



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:18 PM UTC

PDB ID : 4DDX / pdb_00004ddx
Title : Thermotoga maritima reverse gyrase, primitive monoclinic form
Authors : Rudolph, M.G.; Klostermeier, D.
Deposited on : 2012-01-19
Resolution : 4.17 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

<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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

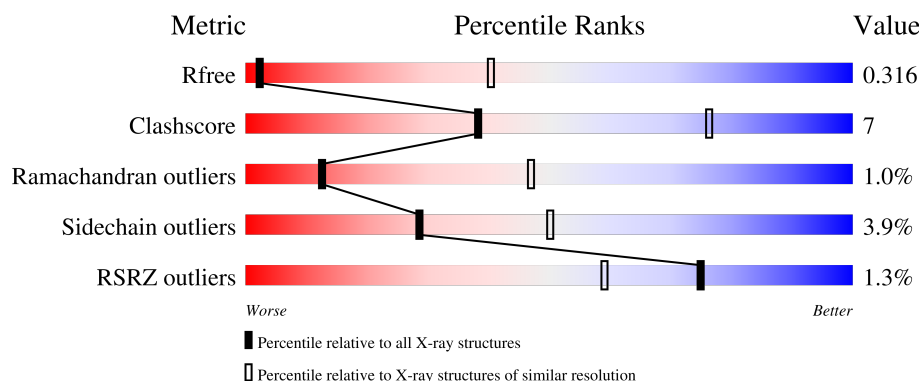
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.17 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}	180053	1026 (4.50-3.86)
Clashscore	190562	1071 (4.50-3.86)
Ramachandran outliers	187476	1014 (4.52-3.84)
Sidechain outliers	187428	1000 (4.52-3.84)
RSRZ outliers	180081	1023 (4.50-3.86)

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 ≥ 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	1104	77% 21% .
1	B	1104	77% 22% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 18064 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 Reverse gyrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1102	Total	C	N	O	S	0	0	0
			9030	5759	1563	1682	26			
1	B	1102	Total	C	N	O	S	0	0	0
			9030	5759	1563	1682	26			

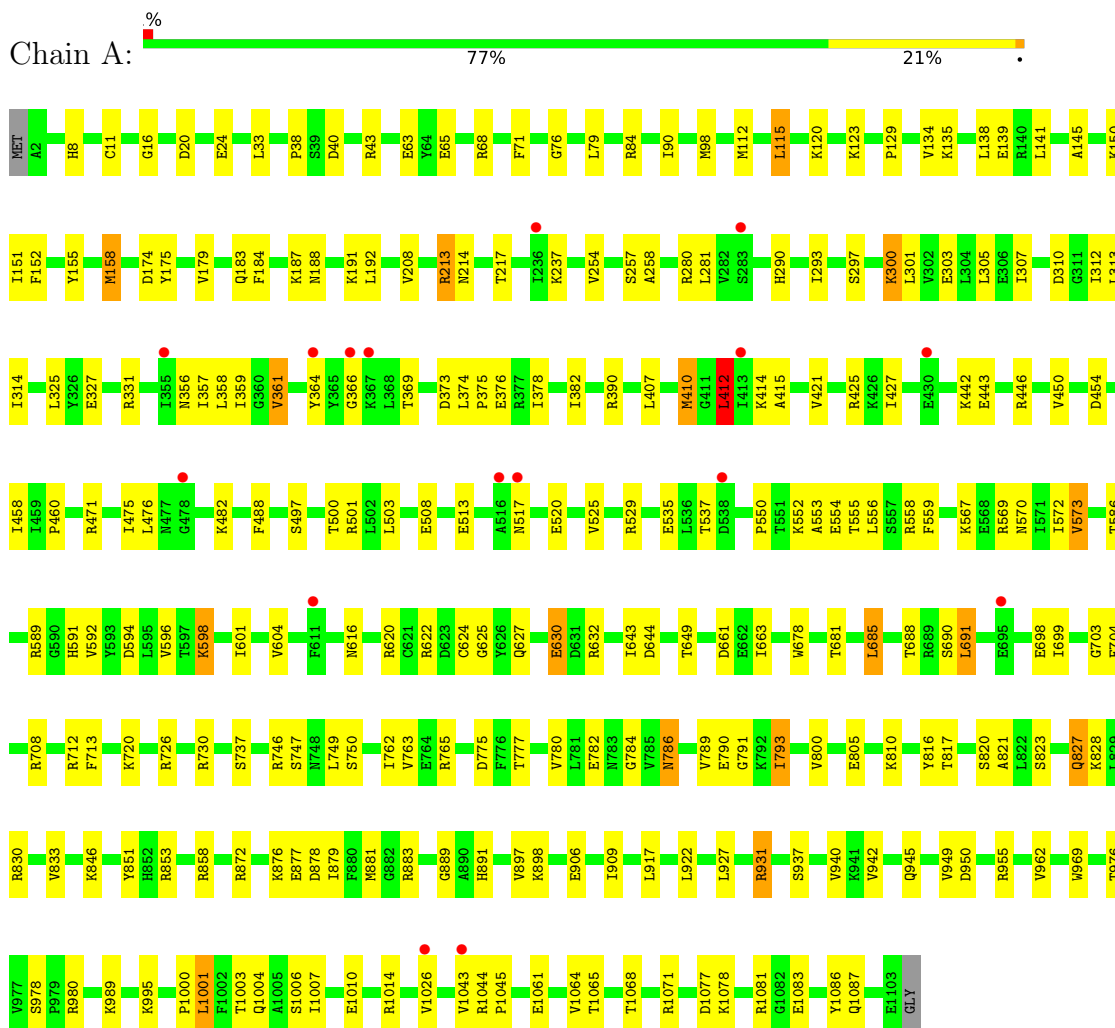
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

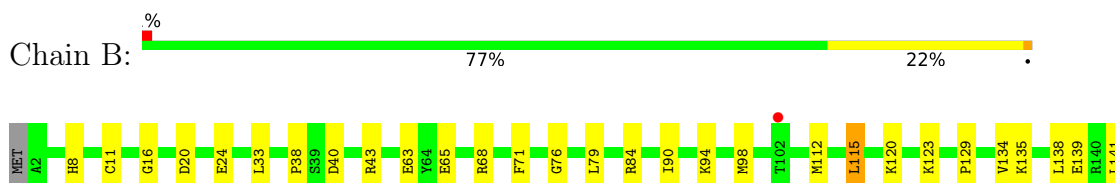
3 Residue-property plots

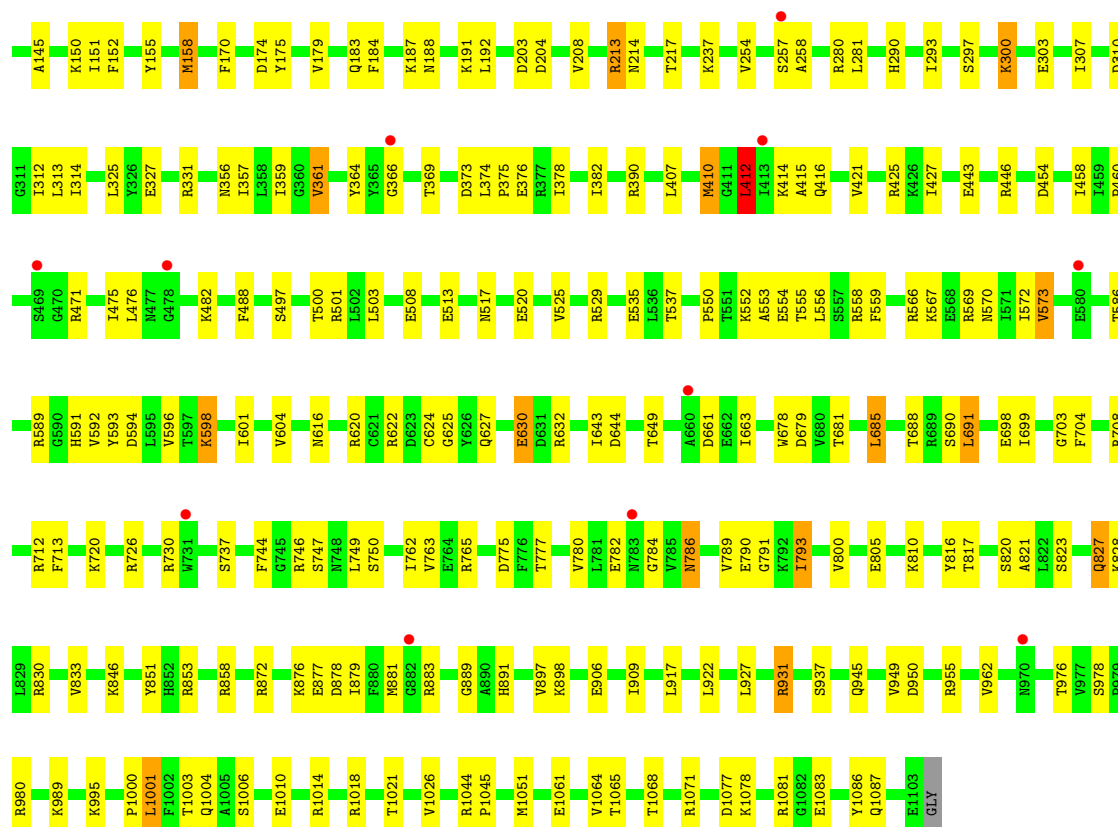
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: Reverse gyrase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	125.81Å 104.90Å 126.18Å 90.00° 91.72° 90.00°	Depositor
Resolution (Å)	48.43 – 4.17 48.43 – 4.17	Depositor EDS
% Data completeness (in resolution range)	96.8 (48.43-4.17) 87.7 (48.43-4.17)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	0.28	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 4.14Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_881)	Depositor
R, R_{free}	0.244 , 0.305 0.257 , 0.316	Depositor DCC
R_{free} test set	1221 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	84.8	Xtriage
Anisotropy	0.684	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 189.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.007 for l,k,-h 0.023 for h,-k,-l 0.239 for l,-k,h	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	18064	wwPDB-VP
Average B, all atoms (Å ²)	141.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.73% 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: ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.26	0/9191	0.70	1/12356 (0.0%)
1	B	0.25	0/9191	0.70	1/12356 (0.0%)
All	All	0.25	0/18382	0.70	2/24712 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	569	ARG	CB-CA-C	-5.28	110.50	116.63
1	B	569	ARG	CB-CA-C	-5.23	110.56	116.63

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	9030	0	9186	126	0
1	B	9030	0	9186	127	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
All	All	18064	0	18372	251	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 7.

All (251) 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:407:LEU:HA	1:B:410:MET:HB2	1.69	0.74
1:A:407:LEU:HA	1:A:410:MET:HB2	1.69	0.73
1:A:897:VAL:HG23	1:A:898:LYS:HG3	1.72	0.72
1:A:375:PRO:HG2	1:A:476:LEU:HD11	1.71	0.72
1:B:375:PRO:HG2	1:B:476:LEU:HD11	1.71	0.71
1:A:780:VAL:HA	1:A:786:ASN:HB3	1.72	0.71
1:B:620:ARG:HG3	1:B:627:GLN:HG2	1.74	0.70
1:B:780:VAL:HA	1:B:786:ASN:HB3	1.72	0.70
1:B:897:VAL:HG23	1:B:898:LYS:HG3	1.72	0.70
1:A:620:ARG:HG3	1:A:627:GLN:HG2	1.73	0.70
1:B:883:ARG:NH2	1:B:937:SER:O	2.25	0.69
1:A:883:ARG:NH2	1:A:937:SER:O	2.25	0.68
1:A:858:ARG:NH1	1:A:891:HIS:O	2.27	0.67
1:B:555:THR:HG21	1:B:846:LYS:HA	1.77	0.67
1:A:555:THR:HG21	1:A:846:LYS:HA	1.76	0.66
1:B:858:ARG:NH1	1:B:891:HIS:O	2.26	0.66
1:A:390:ARG:HG2	1:A:458:ILE:HG12	1.80	0.63
1:A:782:GLU:HA	1:A:989:LYS:HE3	1.80	0.63
1:B:390:ARG:HG2	1:B:458:ILE:HG12	1.79	0.63
1:B:782:GLU:HA	1:B:989:LYS:HE3	1.81	0.62
1:A:237:LYS:HG3	1:B:414:LYS:HE3	1.82	0.62
1:B:20:ASP:OD1	1:B:678:TRP:NE1	2.34	0.60
1:A:314:ILE:HB	1:A:359:ILE:HA	1.84	0.59
1:A:412:LEU:HD13	1:A:427:ILE:HG23	1.84	0.59
1:B:412:LEU:HD13	1:B:427:ILE:HG23	1.84	0.59
1:B:314:ILE:HB	1:B:359:ILE:HA	1.84	0.59
1:B:378:ILE:O	1:B:529:ARG:NH2	2.35	0.59
1:A:688:THR:HG22	1:A:690:SER:H	1.69	0.58
1:A:378:ILE:O	1:A:529:ARG:NH2	2.36	0.58
1:A:115:LEU:HD11	1:A:145:ALA:HB2	1.85	0.58
1:B:115:LEU:HD11	1:B:145:ALA:HB2	1.86	0.57
1:B:290:HIS:NE2	1:B:508:GLU:OE1	2.35	0.57
1:A:65:GLU:HG2	1:A:68:ARG:HH21	1.70	0.57
1:A:482:LYS:O	1:A:529:ARG:NH1	2.38	0.56
1:A:1061:GLU:O	1:A:1065:THR:OG1	2.22	0.56
1:B:482:LYS:O	1:B:529:ARG:NH1	2.38	0.56
1:B:63:GLU:OE2	1:B:120:LYS:NZ	2.38	0.56
1:B:65:GLU:HG2	1:B:68:ARG:HH21	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:688:THR:HG22	1:B:690:SER:H	1.69	0.56
1:A:63:GLU:OE2	1:A:120:LYS:NZ	2.38	0.56
1:B:443:GLU:O	1:B:446:ARG:NH1	2.39	0.55
1:B:663:ILE:HB	1:B:691:LEU:HG	1.88	0.55
1:B:310:ASP:OD1	1:B:356:ASN:ND2	2.38	0.55
1:A:663:ILE:HB	1:A:691:LEU:HG	1.88	0.55
1:A:20:ASP:OD1	1:A:678:TRP:NE1	2.34	0.55
1:A:1001:LEU:HD11	1:A:1044:ARG:HE	1.71	0.55
1:B:762:ILE:HG22	1:B:1045:PRO:HD3	1.88	0.55
1:B:1061:GLU:O	1:B:1065:THR:OG1	2.22	0.55
1:A:290:HIS:NE2	1:A:508:GLU:OE1	2.35	0.55
1:B:878:ASP:OD1	1:B:878:ASP:N	2.36	0.55
1:B:1001:LEU:HD11	1:B:1044:ARG:HE	1.71	0.55
1:A:443:GLU:O	1:A:446:ARG:NH1	2.39	0.54
1:A:817:THR:H	1:A:820:SER:HB2	1.72	0.54
1:B:817:THR:H	1:B:820:SER:HB2	1.73	0.54
1:A:762:ILE:HG22	1:A:1045:PRO:HD3	1.88	0.54
1:A:878:ASP:OD1	1:A:878:ASP:N	2.36	0.54
1:A:300:LYS:NZ	1:A:513:GLU:OE2	2.38	0.54
1:B:594:ASP:OD2	1:B:726:ARG:NH1	2.38	0.54
1:A:33:LEU:HD21	1:A:38:PRO:HG3	1.89	0.53
1:A:310:ASP:OD1	1:A:356:ASN:ND2	2.38	0.53
1:A:589:ARG:NH2	1:A:644:ASP:OD1	2.41	0.53
1:A:950:ASP:OD1	1:A:955:ARG:NE	2.42	0.53
1:B:33:LEU:HD21	1:B:38:PRO:HG3	1.89	0.53
1:A:553:ALA:HA	1:A:556:LEU:HB2	1.91	0.53
1:A:573:VAL:HA	1:A:586:THR:HB	1.91	0.53
1:B:858:ARG:HH11	1:B:889:GLY:HA3	1.73	0.53
1:B:421:VAL:HG12	1:B:425:ARG:HH21	1.74	0.52
1:A:555:THR:HG22	1:A:558:ARG:HH21	1.74	0.52
1:B:553:ALA:HA	1:B:556:LEU:HB2	1.90	0.52
1:A:188:ASN:HB3	1:A:191:LYS:HD2	1.92	0.52
1:B:555:THR:HG22	1:B:558:ARG:HH21	1.74	0.52
1:A:858:ARG:HH11	1:A:889:GLY:HA3	1.73	0.52
1:A:134:VAL:HG13	1:A:179:VAL:HG12	1.92	0.51
1:A:550:PRO:O	1:A:554:GLU:N	2.42	0.51
1:B:950:ASP:OD1	1:B:955:ARG:NE	2.42	0.51
1:A:421:VAL:HG12	1:A:425:ARG:HH21	1.74	0.51
1:B:460:PRO:HB2	1:B:501:ARG:HD2	1.92	0.51
1:A:765:ARG:HG3	1:A:1001:LEU:HD23	1.92	0.51
1:B:188:ASN:HB3	1:B:191:LYS:HD2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:573:VAL:HA	1:B:586:THR:HB	1.91	0.50
1:A:596:VAL:HG22	1:A:616:ASN:HB2	1.92	0.50
1:B:213:ARG:O	1:B:217:THR:OG1	2.25	0.50
1:B:596:VAL:HG22	1:B:616:ASN:HB2	1.92	0.50
1:B:1003:THR:H	1:B:1006:SER:HB2	1.76	0.50
1:A:909:ILE:HD11	1:A:917:LEU:HD12	1.93	0.50
1:B:40:ASP:HB3	1:B:43:ARG:HB3	1.93	0.50
1:B:765:ARG:HG3	1:B:1001:LEU:HD23	1.92	0.50
1:A:1010:GLU:OE1	1:A:1014:ARG:NH1	2.44	0.50
1:B:909:ILE:HD11	1:B:917:LEU:HD12	1.92	0.50
1:A:460:PRO:HB2	1:A:501:ARG:HD2	1.93	0.50
1:B:134:VAL:HG13	1:B:179:VAL:HG12	1.92	0.50
1:B:184:PHE:HE2	1:B:192:LEU:HD21	1.76	0.50
1:A:800:VAL:HA	1:A:949:VAL:HG12	1.94	0.50
1:B:366:GLY:HA2	1:B:471:ARG:HH22	1.77	0.50
1:A:184:PHE:HE2	1:A:192:LEU:HD21	1.76	0.49
1:B:1010:GLU:OE1	1:B:1014:ARG:NH1	2.44	0.49
1:A:40:ASP:HB3	1:A:43:ARG:HB3	1.93	0.49
1:A:141:LEU:HB3	1:A:151:ILE:HD13	1.94	0.49
1:B:141:LEU:HB3	1:B:151:ILE:HD13	1.94	0.49
1:B:550:PRO:O	1:B:554:GLU:N	2.42	0.49
1:A:681:THR:HG22	1:A:691:LEU:HD22	1.94	0.49
1:A:366:GLY:HA2	1:A:471:ARG:HH22	1.77	0.49
1:A:1003:THR:H	1:A:1006:SER:HB2	1.76	0.49
1:B:589:ARG:NH2	1:B:644:ASP:OD1	2.42	0.49
1:B:746:ARG:HB2	1:B:749:LEU:HD11	1.94	0.49
1:A:314:ILE:HD13	1:A:382:ILE:HB	1.95	0.49
1:A:594:ASP:OD2	1:A:726:ARG:NH1	2.37	0.49
1:B:314:ILE:HD13	1:B:382:ILE:HB	1.95	0.49
1:B:79:LEU:HB2	1:B:84:ARG:HG3	1.95	0.49
1:A:98:MET:HE3	1:A:254:VAL:HG11	1.95	0.48
1:A:293:ILE:HD12	1:A:488:PHE:CZ	2.48	0.48
1:B:800:VAL:HA	1:B:949:VAL:HG12	1.95	0.48
1:A:71:PHE:CD2	1:A:112:MET:HG2	2.48	0.48
1:A:79:LEU:HB2	1:A:84:ARG:HG3	1.95	0.48
1:A:777:THR:HG21	1:A:793:ILE:HG12	1.95	0.48
1:B:71:PHE:CD2	1:B:112:MET:HG2	2.49	0.48
1:B:777:THR:HG21	1:B:793:ILE:HG12	1.95	0.48
1:A:213:ARG:O	1:A:217:THR:OG1	2.25	0.48
1:A:828:LYS:HD2	1:A:927:LEU:HD11	1.94	0.48
1:B:517:ASN:HB3	1:B:520:GLU:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:681:THR:HG22	1:B:691:LEU:HD22	1.95	0.48
1:A:517:ASN:HB3	1:A:520:GLU:HB3	1.95	0.48
1:A:746:ARG:HB2	1:A:749:LEU:HD11	1.94	0.48
1:B:816:TYR:HB3	1:B:821:ALA:HB2	1.95	0.48
1:B:300:LYS:NZ	1:B:513:GLU:OE2	2.38	0.47
1:B:777:THR:HG22	1:B:791:GLY:HA2	1.96	0.47
1:B:828:LYS:HD2	1:B:927:LEU:HD11	1.94	0.47
1:B:98:MET:HE3	1:B:254:VAL:HG11	1.95	0.47
1:B:293:ILE:HD12	1:B:488:PHE:CZ	2.48	0.47
1:A:567:LYS:HG3	1:A:572:ILE:HG12	1.97	0.47
1:A:208:VAL:O	1:A:214:ASN:ND2	2.44	0.47
1:A:777:THR:HG22	1:A:791:GLY:HA2	1.96	0.47
1:B:805:GLU:HG2	1:B:945:GLN:HB3	1.97	0.47
1:A:816:TYR:HB3	1:A:821:ALA:HB2	1.96	0.47
1:B:135:LYS:O	1:B:139:GLU:N	2.44	0.47
1:B:567:LYS:HG3	1:B:572:ILE:HG12	1.97	0.47
1:B:823:SER:O	1:B:827:GLN:HB2	2.15	0.47
1:A:726:ARG:HH12	1:A:730:ARG:NH2	2.13	0.46
1:B:258:ALA:H	1:B:280:ARG:HA	1.80	0.46
1:A:823:SER:O	1:A:827:GLN:HB2	2.15	0.46
1:B:150:LYS:HB3	1:B:152:PHE:CE2	2.51	0.46
1:A:258:ALA:H	1:A:280:ARG:HA	1.80	0.46
1:B:622:ARG:NH1	1:B:643:ILE:O	2.48	0.46
1:B:720:LYS:HD3	1:B:1086:TYR:HB2	1.96	0.46
1:A:805:GLU:HG2	1:A:945:GLN:HB3	1.97	0.46
1:B:208:VAL:O	1:B:214:ASN:ND2	2.44	0.46
1:B:726:ARG:HH12	1:B:730:ARG:NH2	2.13	0.46
1:B:624:CYS:SG	1:B:625:GLY:N	2.89	0.46
1:B:737:SER:HB2	1:B:750:SER:HB3	1.97	0.45
1:A:150:LYS:HB3	1:A:152:PHE:CE2	2.51	0.45
1:A:369:THR:HG21	1:A:471:ARG:HH21	1.81	0.45
1:A:622:ARG:NH1	1:A:643:ILE:O	2.48	0.45
1:A:737:SER:HB2	1:A:750:SER:HB3	1.98	0.45
1:B:369:THR:HG21	1:B:471:ARG:HH21	1.81	0.45
1:B:906:GLU:HG2	1:B:922:LEU:HD13	1.99	0.45
1:B:876:LYS:O	1:B:879:ILE:HG12	2.16	0.45
1:A:135:LYS:O	1:A:139:GLU:N	2.44	0.45
1:A:720:LYS:HD3	1:A:1086:TYR:HB2	1.97	0.45
1:A:624:CYS:SG	1:A:625:GLY:N	2.89	0.45
1:A:906:GLU:HG2	1:A:922:LEU:HD13	1.98	0.45
1:B:816:TYR:CE1	1:B:931:ARG:HG3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1068:THR:HG22	1:B:1071:ARG:HH21	1.82	0.45
1:A:876:LYS:O	1:A:879:ILE:HG12	2.16	0.45
1:A:1068:THR:HG22	1:A:1071:ARG:HH21	1.82	0.44
1:B:1078:LYS:HB3	1:B:1083:GLU:HB2	1.99	0.44
1:A:816:TYR:CE1	1:A:931:ARG:HG3	2.52	0.44
1:B:24:GLU:HA	1:B:685:LEU:HD23	1.99	0.44
1:B:203:ASP:HB3	1:B:204:ASP:H	1.67	0.44
1:B:257:SER:HB3	1:B:281:LEU:HB2	2.00	0.44
1:B:775:ASP:HA	1:B:995:LYS:HB3	2.00	0.44
1:B:303:GLU:O	1:B:307:ILE:HG12	2.18	0.44
1:B:816:TYR:CZ	1:B:931:ARG:HG3	2.53	0.44
1:B:833:VAL:HG21	1:B:1026:VAL:HB	2.00	0.44
1:B:698:GLU:HB2	1:B:703:GLY:HA3	2.00	0.44
1:A:712:ARG:NH1	1:A:713:PHE:O	2.49	0.44
1:A:872:ARG:HA	1:A:877:GLU:HB3	2.00	0.44
1:B:872:ARG:HA	1:B:877:GLU:HB3	2.00	0.43
1:A:24:GLU:HA	1:A:685:LEU:HD23	1.99	0.43
1:A:833:VAL:HG21	1:A:1026:VAL:HB	1.99	0.43
1:A:303:GLU:O	1:A:307:ILE:HG12	2.18	0.43
1:A:698:GLU:HB2	1:A:703:GLY:HA3	2.00	0.43
1:A:789:VAL:HG22	1:A:790:GLU:H	1.84	0.43
1:B:135:LYS:HE3	1:B:155:TYR:CZ	2.53	0.43
1:B:123:LYS:HD3	1:B:175:TYR:CE1	2.54	0.43
1:B:570:ASN:OD1	1:B:649:THR:OG1	2.37	0.43
1:A:816:TYR:CZ	1:A:931:ARG:HG3	2.53	0.43
1:A:570:ASN:OD1	1:A:649:THR:OG1	2.37	0.43
1:A:775:ASP:HA	1:A:995:LYS:HB3	2.00	0.43
1:A:1078:LYS:HB3	1:A:1083:GLU:HB2	1.99	0.43
1:B:789:VAL:HG22	1:B:790:GLU:H	1.84	0.43
1:A:135:LYS:HE3	1:A:155:TYR:CZ	2.54	0.43
1:A:257:SER:HB3	1:A:281:LEU:HB2	2.00	0.43
1:A:497:SER:O	1:A:500:THR:OG1	2.37	0.42
1:A:552:LYS:HE2	1:A:699:ILE:HD11	2.01	0.42
1:A:414:LYS:HE3	1:B:237:LYS:HG3	2.00	0.42
1:A:556:LEU:HA	1:A:559:PHE:CE2	2.54	0.42
1:A:589:ARG:HH22	1:A:644:ASP:CG	2.26	0.42
1:A:685:LEU:HA	1:A:685:LEU:HD12	1.80	0.42
1:A:782:GLU:C	1:A:784:GLY:H	2.27	0.42
1:A:962:VAL:HG11	1:A:980:ARG:HG3	2.01	0.42
1:A:123:LYS:HD3	1:A:175:TYR:CE1	2.54	0.42
1:A:833:VAL:HG13	1:A:1004:GLN:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:552:LYS:HE2	1:B:699:ILE:HD11	2.02	0.42
1:B:556:LEU:HA	1:B:559:PHE:CE2	2.54	0.42
1:B:630:GLU:HG3	1:B:632:ARG:NH1	2.35	0.42
1:B:962:VAL:HG11	1:B:980:ARG:HG3	2.01	0.42
1:A:358:LEU:HD23	1:A:358:LEU:HA	1.89	0.42
1:B:497:SER:O	1:B:500:THR:OG1	2.37	0.42
1:A:312:ILE:HD12	1:A:357:ILE:HG12	2.01	0.42
1:A:704:PHE:CZ	1:A:708:ARG:HD2	2.55	0.42
1:B:704:PHE:CZ	1:B:708:ARG:HD2	2.55	0.42
1:B:877:GLU:H	1:B:877:GLU:CD	2.28	0.42
1:A:11:CYS:HB3	1:A:16:GLY:H	1.85	0.41
1:A:374:LEU:HG	1:A:376:GLU:HG2	2.02	0.41
1:A:877:GLU:CD	1:A:877:GLU:H	2.28	0.41
1:B:482:LYS:HG2	1:B:525:VAL:HG22	2.02	0.41
1:B:782:GLU:C	1:B:784:GLY:H	2.27	0.41
1:A:763:VAL:HA	1:A:1045:PRO:HG3	2.03	0.41
1:B:11:CYS:HB3	1:B:16:GLY:H	1.85	0.41
1:A:596:VAL:HG12	1:A:598:LYS:H	1.85	0.41
1:A:940:VAL:HG12	1:A:942:VAL:HG13	2.02	0.41
1:B:327:GLU:O	1:B:331:ARG:HG3	2.21	0.41
1:A:630:GLU:HG3	1:A:632:ARG:NH1	2.35	0.41
1:A:1007:ILE:HD11	1:A:1043:VAL:HG11	2.03	0.41
1:B:312:ILE:HD12	1:B:357:ILE:HG12	2.01	0.41
1:B:596:VAL:HG12	1:B:598:LYS:H	1.86	0.41
1:B:685:LEU:HD12	1:B:685:LEU:HA	1.81	0.41
1:B:851:TYR:CZ	1:B:853:ARG:HB2	2.56	0.41
1:A:482:LYS:HG2	1:A:525:VAL:HG22	2.02	0.41
1:B:155:TYR:CE1	1:B:158:MET:HA	2.56	0.41
1:A:155:TYR:CE1	1:A:158:MET:HA	2.56	0.41
1:A:327:GLU:O	1:A:331:ARG:HG3	2.21	0.41
1:A:940:VAL:HG23	1:A:969:TRP:CD1	2.56	0.41
1:B:833:VAL:HG13	1:B:1004:GLN:HB3	2.02	0.41
1:B:601:ILE:HB	1:B:604:VAL:HB	2.02	0.41
1:B:763:VAL:HA	1:B:1045:PRO:HG3	2.03	0.41
1:B:1018:ARG:O	1:B:1021:THR:N	2.51	0.41
1:A:1077:ASP:OD1	1:A:1081:ARG:NH2	2.54	0.41
1:B:170:PHE:HE1	1:B:191:LYS:HB3	1.86	0.41
1:B:712:ARG:NH1	1:B:713:PHE:O	2.50	0.41
1:B:1077:ASP:OD1	1:B:1081:ARG:NH2	2.54	0.41
1:A:601:ILE:HB	1:A:604:VAL:HB	2.02	0.40
1:A:851:TYR:CZ	1:A:853:ARG:HB2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:LEU:O	1:A:305:LEU:HG	2.21	0.40
1:B:94:LYS:HD3	1:B:566:ARG:HH22	1.86	0.40
1:B:374:LEU:HG	1:B:376:GLU:HG2	2.02	0.40
1:B:589:ARG:HH22	1:B:644:ASP:CG	2.26	0.40
1:B:593:TYR:HD1	1:B:679:ASP:HB3	1.86	0.40
1:B:744:PHE:HE1	1:B:1051:MET:HE2	1.87	0.40
1:B:414:LYS:O	1:B:416:GLN:N	2.54	0.40
1:A:442:LYS:HA	1:A:450:VAL:HG21	2.04	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	1100/1104 (100%)	979 (89%)	110 (10%)	11 (1%)	12	47
1	B	1100/1104 (100%)	979 (89%)	110 (10%)	11 (1%)	12	47
All	All	2200/2208 (100%)	1958 (89%)	220 (10%)	22 (1%)	12	47

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	129	PRO
1	B	129	PRO
1	A	174	ASP
1	A	412	LEU
1	A	535	GLU
1	A	598	LYS
1	B	174	ASP
1	B	412	LEU
1	B	535	GLU
1	B	598	LYS

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Mol	Chain	Res	Type
1	A	415	ALA
1	B	415	ALA
1	A	747	SER
1	B	747	SER
1	A	373	ASP
1	B	373	ASP
1	A	76	GLY
1	B	76	GLY
1	A	1000	PRO
1	B	1000	PRO
1	A	361	VAL
1	B	361	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	992/993 (100%)	953 (96%)	39 (4%)	28	50
1	B	992/993 (100%)	953 (96%)	39 (4%)	28	50
All	All	1984/1986 (100%)	1906 (96%)	78 (4%)	28	50

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	HIS
1	A	90	ILE
1	A	115	LEU
1	A	138	LEU
1	A	158	MET
1	A	183	GLN
1	A	187	LYS
1	A	213	ARG
1	A	297	SER
1	A	300	LYS
1	A	313	LEU

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Mol	Chain	Res	Type
1	A	325	LEU
1	A	361	VAL
1	A	364	TYR
1	A	410	MET
1	A	412	LEU
1	A	454	ASP
1	A	475	ILE
1	A	503	LEU
1	A	537	THR
1	A	573	VAL
1	A	591	HIS
1	A	592	VAL
1	A	630	GLU
1	A	661	ASP
1	A	685	LEU
1	A	691	LEU
1	A	786	ASN
1	A	793	ILE
1	A	810	LYS
1	A	827	GLN
1	A	830	ARG
1	A	881	MET
1	A	931	ARG
1	A	976	THR
1	A	978	SER
1	A	1001	LEU
1	A	1064	VAL
1	A	1087	GLN
1	B	8	HIS
1	B	90	ILE
1	B	115	LEU
1	B	138	LEU
1	B	158	MET
1	B	183	GLN
1	B	187	LYS
1	B	213	ARG
1	B	297	SER
1	B	300	LYS
1	B	313	LEU
1	B	325	LEU
1	B	361	VAL
1	B	364	TYR

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Mol	Chain	Res	Type
1	B	410	MET
1	B	412	LEU
1	B	454	ASP
1	B	475	ILE
1	B	503	LEU
1	B	537	THR
1	B	573	VAL
1	B	591	HIS
1	B	592	VAL
1	B	630	GLU
1	B	661	ASP
1	B	685	LEU
1	B	691	LEU
1	B	786	ASN
1	B	793	ILE
1	B	810	LYS
1	B	827	GLN
1	B	830	ARG
1	B	881	MET
1	B	931	ARG
1	B	976	THR
1	B	978	SER
1	B	1001	LEU
1	B	1064	VAL
1	B	1087	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	434	GLN
1	A	728	GLN
1	A	834	GLN
1	A	852	HIS
1	A	865	ASN
1	B	431	HIS
1	B	434	GLN
1	B	834	GLN
1	B	852	HIS
1	B	865	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 oligosaccharides 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 [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	1102/1104 (99%)	0.50	16 (1%) 72 56	23, 135, 216, 337	0
1	B	1102/1104 (99%)	0.47	12 (1%) 78 62	41, 135, 210, 367	0
All	All	2204/2208 (99%)	0.48	28 (1%) 75 59	23, 135, 212, 367	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	283	SER	3.5
1	A	695	GLU	3.5
1	A	366	GLY	3.0
1	B	731	TRP	2.9
1	B	413	ILE	2.9
1	A	413	ILE	2.8
1	B	102	THR	2.8
1	A	516	ALA	2.7
1	A	1026	VAL	2.5
1	A	364	TYR	2.5
1	B	580	GLU	2.5
1	B	478	GLY	2.4
1	A	430	GLU	2.3
1	B	783	ASN	2.3
1	A	611	PHE	2.3
1	A	478	GLY	2.3
1	B	366	GLY	2.3
1	B	970	ASN	2.2
1	A	517	ASN	2.2
1	A	236	ILE	2.1
1	A	367	LYS	2.1
1	A	1043	VAL	2.1
1	A	538	ASP	2.1
1	B	469	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	355	ILE	2.1
1	B	660	ALA	2.1
1	B	257	SER	2.1
1	B	882	GLY	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	A	1201	1/1	0.93	0.06	203,203,203,203	0
2	ZN	B	1201	1/1	0.96	0.05	189,189,189,189	0
2	ZN	A	1202	1/1	0.97	0.04	125,125,125,125	0
2	ZN	B	1202	1/1	0.99	0.02	101,101,101,101	0

6.5 Other polymers [i](#)

There are no such residues in this entry.