



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:36 AM UTC

PDB ID : 6Z8Y / pdb_00006z8y
Title : Copper transporter OprC
Authors : Bhamidimarri, S.P.; van den Berg, B.
Deposited on : 2020-06-02
Resolution : 2.78 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

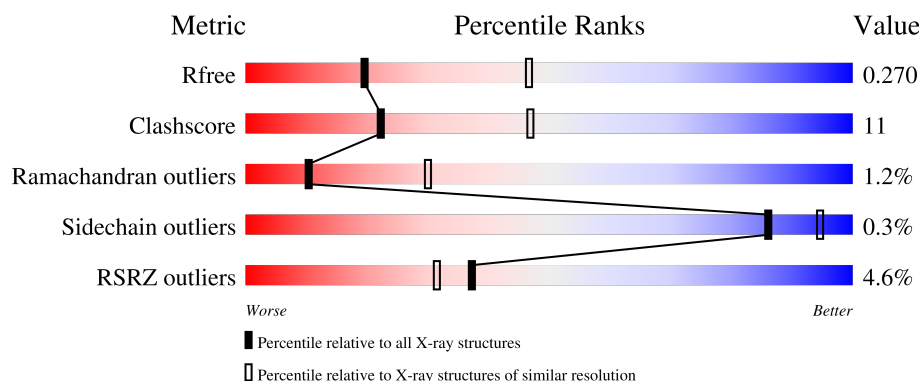
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.78 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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	4% 67% 23% 9%
1	B	723	5% 68% 21% 11%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10142 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	658	Total	C	N	O	S	0	1	0
			5104	3203	902	981	18			
1	B	645	Total	C	N	O	S	0	1	0
			5015	3151	884	965	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	ALA	CYS	engineered mutation	UNP G3XD89
B	143	ALA	CYS	engineered mutation	UNP G3XD89

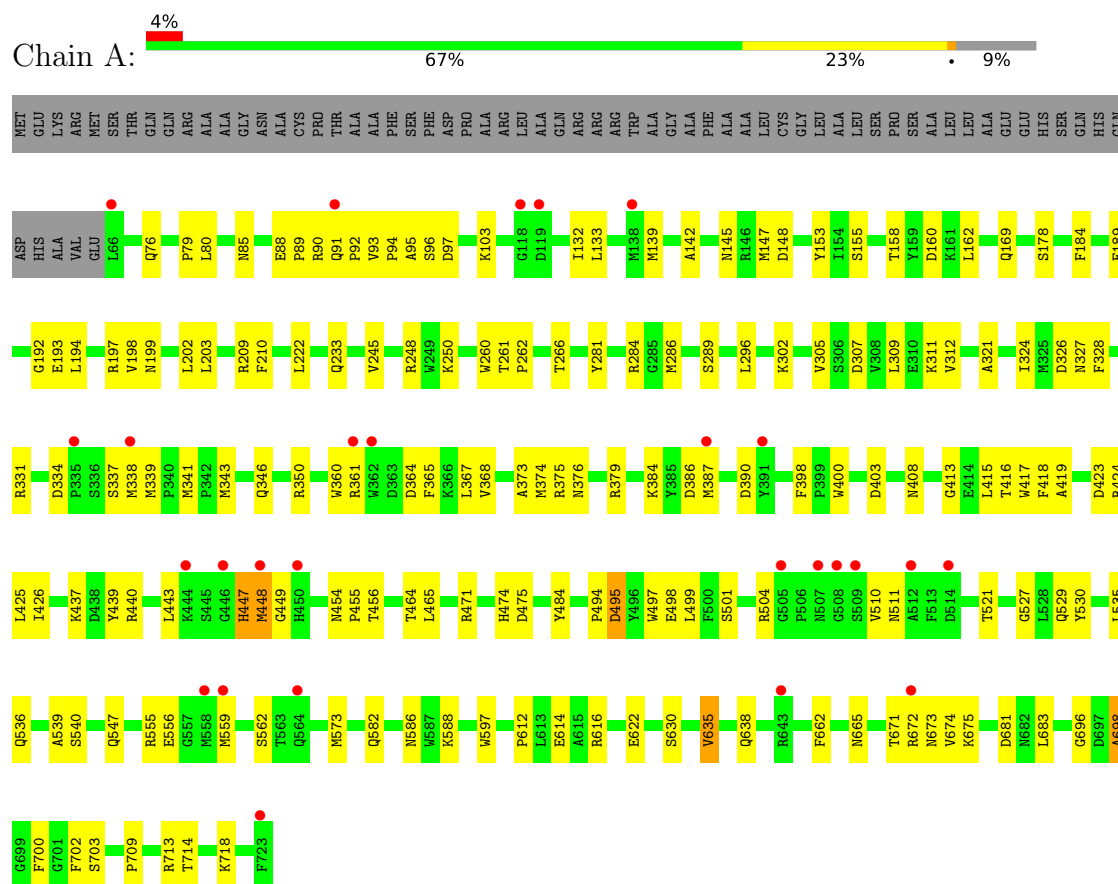
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	12	Total	O	0	0
			12	12		
2	B	11	Total	O	0	0
			11	11		

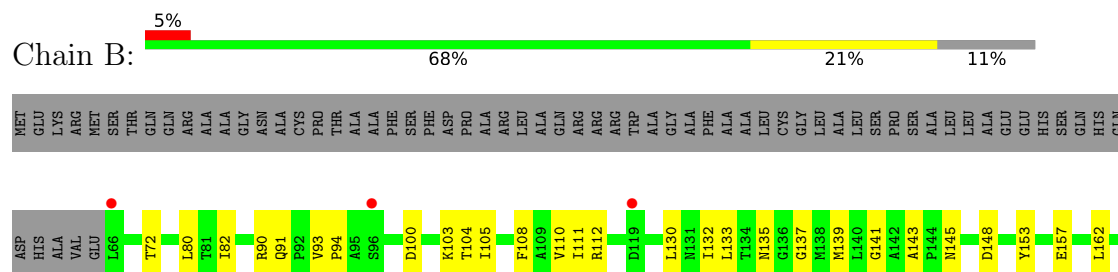
3 Residue-property plots [i](#)

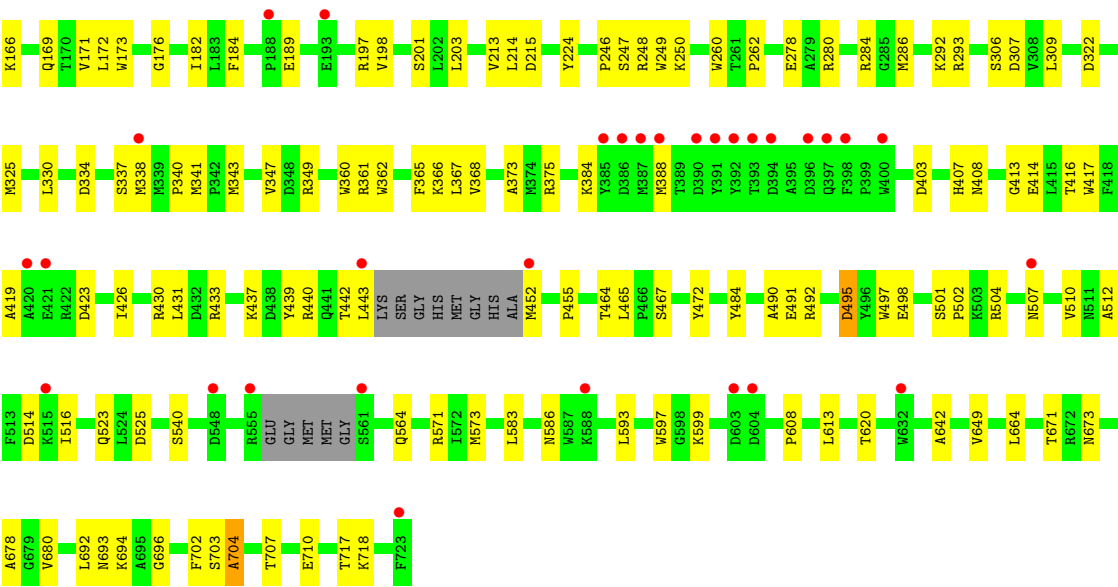
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	67.03Å 170.74Å 198.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	67.03 – 2.78 67.03 – 2.78	Depositor EDS
% Data completeness (in resolution range)	99.8 (67.03-2.78) 100.0 (67.03-2.78)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.06 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.210 , 0.266 0.214 , 0.270	Depositor DCC
R_{free} test set	2894 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	69.3	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 38.0	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	0.93	EDS
Total number of atoms	10142	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.34% 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	0.40	0/5235	0.65	0/7097
1	B	0.40	0/5142	0.67	0/6973
All	All	0.40	0/10377	0.66	0/14070

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	5104	0	4884	112	0
1	B	5015	0	4800	98	0
2	A	12	0	0	0	0
2	B	11	0	0	0	0
All	All	10142	0	9684	209	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 11.

All (209) 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:88:GLU:O	1:A:90:ARG:NH2	1.65	1.28

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:172:LEU:HD13	1:B:414:GLU:OE2	1.75	0.84
1:A:504:ARG:HH12	1:A:562:SER:HB3	1.43	0.84
1:A:289:SER:OG	1:A:326:ASP:OD2	1.95	0.84
1:A:614:GLU:HG3	1:A:635:VAL:HG12	1.64	0.78
1:B:423:ASP:OD2	1:B:472:TYR:OH	2.01	0.78
1:B:671:THR:HG23	1:B:673:ASN:H	1.49	0.78
1:A:324:ILE:HG12	1:A:346:GLN:HG3	1.65	0.76
1:A:681:ASP:HB2	1:A:714:THR:HB	1.69	0.74
1:B:286:MET:HE2	1:B:325:MET:HE2	1.69	0.73
1:A:341:MET:HE2	1:A:343:MET:HE1	1.70	0.73
1:A:671:THR:HG22	1:A:674:VAL:HG12	1.72	0.71
1:A:448:MET:HE3	1:A:559:MET:HE3	1.73	0.69
1:B:361:ARG:HH11	1:B:361:ARG:HG2	1.60	0.66
1:A:334:ASP:HB3	1:A:337:SER:HB2	1.78	0.66
1:A:139:MET:HE3	1:A:321:ALA:HB2	1.77	0.65
1:A:437:LYS:HD3	1:A:439:TYR:CZ	2.32	0.64
1:B:172:LEU:HD11	1:B:414:GLU:HB2	1.79	0.64
1:B:403:ASP:HB2	1:B:440:ARG:HG2	1.78	0.64
1:A:423:ASP:OD1	1:A:474:HIS:ND1	2.26	0.64
1:A:671:THR:HG23	1:A:673:ASN:H	1.63	0.63
1:B:593:LEU:HD11	1:B:613:LEU:HD11	1.79	0.63
1:A:586:ASN:HB2	1:A:622:GLU:O	1.99	0.63
1:A:403:ASP:HB2	1:A:440:ARG:HG2	1.81	0.62
1:A:203:LEU:HD12	1:A:718:LYS:HD3	1.82	0.62
1:A:350:ARG:HH12	1:A:379:ARG:NH2	1.98	0.62
1:B:703:SER:OG	1:B:704:ALA:N	2.33	0.61
1:B:172:LEU:HD13	1:B:414:GLU:CD	2.26	0.61
1:A:612:PRO:HB3	1:A:638:GLN:HB2	1.82	0.60
1:A:91:GLN:OE1	1:A:665:ASN:HB3	2.02	0.60
1:B:417:TRP:CE2	1:B:419:ALA:HB2	2.36	0.60
1:A:417:TRP:CE2	1:A:419:ALA:HB2	2.37	0.60
1:B:82:ILE:HD11	1:B:166:LYS:HE2	1.84	0.60
1:A:447:HIS:C	1:A:449:GLY:H	2.10	0.59
1:A:555:ARG:HG2	1:A:556:GLU:N	2.18	0.59
1:B:608:PRO:HB3	1:B:642:ALA:HB3	1.83	0.59
1:A:103:LYS:O	1:A:616:ARG:NH2	2.35	0.59
1:A:443:LEU:HG	1:A:454:ASN:HB2	1.85	0.59
1:B:306:SER:HB3	1:B:309:LEU:HB3	1.83	0.59
1:A:309:LEU:HD12	1:A:360:TRP:CZ3	2.38	0.59
1:B:497:TRP:O	1:B:501:SER:HB2	2.03	0.59
1:B:80:LEU:HD13	1:B:169:GLN:HG3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:GLN:HB3	1:A:471:ARG:NH1	2.18	0.58
1:A:260:TRP:CE2	1:A:262:PRO:HG3	2.39	0.57
1:A:456:THR:OG1	1:A:511:ASN:ND2	2.35	0.57
1:A:547:GLN:HG3	1:B:433:ARG:HH12	1.68	0.57
1:A:364:ASP:HB3	1:A:418:PHE:HB2	1.87	0.57
1:B:141:GLY:O	1:B:349:ARG:NH1	2.38	0.56
1:A:367:LEU:HB2	1:A:415:LEU:HD12	1.87	0.56
1:B:90:ARG:HG2	1:B:104:THR:HG21	1.88	0.56
1:A:582:GLN:HG3	1:A:588:LYS:HG3	1.87	0.55
1:B:442:THR:HG22	1:B:443:LEU:H	1.71	0.55
1:A:202:LEU:HD11	1:A:210:PHE:CZ	2.43	0.54
1:A:305:VAL:HB	1:A:309:LEU:HD22	1.89	0.54
1:B:112:ARG:NE	1:B:710:GLU:OE1	2.29	0.54
1:A:80:LEU:HD13	1:A:169:GLN:HG3	1.90	0.54
1:A:671:THR:HG22	1:A:674:VAL:CG1	2.38	0.54
1:B:148:ASP:HB2	1:B:153:TYR:CE2	2.43	0.54
1:B:341:MET:HE2	1:B:343:MET:SD	2.49	0.54
1:A:671:THR:OG1	1:A:672:ARG:NH1	2.42	0.53
1:B:361:ARG:HG3	1:B:365:PHE:O	2.08	0.53
1:A:504:ARG:NH1	1:A:562:SER:HB3	2.18	0.53
1:A:94:PRO:O	1:A:209:ARG:NH2	2.41	0.53
1:B:111:ILE:HG12	1:B:692:LEU:HB3	1.91	0.52
1:B:306:SER:OG	1:B:307:ASP:N	2.43	0.52
1:A:696:GLY:HA3	1:A:702:PHE:CG	2.44	0.52
1:B:280:ARG:NH1	1:B:330:LEU:O	2.43	0.52
1:B:189:GLU:OE1	1:B:197:ARG:NH2	2.39	0.52
1:A:91:GLN:HB3	1:A:92:PRO:HD3	1.92	0.52
1:B:145:ASN:HB2	1:B:284:ARG:HH12	1.75	0.52
1:A:169:GLN:HB3	1:A:471:ARG:HH12	1.74	0.51
1:B:416:THR:HG23	1:B:426:ILE:CD1	2.40	0.51
1:A:328:PHE:HA	1:A:331:ARG:O	2.10	0.51
1:A:368:VAL:O	1:A:413:GLY:HA2	2.11	0.51
1:B:214:LEU:HD12	1:B:215:ASP:H	1.75	0.51
1:A:497:TRP:O	1:A:501:SER:HB2	2.11	0.51
1:B:292:LYS:HB3	1:B:322:ASP:HB3	1.92	0.51
1:A:147:MET:HE3	1:A:281:TYR:CE2	2.45	0.51
1:A:266:THR:OG1	1:A:302:LYS:HE2	2.11	0.51
1:A:455:PRO:HG2	1:A:510:VAL:HG12	1.93	0.51
1:B:201:SER:HB3	1:B:213:VAL:CG1	2.40	0.51
1:A:189:GLU:OE2	1:A:197:ARG:NH2	2.44	0.51
1:A:527:GLY:HA3	1:A:539:ALA:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:MET:HE2	1:A:153:TYR:CG	2.46	0.50
1:A:337:SER:OG	1:A:338:MET:N	2.43	0.50
1:B:368:VAL:O	1:B:413:GLY:HA2	2.11	0.50
1:B:375:ARG:HG3	1:B:407:HIS:CD2	2.45	0.50
1:A:76:GLN:HE21	1:A:529:GLN:NE2	2.09	0.50
1:A:374:MET:HG2	1:A:375:ARG:N	2.27	0.50
1:B:696:GLY:HA3	1:B:702:PHE:CG	2.47	0.50
1:A:309:LEU:HD21	1:A:312:VAL:HG23	1.94	0.49
1:A:464:THR:C	1:A:465:LEU:HG	2.36	0.49
1:A:683:LEU:O	1:A:713:ARG:HD3	2.13	0.49
1:B:139:MET:HE1	1:B:293:ARG:HG3	1.94	0.49
1:B:214:LEU:HD12	1:B:215:ASP:N	2.26	0.49
1:A:672:ARG:HH21	1:A:672:ARG:HG3	1.77	0.49
1:B:338:MET:C	1:B:340:PRO:HD3	2.37	0.49
1:B:649:VAL:O	1:B:694:LYS:HE2	2.13	0.49
1:A:334:ASP:CB	1:A:337:SER:HB2	2.42	0.48
1:A:386:ASP:OD2	1:A:387:MET:N	2.47	0.48
1:B:260:TRP:CE2	1:B:262:PRO:HG3	2.48	0.48
1:A:193:GLU:HG3	1:A:194:LEU:H	1.78	0.48
1:B:437:LYS:HD3	1:B:439:TYR:CZ	2.47	0.48
1:B:133:LEU:HD22	1:B:173:TRP:HB3	1.96	0.48
1:B:361:ARG:HH11	1:B:361:ARG:CG	2.26	0.48
1:B:417:TRP:CZ2	1:B:419:ALA:HB2	2.49	0.48
1:B:132:ILE:HG12	1:B:182:ILE:HD12	1.95	0.48
1:B:133:LEU:CD2	1:B:173:TRP:HB3	2.43	0.48
1:A:373:ALA:HA	1:A:408:ASN:O	2.14	0.48
1:A:85:ASN:O	1:A:90:ARG:NH1	2.47	0.47
1:B:495:ASP:HB3	1:B:498:GLU:H	1.77	0.47
1:B:455:PRO:HB2	1:B:510:VAL:HG21	1.97	0.47
1:A:307:ASP:OD1	1:A:307:ASP:N	2.47	0.47
1:A:424:ARG:NH2	1:A:475:ASP:OD1	2.48	0.47
1:B:309:LEU:HD13	1:B:360:TRP:CE3	2.50	0.47
1:A:142:ALA:H	1:A:148:ASP:CG	2.22	0.47
1:A:309:LEU:HD21	1:A:312:VAL:CG2	2.44	0.47
1:A:424:ARG:HH22	1:A:475:ASP:CG	2.23	0.47
1:B:248:ARG:HG2	1:B:249:TRP:N	2.28	0.46
1:B:467:SER:HB3	1:B:490:ALA:HA	1.97	0.46
1:A:573:MET:HE3	1:A:597:TRP:CZ3	2.51	0.46
1:B:100:ASP:HA	1:B:103:LYS:HG3	1.97	0.46
1:A:535:LEU:HD12	1:A:536:GLN:N	2.31	0.46
1:B:130:LEU:HD23	1:B:182:ILE:HD11	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:GLY:HA3	1:B:492:ARG:HG3	1.98	0.45
1:B:135:ASN:C	1:B:137:GLY:H	2.24	0.45
1:B:189:GLU:HB2	1:B:224:TYR:CD2	2.51	0.45
1:B:586:ASN:OD1	1:B:586:ASN:N	2.48	0.45
1:A:178:SER:HB3	1:A:521:THR:HG21	1.98	0.45
1:B:162:LEU:HD13	1:B:184:PHE:CE2	2.53	0.45
1:A:341:MET:HE2	1:A:343:MET:CE	2.44	0.44
1:B:246:PRO:HB3	1:B:280:ARG:HD2	1.98	0.44
1:A:447:HIS:C	1:A:449:GLY:N	2.74	0.44
1:B:203:LEU:HD12	1:B:718:LYS:HB2	2.00	0.44
1:B:452:MET:SD	1:B:452:MET:C	3.00	0.44
1:A:698:ALA:C	1:A:700:PHE:H	2.26	0.44
1:B:334:ASP:HB2	1:B:337:SER:HB2	1.98	0.44
1:A:494:PRO:HB2	1:A:499:LEU:HG	2.00	0.44
1:A:222:LEU:HA	1:A:261:THR:HG23	2.00	0.44
1:A:79:PRO:HB3	1:A:484:TYR:CD1	2.52	0.44
1:B:373:ALA:HA	1:B:408:ASN:O	2.18	0.44
1:B:465:LEU:HB3	1:B:491:GLU:HB2	2.00	0.44
1:A:447:HIS:CG	1:A:448:MET:H	2.34	0.43
1:B:250:LYS:HB3	1:B:278:GLU:HG2	1.98	0.43
1:A:76:GLN:OE1	1:A:540:SER:HB2	2.18	0.43
1:A:96:SER:O	1:A:96:SER:OG	2.28	0.43
1:A:302:LYS:O	1:A:311:LYS:HA	2.18	0.43
1:A:425:LEU:HD12	1:A:425:LEU:HA	1.80	0.43
1:A:197:ARG:CZ	1:A:197:ARG:HB2	2.47	0.43
1:B:367:LEU:HD12	1:B:414:GLU:O	2.19	0.43
1:B:523:GLN:NE2	1:B:525:ASP:OD2	2.48	0.43
1:B:678:ALA:HB2	1:B:717[B]:THR:HG22	2.00	0.43
1:A:286:MET:HG2	1:A:327:ASN:HB3	1.99	0.43
1:A:132:ILE:O	1:A:133:LEU:HD23	2.18	0.43
1:B:338:MET:HE3	1:B:338:MET:H	1.83	0.43
1:A:495:ASP:HB3	1:A:498:GLU:H	1.84	0.43
1:A:95:ALA:O	1:A:97:ASP:N	2.49	0.43
1:A:504:ARG:HH22	1:A:555:ARG:HB3	1.84	0.43
1:B:286:MET:CE	1:B:325:MET:HE2	2.43	0.43
1:B:693:ASN:HB2	1:B:707:THR:HG22	2.01	0.43
1:A:76:GLN:HE21	1:A:529:GLN:HE22	1.67	0.43
1:A:147:MET:HE3	1:A:281:TYR:CD2	2.54	0.43
1:A:379:ARG:HH11	1:A:400:TRP:CD1	2.37	0.43
1:B:105:ILE:HB	1:B:108:PHE:HD2	1.84	0.43
1:B:340:PRO:HD2	1:B:341:MET:SD	2.59	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:ALA:HA	1:B:347:VAL:HG21	2.00	0.42
1:B:145:ASN:HB2	1:B:284:ARG:NH1	2.34	0.42
1:B:384:LYS:HE3	1:B:384:LYS:HB2	1.67	0.42
1:A:233:GLN:HB3	1:A:250:LYS:HG3	2.02	0.42
1:A:417:TRP:CZ2	1:A:419:ALA:HB2	2.55	0.42
1:A:145:ASN:HD22	1:A:284:ARG:HH11	1.67	0.42
1:A:384:LYS:HD3	1:A:398:PHE:CE1	2.54	0.42
1:B:93:VAL:HA	1:B:94:PRO:HD3	1.93	0.42
1:B:157:GLU:H	1:B:157:GLU:CD	2.28	0.42
1:B:172:LEU:CD1	1:B:414:GLU:CD	2.91	0.42
1:B:375:ARG:HG3	1:B:407:HIS:NE2	2.35	0.42
1:A:296:LEU:HA	1:A:296:LEU:HD23	1.85	0.42
1:B:171:VAL:HG22	1:B:414:GLU:HG3	2.02	0.42
1:B:431:LEU:HD11	1:B:464:THR:HG22	2.02	0.42
1:A:88:GLU:C	1:A:90:ARG:HH22	2.27	0.42
1:A:386:ASP:O	1:A:390:ASP:HA	2.20	0.41
1:B:664:LEU:HB2	1:B:680:VAL:HG12	2.02	0.41
1:A:350:ARG:O	1:A:376:ASN:HA	2.20	0.41
1:A:155:SER:O	1:A:158:THR:OG1	2.38	0.41
1:A:403:ASP:HB2	1:A:440:ARG:CG	2.50	0.41
1:B:484:TYR:OH	1:B:540:SER:HB2	2.20	0.41
1:B:502:PRO:HG2	1:B:512:ALA:CB	2.50	0.41
1:A:93:VAL:HA	1:A:94:PRO:HD3	1.87	0.41
1:B:93:VAL:O	1:B:93:VAL:HG13	2.21	0.41
1:B:360:TRP:HB2	1:B:362:TRP:HZ3	1.85	0.41
1:A:416:THR:HG23	1:A:426:ILE:HG13	2.02	0.41
1:B:103:LYS:HE2	1:B:110:VAL:HG21	2.02	0.41
1:B:571:ARG:HH22	1:B:573:MET:CE	2.34	0.41
1:A:160:ASP:N	1:A:160:ASP:OD1	2.53	0.41
1:B:516:ILE:HD13	1:B:516:ILE:HA	1.85	0.41
1:B:597:TRP:CH2	1:B:599:LYS:HB2	2.55	0.41
1:B:72:THR:HG22	1:B:91:GLN:HB3	2.02	0.41
1:B:361:ARG:CG	1:B:361:ARG:NH1	2.84	0.41
1:A:245:VAL:HG21	1:A:709:PRO:HG2	2.03	0.40
1:B:360:TRP:O	1:B:366:LYS:HA	2.22	0.40
1:A:89:PRO:HB3	1:A:718:LYS:HE3	2.02	0.40
1:A:361:ARG:HA	1:A:365:PHE:O	2.21	0.40
1:A:162:LEU:HD13	1:A:184:PHE:CE1	2.56	0.40
1:A:630:SER:HB2	1:A:662:PHE:CE1	2.56	0.40
1:B:504:ARG:HH21	1:B:564:GLN:HB2	1.87	0.40
1:A:198:VAL:C	1:A:199:ASN:HD22	2.30	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:530:TYR:O	1:A:536:GLN:HA	2.22	0.40
1:B:171:VAL:O	1:B:430:ARG:NH1	2.55	0.40
1:B:198:VAL:HG13	1:B:214:LEU:HD11	2.02	0.40

There are no symmetry-related clashes.

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	657/723 (91%)	591 (90%)	58 (9%)	8 (1%)	10	30
1	B	640/723 (88%)	591 (92%)	42 (7%)	7 (1%)	11	32
All	All	1297/1446 (90%)	1182 (91%)	100 (8%)	15 (1%)	10	30

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	495	ASP
1	B	388	MET
1	A	698	ALA
1	B	495	ASP
1	B	507	ASN
1	B	514	ASP
1	A	703	SER
1	B	704	ALA
1	A	248	ARG
1	A	447	HIS
1	A	448	MET
1	B	247	SER
1	B	583	LEU
1	A	192	GLY
1	A	339	MET

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	526/572 (92%)	524 (100%)	2 (0%)	84	93
1	B	518/572 (91%)	517 (100%)	1 (0%)	87	96
All	All	1044/1144 (91%)	1041 (100%)	3 (0%)	86	94

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	635	VAL
1	A	675	LYS
1	B	620	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	ASN
1	A	199	ASN
1	A	207	ASN
1	A	529	GLN
1	A	536	GLN
1	A	582	GLN
1	A	639	ASN
1	B	230	ASN
1	B	253	ASN
1	B	408	ASN
1	B	441	GLN
1	B	458	ASN
1	B	529	GLN

5.3.3 RNA ⓘ

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	658/723 (91%)	0.25	27 (4%) 41 35	36, 67, 107, 142	1 (0%)
1	B	645/723 (89%)	0.26	33 (5%) 33 27	35, 66, 96, 124	1 (0%)
All	All	1303/1446 (90%)	0.25	60 (4%) 37 31	35, 67, 102, 142	2 (0%)

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	66	LEU	6.8
1	B	443	LEU	6.3
1	B	723	PHE	4.9
1	A	514	ASP	4.5
1	B	507	ASN	4.3
1	B	390	ASP	4.3
1	A	362	TRP	4.2
1	A	387	MET	4.1
1	B	338	MET	3.9
1	B	393	THR	3.9
1	B	452	MET	3.9
1	A	723	PHE	3.8
1	B	392	TYR	3.6
1	A	450	HIS	3.5
1	B	119	ASP	3.4
1	B	391	TYR	3.3
1	A	119	ASP	3.3
1	A	338	MET	3.3
1	B	397	GLN	3.2
1	B	385	TYR	3.2
1	A	507	ASN	3.2
1	B	555	ARG	3.0
1	A	505	GLY	3.0
1	B	396	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	448	MET	2.9
1	B	604	ASP	2.9
1	B	632	TRP	2.9
1	B	66	LEU	2.9
1	A	509	SER	2.8
1	B	515	LYS	2.8
1	A	558	MET	2.8
1	B	388	MET	2.8
1	B	603	ASP	2.7
1	A	91	GLN	2.7
1	A	138	MET	2.7
1	B	548	ASP	2.6
1	A	118	GLY	2.5
1	A	335	PRO	2.5
1	B	394	ASP	2.4
1	A	508	GLY	2.4
1	B	398	PHE	2.4
1	B	420	ALA	2.4
1	B	96	SER	2.3
1	A	643	ARG	2.2
1	A	559	MET	2.2
1	A	446	GLY	2.2
1	B	387	MET	2.2
1	A	444	LYS	2.2
1	B	188	PRO	2.1
1	A	391	TYR	2.1
1	B	386	ASP	2.1
1	A	512	ALA	2.1
1	B	193	GLU	2.1
1	A	672	ARG	2.1
1	B	588	LYS	2.1
1	A	564	GLN	2.1
1	B	561	SER	2.1
1	B	400	TRP	2.1
1	A	361	ARG	2.0
1	B	421	GLU	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.