



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:47 AM UTC

PDB ID : 8YJX / pdb_00008yjsx
Title : Crystal structure of penicillin-binding protein 2 (PBP2) from *Campylobacter jejuni*
Authors : Choi, H.J.; Ki, D.W.; Yoon, S.I.
Deposited on : 2024-03-03
Resolution : 2.95 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

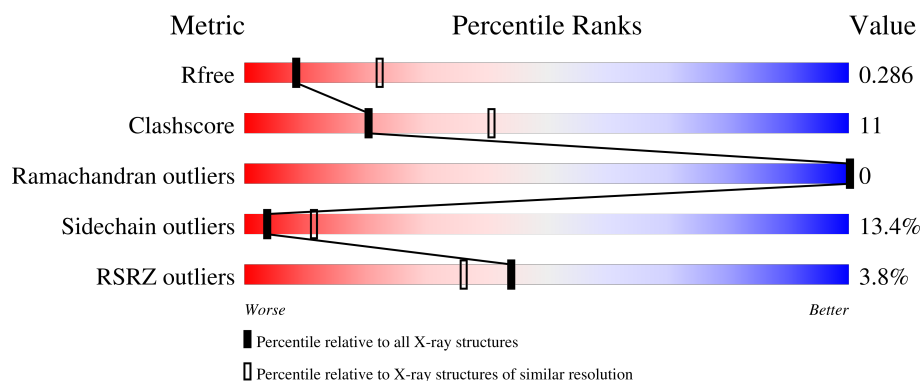
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	568	3% 69% 25% • •
1	B	568	4% 63% 21% • 14%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7642 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 Penicillin-binding protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	558	Total	C	N	O	S	0	0	0
			4126	2634	689	794	9			
1	B	490	Total	C	N	O	S	0	0	0
			3508	2242	592	666	8			

There are 12 discrepancies between the modelled and reference sequences:

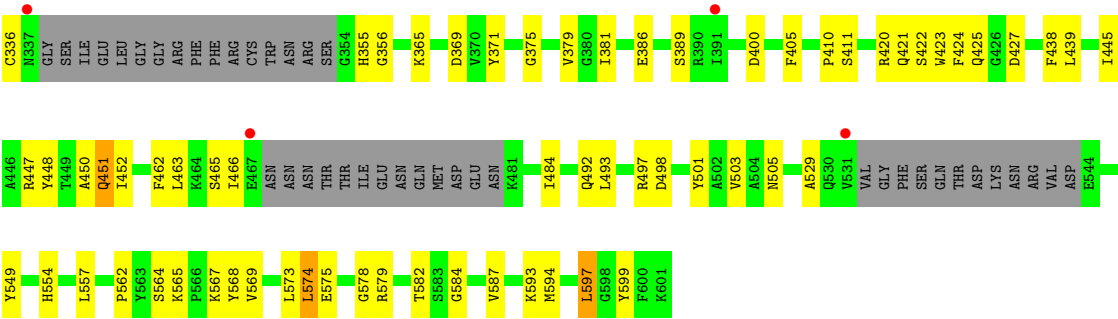
Chain	Residue	Modelled	Actual	Comment	Reference
A	34	GLY	-	expression tag	UNP A0A3X8PTV6
A	35	SER	-	expression tag	UNP A0A3X8PTV6
A	36	ALA	-	expression tag	UNP A0A3X8PTV6
A	37	LYS	-	expression tag	UNP A0A3X8PTV6
A	38	ASP	-	expression tag	UNP A0A3X8PTV6
A	39	PRO	-	expression tag	UNP A0A3X8PTV6
B	34	GLY	-	expression tag	UNP A0A3X8PTV6
B	35	SER	-	expression tag	UNP A0A3X8PTV6
B	36	ALA	-	expression tag	UNP A0A3X8PTV6
B	37	LYS	-	expression tag	UNP A0A3X8PTV6
B	38	ASP	-	expression tag	UNP A0A3X8PTV6
B	39	PRO	-	expression tag	UNP A0A3X8PTV6

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	5	Total 5	O 5	0	0
3	B	1	Total 1	O 1	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	119.87Å 119.87Å 274.06Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.79 – 2.95 29.79 – 2.95	Depositor EDS
% Data completeness (in resolution range)	99.0 (29.79-2.95) 99.4 (29.79-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 2.95Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.235 , 0.279 0.244 , 0.286	Depositor DCC
R_{free} test set	2354 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	79.2	Xtriage
Anisotropy	0.505	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 89.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	0.001 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7642	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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](#)

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

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	0.38	0/4225	0.61	1/5770 (0.0%)
1	B	0.34	0/3588	0.57	0/4904
All	All	0.36	0/7813	0.59	1/10674 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	253	GLY	N-CA-C	6.37	117.84	111.21

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	4126	0	3734	84	0
1	B	3508	0	3103	81	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	5	0	0	0	0
3	B	1	0	0	0	0
All	All	7642	0	6837	162	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 (162) 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:307:TYR:CE2	1:B:422:SER:HA	2.19	0.77
1:B:336:CYS:HB3	1:B:356:GLY:H	1.53	0.74
1:A:349:CYS:HB2	1:A:370:VAL:HG23	1.68	0.74
1:A:43:THR:HG22	1:A:209:VAL:HG13	1.70	0.73
1:A:232:ILE:HG22	1:A:466:ILE:HG12	1.71	0.72
1:A:272:PRO:HG2	1:A:298:PRO:HB3	1.70	0.72
1:A:549:TYR:HA	1:A:552:ARG:HD2	1.73	0.71
1:B:323:LEU:HD12	1:B:492:GLN:HB3	1.73	0.70
1:A:291:LEU:HD22	1:A:299:PHE:HB2	1.74	0.70
1:A:307:TYR:H	1:B:422:SER:HB3	1.57	0.69
1:B:254:ALA:HB1	1:B:303:LEU:HD23	1.74	0.68
1:B:309:PRO:HD2	1:B:529:ALA:HB1	1.74	0.68
1:B:241:GLN:HE22	1:B:268:ALA:H	1.38	0.67
1:B:529:ALA:HB3	1:B:554:HIS:HB2	1.75	0.67
1:A:373:TYR:CD1	1:A:429:LEU:HD13	2.30	0.66
1:B:260:VAL:HG13	1:B:450:ALA:HB1	1.78	0.66
1:A:589:LYS:HA	1:A:592:GLN:HE21	1.60	0.65
1:B:61:LEU:HD21	1:B:238:ILE:HG13	1.77	0.65
1:A:193:ILE:HG23	1:A:197:TYR:HD2	1.62	0.65
1:B:241:GLN:NE2	1:B:268:ALA:O	2.31	0.63
1:A:260:VAL:HG12	1:A:454:LYS:HB2	1.82	0.61
1:B:554:HIS:ND1	1:B:575:GLU:HA	2.17	0.59
1:A:465:SER:HA	1:A:474:GLU:HA	1.85	0.57
1:A:209:VAL:HB	1:A:223:SER:H	1.70	0.57
1:A:164:ILE:HD11	1:A:268:ALA:O	2.04	0.57
1:A:305:ASN:HA	1:A:405:PHE:CD2	2.40	0.57
1:B:325:SER:C	1:B:327:ASN:H	2.11	0.57
1:A:266:LEU:HD11	1:A:599:TYR:CE1	2.39	0.57
1:B:322:PHE:CD1	1:B:379:VAL:HG21	2.41	0.56
1:B:260:VAL:HG22	1:B:569:VAL:HG23	1.87	0.56
1:A:260:VAL:HG22	1:A:569:VAL:HG23	1.87	0.56
1:A:320:LEU:HD13	1:A:496:ILE:HG13	1.87	0.55
1:B:450:ALA:HB2	1:B:569:VAL:HG21	1.89	0.55
1:B:55:ASP:HB3	1:B:61:LEU:HD11	1.89	0.55
1:A:380:GLY:O	1:A:384:ILE:HG13	2.07	0.54
1:A:415:LYS:HE3	1:A:422:SER:O	2.07	0.54
1:A:325:SER:C	1:A:327:ASN:H	2.15	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:438:PHE:O	1:B:439:LEU:HG	2.08	0.54
1:A:370:VAL:HA	1:A:373:TYR:CD2	2.42	0.53
1:B:365:LYS:HA	1:B:503:VAL:HG22	1.91	0.53
1:A:415:LYS:HE2	1:A:427:ASP:OD2	2.08	0.53
1:B:593:LYS:O	1:B:597:LEU:HD12	2.09	0.53
1:A:413:GLU:H	1:A:413:GLU:CD	2.17	0.53
1:B:375:GLY:O	1:B:379:VAL:HG23	2.09	0.52
1:A:45:PHE:CD1	1:A:207:THR:HG22	2.45	0.52
1:A:160:LEU:HD21	1:A:242:SER:HB2	1.91	0.52
1:B:381:ILE:HG12	1:B:423:TRP:CZ2	2.45	0.52
1:B:557:LEU:HD13	1:B:587:VAL:HG21	1.92	0.52
1:A:467:GLU:HA	1:A:472:THR:HA	1.92	0.51
1:B:266:LEU:HD11	1:B:599:TYR:CE1	2.45	0.51
1:B:315:LYS:NZ	1:B:369:ASP:OD1	2.39	0.51
1:A:322:PHE:HA	1:A:379:VAL:HG21	1.93	0.51
1:A:457:GLU:HA	1:A:484:ILE:HD13	1.93	0.51
1:A:557:LEU:HD12	1:A:584:GLY:HA2	1.92	0.51
1:A:550:TYR:HA	1:A:576:HIS:CD2	2.46	0.50
1:B:60:LEU:HD13	1:B:63:ILE:HG21	1.93	0.50
1:B:305:ASN:HA	1:B:405:PHE:CD2	2.46	0.50
1:B:424:PHE:O	1:B:427:ASP:HB2	2.12	0.50
1:A:436:GLY:O	1:B:420:ARG:HD2	2.12	0.49
1:B:241:GLN:NE2	1:B:268:ALA:H	2.09	0.49
1:B:498:ASP:HA	1:B:501:TYR:CD2	2.47	0.49
1:A:55:ASP:HB3	1:A:61:LEU:HD11	1.93	0.49
1:A:522:VAL:HG22	1:A:561:ALA:HB2	1.94	0.49
1:B:265:ILE:HG21	1:B:268:ALA:HB2	1.93	0.49
1:B:197:TYR:CZ	1:B:463:LEU:HD22	2.48	0.49
1:B:71:SER:HA	1:B:122:LYS:HA	1.94	0.49
1:B:334:TYR:HD2	1:B:371:TYR:CD2	2.30	0.49
1:B:69:SER:N	1:B:148:ASP:O	2.42	0.48
1:B:557:LEU:HD22	1:B:587:VAL:HG21	1.95	0.48
1:A:240:LEU:HD12	1:A:240:LEU:HA	1.68	0.48
1:A:370:VAL:HA	1:A:373:TYR:HD2	1.79	0.48
1:B:198:ASN:O	1:B:202:GLN:HB2	2.13	0.48
1:B:311:GLY:O	1:B:314:VAL:HG22	2.13	0.48
1:A:544:GLU:HA	1:A:547:PHE:CD1	2.48	0.48
1:A:270:SER:O	1:A:273:GLU:HG3	2.14	0.47
1:A:544:GLU:HA	1:A:547:PHE:CE1	2.49	0.47
1:A:554:HIS:CE1	1:A:576:HIS:ND1	2.82	0.47
1:B:324:ASN:OD1	1:B:324:ASN:N	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:GLU:O	1:B:227:ALA:HB1	2.13	0.47
1:B:182:ALA:O	1:B:186:ASN:N	2.42	0.47
1:A:159:LYS:HA	1:A:274:TYR:O	2.15	0.47
1:A:525:LYS:HE3	1:A:526:THR:O	2.15	0.47
1:B:322:PHE:CE2	1:B:375:GLY:HA3	2.49	0.47
1:B:270:SER:OG	1:B:302:LYS:HG3	2.15	0.46
1:B:465:SER:O	1:B:466:ILE:HD13	2.15	0.46
1:A:554:HIS:CE1	1:A:576:HIS:CE1	3.03	0.46
1:A:530:GLN:HA	1:A:553:SER:HB3	1.98	0.46
1:A:212:VAL:HA	1:A:217:GLN:O	2.14	0.46
1:A:324:ASN:HB3	1:A:488:PHE:CE1	2.50	0.46
1:A:129:TYR:OH	1:A:133:ILE:HD12	2.16	0.46
1:B:549:TYR:CE2	1:B:579:ARG:HG3	2.52	0.45
1:A:47:PRO:HB3	1:A:184:LEU:HB2	1.97	0.45
1:B:71:SER:O	1:B:145:ILE:HA	2.17	0.45
1:A:213:ASN:OD1	1:A:214:ALA:N	2.49	0.45
1:B:447:ARG:O	1:B:451:GLN:HG2	2.16	0.45
1:B:497:ARG:O	1:B:501:TYR:HD2	2.00	0.45
1:A:211:LYS:N	1:A:220:GLU:O	2.44	0.45
1:B:94:THR:O	1:B:98:PRO:HB3	2.17	0.45
1:A:542:VAL:HG12	1:A:543:ASP:H	1.82	0.45
1:B:241:GLN:HE22	1:B:268:ALA:C	2.23	0.45
1:B:498:ASP:O	1:B:501:TYR:HB2	2.17	0.45
1:B:153:ARG:NH2	1:B:194:GLU:OE1	2.49	0.45
1:A:362:HIS:HA	1:A:365:LYS:HB3	1.98	0.44
1:B:197:TYR:CE1	1:B:463:LEU:HD22	2.52	0.44
1:B:336:CYS:SG	1:B:355:HIS:HB2	2.56	0.44
1:B:557:LEU:HD13	1:B:584:GLY:HA2	2.00	0.44
1:A:133:ILE:N	1:A:134:PRO:HD2	2.32	0.44
1:B:66:LEU:O	1:B:129:TYR:HB2	2.17	0.44
1:A:51:GLY:HA3	1:A:230:ASN:O	2.16	0.44
1:B:180:GLU:HG3	1:B:181:ILE:HD12	1.99	0.44
1:B:210:TYR:H	1:B:210:TYR:HD2	1.64	0.44
1:A:315:LYS:HB3	1:A:372:PHE:CZ	2.52	0.44
1:A:253:GLY:HA2	1:A:575:GLU:HG3	2.00	0.44
1:A:308:TYR:HB2	1:A:556:TRP:HZ2	1.83	0.43
1:A:554:HIS:HE1	1:A:576:HIS:CE1	2.36	0.43
1:A:556:TRP:CZ2	1:A:573:LEU:HD13	2.53	0.43
1:B:400:ASP:OD1	1:B:400:ASP:N	2.48	0.43
1:B:493:LEU:O	1:B:497:ARG:HG3	2.18	0.43
1:B:562:PRO:HG2	1:B:565:LYS:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:316:MET:HE2	1:A:316:MET:HB2	1.76	0.43
1:B:313:VAL:O	1:B:448:TYR:OH	2.32	0.43
1:A:557:LEU:HD13	1:A:587:VAL:HG21	2.00	0.43
1:A:493:LEU:N	1:A:494:PRO:CD	2.82	0.42
1:B:254:ALA:HB1	1:B:303:LEU:CD2	2.46	0.42
1:A:238:ILE:HD13	1:A:238:ILE:HA	1.80	0.42
1:A:302:LYS:HE3	1:A:575:GLU:OE1	2.19	0.42
1:B:244:LEU:HD23	1:B:244:LEU:HA	1.66	0.42
1:A:471:THR:HG23	1:A:471:THR:O	2.20	0.42
1:B:410:PRO:HD3	1:B:438:PHE:HA	2.01	0.42
1:A:171:ALA:HB2	1:A:186:ASN:HA	2.00	0.42
1:A:365:LYS:NZ	1:A:506:GLU:HB3	2.34	0.42
1:A:589:LYS:HA	1:A:592:GLN:NE2	2.32	0.42
1:B:325:SER:C	1:B:327:ASN:N	2.78	0.42
1:B:159:LYS:HG2	1:B:273:GLU:HB2	2.02	0.42
1:A:320:LEU:HD12	1:A:320:LEU:HA	1.79	0.42
1:B:54:THR:HG22	1:B:60:LEU:HD23	2.02	0.41
1:B:60:LEU:HD23	1:B:60:LEU:HA	1.84	0.41
1:B:265:ILE:HD12	1:B:462:PHE:CZ	2.55	0.41
1:A:263:GLY:HA3	1:A:459:ILE:O	2.20	0.41
1:A:352:ARG:H	1:A:352:ARG:HG2	1.63	0.41
1:A:201:LEU:HD21	1:A:232:ILE:CD1	2.51	0.41
1:B:574:LEU:HD12	1:B:578:GLY:CA	2.51	0.41
1:A:305:ASN:HA	1:A:405:PHE:HD2	1.83	0.41
1:A:537:THR:O	1:A:540:ASN:N	2.54	0.41
1:A:314:VAL:HG11	1:A:438:PHE:CE1	2.56	0.41
1:A:320:LEU:CD2	1:A:489:GLU:HG2	2.50	0.41
1:A:591:TYR:O	1:A:595:ILE:HG13	2.21	0.41
1:A:373:TYR:CG	1:A:429:LEU:HD13	2.56	0.41
1:A:381:ILE:HG12	1:A:423:TRP:CZ2	2.55	0.41
1:A:309:PRO:HD2	1:A:529:ALA:HB1	2.02	0.40
1:B:260:VAL:HG23	1:B:567:LYS:O	2.20	0.40
1:B:452:ILE:HD13	1:B:452:ILE:HA	1.87	0.40
1:B:55:ASP:HB2	1:B:236:ILE:O	2.21	0.40
1:B:56:ARG:HG2	1:B:236:ILE:O	2.21	0.40
1:B:213:ASN:OD1	1:B:216:ASN:N	2.54	0.40
1:B:493:LEU:HD23	1:B:493:LEU:HA	1.83	0.40
1:B:568:TYR:CE2	1:B:599:TYR:HB3	2.56	0.40
1:B:594:MET:HE2	1:B:594:MET:HB3	1.86	0.40
1:A:304:ILE:HB	1:A:404:GLU:HG3	2.03	0.40
1:A:87:ASP:OD1	1:A:87:ASP:N	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:450:ALA:HB2	1:A:569:VAL:HG21	2.04	0.40
1:B:296:ASP:OD1	1:B:296:ASP:N	2.54	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	554/568 (98%)	509 (92%)	45 (8%)	0	100	100
1	B	478/568 (84%)	449 (94%)	29 (6%)	0	100	100
All	All	1032/1136 (91%)	958 (93%)	74 (7%)	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	398/491 (81%)	338 (85%)	60 (15%)	3	8
1	B	317/491 (65%)	281 (89%)	36 (11%)	5	16
All	All	715/982 (73%)	619 (87%)	96 (13%)	4	11

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	VAL
1	A	64	ASN
1	A	71	SER
1	A	81	SER
1	A	82	ASN
1	A	87	ASP
1	A	93	LEU
1	A	94	THR
1	A	96	LEU
1	A	118	GLN
1	A	121	ILE
1	A	132	ILE
1	A	133	ILE
1	A	210	TYR
1	A	221	GLN
1	A	230	ASN
1	A	232	ILE
1	A	238	ILE
1	A	291	LEU
1	A	294	SER
1	A	295	LEU
1	A	304	ILE
1	A	307	TYR
1	A	312	SER
1	A	313	VAL
1	A	329	SER
1	A	332	THR
1	A	348	ARG
1	A	352	ARG
1	A	379	VAL
1	A	384	ILE
1	A	385	SER
1	A	388	LEU
1	A	391	ILE
1	A	409	LEU
1	A	413	GLU
1	A	415	LYS
1	A	416	MET
1	A	425	GLN
1	A	429	LEU
1	A	433	ILE
1	A	472	THR
1	A	473	ILE

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Mol	Chain	Res	Type
1	A	483	GLU
1	A	484	ILE
1	A	489	GLU
1	A	507	GLN
1	A	511	SER
1	A	525	LYS
1	A	531	VAL
1	A	532	VAL
1	A	535	SER
1	A	537	THR
1	A	542	VAL
1	A	553	SER
1	A	557	LEU
1	A	564	SER
1	A	573	LEU
1	A	582	THR
1	A	583	SER
1	B	49	VAL
1	B	64	ASN
1	B	66	LEU
1	B	90	LEU
1	B	132	ILE
1	B	176	VAL
1	B	188	THR
1	B	202	GLN
1	B	210	TYR
1	B	220	GLU
1	B	230	ASN
1	B	238	ILE
1	B	244	LEU
1	B	265	ILE
1	B	266	LEU
1	B	276	LEU
1	B	296	ASP
1	B	307	TYR
1	B	313	VAL
1	B	324	ASN
1	B	328	ILE
1	B	331	SER
1	B	386	GLU
1	B	389	SER
1	B	411	SER

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Mol	Chain	Res	Type
1	B	421	GLN
1	B	425	GLN
1	B	445	ILE
1	B	451	GLN
1	B	484	ILE
1	B	505	ASN
1	B	564	SER
1	B	573	LEU
1	B	574	LEU
1	B	582	THR
1	B	597	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	202	GLN
1	A	217	GLN
1	A	297	HIS
1	A	305	ASN
1	A	421	GLN
1	A	425	GLN
1	A	576	HIS
1	A	592	GLN
1	B	163	HIS
1	B	186	ASN
1	B	221	GLN
1	B	230	ASN
1	B	241	GLN
1	B	378	GLN
1	B	425	GLN
1	B	530	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 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

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	558/568 (98%)	0.02	17 (3%)	52 45	52, 93, 147, 178	0
1	B	490/568 (86%)	0.22	23 (4%)	36 30	66, 109, 147, 182	0
All	All	1048/1136 (92%)	0.11	40 (3%)	44 36	52, 102, 147, 182	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	337	ASN	6.5
1	A	39	PRO	5.2
1	A	218	GLU	4.5
1	A	470	ASN	4.4
1	B	172	ASN	4.2
1	A	175	ASP	4.0
1	B	391	ILE	3.9
1	B	175	ASP	3.8
1	B	132	ILE	3.7
1	A	469	ASN	3.7
1	B	467	GLU	3.7
1	B	39	PRO	3.5
1	B	531	VAL	3.5
1	B	136	TYR	3.3
1	B	236	ILE	3.3
1	A	283	ILE	3.2
1	B	210	TYR	2.9
1	B	72	ILE	2.9
1	A	86	LEU	2.8
1	B	294	SER	2.7
1	B	274	TYR	2.6
1	B	100	LEU	2.6
1	A	475	ASN	2.6
1	A	468	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	221	GLN	2.5
1	A	46	LEU	2.5
1	B	229	SER	2.4
1	A	85	ILE	2.4
1	B	144	THR	2.3
1	A	285	PHE	2.3
1	A	41	ILE	2.3
1	B	86	LEU	2.3
1	A	115	TYR	2.3
1	A	145	ILE	2.2
1	B	329	SER	2.2
1	A	467	GLU	2.2
1	B	204	GLU	2.2
1	A	229	SER	2.2
1	B	135	HIS	2.1
1	B	125	ASP	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](#)

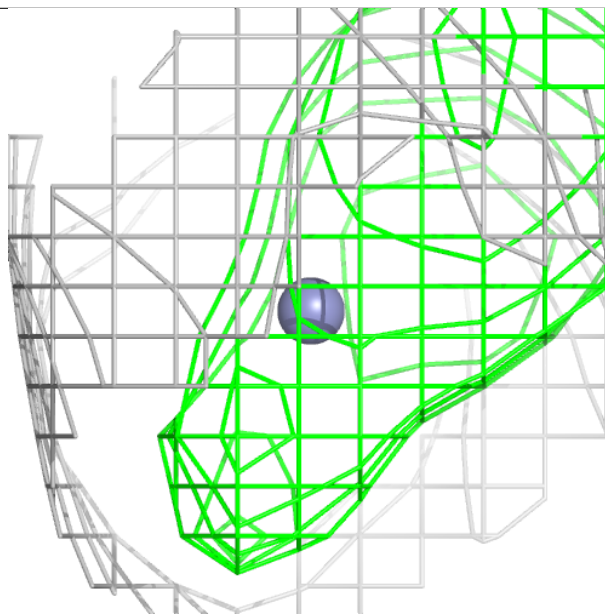
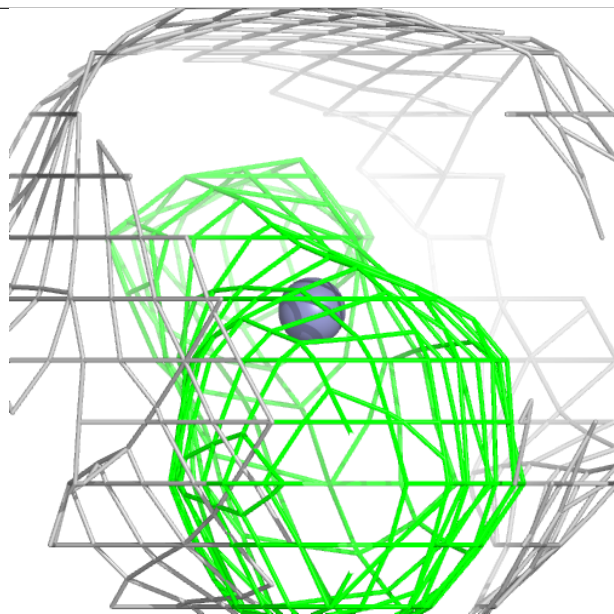
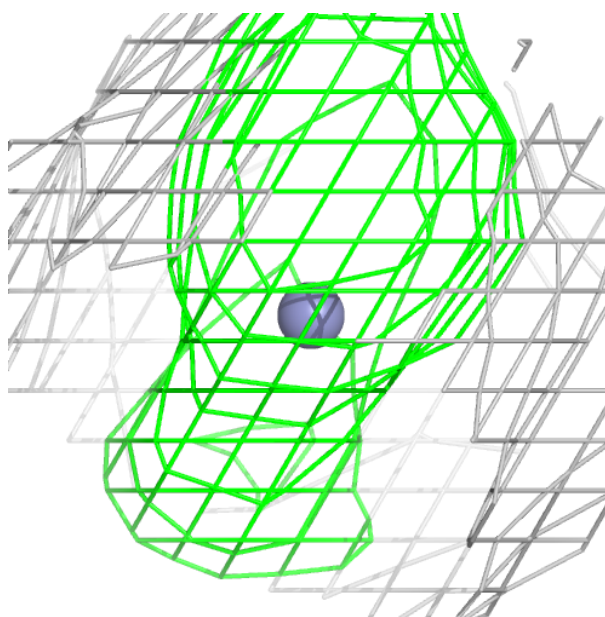
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

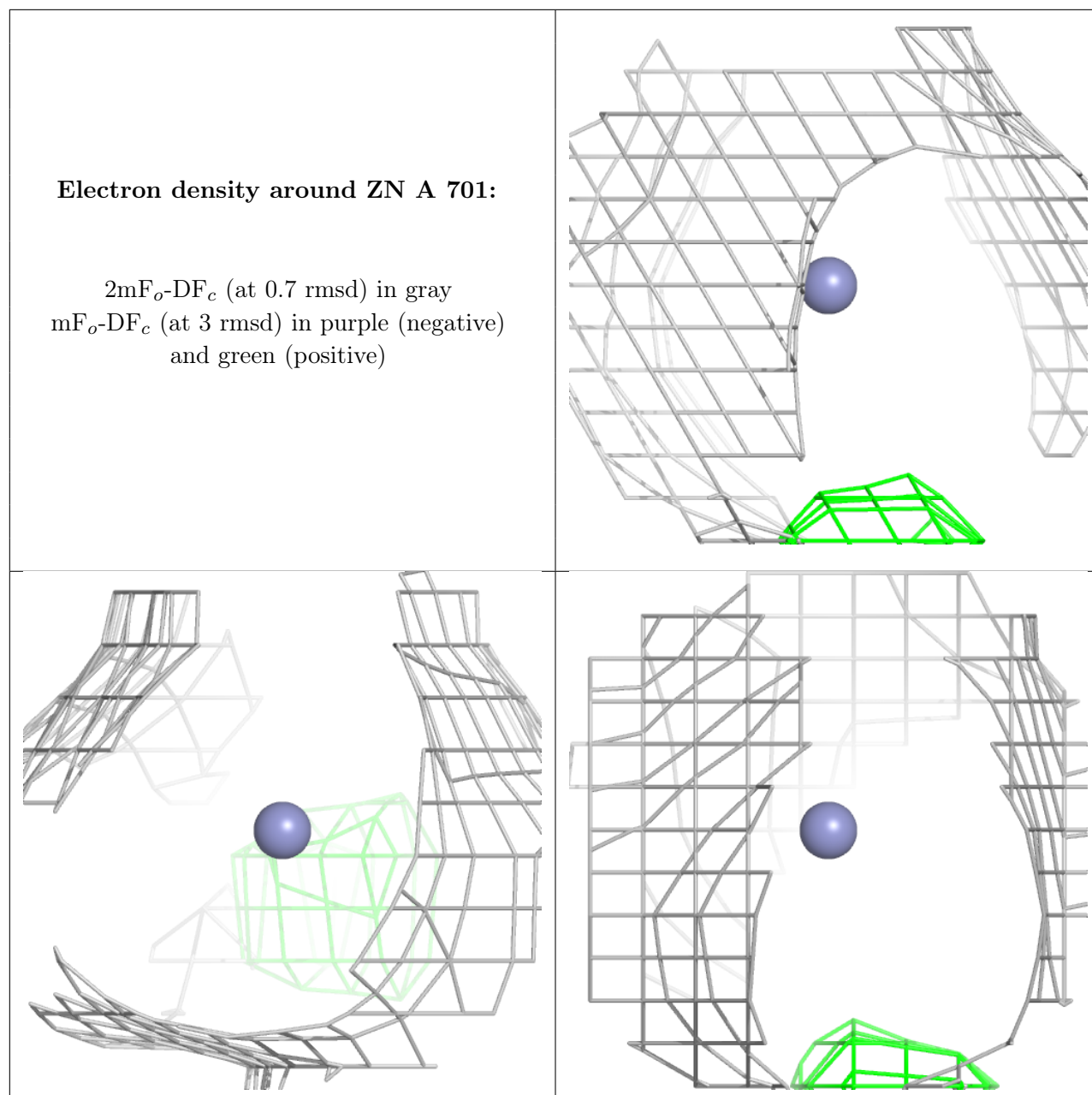
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	B	701	1/1	0.97	0.10	108,108,108,108	0
2	ZN	A	701	1/1	1.00	0.02	68,68,68,68	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around ZN B 701:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers ⓘ

There are no such residues in this entry.