



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 03:09 pm GMT

PDB ID : 3SQI  
Title : DNA binding domain of Ndc10  
Authors : Cho, U.S.; Harrison, S.C.  
Deposited on : 2011-07-05  
Resolution : 2.82 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

<http://wwpdb.org/validation/2016/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) : 1.9-1692  
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

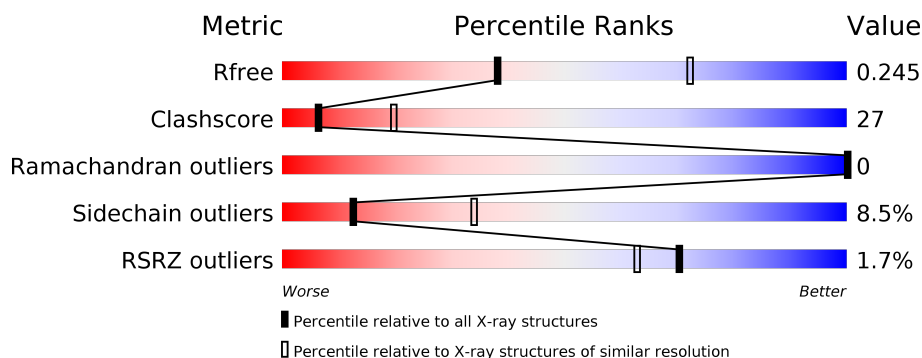
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.82 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	2917 (2.84-2.80)
Clashscore	112137	3382 (2.84-2.80)
Ramachandran outliers	110173	3324 (2.84-2.80)
Sidechain outliers	110143	3326 (2.84-2.80)
RSRZ outliers	101464	2948 (2.84-2.80)

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  $\geq 3$ , 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	534	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0;">%</div> <div style="position: absolute; top: 0; left: 0; width: 38%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 0; left: 38%; width: 30%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 0; left: 68%; width: 1%; height: 10px; background-color: orange;"></div> <div style="position: absolute; top: 0; left: 69%; width: 27%; height: 10px; background-color: grey;"></div> </div> <div>38% 30% 27%</div> </div>
2	B	15	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: 0; left: 0; width: 7%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 0; left: 7%; width: 66%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 0; left: 73%; width: 20%; height: 10px; background-color: orange;"></div> </div> <div>7% 73% 20%</div> </div>
3	C	15	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: 0; left: 0; width: 73%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 0; left: 73%; width: 27%; height: 10px; background-color: orange;"></div> </div> <div>73% 27%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3811 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 KLLA0E03807p.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	388	Total	C	N	O	S	0	0	0
			3196	2086	517	581	12			

- Molecule 2 is a DNA chain called DNA (5'-D(P\*TP\*TP\*AP\*AP\*TP\*TP\*TP\*AP\*TP\*AP\*AP\*AP\*TP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	15	Total	C	N	O	P	0	0	0
			307	150	51	91	15			

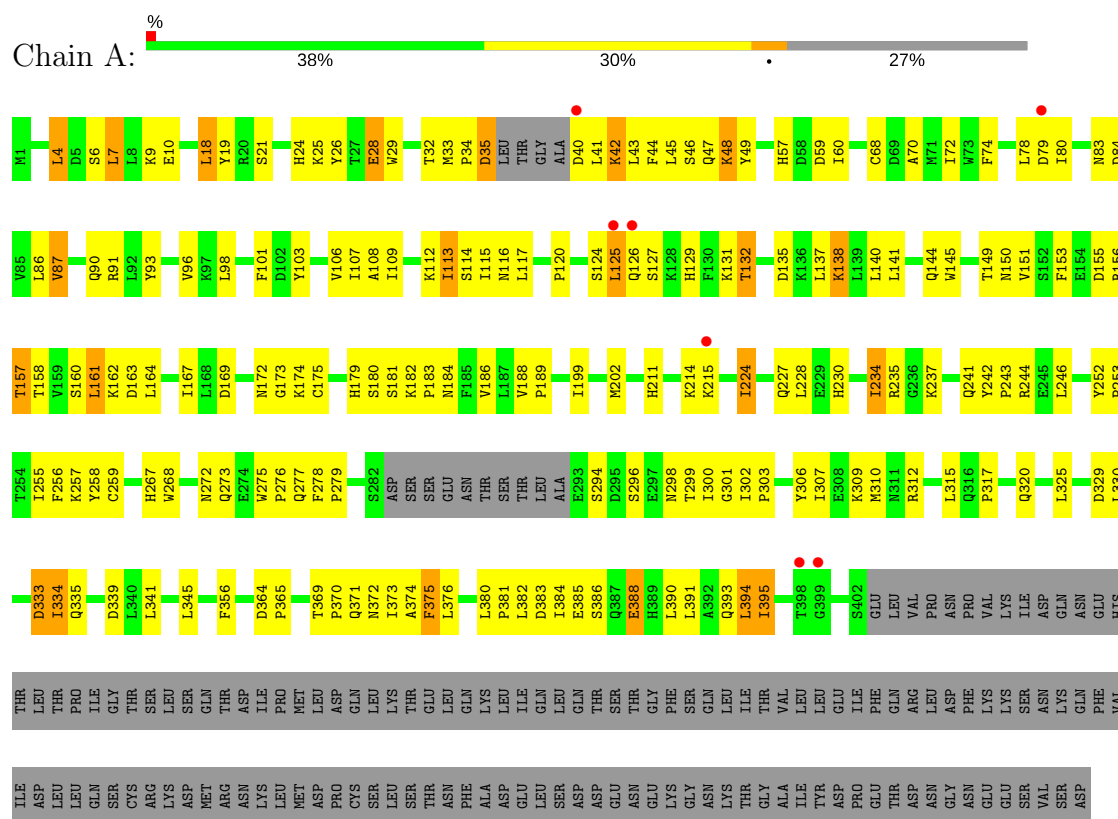
- Molecule 3 is a DNA chain called DNA (5'-D(P\*AP\*AP\*AP\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*AP\*AP\*TP\*TP\*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	15	Total	C	N	O	P	0	0	0
			308	150	54	89	15			

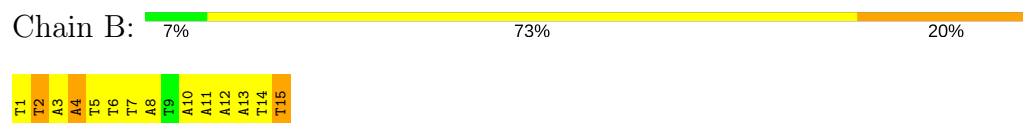
### 3 Residue-property plots

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: KLLA0E03807p



#### • Molecule 2: DNA (5'-D(P\*TP\*TP\*AP\*AP\*TP\*TP\*TP\*AP\*TP\*AP\*AP\*AP\*TP\*T)-3')



#### • Molecule 3: DNA (5'-D(P\*AP\*AP\*AP\*TP\*TP\*TP\*TP\*AP\*TP\*AP\*AP\*AP\*TP\*TP\*A)-3')



A1	A2	A3	T4	T5	T6	T7	A8	T9	A10	A11	A12	T13	T14	A15
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## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	117.30Å 147.54Å 95.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.39 – 2.82 50.00 – 2.82	Depositor EDS
% Data completeness (in resolution range)	91.4 (41.39-2.82) 91.7 (50.00-2.82)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.13 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, $R_{free}$	0.193 , 0.251 0.186 , 0.245	Depositor DCC
$R_{free}$ test set	951 reflections (5.14%)	DCC
Wilson B-factor (Å <sup>2</sup> )	68.6	Xtriage
Anisotropy	1.022	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 81.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3811	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	118.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.35% of the height of the origin peak. No significant pseudotranslation is detected.*

<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

### 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	0.32	0/3287	0.54	0/4473
2	B	0.53	0/343	1.50	4/527 (0.8%)
3	C	0.51	0/345	1.52	6/530 (1.1%)
All	All	0.36	0/3975	0.82	10/5530 (0.2%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	C	6	DT	C1'-O4'-C4'	-9.16	100.94	110.10
2	B	4	DA	O4'-C1'-N9	6.75	112.72	108.00
2	B	15	DT	O4'-C4'-C3'	-6.63	101.85	104.50
3	C	9	DT	C3'-C2'-C1'	-6.41	94.81	102.50
3	C	6	DT	O4'-C1'-N1	6.29	112.41	108.00
2	B	4	DA	C3'-C2'-C1'	-5.42	95.99	102.50
3	C	10	DA	C3'-C2'-C1'	-5.36	96.07	102.50
3	C	9	DT	O4'-C1'-C2'	-5.32	101.65	105.90
3	C	3	DA	C3'-C2'-C1'	-5.12	96.36	102.50
2	B	2	DT	C5-C4-O4	-5.07	121.35	124.90

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	3196	0	3187	161	0
2	B	307	0	174	19	0
3	C	308	0	173	19	0
All	All	3811	0	3534	195	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 27.

All (195) 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:182:LYS:HB2	1:A:183:PRO:HD3	1.54	0.90
1:A:179:HIS:CE1	1:A:181:SER:HB2	2.13	0.83
1:A:214:LYS:HE2	1:A:215:LYS:NZ	1.94	0.83
1:A:86:LEU:O	1:A:90:GLN:NE2	2.12	0.82
2:B:2:DT:H2''	2:B:3:DA:C8	2.15	0.81
1:A:333:ASP:OD2	1:A:333:ASP:N	2.15	0.80
2:B:12:DA:H2''	2:B:13:DA:O5'	1.82	0.78
1:A:160:SER:HB2	1:A:230:HIS:HD2	1.48	0.78
3:C:8:DA:H2''	3:C:9:DT:O5'	1.83	0.77
1:A:214:LYS:HG2	1:A:215:LYS:HD2	1.65	0.76
1:A:156:ARG:NH1	1:A:268:TRP:O	2.19	0.75
1:A:149:THR:HG23	1:A:151:VAL:HG23	1.69	0.74
1:A:169:ASP:OD2	1:A:172:ASN:ND2	2.20	0.73
1:A:125:LEU:HD13	1:A:127:SER:H	1.54	0.73
1:A:155:ASP:O	1:A:158:THR:HG22	1.88	0.73
1:A:371:GLN:O	1:A:374:ALA:N	2.22	0.73
1:A:380:LEU:HG	1:A:381:PRO:HD2	1.72	0.72
1:A:83:ASN:O	1:A:87:VAL:HG13	1.90	0.71
3:C:12:DA:H2''	3:C:13:DT:H72	1.70	0.71
1:A:115:ILE:HD11	1:A:391:LEU:HA	1.72	0.71
1:A:145:TRP:O	1:A:149:THR:HG22	1.90	0.71
1:A:162:LYS:NZ	1:A:228:LEU:O	2.21	0.71
1:A:112:LYS:NZ	1:A:259:CYS:O	2.21	0.70
1:A:160:SER:CB	1:A:230:HIS:HD2	2.03	0.70
1:A:160:SER:HB2	1:A:230:HIS:CD2	2.27	0.69
1:A:333:ASP:HB2	1:A:334:ILE:HD13	1.74	0.69
1:A:6:SER:O	1:A:9:LYS:HG2	1.92	0.69
1:A:18:LEU:HD13	1:A:19:TYR:HD1	1.58	0.69
1:A:18:LEU:HD13	1:A:19:TYR:CD1	2.28	0.68
1:A:115:ILE:HG23	1:A:116:ASN:H	1.58	0.68
1:A:115:ILE:HG23	1:A:116:ASN:N	2.08	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:1:DA:H2''	3:C:2:DA:OP2	1.94	0.68
1:A:370:PRO:O	1:A:373:ILE:HG22	1.94	0.67
1:A:98:LEU:CD1	1:A:309:LYS:HG2	2.25	0.66
1:A:25:LYS:HA	1:A:25:LYS:HE2	1.76	0.65
1:A:214:LYS:HE2	1:A:215:LYS:HZ2	1.59	0.65
3:C:3:DA:H2''	3:C:4:DT:H72	1.78	0.65
1:A:235:ARG:NH2	1:A:237:LYS:O	2.30	0.64
1:A:98:LEU:HD11	1:A:309:LYS:HG2	1.80	0.64
1:A:93:TYR:O	1:A:96:VAL:HG22	1.98	0.64
2:B:2:DT:H2''	2:B:3:DA:H8	1.61	0.64
1:A:42:LYS:HZ2	1:A:42:LYS:H	1.44	0.64
1:A:150:ASN:HD22	1:A:277:GLN:HG2	1.63	0.63
1:A:369:THR:HB	1:A:370:PRO:HD2	1.81	0.62
1:A:252:TYR:O	1:A:255:ILE:HG22	2.00	0.62
2:B:4:DA:H2''	2:B:5:DT:H72	1.82	0.62
1:A:101:PHE:CD2	1:A:276:PRO:HD3	2.35	0.61
1:A:161:LEU:HD12	1:A:161:LEU:C	2.20	0.61
1:A:182:LYS:HB2	1:A:183:PRO:CD	2.29	0.61
1:A:106:VAL:HG11	1:A:275:TRP:CZ2	2.35	0.61
1:A:33:MET:HB3	1:A:34:PRO:HD2	1.83	0.60
1:A:140:LEU:O	1:A:144:GLN:HG3	2.00	0.60
1:A:214:LYS:HE2	1:A:215:LYS:HZ1	1.65	0.60
1:A:237:LYS:O	2:B:15:DT:H5''	2.03	0.59
1:A:115:ILE:HD12	1:A:390:LEU:O	2.03	0.59
1:A:384:ILE:HG13	1:A:385:GLU:N	2.18	0.59
1:A:150:ASN:ND2	1:A:277:GLN:HG2	2.18	0.58
1:A:35:ASP:N	1:A:35:ASP:OD2	2.37	0.58
3:C:4:DT:H2''	3:C:5:DT:H72	1.85	0.58
1:A:372:ASN:O	1:A:376:LEU:HG	2.04	0.57
1:A:29:TRP:O	1:A:32:THR:HB	2.05	0.57
2:B:7:DT:H2''	2:B:8:DA:H8	1.70	0.57
1:A:294:SER:HB2	1:A:299:THR:HA	1.86	0.57
2:B:7:DT:H2''	2:B:8:DA:C8	2.38	0.57
1:A:115:ILE:HB	1:A:390:LEU:HD21	1.87	0.56
1:A:149:THR:CG2	1:A:151:VAL:HG23	2.36	0.56
1:A:83:ASN:O	1:A:86:LEU:HG	2.05	0.56
1:A:24:HIS:O	1:A:28:GLU:HB2	2.05	0.56
1:A:167:ILE:O	1:A:175:CYS:HA	2.05	0.56
3:C:3:DA:H2''	3:C:4:DT:C7	2.35	0.56
1:A:115:ILE:CD1	1:A:391:LEU:HA	2.36	0.55
1:A:124:SER:HB2	1:A:258:TYR:CE1	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:ILE:HB	1:A:390:LEU:CD2	2.35	0.55
1:A:70:ALA:HB1	1:A:74:PHE:HE2	1.71	0.55
1:A:174:LYS:HG2	1:A:315:LEU:HD13	1.88	0.55
1:A:335:GLN:O	1:A:339:ASP:OD2	2.25	0.55
1:A:59:ASP:OD2	1:A:59:ASP:N	2.40	0.55
1:A:385:GLU:HA	1:A:388:GLU:OE1	2.07	0.54
1:A:138:LYS:HG2	1:A:234:ILE:HD12	1.89	0.54
1:A:300:ILE:HG23	1:A:301:GLY:H	1.72	0.54
1:A:120:PRO:HD2	1:A:356:PHE:CD1	2.41	0.54
1:A:179:HIS:NE2	1:A:181:SER:HB2	2.23	0.54
1:A:21:SER:O	1:A:25:LYS:HG2	2.09	0.53
1:A:320:GLN:OE1	1:A:320:GLN:HA	2.07	0.53
1:A:325:LEU:HD12	1:A:345:LEU:HD23	1.90	0.53
1:A:214:LYS:HD3	1:A:365:PRO:HG2	1.89	0.53
1:A:224:ILE:HA	1:A:227:GLN:HG2	1.90	0.53
1:A:383:ASP:O	1:A:386:SER:OG	2.22	0.53
1:A:329:ASP:O	1:A:330:LEU:HG	2.09	0.53
1:A:160:SER:O	1:A:163:ASP:HB2	2.09	0.52
1:A:68:CYS:O	1:A:72:ILE:HG13	2.10	0.52
1:A:124:SER:HB2	1:A:258:TYR:HE1	1.74	0.52
3:C:10:DA:C2	3:C:11:DA:C4	2.97	0.52
1:A:40:ASP:HA	1:A:44:PHE:HE2	1.73	0.52
2:B:8:DA:C2	3:C:10:DA:C2	2.97	0.52
1:A:371:GLN:O	1:A:374:ALA:HB3	2.10	0.51
1:A:106:VAL:HG11	1:A:275:TRP:HZ2	1.73	0.51
1:A:184:ASN:OD1	1:A:302:ILE:HD11	2.10	0.51
3:C:12:DA:H2''	3:C:13:DT:C7	2.38	0.51
1:A:145:TRP:CZ3	1:A:156:ARG:HG2	2.46	0.50
3:C:14:DT:H2''	3:C:15:DA:H5'	1.93	0.50
1:A:115:ILE:CG2	1:A:116:ASN:H	2.22	0.50
1:A:125:LEU:HD13	1:A:127:SER:N	2.25	0.50
1:A:180:SER:HB3	1:A:300:ILE:HG12	1.94	0.50
1:A:70:ALA:O	1:A:74:PHE:HD2	1.95	0.50
1:A:79:ASP:O	1:A:80:ILE:HD12	2.11	0.50
1:A:258:TYR:CE2	1:A:395:ILE:HD11	2.47	0.49
2:B:4:DA:H2''	2:B:5:DT:C7	2.42	0.49
3:C:3:DA:O5'	3:C:3:DA:H2'	2.12	0.49
1:A:4:LEU:O	1:A:7:LEU:HG	2.12	0.49
1:A:278:PHE:CE1	1:A:310:MET:HE3	2.47	0.49
1:A:115:ILE:CG2	1:A:116:ASN:N	2.75	0.49
1:A:188:VAL:HG13	1:A:189:PRO:HD2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:LEU:HD12	1:A:309:LYS:HG2	1.95	0.49
1:A:186:VAL:HG11	1:A:307:ILE:HD13	1.93	0.49
2:B:10:DA:H2''	2:B:11:DA:OP2	2.11	0.49
3:C:3:DA:C2'	3:C:4:DT:H72	2.43	0.49
1:A:103:TYR:CD1	1:A:312:ARG:HD3	2.47	0.49
1:A:46:SER:HB3	1:A:49:TYR:CD2	2.47	0.49
3:C:8:DA:H4'	3:C:9:DT:OP1	2.12	0.49
1:A:112:LYS:O	1:A:116:ASN:ND2	2.45	0.48
1:A:70:ALA:O	1:A:74:PHE:CD2	2.66	0.48
1:A:79:ASP:C	1:A:80:ILE:HD12	2.33	0.48
2:B:6:DT:H2''	2:B:7:DT:O5'	2.13	0.48
1:A:157:THR:HG22	1:A:242:TYR:CD2	2.49	0.48
1:A:242:TYR:CD2	1:A:243:PRO:HD2	2.50	0.47
1:A:141:LEU:HD11	1:A:202:MET:HB2	1.97	0.47
1:A:115:ILE:CD1	1:A:394:LEU:HG	2.45	0.47
1:A:214:LYS:O	1:A:215:LYS:HB2	2.15	0.46
1:A:153:PHE:O	1:A:156:ARG:HB2	2.16	0.46
1:A:26:TYR:HE2	1:A:45:LEU:CD2	2.29	0.46
2:B:12:DA:H4'	2:B:13:DA:OP1	2.15	0.46
1:A:224:ILE:HA	1:A:227:GLN:CG	2.45	0.46
1:A:202:MET:CE	1:A:341:LEU:HD21	2.45	0.46
1:A:241:GLN:CD	1:A:244:ARG:NH1	2.69	0.45
1:A:113:ILE:O	1:A:114:SER:C	2.55	0.45
1:A:155:ASP:HB3	1:A:183:PRO:HG3	1.99	0.45
3:C:4:DT:C2'	3:C:5:DT:H72	2.46	0.45
1:A:32:THR:CG2	1:A:32:THR:O	2.65	0.45
1:A:334:ILE:HD13	1:A:334:ILE:N	2.32	0.44
1:A:279:PRO:HD3	1:A:306:TYR:CE1	2.52	0.44
2:B:14:DT:H2'	2:B:14:DT:H6	1.52	0.44
2:B:1:DT:C2'	2:B:2:DT:H71	2.47	0.44
3:C:1:DA:C2'	3:C:2:DA:OP2	2.64	0.44
1:A:188:VAL:HG21	1:A:315:LEU:HD21	2.00	0.44
3:C:7:DT:H2''	3:C:8:DA:O5'	2.18	0.44
1:A:40:ASP:O	1:A:40:ASP:OD1	2.35	0.44
1:A:241:GLN:CD	1:A:244:ARG:HH12	2.21	0.44
1:A:315:LEU:HD23	1:A:315:LEU:HA	1.83	0.44
1:A:380:LEU:CG	1:A:381:PRO:HD2	2.45	0.44
1:A:25:LYS:HE2	1:A:25:LYS:CA	2.45	0.44
1:A:302:ILE:HG22	1:A:303:PRO:O	2.18	0.43
1:A:48:LYS:HG3	1:A:48:LYS:H	1.46	0.43
1:A:40:ASP:O	1:A:41:LEU:HD12	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:GLN:OE1	1:A:91:ARG:CZ	2.67	0.43
1:A:112:LYS:HD2	1:A:393:GLN:O	2.19	0.43
1:A:137:LEU:HA	1:A:137:LEU:HD23	1.75	0.43
2:B:1:DT:H2''	2:B:2:DT:H71	2.01	0.43
1:A:272:ASN:O	1:A:275:TRP:HB2	2.19	0.43
1:A:57:HIS:HB3	1:A:59:ASP:OD2	2.19	0.43
1:A:160:SER:CB	1:A:230:HIS:CD2	2.93	0.42
2:B:3:DA:H2''	2:B:4:DA:H5'	2.01	0.42
1:A:202:MET:HE3	1:A:202:MET:HB3	1.54	0.42
2:B:2:DT:C2'	2:B:3:DA:C8	2.95	0.42
1:A:173:GLY:HA2	1:A:317:PRO:HB3	2.00	0.42
1:A:234:ILE:O	1:A:243:PRO:HD3	2.19	0.42
1:A:189:PRO:N	1:A:199:ILE:HD12	2.35	0.42
1:A:244:ARG:HG2	3:C:6:DT:H5''	2.00	0.42
1:A:9:LYS:HG3	1:A:10:GLU:N	2.35	0.42
1:A:141:LEU:CD1	1:A:202:MET:HB2	2.50	0.42
1:A:126:GLN:HA	1:A:126:GLN:OE1	2.20	0.42
1:A:132:THR:HG23	1:A:135:ASP:OD2	2.20	0.42
1:A:131:LYS:HG3	1:A:132:THR:N	2.34	0.41
1:A:211:HIS:C	1:A:211:HIS:CD2	2.93	0.41
1:A:300:ILE:CG2	1:A:301:GLY:H	2.33	0.41
1:A:211:HIS:NE2	1:A:364:ASP:OD2	2.53	0.41
1:A:107:ILE:HG23	1:A:108:ALA:N	2.35	0.41
1:A:273:GLN:C	1:A:275:TRP:N	2.73	0.41
1:A:372:ASN:HA	1:A:375:PHE:HD2	1.85	0.41
1:A:117:LEU:HD23	1:A:117:LEU:HA	1.66	0.41
1:A:275:TRP:HA	1:A:276:PRO:HD3	1.94	0.41
1:A:109:ILE:HD12	1:A:109:ILE:N	2.36	0.41
1:A:115:ILE:HD11	1:A:394:LEU:HG	2.03	0.41
1:A:253:PRO:O	1:A:257:LYS:HG3	2.21	0.41
1:A:84:ASP:O	1:A:87:VAL:HG22	2.20	0.41
2:B:12:DA:C2'	2:B:13:DA:O5'	2.63	0.41
1:A:161:LEU:CD1	1:A:161:LEU:C	2.87	0.41
1:A:255:ILE:HG23	1:A:256:PHE:N	2.36	0.41
2:B:12:DA:C2	3:C:5:DT:O4	2.73	0.41
1:A:60:ILE:HD13	1:A:60:ILE:N	2.35	0.41
1:A:300:ILE:HG23	1:A:301:GLY:N	2.35	0.41
3:C:1:DA:H2'	3:C:1:DA:O5'	2.21	0.41
1:A:131:LYS:HG3	1:A:132:THR:HG22	2.02	0.40
1:A:214:LYS:HG2	1:A:215:LYS:CD	2.43	0.40
1:A:373:ILE:HA	1:A:373:ILE:HD12	1.92	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	382/534 (72%)	362 (95%)	20 (5%)	0	100	100

There are no Ramachandran outliers to report.

### 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	366/502 (73%)	335 (92%)	31 (8%)	12	34

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

Mol	Chain	Res	Type
1	A	4	LEU
1	A	7	LEU
1	A	18	LEU
1	A	28	GLU
1	A	35	ASP
1	A	42	LYS
1	A	43	LEU
1	A	48	LYS
1	A	78	LEU
1	A	87	VAL

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Mol	Chain	Res	Type
1	A	113	ILE
1	A	125	LEU
1	A	129	HIS
1	A	132	THR
1	A	138	LYS
1	A	157	THR
1	A	161	LEU
1	A	164	LEU
1	A	224	ILE
1	A	234	ILE
1	A	246	LEU
1	A	267	HIS
1	A	296	SER
1	A	298	ASN
1	A	333	ASP
1	A	334	ILE
1	A	375	PHE
1	A	382	LEU
1	A	388	GLU
1	A	394	LEU
1	A	395	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	ASN
1	A	230	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	388/534 (72%)	-0.11	7 (1%) 69 60	60, 105, 172, 222	0
2	B	15/15 (100%)	-0.93	0 100 100	127, 156, 174, 180	0
3	C	15/15 (100%)	-0.97	0 100 100	118, 158, 181, 199	0
All	All	418/564 (74%)	-0.17	7 (1%) 70 63	60, 108, 174, 222	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	126	GLN	2.6
1	A	399	GLY	2.2
1	A	215	LYS	2.2
1	A	79	ASP	2.2
1	A	398	THR	2.1
1	A	40	ASP	2.1
1	A	125	LEU	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](#)

There are no ligands in this entry.



## 6.5 Other polymers [i](#)

There are no such residues in this entry.