



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 08:35 pm GMT

PDB ID : 2ZXP
Title : Crystal structure of RecJ in complex with Mn²⁺ from *Thermus thermophilus* HB8
Authors : Wakamatsu, T.; Kitamura, Y.; Nakagawa, N.; Masui, R.; Kuramitsu, S.
Deposited on : 2009-01-05
Resolution : 2.30 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

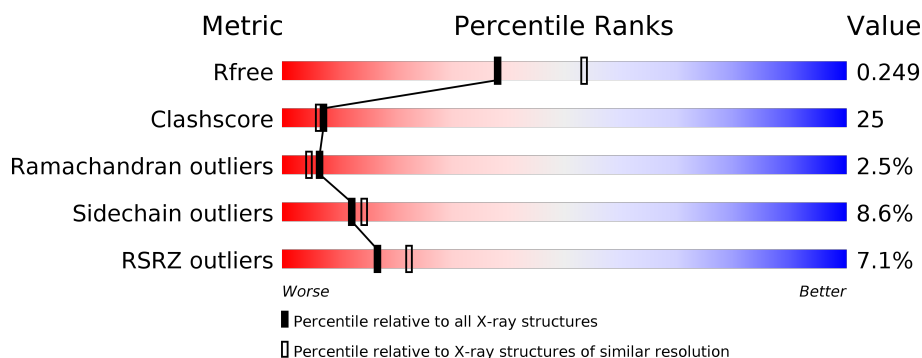
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	666	<div> <div>7%</div> <div>62%</div> <div>30%</div> <div>5% • •</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5315 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 Single-stranded DNA specific exonuclease RecJ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	648	Total	C	N	O	S	0	0	0
			5020	3230	903	880	7			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Mn	0	0
			2	2		

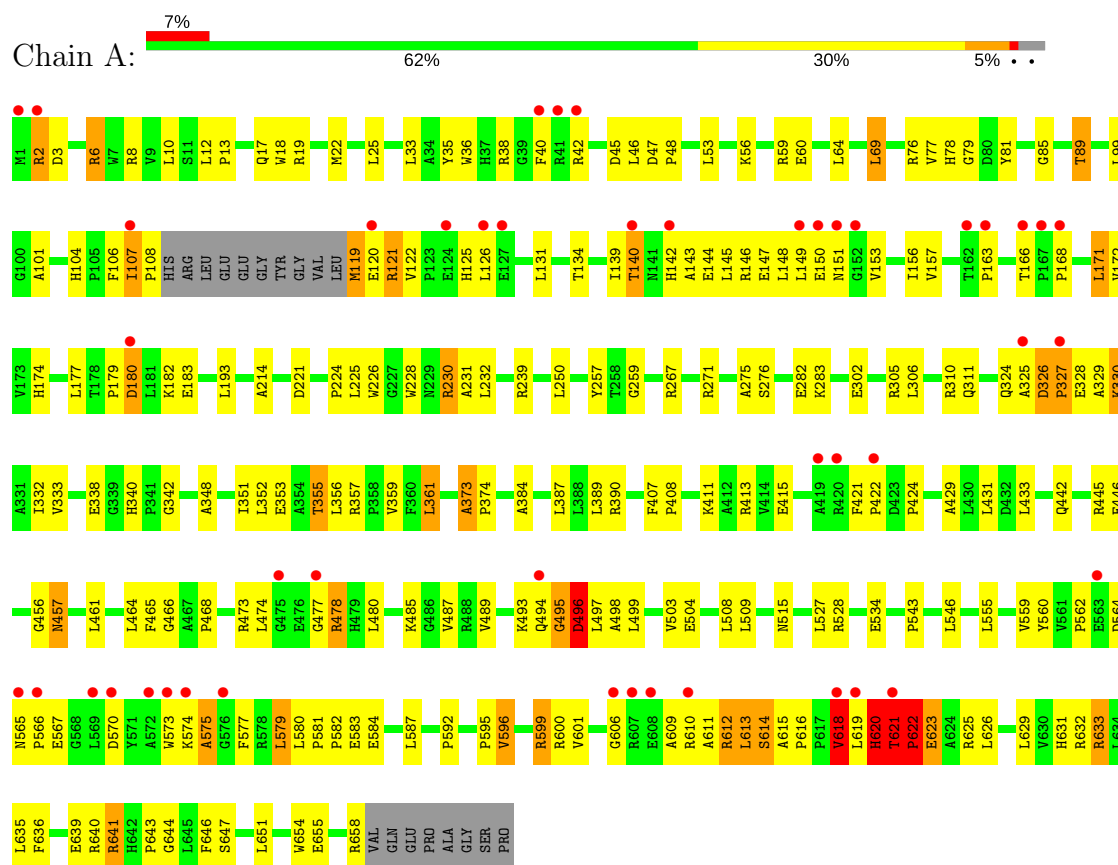
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	293	Total	O	0	0
			293	293		

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: Single-stranded DNA specific exonuclease RecJ



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	83.13Å 83.13Å 249.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.56 – 2.30 49.95 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.4 (41.56-2.30) 99.9 (49.95-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.14 (at 2.29Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.232 , 0.281 0.240 , 0.249	Depositor DCC
R_{free} test set	2005 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	32.7	Xtriage
Anisotropy	0.016	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 50.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5315	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% 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 [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	0.52	0/5145	0.76	6/7012 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	230	ARG	NE-CZ-NH1	-8.60	116.00	120.30
1	A	614	SER	N-CA-C	-6.52	93.40	111.00
1	A	613	LEU	CA-CB-CG	-5.91	101.70	115.30
1	A	239	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	A	239	ARG	NE-CZ-NH2	-5.54	117.53	120.30
1	A	620	HIS	N-CA-C	5.27	125.22	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	5020	0	5130	253	0
2	A	2	0	0	0	0
3	A	293	0	0	29	0
All	All	5315	0	5130	253	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 25.

All (253) 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:614:SER:HB2	1:A:635:LEU:HD21	1.33	1.11
1:A:276:SER:HB2	3:A:920:HOH:O	1.47	1.10
1:A:126:LEU:HD21	1:A:151:ASN:HB3	1.41	1.03
1:A:182:LYS:HA	1:A:182:LYS:HE2	1.47	0.96
1:A:546:LEU:HD22	1:A:606:GLY:HA2	1.47	0.95
1:A:639:GLU:HB3	3:A:959:HOH:O	1.65	0.94
1:A:546:LEU:HD22	1:A:606:GLY:CA	1.98	0.93
1:A:643:PRO:O	3:A:961:HOH:O	1.87	0.93
1:A:457:ASN:HD22	1:A:457:ASN:H	0.93	0.91
1:A:621:THR:HG23	1:A:622:PRO:HD3	1.50	0.91
1:A:329:ALA:O	1:A:330:LYS:HB2	1.71	0.90
1:A:457:ASN:HD22	1:A:457:ASN:N	1.64	0.89
1:A:373:ALA:HB1	1:A:374:PRO:CD	2.03	0.89
1:A:620:HIS:O	1:A:622:PRO:HD2	1.74	0.87
1:A:126:LEU:CD2	1:A:151:ASN:HB3	2.04	0.87
1:A:230:ARG:HD2	3:A:671:HOH:O	1.76	0.85
1:A:442:GLN:NE2	1:A:445:ARG:HH21	1.75	0.84
1:A:340:HIS:HD2	1:A:342:GLY:H	1.21	0.83
1:A:592:PRO:HG2	1:A:654:TRP:CH2	2.14	0.83
1:A:328:GLU:HA	3:A:910:HOH:O	1.78	0.82
1:A:107:ILE:HD13	1:A:108:PRO:HD2	1.60	0.82
1:A:457:ASN:ND2	1:A:457:ASN:H	1.76	0.82
1:A:560:TYR:CE2	1:A:582:PRO:HD3	2.14	0.82
1:A:646:PHE:N	3:A:961:HOH:O	2.14	0.81
1:A:478:ARG:NH1	1:A:495:GLY:HA2	1.97	0.80
1:A:614:SER:HB2	1:A:635:LEU:CD2	2.11	0.80
1:A:477:GLY:HA3	3:A:861:HOH:O	1.82	0.78
1:A:560:TYR:CD2	1:A:582:PRO:HD3	2.18	0.78
1:A:326:ASP:N	1:A:327:PRO:HD2	2.00	0.77
1:A:546:LEU:HB2	1:A:606:GLY:HA2	1.65	0.77
1:A:565:ASN:ND2	1:A:567:GLU:HB3	2.00	0.76
1:A:85:GLY:O	1:A:89:THR:HG23	1.84	0.76
1:A:179:PRO:O	1:A:180:ASP:HB3	1.85	0.76
1:A:565:ASN:CG	1:A:610:ARG:HB2	2.07	0.75
1:A:592:PRO:O	1:A:654:TRP:HZ3	1.72	0.73
1:A:17:GLN:HE21	1:A:36:TRP:HE1	1.34	0.73
1:A:107:ILE:HD13	1:A:108:PRO:CD	2.18	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:546:LEU:CD2	1:A:606:GLY:HA2	2.18	0.72
1:A:126:LEU:HD21	1:A:151:ASN:CB	2.18	0.71
1:A:621:THR:HG23	1:A:622:PRO:CD	2.20	0.71
1:A:139:ILE:HG23	1:A:140:THR:HG22	1.73	0.71
1:A:373:ALA:CB	1:A:374:PRO:CD	2.69	0.70
1:A:581:PRO:HG2	1:A:584:GLU:HG2	1.72	0.70
1:A:442:GLN:HE21	1:A:445:ARG:HH21	1.36	0.70
1:A:17:GLN:NE2	1:A:36:TRP:HE1	1.90	0.69
1:A:326:ASP:H	1:A:327:PRO:HD2	1.55	0.69
1:A:485:LYS:HG3	3:A:955:HOH:O	1.91	0.69
1:A:442:GLN:HE21	1:A:445:ARG:NH2	1.91	0.68
1:A:76:ARG:HE	1:A:104:HIS:HE1	1.40	0.68
1:A:599:ARG:HE	1:A:599:ARG:HA	1.58	0.68
1:A:373:ALA:HB1	1:A:374:PRO:HD2	1.76	0.68
1:A:69:LEU:HD21	1:A:99:LEU:O	1.94	0.68
1:A:478:ARG:HH11	1:A:478:ARG:HG3	1.59	0.68
1:A:559:VAL:HG12	1:A:587:LEU:HD23	1.75	0.67
1:A:592:PRO:HG2	1:A:654:TRP:HH2	1.57	0.67
1:A:351:ILE:O	1:A:355:THR:HB	1.93	0.67
1:A:618:VAL:O	1:A:619:LEU:HD23	1.95	0.67
1:A:327:PRO:O	3:A:910:HOH:O	2.12	0.67
1:A:146:ARG:NH2	1:A:150:GLU:HG2	2.10	0.67
1:A:478:ARG:HH12	1:A:495:GLY:HA2	1.57	0.67
1:A:622:PRO:O	3:A:958:HOH:O	2.13	0.66
1:A:407:PHE:HB3	1:A:408:PRO:HD3	1.78	0.66
1:A:527:LEU:HD23	1:A:528:ARG:N	2.11	0.65
1:A:146:ARG:O	1:A:150:GLU:N	2.29	0.65
1:A:612:ARG:C	1:A:613:LEU:HD22	2.17	0.65
1:A:144:GLU:C	1:A:146:ARG:H	2.00	0.65
1:A:599:ARG:HB3	1:A:601:VAL:HG13	1.79	0.65
1:A:546:LEU:CB	1:A:606:GLY:HA2	2.28	0.64
1:A:559:VAL:CG1	1:A:587:LEU:HD23	2.27	0.64
1:A:565:ASN:HD21	1:A:567:GLU:HB3	1.60	0.63
1:A:620:HIS:O	1:A:622:PRO:CD	2.45	0.63
1:A:324:GLN:O	1:A:326:ASP:OD1	2.16	0.63
1:A:580:LEU:HD22	1:A:584:GLU:OE1	1.99	0.63
1:A:180:ASP:O	1:A:182:LYS:HE3	1.98	0.63
1:A:621:THR:O	1:A:622:PRO:O	2.17	0.62
1:A:411:LYS:O	1:A:415:GLU:HG3	1.99	0.62
1:A:79:GLY:HA3	1:A:89:THR:CG2	2.29	0.62
1:A:25:LEU:HD13	1:A:46:LEU:HD11	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:ARG:CZ	1:A:150:GLU:HG2	2.29	0.62
1:A:611:ALA:C	1:A:613:LEU:H	2.02	0.62
1:A:373:ALA:CB	1:A:374:PRO:HD2	2.30	0.62
1:A:76:ARG:NE	1:A:104:HIS:HE1	1.97	0.61
1:A:621:THR:CG2	1:A:622:PRO:HD3	2.28	0.61
1:A:613:LEU:N	1:A:613:LEU:HD22	2.17	0.60
1:A:182:LYS:HE2	1:A:182:LYS:CA	2.28	0.60
1:A:497:LEU:O	1:A:498:ALA:HB3	2.02	0.60
1:A:25:LEU:HD12	1:A:35:TYR:CD2	2.37	0.59
1:A:230:ARG:CD	3:A:671:HOH:O	2.42	0.59
1:A:457:ASN:ND2	1:A:457:ASN:N	2.38	0.59
1:A:579:LEU:HD12	3:A:777:HOH:O	2.01	0.59
1:A:613:LEU:HD12	1:A:616:PRO:HD3	1.85	0.59
1:A:142:HIS:CE1	1:A:166:THR:HG23	2.38	0.59
1:A:442:GLN:O	1:A:446:GLU:HG3	2.03	0.59
1:A:613:LEU:CD1	1:A:616:PRO:HD3	2.33	0.58
1:A:324:GLN:O	1:A:326:ASP:N	2.36	0.58
1:A:643:PRO:C	3:A:961:HOH:O	2.35	0.58
1:A:373:ALA:HB2	1:A:424:PRO:HB3	1.86	0.57
1:A:106:PHE:CZ	1:A:121:ARG:HG2	2.39	0.57
1:A:53:LEU:HD21	1:A:183:GLU:HG3	1.85	0.57
1:A:326:ASP:O	1:A:327:PRO:O	2.22	0.57
1:A:373:ALA:HB1	1:A:374:PRO:HD3	1.85	0.57
1:A:625:ARG:HD2	3:A:898:HOH:O	2.05	0.57
1:A:13:PRO:HB3	1:A:36:TRP:CD2	2.40	0.57
1:A:282:GLU:H	1:A:282:GLU:CD	2.08	0.56
1:A:373:ALA:CB	1:A:424:PRO:HB3	2.35	0.56
1:A:595:PRO:HD3	1:A:654:TRP:HB3	1.86	0.56
1:A:340:HIS:CD2	1:A:342:GLY:H	2.13	0.56
1:A:76:ARG:HE	1:A:104:HIS:CE1	2.21	0.56
1:A:193:LEU:HD21	1:A:228:TRP:CZ2	2.40	0.56
1:A:583:GLU:HA	1:A:599:ARG:HD2	1.87	0.56
1:A:340:HIS:HE1	3:A:839:HOH:O	1.89	0.55
1:A:60:GLU:CD	1:A:171:LEU:HD11	2.26	0.55
1:A:79:GLY:HA3	1:A:89:THR:HG21	1.88	0.55
1:A:104:HIS:HD2	3:A:901:HOH:O	1.90	0.54
1:A:565:ASN:ND2	1:A:610:ARG:HD2	2.22	0.54
1:A:573:TRP:HE3	3:A:935:HOH:O	1.91	0.54
1:A:633:ARG:NH2	3:A:941:HOH:O	2.40	0.54
1:A:647:SER:N	3:A:961:HOH:O	2.09	0.54
1:A:599:ARG:NE	1:A:599:ARG:HA	2.23	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:546:LEU:HD22	1:A:606:GLY:HA3	1.86	0.53
1:A:126:LEU:HD13	1:A:126:LEU:C	2.29	0.53
1:A:433:LEU:HD21	1:A:636:PHE:HB2	1.91	0.53
1:A:629:LEU:HD23	1:A:629:LEU:C	2.29	0.53
1:A:651:LEU:HA	1:A:654:TRP:CD1	2.44	0.53
1:A:546:LEU:CG	1:A:606:GLY:HA2	2.38	0.52
1:A:144:GLU:C	1:A:146:ARG:N	2.62	0.52
1:A:302:GLU:OE1	1:A:305:ARG:NH1	2.43	0.52
1:A:2:ARG:HD2	1:A:3:ASP:N	2.25	0.52
1:A:355:THR:CG2	1:A:357:ARG:HB2	2.39	0.52
1:A:433:LEU:HD23	1:A:632:ARG:HG2	1.90	0.52
1:A:13:PRO:HB3	1:A:36:TRP:CE3	2.46	0.51
1:A:257:TYR:CZ	1:A:259:GLY:HA2	2.45	0.51
1:A:615:ALA:N	1:A:616:PRO:CD	2.72	0.51
1:A:78:HIS:O	1:A:134:THR:HA	2.10	0.51
1:A:468:PRO:HD3	1:A:503:VAL:HG22	1.93	0.51
1:A:119:MET:HG2	1:A:144:GLU:HG2	1.91	0.51
1:A:180:ASP:OD2	1:A:180:ASP:C	2.49	0.50
1:A:478:ARG:HG2	1:A:493:LYS:O	2.11	0.50
1:A:655:GLU:OE2	1:A:655:GLU:HA	2.11	0.50
1:A:473:ARG:HE	1:A:497:LEU:HD22	1.76	0.50
1:A:139:ILE:O	1:A:168:PRO:HD3	2.12	0.50
1:A:53:LEU:CD2	1:A:183:GLU:HG3	2.41	0.50
1:A:69:LEU:CD1	1:A:101:ALA:HB2	2.41	0.50
1:A:60:GLU:OE2	1:A:171:LEU:HD21	2.12	0.50
1:A:77:VAL:HG12	1:A:89:THR:HB	1.94	0.50
1:A:461:LEU:HD23	1:A:508:LEU:HD23	1.94	0.49
1:A:496:ASP:O	1:A:497:LEU:HB2	2.12	0.49
1:A:560:TYR:OH	1:A:592:PRO:HA	2.12	0.49
1:A:612:ARG:O	1:A:612:ARG:HG2	2.13	0.49
1:A:267:ARG:O	1:A:271:ARG:HD3	2.13	0.48
1:A:38:ARG:HH11	1:A:457:ASN:HB2	1.78	0.48
1:A:478:ARG:CZ	1:A:495:GLY:HA2	2.42	0.48
1:A:565:ASN:HB2	3:A:915:HOH:O	2.12	0.48
1:A:480:LEU:HD12	1:A:480:LEU:C	2.33	0.48
1:A:478:ARG:HH11	1:A:478:ARG:CG	2.26	0.48
1:A:565:ASN:C	1:A:567:GLU:N	2.64	0.48
1:A:139:ILE:CG2	1:A:140:THR:HG22	2.42	0.48
1:A:352:LEU:C	1:A:352:LEU:HD23	2.34	0.48
1:A:611:ALA:C	1:A:613:LEU:N	2.66	0.48
1:A:147:GLU:HB2	1:A:153:VAL:HB	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:ILE:HD12	1:A:355:THR:HG21	1.96	0.47
1:A:348:ALA:HB2	1:A:361:LEU:HD22	1.96	0.47
1:A:48:PRO:HG2	1:A:231:ALA:HB2	1.96	0.47
1:A:76:ARG:NH1	1:A:125:HIS:HD2	2.12	0.47
1:A:610:ARG:O	1:A:613:LEU:HB2	2.14	0.47
1:A:592:PRO:O	1:A:654:TRP:CZ3	2.61	0.47
1:A:149:LEU:HD23	1:A:150:GLU:N	2.29	0.47
1:A:543:PRO:HB2	1:A:641:ARG:HA	1.97	0.47
1:A:79:GLY:HA3	1:A:89:THR:HG22	1.96	0.47
1:A:77:VAL:CG1	1:A:89:THR:HB	2.45	0.47
1:A:565:ASN:N	1:A:566:PRO:HD3	2.29	0.47
1:A:611:ALA:O	1:A:613:LEU:N	2.48	0.47
1:A:338:GLU:HB2	3:A:887:HOH:O	2.14	0.46
1:A:465:PHE:CD2	1:A:504:GLU:HG3	2.50	0.46
1:A:12:LEU:HD22	1:A:18:TRP:CH2	2.51	0.46
1:A:131:LEU:HD11	1:A:156:ILE:HG13	1.98	0.46
1:A:311:GLN:HG2	3:A:952:HOH:O	2.15	0.46
1:A:473:ARG:HG2	1:A:474:LEU:N	2.31	0.46
1:A:19:ARG:HA	1:A:22:MET:HE3	1.98	0.46
1:A:174:HIS:HB3	1:A:177:LEU:HG	1.98	0.46
1:A:581:PRO:CG	1:A:584:GLU:HG2	2.45	0.46
1:A:613:LEU:N	1:A:613:LEU:CD2	2.78	0.46
1:A:142:HIS:CE1	1:A:166:THR:O	2.69	0.46
1:A:134:THR:OG1	1:A:157:VAL:HA	2.15	0.45
1:A:310:ARG:NH1	3:A:952:HOH:O	2.49	0.45
1:A:480:LEU:HD21	1:A:493:LYS:HB2	1.97	0.45
1:A:565:ASN:OD1	1:A:610:ARG:HB2	2.16	0.45
1:A:42:ARG:HH11	1:A:42:ARG:HG3	1.81	0.45
1:A:565:ASN:C	1:A:567:GLU:H	2.20	0.45
1:A:119:MET:HA	1:A:122:VAL:HG23	1.97	0.45
1:A:6:ARG:HB3	1:A:534:GLU:OE1	2.16	0.45
1:A:566:PRO:O	1:A:570:ASP:N	2.50	0.45
1:A:478:ARG:NH1	1:A:478:ARG:HG3	2.29	0.45
1:A:64:LEU:HD22	1:A:171:LEU:HD23	1.99	0.45
1:A:224:PRO:HB3	1:A:226:TRP:CH2	2.52	0.45
1:A:493:LYS:HG2	1:A:497:LEU:CA	2.47	0.45
1:A:493:LYS:HG2	1:A:497:LEU:HA	1.98	0.45
1:A:600:ARG:NH1	3:A:914:HOH:O	2.49	0.45
1:A:142:HIS:HE1	1:A:166:THR:O	1.99	0.44
1:A:565:ASN:HD22	1:A:610:ARG:HD2	1.82	0.44
1:A:635:LEU:O	1:A:639:GLU:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:LEU:H	1:A:456:GLY:N	2.15	0.44
1:A:478:ARG:NH1	1:A:478:ARG:CG	2.81	0.44
1:A:595:PRO:HA	1:A:654:TRP:CE2	2.52	0.44
1:A:559:VAL:HG12	1:A:587:LEU:CD2	2.45	0.44
1:A:631:HIS:O	1:A:635:LEU:HD23	2.17	0.44
1:A:465:PHE:CD1	1:A:466:GLY:N	2.86	0.43
1:A:515:ASN:CG	1:A:515:ASN:O	2.56	0.43
1:A:38:ARG:HB2	1:A:40:PHE:CD1	2.53	0.43
1:A:40:PHE:CE2	1:A:45:ASP:O	2.71	0.43
1:A:464:LEU:HD12	1:A:464:LEU:N	2.34	0.43
1:A:655:GLU:OE1	1:A:658:ARG:NH2	2.52	0.43
1:A:326:ASP:O	1:A:327:PRO:C	2.57	0.43
1:A:497:LEU:O	1:A:498:ALA:CB	2.67	0.43
1:A:172:VAL:O	1:A:172:VAL:HG13	2.17	0.43
1:A:527:LEU:HD23	1:A:527:LEU:C	2.39	0.43
1:A:615:ALA:N	1:A:616:PRO:HD2	2.33	0.43
1:A:625:ARG:HG3	3:A:958:HOH:O	2.18	0.43
1:A:421:PHE:HB3	1:A:422:PRO:HD2	2.00	0.42
1:A:573:TRP:C	1:A:575:ALA:H	2.21	0.42
1:A:596:VAL:HG12	1:A:599:ARG:CG	2.48	0.42
1:A:641:ARG:O	1:A:643:PRO:HD3	2.19	0.42
1:A:384:ALA:HB2	1:A:413:ARG:HG2	2.02	0.42
1:A:389:LEU:C	1:A:390:ARG:HG3	2.40	0.42
1:A:85:GLY:O	1:A:89:THR:CG2	2.62	0.42
1:A:17:GLN:NE2	1:A:36:TRP:NE1	2.64	0.42
1:A:442:GLN:NE2	1:A:445:ARG:NH2	2.50	0.42
1:A:493:LYS:HZ2	1:A:496:ASP:HB2	1.85	0.42
1:A:493:LYS:HG2	1:A:497:LEU:H	1.85	0.42
1:A:275:ALA:HA	1:A:306:LEU:HD13	2.02	0.42
1:A:644:GLY:C	3:A:961:HOH:O	2.58	0.42
1:A:646:PHE:CA	3:A:961:HOH:O	2.65	0.42
1:A:613:LEU:HA	1:A:613:LEU:HD13	1.47	0.42
1:A:38:ARG:HB2	1:A:40:PHE:HD1	1.85	0.42
1:A:225:LEU:H	1:A:456:GLY:H	1.68	0.41
1:A:10:LEU:HD21	1:A:429:ALA:HB1	2.02	0.41
1:A:8:ARG:NH2	1:A:534:GLU:OE2	2.52	0.41
1:A:494:GLN:O	1:A:495:GLY:C	2.58	0.41
1:A:635:LEU:N	1:A:635:LEU:HD22	2.35	0.41
1:A:59:ARG:NH1	3:A:817:HOH:O	2.54	0.41
1:A:623:GLU:O	1:A:623:GLU:HG3	2.20	0.41
1:A:493:LYS:CG	1:A:497:LEU:H	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:VAL:HB	1:A:411:LYS:HG3	2.01	0.41
1:A:38:ARG:NH1	1:A:457:ASN:HB2	2.34	0.41
1:A:606:GLY:HA3	1:A:609:ALA:CB	2.50	0.41
1:A:577:PHE:HB2	1:A:579:LEU:HD22	2.03	0.41
1:A:499:LEU:HD22	1:A:499:LEU:N	2.37	0.40
1:A:581:PRO:HB2	1:A:583:GLU:OE2	2.20	0.40
1:A:214:ALA:HB1	1:A:232:LEU:HD21	2.02	0.40
1:A:359:VAL:HG12	1:A:361:LEU:HD13	2.03	0.40
1:A:562:PRO:HG2	3:A:728:HOH:O	2.22	0.40
1:A:143:ALA:O	1:A:145:LEU:HD23	2.21	0.40
1:A:329:ALA:O	1:A:330:LYS:CB	2.52	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	644/666 (97%)	589 (92%)	39 (6%)	16 (2%)	6 5

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	325	ALA
1	A	327	PRO
1	A	373	ALA
1	A	496	ASP
1	A	620	HIS
1	A	621	THR
1	A	622	PRO
1	A	495	GLY
1	A	612	ARG
1	A	618	VAL

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Mol	Chain	Res	Type
1	A	163	PRO
1	A	180	ASP
1	A	330	LYS
1	A	575	ALA
1	A	574	LYS
1	A	596	VAL

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	502/516 (97%)	459 (91%)	43 (9%)	12	14

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	6	ARG
1	A	33	LEU
1	A	47	ASP
1	A	56	LYS
1	A	69	LEU
1	A	81	TYR
1	A	89	THR
1	A	107	ILE
1	A	119	MET
1	A	120	GLU
1	A	121	ARG
1	A	140	THR
1	A	148	LEU
1	A	171	LEU
1	A	221	ASP
1	A	250	LEU
1	A	283	LYS
1	A	326	ASP
1	A	353	GLU

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Mol	Chain	Res	Type
1	A	355	THR
1	A	356	LEU
1	A	361	LEU
1	A	387	LEU
1	A	431	LEU
1	A	457	ASN
1	A	478	ARG
1	A	487	VAL
1	A	489	VAL
1	A	496	ASP
1	A	509	LEU
1	A	555	LEU
1	A	564	ASP
1	A	579	LEU
1	A	599	ARG
1	A	618	VAL
1	A	621	THR
1	A	622	PRO
1	A	623	GLU
1	A	626	LEU
1	A	633	ARG
1	A	640	ARG
1	A	641	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	71	GLN
1	A	104	HIS
1	A	141	ASN
1	A	142	HIS
1	A	151	ASN
1	A	304	HIS
1	A	311	GLN
1	A	340	HIS
1	A	364	GLN
1	A	394	HIS
1	A	442	GLN
1	A	457	ASN
1	A	494	GLN
1	A	565	ASN

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 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 [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	648/666 (97%)	0.41	46 (7%) 17 22	13, 32, 67, 78	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	7.3
1	A	621	THR	6.3
1	A	607	ARG	6.3
1	A	608	GLU	5.2
1	A	150	GLU	4.9
1	A	325	ALA	4.6
1	A	606	GLY	4.4
1	A	572	ALA	4.3
1	A	565	ASN	4.2
1	A	477	GLY	4.1
1	A	166	THR	4.1
1	A	126	LEU	4.0
1	A	573	TRP	4.0
1	A	167	PRO	3.7
1	A	162	THR	3.6
1	A	574	LYS	3.6
1	A	151	ASN	3.5
1	A	127	GLU	3.4
1	A	107	ILE	3.4
1	A	576	GLY	3.3
1	A	569	LEU	3.2
1	A	42	ARG	3.0
1	A	570	ASP	3.0
1	A	566	PRO	2.9
1	A	422	PRO	2.9
1	A	420	ARG	2.8
1	A	475	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	149	LEU	2.6
1	A	120	GLU	2.6
1	A	41	ARG	2.5
1	A	610	ARG	2.5
1	A	619	LEU	2.4
1	A	142	HIS	2.4
1	A	152	GLY	2.3
1	A	494	GLN	2.3
1	A	168	PRO	2.3
1	A	40	PHE	2.2
1	A	124	GLU	2.2
1	A	140	THR	2.2
1	A	2	ARG	2.2
1	A	327	PRO	2.2
1	A	563	GLU	2.2
1	A	419	ALA	2.1
1	A	163	PRO	2.1
1	A	618	VAL	2.1
1	A	180	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(Å ²)	Q<0.9
2	MN	A	667	1/1	1.00	0.08	-1.70	31,31,31,31	0
2	MN	A	668	1/1	0.93	0.05	-2.44	65,65,65,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.