



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 6, 2021 – 11:02 am BST

PDB ID : 6YL2  
Title : Structural and DNA binding studies of the transcriptional repressor Rv2506 (BkaR) from Mycobacterium tuberculosis supports a role in L-Leucine catabolism  
Authors : Keep, N.H.; Pritchard, J.E.; Sula, A.; Cole, A.R.; Kendall, S.L.  
Deposited on : 2020-04-06  
Resolution : 3.15 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](mailto: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.

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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 : 2.18  
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) : 2.18

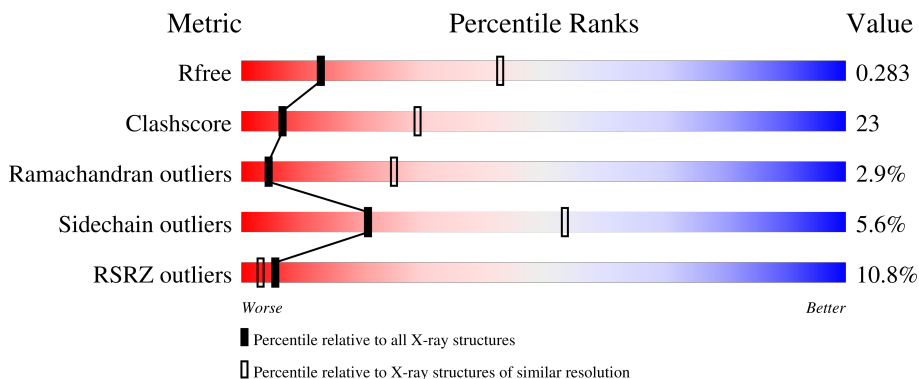
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.15 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	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)
Ramachandran outliers	138981	1770 (3.20-3.12)
Sidechain outliers	138945	1769 (3.20-3.12)
RSRZ outliers	127900	1616 (3.20-3.12)

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  $\geq 3$ , 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	196	
1	B	196	
2	G	20	
3	H	20	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3740 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 Probable transcriptional regulatory protein (Probably TetR-family).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	192	Total	C	N	O	S	0	0	0
			1463	918	279	262	4			
1	A	192	Total	C	N	O	S	0	0	0
			1463	918	279	262	4			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	20	Total	C	N	O	P	0	0	0
			408	196	77	116	19			

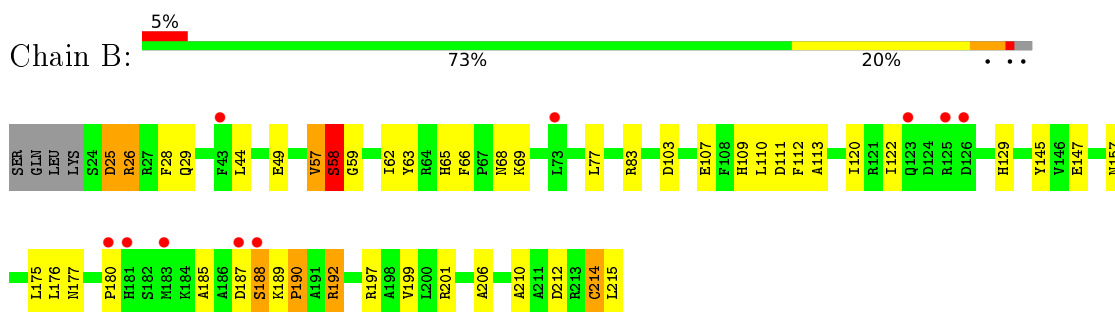
- Molecule 3 is a DNA chain called DNA (5'-D(\*CP\*GP\*GP\*TP\*TP\*AP\*AP\*TP\*CP\*GP\*TP\*CP\*AP\*TP\*TP\*AP\*AP\*CP\*GP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	20	Total	C	N	O	P	0	0	0
			406	196	71	120	19			

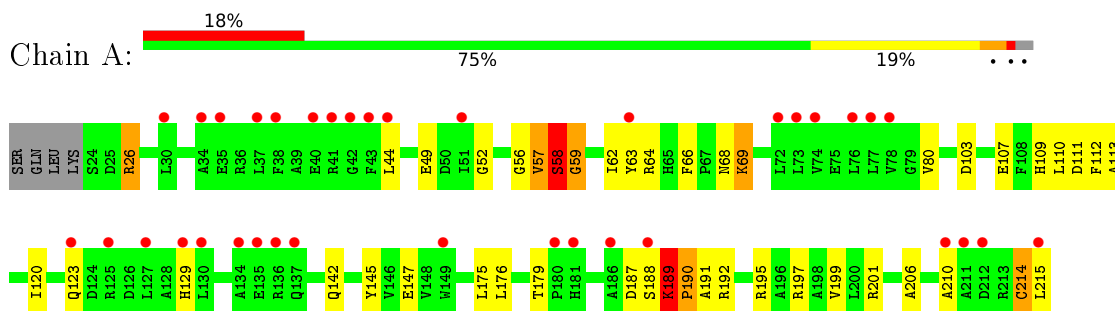
### 3 Residue-property plots [i](#)

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.

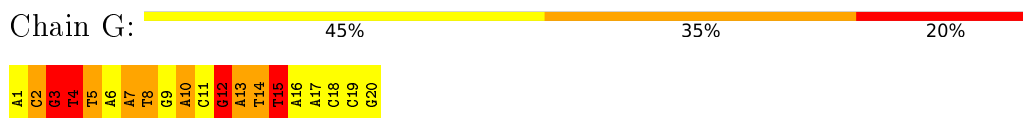
- Molecule 1: Probable transcriptional regulatory protein (Probably TetR-family)



- Molecule 1: Probable transcriptional regulatory protein (Probably TetR-family)



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



- Molecule 3: DNA (5'-D(\*CP\*GP\*GP\*TP\*TP\*AP\*AP\*TP\*CP\*GP\*TP\*CP\*AP\*TP\*TP\*AP\*AP\*CP\*GP\*T)-3')



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.41Å 66.00Å 179.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	62.03 – 3.15 61.95 – 3.15	Depositor EDS
% Data completeness (in resolution range)	97.5 (62.03-3.15) 97.6 (61.95-3.15)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.83 (at 3.13Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.267 , 0.284 0.267 , 0.283	Depositor DCC
$R_{free}$ test set	618 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	90.5	Xtrriage
Anisotropy	0.974	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 92.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3740	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	157.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.30 % 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 1.1679e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

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	0.69	0/1484	0.91	0/2007
1	B	0.66	0/1484	0.90	0/2007
2	G	1.96	13/458 (2.8%)	1.65	13/705 (1.8%)
3	H	1.85	15/454 (3.3%)	1.78	12/699 (1.7%)
All	All	1.10	28/3880 (0.7%)	1.17	25/5418 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
2	G	0	5
3	H	0	3
All	All	0	10

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	3	DG	C8-N7	15.05	1.40	1.30
2	G	15	DT	C5-C7	12.58	1.57	1.50
3	H	9	DC	C4-C5	10.75	1.51	1.43
2	G	3	DG	C5-C4	10.69	1.45	1.38
3	H	15	DT	C5-C7	9.83	1.55	1.50
3	H	9	DC	N1-C2	8.75	1.49	1.40
3	H	13	DA	N9-C4	8.59	1.43	1.37
2	G	4	DT	N1-C2	8.49	1.44	1.38
3	H	12	DC	C4-C5	7.80	1.49	1.43
3	H	2	DG	N9-C8	7.66	1.43	1.37
2	G	8	DT	N1-C2	6.98	1.43	1.38
3	H	15	DT	N1-C2	6.44	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	11	DT	C5-C7	6.32	1.53	1.50
2	G	11	DC	C4-C5	6.12	1.47	1.43
3	H	2	DG	C5-C4	5.93	1.42	1.38
2	G	10	DA	N9-C4	5.80	1.41	1.37
2	G	4	DT	C4-C5	5.67	1.50	1.45
2	G	5	DT	O3'-P	-5.64	1.54	1.61
3	H	8	DT	N1-C2	5.55	1.42	1.38
3	H	2	DG	N1-C2	5.35	1.42	1.37
2	G	12	DG	N1-C2	-5.33	1.33	1.37
2	G	15	DT	C4-C5	5.29	1.49	1.45
2	G	11	DC	N1-C2	5.28	1.45	1.40
3	H	2	DG	O4'-C1'	5.16	1.48	1.42
3	H	6	DA	N9-C4	5.13	1.41	1.37
3	H	19	DG	O3'-P	5.04	1.67	1.61
2	G	5	DT	N1-C2	5.04	1.42	1.38
3	H	8	DT	C4-C5	5.03	1.49	1.45

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	15	DT	O4'-C1'-N1	-16.91	96.16	108.00
2	G	15	DT	C6-C5-C7	-7.70	118.28	122.90
3	H	15	DT	C6-C5-C7	-7.41	118.46	122.90
3	H	2	DG	C8-N9-C4	-7.33	103.47	106.40
3	H	13	DA	O4'-C1'-N9	-7.08	103.05	108.00
3	H	12	DC	N3-C4-C5	-6.98	119.11	121.90
2	G	15	DT	C4'-C3'-O3'	-6.42	93.64	109.70
3	H	2	DG	C2-N3-C4	-6.13	108.83	111.90
2	G	15	DT	C4-C5-C7	6.04	122.63	119.00
2	G	15	DT	C5-C6-N1	-5.96	120.12	123.70
3	H	3	DG	O5'-P-OP1	-5.78	100.50	105.70
2	G	7	DA	O5'-P-OP2	-5.75	100.52	105.70
3	H	2	DG	N3-C4-N9	-5.71	122.57	126.00
2	G	12	DG	N3-C4-C5	-5.67	125.77	128.60
3	H	3	DG	N9-C1'-C2'	-5.61	101.95	112.60
3	H	15	DT	C5-C6-N1	-5.55	120.37	123.70
2	G	8	DT	C6-N1-C2	-5.54	118.53	121.30
2	G	14	DT	P-O5'-C5'	-5.49	112.11	120.90
2	G	13	DA	O4'-C1'-N9	5.45	111.81	108.00
3	H	15	DT	N1-C1'-C2'	5.42	122.90	112.60
2	G	3	DG	C8-N9-C4	-5.38	104.25	106.40
2	G	8	DT	N3-C4-C5	-5.15	112.11	115.20

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Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
2	G	3	DG	N9-C4-C5	5.12	107.45	105.40
3	H	11	DT	C6-C5-C7	-5.09	119.84	122.90
2	G	8	DT	C4-C5-C6	5.04	121.02	118.00

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	58	SER	Peptide
1	B	58	SER	Peptide
2	G	12	DG	Sidechain
2	G	15	DT	Sidechain
2	G	2	DC	Sidechain
2	G	3	DG	Sidechain
2	G	4	DT	Sidechain
3	H	12	DC	Sidechain
3	H	17	DA	Sidechain
3	H	2	DG	Sidechain

## 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	1463	0	1503	47	0
1	B	1463	0	1503	43	0
2	G	408	0	227	49	0
3	H	406	0	229	41	0
All	All	3740	0	3462	166	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 23.

All (166) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance ( $\text{\AA}$ )	Clash overlap ( $\text{\AA}$ )
1:A:44:LEU:HD11	1:A:129:HIS:CE1	1.97	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:GLY:O	1:A:57:VAL:HG22	1.66	0.95
1:A:188:SER:HA	1:A:192:ARG:HB2	1.60	0.81
2:G:14:DT:C6	2:G:15:DT:H71	2.15	0.80
2:G:14:DT:C5	2:G:15:DT:H73	2.17	0.80
3:H:2:DG:H2''	3:H:3:DG:H5'	1.63	0.79
1:B:26:ARG:HB3	1:B:65:HIS:CE1	2.18	0.78
2:G:4:DT:H2'	2:G:5:DT:H71	1.64	0.78
2:G:14:DT:C6	2:G:15:DT:C7	2.66	0.78
1:A:59:GLY:O	1:A:62:ILE:CD1	2.35	0.74
2:G:14:DT:C2	2:G:15:DT:C5	2.76	0.73
2:G:6:DA:H2''	2:G:7:DA:O5'	1.89	0.73
3:H:2:DG:H2''	3:H:3:DG:C5'	2.18	0.72
1:A:44:LEU:CD1	1:A:129:HIS:CE1	2.73	0.72
2:G:1:DA:H3'	2:G:2:DC:H5'	1.70	0.72
3:H:7:DA:OP1	3:H:7:DA:H4'	1.91	0.71
1:A:64:ARG:HD2	3:H:2:DG:C8	2.25	0.71
1:B:59:GLY:O	1:B:62:ILE:CD1	2.40	0.70
1:A:44:LEU:HD11	1:A:129:HIS:NE2	2.07	0.69
1:B:187:ASP:OD2	1:B:192:ARG:HG2	1.92	0.69
1:A:69:LYS:NZ	2:G:13:DA:OP1	2.26	0.68
2:G:7:DA:C2	2:G:8:DT:C2	2.81	0.67
2:G:14:DT:C2	2:G:15:DT:C6	2.82	0.67
3:H:7:DA:H2'	3:H:8:DT:C7	2.25	0.67
2:G:1:DA:H2'	2:G:1:DA:N3	2.09	0.67
3:H:12:DC:H2''	3:H:13:DA:C8	2.30	0.67
3:H:15:DT:H1'	3:H:16:DA:H5''	1.77	0.67
1:B:110:LEU:HD13	1:B:176:LEU:HD22	1.79	0.65
1:A:110:LEU:HD13	1:A:176:LEU:HD22	1.78	0.65
1:A:57:VAL:HG23	1:A:58:SER:H	1.61	0.63
2:G:1:DA:C2'	2:G:2:DC:H5''	2.28	0.63
2:G:19:DC:H2''	2:G:20:DG:C8	2.33	0.63
3:H:11:DT:H2''	3:H:12:DC:O5'	1.99	0.63
2:G:14:DT:N1	2:G:15:DT:H71	2.15	0.61
2:G:1:DA:C2	2:G:2:DC:C5	2.89	0.61
3:H:18:DC:H2''	3:H:19:DG:O5'	2.00	0.61
1:B:26:ARG:HB3	1:B:65:HIS:HE1	1.61	0.61
3:H:6:DA:C4	3:H:7:DA:C8	2.89	0.60
3:H:6:DA:C4	3:H:7:DA:N7	2.70	0.60
1:B:210:ALA:CB	1:A:206:ALA:HB2	2.32	0.60
2:G:1:DA:N1	2:G:2:DC:C4	2.70	0.59
1:B:214:CYS:HB3	1:B:215:LEU:HD13	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:1:DC:C2'	3:H:2:DG:H5''	2.32	0.59
1:B:59:GLY:O	1:B:62:ILE:HD12	2.03	0.58
1:B:210:ALA:HB2	1:A:206:ALA:CB	2.34	0.58
1:B:206:ALA:HB2	1:A:210:ALA:CB	2.34	0.58
2:G:1:DA:H3'	2:G:2:DC:C5'	2.34	0.57
1:B:210:ALA:CB	1:A:206:ALA:CB	2.82	0.57
1:B:214:CYS:O	1:B:215:LEU:HD12	2.05	0.57
2:G:1:DA:C2	2:G:2:DC:C4	2.93	0.56
1:A:214:CYS:C	1:A:215:LEU:HD12	2.26	0.56
1:B:206:ALA:CB	1:A:210:ALA:CB	2.83	0.56
2:G:17:DA:H1'	2:G:18:DC:H5''	1.88	0.55
1:A:80:VAL:CG2	1:A:112:PHE:HZ	2.18	0.55
1:A:113:ALA:HA	1:A:120:ILE:CD1	2.37	0.55
1:A:59:GLY:O	1:A:62:ILE:HD12	2.06	0.55
1:A:214:CYS:O	1:A:215:LEU:HD12	2.05	0.55
1:A:63:TYR:HA	1:A:66:PHE:O	2.07	0.55
1:B:206:ALA:CB	1:A:210:ALA:HB2	2.36	0.54
1:A:111:ASP:OD1	1:A:197:ARG:NH1	2.40	0.54
1:B:111:ASP:OD1	1:B:197:ARG:NH1	2.40	0.54
3:H:6:DA:N3	3:H:7:DA:C8	2.76	0.53
1:B:63:TYR:HA	1:B:66:PHE:O	2.08	0.53
1:B:189:LYS:HB3	1:B:190:PRO:HD2	1.90	0.53
1:A:103:ASP:O	1:A:107:GLU:HG2	2.07	0.53
2:G:14:DT:C5	2:G:15:DT:C7	2.85	0.53
2:G:10:DA:H8	2:G:10:DA:H5'	1.74	0.53
1:B:187:ASP:HA	1:B:188:SER:CB	2.38	0.53
1:B:26:ARG:NH1	1:B:57:VAL:HB	2.24	0.52
3:H:3:DG:H2''	3:H:4:DT:O5'	2.09	0.52
3:H:7:DA:OP1	3:H:7:DA:C4'	2.57	0.52
2:G:1:DA:C2'	2:G:2:DC:C5'	2.87	0.52
3:H:13:DA:C8	3:H:14:DT:C7	2.93	0.52
1:B:214:CYS:C	1:B:215:LEU:HD12	2.30	0.52
2:G:1:DA:H2''	2:G:2:DC:H5''	1.91	0.52
1:A:26:ARG:HG3	3:H:3:DG:OP1	2.10	0.52
1:A:109:HIS:HD2	1:A:145:TYR:OH	1.93	0.52
3:H:7:DA:H2'	3:H:8:DT:H73	1.91	0.52
3:H:6:DA:H2''	3:H:7:DA:O5'	2.09	0.52
2:G:14:DT:C6	2:G:15:DT:H73	2.40	0.51
3:H:12:DC:H4'	3:H:12:DC:OP1	2.11	0.51
1:A:52:GLY:HA3	1:A:58:SER:HA	1.91	0.51
1:B:210:ALA:HB2	1:A:206:ALA:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:13:DA:C5	3:H:14:DT:C4	2.98	0.51
2:G:14:DT:N3	2:G:15:DT:C5	2.79	0.51
2:G:1:DA:C3'	2:G:2:DC:C5'	2.88	0.51
2:G:3:DG:C8	2:G:4:DT:H72	2.46	0.51
1:B:109:HIS:HD2	1:B:145:TYR:OH	1.93	0.50
1:B:103:ASP:O	1:B:107:GLU:HG2	2.11	0.50
3:H:2:DG:C2'	3:H:3:DG:H5''	2.41	0.50
1:B:113:ALA:HA	1:B:120:ILE:CD1	2.41	0.50
1:B:199:VAL:HG22	1:A:215:LEU:HD11	1.92	0.50
1:A:214:CYS:HB3	1:A:215:LEU:HD13	1.92	0.50
2:G:4:DT:H2'	2:G:5:DT:C7	2.37	0.50
2:G:6:DA:C2'	2:G:7:DA:O5'	2.58	0.49
1:A:57:VAL:HG23	1:A:58:SER:N	2.27	0.49
1:B:187:ASP:HA	1:B:188:SER:HB3	1.94	0.49
1:B:206:ALA:HB2	1:A:210:ALA:HB2	1.95	0.49
1:B:26:ARG:HG2	2:G:3:DG:OP1	2.12	0.49
3:H:1:DC:C3'	3:H:2:DG:H5''	2.43	0.48
1:B:215:LEU:HD11	1:A:199:VAL:HG22	1.94	0.48
2:G:12:DG:C2	2:G:13:DA:C6	3.01	0.48
1:A:191:ALA:O	1:A:195:ARG:CB	2.61	0.48
2:G:7:DA:H2'	2:G:8:DT:C7	2.44	0.47
1:A:188:SER:HA	1:A:192:ARG:CB	2.36	0.47
3:H:2:DG:H2''	3:H:3:DG:H5''	1.96	0.47
3:H:2:DG:C2'	3:H:3:DG:C5'	2.91	0.47
2:G:8:DT:H2''	2:G:9:DG:C8	2.49	0.47
1:A:80:VAL:HG23	1:A:112:PHE:HZ	1.79	0.47
1:A:147:GLU:HA	1:A:147:GLU:OE1	2.15	0.47
3:H:12:DC:C2'	3:H:13:DA:C8	2.98	0.47
3:H:2:DG:H1'	3:H:3:DG:H5''	1.97	0.47
3:H:5:DT:H2''	3:H:6:DA:OP2	2.15	0.47
2:G:1:DA:C2	2:G:2:DC:C6	3.03	0.46
3:H:4:DT:H2'	3:H:5:DT:H71	1.96	0.46
2:G:4:DT:H2''	2:G:5:DT:O5'	2.15	0.46
2:G:1:DA:C6	2:G:2:DC:C4	3.03	0.46
1:A:189:LYS:CG	1:A:190:PRO:HD3	2.47	0.45
2:G:9:DG:H1'	2:G:10:DA:C8	2.51	0.45
2:G:15:DT:C2	2:G:16:DA:N7	2.84	0.45
1:B:25:ASP:HB2	1:B:28:PHE:HB3	1.97	0.45
2:G:10:DA:H5'	2:G:10:DA:C8	2.51	0.45
1:A:59:GLY:O	1:A:62:ILE:HD13	2.17	0.44
1:B:44:LEU:HD11	1:B:129:HIS:NE2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:17:DA:H2''	3:H:18:DC:H5''	1.98	0.44
1:B:107:GLU:OE2	1:B:201:ARG:HD3	2.18	0.44
1:B:147:GLU:OE1	1:B:147:GLU:HA	2.17	0.44
1:B:189:LYS:HB3	1:B:190:PRO:CD	2.46	0.44
2:G:1:DA:C4	2:G:2:DC:C5	3.06	0.44
3:H:16:DA:H1'	3:H:17:DA:H5'	1.99	0.44
1:B:187:ASP:CG	1:B:192:ARG:HG2	2.38	0.43
1:A:57:VAL:C	1:A:58:SER:OG	2.56	0.43
1:A:107:GLU:OE2	1:A:201:ARG:HD3	2.18	0.43
1:A:189:LYS:HG3	1:A:190:PRO:HD3	2.00	0.43
2:G:14:DT:H2'	2:G:15:DT:H71	2.00	0.43
2:G:7:DA:H2''	2:G:8:DT:H5'	1.99	0.43
3:H:6:DA:C2	3:H:7:DA:C4	3.07	0.43
2:G:6:DA:H2''	2:G:7:DA:H8	1.84	0.43
3:H:19:DG:H2''	3:H:20:DT:O5'	2.18	0.43
2:G:17:DA:H1'	2:G:18:DC:C5'	2.48	0.43
3:H:2:DG:OP1	3:H:2:DG:H4'	2.19	0.42
1:B:26:ARG:HD3	1:B:26:ARG:HA	1.70	0.42
3:H:1:DC:H3'	3:H:2:DG:C5'	2.48	0.42
3:H:6:DA:C2	3:H:7:DA:C5	3.07	0.42
3:H:7:DA:H2'	3:H:8:DT:C6	2.54	0.42
2:G:12:DG:H1'	2:G:13:DA:C8	2.55	0.42
1:B:175:LEU:C	1:B:175:LEU:HD23	2.40	0.42
3:H:7:DA:C8	3:H:8:DT:H72	2.54	0.42
1:B:109:HIS:CD2	1:B:145:TYR:OH	2.73	0.42
1:A:175:LEU:C	1:A:175:LEU:HD23	2.40	0.42
1:A:187:ASP:HB3	1:A:192:ARG:NH1	2.34	0.42
2:G:7:DA:H2''	2:G:8:DT:C5'	2.50	0.42
1:B:157:ASN:ND2	1:B:212:ASP:OD1	2.53	0.42
1:B:214:CYS:HB3	1:B:215:LEU:CD1	2.48	0.42
2:G:1:DA:C6	2:G:2:DC:N4	2.88	0.42
1:B:83:ARG:HH21	1:B:83:ARG:HG3	1.85	0.41
1:A:56:GLY:O	1:A:57:VAL:CG2	2.54	0.41
1:B:177:ASN:C	1:B:180:PRO:HD2	2.41	0.41
1:A:123:GLN:CD	1:A:142:GLN:HE21	2.24	0.41
2:G:8:DT:H2''	2:G:9:DG:H5'	2.01	0.41
3:H:1:DC:H3'	3:H:2:DG:H5''	2.02	0.41
1:A:191:ALA:O	1:A:195:ARG:HB3	2.20	0.41
3:H:2:DG:N2	3:H:3:DG:C4	2.88	0.41
2:G:4:DT:H2''	2:G:5:DT:C6	2.56	0.40
1:B:77:LEU:HD21	1:B:122:ILE:HG22	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:2:DG:C1'	3:H:3:DG:H5''	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	190/196 (97%)	174 (92%)	11 (6%)	5 (3%)	5	28
1	B	190/196 (97%)	179 (94%)	5 (3%)	6 (3%)	4	22
All	All	380/392 (97%)	353 (93%)	16 (4%)	11 (3%)	4	25

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	25	ASP
1	B	57	VAL
1	B	58	SER
1	A	57	VAL
1	A	190	PRO
1	B	185	ALA
1	A	59	GLY
1	B	190	PRO
1	A	189	LYS
1	B	68	ASN
1	A	68	ASN

### 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	144/148 (97%)	137 (95%)	7 (5%)	25	59
1	B	144/148 (97%)	135 (94%)	9 (6%)	18	49
All	All	288/296 (97%)	272 (94%)	16 (6%)	21	53

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	26	ARG
1	B	29	GLN
1	B	49	GLU
1	B	58	SER
1	B	69	LYS
1	B	112	PHE
1	B	188	SER
1	B	192	ARG
1	B	214	CYS
1	A	26	ARG
1	A	49	GLU
1	A	58	SER
1	A	69	LYS
1	A	179	THR
1	A	189	LYS
1	A	214	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	109	HIS
1	B	142	GLN
1	A	109	HIS
1	A	129	HIS
1	A	142	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](#)

There are no ligands in this entry.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	192/196 (97%)	0.76	36 (18%) <b>1</b> <b>0</b>	127, 159, 214, 238	0
1	B	192/196 (97%)	0.43	10 (5%) <b>27</b> <b>14</b>	126, 148, 220, 241	0
2	G	20/20 (100%)	-0.34	0 <b>100</b> <b>100</b>	119, 146, 175, 182	0
3	H	20/20 (100%)	-0.41	0 <b>100</b> <b>100</b>	111, 139, 173, 175	0
All	All	424/432 (98%)	0.50	46 (10%) <b>5</b> <b>3</b>	111, 152, 215, 241	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	181	HIS	6.9
1	B	181	HIS	6.8
1	A	123	GLN	4.8
1	A	180	PRO	4.6
1	A	43	PHE	4.2
1	A	73	LEU	4.1
1	B	123	GLN	4.1
1	A	38	PHE	3.7
1	A	130	LEU	3.6
1	A	30	LEU	3.5
1	A	125	ARG	3.1
1	B	187	ASP	3.0
1	A	76	LEU	3.0
1	A	188	SER	3.0
1	B	125	ARG	2.9
1	B	180	PRO	2.9
1	A	77	LEU	2.8
1	A	34	ALA	2.8
1	A	215	LEU	2.7
1	A	186	ALA	2.7
1	A	127	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	78	VAL	2.7
1	A	51	ILE	2.7
1	A	72	LEU	2.6
1	A	44	LEU	2.6
1	A	74	VAL	2.6
1	A	210	ALA	2.6
1	A	129	HIS	2.6
1	A	137	GLN	2.6
1	A	211	ALA	2.5
1	B	183	MET	2.4
1	A	212	ASP	2.4
1	A	35	GLU	2.3
1	A	135	GLU	2.3
1	B	188	SER	2.3
1	A	40	GLU	2.2
1	A	136	ARG	2.2
1	A	41	ARG	2.2
1	B	73	LEU	2.1
1	A	149	TRP	2.1
1	B	43	PHE	2.1
1	A	42	GLY	2.1
1	B	126	ASP	2.1
1	A	63	TYR	2.1
1	A	37	LEU	2.1
1	A	134	ALA	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](#)

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.