



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 08:03 PM UTC

PDB ID : 6Z8Q / pdb_00006z8q
Title : Copper transporter OprC
Authors : Bhamidimarri, S.P.; van den Berg, B.
Deposited on : 2020-06-02
Resolution : 2.71 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

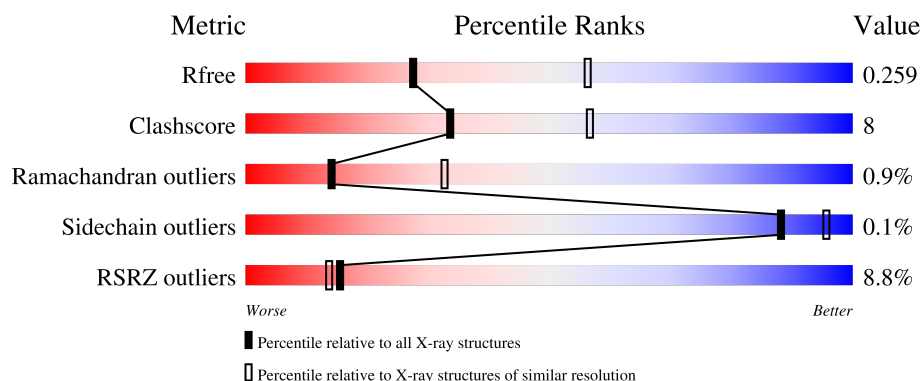
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.71 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	4348 (2.74-2.70)
Clashscore	190562	4665 (2.74-2.70)
Ramachandran outliers	187476	4584 (2.74-2.70)
Sidechain outliers	187428	4585 (2.74-2.70)
RSRZ outliers	180081	4348 (2.74-2.70)

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	723	6% 76% 14% 10%
1	B	723	10% 72% 16% 12%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10073 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 Putative copper transport outer membrane porin OprC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	653	Total	C	N	O	S	4	1	0
			5073	3185	895	975	18			
1	B	639	Total	C	N	O	S	0	1	0
			4981	3126	882	956	17			

- Molecule 2 is SILVER ION (CCD ID: AG) (formula: Ag).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Ag	0	0
			3	3		
2	B	3	Total	Ag	0	0
			3	3		

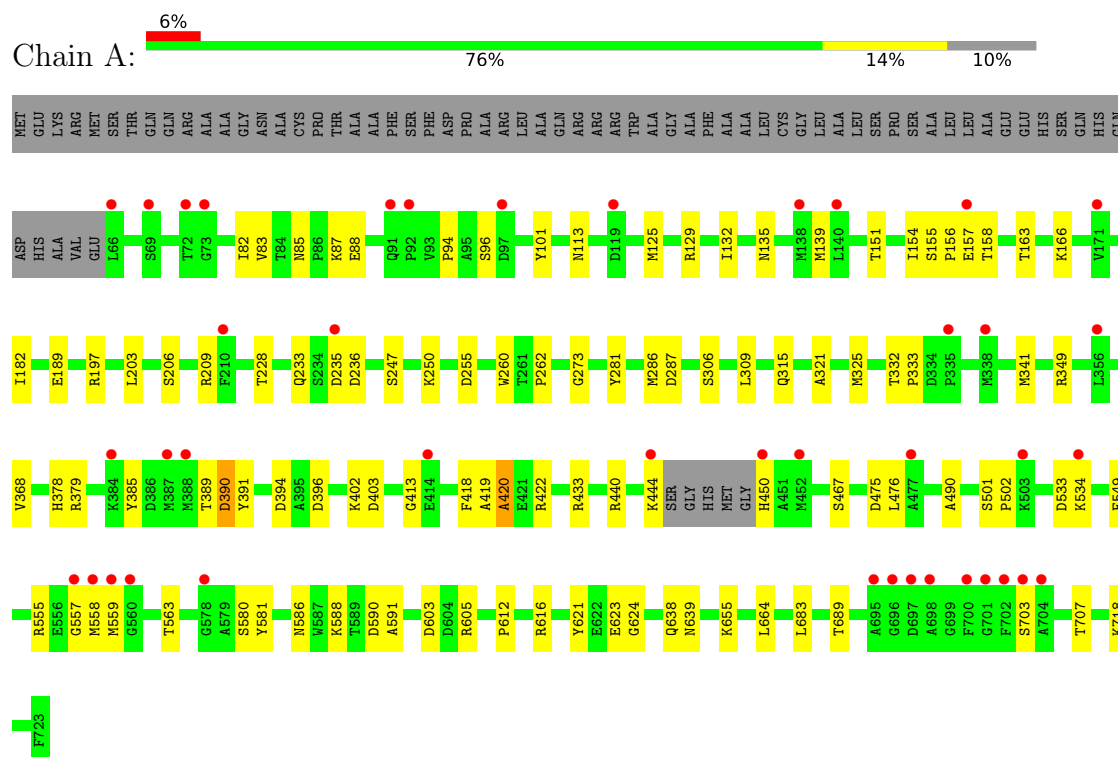
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	2	Total	O	0	0
			2	2		

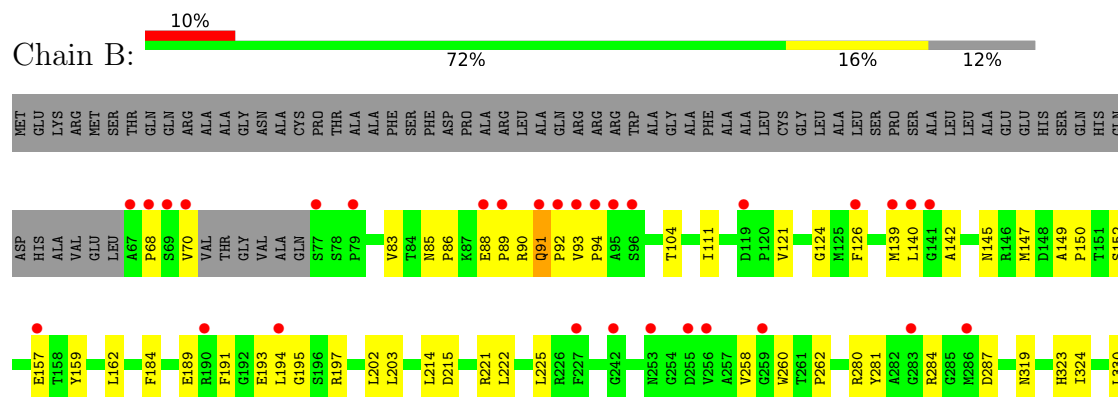
3 Residue-property plots [i](#)

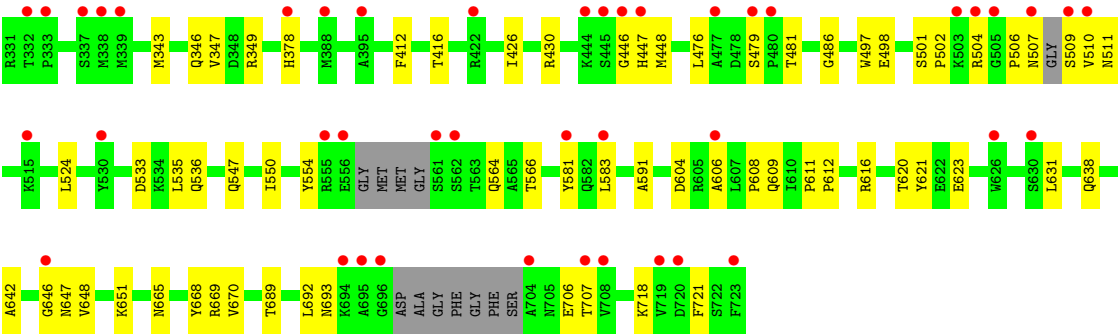
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Putative copper transport outer membrane porin OprC



- Molecule 1: Putative copper transport outer membrane porin OprC





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	156.20Å 195.29Å 167.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	97.65 – 2.71 97.65 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.9 (97.65-2.71) 99.9 (97.65-2.71)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.216 , 0.254 0.226 , 0.259	Depositor DCC
R_{free} test set	3476 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	76.4	Xtriage
Anisotropy	0.397	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10073	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% 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

Bond lengths and bond angles in the following residue types are not validated in this section: AG

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	0.40	0/5202	0.64	0/7052
1	B	0.39	0/5106	0.61	0/6918
All	All	0.39	0/10308	0.62	0/13970

There are no bond length outliers.

There are no bond angle outliers.

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	5073	0	4855	66	0
1	B	4981	0	4764	87	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	11	0	0	0	0
3	B	2	0	0	0	0
All	All	10073	0	9619	152	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 8.

All (152) 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:349:ARG:HH21	1:B:378:HIS:HB2	1.17	1.10
1:B:162:LEU:HD23	1:B:184:PHE:CD1	2.04	0.92
1:A:349:ARG:HH21	1:A:378:HIS:HB2	1.40	0.86
1:B:162:LEU:HD23	1:B:184:PHE:CE1	2.14	0.83
1:A:444:LYS:HE3	1:A:450:HIS:NE2	2.00	0.76
1:A:260:TRP:CE2	1:A:262:PRO:HG3	2.24	0.72
1:B:689:THR:HG21	1:B:707:THR:HB	1.72	0.71
1:A:422:ARG:HB3	1:A:475:ASP:HB2	1.73	0.71
1:A:82:ILE:HD11	1:A:166:LYS:HE3	1.72	0.70
1:A:151:THR:HA	1:A:154:ILE:HD12	1.73	0.70
1:A:580:SER:HB3	1:A:590:ASP:HB3	1.73	0.68
1:A:286:MET:HE3	1:A:325:MET:HE2	1.74	0.68
1:A:389:THR:O	1:A:391:TYR:N	2.24	0.67
1:B:260:TRP:CE2	1:B:262:PRO:HG3	2.30	0.66
1:B:349:ARG:NH2	1:B:378:HIS:HB2	2.01	0.65
1:A:125:MET:HE1	1:A:549:PHE:HE2	1.61	0.65
1:B:145:ASN:O	1:B:284:ARG:NH2	2.30	0.64
1:B:504:ARG:NE	1:B:504:ARG:HA	2.12	0.64
1:A:580:SER:HB2	1:A:588:LYS:HE2	1.79	0.64
1:B:612:PRO:HB3	1:B:638:GLN:HB2	1.81	0.63
1:A:433:ARG:NH1	1:B:547:GLN:OE1	2.31	0.63
1:B:90:ARG:HD3	1:B:104:THR:OG1	1.99	0.62
1:A:83:VAL:HG13	1:A:163:THR:HG22	1.81	0.61
1:B:68:PRO:HG3	1:B:669:ARG:HB2	1.82	0.61
1:A:87:LYS:HA	1:A:157:GLU:HG3	1.83	0.61
1:A:689:THR:HG21	1:A:707:THR:OG1	2.01	0.60
1:B:193:GLU:HG3	1:B:194:LEU:H	1.67	0.59
1:B:157:GLU:OE2	1:B:718:LYS:NZ	2.35	0.57
1:A:555:ARG:HH22	1:A:557:GLY:HA3	1.70	0.57
1:B:533:ASP:N	1:B:533:ASP:OD1	2.38	0.56
1:A:533:ASP:O	1:A:534:LYS:HB2	2.05	0.56
1:B:214:LEU:HD12	1:B:215:ASP:H	1.70	0.56
1:A:255:ASP:OD2	1:A:273:GLY:HA3	2.06	0.55
1:A:612:PRO:HB3	1:A:638:GLN:HB2	1.89	0.55
1:B:162:LEU:CD2	1:B:184:PHE:CE1	2.88	0.55
1:B:225:LEU:HD12	1:B:258[B]:VAL:HG22	1.89	0.55
1:A:379:ARG:HG2	1:A:402:LYS:HA	1.88	0.55
1:B:91:GLN:HG2	1:B:631:LEU:HD13	1.87	0.55
1:B:91:GLN:HB3	1:B:92:PRO:HD3	1.87	0.55
1:A:101:TYR:CD1	1:A:156:PRO:HG2	2.43	0.54
1:A:403:ASP:HB2	1:A:440:ARG:HG3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:ILE:HG12	1:B:692:LEU:HB3	1.89	0.54
1:A:444:LYS:HG3	1:A:450:HIS:HD2	1.72	0.54
1:B:85:ASN:CG	1:B:88:GLU:HG2	2.32	0.54
1:B:214:LEU:HD12	1:B:215:ASP:N	2.22	0.54
1:B:221:ARG:HG3	1:B:222:LEU:HD23	1.91	0.53
1:B:479:SER:HB3	1:B:481:THR:OG1	2.09	0.53
1:B:147:MET:HE3	1:B:281:TYR:CE1	2.43	0.53
1:A:125:MET:HE3	1:A:129:ARG:HD3	1.91	0.52
1:A:534:LYS:HB3	1:A:581:TYR:HE1	1.74	0.52
1:B:280:ARG:HD3	1:B:330:LEU:O	2.09	0.52
1:A:125:MET:HE1	1:A:549:PHE:CE2	2.43	0.52
1:A:664:LEU:HD11	1:A:683:LEU:HD22	1.92	0.52
1:B:606:ALA:O	1:B:642:ALA:HB2	2.10	0.52
1:B:609:GLN:HG3	1:B:647:ASN:ND2	2.24	0.52
1:A:639:ASN:HA	1:A:655:LYS:HG2	1.91	0.51
1:A:203:LEU:HD12	1:A:718:LYS:HB2	1.93	0.51
1:B:497:TRP:O	1:B:502:PRO:HD3	2.10	0.51
1:A:341:MET:HE3	1:A:385:TYR:HB3	1.91	0.51
1:B:535:LEU:HB2	1:B:581:TYR:CE1	2.45	0.51
1:A:385:TYR:OH	1:A:390:ASP:HB2	2.11	0.51
1:A:403:ASP:HB2	1:A:440:ARG:CG	2.41	0.51
1:A:603:ASP:HB3	1:A:605:ARG:HH21	1.76	0.51
1:B:189:GLU:HB3	1:B:191:PHE:CD1	2.46	0.51
1:B:504:ARG:HH22	1:B:511:ASN:HB3	1.75	0.50
1:A:189:GLU:OE2	1:A:197:ARG:NH2	2.35	0.50
1:A:85:ASN:ND2	1:A:88:GLU:OE2	2.43	0.49
1:B:591:ALA:HA	1:B:616:ARG:O	2.12	0.49
1:B:668:TYR:CE2	1:B:670:VAL:HG22	2.47	0.49
1:B:126:PHE:CE1	1:B:647:ASN:HB2	2.47	0.49
1:B:149:ALA:O	1:B:152:SER:HB3	2.13	0.49
1:A:418:PHE:C	1:A:420:ALA:H	2.21	0.49
1:A:418:PHE:O	1:A:420:ALA:N	2.41	0.48
1:B:140:LEU:HG	1:B:150:PRO:HG3	1.95	0.48
1:B:191:PHE:HE2	1:B:195:GLY:C	2.21	0.48
1:B:193:GLU:HG3	1:B:194:LEU:N	2.27	0.48
1:B:225:LEU:HD12	1:B:258[A]:VAL:HG22	1.94	0.48
1:B:189:GLU:CD	1:B:197:ARG:HH12	2.22	0.48
1:B:126:PHE:CZ	1:B:647:ASN:HB2	2.49	0.47
1:A:125:MET:CE	1:A:129:ARG:HD3	2.43	0.47
1:A:139:MET:CE	1:A:321:ALA:HB2	2.45	0.47
1:B:535:LEU:HB2	1:B:581:TYR:CD1	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:591:ALA:HA	1:A:616:ARG:O	2.15	0.47
1:B:91:GLN:HE22	1:B:620:THR:HG22	1.79	0.47
1:A:394:ASP:HB3	1:A:396:ASP:OD1	2.15	0.47
1:B:554:TYR:HB2	1:B:651:LYS:HE2	1.97	0.47
1:B:189:GLU:HB3	1:B:191:PHE:HD1	1.80	0.47
1:A:332:THR:HG22	1:A:333:PRO:HD2	1.97	0.46
1:A:534:LYS:HB3	1:A:581:TYR:CE1	2.50	0.46
1:B:486:GLY:O	1:B:524:LEU:HD12	2.15	0.46
1:B:507:ASN:H	1:B:509:SER:HB2	1.80	0.46
1:B:121:VAL:HG22	1:B:126:PHE:HB3	1.97	0.46
1:B:202:LEU:HD23	1:B:721:PHE:HE1	1.80	0.46
1:B:70:VAL:HA	1:B:89:PRO:HG2	1.97	0.46
1:B:260:TRP:CZ2	1:B:262:PRO:HG3	2.50	0.46
1:A:132:ILE:HD13	1:A:182:ILE:HB	1.98	0.46
1:B:668:TYR:HE2	1:B:670:VAL:HG22	1.80	0.46
1:A:586:ASN:HB3	1:A:621:TYR:CE1	2.51	0.46
1:B:416:THR:HG23	1:B:426:ILE:HG12	1.98	0.45
1:B:147:MET:HE2	1:B:323:HIS:CE1	2.51	0.45
1:B:86:PRO:HG2	1:B:159:TYR:O	2.16	0.45
1:B:498:GLU:HG2	1:B:550:ILE:HG21	1.97	0.45
1:A:555:ARG:HG2	1:A:555:ARG:HH11	1.82	0.45
1:B:343:MET:HE3	1:B:343:MET:HB2	1.69	0.45
1:B:412:PHE:HB3	1:B:430:ARG:HB3	1.97	0.45
1:A:557:GLY:O	1:A:559:MET:N	2.49	0.44
1:A:233:GLN:HB3	1:A:250:LYS:HG3	2.00	0.44
1:B:70:VAL:HG23	1:B:665:ASN:OD1	2.17	0.44
1:B:621:TYR:OH	1:B:623:GLU:OE1	2.35	0.44
1:A:135:ASN:O	1:A:315:GLN:NE2	2.41	0.44
1:B:189:GLU:OE1	1:B:197:ARG:NH1	2.44	0.44
1:B:609:GLN:HG3	1:B:647:ASN:HD22	1.83	0.44
1:B:70:VAL:HA	1:B:89:PRO:CG	2.48	0.43
1:A:113:ASN:HB3	1:A:281:TYR:HD1	1.83	0.43
1:B:203:LEU:HD12	1:B:718:LYS:HB2	2.00	0.43
1:B:446:GLY:O	1:B:448:MET:N	2.51	0.43
1:A:444:LYS:HE3	1:A:450:HIS:CD2	2.54	0.43
1:A:502:PRO:HA	1:A:563:THR:OG1	2.18	0.43
1:A:623:GLU:HB3	1:A:624:GLY:H	1.49	0.43
1:B:324:ILE:HG12	1:B:346:GLN:HG3	2.01	0.43
1:A:501:SER:N	1:A:502:PRO:HD2	2.33	0.43
1:B:535:LEU:HD12	1:B:536:GLN:N	2.34	0.43
1:A:113:ASN:HB3	1:A:281:TYR:CD1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:476:LEU:HD23	1:B:476:LEU:HA	1.82	0.43
1:B:506:PRO:HB3	1:B:566:THR:HA	2.01	0.43
1:B:504:ARG:NH1	1:B:509:SER:O	2.52	0.42
1:B:280:ARG:HG3	1:B:287:ASP:OD1	2.19	0.42
1:A:368:VAL:O	1:A:413:GLY:HA2	2.19	0.42
1:B:89:PRO:HB2	1:B:93:VAL:HG21	2.01	0.42
1:B:142:ALA:HB1	1:B:347:VAL:HG23	2.00	0.42
1:B:139:MET:HE2	1:B:319:ASN:CG	2.45	0.42
1:A:306:SER:HB3	1:A:309:LEU:HB3	2.02	0.42
1:B:139:MET:O	1:B:140:LEU:HD12	2.20	0.42
1:A:235:ASP:CG	1:A:236:ASP:H	2.28	0.42
1:B:91:GLN:HB3	1:B:92:PRO:CD	2.50	0.41
1:B:124:GLY:HA2	1:B:611:PRO:HD3	2.02	0.41
1:B:608:PRO:HB2	1:B:646:GLY:HA2	2.01	0.41
1:B:501:SER:N	1:B:502:PRO:CD	2.84	0.41
1:B:510:VAL:HG23	1:B:511:ASN:H	1.85	0.41
1:A:250:LYS:HB2	1:A:250:LYS:HE3	1.70	0.41
1:B:693:ASN:HB2	1:B:707:THR:HG22	2.03	0.41
1:B:497:TRP:HB3	1:B:648:VAL:HG21	2.03	0.41
1:B:706:GLU:HG2	1:B:707:THR:N	2.36	0.41
1:A:467:SER:HB3	1:A:490:ALA:HA	2.03	0.41
1:B:504:ARG:O	1:B:564:GLN:HA	2.20	0.41
1:A:158:THR:HG22	1:A:228:THR:HG22	2.03	0.40
1:B:162:LEU:C	1:B:162:LEU:HD13	2.45	0.40
1:A:96:SER:O	1:A:96:SER:OG	2.38	0.40
1:A:155:SER:O	1:A:158:THR:OG1	2.30	0.40
1:A:287:ASP:O	1:A:325:MET:HA	2.21	0.40
1:A:94:PRO:O	1:A:209:ARG:NH2	2.44	0.40
1:A:476:LEU:HD23	1:A:476:LEU:HA	1.87	0.40

There are no symmetry-related clashes.

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	650/723 (90%)	611 (94%)	32 (5%)	7 (1%)	11	27
1	B	630/723 (87%)	586 (93%)	39 (6%)	5 (1%)	16	35
All	All	1280/1446 (88%)	1197 (94%)	71 (6%)	12 (1%)	14	33

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	447	HIS
1	B	583	LEU
1	A	206	SER
1	A	390	ASP
1	B	91	GLN
1	A	247	SER
1	A	419	ALA
1	A	420	ALA
1	A	558	MET
1	A	703	SER
1	B	94	PRO
1	B	604	ASP

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	524/573 (91%)	524 (100%)	0	100	100
1	B	516/573 (90%)	515 (100%)	1 (0%)	87	95
All	All	1040/1146 (91%)	1039 (100%)	1 (0%)	88	95

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	83	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	586	ASN
1	A	673	ASN
1	A	705	ASN
1	B	290	GLN
1	B	474	HIS
1	B	673	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](#)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	653/723 (90%)	0.48	41 (6%)	26	23	42, 68, 115, 161	1 (0%)
1	B	639/723 (88%)	0.81	73 (11%)	10	9	51, 80, 125, 152	1 (0%)
All	All	1292/1446 (89%)	0.64	114 (8%)	15	14	42, 73, 122, 161	2 (0%)

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	694	LYS	5.8
1	B	93	VAL	5.8
1	B	92	PRO	5.3
1	A	66	LEU	5.1
1	A	450	HIS	4.8
1	B	70	VAL	4.6
1	A	696	GLY	4.6
1	B	338	MET	4.5
1	B	510	VAL	4.3
1	B	704	ALA	4.3
1	B	339	MET	4.2
1	A	140	LEU	4.1
1	A	92	PRO	4.1
1	B	696	GLY	4.0
1	B	140	LEU	4.0
1	A	695	ALA	3.9
1	A	119	ASP	3.7
1	B	561	SER	3.7
1	A	558	MET	3.7
1	A	338	MET	3.7
1	A	444	LYS	3.6
1	B	445	SER	3.6
1	B	562	SER	3.6
1	B	447	HIS	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	695	ALA	3.5
1	A	335	PRO	3.5
1	B	157	GLU	3.5
1	A	559	MET	3.4
1	A	91	GLN	3.4
1	B	504	ARG	3.4
1	B	723	PHE	3.4
1	B	444	LYS	3.3
1	A	698	ALA	3.3
1	A	701	GLY	3.3
1	A	700	PHE	3.3
1	A	702	PHE	3.3
1	B	581	TYR	3.3
1	B	68	PRO	3.2
1	B	91	GLN	3.2
1	A	388	MET	3.2
1	A	138	MET	3.2
1	B	507	ASN	3.1
1	B	94	PRO	3.1
1	B	332	THR	3.1
1	B	509	SER	3.0
1	B	67	ALA	3.0
1	A	503	LYS	3.0
1	B	89	PRO	2.9
1	B	126	PHE	2.9
1	B	194	LEU	2.8
1	B	555	ARG	2.8
1	B	477	ALA	2.8
1	A	387	MET	2.8
1	B	256	VAL	2.8
1	A	157	GLU	2.8
1	A	384	LYS	2.7
1	B	96	SER	2.7
1	A	210	PHE	2.7
1	B	515	LYS	2.7
1	A	73	GLY	2.6
1	B	69	SER	2.6
1	B	333	PRO	2.6
1	B	446	GLY	2.6
1	B	79	PRO	2.6
1	A	69	SER	2.6
1	A	697	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	141	GLY	2.5
1	A	578	GLY	2.5
1	B	708	VAL	2.5
1	B	95	ALA	2.5
1	B	139	MET	2.5
1	A	704	ALA	2.5
1	A	703	SER	2.4
1	B	583	LEU	2.4
1	B	378	HIS	2.4
1	A	72	THR	2.4
1	B	253	ASN	2.4
1	B	388	MET	2.4
1	B	190	ARG	2.4
1	B	337	SER	2.3
1	B	503	LYS	2.3
1	B	720	ASP	2.3
1	B	480	PRO	2.3
1	B	505	GLY	2.3
1	A	534	LYS	2.3
1	A	235	ASP	2.3
1	A	414	GLU	2.3
1	B	556	GLU	2.3
1	B	606	ALA	2.2
1	B	286	MET	2.2
1	B	259	GLY	2.2
1	B	88	GLU	2.2
1	B	242	GLY	2.2
1	A	477	ALA	2.2
1	B	530	TYR	2.2
1	A	557	GLY	2.2
1	A	171	VAL	2.2
1	A	452	MET	2.2
1	B	719	VAL	2.2
1	B	395	ALA	2.1
1	B	255	ASP	2.1
1	B	479	SER	2.1
1	B	626	TRP	2.1
1	A	560	GLY	2.1
1	A	97	ASP	2.1
1	B	646	GLY	2.1
1	B	119	ASP	2.0
1	B	630	SER	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	707	THR	2.0
1	B	227	PHE	2.0
1	B	422	ARG	2.0
1	B	77	SER	2.0
1	A	356	LEU	2.0
1	B	283	GLY	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	AG	B	802	1/1	0.81	0.28	196,196,196,196	0
2	AG	A	803	1/1	0.92	0.14	166,166,166,166	0
2	AG	A	802	1/1	0.95	0.14	171,171,171,171	0
2	AG	B	803	1/1	0.96	0.10	157,157,157,157	0
2	AG	A	801	1/1	1.00	0.03	82,82,82,82	0
2	AG	B	801	1/1	1.00	0.06	89,89,89,89	0

6.5 Other polymers [i](#)

There are no such residues in this entry.