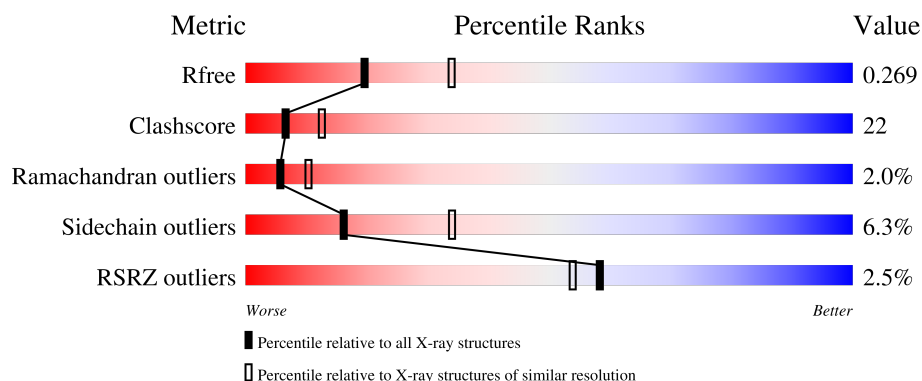
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	385	2% 56% 30% 6% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3019 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 packaging protein Gp17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	357	Total	C	N	O	S	0	0	0
			2904	1855	501	535	13			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	cloning artifact	UNP P17312
A	-23	GLY	-	cloning artifact	UNP P17312
A	-22	SER	-	cloning artifact	UNP P17312
A	-21	SER	-	cloning artifact	UNP P17312
A	-20	HIS	-	cloning artifact	UNP P17312
A	-19	HIS	-	cloning artifact	UNP P17312
A	-18	HIS	-	cloning artifact	UNP P17312
A	-17	HIS	-	cloning artifact	UNP P17312
A	-16	HIS	-	cloning artifact	UNP P17312
A	-15	HIS	-	cloning artifact	UNP P17312
A	-14	SER	-	cloning artifact	UNP P17312
A	-13	SER	-	cloning artifact	UNP P17312
A	-12	GLY	-	cloning artifact	UNP P17312
A	-11	LEU	-	cloning artifact	UNP P17312
A	-10	VAL	-	cloning artifact	UNP P17312
A	-9	PRO	-	cloning artifact	UNP P17312
A	-8	ARG	-	cloning artifact	UNP P17312
A	-7	GLY	-	cloning artifact	UNP P17312
A	-6	SER	-	cloning artifact	UNP P17312
A	-5	HIS	-	cloning artifact	UNP P17312
A	-4	MET	-	cloning artifact	UNP P17312
A	-3	LEU	-	cloning artifact	UNP P17312
A	-2	GLU	-	cloning artifact	UNP P17312
A	-1	ASP	-	cloning artifact	UNP P17312
A	0	PRO	-	cloning artifact	UNP P17312
A	255	GLU	ASP	engineered mutation	UNP P17312
A	256	ASP	GLU	engineered mutation	UNP P17312

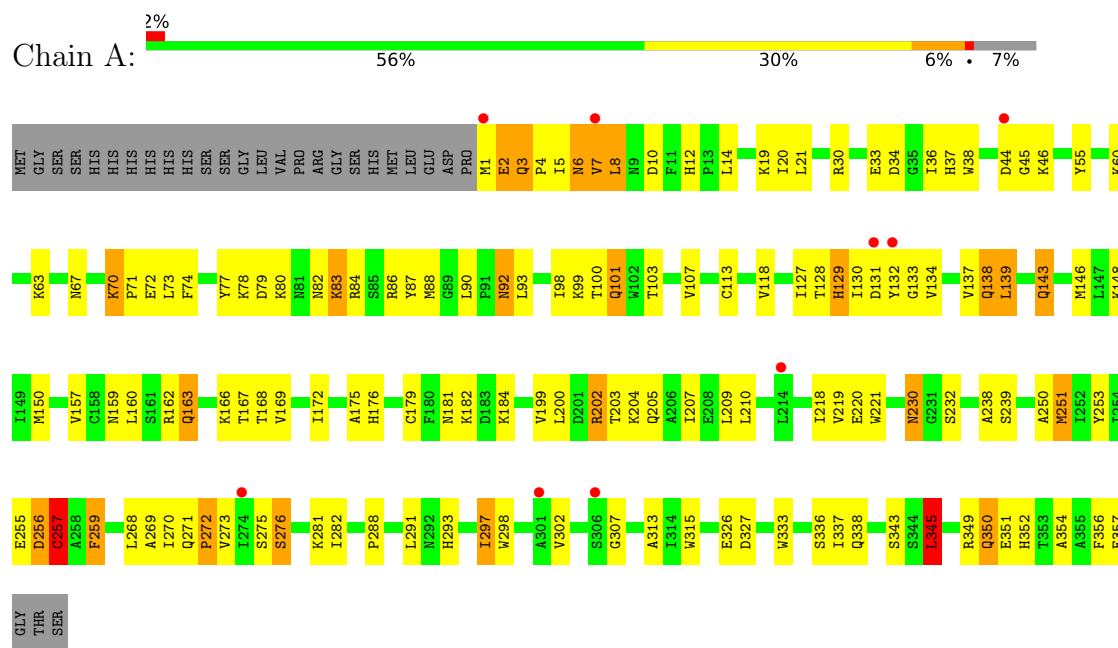
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	115	Total 115	O 115	0	0

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: DNA packaging protein Gp17



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	38.70Å 118.50Å 47.50Å 90.00° 92.50° 90.00°	Depositor
Resolution (Å)	47.40 – 2.50 47.40 – 2.50	Depositor EDS
% Data completeness (in resolution range)	95.9 (47.40-2.50) 95.9 (47.40-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.29Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.246 , 0.274 0.221 , 0.269	Depositor DCC
R_{free} test set	872 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	56.4	Xtriage
Anisotropy	0.306	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.062 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3019	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.76% 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 ⓘ

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	0.52	0/2978	1.15	27/4030 (0.7%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	138	GLN	OE1-CD-NE2	-10.91	111.69	122.60
1	A	163	GLN	OE1-CD-NE2	-10.09	112.51	122.60
1	A	3	GLN	OE1-CD-NE2	-9.82	112.78	122.60
1	A	205	GLN	OE1-CD-NE2	-9.78	112.82	122.60
1	A	45	GLY	N-CA-C	-9.54	102.51	114.92
1	A	350	GLN	OE1-CD-NE2	-9.43	113.17	122.60
1	A	143	GLN	OE1-CD-NE2	-9.34	113.26	122.60
1	A	271	GLN	OE1-CD-NE2	-8.81	113.79	122.60
1	A	338	GLN	OE1-CD-NE2	-8.75	113.85	122.60
1	A	250	ALA	N-CA-C	-8.41	103.15	113.50
1	A	55	TYR	N-CA-C	-7.14	103.58	111.36
1	A	205	GLN	CG-CD-NE2	6.87	126.71	116.40
1	A	138	GLN	CG-CD-NE2	6.82	126.63	116.40
1	A	163	GLN	CG-CD-NE2	6.68	126.43	116.40
1	A	3	GLN	CG-CD-NE2	6.67	126.40	116.40
1	A	350	GLN	CG-CD-NE2	6.44	126.06	116.40
1	A	143	GLN	CG-CD-NE2	6.34	125.91	116.40
1	A	271	GLN	CG-CD-NE2	6.31	125.87	116.40
1	A	92	ASN	N-CA-C	6.18	120.68	113.20
1	A	338	GLN	CG-CD-NE2	6.01	125.41	116.40
1	A	202	ARG	CD-NE-CZ	5.71	132.40	124.40
1	A	307	GLY	N-CA-C	-5.71	107.58	114.66
1	A	6	ASN	N-CA-C	5.52	116.33	108.54
1	A	297	ILE	N-CA-C	-5.43	106.40	111.45
1	A	345	LEU	N-CA-C	-5.22	105.67	111.36
1	A	70	LYS	CA-C-N	5.07	125.10	119.32
1	A	70	LYS	C-N-CA	5.07	125.10	119.32

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	2904	0	2855	125	0
2	A	115	0	0	1	0
All	All	3019	0	2855	125	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 (125) 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:7:VAL:HG12	1:A:8:LEU:H	1.17	1.09
1:A:128:THR:HG22	1:A:129:HIS:H	1.08	1.06
1:A:127:ILE:HG13	1:A:137:VAL:HG22	1.47	0.96
1:A:131:ASP:O	1:A:133:GLY:N	1.99	0.95
1:A:82:ASN:ND2	1:A:84:ARG:HG2	1.81	0.95
1:A:128:THR:HG22	1:A:129:HIS:N	1.83	0.93
1:A:83:LYS:HD3	1:A:83:LYS:H	1.34	0.90
1:A:128:THR:CG2	1:A:129:HIS:H	1.83	0.88
1:A:82:ASN:HD21	1:A:84:ARG:HG2	1.36	0.83
1:A:256:ASP:HB3	1:A:259:PHE:CD2	2.13	0.83
1:A:7:VAL:HG12	1:A:8:LEU:N	1.98	0.79
1:A:315:TRP:HB3	1:A:352:HIS:HB3	1.65	0.78
1:A:256:ASP:HB3	1:A:259:PHE:HD2	1.50	0.77
1:A:345:LEU:HD22	1:A:349:ARG:HH11	1.51	0.75
1:A:253:TYR:CE2	1:A:255:GLU:HB2	2.23	0.73
1:A:204:LYS:HB3	1:A:221:TRP:CZ2	2.23	0.73
1:A:230:ASN:HD22	1:A:230:ASN:C	1.99	0.71
1:A:230:ASN:ND2	1:A:232:SER:H	1.88	0.71
1:A:63:LYS:NZ	1:A:133:GLY:HA3	2.05	0.71
1:A:127:ILE:CG1	1:A:137:VAL:HG22	2.19	0.70
1:A:345:LEU:HD22	1:A:349:ARG:NH1	2.07	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:VAL:O	1:A:8:LEU:HD12	1.93	0.68
1:A:90:LEU:HD12	1:A:93:LEU:HD12	1.77	0.66
1:A:7:VAL:O	1:A:8:LEU:CD1	2.44	0.66
1:A:270:ILE:O	1:A:273:VAL:HG22	1.96	0.65
1:A:80:LYS:NZ	1:A:92:ASN:OD1	2.32	0.63
1:A:7:VAL:CG1	1:A:8:LEU:H	2.01	0.63
1:A:199:VAL:HG21	1:A:255:GLU:OE1	1.99	0.62
1:A:251:MET:HG3	1:A:281:LYS:HB2	1.81	0.62
1:A:36:ILE:HD11	1:A:38:TRP:CZ2	2.35	0.62
1:A:127:ILE:HG22	1:A:128:THR:N	2.14	0.61
1:A:78:LYS:HE2	1:A:210:LEU:O	2.02	0.60
1:A:256:ASP:O	1:A:257:CYS:C	2.44	0.60
1:A:44:ASP:HB3	1:A:46:LYS:H	1.65	0.60
1:A:181:ASN:HB3	1:A:184:LYS:HD3	1.84	0.60
1:A:159:ASN:OD1	1:A:288:PRO:HD3	2.02	0.60
1:A:70:LYS:HG2	1:A:72:GLU:HG2	1.84	0.59
1:A:146:MET:HB3	1:A:150:MET:HE2	1.84	0.58
1:A:230:ASN:HD22	1:A:232:SER:H	1.49	0.58
1:A:160:LEU:HD22	1:A:313:ALA:HB3	1.86	0.58
1:A:90:LEU:CD1	1:A:93:LEU:HD12	2.34	0.58
1:A:70:LYS:HE2	1:A:72:GLU:OE2	2.04	0.57
1:A:130:ILE:HG22	1:A:130:ILE:O	2.05	0.56
1:A:167:THR:CG2	1:A:202:ARG:HD2	2.35	0.56
1:A:218:ILE:HD12	1:A:218:ILE:N	2.20	0.56
1:A:36:ILE:HD11	1:A:38:TRP:CE2	2.41	0.56
1:A:113:CYS:O	1:A:176:HIS:HE1	1.88	0.55
1:A:3:GLN:HB2	1:A:4:PRO:CD	2.36	0.55
1:A:83:LYS:HD3	1:A:83:LYS:N	2.14	0.55
1:A:12:HIS:CE1	1:A:337:ILE:HG13	2.44	0.52
1:A:12:HIS:HE1	1:A:333:TRP:O	1.91	0.52
1:A:199:VAL:CG2	1:A:255:GLU:OE1	2.57	0.51
1:A:259:PHE:O	1:A:259:PHE:CG	2.63	0.51
1:A:83:LYS:H	1:A:83:LYS:CD	2.03	0.51
1:A:182:LYS:O	1:A:184:LYS:HD2	2.10	0.51
1:A:167:THR:HG22	1:A:202:ARG:CD	2.40	0.51
1:A:19:LYS:HG3	1:A:20:ILE:N	2.26	0.50
1:A:128:THR:CG2	1:A:129:HIS:N	2.53	0.50
1:A:70:LYS:HE2	1:A:72:GLU:CD	2.36	0.50
1:A:166:LYS:O	1:A:169:VAL:HG12	2.12	0.50
1:A:275:SER:O	1:A:276:SER:C	2.55	0.50
1:A:12:HIS:CE1	1:A:336:SER:HG	2.30	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:ILE:CG2	1:A:128:THR:N	2.75	0.49
1:A:71:PRO:HG3	1:A:77:TYR:CZ	2.47	0.49
1:A:63:LYS:HZ2	1:A:133:GLY:HA3	1.78	0.48
1:A:168:THR:O	1:A:172:ILE:HG13	2.13	0.48
1:A:288:PRO:HG3	1:A:354:ALA:C	2.39	0.48
1:A:175:ALA:O	1:A:179:CYS:SG	2.72	0.48
1:A:118:VAL:HG21	1:A:148:LYS:HE2	1.96	0.48
1:A:256:ASP:HB3	1:A:259:PHE:CE2	2.47	0.48
1:A:203:THR:O	1:A:207:ILE:HG13	2.15	0.47
1:A:167:THR:HG21	1:A:202:ARG:HD2	1.97	0.47
1:A:163:GLN:HG3	1:A:352:HIS:CE1	2.49	0.47
1:A:356:PHE:O	1:A:357:GLU:HB2	2.14	0.47
1:A:269:ALA:O	1:A:272:PRO:HD2	2.14	0.47
1:A:30:ARG:HB3	1:A:37:HIS:CE1	2.50	0.47
1:A:291:LEU:HD22	1:A:291:LEU:N	2.30	0.47
1:A:146:MET:O	1:A:150:MET:HG3	2.15	0.47
1:A:167:THR:HG22	1:A:202:ARG:HD3	1.97	0.47
1:A:79:ASP:O	1:A:86:ARG:CZ	2.63	0.46
1:A:129:HIS:HA	1:A:134:VAL:HG22	1.97	0.46
1:A:298:TRP:O	1:A:302:VAL:HG23	2.15	0.46
1:A:5:ILE:HG22	1:A:6:ASN:CG	2.41	0.45
1:A:127:ILE:HD11	1:A:137:VAL:HG13	1.97	0.45
1:A:166:LYS:HD2	1:A:167:THR:N	2.31	0.45
1:A:7:VAL:O	1:A:8:LEU:HD13	2.16	0.45
1:A:259:PHE:O	1:A:259:PHE:CD1	2.68	0.45
1:A:33:GLU:HG2	1:A:34:ASP:OD1	2.16	0.45
1:A:293:HIS:O	1:A:297:ILE:HG13	2.17	0.45
1:A:167:THR:CG2	1:A:202:ARG:CD	2.95	0.45
1:A:166:LYS:HZ3	1:A:255:GLU:HG3	1.82	0.45
1:A:253:TYR:HE2	1:A:255:GLU:HB2	1.79	0.45
1:A:166:LYS:HD2	1:A:166:LYS:C	2.42	0.44
1:A:70:LYS:HD3	1:A:73:LEU:CD1	2.47	0.44
1:A:166:LYS:NZ	1:A:255:GLU:OE2	2.45	0.44
1:A:87:TYR:C	1:A:88:MET:HG2	2.41	0.44
1:A:128:THR:O	1:A:129:HIS:HB2	2.17	0.44
1:A:255:GLU:O	1:A:256:ASP:CB	2.65	0.44
1:A:162:ARG:O	1:A:163:GLN:HB2	2.18	0.43
1:A:230:ASN:C	1:A:230:ASN:ND2	2.68	0.43
1:A:251:MET:CG	1:A:281:LYS:HB2	2.47	0.43
1:A:157:VAL:HG13	1:A:157:VAL:O	2.18	0.43
1:A:163:GLN:HG3	1:A:352:HIS:HE1	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:THR:HG22	1:A:101:GLN:N	2.33	0.43
1:A:3:GLN:HB2	1:A:4:PRO:HD2	2.00	0.43
1:A:103:THR:O	1:A:107:VAL:HG23	2.18	0.43
1:A:138:GLN:H	1:A:138:GLN:HG2	1.66	0.43
1:A:14:LEU:O	1:A:21:LEU:HD12	2.18	0.42
1:A:74:PHE:CD1	1:A:74:PHE:C	2.98	0.42
1:A:139:LEU:HD22	1:A:143:GLN:HB2	2.01	0.42
1:A:127:ILE:CD1	1:A:137:VAL:HG22	2.49	0.42
1:A:238:ALA:O	1:A:239:SER:C	2.61	0.42
1:A:30:ARG:HD3	1:A:37:HIS:NE2	2.35	0.42
1:A:200:LEU:O	1:A:204:LYS:HG3	2.20	0.42
1:A:130:ILE:O	1:A:130:ILE:CG2	2.67	0.41
1:A:163:GLN:NE2	2:A:415:HOH:O	2.54	0.41
1:A:12:HIS:CE1	1:A:333:TRP:O	2.72	0.41
1:A:67:ASN:HD22	1:A:77:TYR:HA	1.86	0.41
1:A:2:GLU:H	1:A:2:GLU:HG2	1.49	0.41
1:A:5:ILE:HG22	1:A:6:ASN:ND2	2.35	0.40
1:A:71:PRO:HA	1:A:77:TYR:CG	2.56	0.40
1:A:326:GLU:O	1:A:327:ASP:HB2	2.20	0.40
1:A:269:ALA:C	1:A:272:PRO:HD2	2.46	0.40
1:A:99:LYS:HD2	1:A:99:LYS:N	2.36	0.40
1:A:219:VAL:O	1:A:220:GLU:HG2	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	355/385 (92%)	317 (89%)	31 (9%)	7 (2%)	6 10

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	132	TYR
1	A	129	HIS
1	A	256	ASP
1	A	257	CYS
1	A	60	LYS
1	A	343	SER
1	A	7	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	315/339 (93%)	295 (94%)	20 (6%)	16	34

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	GLU
1	A	8	LEU
1	A	10	ASP
1	A	83	LYS
1	A	98	ILE
1	A	101	GLN
1	A	139	LEU
1	A	209	LEU
1	A	230	ASN
1	A	251	MET
1	A	257	CYS
1	A	259	PHE
1	A	268	LEU
1	A	272	PRO
1	A	276	SER
1	A	282	ILE
1	A	345	LEU
1	A	350	GLN
1	A	351	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	ASN
1	A	82	ASN
1	A	176	HIS
1	A	191	HIS
1	A	230	ASN
1	A	347	GLN
1	A	352	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 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 [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, 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	357/385 (92%)	0.30	9 (2%) 58 54	36, 68, 117, 191	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	7	VAL	3.3
1	A	131	ASP	2.9
1	A	1	MET	2.8
1	A	132	TYR	2.4
1	A	44	ASP	2.3
1	A	274	ILE	2.2
1	A	214	LEU	2.1
1	A	306	SER	2.1
1	A	301	ALA	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.