



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 06:20 am BST

PDB ID : 4YOC  
Title : Crystal Structure of human DNMT1 and USP7/HAUSP complex  
Authors : Cheng, J.; Yang, H.; Fang, J.; Gong, R.; Wang, P.; Li, Z.; Xu, Y.  
Deposited on : 2015-03-11  
Resolution : 2.92 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](mailto: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.

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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  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

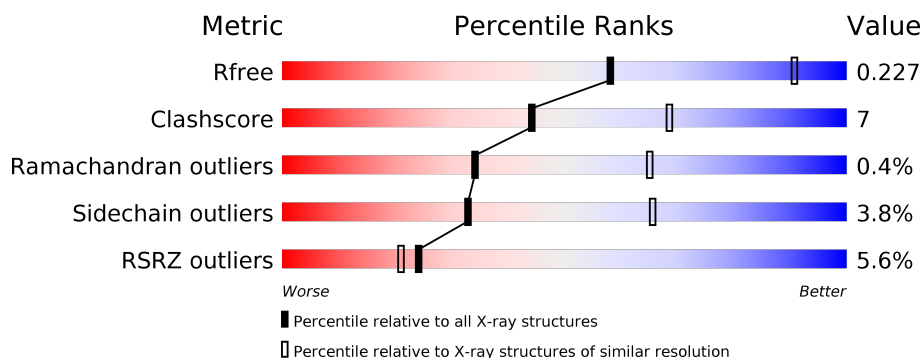
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.92 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))
$R_{free}$	130704	2307 (2.94-2.90)
Clashscore	141614	2531 (2.94-2.90)
Ramachandran outliers	138981	2462 (2.94-2.90)
Sidechain outliers	138945	2464 (2.94-2.90)
RSRZ outliers	127900	2248 (2.94-2.90)

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 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	1004	<div> <div>4%</div> <div> <div></div> <div>71%</div> <div>16%</div> <div>•</div> <div>12%</div> </div> </div>
2	C	548	<div> <div>7%</div> <div> <div></div> <div>78%</div> <div>16%</div> <div>• •</div> </div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11386 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 DNA (cytosine-5)-methyltransferase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	882	Total	C	N	O	S	0	1	0
			7033	4453	1240	1296	44			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	597	SER	-	expression tag	UNP P26358
A	598	GLU	-	expression tag	UNP P26358
A	599	PHE	-	expression tag	UNP P26358

- Molecule 2 is a protein called Ubiquitin carboxyl-terminal hydrolase 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	529	Total	C	N	O	S	0	0	0
			4340	2762	733	818	27			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	555	GLY	-	expression tag	UNP Q93009
C	556	PRO	-	expression tag	UNP Q93009
C	557	LEU	-	expression tag	UNP Q93009
C	558	GLY	-	expression tag	UNP Q93009
C	559	SER	-	expression tag	UNP Q93009

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		

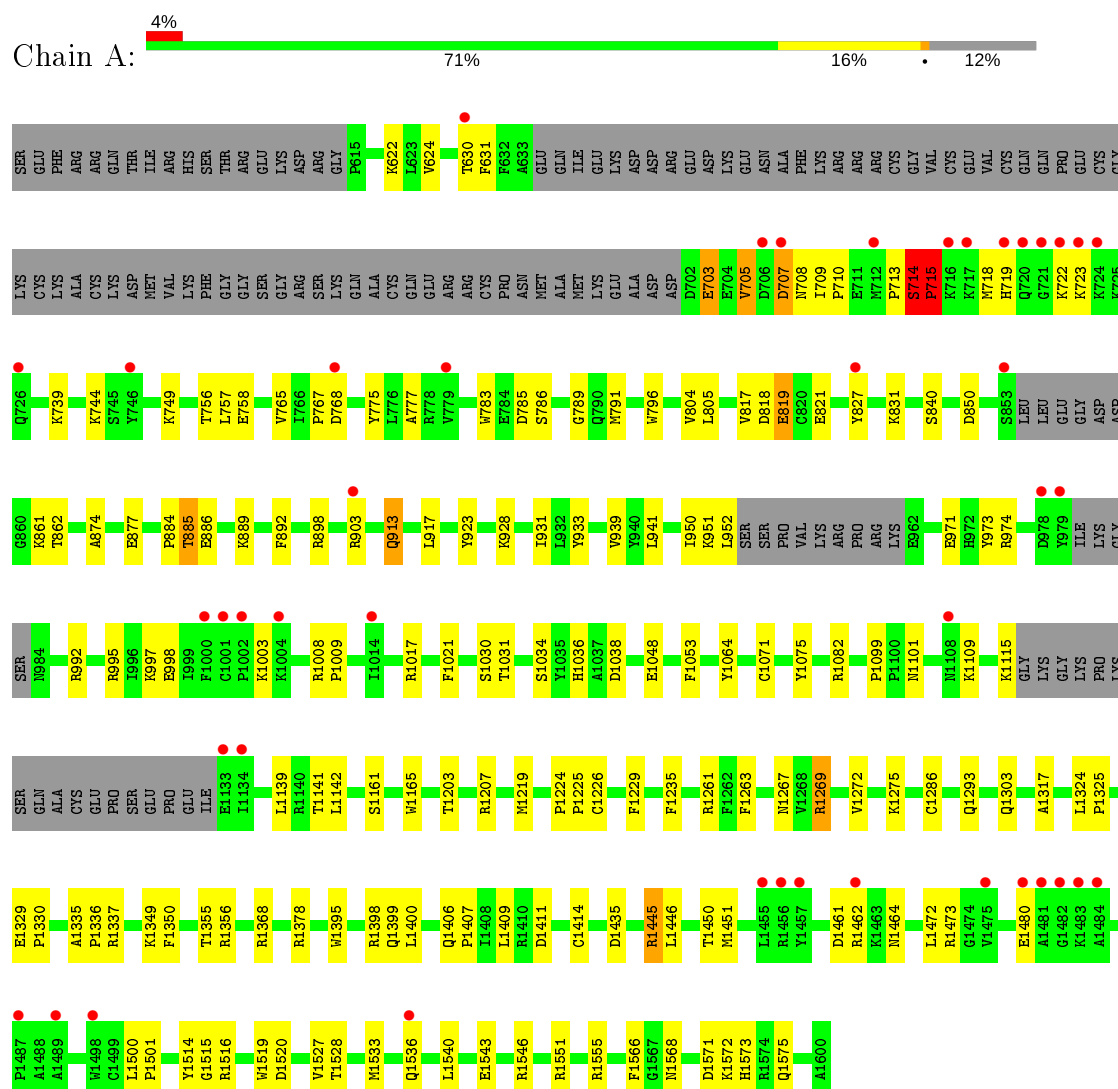
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	9	Total	O	0	0
			9	9		
4	C	2	Total	O	0	0
			2	2		

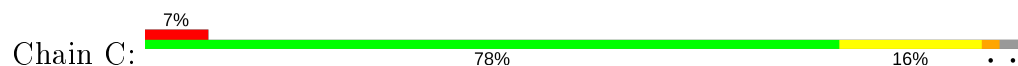
### 3 Residue-property plots [i](#)

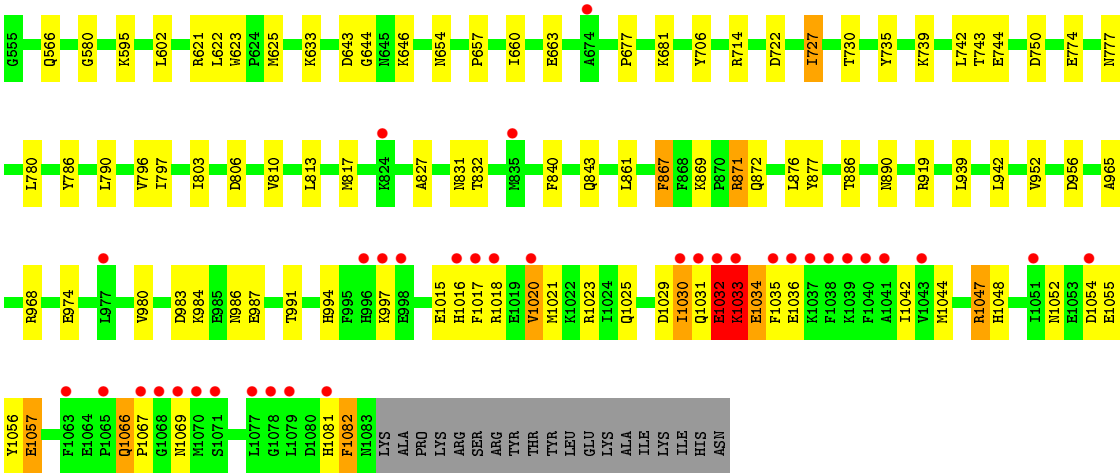
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: DNA (cytosine-5)-methyltransferase 1



- Molecule 2: Ubiquitin carboxyl-terminal hydrolase 7





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	110.23Å 111.62Å 163.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.42 – 2.92 49.42 – 2.92	Depositor EDS
% Data completeness (in resolution range)	99.2 (49.42-2.92) 99.2 (49.42-2.92)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.96 (at 2.91Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.208 , 0.258 0.211 , 0.227	Depositor DCC
$R_{free}$ test set	2247 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	61.7	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 48.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.007 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	11386	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.36% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical 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	0.24	0/7214	0.47	0/9755
2	C	0.24	0/4440	0.46	0/5995
All	All	0.24	0/11654	0.46	0/15750

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	0	5
2	C	0	2
All	All	0	7

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1536	GLN	Peptide
1	A	707	ASP	Peptide
1	A	708	ASN	Peptide
1	A	714	SER	Peptide
1	A	715	PRO	Peptide
2	C	1032	GLU	Peptide
2	C	1033	LYS	Peptide



## 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	7033	0	6901	90	0
2	C	4340	0	4252	59	0
3	A	2	0	0	0	0
4	A	9	0	0	0	0
4	C	2	0	0	0	0
All	All	11386	0	11153	148	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 7.

All (148) 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:1036:HIS:O	1:A:1398:ARG:NH2	2.20	0.74
2:C:1033:LYS:HE3	2:C:1036:GLU:HB2	1.76	0.67
2:C:867:PHE:HD1	2:C:867:PHE:H	1.42	0.67
2:C:625:MET:HG3	2:C:633:LYS:HG2	1.79	0.65
2:C:1016:HIS:HA	2:C:1056:TYR:HB3	1.79	0.64
2:C:919:ARG:NH2	2:C:956:ASP:OD1	2.31	0.64
1:A:622:LYS:HD2	1:A:850:ASP:HB3	1.82	0.62
1:A:1038:ASP:OD1	1:A:1398:ARG:NH1	2.32	0.62
1:A:886:GLU:HA	1:A:889:LYS:HG3	1.82	0.61
2:C:965:ALA:HB3	2:C:968:ARG:HG3	1.83	0.61
1:A:884:PRO:HB3	1:A:892:PHE:CG	2.35	0.61
1:A:1368:ARG:NH2	1:A:1520:ASP:OD1	2.31	0.60
2:C:1017:PHE:O	2:C:1020:VAL:HG12	2.01	0.60
2:C:797:ILE:HG12	2:C:810:VAL:HG22	1.83	0.59
2:C:1015:GLU:HG2	2:C:1020:VAL:HB	1.84	0.59
1:A:997:LYS:NZ	1:A:1048:GLU:OE1	2.33	0.58
1:A:703:GLU:HG3	1:A:705:VAL:H	1.68	0.58
1:A:818:ASP:OD2	1:A:898:ARG:NH1	2.37	0.57
2:C:1031:GLN:HB3	2:C:1034:GLU:HB3	1.87	0.57
1:A:928:LYS:NZ	1:A:1053:PHE:O	2.36	0.57
2:C:1052:ASN:ND2	2:C:1055:GLU:OE2	2.33	0.57
2:C:869:LYS:HB2	2:C:872:GLN:HG3	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1400:LEU:HD23	1:A:1555:ARG:HD3	1.87	0.56
1:A:1411:ASP:OD1	1:A:1572:LYS:NZ	2.34	0.56
1:A:1461:ASP:OD2	1:A:1464:ASN:ND2	2.40	0.55
1:A:707:ASP:OD1	1:A:707:ASP:N	2.40	0.55
1:A:1109:LYS:NZ	2:C:744:GLU:OE1	2.40	0.55
2:C:1044:MET:HG3	2:C:1047:ARG:HG2	1.89	0.54
2:C:983:ASP:HB3	2:C:986:ASN:HB2	1.88	0.54
1:A:998:GLU:HB2	1:A:1017:ARG:HB3	1.89	0.54
1:A:877:GLU:HA	1:A:1293:GLN:HB3	1.89	0.53
1:A:1472:LEU:H	1:A:1472:LEU:HD12	1.73	0.53
1:A:714:SER:CB	1:A:1349:LYS:HB3	2.38	0.53
2:C:643:ASP:OD1	2:C:646:LYS:NZ	2.41	0.53
1:A:1519:TRP:NE1	1:A:1543:GLU:OE2	2.36	0.53
1:A:785:ASP:O	1:A:789:GLY:N	2.35	0.52
2:C:1082:PHE:N	2:C:1082:PHE:HD1	2.07	0.52
1:A:796:TRP:HE1	1:A:821:GLU:HG2	1.73	0.52
1:A:749:LYS:HD3	1:A:758:GLU:HG2	1.91	0.52
1:A:714:SER:HB2	1:A:1349:LYS:HB3	1.92	0.52
1:A:917:LEU:O	1:A:1003:LYS:NZ	2.39	0.52
2:C:1066:GLN:NE2	2:C:1069:ASN:OD1	2.43	0.51
1:A:1235:PHE:CE1	1:A:1275:LYS:HG3	2.45	0.51
2:C:1020:VAL:HG23	2:C:1023:ARG:NH1	2.26	0.51
2:C:774:GLU:OE2	2:C:777:ASN:ND2	2.42	0.51
1:A:1030:SER:OG	1:A:1031:THR:N	2.43	0.51
2:C:840:PHE:HB2	2:C:877:TYR:HB2	1.91	0.51
2:C:663:GLU:OE1	2:C:706:TYR:OH	2.29	0.51
2:C:1082:PHE:N	2:C:1082:PHE:CD1	2.79	0.51
1:A:1446:LEU:HD12	1:A:1450:THR:HB	1.93	0.51
2:C:1031:GLN:OE1	2:C:1032:GLU:N	2.44	0.50
2:C:869:LYS:HB3	2:C:871:ARG:HE	1.77	0.50
1:A:1263:PHE:HB3	1:A:1317:ALA:HB3	1.93	0.50
2:C:1057:GLU:OE1	2:C:1057:GLU:N	2.35	0.50
1:A:913:GLN:HA	1:A:923:TYR:HA	1.94	0.50
1:A:1329:GLU:OE2	1:A:1355:THR:N	2.45	0.49
1:A:1516:ARG:HA	1:A:1540:LEU:HB2	1.94	0.49
1:A:1224:PRO:O	1:A:1267:ASN:ND2	2.45	0.49
2:C:566:GLN:HB2	2:C:657:PRO:HB3	1.93	0.49
1:A:928:LYS:O	1:A:931:ILE:HG12	2.13	0.49
1:A:1528:THR:O	1:A:1573:HIS:HB3	2.12	0.49
1:A:874:ALA:HB2	1:A:1350:PHE:CD1	2.48	0.48
1:A:1008:ARG:HG3	1:A:1009:PRO:HD2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1075:TYR:CZ	1:A:1082:ARG:HB3	2.48	0.48
2:C:942:LEU:HD13	2:C:952:VAL:HG22	1.96	0.48
1:A:709:ILE:HD12	1:A:710:PRO:HD2	1.95	0.48
1:A:1445:ARG:HG2	1:A:1451:MET:SD	2.54	0.48
1:A:1500:LEU:HB2	1:A:1501:PRO:HD3	1.96	0.47
1:A:1414:CYS:SG	1:A:1546:ARG:HD2	2.54	0.47
1:A:1395:TRP:O	1:A:1399:GLN:HG2	2.13	0.47
1:A:1462:ARG:HH12	1:A:1480:GLU:CD	2.18	0.47
2:C:983:ASP:O	2:C:987:GLU:N	2.46	0.47
1:A:1527:VAL:HG22	1:A:1528:THR:H	1.80	0.47
1:A:765:VAL:HG13	1:A:827:TYR:O	2.15	0.47
1:A:885:THR:OG1	1:A:886:GLU:N	2.48	0.47
1:A:1409:LEU:HD11	1:A:1551:ARG:HG2	1.97	0.47
2:C:1025:GLN:HG3	2:C:1035:PHE:CE2	2.50	0.47
2:C:1018:ARG:HA	2:C:1021:MET:HG2	1.97	0.46
2:C:780:LEU:HD13	2:C:786:TYR:HA	1.97	0.46
1:A:1031:THR:O	1:A:1034:SER:OG	2.31	0.46
1:A:1141:THR:HG23	1:A:1219:MET:HG2	1.98	0.46
2:C:997:LYS:HG3	2:C:1082:PHE:CZ	2.51	0.46
2:C:1042:ILE:O	2:C:1048:HIS:HA	2.15	0.46
2:C:939:LEU:HD23	2:C:974:GLU:HA	1.97	0.45
1:A:713:PRO:HG2	1:A:714:SER:H	1.82	0.45
2:C:735:TYR:HB3	2:C:743:THR:HG22	1.98	0.45
1:A:1139:LEU:HD21	1:A:1261:ARG:HH21	1.81	0.45
2:C:714:ARG:NH1	2:C:750:ASP:OD1	2.48	0.45
1:A:1330:PRO:HD2	1:A:1356:ARG:HB2	1.98	0.45
2:C:621:ARG:HG2	2:C:623:TRP:NE1	2.32	0.45
2:C:621:ARG:NH2	2:C:677:PRO:O	2.40	0.45
1:A:624:VAL:HG22	1:A:1286:CYS:SG	2.57	0.45
1:A:971:GLU:OE1	1:A:974:ARG:HD2	2.16	0.44
2:C:739:LYS:HG2	2:C:742:LEU:HB3	1.99	0.44
2:C:803:ILE:HB	2:C:806:ASP:HB2	1.99	0.44
2:C:1030:ILE:HG13	2:C:1031:GLN:H	1.82	0.44
1:A:709:ILE:HA	1:A:710:PRO:HD2	1.78	0.44
1:A:777:ALA:HB2	1:A:796:TRP:CE3	2.52	0.44
2:C:602:LEU:HB3	2:C:644:GLY:HA2	1.99	0.44
2:C:867:PHE:CD1	2:C:867:PHE:N	2.84	0.44
1:A:714:SER:HB3	1:A:715:PRO:HD3	1.99	0.44
2:C:1066:GLN:HB2	2:C:1067:PRO:HD2	1.98	0.44
1:A:1568:ASN:ND2	1:A:1571:ASP:OD2	2.51	0.44
1:A:1269:ARG:O	1:A:1272:VAL:HG22	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:767:PRO:HB3	1:A:775:TYR:HE1	1.82	0.43
2:C:1020:VAL:HG23	2:C:1023:ARG:HH12	1.83	0.43
1:A:1224:PRO:HA	1:A:1225:PRO:HD3	1.88	0.43
1:A:1435:ASP:HA	1:A:1515:GLY:HA2	2.00	0.43
2:C:1015:GLU:O	2:C:1016:HIS:HB3	2.18	0.43
1:A:1335:ALA:HA	1:A:1336:PRO:HD3	1.85	0.43
2:C:1033:LYS:HA	2:C:1033:LYS:HD2	1.68	0.43
2:C:796:VAL:HG21	2:C:861:LEU:HD21	2.00	0.43
1:A:1324:LEU:HD12	1:A:1325:PRO:HD2	2.00	0.43
1:A:723:LYS:HA	1:A:768:ASP:HB3	2.01	0.43
1:A:931:ILE:HG13	1:A:933:TYR:CE1	2.54	0.43
2:C:827:ALA:HB1	2:C:832:THR:O	2.18	0.43
1:A:1064:TYR:CE1	1:A:1099:PRO:HG2	2.54	0.42
1:A:719:HIS:CE1	1:A:796:TRP:HB3	2.53	0.42
1:A:1303:GLN:HB3	1:A:1330:PRO:HB3	2.01	0.42
1:A:757:LEU:HD23	1:A:831:LYS:HD2	2.01	0.42
2:C:580:GLY:HA2	2:C:790:LEU:HD21	2.00	0.42
1:A:973:TYR:CZ	1:A:1473:ARG:HD3	2.55	0.42
1:A:722:LYS:O	1:A:723:LYS:HG3	2.20	0.42
1:A:941:LEU:HB2	1:A:992:ARG:HB2	2.01	0.42
2:C:727:ILE:O	2:C:730:THR:OG1	2.23	0.42
1:A:1003:LYS:HA	1:A:1009:PRO:HA	2.01	0.42
1:A:1141:THR:OG1	1:A:1161:SER:HB2	2.20	0.42
2:C:1031:GLN:OE1	2:C:1034:GLU:N	2.45	0.42
2:C:622:LEU:HG	2:C:660:ILE:HG21	2.02	0.42
2:C:813:LEU:HB3	2:C:817:MET:SD	2.60	0.42
1:A:1514:TYR:CE2	1:A:1533:MET:HG3	2.55	0.41
1:A:631:PHE:O	1:A:631:PHE:CG	2.73	0.41
1:A:1406:GLN:HA	1:A:1407:PRO:HD2	1.91	0.41
1:A:1566:PHE:O	1:A:1575:GLN:NE2	2.53	0.41
1:A:723:LYS:HG2	1:A:767:PRO:HA	2.02	0.41
1:A:995:ARG:HB2	1:A:1021:PHE:HE2	1.84	0.41
1:A:1226:CYS:HA	1:A:1229:PHE:CZ	2.56	0.41
2:C:566:GLN:OE1	2:C:595:LYS:HE2	2.20	0.41
2:C:1029:ASP:HA	2:C:1035:PHE:HZ	1.86	0.41
2:C:886:THR:O	2:C:890:ASN:ND2	2.52	0.41
2:C:869:LYS:HB3	2:C:871:ARG:NE	2.36	0.41
1:A:1142:LEU:HD13	1:A:1165:TRP:HB2	2.03	0.40
1:A:739:LYS:HA	1:A:739:LYS:HD2	1.79	0.40
1:A:744:LYS:HB3	1:A:783:TRP:CD1	2.57	0.40
1:A:783:TRP:CE2	1:A:791:MET:HB2	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:817:VAL:HG23	1:A:819:GLU:HB2	2.03	0.40
1:A:850:ASP:OD1	1:A:850:ASP:N	2.39	0.40
1:A:951:LYS:HE3	1:A:951:LYS:HB3	1.80	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	871/1004 (87%)	830 (95%)	39 (4%)	2 (0%)	47	77
2	C	527/548 (96%)	499 (95%)	25 (5%)	3 (1%)	25	57
All	All	1398/1552 (90%)	1329 (95%)	64 (5%)	5 (0%)	34	65

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	1033	LYS
2	C	1081	HIS
1	A	819	GLU
2	C	980	VAL
1	A	705	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	764/869 (88%)	737 (96%)	27 (4%)	36	68
2	C	484/501 (97%)	463 (96%)	21 (4%)	29	61
All	All	1248/1370 (91%)	1200 (96%)	48 (4%)	33	65

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	630	THR
1	A	703	GLU
1	A	714	SER
1	A	715	PRO
1	A	718	MET
1	A	756	THR
1	A	786	SER
1	A	804	VAL
1	A	805	LEU
1	A	840	SER
1	A	861	LYS
1	A	862	THR
1	A	885	THR
1	A	903	ARG
1	A	913	GLN
1	A	939	VAL
1	A	950	ILE
1	A	952	LEU
1	A	1071	CYS
1	A	1101	ASN
1	A	1115	LYS
1	A	1203	THR
1	A	1207	ARG
1	A	1269	ARG
1	A	1337	ARG
1	A	1378	ARG
1	A	1445	ARG
2	C	654	ASN
2	C	681	LYS
2	C	722	ASP
2	C	727	ILE
2	C	831	ASN
2	C	843	GLN
2	C	867	PHE
2	C	871	ARG

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Mol	Chain	Res	Type
2	C	876	LEU
2	C	984	LYS
2	C	991	THR
2	C	994	HIS
2	C	1020	VAL
2	C	1030	ILE
2	C	1032	GLU
2	C	1034	GLU
2	C	1047	ARG
2	C	1054	ASP
2	C	1057	GLU
2	C	1066	GLN
2	C	1082	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	C	1066	GLN
2	C	1069	ASN
2	C	1072	HIS
2	C	1081	HIS

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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

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	A	882/1004 (87%)	0.30	43 (4%) 29 26	25, 54, 116, 158	0
2	C	529/548 (96%)	0.38	36 (6%) 17 14	31, 61, 136, 173	0
All	All	1411/1552 (90%)	0.33	79 (5%) 24 21	25, 58, 127, 173	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	721	GLY	12.8
2	C	996	HIS	8.5
1	A	720	GLN	5.1
1	A	1483	LYS	5.0
2	C	1067	PRO	4.9
2	C	1071	SER	4.6
1	A	723	LYS	4.5
1	A	979	TYR	4.2
1	A	724	LYS	4.1
1	A	722	LYS	4.1
1	A	978	ASP	4.0
2	C	1070	MET	4.0
2	C	1079	LEU	3.9
1	A	1484	ALA	3.9
2	C	1039	LYS	3.8
2	C	1063	PHE	3.8
1	A	1480	GLU	3.6
2	C	1033	LYS	3.6
1	A	1134	ILE	3.5
1	A	827	TYR	3.5
2	C	674	ALA	3.5
1	A	746	TYR	3.3
2	C	1032	GLU	3.3
2	C	1031	GLN	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	1001	CYS	3.3
2	C	1051	ILE	3.2
1	A	1133	GLU	3.2
1	A	1004	LYS	3.2
2	C	1078	GLY	3.2
1	A	768	ASP	3.1
2	C	1017	PHE	3.1
2	C	1016	HIS	3.1
1	A	1481	ALA	3.0
2	C	1018	ARG	3.0
2	C	1040	PHE	2.9
2	C	998	GLU	2.9
2	C	1037	LYS	2.9
2	C	977	LEU	2.9
2	C	1065	PRO	2.9
1	A	1108	ASN	2.9
2	C	1038	PHE	2.9
2	C	1035	PHE	2.8
1	A	1457	TYR	2.8
1	A	717	LYS	2.7
1	A	1475	VAL	2.7
1	A	712	MET	2.7
1	A	707	ASP	2.7
2	C	1036	GLU	2.7
1	A	716	LYS	2.7
1	A	1462	ARG	2.7
1	A	706	ASP	2.6
1	A	1000	PHE	2.6
2	C	1068	GLY	2.6
1	A	1014	ILE	2.6
1	A	1482	GLY	2.6
1	A	1002	PRO	2.6
2	C	1041	ALA	2.6
2	C	1077	LEU	2.5
2	C	1043	VAL	2.5
1	A	726	GLN	2.5
1	A	853	SER	2.4
2	C	1054	ASP	2.4
2	C	1081	HIS	2.4
2	C	824	LYS	2.4
2	C	1069	ASN	2.4
1	A	1487	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	1456	ARG	2.4
2	C	1030	ILE	2.4
2	C	1020	VAL	2.3
1	A	630	THR	2.3
1	A	1489	ALA	2.2
1	A	779	VAL	2.1
1	A	1455	LEU	2.1
1	A	719	HIS	2.1
1	A	1498	TRP	2.1
1	A	1536	GLN	2.1
1	A	903	ARG	2.0
2	C	997	LYS	2.0
2	C	835	MET	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 carbohydrates 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, 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	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ZN	A	1701	1/1	0.97	0.07	76,76,76,76	0
3	ZN	A	1702	1/1	0.98	0.17	59,59,59,59	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.