



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 12:59 PM GMT

PDB ID : 1KFS  
Title : DNA POLYMERASE I KLENOW FRAGMENT (E.C.2.7.7.7) MUTANT/DNA COMPLEX  
Authors : Brautigam, C.A.; Steitz, T.A.  
Deposited on : 1997-08-18  
Resolution : 2.10 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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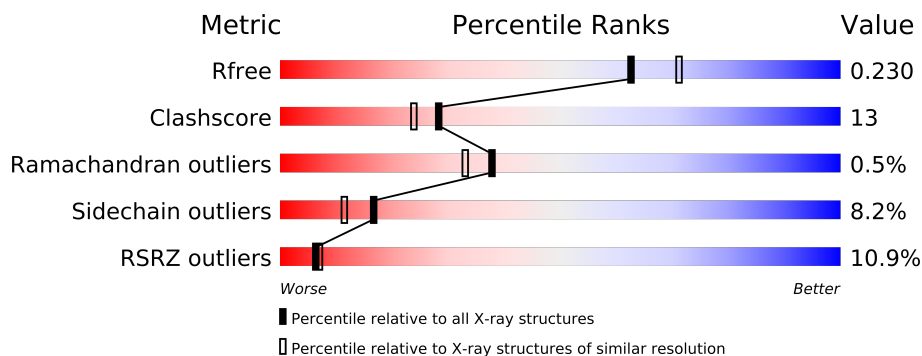
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.10 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))
$R_{free}$	66092	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (2.10-2.10)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	B	7	
2	A	605	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5136 atoms, of which 1 is hydrogen and 0 are deuterium.

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 DNA chain called DNA (5'-D(\*GP\*CP\*TP\*TP\*AP\*CP\*G)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	3	Total	C	H	N	O	P	0	0	0
			63	29	1	13	17	3			

- Molecule 2 is a protein called PROTEIN (DNA POLYMERASE I KLENOW FRAGMENT (E.C.2.7.7.7)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	601	Total	C	N	O	S	0	0	0
			4753	3008	830	899	16			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	324	MET	VAL	ENGINEERED	UNP P00582

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	309	Total	O	0	0
			309	309		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	7	Total	O	0	0
			7	7		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.70Å 101.70Å 85.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.10 19.95 – 2.10	Depositor EDS
% Data completeness (in resolution range)	86.2 (20.00-2.10) 97.5 (19.95-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	5.50	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.58 (at 2.09Å)	Xtriage
Refinement program	X-PLOR 3.8	Depositor
R, $R_{free}$	0.195 , 0.219 0.209 , 0.230	Depositor DCC
$R_{free}$ test set	4057 reflections (8.88%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.8	Xtriage
Anisotropy	0.181	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.42 , 60.9	EDS
Estimated twinning fraction	0.034 for h,-k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 86377 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5136	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 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, 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	B	3.47	7/69 (10.1%)	5.15	25/104 (24.0%)
2	A	0.50	0/4839	0.75	6/6547 (0.1%)
All	All	0.65	7/4908 (0.1%)	0.98	31/6651 (0.5%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1005	DA	C2'-C1'	15.88	1.68	1.52
1	B	1005	DA	C4'-C3'	-10.87	1.41	1.52
1	B	1007	DG	C2-N3	7.57	1.38	1.32
1	B	1005	DA	C2-N3	-5.58	1.28	1.33
1	B	1007	DG	N9-C4	-5.31	1.33	1.38
1	B	1005	DA	N9-C4	-5.25	1.34	1.37
1	B	1005	DA	O4'-C1'	5.23	1.48	1.42

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1005	DA	O4'-C1'-N9	26.65	126.65	108.00
1	B	1005	DA	O4'-C1'-C2'	-18.69	90.95	105.90
1	B	1005	DA	N9-C1'-C2'	-11.71	90.34	112.60
1	B	1007	DG	C5-N7-C8	-9.70	99.45	104.30
1	B	1007	DG	N7-C8-N9	9.51	117.86	113.10
1	B	1005	DA	P-O5'-C5'	9.20	135.62	120.90
1	B	1005	DA	C3'-C2'-C1'	9.18	113.51	102.50
1	B	1007	DG	C8-N9-C4	-8.75	102.90	106.40
1	B	1006	DC	C4'-C3'-C2'	-8.35	95.59	103.10
1	B	1006	DC	P-O3'-C3'	7.99	129.29	119.70
1	B	1005	DA	N3-C4-N9	-7.56	121.35	127.40
1	B	1005	DA	N9-C4-C5	7.26	108.70	105.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1005	DA	C5-C6-N1	-7.12	114.14	117.70
1	B	1007	DG	C6-C5-N7	-7.05	126.17	130.40
1	B	1005	DA	O5'-C5'-C4'	-6.79	94.02	111.00
1	B	1006	DC	O4'-C4'-C3'	6.77	110.06	106.00
1	B	1007	DG	C4-C5-N7	6.69	113.47	110.80
1	B	1005	DA	C1'-O4'-C4'	6.23	116.33	110.10
2	A	821	ARG	NE-CZ-NH2	-6.21	117.20	120.30
1	B	1005	DA	OP2-P-O3'	6.14	118.71	105.20
1	B	1005	DA	C2-N3-C4	-6.09	107.56	110.60
2	A	597	LYS	N-CA-C	5.93	127.02	111.00
1	B	1007	DG	C4'-C3'-C2'	-5.80	97.88	103.10
2	A	607	PRO	N-CA-CB	5.74	110.19	103.30
1	B	1005	DA	C8-N9-C1'	5.61	137.79	127.70
2	A	676	LEU	CA-CB-CG	5.56	128.08	115.30
1	B	1005	DA	C4'-C3'-O3'	-5.39	96.21	109.70
1	B	1005	DA	C4-N9-C1'	-5.16	117.01	126.30
2	A	405	LEU	CA-CB-CG	5.08	126.98	115.30
1	B	1007	DG	O4'-C4'-C3'	-5.01	102.50	104.50
2	A	887	GLU	N-CA-C	-5.00	97.49	111.00

There are no chirality outliers.

There are no planarity outliers.

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	62	1	33	5	0
2	A	4753	0	4753	120	0
3	A	3	0	0	0	0
4	B	1	0	0	0	0
5	A	309	0	0	9	0
5	B	7	0	0	0	0
All	All	5135	1	4786	121	0



Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

All (121) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:740:GLU:HB3	2:A:794:ARG:HG2	1.49	0.95
2:A:828:ILE:HG23	2:A:829:LYS:HE2	1.57	0.86
2:A:782:LYS:HA	2:A:785:GLN:HB3	1.65	0.78
2:A:485:THR:H	2:A:488:GLN:HE21	1.30	0.78
2:A:855:ILE:HG23	2:A:908:THR:HG21	1.66	0.77
2:A:677:GLN:HE21	2:A:881:HIS:H	1.32	0.76
2:A:545:VAL:HG23	2:A:877:ILE:HD12	1.66	0.76
2:A:363:ASN:HD22	2:A:542:ARG:HH11	1.36	0.74
2:A:712:ARG:HD3	2:A:913:PRO:O	1.89	0.73
2:A:696:PRO:HB2	5:A:143:HOH:O	1.89	0.72
2:A:612:GLU:HB3	2:A:615:GLU:HG2	1.72	0.71
2:A:600:LYS:CB	2:A:614:LEU:HG	2.20	0.71
2:A:446:SER:OG	2:A:456:HIS:HD2	1.75	0.69
2:A:336:GLU:HG3	5:A:144:HOH:O	1.94	0.65
2:A:418:GLY:HA3	2:A:421:LEU:HD13	1.77	0.65
2:A:711:LEU:HD13	2:A:765:ILE:HD11	1.78	0.64
2:A:435:LEU:HD13	2:A:438:ILE:HG12	1.81	0.62
2:A:772:GLY:O	2:A:776:GLN:HG2	2.01	0.61
2:A:640:TYR:O	2:A:644:LEU:HB2	1.99	0.61
2:A:746:LEU:O	2:A:749:VAL:HG12	2.01	0.60
2:A:782:LYS:HA	2:A:785:GLN:CB	2.29	0.60
2:A:363:ASN:ND2	2:A:542:ARG:HH11	1.98	0.60
2:A:575:GLY:O	2:A:576:GLU:HG2	2.02	0.60
2:A:872:PRO:HG2	2:A:874:VAL:HG13	1.82	0.60
2:A:717:LEU:HD21	2:A:818:LEU:HD11	1.84	0.59
2:A:346:LEU:CD1	2:A:375:ILE:HG23	2.33	0.59
2:A:858:ARG:HB2	2:A:908:THR:HG23	1.85	0.58
2:A:740:GLU:HG2	5:A:277:HOH:O	2.02	0.58
2:A:828:ILE:CG2	2:A:829:LYS:HE2	2.32	0.58
2:A:556:SER:HB2	2:A:641:THR:HG22	1.86	0.57
2:A:731:LYS:HD2	2:A:746:LEU:HD22	1.87	0.57
2:A:677:GLN:HE21	2:A:881:HIS:N	2.01	0.56
2:A:586:LEU:HD22	2:A:627:ILE:HD13	1.86	0.56
2:A:836:ARG:O	2:A:840:GLU:HG3	2.06	0.56
2:A:674:PRO:HG2	2:A:676:LEU:HD13	1.88	0.56
2:A:421:LEU:HD23	2:A:438:ILE:HG23	1.87	0.56
2:A:677:GLN:NE2	2:A:881:HIS:H	2.02	0.55
1:B:1005:DA:OP1	2:A:455:ARG:NH2	2.40	0.55
2:A:588:THR:O	2:A:592:GLU:HB3	2.07	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:808:GLN:O	2:A:812:GLN:HG2	2.07	0.54
2:A:725:THR:O	2:A:729:GLU:HG2	2.07	0.54
1:B:1005:DA:P	2:A:455:ARG:HH21	2.31	0.54
2:A:684:GLU:O	2:A:687:ARG:HB2	2.08	0.53
2:A:565:GLU:HG2	5:A:157:HOH:O	2.07	0.53
2:A:802:MET:O	2:A:806:ARG:HG3	2.09	0.53
2:A:547:ILE:HD12	2:A:655:VAL:HG21	1.91	0.52
2:A:586:LEU:HD11	2:A:627:ILE:HG21	1.91	0.52
2:A:779:ILE:HD12	2:A:783:GLU:HG3	1.92	0.52
2:A:610:SER:HA	2:A:776:GLN:HE22	1.76	0.51
2:A:556:SER:HB2	2:A:641:THR:CG2	2.41	0.51
2:A:600:LYS:CB	2:A:613:VAL:HB	2.40	0.51
2:A:668:ARG:HE	2:A:849:GLN:HG3	1.76	0.51
2:A:633:LEU:CD2	2:A:685:GLU:HG3	2.41	0.50
2:A:448:ILE:HD11	2:A:530:GLU:HG3	1.92	0.50
2:A:487:ASN:H	2:A:487:ASN:HD22	1.59	0.50
2:A:734:HIS:CD2	2:A:758:LYS:HA	2.47	0.50
2:A:717:LEU:HD21	2:A:818:LEU:CD1	2.42	0.49
2:A:727:PHE:CE1	2:A:733:ILE:HG12	2.47	0.49
2:A:345:LYS:HD2	5:A:268:HOH:O	2.13	0.48
2:A:393:PRO:HD2	2:A:491:LEU:HD22	1.95	0.48
2:A:324:MET:HE2	5:A:247:HOH:O	2.12	0.48
2:A:580:LEU:N	2:A:580:LEU:HD22	2.29	0.47
2:A:617:LEU:HB3	2:A:624:PRO:HG3	1.95	0.47
2:A:487:ASN:ND2	2:A:487:ASN:H	2.11	0.47
2:A:844:ILE:O	2:A:848:MET:HE2	2.14	0.47
2:A:732:ASP:OD2	2:A:754:ARG:NH1	2.47	0.47
2:A:324:MET:HA	2:A:324:MET:CE	2.45	0.47
2:A:586:LEU:CD1	2:A:627:ILE:HG21	2.44	0.46
1:B:1005:DA:N6	2:A:658:SER:HB3	2.31	0.46
2:A:751:SER:HB3	5:A:166:HOH:O	2.16	0.46
2:A:615:GLU:HB3	2:A:628:LEU:HD21	1.96	0.46
2:A:783:GLU:O	2:A:786:LYS:HB3	2.15	0.46
2:A:683:ASN:OD1	2:A:684:GLU:N	2.48	0.46
2:A:677:GLN:HB3	5:A:194:HOH:O	2.16	0.46
2:A:677:GLN:HG2	2:A:880:VAL:HG23	1.97	0.46
2:A:547:ILE:HD13	2:A:644:LEU:HG	1.97	0.46
2:A:909:ARG:HB3	2:A:911:ASP:OD1	2.16	0.46
2:A:486:PHE:CE2	2:A:494:ALA:HB1	2.51	0.45
2:A:418:GLY:O	2:A:441:ASP:HA	2.17	0.45
2:A:756:SER:HA	2:A:777:LEU:HD11	1.99	0.45
2:A:589:ILE:O	2:A:593:LYS:HG2	2.16	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:A:576:GLU:HG3	2:A:577:GLU:O	2.16	0.45
2:A:614:LEU:HB2	2:A:628:LEU:HD13	1.99	0.45
2:A:711:LEU:HD22	2:A:762:PHE:CE1	2.52	0.45
2:A:880:VAL:HG11	2:A:924:TRP:HZ2	1.82	0.44
2:A:580:LEU:HD12	2:A:627:ILE:HG12	1.99	0.44
2:A:513:TRP:HB3	2:A:514:PRO:HD3	2.00	0.44
2:A:472:THR:HG22	2:A:475:GLU:H	1.82	0.44
2:A:481:LYS:HE2	2:A:481:LYS:HB3	1.84	0.44
2:A:418:GLY:HA3	2:A:421:LEU:CD1	2.46	0.44
2:A:346:LEU:HD13	2:A:375:ILE:HG23	2.00	0.44
2:A:808:GLN:OE1	2:A:812:GLN:NE2	2.51	0.43
2:A:472:THR:HG22	2:A:475:GLU:HG3	2.00	0.43
2:A:773:LEU:HD13	2:A:784:ALA:HB1	1.99	0.43
2:A:711:LEU:HD23	2:A:733:ILE:HD13	1.98	0.43
2:A:706:TYR:HB3	2:A:709:ILE:HB	2.00	0.43
2:A:690:ARG:HG2	2:A:924:TRP:CE3	2.53	0.43
2:A:880:VAL:HG11	2:A:924:TRP:CZ2	2.54	0.43
1:B:1005:DA:N3	1:B:1005:DA:C2'	2.79	0.43
2:A:715:ALA:HB1	2:A:724:LEU:HD13	2.00	0.43
2:A:354:PHE:CZ	2:A:428:LEU:HD11	2.54	0.42
2:A:708:GLN:HB3	2:A:708:GLN:HE21	1.61	0.42
2:A:798:VAL:O	2:A:802:MET:HG3	2.19	0.42
2:A:714:MET:HB2	2:A:848:MET:SD	2.59	0.42
2:A:808:GLN:HA	2:A:811:GLU:HB3	2.01	0.42
2:A:760:ILE:HD11	2:A:787:TYR:HB3	2.01	0.41
2:A:705:ASP:O	2:A:916:VAL:HA	2.19	0.41
2:A:710:GLU:OE1	2:A:765:ILE:HG21	2.20	0.41
2:A:623:LEU:N	2:A:624:PRO:HD2	2.35	0.41
2:A:771:PHE:O	2:A:775:ARG:HG3	2.21	0.41
2:A:896:VAL:O	2:A:900:ILE:HG12	2.21	0.41
2:A:780:PRO:O	2:A:783:GLU:HB3	2.20	0.41
1:B:1007:DG:N3	2:A:361:LEU:HG	2.36	0.41
2:A:773:LEU:HD22	2:A:773:LEU:O	2.20	0.41
2:A:619:LEU:HD22	2:A:619:LEU:N	2.36	0.41
2:A:434:GLU:HG3	5:A:303:HOH:O	2.20	0.41
2:A:582:SER:HA	2:A:586:LEU:HB2	2.03	0.41
2:A:552:LEU:HD12	2:A:552:LEU:HA	1.96	0.41
2:A:623:LEU:HG	2:A:627:ILE:CD1	2.52	0.40
2:A:449:LEU:HD13	2:A:516:LEU:HG	2.02	0.40
2:A:449:LEU:HD12	2:A:449:LEU:HA	1.82	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	A	597/605 (99%)	566 (95%)	28 (5%)	3 (0%)	38 33

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	830	SER
2	A	594	GLN
2	A	597	LYS

### 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
2	A	500/510 (98%)	459 (92%)	41 (8%)	17 11

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	324	MET
2	A	334	LEU
2	A	339	LEU
2	A	346	LEU
2	A	375	ILE
2	A	405	LEU
2	A	409	LEU
2	A	421	LEU
2	A	425	ARG
2	A	435	LEU
2	A	449	LEU

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Mol	Chain	Res	Type
2	A	472	THR
2	A	510	LEU
2	A	552	LEU
2	A	578	PHE
2	A	584	LYS
2	A	592	GLU
2	A	597	LYS
2	A	616	GLU
2	A	620	ASP
2	A	629	GLU
2	A	644	LEU
2	A	646	LEU
2	A	676	LEU
2	A	680	PRO
2	A	708	GLN
2	A	711	LEU
2	A	719	ARG
2	A	724	LEU
2	A	754	ARG
2	A	773	LEU
2	A	783	GLU
2	A	794	ARG
2	A	799	LEU
2	A	808	GLN
2	A	821	ARG
2	A	867	LEU
2	A	871	GLN
2	A	874	VAL
2	A	879	GLN
2	A	917	GLU

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

Mol	Chain	Res	Type
2	A	363	ASN
2	A	456	HIS
2	A	487	ASN
2	A	488	GLN
2	A	543	ASN
2	A	571	HIS
2	A	677	GLN
2	A	708	GLN

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Mol	Chain	Res	Type
2	A	716	HIS
2	A	734	HIS
2	A	776	GLN
2	A	812	GLN
2	A	845	ASN
2	A	879	GLN
2	A	899	GLN
2	A	901	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 4 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.

### 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	B	3/7 (42%)	0.17	0	100 100	29, 29, 38, 61	0
2	A	601/605 (99%)	0.49	66 (10%)	6 7	12, 31, 88, 100	0
All	All	604/612 (98%)	0.48	66 (10%)	6 7	12, 31, 88, 100	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	583	THR	16.8
2	A	607	PRO	11.9
2	A	608	SER	11.7
2	A	610	SER	10.0
2	A	601	LYS	9.8
2	A	602	THR	9.1
2	A	596	ILE	9.0
2	A	609	THR	8.4
2	A	597	LYS	7.2
2	A	581	SER	7.2
2	A	598	PRO	6.7
2	A	619	LEU	6.6
2	A	781	ARG	6.6
2	A	594	GLN	6.5
2	A	682	ARG	6.3
2	A	582	SER	6.2
2	A	681	VAL	6.2
2	A	621	TYR	5.9
2	A	620	ASP	5.6
2	A	591	PHE	5.5
2	A	611	GLU	5.4
2	A	599	LEU	5.3
2	A	616	GLU	5.0
2	A	771	PHE	5.0

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Mol	Chain	Res	Type	RSRZ
2	A	588	THR	5.0
2	A	577	GLU	4.8
2	A	575	GLY	4.6
2	A	775	ARG	4.5
2	A	831	SER	4.5
2	A	576	GLU	4.4
2	A	590	LEU	4.4
2	A	584	LYS	4.2
2	A	833	GLY	4.1
2	A	615	GLU	3.8
2	A	778	ASN	3.7
2	A	808	GLN	3.7
2	A	830	SER	3.5
2	A	832	ASN	3.5
2	A	481	LYS	3.4
2	A	834	ALA	3.4
2	A	587	GLN	3.3
2	A	766	TYR	3.3
2	A	592	GLU	3.3
2	A	835	ARG	3.3
2	A	482	ASN	3.1
2	A	612	GLU	3.0
2	A	776	GLN	3.0
2	A	769	SER	3.0
2	A	768	MET	3.0
2	A	782	LYS	2.9
2	A	613	VAL	2.8
2	A	585	GLN	2.7
2	A	579	ASN	2.7
2	A	618	ALA	2.6
2	A	593	LYS	2.5
2	A	480	GLY	2.4
2	A	586	LEU	2.4
2	A	617	LEU	2.4
2	A	783	GLU	2.3
2	A	553	HIS	2.3
2	A	390	LEU	2.2
2	A	772	GLY	2.2
2	A	780	PRO	2.2
2	A	800	GLU	2.1
2	A	793	GLU	2.1
2	A	520	LYS	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 ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	ZN	A	320	1/1	0.11	-0.43	99,99,99,99	0
4	MG	B	2	1/1	0.07	-1.31	36,36,36,36	0
3	ZN	A	3	1/1	0.05	-1.46	66,66,66,66	0
3	ZN	A	1	1/1	0.04	-4.24	21,21,21,21	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.