



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:24 AM UTC

PDB ID : 3LH8 / pdb_00003lh8
Title : Crystal structure of mouse VPS26B in spacegroup P41 21 2
Authors : Collins, B.; Shaw, D.; Norwood, S.
Deposited on : 2010-01-21
Resolution : 2.60 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

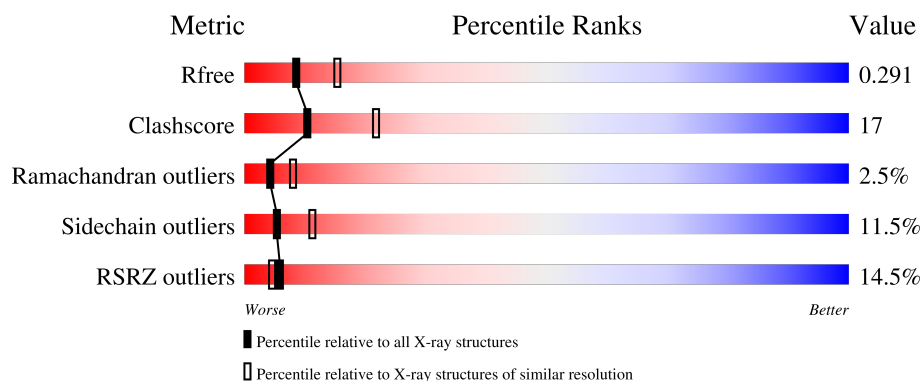
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	340	13% 55% 24% • • 17%
1	B	340	11% 54% 23% 5% • 17%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 4921 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 Vacuolar protein sorting-associated protein 26B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	283	Total	C	N	O	S	0	1	0
			2354	1508	396	444	6			
1	B	282	Total	C	N	O	S	0	1	0
			2355	1509	396	444	6			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	MET	-	expression tag	UNP Q8C0E2
A	-2	GLY	-	expression tag	UNP Q8C0E2
A	-1	SER	-	expression tag	UNP Q8C0E2
A	0	HIS	-	expression tag	UNP Q8C0E2
A	1	HIS	-	expression tag	UNP Q8C0E2
A	2	HIS	-	expression tag	UNP Q8C0E2
A	3	HIS	-	expression tag	UNP Q8C0E2
A	4	HIS	-	expression tag	UNP Q8C0E2
A	5	HIS	-	expression tag	UNP Q8C0E2
A	6	MET	-	expression tag	UNP Q8C0E2
A	197	SER	LEU	engineered mutation	UNP Q8C0E2
A	199	GLU	ARG	engineered mutation	UNP Q8C0E2
B	-3	MET	-	expression tag	UNP Q8C0E2
B	-2	GLY	-	expression tag	UNP Q8C0E2
B	-1	SER	-	expression tag	UNP Q8C0E2
B	0	HIS	-	expression tag	UNP Q8C0E2
B	1	HIS	-	expression tag	UNP Q8C0E2
B	2	HIS	-	expression tag	UNP Q8C0E2
B	3	HIS	-	expression tag	UNP Q8C0E2
B	4	HIS	-	expression tag	UNP Q8C0E2
B	5	HIS	-	expression tag	UNP Q8C0E2
B	6	MET	-	expression tag	UNP Q8C0E2
B	197	SER	LEU	engineered mutation	UNP Q8C0E2
B	199	GLU	ARG	engineered mutation	UNP Q8C0E2

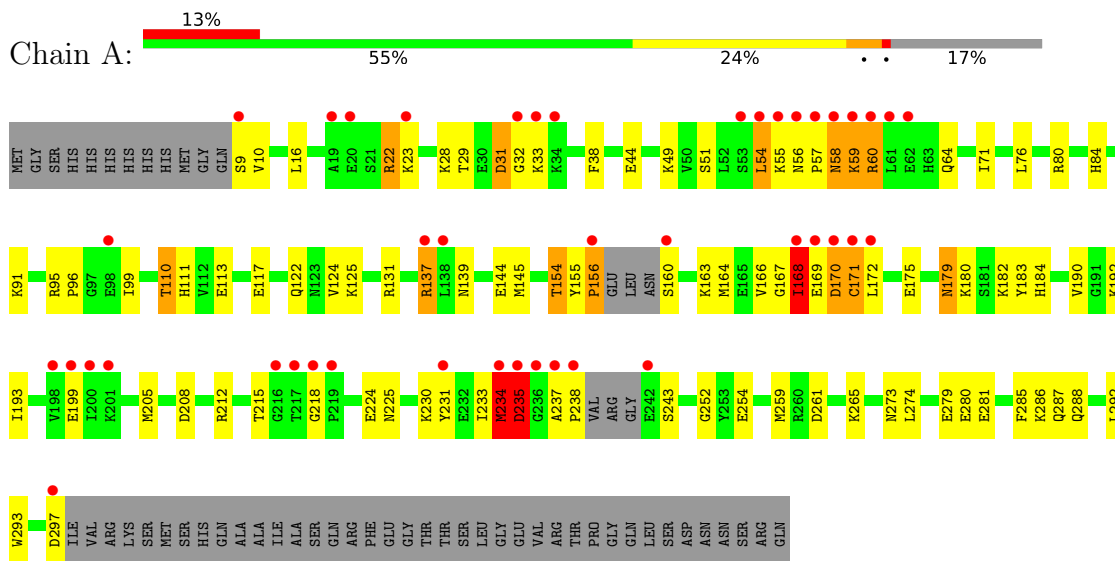
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	117	Total 117	O 117	0	0
2	B	95	Total 95	O 95	0	0

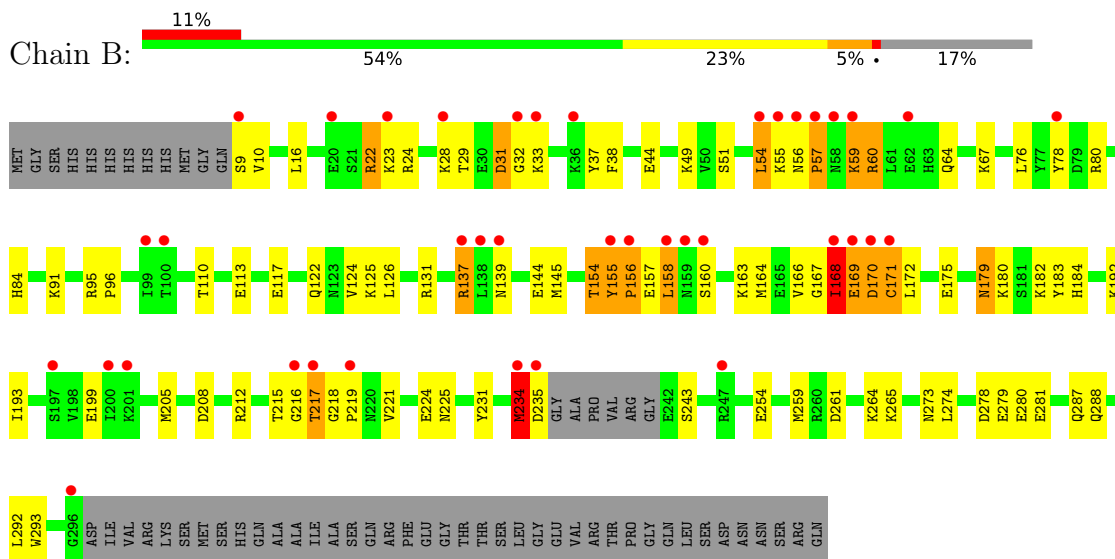
3 Residue-property plots

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: Vacuolar protein sorting-associated protein 26B



- Molecule 1: Vacuolar protein sorting-associated protein 26B



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	118.16Å 118.16Å 193.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.90 – 2.60 40.90 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.7 (40.90-2.60) 98.8 (40.90-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.261 , 0.292 0.259 , 0.291	Depositor DCC
R_{free} test set	2139 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	48.2	Xtriage
Anisotropy	0.022	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	4921	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 2.63% 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	0.55	0/2404	0.91	7/3233 (0.2%)
1	B	0.55	0/2405	0.93	6/3235 (0.2%)
All	All	0.55	0/4809	0.92	13/6468 (0.2%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	155	TYR	CA-C-N	6.06	127.42	119.84
1	B	155	TYR	C-N-CA	6.06	127.42	119.84
1	A	234	MET	N-CA-C	5.93	118.53	110.35
1	B	234	MET	N-CA-C	5.85	118.42	110.35
1	A	58	ASN	CB-CA-C	-5.82	103.05	111.14
1	B	57	PRO	CA-C-N	-5.75	112.31	122.26
1	B	57	PRO	C-N-CA	-5.75	112.31	122.26
1	A	218	GLY	CA-C-N	5.65	125.74	119.87
1	A	218	GLY	C-N-CA	5.65	125.74	119.87
1	B	216	GLY	N-CA-C	5.18	125.47	113.18
1	A	235	ASP	N-CA-C	5.08	121.62	110.80
1	A	95	ARG	CA-C-N	-5.04	114.76	119.90
1	A	95	ARG	C-N-CA	-5.04	114.76	119.90

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	2354	0	2346	82	0
1	B	2355	0	2351	79	1
2	A	117	0	0	7	1
2	B	95	0	0	6	0
All	All	4921	0	4697	160	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 17.

All (160) 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:252:GLY:HA3	2:A:440:HOH:O	1.45	1.17
1:A:60:ARG:HH11	1:A:60:ARG:HG3	0.99	1.16
1:B:60:ARG:HG2	1:B:60:ARG:NH1	1.61	1.05
1:B:84:HIS:HD2	1:B:265:LYS:HE2	1.20	1.05
1:A:110:THR:HA	2:A:428:HOH:O	1.59	1.02
1:B:60:ARG:HG2	1:B:60:ARG:HH11	0.86	1.01
1:A:60:ARG:HG3	1:A:60:ARG:NH1	1.68	0.99
1:B:84:HIS:CD2	1:B:265:LYS:HE2	2.05	0.91
1:B:169:GLU:HG2	1:B:170:ASP:H	1.36	0.90
1:A:168:ILE:HG22	1:A:169:GLU:H	1.37	0.88
1:A:169:GLU:HG2	1:A:170:ASP:H	1.38	0.87
1:B:168:ILE:HG22	1:B:169:GLU:H	1.40	0.86
1:B:156:PRO:HB2	1:B:158:LEU:H	1.43	0.84
1:A:84:HIS:HD2	1:A:265:LYS:HE2	1.42	0.83
1:A:84:HIS:CD2	1:A:265:LYS:HE2	2.13	0.83
1:A:179:ASN:HD22	1:A:180:LYS:N	1.77	0.81
1:B:60:ARG:HH11	1:B:60:ARG:CG	1.82	0.80
1:B:179:ASN:HD22	1:B:180:LYS:N	1.81	0.78
1:B:155:TYR:CD1	1:B:182:LYS:HD3	2.19	0.77
1:A:155:TYR:CD1	1:A:182:LYS:HD3	2.19	0.77
1:A:169:GLU:CG	1:A:170:ASP:H	1.98	0.77
1:B:155:TYR:HE1	1:B:182:LYS:HE3	1.51	0.75
1:B:169:GLU:CG	1:B:170:ASP:H	2.01	0.74
1:A:155:TYR:HE1	1:A:182:LYS:HE3	1.52	0.73
1:B:164:MET:HE3	1:B:274:LEU:HB2	1.71	0.72
1:A:60:ARG:HH11	1:A:60:ARG:CG	1.91	0.71
1:A:155:TYR:HD1	1:A:182:LYS:HD3	1.55	0.71
1:A:154:THR:C	1:A:155:TYR:HD2	2.00	0.70
1:B:155:TYR:HD1	1:B:182:LYS:HD3	1.56	0.69
1:A:111:HIS:N	2:A:428:HOH:O	2.21	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:MET:HE3	1:A:274:LEU:HB2	1.74	0.68
1:A:164:MET:CE	1:A:274:LEU:HB2	2.24	0.67
1:B:154:THR:C	1:B:155:TYR:HD2	2.03	0.65
1:B:164:MET:CE	1:B:274:LEU:HB2	2.26	0.64
1:A:237:ALA:N	1:A:238:PRO:HD3	2.12	0.64
1:A:193:ILE:HG21	1:A:205:MET:HE1	1.80	0.64
1:B:168:ILE:HG22	1:B:169:GLU:N	2.12	0.63
1:A:64:GLN:HA	1:A:96:PRO:HB3	1.82	0.62
1:A:168:ILE:HG22	1:A:169:GLU:N	2.10	0.62
1:B:171:CYS:HA	1:B:199:GLU:HB2	1.81	0.62
1:A:170:ASP:O	1:A:171:CYS:HB3	1.99	0.62
1:B:217:THR:H	1:B:221:VAL:HG22	1.64	0.62
1:A:110:THR:CA	2:A:428:HOH:O	2.32	0.62
1:B:60:ARG:NH1	1:B:60:ARG:CG	2.50	0.62
1:A:171:CYS:HA	1:A:199:GLU:HB2	1.81	0.62
1:A:155:TYR:HE1	1:A:182:LYS:CE	2.13	0.61
1:B:155:TYR:HE1	1:B:182:LYS:CE	2.13	0.61
1:B:64:GLN:HA	1:B:96:PRO:HB3	1.84	0.60
1:B:122:GLN:HG2	1:B:224[A]:GLU:HG3	1.83	0.60
1:B:137:ARG:HG3	2:B:404:HOH:O	2.01	0.60
1:A:22:ARG:NH2	1:A:44:GLU:OE1	2.35	0.59
1:B:163:LYS:HD3	1:B:175:GLU:OE2	2.01	0.59
1:B:170:ASP:O	1:B:171:CYS:HB3	2.01	0.59
1:B:122:GLN:CG	1:B:224[A]:GLU:HG3	2.33	0.59
1:B:155:TYR:HB3	1:B:156:PRO:HD2	1.84	0.58
1:A:212:ARG:HG2	1:A:225:ASN:OD1	2.02	0.58
1:A:169:GLU:CG	1:A:170:ASP:N	2.66	0.58
1:A:163:LYS:HD3	1:A:175:GLU:OE2	2.04	0.58
1:A:155:TYR:HB3	1:A:156:PRO:HD2	1.86	0.58
1:B:234:MET:HG3	1:B:235:ASP:OD2	2.03	0.58
1:A:254:GLU:HA	1:A:254:GLU:OE1	2.04	0.57
1:A:234:MET:HG3	1:A:235:ASP:OD2	2.04	0.56
1:B:160:SER:O	1:B:179:ASN:O	2.24	0.56
1:B:169:GLU:CG	1:B:170:ASP:N	2.68	0.56
1:B:193:ILE:HG21	1:B:205:MET:HE1	1.86	0.56
1:A:179:ASN:HD22	1:A:179:ASN:C	2.14	0.56
1:A:170:ASP:O	1:A:171:CYS:CB	2.54	0.55
1:B:156:PRO:HB2	1:B:158:LEU:N	2.18	0.54
1:A:131:ARG:HD3	1:A:144:GLU:OE2	2.06	0.54
1:B:208:ASP:HB2	1:B:273:ASN:HB3	1.90	0.54
1:A:286:LYS:HD3	2:B:364:HOH:O	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:GLN:HG2	1:A:224[A]:GLU:HG3	1.90	0.53
1:B:280:GLU:O	1:B:281:GLU:HB2	2.09	0.53
1:A:122:GLN:CG	1:A:224[A]:GLU:HG3	2.37	0.53
1:A:208:ASP:HB2	1:A:273:ASN:HB3	1.89	0.53
1:B:179:ASN:HD21	1:B:183:TYR:HE2	1.56	0.53
1:B:212:ARG:HG2	1:B:225:ASN:OD1	2.09	0.53
1:B:67:LYS:HG2	2:B:346:HOH:O	2.08	0.53
1:B:164:MET:HE3	1:B:274:LEU:CB	2.38	0.53
1:A:160:SER:O	1:A:179:ASN:O	2.26	0.52
1:B:22:ARG:NH2	1:B:44:GLU:OE1	2.43	0.52
1:B:254:GLU:HA	1:B:254:GLU:OE1	2.09	0.52
1:B:170:ASP:O	1:B:171:CYS:CB	2.57	0.52
1:B:155:TYR:CE1	1:B:182:LYS:HD3	2.43	0.52
1:A:155:TYR:CE1	1:A:182:LYS:HD3	2.44	0.52
1:A:56:ASN:HB3	1:A:59:LYS:HB3	1.92	0.52
1:A:280:GLU:O	1:A:281:GLU:HB2	2.10	0.51
1:A:168:ILE:HD12	1:A:171:CYS:SG	2.50	0.51
1:B:231:TYR:CD2	1:B:231:TYR:C	2.89	0.50
1:B:124:VAL:HG22	1:B:125:LYS:N	2.27	0.49
1:A:164:MET:HE3	1:A:274:LEU:CB	2.43	0.49
1:A:124:VAL:HG22	1:A:125:LYS:N	2.27	0.49
1:B:22:ARG:NH1	2:B:378:HOH:O	2.46	0.49
1:B:56:ASN:HB3	1:B:59:LYS:HB3	1.93	0.49
1:B:179:ASN:HD22	1:B:179:ASN:C	2.17	0.49
1:A:184:HIS:HA	1:A:293:TRP:O	2.13	0.49
1:A:193:ILE:CG2	1:A:205:MET:HE1	2.43	0.48
1:B:16:LEU:HG	1:B:145:MET:HE1	1.94	0.48
1:A:154:THR:C	1:A:155:TYR:CD2	2.87	0.48
1:A:54:LEU:HD12	1:A:54:LEU:HA	1.73	0.48
1:A:212:ARG:NH2	1:A:287:GLN:OE1	2.42	0.48
1:A:117:GLU:HB3	1:A:259:MET:HE2	1.96	0.48
1:B:193:ILE:CG2	1:B:205:MET:HE1	2.43	0.48
1:B:217:THR:HG21	1:B:264:LYS:O	2.13	0.48
1:A:155:TYR:CE1	1:A:182:LYS:CE	2.96	0.47
1:A:71:ILE:HD13	1:A:131:ARG:NH2	2.30	0.47
1:A:91:LYS:HB2	1:A:91:LYS:HE2	1.65	0.47
1:B:184:HIS:HA	1:B:293:TRP:O	2.15	0.47
1:B:117:GLU:HB3	1:B:259:MET:HE2	1.95	0.47
1:B:131:ARG:HD3	1:B:144:GLU:OE2	2.15	0.47
1:B:155:TYR:CE1	1:B:182:LYS:CE	2.96	0.47
1:A:59:LYS:HG2	1:A:60:ARG:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:TYR:O	1:A:292:LEU:HA	2.15	0.46
1:A:231:TYR:CD2	1:A:231:TYR:C	2.93	0.46
1:A:179:ASN:OD1	1:A:190:VAL:HB	2.17	0.45
1:A:16:LEU:HG	1:A:145:MET:HE1	1.97	0.45
1:B:278:ASP:OD1	1:B:278:ASP:C	2.60	0.45
1:B:154:THR:C	1:B:155:TYR:CD2	2.90	0.45
1:A:80:ARG:HA	1:A:80:ARG:HD2	1.74	0.45
1:B:168:ILE:HD12	1:B:171:CYS:SG	2.56	0.45
1:B:126:LEU:C	1:B:126:LEU:HD23	2.41	0.45
1:B:167:GLY:O	1:B:168:ILE:O	2.35	0.45
1:A:168:ILE:CG2	1:A:169:GLU:H	2.19	0.44
1:A:137:ARG:CG	2:A:419:HOH:O	2.65	0.44
1:B:55:LYS:O	1:B:57:PRO:HD3	2.18	0.44
1:A:179:ASN:HD21	1:A:183:TYR:HE2	1.64	0.44
1:A:56:ASN:C	1:A:58:ASN:N	2.74	0.44
1:B:225:ASN:O	2:B:413:HOH:O	2.20	0.44
1:A:137:ARG:HG2	2:A:419:HOH:O	2.17	0.44
1:B:59:LYS:HG2	1:B:60:ARG:O	2.18	0.44
1:A:285:PHE:HB3	1:B:78:TYR:CE2	2.53	0.43
1:B:168:ILE:HB	1:B:171:CYS:SG	2.59	0.43
1:B:155:TYR:CE1	1:B:182:LYS:HE3	2.42	0.43
1:B:22:ARG:HD3	1:B:38:PHE:HB3	2.00	0.43
1:B:137:ARG:HG2	2:B:403:HOH:O	2.19	0.43
1:A:155:TYR:CE1	1:A:182:LYS:CD	3.02	0.43
1:A:167:GLY:O	1:A:168:ILE:O	2.37	0.43
1:A:230:LYS:HE2	1:A:230:LYS:HB2	1.78	0.43
1:B:80:ARG:HA	1:B:80:ARG:HD2	1.76	0.43
1:A:56:ASN:C	1:A:58:ASN:H	2.27	0.42
1:A:233:ILE:HD12	1:A:233:ILE:N	2.34	0.42
1:B:24:ARG:HA	1:B:37:TYR:O	2.18	0.42
1:B:212:ARG:NH2	1:B:287:GLN:OE1	2.42	0.42
1:B:155:TYR:CE1	1:B:182:LYS:CD	3.02	0.42
1:A:55:LYS:O	1:A:57:PRO:HD3	2.20	0.42
1:A:16:LEU:HD12	1:A:38:PHE:CD2	2.55	0.42
1:A:22:ARG:NH1	2:A:411:HOH:O	2.48	0.42
1:A:22:ARG:HD3	1:A:38:PHE:HB3	2.02	0.42
1:B:16:LEU:CG	1:B:145:MET:HE1	2.50	0.42
1:B:54:LEU:HD12	1:B:54:LEU:HA	1.72	0.42
1:A:99:ILE:HD13	1:A:99:ILE:HA	1.92	0.41
1:B:76:LEU:HD22	1:B:215:THR:CG2	2.50	0.41
1:A:76:LEU:HD22	1:A:215:THR:CG2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:GLU:OE1	1:A:254:GLU:CA	2.69	0.41
1:B:91:LYS:HB2	1:B:91:LYS:HE2	1.64	0.41
1:B:218:GLY:HA2	1:B:219:PRO:HA	1.59	0.40
1:A:154:THR:O	1:A:155:TYR:CD2	2.74	0.40
1:B:95:ARG:O	1:B:96:PRO:C	2.64	0.40
1:A:16:LEU:CG	1:A:145:MET:HE1	2.51	0.40
1:B:183:TYR:O	1:B:292:LEU:HA	2.20	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:95:ARG:NH2	2:A:375:HOH:O[6_444]	1.92	0.28

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	278/340 (82%)	254 (91%)	18 (6%)	6 (2%)	5	10
1	B	279/340 (82%)	256 (92%)	15 (5%)	8 (3%)	3	6
All	All	557/680 (82%)	510 (92%)	33 (6%)	14 (2%)	4	8

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	168	ILE
1	A	171	CYS
1	A	235	ASP
1	B	168	ILE
1	B	171	CYS
1	B	217	THR

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Mol	Chain	Res	Type
1	B	31	ASP
1	A	31	ASP
1	A	170	ASP
1	B	170	ASP
1	A	32	GLY
1	B	32	GLY
1	B	169	GLU
1	B	156	PRO

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	261/308 (85%)	231 (88%)	30 (12%)	5	11
1	B	262/308 (85%)	232 (88%)	30 (12%)	5	11
All	All	523/616 (85%)	463 (88%)	60 (12%)	5	11

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	10	VAL
1	A	22	ARG
1	A	23	LYS
1	A	28	LYS
1	A	29	THR
1	A	31	ASP
1	A	33	LYS
1	A	49	LYS
1	A	51	SER
1	A	54	LEU
1	A	59	LYS
1	A	60	ARG
1	A	110	THR
1	A	113	GLU

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Mol	Chain	Res	Type
1	A	137	ARG
1	A	139	ASN
1	A	154	THR
1	A	156	PRO
1	A	166	VAL
1	A	168	ILE
1	A	172	LEU
1	A	179	ASN
1	A	192	LYS
1	A	234	MET
1	A	243	SER
1	A	261	ASP
1	A	279	GLU
1	A	288	GLN
1	A	297	ASP
1	B	9	SER
1	B	10	VAL
1	B	22	ARG
1	B	23	LYS
1	B	28	LYS
1	B	29	THR
1	B	31	ASP
1	B	33	LYS
1	B	49	LYS
1	B	51	SER
1	B	54	LEU
1	B	59	LYS
1	B	60	ARG
1	B	110	THR
1	B	113	GLU
1	B	137	ARG
1	B	139	ASN
1	B	154	THR
1	B	157	GLU
1	B	158	LEU
1	B	166	VAL
1	B	168	ILE
1	B	172	LEU
1	B	179	ASN
1	B	192	LYS
1	B	234	MET
1	B	243	SER

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Mol	Chain	Res	Type
1	B	261	ASP
1	B	279	GLU
1	B	288	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	ASN
1	A	82	ASN
1	A	84	HIS
1	A	179	ASN
1	A	220	ASN
1	A	273	ASN
1	B	56	ASN
1	B	84	HIS
1	B	179	ASN
1	B	220	ASN
1	B	273	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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 ⓘ

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	283/340 (83%)	0.67	43 (15%) 5 4	21, 46, 93, 127	1 (0%)
1	B	282/340 (82%)	0.55	39 (13%) 6 5	20, 45, 90, 127	1 (0%)
All	All	565/680 (83%)	0.61	82 (14%) 6 4	20, 46, 93, 127	2 (0%)

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	156	PRO	6.3
1	A	236	GLY	5.9
1	A	33	LYS	5.8
1	A	234	MET	5.5
1	B	156	PRO	5.4
1	B	158	LEU	5.3
1	B	234	MET	5.1
1	A	59	LYS	4.9
1	B	59	LYS	4.9
1	B	33	LYS	4.8
1	A	231	TYR	4.6
1	B	9	SER	4.6
1	B	235	ASP	4.5
1	A	170	ASP	4.4
1	A	235	ASP	4.4
1	A	9	SER	4.3
1	A	57	PRO	4.3
1	A	160	SER	4.3
1	A	55	LYS	4.2
1	B	58	ASN	4.1
1	A	168	ILE	4.0
1	A	60	ARG	3.9
1	A	297	ASP	3.9
1	A	138	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	296	GLY	3.7
1	A	23	LYS	3.7
1	B	169	GLU	3.6
1	A	169	GLU	3.6
1	A	237	ALA	3.6
1	A	201	LYS	3.6
1	A	216	GLY	3.5
1	B	57	PRO	3.5
1	B	23	LYS	3.4
1	A	137	ARG	3.4
1	B	20	GLU	3.4
1	B	159	ASN	3.4
1	B	170	ASP	3.3
1	B	55	LYS	3.3
1	A	242	GLU	3.3
1	A	32	GLY	3.2
1	A	200	ILE	3.2
1	A	199	GLU	3.2
1	B	217	THR	3.2
1	B	99	ILE	3.1
1	B	168	ILE	3.1
1	B	171	CYS	3.1
1	B	138	LEU	3.0
1	A	171	CYS	3.0
1	B	197	SER	3.0
1	A	56	ASN	2.9
1	A	198	VAL	2.8
1	A	238	PRO	2.8
1	A	218	GLY	2.7
1	B	247	ARG	2.7
1	B	219	PRO	2.7
1	B	62	GLU	2.7
1	A	58	ASN	2.7
1	B	201	LYS	2.6
1	A	34	LYS	2.6
1	B	137	ARG	2.4
1	A	217	THR	2.4
1	B	160	SER	2.4
1	A	20	GLU	2.4
1	A	219	PRO	2.4
1	A	98	GLU	2.3
1	B	216	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	54	LEU	2.3
1	A	19	ALA	2.2
1	B	139	ASN	2.2
1	B	32	GLY	2.2
1	A	172	LEU	2.2
1	A	62	GLU	2.1
1	B	155	TYR	2.1
1	A	54	LEU	2.1
1	B	100	THR	2.1
1	B	56	ASN	2.1
1	B	28	LYS	2.1
1	B	36	LYS	2.1
1	A	61	LEU	2.1
1	B	200	ILE	2.1
1	A	53	SER	2.0
1	B	78	TYR	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.