



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:32 PM GMT

PDB ID : 4FCX
Title : S.pombe Mre11 apoenzym
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Deposited on : 2012-05-25
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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

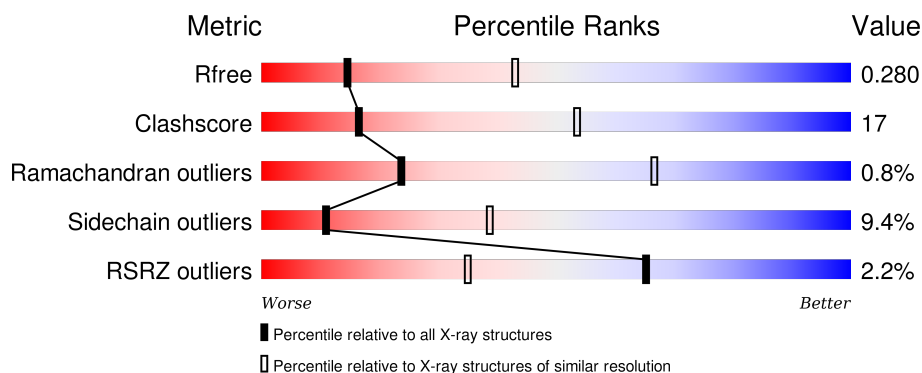
1 Overall quality at a glance

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}	91344	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (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	404	
1	B	404	

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	MN	A	502	-	-	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5900 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 repair protein rad32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	369	Total	C	N	O	S	115	0	0
			2973	1899	515	552	7			
1	A	366	Total	C	N	O	S	38	0	0
			2923	1870	504	542	7			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	10	GLY	-	EXPRESSION TAG	UNP Q09683
B	11	PRO	-	EXPRESSION TAG	UNP Q09683
B	12	LEU	-	EXPRESSION TAG	UNP Q09683
B	13	GLY	-	EXPRESSION TAG	UNP Q09683
B	14	SER	-	EXPRESSION TAG	UNP Q09683
A	10	GLY	-	EXPRESSION TAG	UNP Q09683
A	11	PRO	-	EXPRESSION TAG	UNP Q09683
A	12	LEU	-	EXPRESSION TAG	UNP Q09683
A	13	GLY	-	EXPRESSION TAG	UNP Q09683
A	14	SER	-	EXPRESSION TAG	UNP Q09683

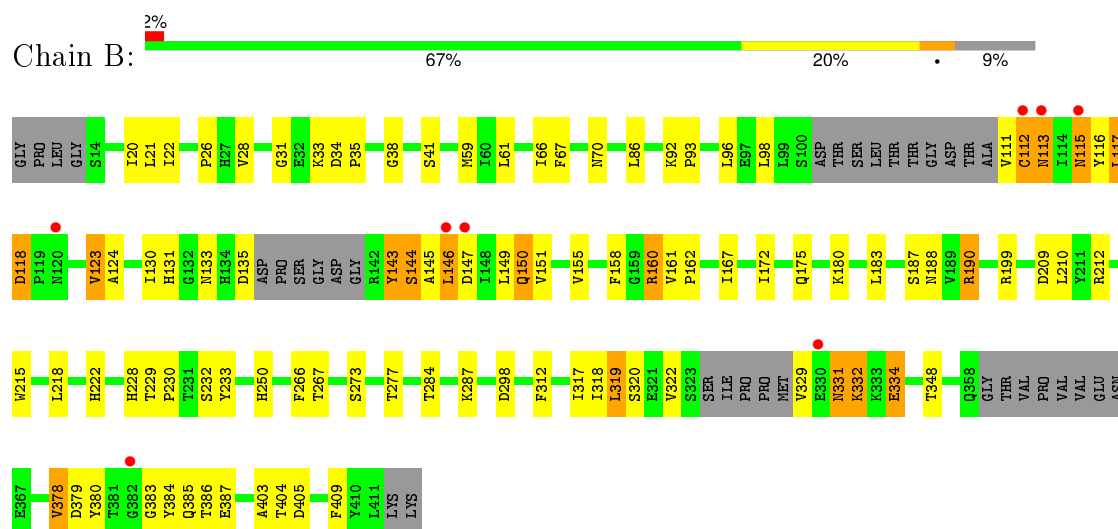
- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Mn	0	0
			2	2		
2	A	2	Total	Mn	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 errors 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 repair protein rad32



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.41 Å 82.26 Å 164.30 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.84 – 3.00 21.84 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.1 (21.84-3.00) 98.2 (21.84-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.99 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.227 , 0.286 0.222 , 0.280	Depositor DCC
R_{free} test set	1978 reflections (9.45%)	DCC
Wilson B-factor (Å ²)	76.6	Xtriage
Anisotropy	0.424	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 66.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 20950 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5900	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 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	0.41	0/2990	0.62	1/4057 (0.0%)
1	B	0.40	0/3039	0.57	0/4120
All	All	0.41	0/6029	0.60	1/8177 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	251	GLU	N-CA-C	-5.27	96.78	111.00

There are no chirality outliers.

There are no planarity outliers.

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	2923	0	2882	81	0
1	B	2973	0	2932	111	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
All	All	5900	0	5814	188	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 17.

All (188) 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:131:HIS:CE1	1:B:146:LEU:HD13	1.46	1.50
1:B:143:TYR:HA	1:B:144:SER:CB	1.51	1.39
1:B:319:LEU:HD13	1:B:384:TYR:CG	1.59	1.36
1:B:319:LEU:CD1	1:B:384:TYR:CG	2.18	1.24
1:B:143:TYR:CA	1:B:144:SER:HB3	1.75	1.17
1:B:319:LEU:HD13	1:B:384:TYR:CD2	1.79	1.16
1:B:319:LEU:CD1	1:B:384:TYR:CD2	2.27	1.16
1:A:167:ILE:HG23	1:A:203:VAL:HG22	1.18	1.12
1:B:319:LEU:HD11	1:B:384:TYR:HB2	1.27	1.09
1:B:111:VAL:HA	1:B:112:CYS:SG	1.93	1.08
1:B:143:TYR:HA	1:B:144:SER:HB3	1.09	1.08
1:B:131:HIS:CE1	1:B:146:LEU:CD1	2.35	1.08
1:A:393:SER:HB3	1:A:403:ALA:O	1.56	1.05
1:B:319:LEU:HD11	1:B:384:TYR:CB	1.86	1.04
1:A:167:ILE:CG2	1:A:203:VAL:HG22	1.86	1.04
1:B:131:HIS:HE1	1:B:146:LEU:CD1	1.69	1.03
1:B:319:LEU:HG	1:B:380:TYR:HA	1.03	1.03
1:B:319:LEU:CD1	1:B:384:TYR:HB2	1.87	1.03
1:B:319:LEU:CD1	1:B:384:TYR:CB	2.36	1.03
1:B:143:TYR:CA	1:B:144:SER:CB	2.33	1.02
1:A:166:ASN:OD1	1:A:202:LYS:HB3	1.60	1.02
1:A:167:ILE:HG12	1:A:203:VAL:HG13	1.38	1.02
1:B:143:TYR:HA	1:B:144:SER:HB2	1.38	1.01
1:A:145:ALA:HA	1:A:148:ILE:HD12	1.42	1.00
1:B:111:VAL:N	1:B:112:CYS:HG	1.60	0.99
1:B:319:LEU:CG	1:B:380:TYR:HA	1.91	0.99
1:A:375:ARG:HH12	1:A:405:ASP:CG	1.69	0.96
1:B:111:VAL:HB	1:B:113:ASN:HB2	1.47	0.96
1:B:331:ASN:O	1:B:385:GLN:HB3	1.65	0.94
1:B:319:LEU:HG	1:B:380:TYR:CA	1.95	0.93
1:B:319:LEU:HD12	1:B:384:TYR:CD2	2.02	0.93
1:A:375:ARG:NH1	1:A:405:ASP:OD2	2.05	0.89
1:A:393:SER:CB	1:A:403:ALA:O	2.21	0.88
1:B:111:VAL:HG11	1:B:113:ASN:HD22	1.39	0.88
1:B:111:VAL:CA	1:B:112:CYS:SG	2.65	0.84
1:B:93:PRO:HA	1:B:115:ASN:OD1	1.78	0.84
1:A:14:SER:H	1:A:15:GLU:HG2	1.42	0.83
1:A:93:PRO:HB2	1:A:94:CYS:HA	1.59	0.82
1:A:393:SER:HA	1:A:406:VAL:CG2	2.11	0.81
1:B:98:LEU:HD13	1:B:112:CYS:SG	2.21	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:VAL:HG12	1:B:329:VAL:O	1.82	0.79
1:A:55:ARG:NH2	1:A:302:GLU:OE2	2.19	0.75
1:B:111:VAL:HA	1:B:112:CYS:C	2.05	0.75
1:A:344:GLU:OE2	1:A:399:ARG:NH1	2.20	0.75
1:A:393:SER:HA	1:A:406:VAL:HG21	1.69	0.74
1:B:329:VAL:CG2	1:B:383:GLY:O	2.37	0.73
1:B:112:CYS:O	1:B:112:CYS:SG	2.47	0.72
1:B:143:TYR:CB	1:B:144:SER:HB3	2.18	0.72
1:B:229:THR:OG1	1:B:232:SER:HB2	1.90	0.71
1:B:319:LEU:HG	1:B:379:ASP:O	1.91	0.71
1:B:319:LEU:HB2	1:B:379:ASP:O	1.90	0.70
1:B:111:VAL:N	1:B:112:CYS:SG	2.63	0.70
1:B:111:VAL:CG1	1:B:113:ASN:HD22	2.05	0.69
1:B:115:ASN:ND2	1:B:115:ASN:N	2.40	0.69
1:B:111:VAL:HG11	1:B:113:ASN:ND2	2.08	0.68
1:A:393:SER:OG	1:A:406:VAL:N	2.26	0.68
1:A:145:ALA:HA	1:A:148:ILE:CD1	2.22	0.68
1:B:143:TYR:N	1:B:143:TYR:CD2	2.60	0.67
1:A:323:SER:HA	1:A:325:ILE:H	1.58	0.67
1:B:319:LEU:HD13	1:B:384:TYR:CD1	2.25	0.66
1:A:167:ILE:HD11	1:A:186:ILE:HG12	1.78	0.66
1:B:93:PRO:CA	1:B:115:ASN:OD1	2.42	0.66
1:B:319:LEU:CB	1:B:379:ASP:O	2.46	0.64
1:B:319:LEU:CG	1:B:379:ASP:O	2.47	0.63
1:A:165:ASP:O	1:A:197:SER:OG	2.13	0.62
1:A:167:ILE:HG23	1:A:203:VAL:CG2	2.12	0.61
1:B:131:HIS:HE1	1:B:146:LEU:HD13	0.88	0.61
1:A:393:SER:HB3	1:A:406:VAL:HG23	1.82	0.60
1:B:123:VAL:HG12	1:A:93:PRO:HB3	1.84	0.60
1:B:329:VAL:O	1:B:329:VAL:CG1	2.48	0.60
1:B:143:TYR:HB3	1:B:147:ASP:OD1	2.02	0.60
1:B:131:HIS:CE1	1:B:146:LEU:HD22	2.37	0.59
1:B:318:ILE:O	1:B:322:VAL:HG23	2.02	0.59
1:B:329:VAL:HG23	1:B:383:GLY:O	2.03	0.58
1:A:407:VAL:HG22	1:A:408:GLN:N	2.18	0.58
1:B:115:ASN:N	1:B:115:ASN:HD22	2.02	0.58
1:B:147:ASP:HB3	1:A:78:TYR:CE1	2.39	0.58
1:A:250:HIS:CG	1:A:250:HIS:O	2.58	0.56
1:A:256:ILE:HD13	1:A:303:LYS:HB2	1.87	0.56
1:A:145:ALA:CA	1:A:148:ILE:HD12	2.26	0.56
1:A:19:ARG:NH1	1:A:55:ARG:O	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:VAL:O	1:B:111:VAL:HG23	2.06	0.55
1:B:118:ASP:OD2	1:B:118:ASP:N	2.39	0.55
1:B:334:GLU:CD	1:B:334:GLU:H	2.06	0.55
1:A:407:VAL:CG2	1:A:408:GLN:N	2.68	0.55
1:A:225:HIS:NE2	1:A:236:GLU:OE2	2.36	0.55
1:B:405:ASP:OD1	1:B:405:ASP:N	2.39	0.55
1:A:167:ILE:CD1	1:A:186:ILE:HG12	2.37	0.54
1:B:144:SER:HG	1:B:147:ASP:CG	2.11	0.54
1:B:124:ALA:H	1:A:93:PRO:HG2	1.71	0.54
1:A:145:ALA:O	1:A:148:ILE:HB	2.08	0.54
1:A:393:SER:CB	1:A:406:VAL:HG23	2.38	0.53
1:B:329:VAL:HG21	1:B:383:GLY:O	2.07	0.53
1:A:165:ASP:OD1	1:A:166:ASN:ND2	2.41	0.53
1:A:392:PHE:CE2	1:A:406:VAL:HG12	2.44	0.53
1:B:131:HIS:HE1	1:B:146:LEU:CG	2.21	0.52
1:B:22:ILE:HG22	1:B:61:LEU:HB3	1.92	0.52
1:B:143:TYR:HB3	1:B:144:SER:HB3	1.89	0.52
1:A:91:ASP:OD1	1:A:122:ASN:ND2	2.40	0.52
1:B:31:GLY:O	1:B:38:GLY:HA2	2.09	0.52
1:B:386:THR:OG1	1:B:387:GLU:N	2.41	0.52
1:A:77:LEU:HD13	1:A:145:ALA:HB1	1.90	0.51
1:B:96:LEU:HD12	1:A:58:ASP:HB3	1.93	0.51
1:B:111:VAL:CG1	1:B:113:ASN:ND2	2.71	0.51
1:B:172:ILE:HB	1:B:183:LEU:HB3	1.91	0.51
1:B:319:LEU:HD12	1:B:384:TYR:HD2	1.69	0.51
1:A:392:PHE:CE2	1:A:406:VAL:CG1	2.94	0.50
1:B:67:PHE:CD2	1:B:145:ALA:HB1	2.46	0.50
1:A:338:TYR:O	1:A:342:LYS:HG2	2.10	0.50
1:A:392:PHE:HE2	1:A:406:VAL:HG12	1.75	0.50
1:B:26:PRO:HG2	1:B:66:ILE:HA	1.92	0.50
1:A:393:SER:CA	1:A:406:VAL:CG2	2.86	0.50
1:A:14:SER:HB2	1:A:15:GLU:HA	1.94	0.49
1:A:343:VAL:O	1:A:347:ILE:HG13	2.12	0.49
1:B:378:VAL:HG22	1:B:409:PHE:HD1	1.77	0.49
1:A:167:ILE:O	1:A:203:VAL:HA	2.12	0.49
1:B:149:LEU:HD13	1:B:155:VAL:HG21	1.94	0.49
1:B:284:THR:HG22	1:B:312:PHE:CE2	2.48	0.49
1:B:180:LYS:HB3	1:B:215:TRP:CE3	2.48	0.49
1:A:319:LEU:HD12	1:A:380:TYR:HA	1.94	0.49
1:A:287:LYS:HG3	1:A:309:VAL:HG11	1.94	0.48
1:B:331:ASN:O	1:B:385:GLN:CB	2.50	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:PRO:CB	1:A:94:CYS:HA	2.36	0.48
1:B:111:VAL:CA	1:B:112:CYS:CB	2.91	0.48
1:B:111:VAL:HA	1:B:112:CYS:CB	2.33	0.48
1:A:26:PRO:HG2	1:A:66:ILE:HA	1.96	0.48
1:B:230:PRO:HA	1:B:233:TYR:HE1	1.78	0.48
1:A:379:ASP:OD1	1:A:381:THR:OG1	2.30	0.48
1:A:195:TYR:CD2	1:A:231:THR:HG22	2.48	0.48
1:B:116:TYR:HB2	1:B:118:ASP:OD2	2.14	0.48
1:A:167:ILE:HD11	1:A:169:VAL:CG2	2.44	0.47
1:B:331:ASN:HB3	1:B:334:GLU:CD	2.34	0.47
1:B:143:TYR:HB3	1:B:147:ASP:CB	2.44	0.47
1:B:130:ILE:HG22	1:B:158:PHE:O	2.14	0.47
1:B:111:VAL:CA	1:B:112:CYS:C	2.80	0.47
1:B:135:ASP:HB2	1:B:146:LEU:HD21	1.96	0.47
1:B:146:LEU:O	1:B:150:GLN:HB3	2.15	0.46
1:A:66:ILE:HD11	1:A:129:SER:OG	2.16	0.46
1:A:393:SER:HA	1:A:406:VAL:CB	2.45	0.46
1:B:131:HIS:CE1	1:B:146:LEU:CD2	2.98	0.46
1:A:373:LEU:HA	1:A:402:ASN:OD1	2.16	0.46
1:A:148:ILE:O	1:A:152:THR:HG23	2.16	0.45
1:A:333:LYS:O	1:A:337:THR:HG23	2.16	0.45
1:A:375:ARG:NH1	1:A:405:ASP:CG	2.47	0.45
1:A:20:ILE:HG23	1:A:59:MET:HG3	1.99	0.45
1:B:98:LEU:CD1	1:B:112:CYS:SG	3.01	0.45
1:A:31:GLY:O	1:A:38:GLY:HA2	2.16	0.45
1:B:144:SER:OG	1:B:147:ASP:OD2	2.32	0.45
1:A:393:SER:OG	1:A:406:VAL:HB	2.17	0.45
1:B:319:LEU:HD22	1:B:319:LEU:HA	1.85	0.45
1:B:188:ASN:OD1	1:B:190:ARG:NH2	2.50	0.44
1:A:14:SER:CB	1:A:15:GLU:HA	2.46	0.44
1:B:20:ILE:HG23	1:B:59:MET:HG3	2.00	0.44
1:B:331:ASN:HB3	1:B:334:GLU:CG	2.48	0.44
1:B:222:HIS:CE1	1:B:250:HIS:H	2.36	0.44
1:A:354:TRP:O	1:A:358:GLN:HG2	2.18	0.44
1:A:222:HIS:ND1	1:A:250:HIS:HB2	2.32	0.44
1:A:392:PHE:HE2	1:A:406:VAL:CG1	2.31	0.44
1:B:131:HIS:ND1	1:B:146:LEU:HD22	2.32	0.43
1:A:314:MET:HE2	1:A:375:ARG:HD3	2.00	0.43
1:A:326:PRO:HA	1:A:327:PRO:HD3	1.87	0.43
1:A:98:LEU:HD12	1:A:100:SER:H	1.83	0.43
1:B:319:LEU:CD1	1:B:380:TYR:HA	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:ASN:ND2	1:A:405:ASP:O	2.40	0.43
1:B:334:GLU:OE1	1:B:334:GLU:N	2.30	0.43
1:A:393:SER:HA	1:A:406:VAL:HB	2.00	0.42
1:A:167:ILE:HD11	1:A:169:VAL:HG22	2.00	0.42
1:A:323:SER:HA	1:A:325:ILE:N	2.31	0.42
1:B:92:LYS:HA	1:B:93:PRO:HD3	1.61	0.42
1:B:116:TYR:HB3	1:B:117:LEU:H	1.63	0.42
1:B:273:SER:O	1:B:287:LYS:HD2	2.19	0.42
1:B:332:LYS:HE2	1:B:332:LYS:HB3	1.28	0.42
1:B:34:ASP:HA	1:B:35:PRO:HD3	1.90	0.41
1:A:204:LYS:HA	1:A:204:LYS:HD2	1.86	0.41
1:A:190:ARG:NH1	1:A:190:ARG:H	2.18	0.41
1:B:160:ARG:HA	1:B:187:SER:OG	2.21	0.41
1:B:317:ILE:O	1:B:378:VAL:HA	2.21	0.41
1:A:48:ILE:HA	1:A:48:ILE:HD13	1.80	0.41
1:A:339:LEU:O	1:A:343:VAL:HG23	2.21	0.41
1:A:269:VAL:O	1:A:271:PRO:HD3	2.21	0.41
1:B:28:VAL:HA	1:B:41:SER:OG	2.21	0.41
1:B:161:VAL:HG11	1:B:167:ILE:HG12	2.03	0.40
1:A:52:ALA:O	1:A:57:VAL:HG22	2.21	0.40
1:B:143:TYR:H	1:B:143:TYR:HD2	1.67	0.40
1:B:115:ASN:ND2	1:B:115:ASN:H	2.14	0.40
1:B:161:VAL:CG1	1:B:167:ILE:HG12	2.51	0.40
1:B:380:TYR:CD1	1:B:409:PHE:HB3	2.57	0.40
1:A:260:TYR:OH	1:A:265:LYS:HG2	2.21	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	358/404 (89%)	334 (93%)	21 (6%)	3 (1%)	24 66

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	359/404 (89%)	338 (94%)	18 (5%)	3 (1%)	24	66
All	All	717/808 (89%)	672 (94%)	39 (5%)	6 (1%)	24	66

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	144	SER
1	A	93	PRO
1	B	162	PRO
1	B	403	ALA
1	A	330	GLU
1	A	118	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	324/362 (90%)	298 (92%)	26 (8%)	15	47
1	B	334/362 (92%)	298 (89%)	36 (11%)	8	30
All	All	658/724 (91%)	596 (91%)	62 (9%)	11	39

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	21	LEU
1	B	33	LYS
1	B	70	ASN
1	B	86	LEU
1	B	112	CYS
1	B	113	ASN
1	B	115	ASN
1	B	117	LEU
1	B	118	ASP
1	B	123	VAL

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Mol	Chain	Res	Type
1	B	133	ASN
1	B	143	TYR
1	B	146	LEU
1	B	150	GLN
1	B	151	VAL
1	B	160	ARG
1	B	175	GLN
1	B	190	ARG
1	B	199	ARG
1	B	209	ASP
1	B	210	LEU
1	B	212	ARG
1	B	218	LEU
1	B	228	HIS
1	B	266	PHE
1	B	267	THR
1	B	277	THR
1	B	298	ASP
1	B	319	LEU
1	B	320	SER
1	B	331	ASN
1	B	332	LYS
1	B	334	GLU
1	B	348	THR
1	B	378	VAL
1	B	404	THR
1	A	14	SER
1	A	21	LEU
1	A	37	ARG
1	A	41	SER
1	A	85	ARG
1	A	94	CYS
1	A	98	LEU
1	A	146	LEU
1	A	149	LEU
1	A	164	ASN
1	A	165	ASP
1	A	167	ILE
1	A	190	ARG
1	A	211	TYR
1	A	229	THR
1	A	231	THR

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Mol	Chain	Res	Type
1	A	232	SER
1	A	266	PHE
1	A	275	ILE
1	A	277	THR
1	A	298	ASP
1	A	300	HIS
1	A	320	SER
1	A	371	LEU
1	A	381	THR
1	A	404	THR

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

Mol	Chain	Res	Type
1	B	113	ASN
1	B	131	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	366/404 (90%)	-0.34	8 (2%) 65 35	30, 72, 101, 125	7 (1%)
1	B	368/404 (91%)	-0.19	8 (2%) 65 35	26, 77, 131, 153	23 (6%)
All	All	734/808 (90%)	-0.26	16 (2%) 65 35	26, 75, 119, 153	30 (4%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	164	ASN	5.2
1	B	115	ASN	4.9
1	B	113	ASN	4.0
1	B	120	ASN	3.5
1	B	146	LEU	3.0
1	B	382	GLY	2.6
1	B	330	GLU	2.6
1	A	228	HIS	2.4
1	A	122	ASN	2.3
1	A	166	ASN	2.3
1	A	91	ASP	2.3
1	A	165	ASP	2.2
1	A	404	THR	2.2
1	B	112	CYS	2.1
1	A	330	GLU	2.0
1	B	147	ASP	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(\AA^2)	Q<0.9
2	MN	A	502	1/1	0.94	0.21	2.13	64,64,64,64	0
2	MN	A	501	1/1	0.98	0.19	0.48	58,58,58,58	0
2	MN	B	501	1/1	0.99	0.18	0.29	49,49,49,49	0
2	MN	B	502	1/1	0.99	0.14	-1.15	49,49,49,49	0

6.5 Other polymers [i](#)

There are no such residues in this entry.