



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 01:56 am GMT

PDB ID : 4W8Y
Title : Structure of full length Cmr2 from *Pyrococcus furiosus* (Manganese bound form)
Authors : Benda, C.; Ebert, J.; Baumgaertner, M.; Conti, E.
Deposited on : 2014-08-26
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

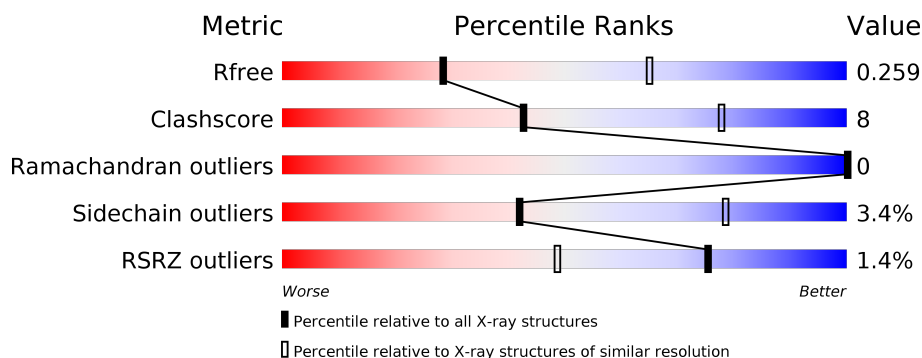
1 Overall quality at a glance

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}	100719	1692 (3.00-3.00)
Clashscore	112137	2037 (3.00-3.00)
Ramachandran outliers	110173	1973 (3.00-3.00)
Sidechain outliers	110143	1976 (3.00-3.00)
RSRZ outliers	101464	1716 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	888	<div> <div> <div></div> <div>75%</div> <div>18%</div> <div>• 6%</div> </div> <div>2%</div> </div>
1	B	888	<div> <div> <div></div> <div>73%</div> <div>20%</div> <div>• 7%</div> </div> <div>2%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ZN	B	901	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13055 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 CRISPR system Cmr subunit Cmr2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	835	Total	C	N	O	S	0	0	0
			6498	4211	1076	1198	13			
1	B	828	Total	C	N	O	S	0	0	0
			6517	4235	1069	1199	14			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	MET	-	initiating methionine	UNP Q8U1S6
A	-15	HIS	-	expression tag	UNP Q8U1S6
A	-14	HIS	-	expression tag	UNP Q8U1S6
A	-13	HIS	-	expression tag	UNP Q8U1S6
A	-12	HIS	-	expression tag	UNP Q8U1S6
A	-11	HIS	-	expression tag	UNP Q8U1S6
A	-10	HIS	-	expression tag	UNP Q8U1S6
A	-9	GLU	-	expression tag	UNP Q8U1S6
A	-8	ASN	-	expression tag	UNP Q8U1S6
A	-7	LEU	-	expression tag	UNP Q8U1S6
A	-6	TYR	-	expression tag	UNP Q8U1S6
A	-5	PHE	-	expression tag	UNP Q8U1S6
A	-4	GLN	-	expression tag	UNP Q8U1S6
A	-3	GLY	-	expression tag	UNP Q8U1S6
A	-2	ALA	-	expression tag	UNP Q8U1S6
A	-1	ALA	-	expression tag	UNP Q8U1S6
A	0	SER	-	expression tag	UNP Q8U1S6
B	-16	MET	-	initiating methionine	UNP Q8U1S6
B	-15	HIS	-	expression tag	UNP Q8U1S6
B	-14	HIS	-	expression tag	UNP Q8U1S6
B	-13	HIS	-	expression tag	UNP Q8U1S6
B	-12	HIS	-	expression tag	UNP Q8U1S6
B	-11	HIS	-	expression tag	UNP Q8U1S6
B	-10	HIS	-	expression tag	UNP Q8U1S6
B	-9	GLU	-	expression tag	UNP Q8U1S6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-8	ASN	-	expression tag	UNP Q8U1S6
B	-7	LEU	-	expression tag	UNP Q8U1S6
B	-6	TYR	-	expression tag	UNP Q8U1S6
B	-5	PHE	-	expression tag	UNP Q8U1S6
B	-4	GLN	-	expression tag	UNP Q8U1S6
B	-3	GLY	-	expression tag	UNP Q8U1S6
B	-2	ALA	-	expression tag	UNP Q8U1S6
B	-1	ALA	-	expression tag	UNP Q8U1S6
B	0	SER	-	expression tag	UNP Q8U1S6

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	4	Total	Mn	0	0
			4	4		
3	A	4	Total	Mn	0	0
			4	4		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	16	Total	O	0	0
			16	16		
4	B	14	Total	O	0	0
			14	14		

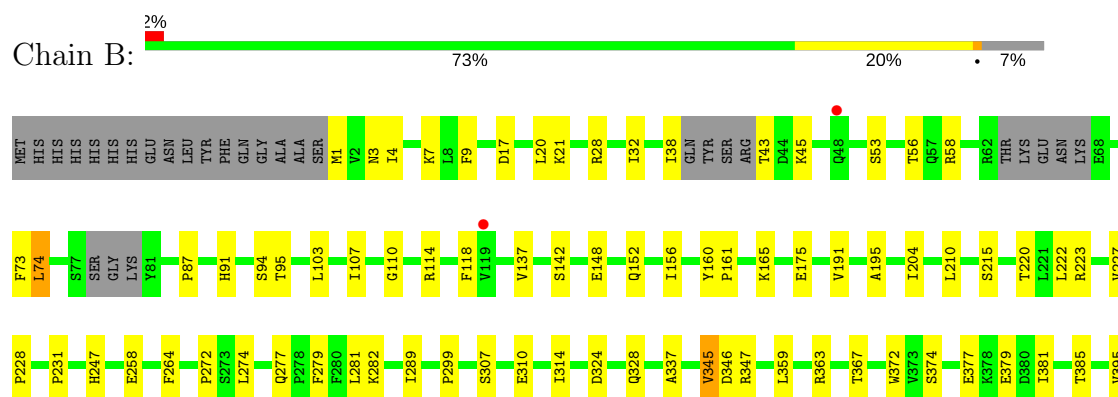
3 Residue-property plots

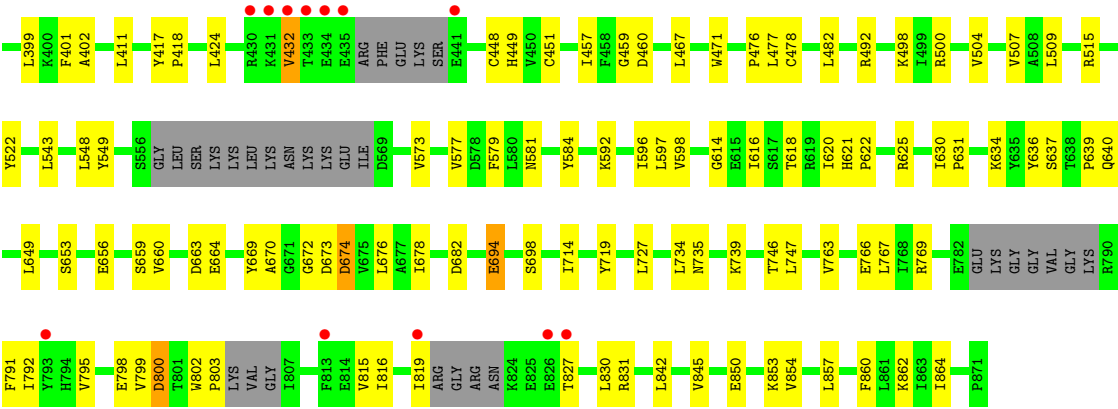
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: CRISPR system Cmr subunit Cmr2



• Molecule 1: CRISPR system Cmr subunit Cmr2





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.40Å 167.93Å 100.91Å 90.00° 98.91° 90.00°	Depositor
Resolution (Å)	85.72 – 3.00 61.65 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.4 (85.72-3.00) 98.5 (61.65-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1678)	Depositor
R, R_{free}	0.223 , 0.259 0.225 , 0.259	Depositor DCC
R_{free} test set	3949 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	90.0	Xtriage
Anisotropy	0.282	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 57.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13055	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% 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, MN

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	0/6634	0.42	0/9002
1	B	0.24	0/6656	0.41	0/9024
All	All	0.24	0/13290	0.41	0/18026

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

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	6498	0	6297	94	0
1	B	6517	0	6365	107	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	4	0	0	0	0
3	B	4	0	0	0	0
4	A	16	0	0	1	0
4	B	14	0	0	2	0
All	All	13055	0	12662	201	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 8.

All (201) 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:360:LYS:NZ	1:A:440:SER:OG	2.07	0.86
1:A:210:LEU:HD11	1:A:639:PRO:HB2	1.63	0.80
1:B:21:LYS:O	1:B:28:ARG:NH2	2.19	0.73
1:A:21:LYS:O	1:A:28:ARG:NH2	2.21	0.72
1:B:656:GLU:HB3	1:B:694:GLU:HG2	1.71	0.71
1:B:17:ASP:OD2	1:B:28:ARG:NH1	2.26	0.68
1:A:656:GLU:HB3	1:A:694:GLU:HG2	1.76	0.68
1:A:381:ILE:HD11	1:A:411:LEU:HD12	1.75	0.67
1:A:790:ARG:O	1:A:794:HIS:ND1	2.27	0.67
1:A:529:LYS:HB3	1:A:576:VAL:HG22	1.78	0.66
1:B:800:ASP:N	1:B:800:ASP:OD1	2.30	0.64
1:A:228:PRO:HB2	1:A:231:PRO:HD2	1.80	0.63
1:B:381:ILE:HD11	1:B:411:LEU:HD12	1.79	0.63
1:B:9:PHE:HB2	1:B:38:ILE:HG21	1.81	0.62
1:B:620:ILE:HD11	1:B:630:ILE:HD12	1.80	0.62
1:B:795:VAL:O	1:B:799:VAL:HG23	1.99	0.62
1:A:845:VAL:HG21	1:A:857:LEU:HB2	1.82	0.62
1:B:227:VAL:HG21	1:B:432:VAL:HG23	1.82	0.61
1:A:394:ASP:OD2	1:A:634:LYS:N	2.33	0.61
1:B:4:ILE:HD11	1:B:156:ILE:HD12	1.82	0.61
1:B:74:LEU:HD11	1:B:107:ILE:HD12	1.84	0.59
1:B:735:ASN:OD1	1:B:739:LYS:NZ	2.36	0.59
1:A:222:LEU:HD12	1:A:372:TRP:HB3	1.85	0.59
1:A:125:ARG:NH2	1:A:175:GLU:OE1	2.35	0.58
1:A:28:ARG:O	1:A:32:ILE:HG12	2.04	0.58
1:B:28:ARG:O	1:B:32:ILE:HG12	2.04	0.57
1:A:418:PRO:HG2	1:A:636:TYR:HB3	1.87	0.57
1:B:845:VAL:HG21	1:B:857:LEU:HB2	1.86	0.57
1:A:377:GLU:OE1	1:A:413:ARG:NH1	2.38	0.57
1:B:281:LEU:HB2	1:B:289:ILE:HD11	1.87	0.57
1:B:222:LEU:HD12	1:B:372:TRP:HB3	1.86	0.57
1:A:827:THR:OG1	1:A:831:ARG:NH1	2.38	0.56
1:A:617:SER:O	1:A:625:ARG:NH2	2.39	0.56
1:A:220:THR:HG22	1:A:374:SER:HA	1.87	0.56
1:A:735:ASN:OD1	1:A:739:LYS:NZ	2.39	0.55
1:B:228:PRO:HB2	1:B:231:PRO:HD2	1.88	0.55
1:A:2:VAL:HG13	1:A:7:LYS:HD2	1.88	0.55
1:B:195:ALA:HB2	1:B:204:ILE:HA	1.87	0.55
1:B:592:LYS:NZ	4:B:1010:HOH:O	2.35	0.55
1:B:674:ASP:N	1:B:674:ASP:OD1	2.36	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:816:ILE:HA	1:A:831:ARG:HE	1.72	0.55
1:B:395:VAL:O	1:B:399:LEU:HG	2.07	0.55
1:A:195:ALA:HB2	1:A:204:ILE:HA	1.89	0.54
1:B:381:ILE:HG22	1:B:399:LEU:HD13	1.90	0.54
1:A:714:ILE:HG13	1:A:747:LEU:HD11	1.90	0.54
1:A:860:PHE:CZ	1:A:864:ILE:HD11	2.42	0.54
1:A:448:CYS:HB2	1:A:478:CYS:HB3	1.90	0.54
1:B:507:VAL:HG11	1:B:678:ILE:HG22	1.90	0.54
1:B:418:PRO:HG3	1:B:637:SER:O	2.07	0.53
1:A:401:PHE:HB2	1:A:631:PRO:HB3	1.89	0.53
1:A:500:ARG:NH1	1:A:663:ASP:OD1	2.41	0.53
1:A:443:LEU:HB2	1:A:455:LEU:HD22	1.89	0.53
1:A:91:HIS:HB3	1:A:94:SER:HB2	1.90	0.53
1:B:94:SER:HG	1:B:698:SER:HG	1.57	0.53
1:B:448:CYS:HB3	1:B:451:CYS:SG	2.49	0.52
1:A:379:GLU:OE2	1:A:413:ARG:NH2	2.43	0.52
1:A:459:GLY:HA2	1:A:467:LEU:HD22	1.92	0.52
1:A:395:VAL:O	1:A:399:LEU:HG	2.09	0.52
1:B:220:THR:HG22	1:B:374:SER:HA	1.90	0.52
1:A:330:TYR:HB2	1:A:483:ILE:HD11	1.92	0.52
1:B:7:LYS:HE2	1:B:264:PHE:O	2.09	0.52
1:A:460:ASP:OD1	1:A:460:ASP:N	2.37	0.52
1:A:791:PHE:O	1:A:795:VAL:HG23	2.09	0.52
1:A:346:ASP:HB2	1:A:347:ARG:HE	1.75	0.51
1:B:142:SER:HA	1:B:152:GLN:HE21	1.76	0.51
1:A:649:LEU:HB3	1:A:672:GLY:HA2	1.93	0.51
1:B:210:LEU:HD12	1:B:640:GLN:HG3	1.92	0.51
1:B:792:ILE:HG21	1:B:862:LYS:HB2	1.91	0.51
1:A:74:LEU:HG	1:A:119:VAL:HG13	1.93	0.51
1:A:607:VAL:HG13	1:A:613:LEU:HD21	1.93	0.51
1:B:460:ASP:N	1:B:460:ASP:OD1	2.38	0.51
1:B:247:HIS:CD2	1:B:279:PHE:HZ	2.28	0.51
1:B:714:ILE:HG13	1:B:747:LEU:HD11	1.94	0.50
1:A:148:GLU:HG3	1:A:151:LYS:HG3	1.93	0.50
1:A:674:ASP:N	1:A:674:ASP:OD1	2.32	0.50
1:B:477:LEU:HD23	1:B:482:LEU:HA	1.94	0.50
1:B:94:SER:O	1:B:95:THR:OG1	2.27	0.50
1:A:94:SER:O	1:A:95:THR:OG1	2.25	0.49
1:B:500:ARG:NH1	1:B:663:ASP:OD1	2.45	0.49
1:A:227:VAL:HG21	1:A:432:VAL:HG13	1.95	0.49
1:B:114:ARG:HB2	1:B:118:PHE:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:PHE:CE2	1:B:87:PRO:HG2	2.48	0.49
1:A:100:TYR:HD1	1:A:189:GLU:HB3	1.78	0.48
1:B:103:LEU:O	1:B:107:ILE:HG12	2.13	0.48
1:B:815:VAL:O	1:B:819:ILE:HG22	2.13	0.48
1:B:682:ASP:OD1	1:B:682:ASP:N	2.47	0.48
1:A:247:HIS:CD2	1:A:279:PHE:HZ	2.31	0.48
1:A:73:PHE:CE2	1:A:87:PRO:HG2	2.48	0.48
1:B:860:PHE:CZ	1:B:864:ILE:HD11	2.48	0.48
1:A:746:THR:HG22	1:A:846:ARG:HE	1.78	0.48
1:A:385:THR:HG21	1:A:399:LEU:HD11	1.95	0.48
1:B:769:ARG:NH1	4:B:1004:HOH:O	2.46	0.48
1:A:545:ASP:OD1	4:A:1003:HOH:O	2.20	0.48
1:A:850:GLU:O	1:A:854:VAL:HG23	2.14	0.48
1:B:791:PHE:O	1:B:795:VAL:HG23	2.13	0.48
1:B:272:PRO:HG3	1:B:299:PRO:HD2	1.96	0.47
1:B:282:LYS:HB2	1:B:289:ILE:HD12	1.95	0.47
1:B:448:CYS:HB2	1:B:478:CYS:HB3	1.95	0.47
1:A:457:ILE:HB	1:A:471:TRP:HZ2	1.79	0.47
1:B:53:SER:O	1:B:56:THR:OG1	2.28	0.47
1:B:842:LEU:O	1:B:853:LYS:HD3	2.15	0.47
1:A:771:PHE:CE2	1:A:777:ARG:HG3	2.50	0.47
1:B:142:SER:HA	1:B:152:GLN:NE2	2.30	0.47
1:A:815:VAL:HG13	1:A:819:ILE:HD12	1.97	0.47
1:A:272:PRO:HG3	1:A:299:PRO:HD2	1.97	0.47
1:A:86:TYR:CD2	1:A:104:GLU:HG3	2.50	0.47
1:B:45:LYS:HB3	1:B:215:SER:HB2	1.97	0.47
1:B:596:ILE:HD13	1:B:727:LEU:HD23	1.97	0.47
1:A:160:TYR:HB3	1:A:161:PRO:HD3	1.96	0.46
1:A:84:VAL:HG12	1:A:623:ASN:HB3	1.96	0.46
1:B:160:TYR:HB3	1:B:161:PRO:HD3	1.98	0.46
1:B:543:LEU:HB3	1:B:548:LEU:HD12	1.97	0.46
1:A:842:LEU:O	1:A:853:LYS:HD3	2.14	0.46
1:B:614:GLY:HA2	1:B:634:LYS:HE2	1.98	0.46
1:A:377:GLU:O	1:A:413:ARG:HD2	2.16	0.46
1:B:165:LYS:HB2	1:B:191:VAL:HG21	1.97	0.46
1:A:507:VAL:HG11	1:A:678:ILE:HG22	1.97	0.46
1:B:417:TYR:CG	1:B:639:PRO:HG3	2.51	0.46
1:A:446:TRP:CZ2	1:A:464:HIS:HB2	2.51	0.46
1:A:245:SER:OG	1:A:484:LYS:HD2	2.16	0.46
1:A:673:ASP:OD1	1:A:673:ASP:N	2.48	0.45
1:B:850:GLU:O	1:B:854:VAL:HG23	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:816:ILE:HA	1:B:831:ARG:HE	1.81	0.45
1:A:511:TYR:HB3	1:A:591:PRO:HG3	1.99	0.45
1:A:792:ILE:O	1:A:796:LEU:HD22	2.17	0.45
1:A:841:LEU:O	1:A:845:VAL:HG23	2.17	0.45
1:B:449:HIS:CG	1:B:476:PRO:HD2	2.52	0.44
1:B:543:LEU:HD13	1:B:548:LEU:HB2	1.99	0.44
1:B:597:LEU:HD13	1:B:714:ILE:HG12	1.99	0.44
1:A:807:ILE:HA	1:A:810:MET:HB3	2.00	0.44
1:B:110:GLY:HA3	1:B:118:PHE:CE2	2.53	0.44
1:B:673:ASP:OD1	1:B:673:ASP:N	2.45	0.44
1:B:598:VAL:HG12	1:B:734:LEU:HD22	1.99	0.44
1:A:74:LEU:HD23	1:A:123:LEU:HD21	2.00	0.44
1:B:359:LEU:O	1:B:363:ARG:HB2	2.18	0.44
1:B:310:GLU:O	1:B:314:ILE:HG13	2.18	0.44
1:B:324:ASP:O	1:B:328:GLN:HG3	2.18	0.44
1:B:337:ALA:HB2	1:B:482:LEU:HD11	2.00	0.44
1:B:58:ARG:CZ	1:B:402:ALA:HB2	2.48	0.44
1:B:424:LEU:HA	1:B:424:LEU:HD12	1.85	0.44
1:B:549:TYR:O	1:B:584:TYR:OH	2.35	0.44
1:B:577:VAL:O	1:B:581:ASN:ND2	2.47	0.44
1:B:792:ILE:H	1:B:792:ILE:HD12	1.83	0.44
1:A:356:GLU:O	1:A:360:LYS:HB2	2.17	0.43
1:B:385:THR:HG21	1:B:399:LEU:HD11	2.00	0.43
1:B:649:LEU:HB3	1:B:672:GLY:HA2	1.99	0.43
1:A:839:LYS:HA	1:A:839:LYS:HE2	2.00	0.43
1:B:459:GLY:HA2	1:B:467:LEU:HD22	2.00	0.43
1:B:592:LYS:HD3	1:B:719:TYR:HE2	1.83	0.43
1:B:653:SER:HB2	1:B:670:ALA:O	2.18	0.43
1:A:624:ILE:O	1:A:628:VAL:HG12	2.18	0.43
1:B:227:VAL:HG21	1:B:432:VAL:CG2	2.48	0.43
1:B:457:ILE:HB	1:B:471:TRP:HZ2	1.83	0.43
1:A:4:ILE:HD11	1:A:156:ILE:HD12	1.99	0.43
1:B:274:LEU:HA	1:B:277:GLN:HG3	2.01	0.43
1:B:91:HIS:HB3	1:B:94:SER:HB2	2.00	0.43
1:A:324:ASP:O	1:A:328:GLN:HG3	2.18	0.43
1:B:622:PRO:HA	1:B:625:ARG:HG3	2.01	0.43
1:A:23:GLU:O	1:A:23:GLU:HG2	2.18	0.43
1:A:310:GLU:O	1:A:314:ILE:HG13	2.19	0.43
1:B:3:ASN:OD1	1:B:3:ASN:N	2.51	0.43
1:B:763:VAL:HG22	1:B:767:LEU:HD13	2.00	0.43
1:A:416:ILE:HG22	1:A:419:LEU:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:656:GLU:O	1:B:659:SER:OG	2.23	0.42
1:A:274:LEU:HA	1:A:277:GLN:HG3	2.01	0.42
1:A:291:ASP:OD2	1:A:291:ASP:N	2.52	0.42
1:A:51:ALA:O	1:A:55:LYS:HG2	2.20	0.42
1:B:418:PRO:HG2	1:B:636:TYR:HB3	2.01	0.42
1:B:492:ARG:NH1	1:B:498:LYS:HD3	2.35	0.42
1:A:32:ILE:HD12	1:A:130:GLU:HB2	2.00	0.42
1:A:420:LEU:HA	1:A:420:LEU:HD12	1.93	0.42
1:B:621:HIS:HA	1:B:622:PRO:HD3	1.92	0.42
1:A:270:ILE:HD11	1:A:305:ILE:HD11	2.00	0.42
1:B:766:GLU:H	1:B:766:GLU:CD	2.23	0.42
1:A:145:PHE:O	1:A:152:GLN:NE2	2.53	0.42
1:B:802:TRP:HA	1:B:803:PRO:HD3	1.90	0.42
1:A:827:THR:O	1:A:831:ARG:HG3	2.20	0.42
1:B:345:VAL:HG13	1:B:346:ASP:O	2.19	0.42
1:B:798:GLU:O	1:B:802:TRP:N	2.52	0.42
1:A:145:PHE:HB2	1:A:152:GLN:HG3	2.01	0.41
1:B:401:PHE:HB2	1:B:631:PRO:HB3	2.02	0.41
1:B:819:ILE:HD12	1:B:819:ILE:HA	1.94	0.41
1:A:558:LEU:HD12	1:A:558:LEU:HA	1.90	0.41
1:B:509:LEU:HB3	1:B:515:ARG:HB2	2.02	0.41
1:A:573:VAL:O	1:A:577:VAL:HG23	2.20	0.41
1:B:616:ILE:HG12	1:B:630:ILE:HD13	2.02	0.41
1:B:573:VAL:O	1:B:577:VAL:HG23	2.21	0.41
1:A:763:VAL:HG11	1:A:767:LEU:HD22	2.02	0.41
1:B:669:TYR:HB3	1:B:676:LEU:HB3	2.03	0.41
1:B:20:LEU:HD23	1:B:73:PHE:CD2	2.56	0.41
1:A:369:SER:HB3	1:A:431:LYS:HD3	2.02	0.41
1:B:137:VAL:HG12	1:B:156:ILE:HG12	2.03	0.41
1:A:603:ASP:O	1:A:607:VAL:HG23	2.21	0.41
1:B:9:PHE:HE2	1:B:43:THR:HG21	1.85	0.41
1:B:227:VAL:HG23	1:B:367:THR:HB	2.02	0.41
1:B:522:TYR:HB3	1:B:579:PHE:HZ	1.86	0.41
1:B:656:GLU:O	1:B:660:VAL:HG23	2.21	0.40
1:A:598:VAL:HG12	1:A:734:LEU:HD22	2.03	0.40
1:A:603:ASP:OD1	1:A:744:LYS:NZ	2.48	0.40
1:A:523:GLY:O	1:A:527:VAL:HG23	2.21	0.40
1:A:223:ARG:NH1	1:A:225:LYS:HB2	2.36	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	
1	A	819/888 (92%)	797 (97%)	22 (3%)	0	100	100
1	B	810/888 (91%)	789 (97%)	21 (3%)	0	100	100
All	All	1629/1776 (92%)	1586 (97%)	43 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	666/795 (84%)	641 (96%)	25 (4%)	38	75
1	B	677/795 (85%)	656 (97%)	21 (3%)	45	80
All	All	1343/1590 (84%)	1297 (97%)	46 (3%)	42	78

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	LEU
1	A	148	GLU
1	A	210	LEU
1	A	223	ARG
1	A	258	GLU
1	A	345	VAL
1	A	347	ARG
1	A	394	ASP

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Mol	Chain	Res	Type
1	A	413	ARG
1	A	416	ILE
1	A	443	LEU
1	A	504	VAL
1	A	532	GLU
1	A	541	ASN
1	A	575	GLU
1	A	624	ILE
1	A	626	ASP
1	A	634	LYS
1	A	674	ASP
1	A	746	THR
1	A	796	LEU
1	A	800	ASP
1	A	801	THR
1	A	810	MET
1	A	825	GLU
1	B	1	MET
1	B	74	LEU
1	B	148	GLU
1	B	175	GLU
1	B	223	ARG
1	B	258	GLU
1	B	307	SER
1	B	345	VAL
1	B	347	ARG
1	B	377	GLU
1	B	379	GLU
1	B	432	VAL
1	B	504	VAL
1	B	618	THR
1	B	664	GLU
1	B	674	ASP
1	B	694	GLU
1	B	746	THR
1	B	800	ASP
1	B	827	THR
1	B	830	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	152	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 10 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	835/888 (94%)	-0.14	9 (1%)	80 55	53, 89, 134, 213	0
1	B	828/888 (93%)	-0.11	14 (1%)	70 42	58, 92, 129, 192	0
All	All	1663/1776 (93%)	-0.12	23 (1%)	75 49	53, 91, 131, 213	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	435	GLU	13.2
1	B	432	VAL	6.6
1	B	434	GLU	6.6
1	B	431	LYS	5.7
1	A	627	TYR	5.1
1	A	825	GLU	4.7
1	B	433	THR	4.2
1	B	48	GLN	3.7
1	A	532	GLU	3.5
1	B	119	VAL	3.0
1	B	819	ILE	2.6
1	A	62	ARG	2.5
1	B	793	TYR	2.4
1	A	826	GLU	2.4
1	B	813	PHE	2.3
1	A	23	GLU	2.3
1	A	828	LYS	2.2
1	B	430	ARG	2.2
1	B	441	GLU	2.1
1	A	22	ILE	2.1
1	B	827	THR	2.1
1	A	116	GLU	2.0
1	B	826	GLU	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. 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, 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	LLDF	B-factors(Å ²)	Q<0.9
2	ZN	B	901	1/1	0.96	0.19	2.84	85,85,85,85	0
3	MN	B	902	1/1	0.91	0.17	-0.63	143,143,143,143	0
2	ZN	A	901	1/1	0.99	0.15	-0.99	64,64,64,64	0
3	MN	A	905	1/1	0.58	0.13	-	132,132,132,132	0
3	MN	A	904	1/1	0.96	0.19	-	76,76,76,76	0
3	MN	B	905	1/1	0.81	0.23	-	191,191,191,191	0
3	MN	A	902	1/1	0.81	0.23	-	172,172,172,172	0
3	MN	A	903	1/1	0.93	0.18	-	153,153,153,153	0
3	MN	B	903	1/1	0.24	0.14	-	168,168,168,168	0
3	MN	B	904	1/1	0.98	0.23	-	82,82,82,82	0

6.5 Other polymers [i](#)

There are no such residues in this entry.