



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 05:28 PM UTC

PDB ID : 5H61 / pdb_00005h61
Title : Structure of Transferase mutant-C23S,C199S
Authors : Park, J.B.; Yoo, Y.; Kim, J.
Deposited on : 2016-11-10
Resolution : 1.86 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

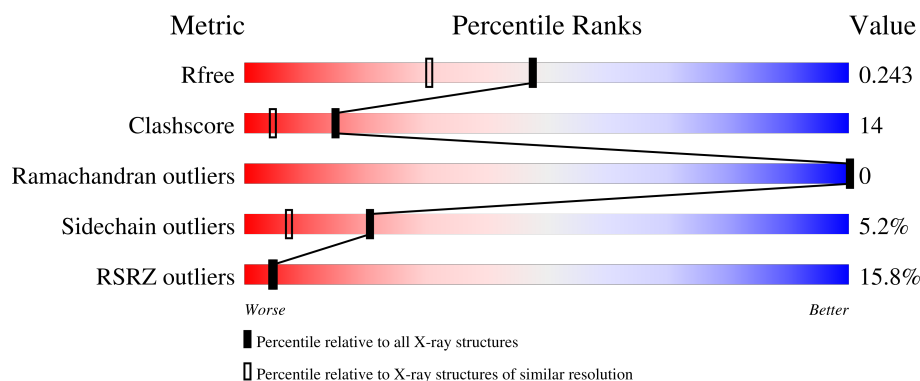
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.86 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)

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	348	12% 66% 17% • 15%
1	B	348	15% 64% 18% • 15%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5030 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 Transferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	295	Total	C	N	O	S	0	0	0
			2382	1525	406	444	7			
1	B	295	Total	C	N	O	S	0	0	0
			2383	1527	406	443	7			

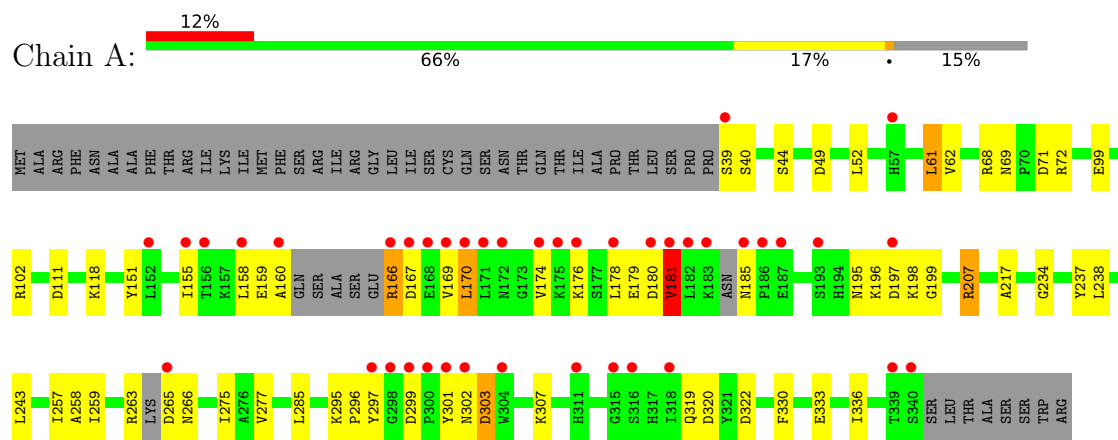
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	141	Total	O	0	0
			141	141		
2	B	124	Total	O	0	0
			124	124		

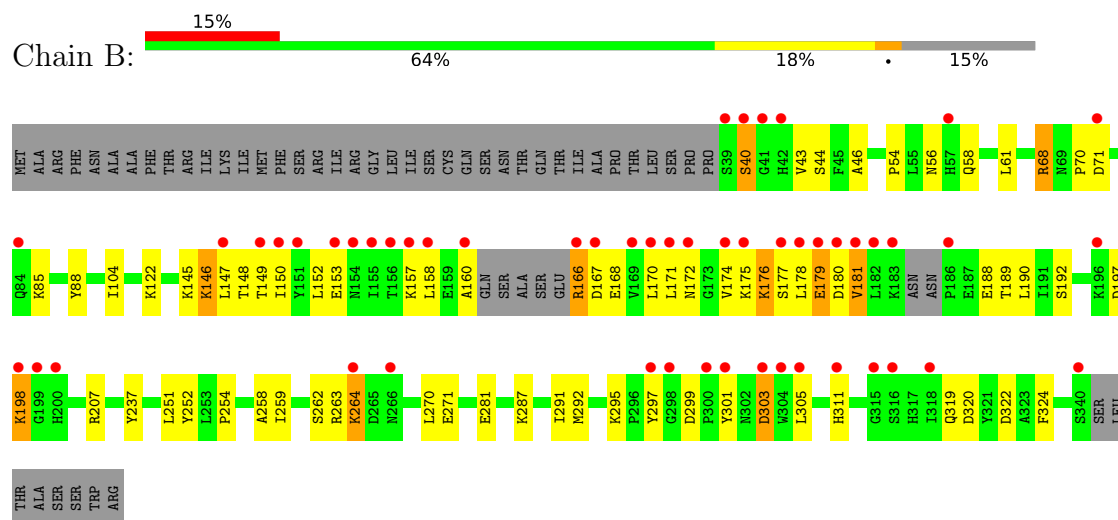
3 Residue-property plots [i](#)

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: Transferase



• Molecule 1: Transferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.61Å 145.58Å 55.48Å 90.00° 105.12° 90.00°	Depositor
Resolution (Å)	72.79 – 1.86 72.79 – 1.86	Depositor EDS
% Data completeness (in resolution range)	97.8 (72.79-1.86) 97.8 (72.79-1.86)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 1.86Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.198 , 0.233 0.206 , 0.243	Depositor DCC
R_{free} test set	2856 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	26.8	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5030	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.56% 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	1.40	8/2438 (0.3%)	1.08	7/3294 (0.2%)
1	B	1.36	2/2440 (0.1%)	1.08	2/3296 (0.1%)
All	All	1.38	10/4878 (0.2%)	1.08	9/6590 (0.1%)

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	0	1
1	B	0	1
All	All	0	2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	118	LYS	C-O	-6.26	1.16	1.24
1	A	234	GLY	C-O	-6.14	1.18	1.24
1	B	122	LYS	C-O	-5.87	1.17	1.24
1	A	207	ARG	CA-C	-5.74	1.45	1.52
1	A	62	VAL	C-O	-5.53	1.18	1.24
1	A	99	GLU	CA-C	-5.45	1.47	1.53
1	A	217	ALA	N-CA	-5.42	1.39	1.46
1	A	62	VAL	N-CA	-5.14	1.40	1.46
1	A	238	LEU	CA-C	-5.01	1.46	1.52
1	B	281	GLU	CA-C	-5.00	1.46	1.53

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	185	ASN	CA-C-N	-6.22	113.40	119.87
1	A	185	ASN	C-N-CA	-6.22	113.40	119.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	196	LYS	N-CA-C	5.93	118.99	111.69
1	B	299	ASP	CA-C-N	-5.85	112.48	118.85
1	B	299	ASP	C-N-CA	-5.85	112.48	118.85
1	A	299	ASP	CA-C-N	-5.75	112.58	118.85
1	A	299	ASP	C-N-CA	-5.75	112.58	118.85
1	A	181	VAL	CB-CA-C	-5.68	104.47	112.14
1	A	40	SER	N-CA-C	-5.58	106.27	112.57

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	39	SER	Peptide
1	B	40	SER	Peptide

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	2382	0	2330	58	2
1	B	2383	0	2339	85	2
2	A	141	0	0	1	0
2	B	124	0	0	2	0
All	All	5030	0	4669	129	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 14.

All (129) 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:176:LYS:CB	1:B:301:TYR:HD1	1.52	1.22
1:A:176:LYS:HB3	1:B:301:TYR:HD1	1.04	1.15
1:A:301:TYR:HB3	1:B:176:LYS:HD3	1.28	1.15
1:A:167:ASP:N	1:A:170:LEU:HD21	1.63	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:ASP:H	1:A:170:LEU:CD2	1.61	1.12
1:B:176:LYS:O	1:B:179:GLU:HG3	1.49	1.10
1:A:176:LYS:HB3	1:B:301:TYR:CD1	1.93	1.04
1:A:176:LYS:CB	1:B:301:TYR:CD1	2.40	1.04
1:A:167:ASP:H	1:A:170:LEU:HD21	0.87	1.02
1:B:167:ASP:OD1	1:B:168:GLU:N	1.97	0.96
1:A:166:ARG:N	1:A:170:LEU:CD2	2.30	0.94
1:A:176:LYS:HB2	1:B:301:TYR:CD1	2.03	0.93
1:A:320:ASP:OD2	1:A:322:ASP:HB2	1.70	0.92
1:A:167:ASP:N	1:A:170:LEU:CD2	2.28	0.92
1:A:158:LEU:HB3	1:A:174:VAL:HG22	1.53	0.91
1:A:166:ARG:N	1:A:170:LEU:HD21	1.89	0.88
1:B:176:LYS:O	1:B:179:GLU:CG	2.23	0.85
1:A:301:TYR:HB3	1:B:176:LYS:CD	2.07	0.85
1:A:68:ARG:HD2	1:A:198:LYS:O	1.78	0.84
1:A:265:ASP:OD1	1:A:266:ASN:N	2.09	0.84
1:B:176:LYS:HA	1:B:179:GLU:HG2	1.59	0.83
1:B:160:ALA:O	1:B:166:ARG:NE	2.10	0.83
1:A:166:ARG:N	1:A:170:LEU:HD22	1.94	0.81
1:A:167:ASP:O	1:A:170:LEU:HG	1.85	0.76
1:B:303:ASP:OD1	1:B:305:LEU:N	2.17	0.76
1:B:148:THR:O	1:B:152:LEU:HD12	1.89	0.73
1:A:170:LEU:O	1:A:174:VAL:HG23	1.88	0.72
1:B:178:LEU:HA	1:B:181:VAL:HG23	1.71	0.72
1:B:179:GLU:HG3	1:B:180:ASP:N	2.05	0.72
1:A:303:ASP:HB2	2:A:451:HOH:O	1.88	0.72
1:A:166:ARG:N	1:A:166:ARG:HD3	2.04	0.71
1:A:302:ASN:O	1:A:307:LYS:HE3	1.92	0.70
1:A:68:ARG:CD	1:A:198:LYS:O	2.42	0.68
1:B:168:GLU:O	1:B:172:ASN:CG	2.38	0.66
1:B:177:SER:O	1:B:181:VAL:HG23	1.96	0.65
1:B:153:GLU:O	1:B:157:LYS:HG2	1.97	0.65
1:B:146:LYS:CD	1:B:146:LYS:H	2.10	0.64
1:B:179:GLU:HG3	1:B:180:ASP:H	1.62	0.64
1:A:301:TYR:CB	1:B:176:LYS:HD3	2.19	0.63
1:B:167:ASP:N	1:B:170:LEU:HD13	2.14	0.63
1:A:180:ASP:OD2	1:B:301:TYR:OH	2.14	0.62
1:B:146:LYS:H	1:B:146:LYS:HE2	1.65	0.61
1:B:46:ALA:H	1:B:319:GLN:HE21	1.49	0.60
1:B:197:ASP:O	1:B:198:LYS:HE2	2.02	0.60
1:B:291:ILE:HG23	1:B:295:LYS:HD2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:145:LYS:HE2	1:B:150:ILE:HG12	1.85	0.58
1:B:170:LEU:O	1:B:174:VAL:HG23	2.04	0.58
1:B:178:LEU:CA	1:B:181:VAL:HG23	2.34	0.56
1:B:68:ARG:CG	1:B:198:LYS:HB3	2.35	0.55
1:B:262:SER:OG	1:B:271:GLU:OE1	2.18	0.55
1:B:146:LYS:H	1:B:146:LYS:CE	2.18	0.55
1:A:258:ALA:O	1:A:259:ILE:HD13	2.07	0.55
1:B:190:LEU:HD13	1:B:292:MET:HE3	1.87	0.55
1:B:178:LEU:HA	1:B:181:VAL:CG2	2.36	0.54
1:A:158:LEU:HB3	1:A:174:VAL:CG2	2.33	0.54
1:A:167:ASP:N	1:A:170:LEU:HD23	2.18	0.54
1:B:263:ARG:HG2	1:B:263:ARG:HH11	1.73	0.54
1:B:158:LEU:O	1:B:160:ALA:N	2.41	0.53
1:A:259:ILE:CD1	1:A:275:ILE:HG23	2.39	0.52
1:B:320:ASP:OD2	1:B:322:ASP:HB3	2.09	0.52
1:B:258:ALA:O	1:B:259:ILE:HD13	2.10	0.52
1:A:195:ASN:O	1:A:198:LYS:NZ	2.42	0.52
1:A:257:ILE:HD12	1:A:277:VAL:HG12	1.93	0.51
1:A:302:ASN:O	1:A:307:LYS:CE	2.58	0.51
1:A:301:TYR:CE2	1:B:172:ASN:HB2	2.45	0.51
1:B:168:GLU:O	1:B:172:ASN:OD1	2.29	0.51
1:A:167:ASP:O	1:A:170:LEU:CG	2.57	0.50
1:B:147:LEU:CD1	1:B:188:GLU:HG2	2.41	0.50
1:B:146:LYS:H	1:B:146:LYS:HD3	1.75	0.50
1:B:146:LYS:HE3	1:B:149:THR:OG1	2.11	0.49
1:B:167:ASP:H	1:B:170:LEU:HD13	1.76	0.49
1:A:69:ASN:O	1:A:72:ARG:HG2	2.11	0.49
1:B:166:ARG:HA	1:B:170:LEU:HD13	1.93	0.49
1:B:54:PRO:HA	1:B:251:LEU:O	2.12	0.49
1:A:296:PRO:HG2	1:A:297:TYR:CD2	2.48	0.49
1:B:158:LEU:HD23	1:B:170:LEU:HD23	1.95	0.49
1:A:180:ASP:OD2	1:B:301:TYR:CE1	2.68	0.47
1:A:158:LEU:CB	1:A:174:VAL:HG22	2.33	0.47
1:A:303:ASP:OD1	1:A:303:ASP:N	2.37	0.47
1:B:167:ASP:N	1:B:170:LEU:CD1	2.78	0.47
1:A:197:ASP:OD1	1:A:199:GLY:N	2.37	0.47
1:B:177:SER:O	1:B:181:VAL:N	2.38	0.47
1:A:166:ARG:C	1:A:170:LEU:HD21	2.35	0.46
1:B:54:PRO:HB3	1:B:252:TYR:CE1	2.51	0.46
1:A:151:TYR:HB3	1:A:181:VAL:HG22	1.97	0.46
1:A:176:LYS:HE2	1:B:301:TYR:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:LYS:HG3	1:B:301:TYR:HB3	1.97	0.46
1:B:167:ASP:H	1:B:170:LEU:CD1	2.29	0.46
1:B:179:GLU:CG	1:B:180:ASP:N	2.77	0.46
1:B:158:LEU:O	1:B:166:ARG:HD2	2.16	0.46
1:B:291:ILE:CG2	1:B:295:LYS:HD2	2.45	0.45
1:A:61:LEU:HD13	1:A:102:ARG:CD	2.46	0.45
1:B:68:ARG:C	1:B:70:PRO:HD3	2.42	0.45
1:B:157:LYS:N	1:B:157:LYS:HD3	2.32	0.45
1:B:175:LYS:HD2	1:B:175:LYS:HA	1.73	0.45
1:B:270:LEU:HD11	1:B:324:PHE:HE2	1.81	0.45
1:A:44:SER:HB3	1:A:49:ASP:OD1	2.17	0.45
1:B:147:LEU:HD11	1:B:188:GLU:HG2	1.98	0.44
1:A:301:TYR:HE1	1:B:297:TYR:OH	2.00	0.44
1:B:189:THR:H	1:B:192:SER:HG	1.62	0.44
1:A:243:LEU:HD23	1:A:336:ILE:HD12	1.99	0.44
1:B:158:LEU:CD2	1:B:170:LEU:HD23	2.48	0.44
1:B:287:LYS:HE2	2:B:438:HOH:O	2.18	0.43
1:B:56:ASN:OD1	1:B:58:GLN:HB2	2.18	0.43
1:A:167:ASP:OD2	1:A:169:VAL:HG23	2.18	0.43
1:A:68:ARG:CG	1:A:198:LYS:O	2.67	0.43
1:B:68:ARG:HG3	1:B:198:LYS:HB3	2.00	0.43
1:B:166:ARG:CA	1:B:170:LEU:HD13	2.49	0.42
1:B:166:ARG:C	1:B:170:LEU:HD13	2.44	0.42
1:B:168:GLU:O	1:B:172:ASN:ND2	2.52	0.42
1:A:61:LEU:HD13	1:A:102:ARG:HD3	2.02	0.42
1:B:148:THR:HG23	2:B:425:HOH:O	2.19	0.42
1:A:207:ARG:HG3	1:A:237:TYR:CZ	2.55	0.42
1:B:146:LYS:CD	1:B:146:LYS:N	2.79	0.42
1:B:207:ARG:HG3	1:B:237:TYR:CE2	2.55	0.41
1:B:263:ARG:HG2	1:B:263:ARG:NH1	2.35	0.41
1:B:287:LYS:HE3	1:B:311:HIS:ND1	2.36	0.41
1:B:176:LYS:HE3	1:B:176:LYS:HB3	1.64	0.41
1:B:43:VAL:HG21	1:B:254:PRO:HB3	2.02	0.41
1:B:146:LYS:HD3	1:B:146:LYS:N	2.36	0.41
1:B:188:GLU:HG3	1:B:189:THR:N	2.35	0.41
1:A:330:PHE:C	1:A:330:PHE:CD1	2.98	0.41
1:B:88:TYR:N	1:B:88:TYR:CD1	2.88	0.40
1:A:155:ILE:CD1	1:A:178:LEU:HD23	2.51	0.40
1:A:159:GLU:O	1:A:160:ALA:HB2	2.22	0.40
1:A:166:ARG:CA	1:A:170:LEU:HD21	2.50	0.40
1:B:68:ARG:HG3	1:B:198:LYS:O	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:ARG:O	1:B:264:LYS:HG3	2.21	0.40
1:A:295:LYS:HA	1:A:296:PRO:HD3	2.01	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:A:333:GLU:OE2	1:B:40:SER:OG[2_545]	2.09	0.11
1:A:333:GLU:OE2	1:B:40:SER:CB[2_545]	2.18	0.02

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	287/348 (82%)	284 (99%)	3 (1%)	0	100	100
1	B	289/348 (83%)	285 (99%)	4 (1%)	0	100	100
All	All	576/696 (83%)	569 (99%)	7 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	261/307 (85%)	249 (95%)	12 (5%)	24	9

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	261/307 (85%)	246 (94%)	15 (6%)	18	6
All	All	522/614 (85%)	495 (95%)	27 (5%)	21	7

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	LEU
1	A	61	LEU
1	A	71	ASP
1	A	111	ASP
1	A	166	ARG
1	A	170	LEU
1	A	179	GLU
1	A	181	VAL
1	A	263	ARG
1	A	285	LEU
1	A	303	ASP
1	A	319	GLN
1	B	44	SER
1	B	61	LEU
1	B	68	ARG
1	B	71	ASP
1	B	85	LYS
1	B	104	ILE
1	B	146	LYS
1	B	166	ARG
1	B	171	LEU
1	B	176	LYS
1	B	179	GLU
1	B	181	VAL
1	B	198	LYS
1	B	264	LYS
1	B	303	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	125	GLN
1	A	228	HIS
1	A	338	ASN
1	B	154	ASN

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Mol	Chain	Res	Type
1	B	319	GLN

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	295/348 (84%)	0.63	41 (13%) 6 6	15, 31, 66, 88	0
1	B	295/348 (84%)	0.80	52 (17%) 4 3	14, 33, 71, 94	0
All	All	590/696 (84%)	0.71	93 (15%) 5 4	14, 32, 69, 94	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	186	PRO	6.4
1	A	160	ALA	6.1
1	B	182	LEU	5.6
1	B	181	VAL	5.5
1	A	297	TYR	5.2
1	B	171	LEU	5.0
1	B	301	TYR	5.0
1	A	182	LEU	4.8
1	B	178	LEU	4.4
1	A	183	LYS	4.4
1	B	200	HIS	4.3
1	B	158	LEU	4.1
1	B	169	VAL	4.0
1	B	40	SER	3.9
1	A	171	LEU	3.9
1	A	170	LEU	3.9
1	A	175	LYS	3.9
1	B	297	TYR	3.8
1	B	160	ALA	3.8
1	A	301	TYR	3.8
1	B	315	GLY	3.7
1	B	39	SER	3.6
1	B	167	ASP	3.6
1	B	340	SER	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	174	VAL	3.5
1	B	174	VAL	3.4
1	A	169	VAL	3.4
1	A	340	SER	3.3
1	A	300	PRO	3.3
1	B	183	LYS	3.3
1	B	304	TRP	3.3
1	B	175	LYS	3.2
1	A	155	ILE	3.2
1	B	318	ILE	3.1
1	A	181	VAL	3.1
1	B	41	GLY	3.1
1	B	155	ILE	3.1
1	B	147	LEU	3.1
1	A	304	TRP	3.0
1	B	156	THR	3.0
1	B	150	ILE	2.9
1	B	177	SER	2.9
1	A	176	LYS	2.8
1	B	153	GLU	2.8
1	B	170	LEU	2.8
1	B	266	ASN	2.8
1	A	298	GLY	2.8
1	B	166	ARG	2.7
1	B	151	TYR	2.7
1	B	311	HIS	2.7
1	B	298	GLY	2.6
1	A	265	ASP	2.6
1	A	315	GLY	2.5
1	A	156	THR	2.5
1	A	339	THR	2.5
1	B	198	LYS	2.4
1	A	167	ASP	2.4
1	A	168	GLU	2.4
1	B	57	HIS	2.4
1	A	178	LEU	2.4
1	A	185	ASN	2.4
1	A	152	LEU	2.4
1	A	166	ARG	2.4
1	B	300	PRO	2.4
1	B	180	ASP	2.3
1	A	187	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	172	ASN	2.3
1	A	193	SER	2.3
1	B	154	ASN	2.3
1	B	196	LYS	2.3
1	B	303	ASP	2.2
1	B	157	LYS	2.2
1	B	149	THR	2.2
1	A	180	ASP	2.2
1	A	186	PRO	2.2
1	B	305	LEU	2.2
1	A	299	ASP	2.2
1	B	84	GLN	2.2
1	A	302	ASN	2.1
1	A	39	SER	2.1
1	B	199	GLY	2.1
1	B	172	ASN	2.1
1	A	318	ILE	2.1
1	B	71	ASP	2.1
1	A	316	SER	2.1
1	B	316	SER	2.1
1	A	57	HIS	2.1
1	A	311	HIS	2.1
1	B	42	HIS	2.1
1	A	158	LEU	2.1
1	B	179	GLU	2.0
1	A	197	ASP	2.0
1	B	264	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.