



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:19 PM UTC

PDB ID : 5IXA / pdb_00005ixa
Title : HCMV DNA polymerase processivity subunit UL44 at neutral pH and low salt
Authors : Chen, H.; Coen, D.M.; Hogle, J.M.; Filman, D.J.
Deposited on : 2016-03-23
Resolution : 2.68 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

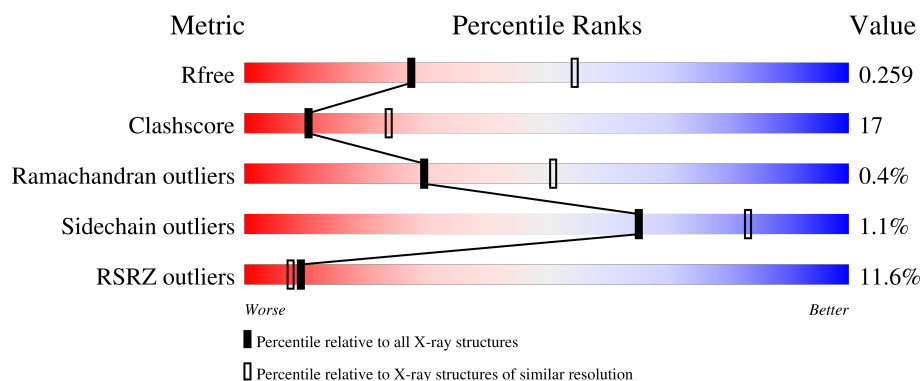
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.68 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	5070 (2.70-2.66)
Clashscore	190562	5409 (2.70-2.66)
Ramachandran outliers	187476	5324 (2.70-2.66)
Sidechain outliers	187428	5324 (2.70-2.66)
RSRZ outliers	180081	5070 (2.70-2.66)

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	290	11% 59% 27% 14%
1	B	290	9% 57% 29% 13%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4023 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 processivity factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	250	Total	C	N	O	S	0	0	0
			1956	1249	329	368	10			
1	B	251	Total	C	N	O	S	0	0	0
			1969	1260	330	369	10			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	205	SER	ALA	SEE REMARK 999	UNP P16790
B	205	SER	ALA	SEE REMARK 999	UNP P16790

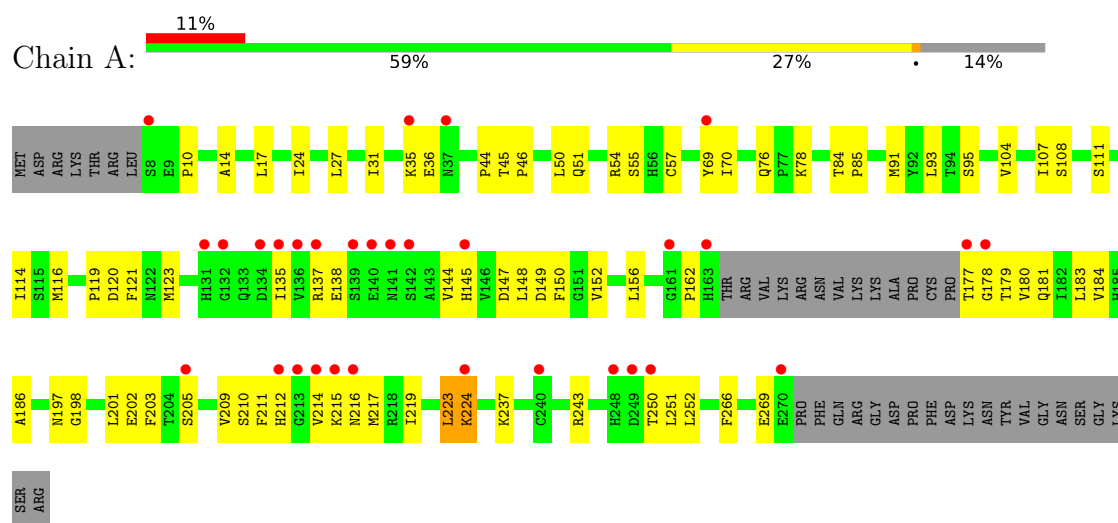
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	44	Total	O	0	0
			44	44		
2	B	54	Total	O	0	0
			54	54		

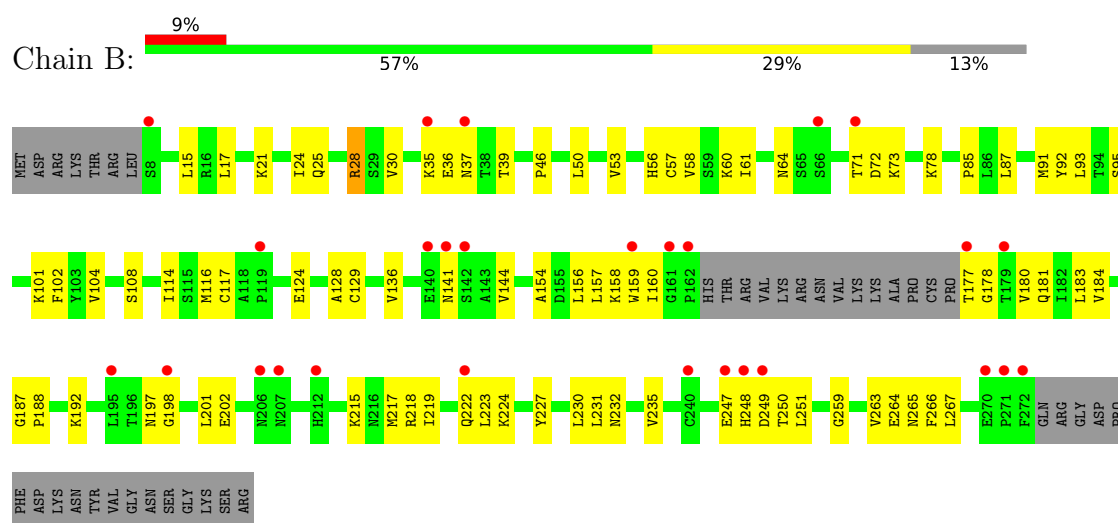
3 Residue-property plots

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 polymerase processivity factor



• Molecule 1: DNA polymerase processivity factor



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.03Å 100.83Å 109.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.11 – 2.68 37.11 – 2.68	Depositor EDS
% Data completeness (in resolution range)	97.7 (37.11-2.68) 97.3 (37.11-2.68)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.210 , 0.259 0.209 , 0.259	Depositor DCC
R_{free} test set	1187 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	43.7	Xtriage
Anisotropy	0.185	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 58.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4023	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.24% 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	0.53	0/1994	0.94	4/2710 (0.1%)
1	B	0.47	0/2009	0.82	1/2731 (0.0%)
All	All	0.50	0/4003	0.88	5/5441 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	224	LYS	CG-CD-CE	10.90	136.37	111.30
1	A	223	LEU	CA-C-N	-7.61	107.82	121.14
1	A	223	LEU	C-N-CA	-7.61	107.82	121.14
1	A	224	LYS	CB-CG-CD	-7.60	93.82	111.30
1	B	28	ARG	NE-CZ-NH2	-5.75	114.03	119.20

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	1956	0	1982	68	0
1	B	1969	0	1996	68	0
2	A	44	0	0	12	2
2	B	54	0	0	15	2
All	All	4023	0	3978	133	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 17.

All (133) 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:35:LYS:CB	2:B:302:HOH:O	2.08	1.00
1:B:57:CYS:SG	2:B:353:HOH:O	2.20	1.00
1:B:35:LYS:N	2:B:302:HOH:O	1.97	0.96
1:A:54:ARG:HG2	1:A:55:SER:H	1.28	0.96
1:A:57:CYS:SG	2:A:342:HOH:O	2.27	0.93
1:A:120:ASP:N	2:A:301:HOH:O	1.81	0.92
1:B:177:THR:HG22	1:B:178:GLY:H	1.37	0.88
1:B:35:LYS:CA	2:B:302:HOH:O	2.22	0.88
1:A:24:ILE:HD12	1:A:95:SER:HA	1.57	0.86
1:B:249:ASP:OD1	2:B:301:HOH:O	1.92	0.86
1:B:24:ILE:HD12	1:B:95:SER:HA	1.63	0.81
1:B:35:LYS:HB3	2:B:302:HOH:O	1.76	0.79
1:B:219:ILE:HD12	1:B:250:THR:HG22	1.63	0.78
1:A:211:PHE:HD1	1:A:214:VAL:HG11	1.49	0.78
1:A:119:PRO:HB2	2:A:301:HOH:O	1.84	0.78
1:A:76:GLN:OE1	2:A:302:HOH:O	2.01	0.78
1:B:37:ASN:ND2	1:B:53:VAL:O	2.17	0.77
1:A:119:PRO:CA	2:A:301:HOH:O	2.34	0.75
1:B:264:GLU:OE2	2:B:303:HOH:O	2.04	0.75
1:A:119:PRO:CB	2:A:301:HOH:O	2.36	0.72
1:A:14:ALA:HB3	1:A:69:TYR:HD1	1.55	0.71
1:B:144:VAL:HG21	1:B:184:VAL:HG11	1.71	0.71
1:B:21:LYS:O	1:B:25:GLN:HG2	1.91	0.70
1:B:156:LEU:HD23	1:B:201:LEU:HD23	1.72	0.70
1:B:215:LYS:NZ	2:B:308:HOH:O	2.23	0.70
1:B:35:LYS:HG2	1:B:36:GLU:H	1.56	0.70
1:B:181:GLN:NE2	1:B:218:ARG:HG2	2.07	0.70
1:A:137:ARG:NH1	2:A:305:HOH:O	2.24	0.69
1:A:145:HIS:CE1	1:A:212:HIS:HD2	2.11	0.69
1:B:192:LYS:HG3	1:B:202:GLU:HB2	1.75	0.68
1:A:93:LEU:HD22	1:A:116:MET:HE1	1.75	0.67
1:A:138:GLU:OE2	1:A:243:ARG:NH1	2.28	0.67
1:B:60:LYS:NZ	1:B:136:VAL:O	2.26	0.66
1:A:14:ALA:HB3	1:A:69:TYR:CD1	2.30	0.66
1:A:54:ARG:HG2	1:A:55:SER:N	2.05	0.66
1:B:128:ALA:O	2:B:304:HOH:O	2.12	0.65
1:B:177:THR:CG2	1:B:178:GLY:H	2.09	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:ILE:HD12	1:A:250:THR:HG22	1.79	0.64
1:A:35:LYS:HG2	1:A:36:GLU:H	1.63	0.63
1:B:180:VAL:HB	1:B:223:LEU:HD13	1.81	0.63
1:B:30:VAL:HG11	1:B:61:ILE:HD11	1.80	0.62
1:A:224:LYS:NZ	2:A:303:HOH:O	2.01	0.62
1:B:25:GLN:O	1:B:28:ARG:HG3	1.98	0.62
1:A:144:VAL:HG13	1:A:214:VAL:HG12	1.81	0.61
1:B:197:ASN:OD1	1:B:198:GLY:N	2.33	0.61
1:A:50:LEU:HD21	1:A:91:MET:HE2	1.83	0.61
1:B:177:THR:HG22	1:B:178:GLY:N	2.15	0.60
1:A:145:HIS:NE2	1:A:212:HIS:HD2	1.99	0.60
1:A:119:PRO:N	2:A:301:HOH:O	2.35	0.59
1:A:145:HIS:CE1	1:A:212:HIS:CD2	2.90	0.59
1:B:104:VAL:HG22	1:B:114:ILE:HG13	1.85	0.59
1:A:219:ILE:HD13	1:A:252:LEU:HB2	1.86	0.58
1:A:147:ASP:OD2	1:A:212:HIS:NE2	2.29	0.58
1:B:73:LYS:NZ	2:B:305:HOH:O	2.18	0.57
1:B:181:GLN:HE21	1:B:218:ARG:HG2	1.69	0.57
1:A:181:GLN:HG2	1:A:183:LEU:HD12	1.87	0.57
1:A:111:SER:O	2:A:304:HOH:O	2.18	0.56
1:B:141:ASN:ND2	1:B:247:GLU:OE2	2.40	0.55
1:B:64:ASN:ND2	1:B:259:GLY:O	2.36	0.54
1:B:93:LEU:HG	1:B:116:MET:HE1	1.88	0.53
1:A:27:LEU:O	1:A:31:ILE:HG12	2.08	0.53
1:A:138:GLU:HG3	2:A:317:HOH:O	2.08	0.53
1:A:177:THR:HG23	1:A:178:GLY:H	1.74	0.53
1:A:121:PHE:HB2	1:B:85:PRO:HG2	1.90	0.52
1:A:156:LEU:HD23	1:A:201:LEU:HD23	1.91	0.52
1:B:227:TYR:CE2	1:B:231:LEU:HD11	2.45	0.52
1:A:197:ASN:OD1	1:A:198:GLY:N	2.43	0.51
1:B:24:ILE:CD1	1:B:95:SER:HA	2.39	0.50
1:B:154:ALA:O	1:B:158:LYS:HG2	2.10	0.50
1:B:248:HIS:HA	2:B:309:HOH:O	2.11	0.50
1:A:224:LYS:CE	2:A:303:HOH:O	2.54	0.50
1:A:219:ILE:HG13	1:A:269:GLU:HG2	1.94	0.50
1:B:78:LYS:NZ	1:B:108:SER:O	2.45	0.49
1:B:157:LEU:C	1:B:159:TRP:H	2.19	0.49
1:A:104:VAL:HG22	1:A:114:ILE:HG13	1.95	0.48
1:A:149:ASP:HB3	1:A:152:VAL:HG23	1.96	0.48
1:A:180:VAL:HB	1:A:223:LEU:HD13	1.94	0.48
1:B:248:HIS:CA	2:B:309:HOH:O	2.61	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:PHE:CE1	1:A:237:LYS:HA	2.49	0.47
1:A:215:LYS:HE2	1:A:215:LYS:N	2.29	0.47
1:A:93:LEU:CD2	1:A:116:MET:HE1	2.41	0.47
1:A:123:MET:SD	1:B:87:LEU:HD21	2.55	0.47
1:A:144:VAL:HG22	1:A:214:VAL:HA	1.96	0.47
1:A:186:ALA:N	1:A:216:ASN:OD1	2.48	0.47
1:B:71:THR:HG22	1:B:72:ASP:CB	2.44	0.46
1:B:58:VAL:HG11	1:B:136:VAL:HG21	1.97	0.46
1:A:251:LEU:HD11	1:A:266:PHE:HB3	1.96	0.46
1:A:78:LYS:NZ	1:A:108:SER:O	2.49	0.46
1:B:92:TYR:O	1:B:95:SER:OG	2.31	0.46
1:A:210:SER:OG	1:A:212:HIS:CE1	2.69	0.46
1:B:251:LEU:HD21	1:B:266:PHE:HD2	1.80	0.46
1:B:217:MET:HB3	1:B:217:MET:HE2	1.56	0.45
1:A:44:PRO:HA	1:A:70:ILE:HD12	1.97	0.45
1:B:15:LEU:HD13	1:B:102:PHE:CD2	2.51	0.45
1:A:147:ASP:CG	1:A:212:HIS:HE2	2.21	0.45
1:B:183:LEU:HD12	1:B:218:ARG:HH11	1.81	0.45
1:A:145:HIS:CD2	1:A:212:HIS:HD2	2.35	0.44
1:A:45:THR:N	1:A:46:PRO:HA	2.32	0.44
1:A:51:GLN:OE1	1:A:135:ILE:HA	2.17	0.44
1:A:217:MET:HB2	1:A:217:MET:HE3	1.34	0.44
1:B:160:ILE:CG2	1:B:223:LEU:HD11	2.47	0.44
1:A:144:VAL:HG11	1:A:184:VAL:HG21	2.00	0.44
1:B:157:LEU:HD21	1:B:230:LEU:HB2	2.00	0.44
1:A:45:THR:HB	1:A:46:PRO:C	2.43	0.43
1:A:145:HIS:CE1	1:A:243:ARG:HH21	2.35	0.43
1:A:17:LEU:HD23	1:A:17:LEU:HA	1.66	0.43
1:A:186:ALA:HA	1:A:214:VAL:O	2.18	0.43
1:B:184:VAL:HG13	1:B:215:LYS:O	2.19	0.43
1:B:232:ASN:HD22	1:B:265:ASN:HD21	1.67	0.43
1:B:50:LEU:HD11	1:B:91:MET:HE2	2.01	0.43
1:B:93:LEU:CD2	1:B:116:MET:HE1	2.49	0.43
1:B:56:HIS:O	2:B:306:HOH:O	2.20	0.42
1:A:10:PRO:HG2	1:A:107:ILE:HD11	2.01	0.42
1:B:267:LEU:HD12	1:B:267:LEU:N	2.34	0.42
1:B:17:LEU:HD23	1:B:17:LEU:HA	1.72	0.42
1:B:93:LEU:CG	1:B:116:MET:HE1	2.50	0.42
1:B:56:HIS:C	2:B:306:HOH:O	2.63	0.42
1:A:148:LEU:HD23	1:A:209:VAL:HG13	2.01	0.42
1:A:84:THR:HA	1:A:85:PRO:HD3	1.92	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:GLN:OE1	1:B:224:LYS:HB3	2.19	0.42
1:A:144:VAL:HG11	1:A:184:VAL:HG11	2.02	0.41
1:B:232:ASN:HB3	1:B:263:VAL:HG11	2.01	0.41
1:B:71:THR:HG22	1:B:72:ASP:HB2	2.02	0.41
1:B:187:GLY:HA2	1:B:188:PRO:HA	1.83	0.41
1:A:224:LYS:O	1:A:224:LYS:HD2	2.19	0.41
1:A:202:GLU:HG2	1:A:203:PHE:N	2.36	0.41
1:A:181:GLN:HG2	1:A:183:LEU:CD1	2.50	0.40
1:B:160:ILE:HG23	1:B:223:LEU:HD11	2.03	0.40
1:A:120:ASP:HB3	1:B:128:ALA:HB2	2.03	0.40
1:B:102:PHE:HE1	1:B:114:ILE:HG23	1.86	0.40
1:A:147:ASP:HB2	1:A:210:SER:HB3	2.02	0.40
1:B:46:PRO:CG	2:B:320:HOH:O	2.69	0.40
1:B:101:LYS:HB3	1:B:117:CYS:HB2	2.03	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:339:HOH:O	2:B:317:HOH:O[3_656]	2.12	0.08
2:A:336:HOH:O	2:B:350:HOH:O[4_466]	2.17	0.03

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	246/290 (85%)	227 (92%)	17 (7%)	2 (1%)	16	34
1	B	247/290 (85%)	230 (93%)	17 (7%)	0	100	100
All	All	493/580 (85%)	457 (93%)	34 (7%)	2 (0%)	30	51

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	162	PRO
1	A	205	SER

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	227/264 (86%)	226 (100%)	1 (0%)	84	93
1	B	229/264 (87%)	225 (98%)	4 (2%)	53	77
All	All	456/528 (86%)	451 (99%)	5 (1%)	65	83

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	179	THR
1	B	39	THR
1	B	124	GLU
1	B	129	CYS
1	B	235	VAL

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

Mol	Chain	Res	Type
1	A	25	GLN
1	A	37	ASN
1	A	131	HIS
1	A	212	HIS
1	B	25	GLN
1	B	56	HIS
1	B	131	HIS
1	B	181	GLN
1	B	212	HIS
1	B	265	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 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

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, 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	250/290 (86%)	0.42	31 (12%) 8 7	25, 44, 102, 131	0
1	B	251/290 (86%)	0.35	27 (10%) 11 9	20, 48, 87, 124	1 (0%)
All	All	501/580 (86%)	0.38	58 (11%) 9 7	20, 46, 92, 131	1 (0%)

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	214	VAL	6.5
1	A	139	SER	5.4
1	A	137	ARG	4.7
1	A	178	GLY	4.5
1	A	213	GLY	4.1
1	B	177	THR	3.9
1	A	163	HIS	3.8
1	B	272	PHE	3.8
1	B	162	PRO	3.7
1	A	215	LYS	3.6
1	B	140	GLU	3.6
1	A	132	GLY	3.4
1	B	141	ASN	3.3
1	A	69	TYR	3.2
1	B	37	ASN	3.2
1	B	179	THR	3.1
1	A	131	HIS	3.1
1	A	141	ASN	2.9
1	A	248	HIS	2.8
1	A	8	SER	2.8
1	A	205	SER	2.7
1	A	224	LYS	2.7
1	B	71	THR	2.6
1	B	271	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	134	ASP	2.6
1	A	240	CYS	2.6
1	A	212	HIS	2.5
1	A	140	GLU	2.5
1	B	249	ASP	2.5
1	A	177	THR	2.5
1	B	207	ASN	2.5
1	A	249	ASP	2.5
1	A	135	ILE	2.4
1	B	206	ASN	2.4
1	B	8	SER	2.4
1	A	35	LYS	2.4
1	A	250	THR	2.3
1	B	270	GLU	2.3
1	B	35	LYS	2.3
1	B	142	SER	2.3
1	A	161	GLY	2.3
1	A	136	VAL	2.3
1	B	161	GLY	2.3
1	B	222	GLN	2.3
1	B	247	GLU	2.3
1	A	270	GLU	2.2
1	B	240	CYS	2.2
1	A	142	SER	2.2
1	A	145	HIS	2.1
1	B	248	HIS	2.1
1	B	212	HIS	2.1
1	B	195	LEU	2.1
1	B	119	PRO	2.1
1	A	216	ASN	2.1
1	B	198	GLY	2.1
1	B	66	SER	2.0
1	B	159	TRP	2.0
1	A	37	ASN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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.