



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:12 PM UTC

PDB ID : 3LHA / pdb_00003lha
Title : Crystal structure of mouse VPS26B(R240S/G241A/E242S) in spacegroup P41 21 2
Authors : Collins, B.; Shaw, D.; Norwood, S.
Deposited on : 2010-01-21
Resolution : 2.80 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

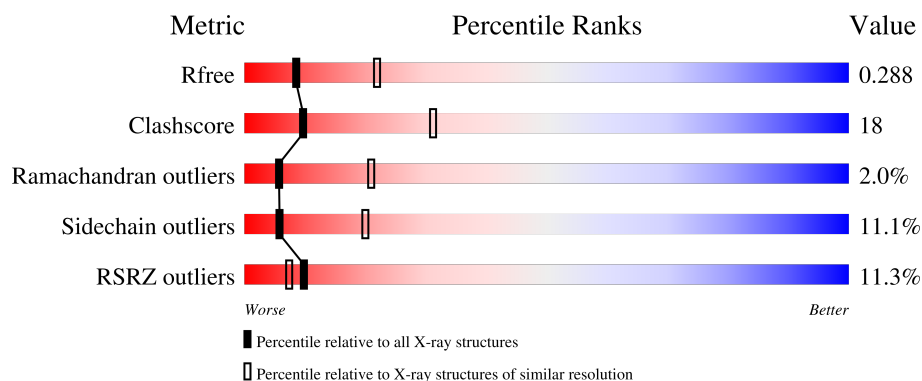
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	11% 55% 24% • • 17%
1	B	340	8% 53% 24% 5% • 17%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4855 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
			2351	1506	396	443	6			
1	B	282	Total	C	N	O	S	0	1	0
			2352	1507	396	443	6			

There are 26 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
A	242	SER	GLU	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

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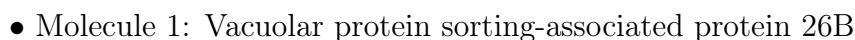
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Chain	Residue	Modelled	Actual	Comment	Reference
B	242	SER	GLU	engineered mutation	UNP Q8C0E2

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	93	Total O 93 93	0	0
2	B	59	Total O 59 59	0	0

- 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	118.14Å 118.14Å 193.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.85 – 2.80 40.85 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (40.85-2.80) 99.9 (40.85-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.92 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.252 , 0.284 0.254 , 0.288	Depositor DCC
R_{free} test set	1734 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	52.3	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	4855	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 2.62% 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 [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	0.60	0/2401	0.90	3/3229 (0.1%)
1	B	0.60	0/2402	0.95	5/3231 (0.2%)
All	All	0.60	0/4803	0.92	8/6460 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	155	TYR	CA-C-N	7.45	129.15	119.84
1	B	155	TYR	C-N-CA	7.45	129.15	119.84
1	B	60	ARG	CB-CG-CD	5.95	124.99	111.30
1	B	234	MET	N-CA-C	5.91	118.52	110.55
1	A	234	MET	N-CA-C	5.84	118.44	110.55
1	A	60	ARG	CB-CG-CD	5.54	124.04	111.30
1	A	179	ASN	N-CA-C	5.31	117.15	111.36
1	B	179	ASN	N-CA-C	5.20	117.03	111.36

There are no chirality outliers.

There are no planarity outliers.

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	2351	0	2345	81	0
1	B	2352	0	2350	88	2
2	A	93	0	0	8	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	59	0	0	3	0
All	All	4855	0	4695	166	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 18.

All (166) 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:179:ASN:ND2	1:B:190:VAL:HB	1.72	1.05
1:B:179:ASN:HD22	1:B:190:VAL:HB	1.22	1.04
1:A:259:MET:HE3	2:A:367:HOH:O	1.57	1.02
1:A:179:ASN:ND2	1:A:190:VAL:HB	1.74	1.02
1:A:179:ASN:HD22	1:A:190:VAL:HB	1.25	0.99
1:B:259:MET:HE3	2:B:358:HOH:O	1.67	0.94
1:B:164:MET:HE3	1:B:274:LEU:HB2	1.50	0.93
1:A:164:MET:HE3	1:A:274:LEU:HB2	1.54	0.87
1:B:171:CYS:HA	1:B:199:GLU:HB2	1.58	0.86
1:A:171:CYS:HA	1:A:199:GLU:HB2	1.59	0.84
1:A:119:TYR:HB2	1:A:259:MET:HE1	1.58	0.82
1:A:155:TYR:CD1	1:A:182:LYS:HD3	2.17	0.80
1:B:155:TYR:CD1	1:B:182:LYS:HD3	2.18	0.79
1:B:119:TYR:HB2	1:B:259:MET:HE1	1.66	0.76
1:A:155:TYR:HB3	1:A:156:PRO:HD2	1.68	0.76
1:B:155:TYR:HB3	1:B:156:PRO:HD2	1.66	0.76
1:B:217:THR:HG21	1:B:264:LYS:O	1.88	0.74
1:A:154:THR:C	1:A:155:TYR:HD2	1.96	0.73
1:A:110:THR:HA	2:A:409:HOH:O	1.89	0.71
1:A:155:TYR:CE1	1:A:182:LYS:HD3	2.26	0.70
1:B:155:TYR:CE1	1:B:182:LYS:HD3	2.26	0.70
1:B:154:THR:C	1:B:155:TYR:HD2	1.98	0.70
1:B:156:PRO:HB2	1:B:158:LEU:H	1.57	0.69
1:A:168:ILE:HD12	1:A:171:CYS:SG	2.33	0.69
1:A:122:GLN:HG2	1:A:224[A]:GLU:HG3	1.76	0.68
1:B:164:MET:HE3	1:B:274:LEU:CB	2.24	0.67
1:B:122:GLN:HG2	1:B:224[A]:GLU:HG3	1.76	0.67
1:B:217:THR:H	1:B:221:VAL:HG22	1.59	0.66
1:A:243:SER:O	1:A:244:ILE:HG13	1.96	0.65
1:B:243:SER:O	1:B:244:ILE:HG13	1.97	0.64
1:A:164:MET:HE3	1:A:274:LEU:CB	2.27	0.64
1:A:60:ARG:NH1	1:A:98:GLU:OE2	2.32	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:PRO:HB2	1:B:157:GLU:HA	1.79	0.63
1:A:210:ILE:HD12	1:A:273:ASN:HB2	1.81	0.63
1:B:60:ARG:NH1	1:B:98:GLU:OE2	2.31	0.63
1:B:182:LYS:HG2	1:B:291:VAL:HB	1.80	0.62
1:B:169:GLU:O	1:B:169:GLU:HG3	1.99	0.62
1:B:210:ILE:HD12	1:B:273:ASN:HB2	1.81	0.62
1:B:168:ILE:HD12	1:B:171:CYS:SG	2.40	0.62
1:A:280:GLU:O	1:A:281:GLU:HB2	1.99	0.61
1:B:155:TYR:HD1	1:B:182:LYS:HD3	1.64	0.61
1:A:119:TYR:CB	1:A:259:MET:HE1	2.28	0.61
1:A:154:THR:O	1:A:155:TYR:HD2	1.82	0.61
1:A:155:TYR:HD1	1:A:182:LYS:HD3	1.63	0.61
1:B:119:TYR:CB	1:B:259:MET:HE1	2.29	0.61
1:A:16:LEU:HG	1:A:145:MET:HE1	1.85	0.59
1:A:182:LYS:HG2	1:A:291:VAL:HB	1.83	0.59
1:B:60:ARG:CG	1:B:60:ARG:HH11	2.16	0.59
1:B:280:GLU:O	1:B:281:GLU:HB2	2.02	0.59
1:A:169:GLU:O	1:A:169:GLU:HG3	2.04	0.57
1:B:122:GLN:CG	1:B:224[A]:GLU:HG3	2.35	0.57
1:B:16:LEU:HG	1:B:145:MET:HE1	1.86	0.57
1:A:111:HIS:N	2:A:409:HOH:O	2.14	0.56
1:A:91:LYS:HD3	2:A:365:HOH:O	2.05	0.56
1:B:156:PRO:CB	1:B:158:LEU:H	2.18	0.56
1:B:154:THR:O	1:B:155:TYR:HD2	1.89	0.55
1:B:24:ARG:HA	1:B:37:TYR:O	2.06	0.55
1:B:217:THR:CG2	1:B:264:LYS:O	2.55	0.55
1:B:155:TYR:CB	1:B:156:PRO:HD2	2.34	0.55
1:A:122:GLN:CG	1:A:224[A]:GLU:HG3	2.35	0.54
1:B:22:ARG:HD2	2:B:365:HOH:O	2.08	0.54
1:B:156:PRO:HB2	1:B:157:GLU:CA	2.37	0.54
1:B:166:VAL:HG12	1:B:174:ILE:HB	1.90	0.54
1:A:137:ARG:HG3	2:A:402:HOH:O	2.07	0.53
1:B:225:ASN:O	2:B:385:HOH:O	2.19	0.53
1:B:60:ARG:HH11	1:B:60:ARG:HG2	1.74	0.52
1:A:20:GLU:CD	1:A:20:GLU:H	2.18	0.52
1:B:69:GLU:HG2	1:B:90:VAL:HG22	1.90	0.52
1:B:171:CYS:CA	1:B:199:GLU:HB2	2.37	0.52
1:B:179:ASN:ND2	1:B:190:VAL:CB	2.61	0.52
1:B:156:PRO:HB2	1:B:158:LEU:N	2.25	0.52
1:B:208:ASP:HB2	1:B:273:ASN:HB3	1.93	0.51
1:B:231:TYR:CD2	1:B:231:TYR:C	2.88	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:GLU:C	1:A:233:ILE:HD12	2.34	0.51
1:A:166:VAL:HG12	1:A:174:ILE:HB	1.92	0.51
1:A:155:TYR:CB	1:A:156:PRO:HD2	2.34	0.50
1:A:285:PHE:HB3	1:B:78:TYR:CE2	2.46	0.50
1:B:20:GLU:H	1:B:20:GLU:CD	2.17	0.50
1:B:254:GLU:HA	1:B:254:GLU:OE1	2.12	0.50
1:A:254:GLU:HA	1:A:254:GLU:OE1	2.11	0.50
1:A:64:GLN:HA	1:A:96:PRO:HB3	1.94	0.49
1:B:171:CYS:HB2	1:B:200:ILE:HG13	1.95	0.49
1:B:278:ASP:OD1	1:B:278:ASP:C	2.56	0.49
1:A:170:ASP:O	1:A:171:CYS:CB	2.61	0.48
1:B:164:MET:CE	1:B:274:LEU:HB2	2.35	0.48
1:A:22:ARG:NH2	1:A:44:GLU:OE1	2.41	0.48
1:B:124:VAL:HG22	1:B:125:LYS:N	2.29	0.48
1:A:110:THR:CA	2:A:409:HOH:O	2.53	0.47
1:A:122:GLN:HG2	1:A:224[B]:GLU:CD	2.40	0.47
1:B:76:LEU:HD23	1:B:124:VAL:HG23	1.97	0.47
1:A:124:VAL:HG22	1:A:125:LYS:N	2.29	0.47
1:B:234:MET:HG3	1:B:235:ASP:OD2	2.14	0.47
1:A:208:ASP:HB2	1:A:273:ASN:HB3	1.96	0.47
1:B:137:ARG:H	1:B:137:ARG:HG3	1.34	0.47
1:B:170:ASP:O	1:B:171:CYS:CB	2.62	0.47
1:B:170:ASP:O	1:B:171:CYS:HB3	2.15	0.47
1:A:123:ASN:HD22	1:B:285:PHE:HE2	1.62	0.46
1:B:180:LYS:HB2	1:B:183:TYR:CZ	2.50	0.46
1:A:193:ILE:HG21	1:A:205:MET:HE1	1.97	0.46
1:B:122:GLN:HG2	1:B:224[B]:GLU:CD	2.40	0.46
1:A:137:ARG:HG3	1:A:137:ARG:H	1.37	0.46
1:A:180:LYS:HB2	1:A:183:TYR:CZ	2.51	0.46
1:A:171:CYS:CA	1:A:199:GLU:HB2	2.39	0.46
1:A:206:GLU:OE2	1:A:230:LYS:HD3	2.16	0.46
1:A:69:GLU:HG2	1:A:90:VAL:HG22	1.98	0.46
1:A:170:ASP:O	1:A:171:CYS:HB3	2.15	0.46
1:A:171:CYS:HB2	1:A:200:ILE:HG13	1.97	0.46
1:B:230:LYS:HB2	1:B:230:LYS:HE2	1.69	0.46
1:A:233:ILE:HD12	1:A:233:ILE:N	2.31	0.45
1:B:154:THR:C	1:B:155:TYR:CD2	2.88	0.45
1:B:228:ILE:HG21	1:B:255:LEU:HD21	1.97	0.45
1:A:169:GLU:O	1:A:170:ASP:OD1	2.34	0.45
1:B:206:GLU:OE2	1:B:230:LYS:HD3	2.15	0.45
1:A:228:ILE:HG21	1:A:255:LEU:HD21	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:TYR:CD2	1:A:231:TYR:C	2.95	0.45
1:A:234:MET:HG3	1:A:235:ASP:OD2	2.16	0.45
1:B:168:ILE:HB	1:B:171:CYS:SG	2.56	0.45
1:A:24:ARG:HA	1:A:37:TYR:O	2.16	0.45
1:A:126:LEU:C	1:A:126:LEU:HD23	2.42	0.45
1:B:232:GLU:C	1:B:233:ILE:HD12	2.41	0.45
1:B:49:LYS:HA	1:B:105:PHE:O	2.16	0.45
1:A:59:LYS:HG2	1:A:60:ARG:O	2.18	0.44
1:A:122:GLN:HG2	1:A:224[B]:GLU:CG	2.47	0.44
1:A:95:ARG:O	1:A:96:PRO:C	2.61	0.44
1:A:154:THR:C	1:A:155:TYR:CD2	2.86	0.44
1:B:122:GLN:HG2	1:B:224[B]:GLU:CG	2.48	0.44
1:A:183:TYR:O	1:A:292:LEU:HA	2.17	0.44
1:A:154:THR:O	1:A:155:TYR:CD2	2.69	0.44
1:A:285:PHE:HE2	1:B:123:ASN:HD22	1.66	0.44
1:A:80:ARG:HD2	1:A:80:ARG:HA	1.66	0.43
1:A:278:ASP:OD1	1:A:278:ASP:C	2.60	0.43
1:A:49:LYS:NZ	1:A:49:LYS:HB3	2.33	0.43
1:B:59:LYS:HG2	1:B:60:ARG:O	2.17	0.43
1:B:64:GLN:HA	1:B:96:PRO:HB3	2.00	0.43
1:A:168:ILE:HB	1:A:171:CYS:SG	2.58	0.43
1:B:80:ARG:HA	1:B:80:ARG:HD2	1.68	0.43
1:B:218:GLY:HA2	1:B:219:PRO:HA	1.68	0.43
1:A:95:ARG:NH2	2:A:379:HOH:O	2.51	0.43
1:A:247:ARG:NH1	2:A:377:HOH:O	2.42	0.43
1:A:49:LYS:HA	1:A:105:PHE:O	2.18	0.42
1:B:60:ARG:HG2	1:B:98:GLU:OE2	2.17	0.42
1:B:119:TYR:CZ	1:B:268:VAL:HG21	2.54	0.42
1:A:131:ARG:HD3	1:A:144:GLU:OE2	2.19	0.42
1:B:124:VAL:HG22	1:B:125:LYS:H	1.84	0.42
1:B:233:ILE:HD12	1:B:233:ILE:N	2.35	0.42
1:A:69:GLU:O	1:A:130:LEU:HA	2.19	0.42
1:B:95:ARG:O	1:B:96:PRO:C	2.62	0.42
1:A:163:LYS:HD3	1:A:175:GLU:OE2	2.20	0.42
1:B:108:GLU:OE1	1:B:110:THR:HG23	2.20	0.42
1:B:216:GLY:C	1:B:217:THR:HG23	2.44	0.42
1:B:155:TYR:HB3	1:B:156:PRO:CD	2.43	0.42
1:B:193:ILE:HG21	1:B:205:MET:HE1	2.00	0.42
1:B:60:ARG:HH11	1:B:60:ARG:CB	2.32	0.42
1:A:230:LYS:HE2	1:A:230:LYS:HB2	1.69	0.41
1:A:171:CYS:HA	1:A:199:GLU:CB	2.42	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:16:LEU:CG	1:B:145:MET:HE1	2.50	0.41
1:B:35:GLU:OE2	1:B:35:GLU:HA	2.19	0.41
1:A:16:LEU:CG	1:A:145:MET:HE1	2.49	0.41
1:A:35:GLU:OE2	1:A:35:GLU:HA	2.20	0.41
1:A:119:TYR:CZ	1:A:268:VAL:HG21	2.55	0.41
1:A:272:LEU:O	1:A:287:GLN:HA	2.21	0.41
1:B:271:TYR:CE1	1:B:289:GLU:HB2	2.56	0.41
1:B:22:ARG:NH2	1:B:44:GLU:OE1	2.45	0.41
1:A:210:ILE:HD12	1:A:273:ASN:CB	2.49	0.41
1:B:210:ILE:HD12	1:B:273:ASN:CB	2.50	0.40
1:B:169:GLU:O	1:B:170:ASP:OD1	2.39	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)
1:B:95:ARG:NH2	2:A:366:HOH:O[6_444]	2.11	0.09
1:B:196:LEU:O	1:B:196:LEU:O[7_555]	2.14	0.06

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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%)	256 (92%)	17 (6%)	5 (2%)	6	23
1	B	279/340 (82%)	254 (91%)	19 (7%)	6 (2%)	5	19
All	All	557/680 (82%)	510 (92%)	36 (6%)	11 (2%)	6	21

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	168	ILE

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Mol	Chain	Res	Type
1	A	171	CYS
1	A	235	ASP
1	B	168	ILE
1	B	171	CYS
1	B	217	THR
1	A	138	LEU
1	B	138	LEU
1	B	156	PRO
1	B	57	PRO
1	A	32	GLY

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	261/308 (85%)	233 (89%)	28 (11%)	6	21
1	B	262/308 (85%)	232 (88%)	30 (12%)	5	18
All	All	523/616 (85%)	465 (89%)	58 (11%)	6	20

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	10	VAL
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	54	LEU
1	A	55	LYS
1	A	59	LYS
1	A	60	ARG
1	A	110	THR

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Mol	Chain	Res	Type
1	A	113	GLU
1	A	137	ARG
1	A	154	THR
1	A	156	PRO
1	A	166	VAL
1	A	168	ILE
1	A	169	GLU
1	A	170	ASP
1	A	172	LEU
1	A	230	LYS
1	A	234	MET
1	A	243	SER
1	A	261	ASP
1	A	286	LYS
1	A	288	GLN
1	B	9	SER
1	B	10	VAL
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	54	LEU
1	B	55	LYS
1	B	59	LYS
1	B	60	ARG
1	B	110	THR
1	B	113	GLU
1	B	137	ARG
1	B	154	THR
1	B	157	GLU
1	B	159	ASN
1	B	166	VAL
1	B	168	ILE
1	B	169	GLU
1	B	170	ASP
1	B	172	LEU
1	B	181	SER
1	B	230	LYS
1	B	234	MET
1	B	243	SER

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Mol	Chain	Res	Type
1	B	261	ASP
1	B	286	LYS
1	B	288	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	HIS
1	A	84	HIS
1	A	123	ASN
1	A	273	ASN
1	B	83	HIS
1	B	84	HIS
1	B	123	ASN
1	B	179	ASN
1	B	273	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 ⓘ

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%)	0.50	37 (13%) 7 6	21, 45, 93, 133	1 (0%)
1	B	282/340 (82%)	0.32	27 (9%) 13 10	20, 45, 92, 132	1 (0%)
All	All	565/680 (83%)	0.41	64 (11%) 10 7	20, 45, 92, 133	2 (0%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	156	PRO	5.4
1	A	160	SER	4.9
1	B	158	LEU	4.8
1	A	234	MET	4.8
1	A	59	LYS	4.6
1	B	156	PRO	4.5
1	A	9	SER	4.4
1	B	59	LYS	4.4
1	A	236	GLY	4.4
1	A	169	GLU	4.3
1	A	138	LEU	4.1
1	A	33	LYS	4.0
1	A	231	TYR	4.0
1	B	234	MET	3.9
1	A	168	ILE	3.9
1	A	235	ASP	3.9
1	B	9	SER	3.8
1	B	58	ASN	3.7
1	A	170	ASP	3.7
1	B	62	GLU	3.7
1	A	55	LYS	3.6
1	A	242	SER	3.6
1	B	235	ASP	3.6
1	B	160	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	169	GLU	3.4
1	A	297	ASP	3.4
1	A	167	GLY	3.3
1	A	216	GLY	3.2
1	B	138	LEU	3.1
1	A	171	CYS	3.1
1	B	20	GLU	3.0
1	A	57	PRO	3.0
1	B	33	LYS	2.9
1	B	296	GLY	2.9
1	A	32	GLY	2.8
1	B	23	LYS	2.8
1	A	60	ARG	2.8
1	A	172	LEU	2.7
1	A	237	ALA	2.6
1	B	100	THR	2.5
1	B	57	PRO	2.5
1	A	137	ARG	2.5
1	B	171	CYS	2.5
1	A	20	GLU	2.5
1	A	56	ASN	2.4
1	A	23	LYS	2.4
1	A	35	GLU	2.4
1	B	197	SER	2.4
1	A	201	LYS	2.3
1	B	137	ARG	2.3
1	A	198	VAL	2.3
1	A	249	PHE	2.3
1	A	199	GLU	2.3
1	B	55	LYS	2.3
1	A	238	PRO	2.3
1	A	218	GLY	2.3
1	B	201	LYS	2.2
1	A	200	ILE	2.2
1	B	168	ILE	2.2
1	A	34	LYS	2.2
1	B	99	ILE	2.1
1	B	32	GLY	2.1
1	B	159	ASN	2.0
1	B	34	LYS	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.