



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 16, 2024 – 07:14 AM EDT

PDB ID : 4ZHJ  
Title : Crystal Structure of the Catalytic Subunit of Magnesium Chelatase  
Authors : Chen, X.; Pu, H.; Fang, Y.; Liu, L.  
Deposited on : 2015-04-25  
Resolution : 2.50 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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	:	2.37.1
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)	:	2.37.1

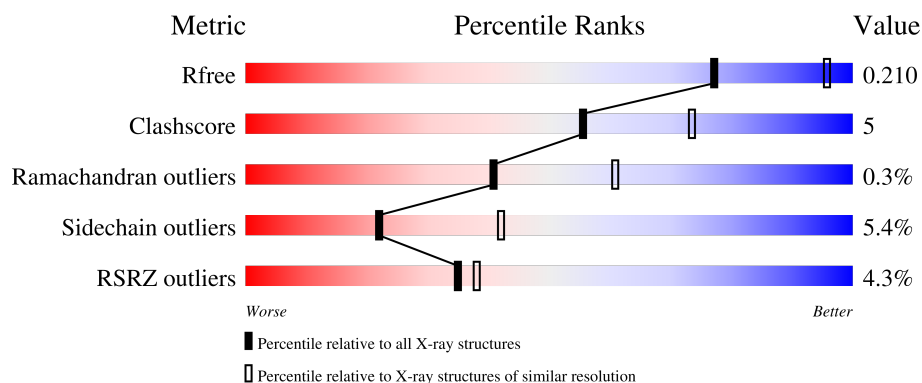
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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  $\geq 3$ , 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	1351	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>12%</div> <div>• 7%</div> </div> </div>
1	B	1351	<div> <div>6%</div> <div> <div></div> <div>76%</div> <div>14%</div> <div>• 9%</div> </div> </div>

## 2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 20168 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 Mg-chelatase subunit ChlH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1258	Total	C	N	O	S	0	3	0
			9817	6230	1664	1875	48			
1	B	1233	Total	C	N	O	S	0	5	0
			9471	6019	1614	1790	48			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P73020
A	-18	GLY	-	expression tag	UNP P73020
A	-17	SER	-	expression tag	UNP P73020
A	-16	SER	-	expression tag	UNP P73020
A	-15	HIS	-	expression tag	UNP P73020
A	-14	HIS	-	expression tag	UNP P73020
A	-13	HIS	-	expression tag	UNP P73020
A	-12	HIS	-	expression tag	UNP P73020
A	-11	HIS	-	expression tag	UNP P73020
A	-10	HIS	-	expression tag	UNP P73020
A	-9	SER	-	expression tag	UNP P73020
A	-8	SER	-	expression tag	UNP P73020
A	-7	GLY	-	expression tag	UNP P73020
A	-6	LEU	-	expression tag	UNP P73020
A	-5	VAL	-	expression tag	UNP P73020
A	-4	PRO	-	expression tag	UNP P73020
A	-3	ARG	-	expression tag	UNP P73020
A	-2	GLY	-	expression tag	UNP P73020
A	-1	SER	-	expression tag	UNP P73020
A	0	HIS	-	expression tag	UNP P73020
B	-19	MET	-	expression tag	UNP P73020
B	-18	GLY	-	expression tag	UNP P73020
B	-17	SER	-	expression tag	UNP P73020
B	-16	SER	-	expression tag	UNP P73020
B	-15	HIS	-	expression tag	UNP P73020

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP P73020
B	-13	HIS	-	expression tag	UNP P73020
B	-12	HIS	-	expression tag	UNP P73020
B	-11	HIS	-	expression tag	UNP P73020
B	-10	HIS	-	expression tag	UNP P73020
B	-9	SER	-	expression tag	UNP P73020
B	-8	SER	-	expression tag	UNP P73020
B	-7	GLY	-	expression tag	UNP P73020
B	-6	LEU	-	expression tag	UNP P73020
B	-5	VAL	-	expression tag	UNP P73020
B	-4	PRO	-	expression tag	UNP P73020
B	-3	ARG	-	expression tag	UNP P73020
B	-2	GLY	-	expression tag	UNP P73020
B	-1	SER	-	expression tag	UNP P73020
B	0	HIS	-	expression tag	UNP P73020

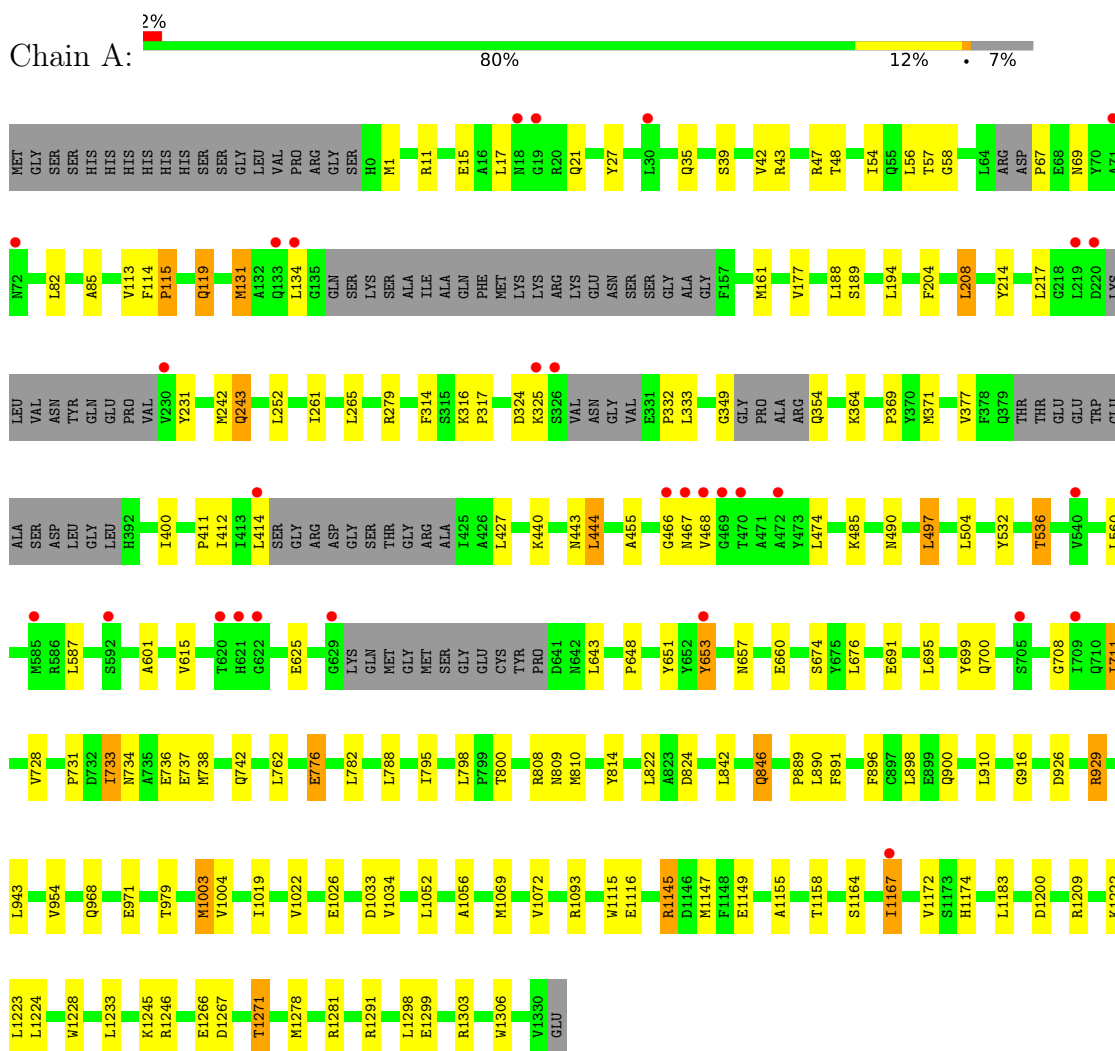
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	603	Total O 603 603	0	0
2	B	277	Total O 277 277	0	0

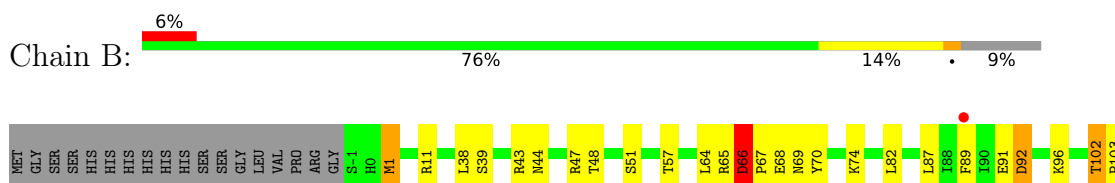
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Mg-chelatase subunit ChIH



#### • Molecule 1: Mg-chelatase subunit ChIH





## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	319.72Å 319.72Å 105.19Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.96 – 2.50 44.96 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.1 (44.96-2.50) 97.1 (44.96-2.50)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.51 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, $R_{free}$	0.177 , 0.209 0.179 , 0.210	Depositor DCC
$R_{free}$ test set	6735 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.7	Xtriage
Anisotropy	0.083	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 50.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	20168	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.10% 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](#)

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	0/10021	0.44	1/13604 (0.0%)
1	B	0.23	0/9675	0.47	10/13164 (0.1%)
All	All	0.24	0/19696	0.46	11/26768 (0.0%)

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	B	0	2

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	138	LYS	CB-CA-C	7.05	124.51	110.40
1	B	792	ASP	N-CA-C	6.59	128.80	111.00
1	B	1141	MET	N-CA-C	-5.68	95.67	111.00
1	B	1222	LYS	N-CA-C	5.59	126.10	111.00
1	B	792	ASP	CB-CA-C	-5.42	99.57	110.40
1	B	1222	LYS	CB-CA-C	-5.34	99.73	110.40
1	B	582	GLY	N-CA-C	5.32	126.41	113.10
1	B	1168	SER	N-CA-CB	5.28	118.42	110.50
1	B	516	TYR	CB-CA-C	-5.22	99.95	110.40
1	B	793	GLU	N-CA-C	5.20	125.04	111.00
1	A	1167	ILE	N-CA-C	-5.10	97.23	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:



Mol	Chain	Res	Type	Group
1	B	138	LYS	Peptide
1	B	792	ASP	Peptide

## 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	9817	0	9630	83	0
1	B	9471	0	9182	112	0
2	A	603	0	0	10	0
2	B	277	0	0	5	0
All	All	20168	0	18812	195	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 5.

All (195) 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:B:1222:LYS:O	1:B:1228:TRP:CE3	2.11	1.02
1:B:66:ASP:CB	1:B:67:PRO:HA	2.00	0.90
1:B:66:ASP:CG	1:B:67:PRO:HA	2.00	0.82
1:B:67:PRO:O	1:B:69:ASN:N	2.13	0.81
1:B:883:ASN:O	1:B:887:LEU:HD23	1.80	0.80
1:A:1299:GLU:OE2	1:A:1303:ARG:NH1	2.16	0.78
1:A:1003:MET:SD	2:A:1958:HOH:O	2.43	0.77
1:B:679:PRO:HG3	1:B:773:PRO:HB3	1.68	0.76
1:B:1003:MET:SD	2:B:1657:HOH:O	2.44	0.75
1:B:126:MET:SD	2:B:1659:HOH:O	2.47	0.71
1:A:1115:TRP:O	1:A:1303:ARG:NH2	2.25	0.70
1:B:66:ASP:HB3	1:B:67:PRO:HA	1.73	0.69
1:B:1115:TRP:O	1:B:1303:ARG:NH2	2.26	0.69
1:A:467:ASN:HB2	1:A:1245:LYS:HG3	1.77	0.67
1:A:1116:GLU:O	2:A:1401:HOH:O	2.13	0.66
1:B:340:VAL:HG22	1:B:371:MET:HE2	1.78	0.66
1:B:738:MET:HB3	1:B:742:GLN:HB3	1.77	0.66
1:B:973:GLY:HA3	1:B:975:ASN:H	1.61	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:PRO:C	1:B:69:ASN:H	1.97	0.65
1:B:883:ASN:O	1:B:887:LEU:CD2	2.46	0.64
1:A:231:TYR:HB3	1:A:279:ARG:HE	1.61	0.64
1:A:808:ARG:NH2	1:A:824:ASP:OD1	2.29	0.64
1:A:1222:LYS:O	1:A:1228:TRP:CE3	2.52	0.63
1:B:808:ARG:NH2	1:B:824:ASP:OD1	2.28	0.62
1:A:261:ILE:HB	1:A:265:LEU:HD12	1.81	0.61
1:B:66:ASP:CB	1:B:67:PRO:CA	2.78	0.61
1:A:1164:SER:O	1:A:1167:ILE:O	2.19	0.60
1:A:1200:ASP:OD1	1:A:1209:ARG:NH1	2.35	0.60
1:B:1222:LYS:O	1:B:1228:TRP:CD2	2.55	0.60
1:A:455:ALA:HB3	1:A:615[A]:VAL:HG22	1.85	0.59
1:B:926:ASP:HB2	1:B:991:LYS:HD3	1.85	0.59
1:B:89:PHE:HA	1:B:117:MET:HG3	1.85	0.59
1:B:1222:LYS:O	1:B:1228:TRP:HE3	1.80	0.58
1:B:1299:GLU:OE2	1:B:1303:ARG:NH1	2.36	0.57
1:B:1164:SER:O	1:B:1167:ILE:O	2.22	0.57
1:A:485:LYS:HG2	1:A:497:LEU:HD11	1.86	0.57
1:B:699:TYR:O	1:B:701:THR:N	2.37	0.57
1:A:532:TYR:O	1:A:536:THR:HG23	2.05	0.56
1:B:1066:PRO:HG2	1:B:1069:MET:HG3	1.87	0.56
1:A:708:GLY:HA2	1:A:711:ILE:HD11	1.87	0.56
1:A:466:GLY:O	2:A:1402:HOH:O	2.18	0.56
1:B:588:LEU:HD21	1:B:1124:MET:HG3	1.88	0.56
1:B:1033:ASP:HA	1:B:1072:VAL:HG22	1.88	0.56
1:B:217:LEU:O	1:B:219:LEU:N	2.38	0.55
1:B:779:ILE:HG12	1:B:843:VAL:HG21	1.89	0.55
1:A:846:GLN:NE2	2:A:1431:HOH:O	2.40	0.54
1:B:797:ALA:HB3	1:B:800:THR:HG23	1.90	0.54
1:A:324:ASP:HB2	1:A:333:LEU:H	1.73	0.54
1:A:1267:ASP:O	1:A:1271:THR:HG23	2.08	0.54
1:A:214:TYR:HB3	1:A:217:LEU:HD13	1.90	0.53
1:B:1299:GLU:O	1:B:1303:ARG:HG3	2.09	0.53
1:B:82:LEU:HD21	1:B:208:LEU:HD23	1.91	0.53
1:B:973:GLY:HA3	1:B:975:ASN:N	2.23	0.53
1:A:371:MET:HE2	1:A:411:PRO:HB3	1.91	0.52
1:B:1169:LEU:HG	1:B:1253:TRP:CZ2	2.45	0.52
1:A:1266:GLU:HG2	1:A:1306:TRP:HE1	1.74	0.51
1:A:11:ARG:HG3	1:A:57:THR:HG22	1.93	0.51
1:B:377:VAL:HG21	1:B:412:ILE:HG21	1.93	0.51
1:A:114:PHE:HB3	2:A:1957:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:658:PRO:O	1:B:662:THR:HG22	2.09	0.51
1:B:47:ARG:HG3	1:B:48:THR:HG23	1.91	0.51
1:B:1003:MET:O	1:B:1029:ARG:NH2	2.44	0.51
1:A:377:VAL:HG21	1:A:412:ILE:HG21	1.92	0.51
1:A:653:TYR:O	1:A:674:SER:HA	2.11	0.51
1:B:579:GLY:H	1:B:580:TYR:HA	1.74	0.51
1:B:87:LEU:HD22	1:B:158:GLN:HG3	1.93	0.51
1:A:731:PRO:HB2	1:A:733:ILE:HG23	1.93	0.50
1:A:27:TYR:CZ	1:A:58:GLY:HA3	2.47	0.50
1:A:1022:VAL:HG13	1:A:1026:GLU:HB3	1.94	0.50
1:A:316:LYS:HB2	1:A:317:PRO:HD3	1.94	0.50
1:B:532:TYR:O	1:B:536:THR:HG23	2.12	0.50
1:A:115:PRO:HG2	1:A:161:MET:HG2	1.92	0.50
1:B:92:ASP:OD1	1:B:92:ASP:N	2.43	0.50
1:A:1145:ARG:HD3	1:A:1149:GLU:OE2	2.11	0.50
1:B:129:PHE:HA	1:B:133:GLN:HG3	1.92	0.50
1:A:377:VAL:HG22	1:A:414:LEU:HD21	1.93	0.49
1:A:691:GLU:OE2	2:A:1403:HOH:O	2.20	0.49
1:B:82:LEU:HD11	1:B:112:ILE:HD13	1.94	0.49
1:B:70:TYR:OH	1:B:96:LYS:O	2.26	0.49
1:B:44:ASN:HD22	1:B:202:GLU:CD	2.16	0.49
1:B:65:ARG:C	1:B:66:ASP:O	2.49	0.49
1:A:625:GLU:OE1	1:A:651:TYR:OH	2.21	0.49
1:B:1022:VAL:HG13	1:B:1026:GLU:HB3	1.94	0.48
1:B:1172:VAL:HG21	1:B:1174:HIS:CE1	2.48	0.48
1:A:85:ALA:HB3	1:A:113:VAL:HG22	1.96	0.48
1:A:189:SER:HB2	1:A:204:PHE:CE1	2.48	0.48
1:B:625:GLU:OE2	1:B:651:TYR:OH	2.23	0.48
1:B:684:GLY:HA2	1:B:758:GLU:HG2	1.95	0.48
1:B:685:LEU:HD23	1:B:754:LEU:HD12	1.95	0.48
1:A:440:LYS:HB3	1:A:648:PRO:HD3	1.94	0.47
1:B:11:ARG:HG3	1:B:57:THR:HG22	1.95	0.47
1:A:926:ASP:HB3	1:A:929:ARG:HG2	1.97	0.47
1:A:1019:ILE:HD13	1:A:1056:ALA:HB2	1.96	0.47
1:A:325:LYS:H	1:A:332:PRO:HB3	1.80	0.47
1:B:653:TYR:O	1:B:674:SER:HA	2.15	0.47
1:A:738:MET:HB3	1:A:742:GLN:HG3	1.96	0.47
1:B:1164:SER:HA	1:B:1202:THR:HG21	1.97	0.47
1:B:1003:MET:HE2	1:B:1034:VAL:HG12	1.97	0.47
1:B:1200:ASP:OD1	1:B:1209:ARG:NH1	2.48	0.47
1:B:138:LYS:HA	1:B:139:SER:HA	1.54	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:969:ARG:O	1:B:974:GLY:HA2	2.15	0.46
1:A:349:GLY:HA2	1:A:354:GLN:HA	1.97	0.46
1:B:1278:MET:HE1	1:B:1281:ARG:NE	2.31	0.46
2:A:2002:HOH:O	1:B:1:MET:HE1	2.15	0.46
1:A:657:ASN:HD21	1:A:660:GLU:HG3	1.81	0.46
1:A:810:MET:HG3	1:A:814:TYR:CE2	2.50	0.46
1:B:840:ALA:N	2:B:1404:HOH:O	2.26	0.46
1:B:739:ASP:OD1	1:B:739:ASP:N	2.47	0.45
1:B:555:SER:HB3	1:B:556:ASP:C	2.36	0.45
1:B:51:SER:HB3	1:B:217:LEU:HD11	1.98	0.45
1:B:665:LYS:NZ	1:B:672:THR:OG1	2.49	0.45
1:B:102:THR:HG22	1:B:103:PRO:HD3	1.99	0.45
1:B:698:SER:HA	1:B:699:TYR:C	2.36	0.45
1:A:231:TYR:HB2	1:A:279:ARG:HH21	1.81	0.45
1:A:177:VAL:HG22	1:A:427:LEU:HD21	1.98	0.45
1:B:236:ILE:HB	1:B:245:PHE:HB2	1.98	0.44
1:B:678:PRO:HA	1:B:679:PRO:HD3	1.88	0.44
1:B:897:CYS:O	1:B:901:VAL:HG23	2.17	0.44
1:B:70:TYR:CE1	1:B:74:LYS:HE3	2.52	0.44
1:B:780:ALA:O	1:B:783:VAL:HG12	2.17	0.44
1:A:27:TYR:OH	1:A:35:GLN:HG3	2.18	0.44
1:B:189:SER:HB2	1:B:204:PHE:CE1	2.52	0.44
1:B:605:TYR:HA	1:B:609:ILE:HD13	2.00	0.44
1:A:369:PRO:HD3	1:A:916:GLY:O	2.17	0.44
1:B:350:GLY:O	1:B:352:ALA:HA	2.18	0.44
1:A:1222:LYS:O	1:A:1228:TRP:HB3	2.18	0.43
1:B:369:PRO:HD3	1:B:916:GLY:O	2.18	0.43
1:A:188:LEU:HD12	1:A:188:LEU:HA	1.87	0.43
1:A:1278:MET:HE3	1:A:1281:ARG:NE	2.33	0.43
1:B:379:GLN:NE2	2:B:1405:HOH:O	2.26	0.43
1:A:1033:ASP:HA	1:A:1072:VAL:HG22	1.98	0.43
1:B:1291:ARG:HD3	1:B:1327:ILE:HD12	1.99	0.43
1:B:967:ARG:HH22	1:B:1213:GLU:CD	2.22	0.43
1:A:21:GLN:HG3	1:A:217:LEU:HD11	1.99	0.43
1:A:560:LEU:HD13	1:A:601:ALA:HB2	2.01	0.43
1:B:523:ILE:HD13	1:B:561:LEU:HD22	2.01	0.43
1:B:1018:LYS:HB2	1:B:1018:LYS:HE3	1.77	0.43
1:A:42:VAL:HG13	1:A:54:ILE:HB	2.00	0.43
1:A:82:LEU:HD21	1:A:208:LEU:HB3	2.00	0.43
1:B:211:LYS:HB3	1:B:212:TYR:CD2	2.54	0.43
1:A:896:PHE:O	1:A:900:GLN:HG2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1156:ASP:OD2	1:B:1188:ARG:NH2	2.52	0.42
1:A:47:ARG:HG3	1:A:48:THR:HG23	2.01	0.42
1:B:676:LEU:HD23	1:B:676:LEU:HA	1.92	0.42
1:B:692:LEU:HD23	1:B:754:LEU:HD21	2.01	0.42
1:A:27:TYR:CE2	1:A:58:GLY:HA3	2.54	0.42
1:A:131:MET:HE2	1:A:134:LEU:HD11	2.01	0.42
1:B:699:TYR:H	1:B:711:ILE:HD12	1.85	0.42
1:A:443:ASN:ND2	2:A:1414:HOH:O	2.31	0.42
1:B:529:VAL:HG13	1:B:549:PRO:HB2	2.00	0.42
1:B:548:LYS:HA	1:B:549:PRO:HD3	1.91	0.42
1:B:950:THR:O	1:B:954:VAL:HG13	2.20	0.42
1:A:810:MET:HB2	1:A:810:MET:HE2	1.95	0.42
1:A:1172:VAL:HG21	1:A:1174:HIS:CE1	2.55	0.42
1:B:484:MET:HE3	1:B:497:LEU:HD23	2.02	0.42
1:B:795:ILE:HG22	1:B:889:PRO:HB2	2.01	0.42
1:B:39:SER:O	1:B:43:ARG:HG3	2.19	0.41
1:B:901:VAL:HG22	1:B:934:LEU:HD13	2.02	0.41
1:A:67:PRO:O	1:A:69:ASN:N	2.51	0.41
1:A:364:LYS:HA	1:A:364:LYS:HD2	1.88	0.41
1:B:973:GLY:CA	1:B:975:ASN:H	2.31	0.41
1:B:1226:PRO:HA	1:B:1229:TYR:CZ	2.55	0.41
1:A:979:THR:HB	1:A:1155:ALA:HA	2.03	0.41
1:B:64:LEU:HD23	1:B:64:LEU:HA	1.82	0.41
1:B:679:PRO:HB3	1:B:777:GLU:O	2.19	0.41
1:B:774:SER:OG	1:B:777:GLU:HG2	2.20	0.41
1:A:699:TYR:HD1	1:A:711:ILE:HD13	1.85	0.41
1:A:776:GLU:HB2	2:A:1797:HOH:O	2.21	0.41
1:A:1034:VAL:O	1:A:1093:ARG:HD2	2.20	0.41
1:B:882:VAL:HB	1:B:887:LEU:HD21	2.00	0.41
1:B:273:ILE:HD12	1:B:297:LEU:HD22	2.03	0.41
1:B:398:LEU:HD22	1:B:946:GLN:HG3	2.03	0.41
1:A:119:GLN:H	1:A:119:GLN:HG3	1.52	0.41
1:A:400:ILE:HD12	1:A:414:LEU:HD12	2.03	0.41
1:B:1298:LEU:HD12	1:B:1298:LEU:HA	1.90	0.41
1:A:243:GLN:NE2	2:A:1478:HOH:O	2.54	0.41
1:A:371:MET:CE	1:A:411:PRO:HB3	2.51	0.41
1:A:734:ASN:HB3	1:A:736:GLU:OE1	2.21	0.41
1:A:795:ILE:HG22	1:A:889:PRO:HB2	2.03	0.41
1:B:896:PHE:O	1:B:900:GLN:HG2	2.21	0.41
1:B:963:ARG:HD2	1:B:1208:VAL:HB	2.03	0.41
1:B:1019:ILE:HD13	1:B:1056:ALA:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:699:TYR:CD1	1:A:711:ILE:HD13	2.55	0.41
1:B:925:GLY:O	1:B:927:PRO:HD3	2.21	0.41
1:A:842:LEU:HG	1:A:891:PHE:HD1	1.85	0.40
1:B:980:ILE:HB	1:B:1003:MET:HE1	2.02	0.40
1:A:27:TYR:HD2	1:A:56:LEU:HD11	1.87	0.40
1:A:39:SER:O	1:A:43:ARG:HG3	2.21	0.40
1:A:314:PHE:C	1:A:317:PRO:HD2	2.42	0.40
1:A:208:LEU:HD12	1:A:208:LEU:HA	1.88	0.40
1:B:231:TYR:N	2:B:1453:HOH:O	2.54	0.40
1:B:578:PHE:HD2	1:B:580:TYR:HE1	1.70	0.40
1:B:772:PRO:HG3	1:B:904:ASP:HB2	2.02	0.40
1:A:444:LEU:HD12	1:A:444:LEU:HA	1.92	0.40
1:B:1030:PRO:HG3	1:B:1069:MET:HE1	2.02	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	1242/1351 (92%)	1216 (98%)	25 (2%)	1 (0%)	51	73
1	B	1220/1351 (90%)	1159 (95%)	54 (4%)	7 (1%)	25	43
All	All	2462/2702 (91%)	2375 (96%)	79 (3%)	8 (0%)	41	61

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	68	GLU
1	B	91	GLU
1	B	700	GLN
1	A	115	PRO
1	B	115	PRO

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Mol	Chain	Res	Type
1	B	351	PRO
1	B	353	ARG
1	B	66	ASP

### 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	
1	A	1050/1147 (92%)	991 (94%)	59 (6%)	21	40
1	B	988/1147 (86%)	938 (95%)	50 (5%)	24	45
All	All	2038/2294 (89%)	1929 (95%)	109 (5%)	22	43

All (109) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	15	GLU
1	A	17	LEU
1	A	119	GLN
1	A	131	MET
1	A	194	LEU
1	A	208	LEU
1	A	242	MET
1	A	243	GLN
1	A	252	LEU
1	A	444	LEU
1	A	468	VAL
1	A	474	LEU
1	A	490	ASN
1	A	497	LEU
1	A	504	LEU
1	A	536	THR
1	A	587	LEU
1	A	643	LEU
1	A	653	TYR

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Mol	Chain	Res	Type
1	A	676	LEU
1	A	695	LEU
1	A	700	GLN
1	A	711	ILE
1	A	728	VAL
1	A	733	ILE
1	A	737	GLU
1	A	762	LEU
1	A	776	GLU
1	A	782	LEU
1	A	788	LEU
1	A	798	LEU
1	A	800	THR
1	A	809	ASN
1	A	822	LEU
1	A	846	GLN
1	A	890	LEU
1	A	898	LEU
1	A	910	LEU
1	A	929	ARG
1	A	943	LEU
1	A	954	VAL
1	A	968	GLN
1	A	971	GLU
1	A	1003	MET
1	A	1004	VAL
1	A	1052	LEU
1	A	1069	MET
1	A	1145	ARG
1	A	1147	MET
1	A	1158	THR
1	A	1183	LEU
1	A	1223	LEU
1	A	1224	LEU
1	A	1233	LEU
1	A	1246	ARG
1	A	1271	THR
1	A	1291	ARG
1	A	1298	LEU
1	B	1	MET
1	B	38	LEU
1	B	66	ASP

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Mol	Chain	Res	Type
1	B	92	ASP
1	B	102	THR
1	B	158	GLN
1	B	177	VAL
1	B	219	LEU
1	B	220	ASP
1	B	269	LEU
1	B	457	THR
1	B	474	LEU
1	B	490	ASN
1	B	504	LEU
1	B	536	THR
1	B	647	ILE
1	B	653	TYR
1	B	662	THR
1	B	676	LEU
1	B	685	LEU
1	B	699	TYR
1	B	739	ASP
1	B	742	GLN
1	B	793	GLU
1	B	798	LEU
1	B	808	ARG
1	B	811	GLU
1	B	822	LEU
1	B	828	LEU
1	B	846	GLN
1	B	890	LEU
1	B	898	LEU
1	B	910	LEU
1	B	954	VAL
1	B	963	ARG
1	B	968	GLN
1	B	997	LEU
1	B	1004	VAL
1	B	1022	VAL
1	B	1029	ARG
1	B	1052	LEU
1	B	1147	MET
1	B	1169	LEU
1	B	1183	LEU
1	B	1187	LEU

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Mol	Chain	Res	Type
1	B	1224	LEU
1	B	1246	ARG
1	B	1281	ARG
1	B	1291	ARG
1	B	1292	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1174	HIS
1	B	1174	HIS

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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, 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	1258/1351 (93%)	-0.12	30 (2%) 59 62	21, 37, 76, 119	38 (3%)
1	B	1233/1351 (91%)	0.13	78 (6%) 20 21	24, 55, 95, 123	3 (0%)
All	All	2491/2702 (92%)	0.00	108 (4%) 35 38	21, 46, 88, 123	41 (1%)

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	547	GLY	5.7
1	A	325	LYS	5.4
1	B	468	VAL	5.4
1	A	468	VAL	4.8
1	B	469	GLY	4.7
1	B	701	THR	4.7
1	B	529	VAL	4.7
1	A	467	ASN	4.6
1	B	470	THR	4.5
1	A	326	SER	4.2
1	B	516	TYR	3.9
1	B	747	VAL	3.7
1	B	879	TYR	3.7
1	A	466	GLY	3.7
1	B	538	TYR	3.6
1	B	548	LYS	3.6
1	B	733	ILE	3.6
1	B	555	SER	3.6
1	B	870	TRP	3.6
1	B	467	ASN	3.6
1	B	556	ASP	3.6
1	A	540	VAL	3.6
1	B	550	PRO	3.6
1	B	532	TYR	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	133	GLN	3.5
1	B	848	ASN	3.4
1	B	559	ASN	3.4
1	B	741	GLY	3.3
1	A	705	SER	3.3
1	B	560	LEU	3.3
1	A	470	THR	3.2
1	B	805	SER	3.1
1	B	711	ILE	3.0
1	B	882	VAL	3.0
1	B	558	GLN	3.0
1	A	592	SER	3.0
1	B	887	LEU	3.0
1	B	623	SER	3.0
1	A	1167	ILE	3.0
1	B	1172	VAL	2.9
1	B	551	GLY	2.9
1	B	809	ASN	2.9
1	B	580	TYR	2.9
1	B	740	GLN	2.9
1	A	621	HIS	2.9
1	B	549	PRO	2.8
1	B	561	LEU	2.8
1	B	546	TRP	2.8
1	B	535	LEU	2.8
1	A	71	ALA	2.8
1	B	877	SER	2.8
1	B	530	GLU	2.7
1	B	1167	ILE	2.7
1	B	533	GLU	2.6
1	A	709	ILE	2.6
1	B	698	SER	2.6
1	A	585	MET	2.6
1	A	629	GLY	2.6
1	A	469	GLY	2.5
1	B	581	GLU	2.5
1	B	609	ILE	2.4
1	B	643	LEU	2.4
1	B	540	VAL	2.4
1	B	621	HIS	2.4
1	B	255	TYR	2.4
1	B	709	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	881[A]	ASN	2.4
1	B	352	ALA	2.4
1	A	30	LEU	2.4
1	B	823	ALA	2.3
1	B	590	SER	2.3
1	B	1168	SER	2.3
1	A	622	GLY	2.3
1	B	351	PRO	2.3
1	A	72	ASN	2.3
1	B	541	ARG	2.3
1	B	537	PRO	2.2
1	B	743	ARG	2.2
1	B	527	MET	2.2
1	A	653	TYR	2.2
1	B	557	GLY	2.2
1	B	531	GLN	2.2
1	B	1274	LYS	2.2
1	A	18	ASN	2.2
1	A	472	ALA	2.2
1	B	601	ALA	2.2
1	B	716	MET	2.1
1	B	802	ILE	2.1
1	B	699	TYR	2.1
1	A	220	ASP	2.1
1	A	19	GLY	2.1
1	B	471	ALA	2.1
1	B	515	GLN	2.1
1	B	746	ILE	2.1
1	A	414	LEU	2.1
1	B	465	LYS	2.1
1	B	218	GLY	2.1
1	A	134	LEU	2.1
1	A	219	LEU	2.1
1	B	89	PHE	2.1
1	B	796	TRP	2.1
1	B	738	MET	2.1
1	B	552	HIS	2.1
1	A	230	VAL	2.0
1	B	135	GLY	2.0
1	B	526	ARG	2.0
1	A	620	THR	2.0
1	B	136	GLN	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.