



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 04:29 AM GMT

PDB ID : 2GAJ
Title : Structure of Full Length Topoisomerase I from *Thermotoga maritima* in monoclinic crystal form
Authors : Hansen, G.
Deposited on : 2006-03-09
Resolution : 1.95 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

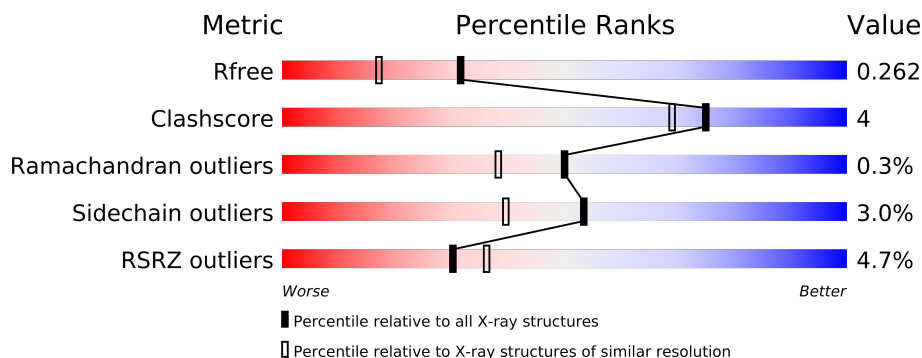
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.95 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}	66092	1321 (1.96-1.96)
Clashscore	79885	1488 (1.96-1.96)
Ramachandran outliers	78287	1475 (1.96-1.96)
Sidechain outliers	78261	1475 (1.96-1.96)
RSRZ outliers	66119	1321 (1.96-1.96)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	633	
1	B	633	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10305 atoms, of which 0 are hydrogen and 0 are deuterium.

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 topoisomerase I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	581	Total	C	N	O	S	0	0	0
			4714	3017	799	881	17			
1	B	581	Total	C	N	O	S	0	0	0
			4714	3017	799	881	17			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	INITIATING METHIONINE	UNP P46799
A	2	ALA	-	CLONING ARTIFACT	UNP P46799
B	1	MET	-	INITIATING METHIONINE	UNP P46799
B	2	ALA	-	CLONING ARTIFACT	UNP P46799

- Molecule 2 is water.

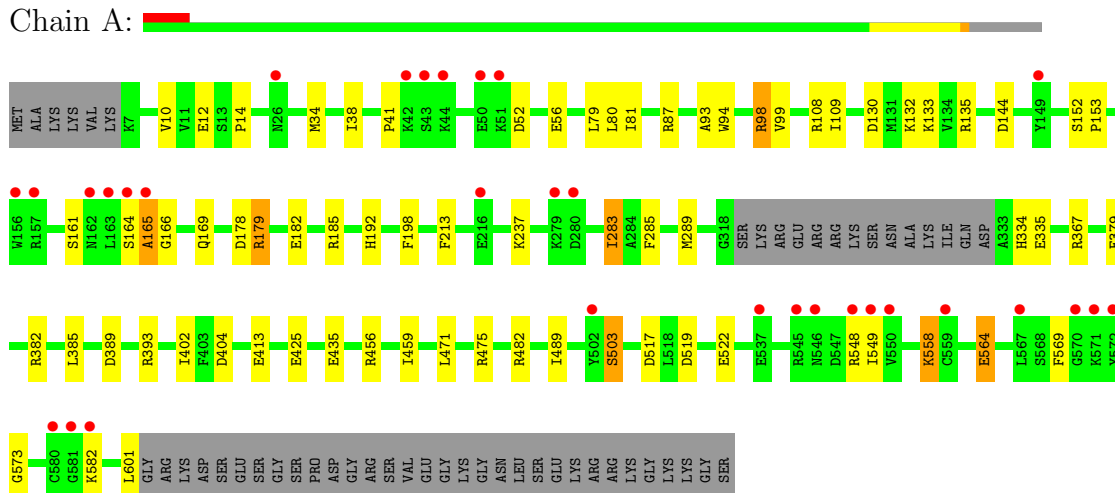
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	427	Total	O	0	0
			427	427		
2	B	450	Total	O	0	0
			450	450		

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 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.

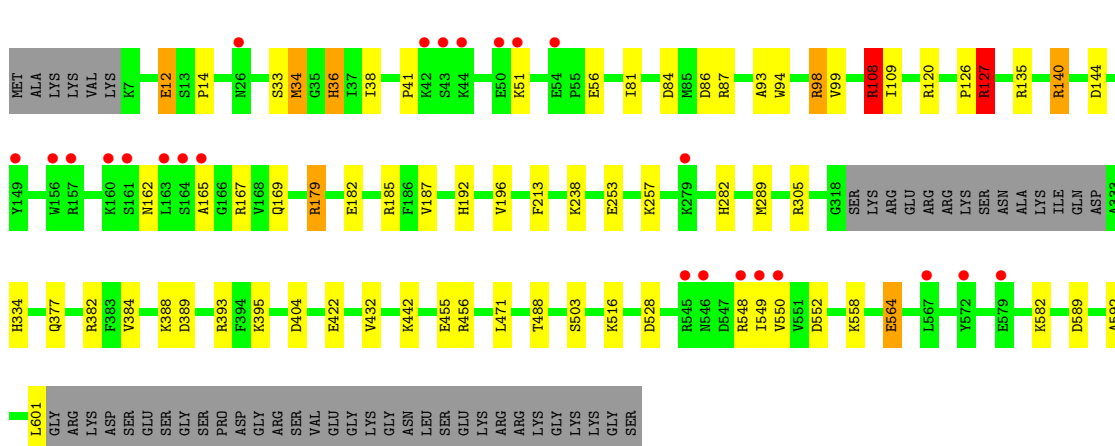
• Molecule 1: DNA topoisomerase I

Chain A:



• Molecule 1: DNA topoisomerase I

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	45.05Å 142.71Å 127.64Å 90.00° 90.84° 90.00°	Depositor
Resolution (Å)	15.00 – 1.95 14.99 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.6 (15.00-1.95) 99.8 (14.99-1.95)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.195 , 0.246 0.213 , 0.262	Depositor DCC
R_{free} test set	5837 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	20.5	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 41.0	EDS
Estimated twinning fraction	0.000 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 116577 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10305	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.74 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.2655e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	0.67	0/4799	1.07	23/6441 (0.4%)
1	B	0.69	0/4799	1.16	27/6441 (0.4%)
All	All	0.68	0/9598	1.12	50/12882 (0.4%)

There are no bond length outliers.

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	179	ARG	NE-CZ-NH2	-25.81	107.40	120.30
1	B	179	ARG	NE-CZ-NH1	18.68	129.64	120.30
1	A	456	ARG	NE-CZ-NH2	-11.42	114.59	120.30
1	B	127	ARG	NE-CZ-NH1	10.76	125.68	120.30
1	B	86	ASP	CB-CG-OD1	10.51	127.76	118.30
1	A	519	ASP	CB-CG-OD1	10.44	127.69	118.30
1	B	86	ASP	CB-CG-OD2	-10.25	109.07	118.30
1	B	127	ARG	NE-CZ-NH2	-8.66	115.97	120.30
1	A	456	ARG	NE-CZ-NH1	8.47	124.54	120.30
1	A	98	ARG	NE-CZ-NH2	-8.31	116.14	120.30
1	B	404	ASP	CB-CG-OD1	8.23	125.71	118.30
1	B	179	ARG	CG-CD-NE	-8.10	94.79	111.80
1	B	179	ARG	CD-NE-CZ	8.00	134.80	123.60
1	B	404	ASP	CB-CG-OD2	-7.89	111.20	118.30
1	A	393	ARG	NE-CZ-NH1	6.95	123.78	120.30
1	A	367	ARG	NE-CZ-NH2	-6.89	116.85	120.30
1	A	179	ARG	NE-CZ-NH2	-6.87	116.87	120.30
1	A	144	ASP	CB-CG-OD1	6.86	124.48	118.30
1	A	519	ASP	CB-CG-OD2	-6.78	112.19	118.30
1	B	185	ARG	NE-CZ-NH2	-6.73	116.93	120.30
1	B	98	ARG	NE-CZ-NH2	-6.64	116.98	120.30
1	B	84	ASP	CB-CG-OD2	-6.57	112.39	118.30
1	A	404	ASP	CB-CG-OD2	-6.56	112.40	118.30
1	A	475	ARG	NE-CZ-NH2	-6.52	117.04	120.30
1	B	135	ARG	NE-CZ-NH1	6.36	123.48	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	98	ARG	NE-CZ-NH1	6.33	123.46	120.30
1	A	178	ASP	CB-CG-OD1	6.30	123.97	118.30
1	B	389	ASP	CB-CG-OD1	6.30	123.97	118.30
1	A	482	ARG	NE-CZ-NH1	6.19	123.39	120.30
1	B	382	ARG	NE-CZ-NH1	6.17	123.39	120.30
1	B	140	ARG	NE-CZ-NH2	-6.16	117.22	120.30
1	B	305	ARG	NE-CZ-NH1	6.16	123.38	120.30
1	A	130	ASP	CB-CG-OD1	6.09	123.78	118.30
1	B	305	ARG	NE-CZ-NH2	-6.01	117.29	120.30
1	A	389	ASP	CB-CG-OD1	5.87	123.58	118.30
1	A	517	ASP	CB-CG-OD1	5.85	123.56	118.30
1	A	135	ARG	NE-CZ-NH2	-5.83	117.39	120.30
1	A	482	ARG	NE-CZ-NH2	-5.69	117.45	120.30
1	B	528	ASP	CB-CG-OD1	5.64	123.37	118.30
1	B	140	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	B	12	GLU	CA-CB-CG	5.62	125.76	113.40
1	A	185	ARG	NE-CZ-NH2	-5.55	117.53	120.30
1	B	456	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	B	120	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	B	87	ARG	NE-CZ-NH2	-5.49	117.56	120.30
1	A	404	ASP	CB-CG-OD1	5.41	123.17	118.30
1	A	335	GLU	OE1-CD-OE2	5.36	129.74	123.30
1	B	144	ASP	CB-CG-OD1	5.24	123.02	118.30
1	A	382	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	B	108	ARG	NE-CZ-NH1	5.12	122.86	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4714	0	4821	35	0
1	B	4714	0	4821	42	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	427	0	0	7	0
2	B	450	0	0	13	0
All	All	10305	0	9642	77	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (77) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:489:ILE:HD11	1:A:601:LEU:HD21	1.56	0.86
1:A:80:LEU:HD13	1:A:109:ILE:HD11	1.58	0.84
1:A:413:GLU:O	2:A:892:HOH:O	1.96	0.82
1:A:334:HIS:HD2	2:A:729:HOH:O	1.65	0.80
1:B:192:HIS:HD2	1:B:213:PHE:H	1.31	0.77
1:A:334:HIS:CD2	2:A:729:HOH:O	2.37	0.77
1:B:253:GLU:OE1	1:B:257:LYS:HE2	1.88	0.73
1:B:94:TRP:CE2	1:B:98:ARG:HD2	2.25	0.71
1:B:14:PRO:HG3	1:B:34:MET:HE1	1.73	0.70
1:A:80:LEU:HB3	1:A:109:ILE:HD12	1.73	0.70
1:B:187:VAL:O	2:B:1009:HOH:O	2.10	0.69
1:B:282:HIS:HE1	2:B:802:HOH:O	1.78	0.66
1:A:41:PRO:HD3	1:A:56:GLU:O	1.98	0.62
1:A:169:GLN:NE2	1:A:471:LEU:HD21	2.14	0.62
1:B:94:TRP:CZ2	1:B:98:ARG:HD2	2.37	0.60
1:B:592:ALA:HB2	1:B:601:LEU:HD23	1.84	0.60
1:A:379:GLU:HG3	1:A:402:ILE:HD11	1.84	0.59
1:A:10:VAL:CG2	1:A:79:LEU:HD11	2.32	0.59
1:A:192:HIS:HD2	1:A:213:PHE:H	1.50	0.58
1:B:14:PRO:HG3	1:B:34:MET:CE	2.35	0.56
1:A:334:HIS:CD2	2:A:1024:HOH:O	2.57	0.55
1:B:41:PRO:HD3	1:B:56:GLU:O	2.07	0.55
1:B:38:ILE:HD11	1:B:99:VAL:HG11	1.89	0.55
1:B:140:ARG:HD2	2:B:761:HOH:O	2.06	0.54
1:B:167:ARG:HD3	2:B:635:HOH:O	2.08	0.54
1:A:165:ALA:HB2	1:A:471:LEU:HD13	1.90	0.54
1:B:127:ARG:HG3	1:B:127:ARG:HH11	1.73	0.53
1:B:34:MET:CA	1:B:34:MET:HE3	2.39	0.52
1:A:425:GLU:OE2	2:A:847:HOH:O	2.19	0.51
1:B:81:ILE:HG21	1:B:93:ALA:HA	1.92	0.50
1:B:516:LYS:HG2	2:B:907:HOH:O	2.11	0.50
1:B:34:MET:HA	1:B:34:MET:HE3	1.91	0.50
1:B:552:ASP:OD2	2:B:945:HOH:O	2.19	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:182:GLU:OE2	1:B:442:LYS:NZ	2.40	0.50
1:B:282:HIS:CE1	2:B:802:HOH:O	2.59	0.49
1:A:94:TRP:CE2	1:A:98:ARG:HD2	2.49	0.48
1:B:169:GLN:NE2	1:B:471:LEU:CD2	2.77	0.48
1:B:334:HIS:HD2	2:B:661:HOH:O	1.97	0.47
1:B:14:PRO:HD3	1:B:34:MET:CE	2.44	0.47
1:B:388:LYS:NZ	2:B:972:HOH:O	2.25	0.47
1:A:80:LEU:HD13	1:A:109:ILE:CD1	2.38	0.47
1:A:94:TRP:CZ2	1:A:98:ARG:HD2	2.50	0.47
1:B:238:LYS:HG2	1:B:377:GLN:HG2	1.96	0.47
1:A:14:PRO:HG3	1:A:34:MET:SD	2.55	0.47
1:A:165:ALA:CB	1:A:471:LEU:HD13	2.43	0.47
1:A:198:PHE:CZ	1:A:385:LEU:HD11	2.49	0.47
1:A:38:ILE:HD11	1:A:99:VAL:CG1	2.46	0.46
1:A:152:SER:N	1:A:153:PRO:CD	2.80	0.45
1:B:488:THR:HB	1:B:589:ASP:HA	1.98	0.45
1:B:384:VAL:HG22	1:B:395:LYS:HG2	1.97	0.45
1:A:237:LYS:NZ	2:A:943:HOH:O	2.49	0.44
1:B:34:MET:CE	1:B:34:MET:HA	2.47	0.44
1:B:165:ALA:HB3	1:B:471:LEU:HD13	1.99	0.44
1:B:33:SER:C	1:B:34:MET:HG2	2.39	0.43
1:A:179:ARG:NH1	1:A:182:GLU:OE1	2.51	0.43
1:B:558:LYS:HE3	1:B:564:GLU:HB3	2.01	0.43
1:A:132:LYS:HE2	2:A:866:HOH:O	2.18	0.43
1:A:569:PHE:CE1	1:A:573:GLY:HA2	2.53	0.43
1:A:38:ILE:HD11	1:A:99:VAL:HG11	2.01	0.43
1:A:558:LYS:NZ	1:A:564:GLU:HB2	2.33	0.43
1:B:108:ARG:HD2	1:B:127:ARG:NH1	2.35	0.42
1:B:282:HIS:HD2	2:B:678:HOH:O	2.03	0.42
1:B:109:ILE:CD1	1:B:126:PRO:HB3	2.50	0.42
1:B:196:VAL:HG12	1:B:432:VAL:HG22	2.00	0.42
1:B:282:HIS:CD2	2:B:678:HOH:O	2.73	0.42
1:B:38:ILE:HD11	1:B:99:VAL:CG1	2.48	0.42
1:A:164:SER:C	1:A:166:GLY:H	2.22	0.41
1:B:169:GLN:NE2	1:B:471:LEU:HD21	2.35	0.41
1:A:81:ILE:HG21	1:A:93:ALA:HA	2.02	0.41
1:A:283:ILE:HD11	1:A:285:PHE:CZ	2.54	0.41
1:A:80:LEU:HB3	1:A:109:ILE:CD1	2.47	0.41
1:B:34:MET:CE	1:B:34:MET:CA	2.98	0.41
1:B:36:HIS:CD2	2:B:942:HOH:O	2.74	0.41
1:A:459:ILE:HA	1:A:459:ILE:HD13	1.78	0.41
1:B:516:LYS:HG3	2:B:932:HOH:O	2.21	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:10:VAL:HG22	1:A:79:LEU:HD11	2.04	0.40
1:A:87:ARG:NE	1:A:522:GLU:OE1	2.45	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	577/633 (91%)	561 (97%)	14 (2%)	2 (0%)	50 38
1	B	577/633 (91%)	562 (97%)	14 (2%)	1 (0%)	56 46
All	All	1154/1266 (91%)	1123 (97%)	28 (2%)	3 (0%)	50 38

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	550	VAL
1	A	503	SER
1	A	165	ALA

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	519/562 (92%)	505 (97%)	14 (3%)	57 45
1	B	519/562 (92%)	502 (97%)	17 (3%)	50 35
All	All	1038/1124 (92%)	1007 (97%)	31 (3%)	53 39

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	GLU
1	A	52	ASP
1	A	108	ARG
1	A	133	LYS
1	A	161	SER
1	A	283	ILE
1	A	289	MET
1	A	435	GLU
1	A	503	SER
1	A	548	ARG
1	A	549	ILE
1	A	558	LYS
1	A	564	GLU
1	A	582	LYS
1	B	12	GLU
1	B	34	MET
1	B	36	HIS
1	B	51	LYS
1	B	108	ARG
1	B	127	ARG
1	B	162	ASN
1	B	179	ARG
1	B	289	MET
1	B	393	ARG
1	B	422	GLU
1	B	455	GLU
1	B	503	SER
1	B	548	ARG
1	B	549	ILE
1	B	564	GLU
1	B	582	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	ASN
1	A	158	ASN
1	A	192	HIS
1	A	334	HIS
1	A	523	GLN
1	A	557	GLN
1	B	26	ASN
1	B	36	HIS

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Mol	Chain	Res	Type
1	B	158	ASN
1	B	169	GLN
1	B	192	HIS
1	B	282	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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 ⓘ

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	581/633 (91%)	0.23	31 (5%)	25 31	2, 10, 24, 34	0
1	B	581/633 (91%)	0.17	24 (4%)	35 40	2, 9, 24, 36	0
All	All	1162/1266 (91%)	0.20	55 (4%)	30 36	2, 9, 24, 36	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	165	ALA	5.6
1	B	549	ILE	5.5
1	A	43	SER	5.3
1	B	165	ALA	5.2
1	A	572	TYR	5.1
1	B	43	SER	4.9
1	B	164	SER	4.8
1	A	280	ASP	4.5
1	A	156	TRP	4.3
1	B	156	TRP	4.2
1	A	549	ILE	4.2
1	A	164	SER	4.2
1	A	570	GLY	3.9
1	A	279	LYS	3.8
1	A	545	ARG	3.8
1	B	163	LEU	3.6
1	A	580	CYS	3.6
1	B	546	ASN	3.5
1	B	157	ARG	3.5
1	B	50	GLU	3.3
1	B	548	ARG	3.3
1	B	550	VAL	3.3
1	A	51	LYS	3.3
1	A	50	GLU	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	548	ARG	3.2
1	B	42	LYS	3.2
1	A	157	ARG	3.2
1	A	550	VAL	3.1
1	A	149	TYR	3.1
1	B	545	ARG	3.1
1	A	26	ASN	2.8
1	B	26	ASN	2.7
1	B	160	LYS	2.7
1	A	582	LYS	2.7
1	A	163	LEU	2.7
1	A	546	ASN	2.6
1	B	44	LYS	2.6
1	B	572	TYR	2.6
1	B	51	LYS	2.5
1	B	279	LYS	2.5
1	A	44	LYS	2.5
1	A	571	LYS	2.4
1	B	579	GLU	2.4
1	A	567	LEU	2.4
1	A	216	GLU	2.3
1	B	149	TYR	2.3
1	A	537	GLU	2.3
1	A	502	TYR	2.2
1	A	559	CYS	2.2
1	A	42	LYS	2.1
1	B	567	LEU	2.1
1	B	161	SER	2.0
1	A	162	ASN	2.0
1	A	581	GLY	2.0
1	B	54	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.