



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:01 PM UTC

PDB ID : 4MNA / pdb_00004mna
Title : Crystal structure of the free FLS2 ectodomains
Authors : Chai, J.; Han, Z.; Sun, Y.
Deposited on : 2013-09-10
Resolution : 4.00 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
Mogul : 2022.3.0, CSD as543be (2022)
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

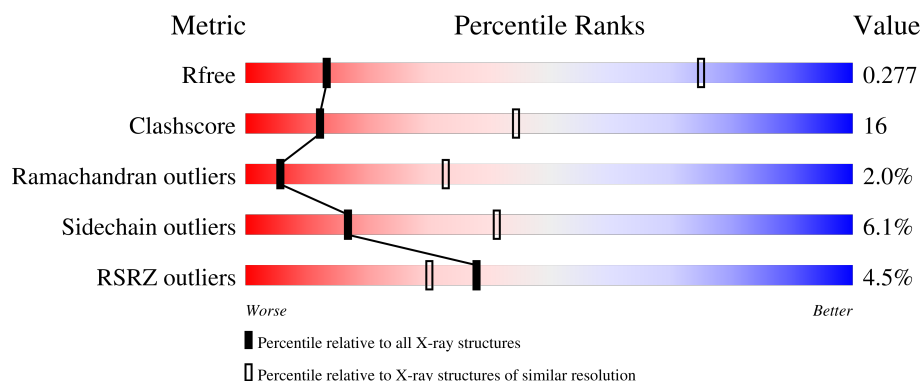
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.00 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	1082 (4.20-3.80)
Clashscore	190562	1129 (4.20-3.80)
Ramachandran outliers	187476	1064 (4.20-3.80)
Sidechain outliers	187428	1055 (4.20-3.80)
RSRZ outliers	180081	1082 (4.20-3.80)

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	662	

2 Entry composition [i](#)

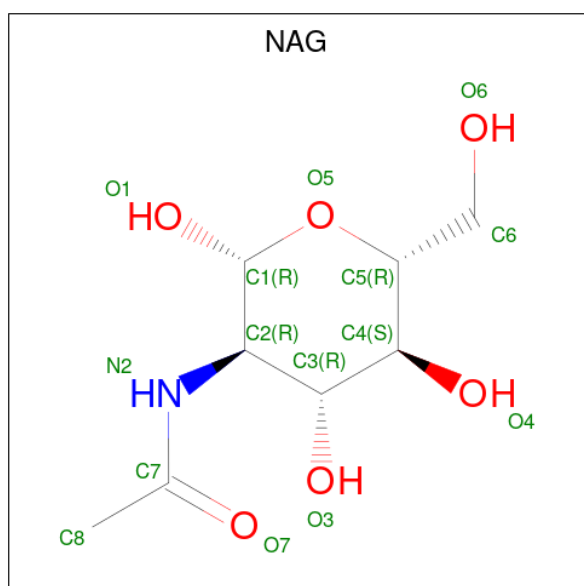
There are 3 unique types of molecules in this entry. The entry contains 4964 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 LRR receptor-like serine/threonine-protein kinase FLS2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	639	Total	C	N	O	S	0	0	0
			4890	3090	819	962	19			

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total 4	Zn 4	0	0

- Molecule 1: LRR receptor-like serine/threonine-protein kinase FLS2



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	189.54Å 189.54Å 117.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.73 – 4.00 29.73 – 4.00	Depositor EDS
% Data completeness (in resolution range)	98.9 (29.73-4.00) 98.5 (29.73-4.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.74 (at 3.98Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.249 , 0.279 0.253 , 0.277	Depositor DCC
R_{free} test set	1057 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	94.2	Xtriage
Anisotropy	0.681	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 78.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.031 for -h,-k,l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4964	wwPDB-VP
Average B, all atoms (Å ²)	105.0	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: NAG, 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	0.52	0/4969	0.92	8/6743 (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	A	119	LEU	N-CA-C	-7.75	98.84	110.24
1	A	431	MET	N-CA-C	7.34	122.40	113.38
1	A	89	SER	CA-C-N	5.75	125.37	119.56
1	A	89	SER	C-N-CA	5.75	125.37	119.56
1	A	243	GLN	N-CA-C	-5.71	104.31	112.13
1	A	754	ALA	CB-CA-C	-5.60	110.09	116.54
1	A	447	ILE	CA-C-N	5.29	125.55	119.83
1	A	447	ILE	C-N-CA	5.29	125.55	119.83

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	4890	0	4944	155	1
2	A	70	0	65	3	0
3	A	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4964	0	5009	155	1

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 16.

All (155) 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:245:LEU:HD12	1:A:246:VAL:N	1.37	1.38
1:A:245:LEU:HD12	1:A:245:LEU:C	1.52	1.30
1:A:246:VAL:O	1:A:247:LEU:CD2	1.79	1.29
1:A:246:VAL:O	1:A:247:LEU:HD23	1.03	1.20
1:A:433:LEU:HD12	1:A:433:LEU:N	1.52	1.20
1:A:433:LEU:H	1:A:433:LEU:CD1	1.62	1.11
1:A:428:PHE:HA	1:A:431:MET:HE3	1.36	1.06
1:A:121:GLU:OE2	1:A:121:GLU:HA	1.45	1.06
1:A:411:LYS:HE3	2:A:906:NAG:O6	1.59	1.02
1:A:245:LEU:CD1	1:A:246:VAL:N	2.30	0.95
1:A:245:LEU:C	1:A:245:LEU:CD1	2.30	0.89
1:A:99:GLN:O	1:A:242:LEU:CD1	2.21	0.88
1:A:428:PHE:HA	1:A:431:MET:CE	2.07	0.84
1:A:603:THR:HG23	1:A:629:TYR:HB3	1.61	0.83
1:A:532:LEU:HD11	1:A:534:MET:HE3	1.59	0.82
1:A:120:THR:O	1:A:120:THR:HG23	1.80	0.81
1:A:433:LEU:HD12	1:A:433:LEU:H	0.70	0.80
1:A:120:THR:O	1:A:120:THR:CG2	2.30	0.77
1:A:245:LEU:HD12	1:A:246:VAL:CA	2.16	0.76
1:A:246:VAL:O	1:A:247:LEU:CG	2.37	0.73
1:A:99:GLN:O	1:A:242:LEU:HD12	1.90	0.71
1:A:121:GLU:OE2	1:A:121:GLU:CA	2.30	0.71
1:A:411:LYS:HE3	2:A:906:NAG:HO6	1.54	0.69
1:A:98:LEU:HB3	1:A:119:LEU:CD1	2.23	0.68
1:A:436:ILE:HD12	1:A:451:ILE:HD11	1.76	0.68
1:A:436:ILE:HG13	1:A:457:LEU:HD11	1.75	0.68
1:A:377:PRO:HG2	1:A:380:LEU:HD23	1.76	0.68
1:A:311:LEU:HD12	1:A:314:LEU:HD11	1.74	0.67
1:A:333:GLY:O	1:A:358:ASN:ND2	2.25	0.67
1:A:285:GLY:HA2	1:A:311:LEU:HD21	1.77	0.67
1:A:99:GLN:HA	1:A:121:GLU:O	1.95	0.66
1:A:271:LEU:HB2	1:A:295:ILE:HG22	1.77	0.66
1:A:433:LEU:N	1:A:433:LEU:CD1	2.34	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:PHE:HA	1:A:335:LEU:HD21	1.78	0.65
1:A:118:LYS:O	1:A:119:LEU:C	2.38	0.65
1:A:697:MET:O	1:A:699:MET:N	2.29	0.65
1:A:99:GLN:O	1:A:242:LEU:HD13	1.98	0.64
1:A:559:SER:HB2	1:A:583:GLN:HG3	1.78	0.64
1:A:360:ARG:HH21	1:A:382:LEU:HD13	1.63	0.63
1:A:779:ASN:O	1:A:781:ASP:N	2.33	0.62
1:A:266:LEU:HD23	1:A:287:LEU:HD13	1.82	0.62
1:A:483:ILE:HG22	1:A:507:ILE:HB	1.82	0.61
1:A:359:LEU:H	1:A:359:LEU:HD23	1.65	0.61
1:A:242:LEU:HD12	1:A:243:GLN:N	2.15	0.60
1:A:473:PRO:HB3	1:A:498:GLU:HG3	1.84	0.59
1:A:246:VAL:HG12	1:A:247:LEU:N	2.16	0.59
1:A:645:LEU:HD23	1:A:669:LEU:HD22	1.85	0.59
1:A:38:LYS:HG3	1:A:49:LEU:HD13	1.83	0.59
1:A:415:LEU:HB2	1:A:438:ILE:HG22	1.85	0.58
1:A:753:LEU:O	1:A:756:ASN:ND2	2.37	0.58
1:A:268:GLN:HG3	1:A:292:ALA:HB3	1.86	0.57
1:A:582:LEU:HB2	1:A:606:ILE:HD12	1.88	0.56
1:A:243:GLN:O	1:A:267:VAL:HG12	2.05	0.56
1:A:81:GLU:HG3	1:A:105:SER:HB3	1.88	0.55
1:A:246:VAL:CG1	1:A:247:LEU:N	2.69	0.55
1:A:518:ARG:NH1	1:A:542:PRO:HD2	2.21	0.55
1:A:607:SER:O	1:A:609:ASN:ND2	2.40	0.55
1:A:678:LEU:HD22	1:A:697:MET:HE2	1.89	0.55
1:A:433:LEU:CD1	1:A:454:CYS:SG	2.96	0.55
1:A:433:LEU:HD22	1:A:436:ILE:HD11	1.88	0.54
1:A:719:GLY:HA3	1:A:741:SER:HB3	1.88	0.54
1:A:52:TRP:CH2	1:A:66:ILE:HD11	2.42	0.54
1:A:392:HIS:HD2	1:A:393:ASP:OD1	1.89	0.54
1:A:433:LEU:HD11	1:A:454:CYS:SG	2.48	0.54
1:A:242:LEU:HD12	1:A:243:GLN:H	1.72	0.53
1:A:370:ASN:HB3	1:A:372:ILE:HG13	1.91	0.53
1:A:606:ILE:HG22	1:A:632:PHE:HD1	1.74	0.53
1:A:626:MET:HE1	1:A:630:LEU:HB2	1.91	0.53
1:A:411:LYS:HA	1:A:433:LEU:HA	1.91	0.52
1:A:245:LEU:HD11	1:A:247:LEU:HG	1.91	0.52
1:A:246:VAL:C	1:A:247:LEU:CG	2.83	0.52
1:A:280:ILE:HG22	1:A:305:PRO:HG2	1.92	0.51
1:A:763:PRO:HG2	1:A:768:PHE:CZ	2.46	0.51
1:A:296:TYR:O	1:A:298:ASN:ND2	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:SER:HA	1:A:310:ARG:HH11	1.76	0.51
1:A:582:LEU:O	1:A:585:ASN:ND2	2.44	0.50
1:A:758:LEU:HD12	1:A:779:ASN:HD21	1.77	0.49
1:A:294:ARG:HH21	1:A:316:HIS:CD2	2.31	0.49
1:A:427:GLY:N	1:A:450:ASP:OD2	2.46	0.48
1:A:411:LYS:CE	2:A:906:NAG:O6	2.48	0.48
1:A:477:LYS:O	1:A:479:GLN:HG3	2.13	0.48
1:A:436:ILE:CD1	1:A:451:ILE:HD11	2.42	0.48
1:A:246:VAL:C	1:A:247:LEU:HG	2.39	0.48
1:A:625:ASN:OD1	1:A:625:ASN:N	2.41	0.48
1:A:417:HIS:H	1:A:440:ARG:HB3	1.79	0.47
1:A:285:GLY:HA3	1:A:307:SER:HB2	1.95	0.46
1:A:98:LEU:HD23	1:A:119:LEU:HD11	1.97	0.46
1:A:424:ILE:HD11	1:A:438:ILE:HD13	1.97	0.46
1:A:539:LEU:HD23	1:A:539:LEU:HA	1.78	0.46
1:A:629:TYR:CE2	1:A:631:ASN:HB2	2.51	0.46
1:A:60:HIS:HA	1:A:63:TRP:CE2	2.51	0.46
1:A:739:PRO:HB2	1:A:742:LEU:HD23	1.98	0.46
1:A:87:VAL:HG23	1:A:110:GLY:HA3	1.98	0.45
1:A:48:VAL:HG21	1:A:82:LYS:HD2	1.98	0.45
1:A:92:ILE:O	1:A:119:LEU:HD21	2.17	0.45
1:A:707:ARG:NH1	1:A:730:SER:OG	2.49	0.45
1:A:386:LEU:HD23	1:A:386:LEU:HA	1.83	0.45
1:A:242:LEU:HD13	1:A:242:LEU:HA	1.42	0.45
1:A:449:ASP:OD1	1:A:472:LYS:HD2	2.17	0.45
1:A:356:ILE:HD12	1:A:356:ILE:HA	1.87	0.45
1:A:48:VAL:HG21	1:A:82:LYS:HB3	1.99	0.45
1:A:768:PHE:HA	1:A:771:ILE:HB	1.98	0.45
1:A:41:ILE:HD12	1:A:49:LEU:HD11	2.00	0.44
1:A:633:SER:O	1:A:635:ASN:ND2	2.50	0.44
1:A:777:MET:HA	1:A:778:GLY:HA2	1.46	0.44
1:A:330:GLU:OE1	1:A:354:GLN:N	2.49	0.44
1:A:724:LEU:HD22	1:A:748:LEU:HD13	1.99	0.44
1:A:626:MET:HE2	1:A:648:LEU:HD22	1.99	0.44
1:A:309:PHE:HE1	1:A:332:ILE:HG23	1.83	0.44
1:A:523:MET:HE3	1:A:523:MET:HB2	1.75	0.44
1:A:502:LEU:O	1:A:504:ASP:N	2.51	0.44
1:A:626:MET:HE3	1:A:651:VAL:HG21	2.00	0.43
1:A:287:LEU:HB2	1:A:290:LEU:HD12	2.00	0.43
1:A:574:LEU:C	1:A:576:SER:H	2.26	0.43
1:A:658:ASN:H	1:A:682:GLN:HB3	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:487:SER:O	1:A:489:ASN:ND2	2.52	0.43
1:A:528:LEU:HD13	1:A:528:LEU:HA	1.77	0.43
1:A:602:ASN:O	1:A:628:LEU:HB2	2.18	0.43
1:A:623:LEU:HD13	1:A:626:MET:SD	2.58	0.43
1:A:758:LEU:HD12	1:A:779:ASN:ND2	2.34	0.43
1:A:346:ASN:HB3	1:A:347:ASN:H	1.59	0.43
1:A:561:ASN:HB3	1:A:562:LYS:H	1.69	0.43
1:A:433:LEU:HD22	1:A:436:ILE:CD1	2.48	0.43
1:A:261:GLY:HA2	1:A:287:LEU:HD21	2.00	0.43
1:A:268:GLN:NE2	1:A:270:GLU:OE1	2.42	0.42
1:A:626:MET:CE	1:A:630:LEU:HB2	2.49	0.42
1:A:495:ILE:HD11	1:A:510:LEU:HD11	2.01	0.42
1:A:381:GLY:HA3	1:A:406:ASN:HB2	2.02	0.42
1:A:340:VAL:HG22	1:A:364:VAL:HB	2.02	0.42
1:A:738:ILE:HA	1:A:739:PRO:HD3	1.90	0.42
1:A:629:TYR:OH	1:A:631:ASN:ND2	2.53	0.42
1:A:545:GLU:HG2	1:A:570:LEU:HD13	2.02	0.42
1:A:98:LEU:HB3	1:A:119:LEU:HD11	1.99	0.42
1:A:678:LEU:HD22	1:A:697:MET:CE	2.50	0.41
1:A:302:SER:OG	1:A:303:SER:N	2.41	0.41
1:A:507:ILE:HG23	1:A:509:TYR:CE1	2.56	0.41
1:A:592:PRO:HG2	1:A:595:LEU:HG	2.01	0.41
1:A:574:LEU:HD12	1:A:577:LEU:HD11	2.02	0.41
1:A:111:LYS:HE2	1:A:111:LYS:HB3	1.83	0.41
1:A:89:SER:HA	1:A:90:PRO:HD3	1.82	0.41
1:A:245:LEU:HB3	1:A:266:LEU:HD11	2.03	0.41
1:A:310:ARG:O	1:A:312:THR:HG23	2.21	0.41
1:A:523:MET:HG3	1:A:526:LEU:HD12	2.02	0.41
1:A:529:LEU:HD23	1:A:550:MET:HE1	2.02	0.41
1:A:672:CYS:HB2	1:A:675:VAL:HG23	2.03	0.41
1:A:531:GLY:HA2	1:A:555:VAL:HG13	2.03	0.41
1:A:646:GLY:HA2	1:A:669:LEU:HA	2.03	0.41
1:A:738:ILE:HA	1:A:738:ILE:HD13	1.83	0.41
1:A:383:LEU:H	1:A:383:LEU:HG	1.65	0.40
1:A:433:LEU:HD22	1:A:436:ILE:CG1	2.51	0.40
1:A:547:MET:HE2	1:A:547:MET:HB3	1.94	0.40
1:A:360:ARG:NH2	1:A:382:LEU:HD13	2.32	0.40
1:A:304:ILE:HA	1:A:305:PRO:HD3	1.95	0.40
1:A:666:PRO:HG2	1:A:669:LEU:HD23	2.03	0.40
1:A:478:LEU:HD13	1:A:481:LEU:HD22	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:GLU:OE1	1:A:497:ARG:NH1[3_655]	2.19	0.01

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	637/662 (96%)	543 (85%)	81 (13%)	13 (2%)	6	33

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	698	ASP
1	A	302	SER
1	A	407	CYS
1	A	503	LYS
1	A	311	LEU
1	A	455	SER
1	A	780	THR
1	A	575	GLU
1	A	738	ILE
1	A	743	ALA
1	A	247	LEU
1	A	744	ASN
1	A	288	VAL

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	571/594 (96%)	536 (94%)	35 (6%)	17	41

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	LEU
1	A	118	LYS
1	A	119	LEU
1	A	120	THR
1	A	121	GLU
1	A	242	LEU
1	A	244	SER
1	A	245	LEU
1	A	289	GLN
1	A	290	LEU
1	A	330	GLU
1	A	332	ILE
1	A	341	LEU
1	A	356	ILE
1	A	359	LEU
1	A	380	LEU
1	A	383	LEU
1	A	402	SER
1	A	408	THR
1	A	433	LEU
1	A	434	THR
1	A	438	ILE
1	A	451	ILE
1	A	478	LEU
1	A	490	SER
1	A	507	ILE
1	A	510	LEU
1	A	555	VAL
1	A	598	LEU
1	A	625	ASN
1	A	627	GLN
1	A	693	VAL
1	A	700	ILE
1	A	762	VAL
1	A	764	GLU

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	286	ASN
1	A	289	GLN
1	A	347	ASN
1	A	417	HIS
1	A	761	HIS
1	A	772	ASN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	A	904	1	14,14,15	0.51	0	17,19,21	1.12	1 (5%)
2	NAG	A	901	1	14,14,15	0.57	0	17,19,21	1.05	1 (5%)
2	NAG	A	909	1	14,14,15	0.52	0	17,19,21	0.76	0
2	NAG	A	906	1	14,14,15	0.54	0	17,19,21	1.39	3 (17%)
2	NAG	A	902	1	14,14,15	0.66	1 (7%)	17,19,21	1.60	3 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	904	1	-	3/6/23/26	0/1/1/1
2	NAG	A	901	1	-	4/6/23/26	0/1/1/1
2	NAG	A	909	1	-	0/6/23/26	0/1/1/1
2	NAG	A	906	1	-	4/6/23/26	0/1/1/1
2	NAG	A	902	1	-	3/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	902	NAG	C1-C2	2.02	1.55	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	906	NAG	C1-O5-C5	3.93	117.45	112.19
2	A	902	NAG	C2-N2-C7	3.83	128.03	122.90
2	A	902	NAG	C1-O5-C5	3.17	116.43	112.19
2	A	904	NAG	C1-O5-C5	2.78	115.92	112.19
2	A	906	NAG	O5-C1-C2	-2.50	107.42	111.29
2	A	902	NAG	C3-C4-C5	-2.17	106.29	110.23
2	A	901	NAG	C1-C2-N2	-2.17	107.01	110.43
2	A	906	NAG	C1-C2-N2	2.07	113.69	110.43

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	901	NAG	C8-C7-N2-C2
2	A	901	NAG	O7-C7-N2-C2
2	A	904	NAG	C8-C7-N2-C2
2	A	904	NAG	O7-C7-N2-C2
2	A	906	NAG	C8-C7-N2-C2
2	A	906	NAG	O7-C7-N2-C2
2	A	902	NAG	O7-C7-N2-C2
2	A	902	NAG	C8-C7-N2-C2
2	A	901	NAG	C4-C5-C6-O6
2	A	901	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	A	906	NAG	C4-C5-C6-O6
2	A	902	NAG	O5-C5-C6-O6
2	A	904	NAG	O5-C5-C6-O6
2	A	906	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	906	NAG	3	0

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	639/662 (96%)	0.17	29 (4%) 38 30	46, 99, 181, 324	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	27	PHE	5.6
1	A	28	GLU	5.4
1	A	764	GLU	5.2
1	A	765	SER	4.6
1	A	55	ILE	4.0
1	A	379	ASP	3.9
1	A	382	LEU	3.8
1	A	72	GLY	3.6
1	A	750	HIS	3.6
1	A	58	LEU	3.5
1	A	435	PHE	3.4
1	A	26	SER	3.3
1	A	649	GLU	3.1
1	A	761	HIS	3.0
1	A	763	PRO	3.0
1	A	59	ARG	2.9
1	A	740	GLU	2.8
1	A	698	ASP	2.6
1	A	684	ASN	2.6
1	A	313	GLN	2.5
1	A	96	THR	2.3
1	A	68	CYS	2.3
1	A	430	ARG	2.2
1	A	562	LYS	2.2
1	A	752	LYS	2.2
1	A	319	LEU	2.2
1	A	784	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	310	ARG	2.1
1	A	31	ILE	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	NAG	A	904	14/15	0.74	0.20	112,114,115,117	0
3	ZN	A	908	1/1	0.81	0.10	267,267,267,267	0
2	NAG	A	906	14/15	0.85	0.13	92,94,97,97	0
2	NAG	A	902	14/15	0.85	0.10	82,82,83,83	0
2	NAG	A	909	14/15	0.87	0.15	20,20,20,20	0
3	ZN	A	907	1/1	0.90	0.12	143,143,143,143	0
2	NAG	A	901	14/15	0.92	0.10	108,110,111,112	0
3	ZN	A	903	1/1	0.96	0.19	50,50,50,50	0
3	ZN	A	905	1/1	0.97	0.17	93,93,93,93	0

6.5 Other polymers [i](#)

There are no such residues in this entry.