



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:27 AM UTC

PDB ID : 2HFZ / pdb_00002hfz
Title : Crystal structure of RNA dependent RNA polymerase domain from West Nile virus
Authors : Egloff, M.P.; Malet, H.; Marseilles Structural Genomics Program @ AFMB (MSGP)
Deposited on : 2006-06-26
Resolution : 3.00 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
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

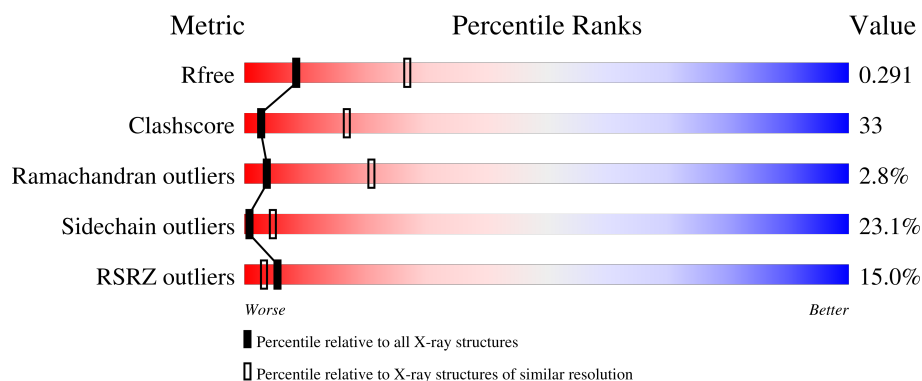
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	639	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4929 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 RNA-directed RNA polymerase(NS5).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	608	Total	C	N	O	S	0	0	0
			4915	3083	888	915	29			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	267	HIS	-	expression tag	UNP P14335
A	268	HIS	-	expression tag	UNP P14335
A	269	HIS	-	expression tag	UNP P14335
A	270	HIS	-	expression tag	UNP P14335
A	271	HIS	-	expression tag	UNP P14335
A	272	HIS	-	expression tag	UNP P14335
A	273	LYS	-	SEE REMARK 999	UNP P14335

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total	O	0	0
			12	12		

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	73.33Å 103.45Å 190.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	95.35 – 3.00 95.04 – 3.00	Depositor EDS
% Data completeness (in resolution range)	96.1 (95.35-3.00) 96.2 (95.04-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.99Å)	Xtriage
Refinement program	BUSTER-TNT 1.3.1	Depositor
R, R_{free}	0.260 , 0.265 0.276 , 0.291	Depositor DCC
R_{free} test set	1440 reflections (10.05%)	wwPDB-VP
Wilson B-factor (Å ²)	48.3	Xtriage
Anisotropy	0.528	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 55.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	4929	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.42% 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: ZN, MG

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	0.45	3/5029 (0.1%)	0.89	21/6807 (0.3%)

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	1	0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	799	THR	CA-CB	5.97	1.60	1.53
1	A	641	THR	CA-C	-5.70	1.45	1.52
1	A	733	CYS	CB-SG	5.63	1.99	1.81

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	420	GLN	N-CA-C	-9.71	101.37	113.02
1	A	845	GLY	N-CA-C	-8.04	101.77	112.14
1	A	361	VAL	N-CA-C	7.39	119.11	109.58
1	A	898	THR	N-CA-C	7.25	126.23	110.80
1	A	852	CYS	N-CA-C	6.88	120.56	112.72
1	A	501	GLY	N-CA-C	-6.84	103.78	112.54
1	A	811	THR	N-CA-C	-6.69	104.34	113.30
1	A	880	GLY	N-CA-C	6.60	121.56	112.52
1	A	841	VAL	CA-C-N	6.41	127.85	119.84
1	A	841	VAL	C-N-CA	6.41	127.85	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	444	HIS	CB-CA-C	-6.37	97.39	110.38
1	A	798	THR	N-CA-C	-6.01	105.19	113.18
1	A	706	TYR	N-CA-C	-5.79	106.25	113.55
1	A	751	TRP	N-CA-C	-5.76	106.42	113.50
1	A	633	GLY	CA-C-N	5.57	125.93	119.47
1	A	633	GLY	C-N-CA	5.57	125.93	119.47
1	A	345	MET	N-CA-C	5.26	116.73	109.15
1	A	852	CYS	N-CA-CB	5.25	118.52	110.70
1	A	844	SER	N-CA-C	-5.12	104.11	110.41
1	A	678	ARG	N-CA-C	-5.03	106.41	112.54
1	A	298	GLU	N-CA-C	-5.02	107.50	113.38

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	898	THR	CA

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	4915	0	4784	321	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	12	0	0	1	0
All	All	4929	0	4784	321	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 33.

All (321) 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:321:SER:HB3	1:A:345:MET:HA	1.21	1.12

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:GLY:HA3	1:A:349:THR:HA	1.29	1.11
1:A:581:LYS:HB3	1:A:593:MET:HE2	1.14	1.10
1:A:867:ASN:HD22	1:A:867:ASN:N	1.56	0.99
1:A:279:ILE:HG22	1:A:283:ILE:HD11	1.46	0.97
1:A:293:THR:HG21	1:A:312:GLU:HB2	1.49	0.94
1:A:846:LYS:HZ1	1:A:899:LEU:HD22	1.32	0.93
1:A:899:LEU:HD23	1:A:899:LEU:H	1.33	0.92
1:A:334:TRP:HA	1:A:337:ILE:HD13	1.57	0.87
1:A:444:HIS:CE1	1:A:486:LEU:HD13	2.08	0.87
1:A:455:ASN:ND2	1:A:580:VAL:HG13	1.91	0.86
1:A:293:THR:CG2	1:A:312:GLU:HB2	2.06	0.84
1:A:384:TRP:CH2	1:A:555:VAL:HG13	2.13	0.84
1:A:797:ARG:HD3	1:A:800:TRP:CH2	2.14	0.83
1:A:797:ARG:HD3	1:A:800:TRP:CZ2	2.15	0.82
1:A:279:ILE:HG22	1:A:283:ILE:CD1	2.11	0.81
1:A:581:LYS:CB	1:A:593:MET:HE2	2.06	0.81
1:A:601:GLN:HA	1:A:601:GLN:NE2	1.95	0.81
1:A:355:ARG:HG3	1:A:355:ARG:HH11	1.45	0.81
1:A:457:MET:SD	1:A:593:MET:HE1	2.21	0.81
1:A:867:ASN:N	1:A:867:ASN:ND2	2.28	0.81
1:A:727:ARG:HD2	1:A:829:MET:HE2	1.66	0.78
1:A:846:LYS:NZ	1:A:899:LEU:HB3	2.00	0.77
1:A:354:GLN:CD	1:A:584:ARG:HG3	2.10	0.77
1:A:846:LYS:HZ3	1:A:899:LEU:HD13	1.49	0.77
1:A:587:ALA:O	1:A:590:ARG:HG2	1.85	0.77
1:A:773:HIS:H	1:A:773:HIS:CD2	2.03	0.77
1:A:709:GLN:OE1	1:A:720:GLU:HB2	1.84	0.76
1:A:312:GLU:OE2	1:A:590:ARG:HB2	1.87	0.75
1:A:431:PRO:O	1:A:435:GLU:HG3	1.86	0.75
1:A:451:THR:HB	1:A:479:MET:SD	2.26	0.75
1:A:456:MET:HE3	1:A:480:TRP:CB	2.17	0.74
1:A:453:ILE:CD1	1:A:581:LYS:HD2	2.18	0.73
1:A:867:ASN:HD22	1:A:867:ASN:H	1.36	0.73
1:A:681:THR:HG22	1:A:701:PRO:HG3	1.70	0.73
1:A:317:GLY:O	1:A:349:THR:HG22	1.89	0.73
1:A:641:THR:HG21	1:A:642:LYS:HE2	1.71	0.73
1:A:846:LYS:HZ3	1:A:899:LEU:HB3	1.52	0.72
1:A:454:TYR:HE1	1:A:580:VAL:HG22	1.53	0.72
1:A:872:ILE:O	1:A:876:ARG:HG3	1.90	0.72
1:A:622:LEU:HD13	1:A:663:MET:HE1	1.72	0.72
1:A:598:ARG:HB3	1:A:601:GLN:HG2	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:293:THR:CG2	1:A:312:GLU:H	2.02	0.71
1:A:321:SER:CB	1:A:345:MET:HA	2.12	0.71
1:A:581:LYS:HB3	1:A:593:MET:CE	2.09	0.70
1:A:775:ARG:HD2	1:A:856:ILE:HD12	1.73	0.70
1:A:320:SER:HB3	1:A:345:MET:HG3	1.73	0.69
1:A:552:GLU:O	1:A:555:VAL:HG22	1.93	0.69
1:A:846:LYS:NZ	1:A:899:LEU:HD22	2.06	0.69
1:A:694:LYS:HE2	1:A:695:ASP:OD1	1.92	0.69
1:A:641:THR:OG1	1:A:642:LYS:HG2	1.93	0.69
1:A:800:TRP:O	1:A:802:ILE:N	2.26	0.69
1:A:569:ALA:O	1:A:573:LEU:HB2	1.94	0.68
1:A:719:THR:HG21	1:A:848:GLU:OE2	1.92	0.68
1:A:846:LYS:NZ	1:A:899:LEU:HD13	2.07	0.68
1:A:389:ARG:HG3	1:A:389:ARG:HH11	1.58	0.68
1:A:322:LEU:N	1:A:322:LEU:HD22	2.09	0.68
1:A:453:ILE:HD11	1:A:581:LYS:HD2	1.75	0.67
1:A:578:LYS:HE2	1:A:602:ARG:HD3	1.75	0.67
1:A:389:ARG:HH11	1:A:389:ARG:CG	2.09	0.66
1:A:846:LYS:CE	1:A:899:LEU:HD13	2.26	0.66
1:A:386:PHE:O	1:A:389:ARG:HG2	1.95	0.66
1:A:455:ASN:HD21	1:A:580:VAL:HG13	1.60	0.66
1:A:316:THR:HG23	1:A:583:MET:HE1	1.77	0.66
1:A:441:ARG:NH1	1:A:489:GLU:OE1	2.29	0.66
1:A:727:ARG:HD2	1:A:829:MET:CE	2.25	0.66
1:A:456:MET:HE3	1:A:480:TRP:HB3	1.77	0.65
1:A:624:ARG:HG2	1:A:679:PHE:CE1	2.31	0.65
1:A:879:ILE:C	1:A:879:ILE:HD13	2.22	0.65
1:A:355:ARG:HG3	1:A:355:ARG:NH1	2.11	0.65
1:A:863:THR:O	1:A:867:ASN:ND2	2.30	0.65
1:A:474:ARG:HH11	1:A:474:ARG:HG2	1.61	0.65
1:A:719:THR:HG21	1:A:848:GLU:CD	2.21	0.65
1:A:320:SER:CB	1:A:345:MET:HE2	2.26	0.65
1:A:605:GLY:O	1:A:607:VAL:N	2.30	0.65
1:A:297:ASP:O	1:A:299:ASN:ND2	2.30	0.64
1:A:322:LEU:HD22	1:A:322:LEU:H	1.61	0.64
1:A:353:GLN:HE22	1:A:457:MET:HE2	1.62	0.64
1:A:311:TYR:OH	1:A:581:LYS:HE2	1.97	0.64
1:A:776:ASP:OD1	1:A:861:ARG:NH1	2.30	0.63
1:A:282:ARG:NH2	1:A:573:LEU:O	2.29	0.63
1:A:868:ILE:HD13	1:A:868:ILE:O	1.98	0.63
1:A:641:THR:O	1:A:643:GLY:N	2.30	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:893:ARG:NH1	1:A:893:ARG:HG2	2.13	0.62
1:A:279:ILE:CG2	1:A:283:ILE:HD11	2.26	0.62
1:A:317:GLY:CA	1:A:349:THR:HA	2.17	0.62
1:A:836:GLU:O	1:A:837:LYS:HG3	2.00	0.62
1:A:837:LYS:NZ	1:A:837:LYS:HB2	2.14	0.62
1:A:343:MET:HE1	1:A:738:GLU:HA	1.82	0.62
1:A:754:ARG:HG2	1:A:754:ARG:HH11	1.65	0.62
1:A:899:LEU:H	1:A:899:LEU:CD2	2.11	0.61
1:A:766:MET:HE2	1:A:798:THR:CG2	2.29	0.61
1:A:320:SER:HB3	1:A:345:MET:CE	2.30	0.61
1:A:354:GLN:CG	1:A:584:ARG:HG3	2.30	0.61
1:A:527:ARG:O	1:A:673:LYS:HD3	2.00	0.61
1:A:748:GLY:O	1:A:750:GLY:N	2.33	0.61
1:A:404:LYS:O	1:A:408:ASN:HB3	2.01	0.61
1:A:551:ASN:ND2	1:A:554:LYS:HZ1	1.97	0.61
1:A:641:THR:OG1	1:A:642:LYS:HE3	2.00	0.61
1:A:320:SER:HB3	1:A:345:MET:HE2	1.82	0.60
1:A:318:SER:O	1:A:349:THR:HG23	2.01	0.60
1:A:582:VAL:N	1:A:593:MET:HE3	2.16	0.60
1:A:445:LEU:CD2	1:A:562:GLU:HG3	2.32	0.60
1:A:418:GLN:HG3	1:A:747:PRO:HB3	1.83	0.59
1:A:454:TYR:CE1	1:A:580:VAL:HG22	2.37	0.59
1:A:355:ARG:HG3	1:A:356:VAL:N	2.17	0.59
1:A:454:TYR:HE1	1:A:580:VAL:CG2	2.15	0.59
1:A:320:SER:HB3	1:A:345:MET:CG	2.33	0.59
1:A:893:ARG:HG2	1:A:893:ARG:HH11	1.68	0.59
1:A:775:ARG:NH2	1:A:844:SER:O	2.36	0.59
1:A:861:ARG:O	1:A:864:TRP:HD1	1.84	0.59
1:A:502:ARG:NH2	1:A:663:MET:O	2.28	0.59
1:A:879:ILE:HD12	1:A:884:TYR:OH	2.03	0.58
1:A:578:LYS:NZ	1:A:603:GLY:O	2.29	0.58
1:A:293:THR:HG21	1:A:312:GLU:H	1.67	0.58
1:A:305:TRP:HE1	1:A:360:LYS:HD3	1.69	0.58
1:A:407:SER:O	1:A:409:ALA:N	2.36	0.58
1:A:707:ASP:OD1	1:A:709:GLN:HB2	2.03	0.58
1:A:453:ILE:HD12	1:A:581:LYS:HD2	1.85	0.57
1:A:817:VAL:O	1:A:821:VAL:HG13	2.04	0.57
1:A:650:THR:O	1:A:654:GLU:HG2	2.04	0.57
1:A:578:LYS:CE	1:A:602:ARG:HD3	2.35	0.57
1:A:867:ASN:ND2	1:A:867:ASN:H	1.97	0.57
1:A:337:ILE:C	1:A:339:ASN:H	2.13	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:457:MET:HG2	1:A:583:MET:HE3	1.86	0.56
1:A:578:LYS:HZ3	1:A:603:GLY:N	2.04	0.56
1:A:786:SER:O	1:A:876:ARG:NH2	2.30	0.56
1:A:288:ARG:NH1	1:A:288:ARG:HG3	2.21	0.56
1:A:340:VAL:HG12	1:A:737:ASP:O	2.06	0.56
1:A:543:ARG:HD2	1:A:690:SER:O	2.05	0.56
1:A:651:TRP:CZ2	1:A:659:ARG:HD2	2.41	0.56
1:A:402:ILE:HG12	1:A:428:VAL:HB	1.86	0.55
1:A:284:GLU:O	1:A:288:ARG:HG2	2.07	0.55
1:A:436:MET:HE1	1:A:475:ALA:HB1	1.88	0.55
1:A:276:THR:O	1:A:276:THR:HG22	2.07	0.54
1:A:721:LEU:HD11	1:A:773:HIS:HB3	1.89	0.54
1:A:752:ASN:O	1:A:756:THR:HG23	2.07	0.54
1:A:888:MET:O	1:A:891:LEU:HB3	2.08	0.54
1:A:337:ILE:O	1:A:339:ASN:N	2.40	0.54
1:A:636:ASP:HB3	1:A:640:LEU:HD12	1.88	0.54
1:A:689:MET:O	1:A:690:SER:HB2	2.08	0.54
1:A:293:THR:HG21	1:A:312:GLU:CB	2.31	0.54
1:A:293:THR:HG22	1:A:312:GLU:H	1.72	0.54
1:A:445:LEU:HD22	1:A:562:GLU:HG3	1.90	0.54
1:A:893:ARG:HH11	1:A:893:ARG:CG	2.21	0.54
1:A:313:VAL:HG11	1:A:593:MET:HB2	1.90	0.53
1:A:846:LYS:HZ3	1:A:899:LEU:CD1	2.20	0.53
1:A:641:THR:OG1	1:A:644:LYS:HD3	2.08	0.53
1:A:288:ARG:HG3	1:A:288:ARG:HH11	1.73	0.53
1:A:830:GLU:CD	1:A:830:GLU:H	2.17	0.53
1:A:846:LYS:HZ1	1:A:899:LEU:CD2	2.11	0.53
1:A:555:VAL:O	1:A:558:LEU:HB2	2.09	0.53
1:A:602:ARG:HH21	1:A:609:THR:HG21	1.74	0.53
1:A:317:GLY:HA2	1:A:348:THR:O	2.09	0.53
1:A:456:MET:HE3	1:A:480:TRP:HB2	1.89	0.53
1:A:353:GLN:NE2	1:A:457:MET:HE2	2.24	0.52
1:A:622:LEU:HD13	1:A:663:MET:CE	2.37	0.52
1:A:602:ARG:NH2	1:A:606:GLN:OE1	2.43	0.52
1:A:879:ILE:HD13	1:A:879:ILE:O	2.10	0.52
1:A:354:GLN:HG3	1:A:584:ARG:HG3	1.91	0.52
1:A:553:ALA:HB2	1:A:576:ARG:NH2	2.24	0.51
1:A:436:MET:HE1	1:A:475:ALA:CB	2.40	0.51
1:A:330:LEU:O	1:A:864:TRP:HZ3	1.91	0.51
1:A:856:ILE:HA	1:A:861:ARG:HD3	1.92	0.51
1:A:524:VAL:C	1:A:526:THR:H	2.18	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:SER:HA	1:A:346:THR:O	2.10	0.51
1:A:420:GLN:O	1:A:423:SER:HB2	2.10	0.51
1:A:297:ASP:OD2	1:A:584:ARG:NH2	2.44	0.51
1:A:694:LYS:HG3	1:A:695:ASP:N	2.25	0.51
1:A:766:MET:HG3	1:A:798:THR:HG21	1.93	0.51
1:A:320:SER:HB2	1:A:345:MET:HE2	1.93	0.50
1:A:440:GLU:OE1	1:A:450:HIS:HB2	2.12	0.50
1:A:444:HIS:HE1	1:A:486:LEU:HD13	1.70	0.50
1:A:287:ARG:O	1:A:291:SER:HB2	2.10	0.50
1:A:456:MET:HE2	1:A:456:MET:HA	1.93	0.50
1:A:449:CYS:HB2	1:A:573:LEU:HD13	1.93	0.50
1:A:814:MET:HA	1:A:817:VAL:HG13	1.93	0.50
1:A:719:THR:HG21	1:A:848:GLU:OE1	2.11	0.49
1:A:626:MET:SD	1:A:648:VAL:HG13	2.52	0.49
1:A:766:MET:HE2	1:A:798:THR:HG23	1.94	0.49
1:A:432:LYS:NZ	1:A:474:ARG:HH22	2.10	0.49
1:A:773:HIS:H	1:A:773:HIS:HD2	1.55	0.49
1:A:823:ILE:HD11	1:A:835:VAL:HG23	1.95	0.49
1:A:899:LEU:HD23	1:A:899:LEU:N	2.14	0.49
1:A:599:GLU:HG3	1:A:600:ASP:OD2	2.13	0.49
1:A:846:LYS:HZ3	1:A:899:LEU:CB	2.25	0.49
1:A:337:ILE:O	1:A:341:THR:HG23	2.12	0.48
1:A:443:ALA:O	1:A:446:ARG:HG2	2.12	0.48
1:A:641:THR:HB	1:A:642:LYS:H	1.28	0.48
1:A:543:ARG:HG3	1:A:543:ARG:HH11	1.77	0.48
1:A:754:ARG:HH11	1:A:754:ARG:CG	2.24	0.48
1:A:759:LEU:HD13	1:A:796:GLY:HA3	1.96	0.48
1:A:605:GLY:C	1:A:607:VAL:H	2.21	0.48
1:A:419:ASN:C	1:A:421:TRP:H	2.17	0.48
1:A:897:THR:OG1	1:A:898:THR:N	2.46	0.48
1:A:332:LYS:HB3	1:A:333:PRO:HD3	1.95	0.48
1:A:347:ASP:CG	1:A:352:GLY:HA3	2.38	0.48
1:A:401:PHE:CE1	1:A:485:PHE:HA	2.49	0.48
1:A:418:GLN:CG	1:A:419:ASN:H	2.25	0.48
1:A:281:ASN:HB2	4:A:2:HOH:O	2.14	0.47
1:A:500:LEU:HD13	1:A:611:ALA:HA	1.95	0.47
1:A:708:TRP:C	1:A:710:GLN:H	2.22	0.47
1:A:875:VAL:O	1:A:879:ILE:HG23	2.14	0.47
1:A:402:ILE:HG22	1:A:406:ASN:HD22	1.79	0.47
1:A:896:ASP:N	1:A:896:ASP:OD1	2.46	0.47
1:A:773:HIS:O	1:A:843:TYR:HD2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:ILE:HG22	1:A:406:ASN:ND2	2.29	0.47
1:A:405:VAL:O	1:A:421:TRP:NE1	2.47	0.47
1:A:645:GLY:N	1:A:646:PRO:CD	2.78	0.47
1:A:716:ASN:HD21	1:A:734:ARG:HD2	1.80	0.47
1:A:747:PRO:O	1:A:748:GLY:O	2.32	0.47
1:A:294:TRP:O	1:A:295:HIS:HB3	2.14	0.47
1:A:772:PHE:O	1:A:778:ARG:HD3	2.15	0.47
1:A:330:LEU:CD2	1:A:868:ILE:HD11	2.45	0.47
1:A:752:ASN:HD22	1:A:752:ASN:C	2.22	0.47
1:A:801:SER:OG	1:A:802:ILE:N	2.48	0.47
1:A:337:ILE:O	1:A:337:ILE:HG22	2.14	0.46
1:A:578:LYS:O	1:A:597:SER:HB2	2.15	0.46
1:A:610:TYR:CE1	1:A:803:HIS:CE1	3.03	0.46
1:A:300:HIS:O	1:A:300:HIS:ND1	2.49	0.46
1:A:275:ASP:O	1:A:276:THR:OG1	2.27	0.46
1:A:416:GLU:OE1	1:A:416:GLU:HA	2.15	0.46
1:A:810:THR:HB	1:A:812:GLU:H	1.79	0.46
1:A:293:THR:HG21	1:A:312:GLU:N	2.30	0.46
1:A:402:ILE:CG2	1:A:406:ASN:ND2	2.79	0.46
1:A:874:GLN:O	1:A:878:ILE:HG13	2.15	0.46
1:A:332:LYS:O	1:A:335:ASP:HB2	2.15	0.46
1:A:617:ASN:ND2	1:A:621:GLN:HE21	2.14	0.46
1:A:578:LYS:HG3	1:A:601:GLN:CB	2.46	0.46
1:A:814:MET:CE	1:A:817:VAL:HG11	2.46	0.45
1:A:355:ARG:CG	1:A:356:VAL:N	2.79	0.45
1:A:532:ILE:O	1:A:705:TRP:HE3	1.98	0.45
1:A:315:PRO:HG3	1:A:591:THR:OG1	2.15	0.45
1:A:322:LEU:HB2	1:A:746:SER:O	2.16	0.45
1:A:754:ARG:CG	1:A:754:ARG:NH1	2.80	0.45
1:A:278:LYS:HB3	1:A:573:LEU:CD2	2.46	0.45
1:A:456:MET:HE1	1:A:476:ILE:O	2.16	0.45
1:A:888:MET:O	1:A:891:LEU:N	2.44	0.45
1:A:419:ASN:O	1:A:422:ARG:HG2	2.17	0.45
1:A:723:MET:HG2	1:A:842:PRO:HG3	1.99	0.45
1:A:549:LEU:O	1:A:576:ARG:NH2	2.50	0.45
1:A:775:ARG:NE	1:A:849:ASP:OD2	2.46	0.45
1:A:444:HIS:C	1:A:446:ARG:N	2.74	0.45
1:A:306:ASN:O	1:A:596:ILE:HA	2.17	0.45
1:A:846:LYS:NZ	1:A:899:LEU:CB	2.76	0.45
1:A:712:PRO:HG3	1:A:717:HIS:HE1	1.81	0.44
1:A:779:LEU:HD23	1:A:779:LEU:C	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:866:GLU:OE2	1:A:866:GLU:HA	2.17	0.44
1:A:893:ARG:NH1	1:A:893:ARG:CG	2.78	0.44
1:A:290:TYR:O	1:A:294:TRP:N	2.47	0.44
1:A:343:MET:HG2	1:A:343:MET:O	2.18	0.44
1:A:321:SER:HB3	1:A:345:MET:CA	2.15	0.44
1:A:697:GLN:OE1	1:A:697:GLN:HA	2.09	0.44
1:A:330:LEU:HD22	1:A:868:ILE:HD11	2.00	0.44
1:A:475:ALA:HB1	1:A:479:MET:HE2	1.99	0.44
1:A:681:THR:HG22	1:A:701:PRO:CG	2.45	0.44
1:A:348:THR:O	1:A:348:THR:HG22	2.16	0.44
1:A:418:GLN:HG3	1:A:747:PRO:CB	2.46	0.44
1:A:531:ARG:HG2	1:A:706:TYR:CE1	2.52	0.44
1:A:606:GLN:HA	1:A:609:THR:HB	2.00	0.44
1:A:562:GLU:OE2	1:A:565:ARG:NH2	2.51	0.44
1:A:858:THR:HG22	1:A:859:ARG:N	2.32	0.44
1:A:297:ASP:C	1:A:299:ASN:N	2.76	0.44
1:A:311:TYR:O	1:A:313:VAL:HG13	2.17	0.43
1:A:330:LEU:O	1:A:864:TRP:CZ3	2.71	0.43
1:A:475:ALA:O	1:A:479:MET:HG3	2.17	0.43
1:A:578:LYS:HZ3	1:A:603:GLY:H	1.64	0.43
1:A:444:HIS:C	1:A:446:ARG:H	2.24	0.43
1:A:281:ASN:HB3	1:A:449:CYS:O	2.19	0.43
1:A:602:ARG:HE	1:A:602:ARG:HB2	1.64	0.43
1:A:310:SER:HA	1:A:593:MET:O	2.18	0.43
1:A:419:ASN:OD1	1:A:755:ASP:OD2	2.37	0.43
1:A:445:LEU:HD21	1:A:562:GLU:HG3	1.98	0.43
1:A:685:PHE:O	1:A:689:MET:HG3	2.19	0.43
1:A:602:ARG:H	1:A:602:ARG:HG3	1.50	0.42
1:A:802:ILE:HG12	1:A:803:HIS:H	1.83	0.42
1:A:433:PHE:HB2	1:A:478:PHE:CE2	2.54	0.42
1:A:645:GLY:N	1:A:646:PRO:HD2	2.34	0.42
1:A:778:ARG:NH2	1:A:838:TRP:O	2.53	0.42
1:A:708:TRP:C	1:A:710:GLN:N	2.78	0.42
1:A:797:ARG:HG2	1:A:800:TRP:CE2	2.54	0.42
1:A:849:ASP:O	1:A:854:SER:HB2	2.19	0.42
1:A:389:ARG:CG	1:A:389:ARG:NH1	2.74	0.42
1:A:819:ASN:OD1	1:A:823:ILE:HD12	2.20	0.42
1:A:830:GLU:CD	1:A:830:GLU:N	2.78	0.42
1:A:868:ILE:HD13	1:A:871:ALA:HB3	2.02	0.42
1:A:474:ARG:HG2	1:A:474:ARG:NH1	2.28	0.42
1:A:310:SER:HB3	1:A:594:ASP:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:TRP:CZ2	1:A:555:VAL:HG13	2.52	0.42
1:A:526:THR:O	1:A:526:THR:HG22	2.20	0.42
1:A:520:ILE:HD13	1:A:520:ILE:N	2.34	0.42
1:A:727:ARG:HB3	1:A:829:MET:HE1	2.02	0.42
1:A:766:MET:CE	1:A:798:THR:CG2	2.98	0.42
1:A:456:MET:HE3	1:A:456:MET:HB2	1.95	0.41
1:A:833:THR:HA	1:A:834:PRO:HD3	1.80	0.41
1:A:297:ASP:C	1:A:299:ASN:H	2.28	0.41
1:A:456:MET:CE	1:A:480:TRP:HB2	2.49	0.41
1:A:727:ARG:HD3	1:A:831:ASP:HB3	2.01	0.41
1:A:484:ARG:HD3	1:A:484:ARG:HA	1.74	0.41
1:A:293:THR:HB	1:A:311:TYR:HB2	2.02	0.41
1:A:337:ILE:C	1:A:339:ASN:N	2.78	0.41
1:A:858:THR:CG2	1:A:859:ARG:N	2.83	0.41
1:A:559:LEU:HD22	1:A:563:HIS:CE1	2.56	0.41
1:A:584:ARG:HE	1:A:584:ARG:HB3	1.55	0.41
1:A:347:ASP:OD1	1:A:349:THR:OG1	2.30	0.41
1:A:347:ASP:HB3	1:A:352:GLY:HA3	2.03	0.41
1:A:846:LYS:O	1:A:850:ILE:HD12	2.21	0.41
1:A:441:ARG:HH11	1:A:489:GLU:CD	2.28	0.41
1:A:846:LYS:NZ	1:A:899:LEU:CG	2.84	0.41
1:A:846:LYS:NZ	1:A:899:LEU:CD1	2.80	0.41
1:A:886:ASP:OD1	1:A:888:MET:HB2	2.20	0.41
1:A:551:ASN:HD22	1:A:554:LYS:HE3	1.85	0.40
1:A:582:VAL:C	1:A:593:MET:HE3	2.46	0.40
1:A:879:ILE:HD11	1:A:882:GLU:HG3	2.02	0.40
1:A:336:THR:O	1:A:336:THR:OG1	2.33	0.40
1:A:495:ASN:ND2	1:A:607:VAL:O	2.54	0.40
1:A:606:GLN:H	1:A:606:GLN:HE21	1.67	0.40
1:A:766:MET:HG3	1:A:798:THR:CG2	2.51	0.40
1:A:641:THR:CB	1:A:642:LYS:HG2	2.52	0.40
1:A:578:LYS:HZ1	1:A:603:GLY:C	2.24	0.40
1:A:712:PRO:HG3	1:A:717:HIS:CE1	2.56	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	602/639 (94%)	546 (91%)	39 (6%)	17 (3%)	4	21

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	276	THR
1	A	320	SER
1	A	338	THR
1	A	363	THR
1	A	408	ASN
1	A	606	GLN
1	A	642	LYS
1	A	748	GLY
1	A	749	ALA
1	A	750	GLY
1	A	898	THR
1	A	364	LYS
1	A	472	GLY
1	A	604	SER
1	A	317	GLY
1	A	709	GLN
1	A	802	ILE

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	523/550 (95%)	402 (77%)	121 (23%)	1 5

All (121) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	274	SER
1	A	275	ASP
1	A	280	LYS
1	A	284	GLU
1	A	288	ARG
1	A	293	THR
1	A	297	ASP
1	A	303	ARG
1	A	306	ASN
1	A	312	GLU
1	A	313	VAL
1	A	318	SER
1	A	320	SER
1	A	322	LEU
1	A	323	VAL
1	A	332	LYS
1	A	336	THR
1	A	342	THR
1	A	343	MET
1	A	345	MET
1	A	360	LYS
1	A	361	VAL
1	A	364	LYS
1	A	376	LEU
1	A	389	ARG
1	A	404	LYS
1	A	405	VAL
1	A	416	GLU
1	A	417	GLU
1	A	418	GLN
1	A	419	ASN
1	A	420	GLN
1	A	421	TRP
1	A	422	ARG
1	A	423	SER
1	A	436	MET
1	A	441	ARG
1	A	442	GLU

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Mol	Chain	Res	Type
1	A	454	TYR
1	A	455	ASN
1	A	456	MET
1	A	457	MET
1	A	484	ARG
1	A	503	LYS
1	A	514	LEU
1	A	515	GLN
1	A	520	ILE
1	A	521	LEU
1	A	527	ARG
1	A	555	VAL
1	A	558	LEU
1	A	562	GLU
1	A	564	ARG
1	A	572	GLU
1	A	582	VAL
1	A	588	ASP
1	A	592	VAL
1	A	596	ILE
1	A	598	ARG
1	A	601	GLN
1	A	602	ARG
1	A	604	SER
1	A	606	GLN
1	A	617	ASN
1	A	622	LEU
1	A	639	LYS
1	A	642	LYS
1	A	659	ARG
1	A	660	LEU
1	A	675	LEU
1	A	694	LYS
1	A	697	GLN
1	A	705	TRP
1	A	709	GLN
1	A	719	THR
1	A	724	LYS
1	A	733	CYS
1	A	740	VAL
1	A	746	SER
1	A	754	ARG

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Mol	Chain	Res	Type
1	A	756	THR
1	A	762	SER
1	A	769	LEU
1	A	773	HIS
1	A	778	ARG
1	A	790	VAL
1	A	797	ARG
1	A	799	THR
1	A	802	ILE
1	A	810	THR
1	A	811	THR
1	A	815	LEU
1	A	816	GLU
1	A	817	VAL
1	A	821	VAL
1	A	830	GLU
1	A	836	GLU
1	A	837	LYS
1	A	844	SER
1	A	849	ASP
1	A	850	ILE
1	A	852	CYS
1	A	855	LEU
1	A	856	ILE
1	A	859	ARG
1	A	861	ARG
1	A	864	TRP
1	A	867	ASN
1	A	868	ILE
1	A	869	GLN
1	A	872	ILE
1	A	874	GLN
1	A	877	SER
1	A	879	ILE
1	A	890	SER
1	A	892	LYS
1	A	893	ARG
1	A	896	ASP
1	A	897	THR
1	A	898	THR
1	A	899	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	377	ASN
1	A	455	ASN
1	A	495	ASN
1	A	551	ASN
1	A	601	GLN
1	A	617	ASN
1	A	621	GLN
1	A	716	ASN
1	A	717	HIS
1	A	752	ASN
1	A	773	HIS
1	A	803	HIS
1	A	867	ASN
1	A	869	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](#)

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	608/639 (95%)	1.17	91 (14%) 5 3	67, 89, 111, 133	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	321	SER	4.5
1	A	320	SER	3.9
1	A	361	VAL	3.9
1	A	801	SER	3.8
1	A	642	LYS	3.7
1	A	710	GLN	3.7
1	A	408	ASN	3.7
1	A	345	MET	3.4
1	A	719	THR	3.4
1	A	746	SER	3.3
1	A	560	ASP	3.2
1	A	749	ALA	3.1
1	A	318	SER	3.1
1	A	842	PRO	3.1
1	A	866	GLU	3.0
1	A	725	ASP	3.0
1	A	895	GLU	3.0
1	A	274	SER	3.0
1	A	458	GLY	2.9
1	A	453	ILE	2.9
1	A	686	LEU	2.9
1	A	638	GLU	2.8
1	A	602	ARG	2.8
1	A	886	ASP	2.8
1	A	898	THR	2.8
1	A	791	ASN	2.8
1	A	803	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	896	ASP	2.7
1	A	406	ASN	2.7
1	A	317	GLY	2.7
1	A	527	ARG	2.7
1	A	455	ASN	2.7
1	A	799	THR	2.6
1	A	409	ALA	2.6
1	A	606	GLN	2.6
1	A	454	TYR	2.6
1	A	365	ALA	2.6
1	A	640	LEU	2.5
1	A	887	TYR	2.5
1	A	416	GLU	2.5
1	A	528	PRO	2.5
1	A	703	THR	2.5
1	A	442	GLU	2.4
1	A	445	LEU	2.4
1	A	811	THR	2.4
1	A	407	SER	2.4
1	A	846	LYS	2.4
1	A	864	TRP	2.4
1	A	675	LEU	2.4
1	A	439	GLU	2.4
1	A	326	VAL	2.4
1	A	601	GLN	2.3
1	A	557	GLU	2.3
1	A	572	GLU	2.3
1	A	363	THR	2.3
1	A	639	LYS	2.3
1	A	319	ALA	2.3
1	A	843	TYR	2.3
1	A	515	GLN	2.3
1	A	607	VAL	2.3
1	A	798	THR	2.2
1	A	322	LEU	2.2
1	A	690	SER	2.2
1	A	844	SER	2.2
1	A	733	CYS	2.2
1	A	362	ASP	2.2
1	A	569	ALA	2.2
1	A	338	THR	2.2
1	A	682	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	475	ALA	2.2
1	A	534	ALA	2.2
1	A	600	ASP	2.2
1	A	637	VAL	2.1
1	A	432	LYS	2.1
1	A	390	GLU	2.1
1	A	275	ASP	2.1
1	A	335	ASP	2.1
1	A	296	HIS	2.1
1	A	300	HIS	2.1
1	A	476	ILE	2.1
1	A	532	ILE	2.1
1	A	731	THR	2.1
1	A	495	ASN	2.1
1	A	329	LEU	2.1
1	A	695	ASP	2.1
1	A	473	SER	2.0
1	A	278	LYS	2.0
1	A	399	GLU	2.0
1	A	531	ARG	2.0
1	A	748	GLY	2.0
1	A	856	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	MG	A	906	1/1	0.87	0.28	79,79,79,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ZN	A	907	1/1	0.94	0.07	77,77,77,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.