



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2020 – 04:48 PM BST

PDB ID : 6IS7
Title : Structure of 9N-I DNA polymerase incorporation with dA in the active site
Authors : Linwu, S.W.; Maestre-Reyna, M.; Tsai, M.D.; Tu, Y.H.; Chang, W.H.
Deposited on : 2018-11-15
Resolution : 2.80 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.13
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13

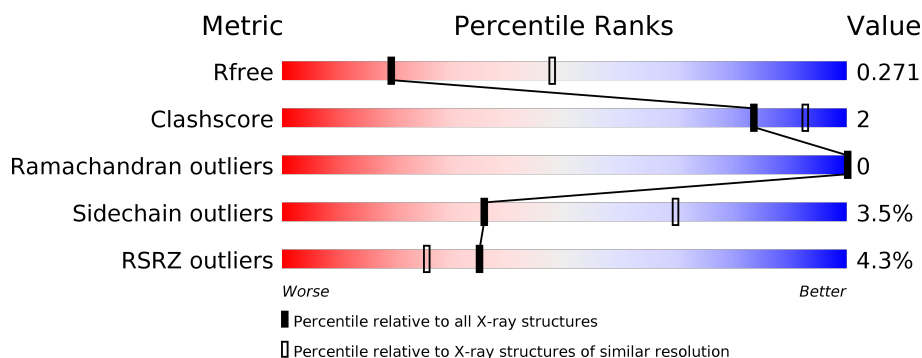
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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	783	<div> <div>5%</div> <div> <div></div> <div>90%</div> <div>6%</div> <div>••</div> </div> </div>
1	B	783	<div> <div>3%</div> <div> <div></div> <div>90%</div> <div>6%</div> <div>••</div> </div> </div>
2	C	15	<div> <div></div> <div> <div>80%</div> <div>20%</div> </div> </div>
2	E	15	<div> <div></div> <div> <div>60%</div> <div>40%</div> </div> </div>
2	K	15	<div> <div>20%</div> <div>80%</div> </div>
2	L	15	<div> <div>27%</div> <div>73%</div> </div>

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Mol	Chain	Length	Quality of chain
3	D	18	 6% 94% 6%
3	F	18	 6% 94% 6%
3	I	18	 11% 11% 78%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13390 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 polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	757	Total	C	N	O	S	0	0	0
			5883	3812	997	1059	15			
1	B	756	Total	C	N	O	S	0	0	0
			5723	3711	955	1042	15			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	141	ALA	ASP	engineered mutation	UNP Q56366
A	143	ALA	GLU	engineered mutation	UNP Q56366
A	485	LEU	ALA	engineered mutation	UNP Q56366
A	776	LEU	-	expression tag	UNP Q56366
A	777	GLU	-	expression tag	UNP Q56366
A	778	HIS	-	expression tag	UNP Q56366
A	779	HIS	-	expression tag	UNP Q56366
A	780	HIS	-	expression tag	UNP Q56366
A	781	HIS	-	expression tag	UNP Q56366
A	782	HIS	-	expression tag	UNP Q56366
A	783	HIS	-	expression tag	UNP Q56366
B	141	ALA	ASP	engineered mutation	UNP Q56366
B	143	ALA	GLU	engineered mutation	UNP Q56366
B	485	LEU	ALA	engineered mutation	UNP Q56366
B	776	LEU	-	expression tag	UNP Q56366
B	777	GLU	-	expression tag	UNP Q56366
B	778	HIS	-	expression tag	UNP Q56366
B	779	HIS	-	expression tag	UNP Q56366
B	780	HIS	-	expression tag	UNP Q56366
B	781	HIS	-	expression tag	UNP Q56366
B	782	HIS	-	expression tag	UNP Q56366
B	783	HIS	-	expression tag	UNP Q56366

- Molecule 2 is a DNA chain called DNA (5'-D(P*GP*CP*GP*GP*AP*CP*TP*GP*CP*TP*TP*AP*CP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	15	Total	C	N	O	P	0	0	0
			303	145	56	88	14			
2	E	15	Total	C	N	O	P	0	0	0
			302	145	56	87	14			
2	K	3	Total	C	N	O	P	0	0	0
			59	29	7	20	3			
2	L	4	Total	C	N	O	P	0	0	0
			81	39	12	26	4			

- Molecule 3 is a DNA chain called DNA (5'-D(*AP*CP*TP*GP*GP*TP*AP*AP*GP*CP*AP*GP*TP*CP*CP*GP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	18	Total	C	N	O	P	0	0	0
			368	175	71	105	17			
3	F	18	Total	C	N	O	P	0	0	0
			368	175	71	105	17			
3	I	4	Total	C	N	O	P	0	0	0
			83	39	18	22	4			

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	5	Total	Ca	0	0
			5	5		
4	A	2	Total	Ca	0	0
			2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	93	Total	O	0	0
			93	93		
5	B	76	Total	O	0	1
			77	77		
5	C	8	Total	O	0	0
			8	8		
5	D	8	Total	O	0	0
			8	8		
5	E	7	Total	O	0	0
			7	7		
5	F	8	Total	O	0	0
			8	8		

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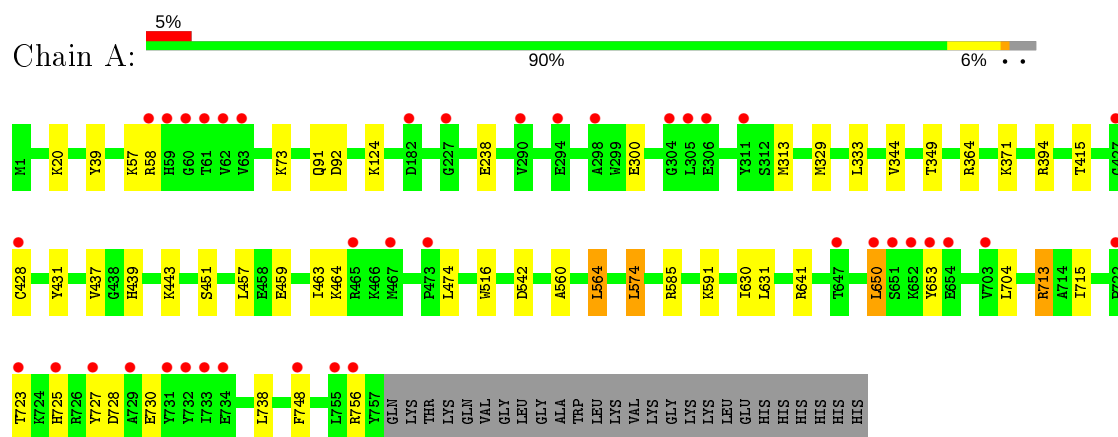
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	I	3	Total 3	O 3	0	0
5	K	3	Total 3	O 3	0	0
5	L	6	Total 6	O 6	0	0

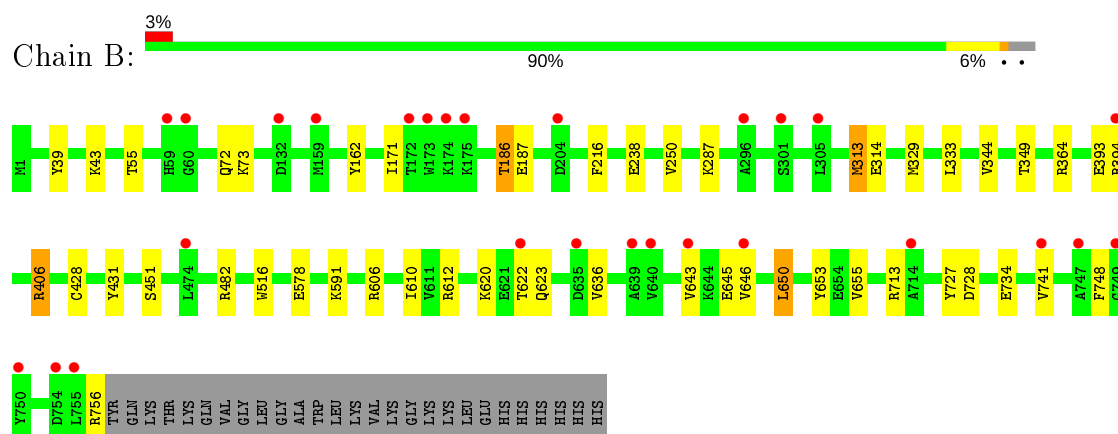
3 Residue-property plots [i](#)

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.

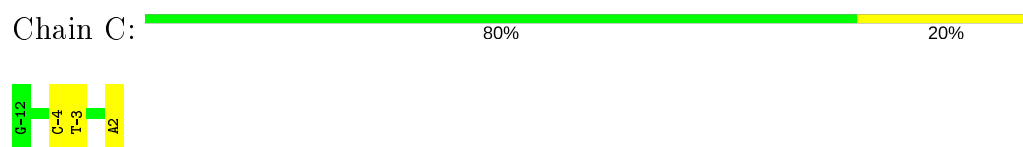
• Molecule 1: DNA polymerase



• Molecule 1: DNA polymerase



• Molecule 2: DNA (5'-D(P*GP*CP*GP*GP*AP*CP*TP*GP*CP*TP*TP*AP*CP*CP*A)-3')



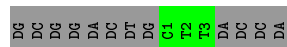
- Molecule 2: DNA (5'-D(P*GP*CP*GP*GP*AP*CP*TP*GP*CP*TP*TP*AP*CP*CP*A)-3')

Chain E: 



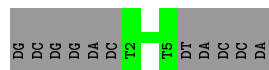
- Molecule 2: DNA (5'-D(P*GP*CP*GP*GP*AP*CP*TP*GP*CP*TP*TP*AP*CP*CP*A)-3')

Chain K: 



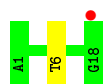
- Molecule 2: DNA (5'-D(P*GP*CP*GP*GP*AP*CP*TP*GP*CP*TP*TP*AP*CP*CP*A)-3')

Chain L: 



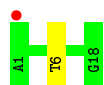
- Molecule 3: DNA (5'-D(*AP*CP*TP*GP*GP*TP*AP*AP*GP*CP*AP*GP*TP*CP*CP*GP*CP*G)-3')

Chain D: 



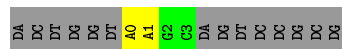
- Molecule 3: DNA (5'-D(*AP*CP*TP*GP*GP*TP*AP*AP*GP*CP*AP*GP*TP*CP*CP*GP*CP*G)-3')

Chain F: 



- Molecule 3: DNA (5'-D(*AP*CP*TP*GP*GP*TP*AP*AP*GP*CP*AP*GP*TP*CP*CP*GP*CP*G)-3')

Chain I: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	92.95Å 106.77Å 255.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.29 – 2.80 47.26 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.4 (46.29-2.80) 99.5 (47.26-2.80)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.225 , 0.276 0.225 , 0.271	Depositor DCC
R_{free} test set	3139 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	58.1	Xtriage
Anisotropy	0.272	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 66.2	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	0.91	EDS
Total number of atoms	13390	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% 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: CA

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	0.64	0/6022	0.73	0/8171
1	B	0.65	0/5854	0.74	0/7959
2	C	0.47	0/339	0.84	0/521
2	E	0.48	0/338	0.88	2/520 (0.4%)
2	K	0.20	0/64	0.64	0/96
2	L	0.20	0/89	0.68	0/135
3	D	0.40	0/413	0.77	0/636
3	F	0.43	0/413	0.76	0/636
3	I	0.23	0/93	0.84	0/141
All	All	0.62	0/13625	0.74	2/18815 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	0	DC	O4'-C4'-C3'	-5.88	102.15	104.50
2	E	0	DC	C1'-O4'-C4'	-5.21	104.89	110.10

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	5883	0	5613	27	0
1	B	5723	0	5342	26	0
2	C	303	0	170	3	0
2	E	302	0	167	4	0
2	K	59	0	36	0	0
2	L	81	0	47	0	0
3	D	368	0	203	1	0
3	F	368	0	203	1	0
3	I	83	0	45	1	0
4	A	2	0	0	0	0
4	B	5	0	0	0	0
5	A	93	0	0	2	0
5	B	77	0	0	0	0
5	C	8	0	0	0	0
5	D	8	0	0	0	0
5	E	7	0	0	0	0
5	F	8	0	0	0	0
5	I	3	0	0	0	0
5	K	3	0	0	0	0
5	L	6	0	0	0	0
All	All	13390	0	11826	59	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 2.

All (59) 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:610:ILE:HG13	1:B:623:GLN:HE21	1.44	0.82
1:B:636:VAL:HG21	1:B:748:PHE:CD2	2.20	0.76
1:B:643:VAL:HG21	1:B:741:VAL:HG21	1.70	0.73
1:A:730:GLU:OE2	1:A:756:ARG:NH2	2.21	0.73
1:B:364:ARG:NH2	1:B:451:SER:OG	2.22	0.73
1:B:713:ARG:NH1	1:B:728:ASP:OD2	2.24	0.70
1:A:713:ARG:NH1	1:A:728:ASP:OD2	2.25	0.69
1:A:437:VAL:CG1	1:A:439:HIS:CD2	2.75	0.69
1:A:437:VAL:HG12	1:A:439:HIS:CD2	2.31	0.66
1:A:58:ARG:NH1	1:A:92:ASP:OD1	2.31	0.64
1:A:459:GLU:O	1:A:463:ILE:HG13	2.00	0.62
1:A:574:LEU:H	1:A:574:LEU:HD23	1.66	0.61
1:A:437:VAL:HG11	1:A:439:HIS:CD2	2.37	0.59
1:A:574:LEU:HD23	1:A:574:LEU:N	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:591:LYS:HD3	3:D:6:DT:H5"	1.89	0.55
2:C:-4:DC:H2"	2:C:-3:DT:H5"	1.90	0.54
1:B:643:VAL:CG2	1:B:741:VAL:HG21	2.37	0.54
1:B:636:VAL:HG21	1:B:748:PHE:HD2	1.69	0.54
1:B:734:GLU:OE2	1:B:756:ARG:NH1	2.40	0.54
2:E:-4:DC:H2"	2:E:-3:DT:H5"	1.90	0.52
1:A:333:LEU:HD23	1:A:344:VAL:HG13	1.93	0.51
1:B:622:THR:HG21	1:B:646:VAL:HG11	1.92	0.51
1:B:333:LEU:HD23	1:B:344:VAL:HG13	1.93	0.51
1:B:650:LEU:HD13	1:B:655:VAL:HG21	1.92	0.51
1:B:591:LYS:HD3	3:F:6:DT:H5"	1.95	0.49
1:A:39:TYR:CZ	1:A:73:LYS:HE3	2.48	0.49
1:A:364:ARG:NH2	1:A:451:SER:OG	2.43	0.48
1:A:738:LEU:HD11	1:A:756:ARG:HG3	1.95	0.48
1:A:58:ARG:HD2	1:A:91:GLN:HB3	1.96	0.48
1:A:630:ILE:HD13	1:A:748:PHE:HE2	1.79	0.47
1:A:371:LYS:NZ	5:A:901:HOH:O	2.47	0.47
2:E:-11:DC:H2'	2:E:-10:DG:C8	2.49	0.47
1:B:428:CYS:HB2	1:B:431:TYR:CE1	2.51	0.46
1:A:428:CYS:HB2	1:A:431:TYR:CE1	2.50	0.46
1:A:653:TYR:HA	1:A:727:TYR:OH	2.15	0.46
1:B:186:THR:OG1	1:B:187:GLU:N	2.48	0.46
1:B:39:TYR:CZ	1:B:73:LYS:HE3	2.52	0.45
3:I:0:DA:H2"	3:I:1:DA:H5"	1.99	0.45
1:B:313:MET:HG3	1:B:314:GLU:N	2.33	0.44
1:A:560:ALA:O	1:A:564:LEU:HD22	2.18	0.44
1:B:653:TYR:HA	1:B:727:TYR:OH	2.17	0.44
1:B:406:ARG:NH2	1:B:578:GLU:OE2	2.50	0.44
1:B:643:VAL:O	1:B:646:VAL:HG12	2.19	0.43
1:A:704:LEU:HD22	1:A:725:HIS:HB3	2.01	0.43
1:B:612:ARG:O	1:B:620:LYS:HD3	2.20	0.42
1:A:542:ASP:OD1	2:C:2:DA:O3'	2.37	0.42
1:A:437:VAL:CG1	1:A:439:HIS:NE2	2.82	0.42
1:B:636:VAL:CG2	1:B:748:PHE:CD2	2.99	0.42
1:B:606:ARG:HD2	2:E:2:DA:OP2	2.20	0.42
1:A:415:THR:O	1:A:443:LYS:NZ	2.53	0.42
1:B:610:ILE:HG13	1:B:623:GLN:NE2	2.24	0.42
1:A:631:LEU:HD23	1:A:631:LEU:HA	1.90	0.41
1:B:162:TYR:CE2	1:B:171:ILE:HD11	2.55	0.41
1:B:636:VAL:HG21	1:B:748:PHE:CE2	2.54	0.41
1:A:650:LEU:HA	1:A:650:LEU:HD12	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:PHE:HB2	1:B:250:VAL:HG21	2.03	0.40
2:E:-4:DC:H2''	2:E:-3:DT:C5'	2.52	0.40
2:C:-4:DC:H2''	2:C:-3:DT:C5'	2.51	0.40
1:A:725:HIS:HA	5:A:955:HOH:O	2.20	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	755/783 (96%)	736 (98%)	19 (2%)	0	100	100
1	B	754/783 (96%)	735 (98%)	19 (2%)	0	100	100
All	All	1509/1566 (96%)	1471 (98%)	38 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	550/675 (82%)	529 (96%)	21 (4%)	33	67
1	B	515/675 (76%)	499 (97%)	16 (3%)	40	74
All	All	1065/1350 (79%)	1028 (96%)	37 (4%)	36	70

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LYS
1	A	57	LYS
1	A	124	LYS
1	A	238	GLU
1	A	300	GLU
1	A	313	MET
1	A	329	MET
1	A	349	THR
1	A	394	ARG
1	A	457	LEU
1	A	464	LYS
1	A	474	LEU
1	A	516	TRP
1	A	564	LEU
1	A	574	LEU
1	A	585	ARG
1	A	641	ARG
1	A	650	LEU
1	A	713	ARG
1	A	715	ILE
1	A	723	THR
1	B	43	LYS
1	B	55	THR
1	B	72	GLN
1	B	186	THR
1	B	238	GLU
1	B	287	LYS
1	B	313	MET
1	B	329	MET
1	B	349	THR
1	B	393	GLU
1	B	394	ARG
1	B	406	ARG
1	B	482	ARG
1	B	516	TRP
1	B	645	GLU
1	B	650	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	297	GLN
1	A	351	ASN
1	A	461	GLN
1	A	679	HIS
1	B	351	ASN
1	B	623	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 7 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	757/783 (96%)	0.25	39 (5%) 27 18	55, 89, 135, 161	3 (0%)
1	B	756/783 (96%)	0.20	27 (3%) 42 32	55, 89, 162, 232	1 (0%)
2	C	15/15 (100%)	0.03	0 100 100	60, 71, 88, 88	0
2	E	15/15 (100%)	-0.13	0 100 100	51, 62, 75, 78	0
2	K	3/15 (20%)	1.43	0 100 100	117, 117, 118, 164	2 (66%)
2	L	4/15 (26%)	0.73	0 100 100	112, 113, 114, 156	3 (75%)
3	D	18/18 (100%)	0.56	1 (5%) 24 16	56, 71, 119, 142	1 (5%)
3	F	18/18 (100%)	0.19	1 (5%) 24 16	57, 72, 106, 121	0
3	I	4/18 (22%)	1.03	0 100 100	116, 154, 156, 158	1 (25%)
All	All	1590/1680 (94%)	0.23	68 (4%) 35 25	51, 88, 146, 232	11 (0%)

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	18	DG	4.9
1	B	749	GLY	4.9
1	A	428	CYS	4.4
1	A	650	LEU	4.2
1	A	727	TYR	4.2
1	B	305	LEU	4.0
1	B	755	LEU	3.9
1	B	174	LYS	3.9
1	A	427	GLY	3.7
1	B	172	THR	3.7
1	A	653	TYR	3.6
1	A	651	SER	3.6
1	B	173	TRP	3.3
1	B	474	LEU	3.1
1	B	754	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	714	ALA	3.1
1	B	643	VAL	3.1
1	A	755	LEU	3.0
1	A	703	VAL	3.0
1	A	722	PRO	2.9
1	B	639	ALA	2.9
1	A	59	HIS	2.9
1	B	59	HIS	2.9
1	A	473	PRO	2.8
1	A	304	GLY	2.8
1	A	60	GLY	2.8
1	A	62	VAL	2.8
1	B	159	MET	2.8
1	A	733	ILE	2.8
1	A	465	ARG	2.7
1	A	227	GLY	2.7
1	A	756	ARG	2.7
1	A	63	VAL	2.6
1	A	654	GLU	2.6
1	A	647	THR	2.6
1	A	305	LEU	2.5
1	B	60	GLY	2.5
1	A	58	ARG	2.5
1	B	640	VAL	2.5
1	A	652	LYS	2.5
1	A	731	TYR	2.5
1	B	747	ALA	2.5
1	B	622	THR	2.4
1	A	723	THR	2.4
1	A	748	PHE	2.4
1	A	61	THR	2.4
1	B	132	ASP	2.3
1	A	306	GLU	2.3
1	B	296	ALA	2.3
1	A	290	VAL	2.2
1	A	734	GLU	2.2
3	F	1	DA	2.2
1	B	204	ASP	2.2
1	B	301	SER	2.2
1	A	729	ALA	2.2
1	B	394	ARG	2.2
1	B	175	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	725	HIS	2.2
1	B	750	TYR	2.2
1	A	298	ALA	2.2
1	B	646	VAL	2.2
1	B	635	ASP	2.2
1	A	467	MET	2.1
1	B	741	VAL	2.0
1	A	294	GLU	2.0
1	A	311	TYR	2.0
1	A	732	TYR	2.0
1	A	182	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 monosaccharides 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(Å ²)	Q<0.9
4	CA	B	805	1/1	0.81	0.33	102,102,102,102	0
4	CA	B	801	1/1	0.84	0.30	109,109,109,109	0
4	CA	A	802	1/1	0.84	0.16	107,107,107,107	0
4	CA	A	801	1/1	0.89	0.21	109,109,109,109	0
4	CA	B	804	1/1	0.91	0.10	117,117,117,117	0
4	CA	B	802	1/1	0.91	0.13	126,126,126,126	0
4	CA	B	803	1/1	0.93	0.09	77,77,77,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.