



Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 05:48 am BST

PDB ID : 5I6T
Title : Crystal structure of color device state C
Authors : Hao, Y.; Birktoft, J.; Seeman, N.C.
Deposited on : 2016-02-16
Resolution : 5.28 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 : 2.11
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) : 2.11

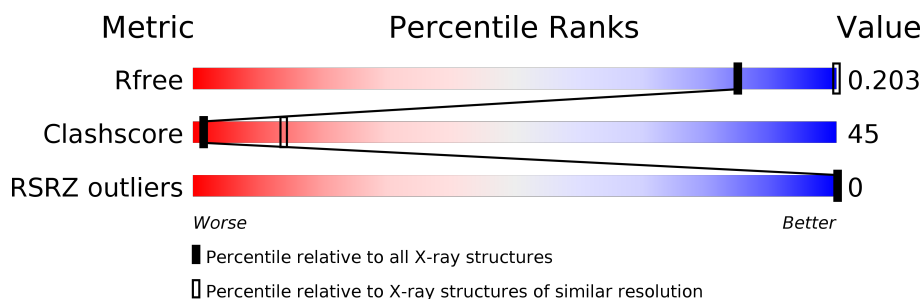
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 5.28 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	1177 (6.76-3.80)
Clashscore	141614	1004 (6.70-3.86)
RSRZ outliers	127900	1003 (6.80-3.76)

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	21	<div> <div>62%</div> <div>38%</div> </div>
2	B	26	<div> <div>23%</div> <div>73%</div> <div>.</div> </div>
3	C	14	<div> <div>7%</div> <div>79%</div> <div>14%</div> </div>
4	D	14	<div> <div>21%</div> <div>71%</div> <div>7%</div> </div>
5	E	14	<div> <div>7%</div> <div>86%</div> <div>7%</div> </div>
6	F	21	<div> <div>90%</div> <div>10%</div> </div>
7	G	21	<div> <div>19%</div> <div>76%</div> <div>5%</div> </div>
8	H	5	<div> <div>100%</div> </div>

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 2767 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 DNA chain called DNA (5'-D(*GP*TP*CP*CP*TP*AP*CP*CP*TP*GP*GP*CP*AP*GP*GP*AP*CP*GP*AP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	21	Total	C	N	O	P	0	0	0
			426	203	79	124	20			

- Molecule 2 is a DNA chain called DNA (26-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	26	Total	C	N	O	P	0	0	0
			520	249	93	153	25			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	14	Total	C	N	O	P	0	0	0
			290	139	53	85	13			

- Molecule 4 is a DNA chain called DNA (5'-D(*AP*CP*AP*GP*TP*CP*GP*TP*GP*GP*TP*AP*TP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	14	Total	C	N	O	P	0	0	0
			285	137	52	83	13			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	14	Total	C	N	O	P	0	0	0
			287	137	52	85	13			

- Molecule 6 is a DNA chain called DNA (5'-D(*CP*AP*GP*AP*TP*AP*CP*CP*TP*GP*

AP*TP*CP*GP*GP*AP*CP*TP*AP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	21	Total	C	N	O	P	0	0	0
			427	204	81	122	20			

- Molecule 7 is a DNA chain called DNA (5'-D(*GP*AP*GP*CP*GP*AP*CP*CP*TP*GP*TP*AP*CP*GP*GP*AP*CP*AP*TP*CP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	G	21	Total	C	N	O	P	0	0	0
			429	204	84	121	20			

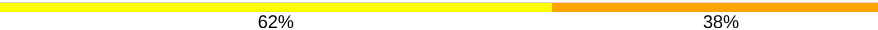
- Molecule 8 is a DNA chain called DNA (5'-D(P*TP*AP*GP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	5	Total	C	N	O	P	0	0	0
			103	49	20	29	5			

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

Chain A: 



- Molecule 2: DNA (26-MER)

Chain B: 



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

Chain C: 



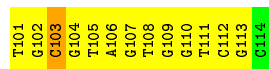
- Molecule 4: DNA (5'-D(*AP*CP*AP*GP*TP*CP*GP*TP*GP*GP*TP*AP*TP*C)-3')

Chain D: 

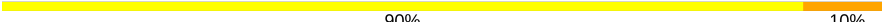


- Molecule 5: DNA (5'-D(*TP*GP*CP*GP*TP*AP*GP*TP*GP*GP*TP*CP*GP*C)-3')

Chain E: 

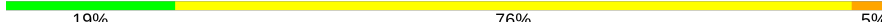


- Molecule 6: DNA (5'-D(*CP*AP*GP*AP*TP*AP*CP*CP*TP*GP*AP*TP*CP*GP*GP*AP*CP*TP*AP*CP*G)-3')

Chain F:  90% 10%

G201
A202
G203
A204
T205
A206
C207
C208
T209
G210
A211
T212
C213
G214
G215
A216
C217
T218
A219
C220
G221

- Molecule 7: DNA (5'-D(*GP*AP*GP*CP*GP*AP*CP*CP*TP*GP*TP*AP*CP*GP*GP*AP*CP*AP*TP*CP*A)-3')

Chain G:  19% 76% 5%

G301
A302
G303
C304
G305
A306
C307
C308
T309
G310
T311
A312
C313
G314
G315
A316
C317
A318
T319
C320
A321

- Molecule 8: DNA (5'-D(P*TP*AP*GP*AP*C)-3')

Chain H:  100%

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	68.06Å 69.41Å 69.86Å 98.73° 97.58° 99.27°	Depositor
Resolution (Å)	41.74 – 5.28 41.74 – 5.28	Depositor EDS
% Data completeness (in resolution range)	98.0 (41.74-5.28) 90.1 (41.74-5.28)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 5.38Å)	Xtriage
Refinement program	PHENIX dev_2276	Depositor
R, R_{free}	0.155 , 0.203 0.155 , 0.203	Depositor DCC
R_{free} test set	437 reflections (9.89%)	wwPDB-VP
Wilson B-factor (Å ²)	247.5	Xtriage
Anisotropy	0.195	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.08 , 109.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.078 for k,l,h 0.078 for l,h,k 0.044 for -k,-h,-l 0.035 for -l,-k,-h 0.053 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	2767	wwPDB-VP
Average B, all atoms (Å ²)	358.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.08% 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](#)

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	1.35	3/477 (0.6%)	1.30	5/734 (0.7%)
2	B	1.10	0/581	1.17	1/892 (0.1%)
3	C	1.10	0/325	1.24	2/502 (0.4%)
4	D	1.17	1/319 (0.3%)	1.23	1/491 (0.2%)
5	E	0.95	0/321	1.12	1/495 (0.2%)
6	F	1.22	2/479 (0.4%)	1.21	1/737 (0.1%)
7	G	1.08	1/482 (0.2%)	1.06	0/742
8	H	1.04	0/115	0.89	0/175
All	All	1.15	7/3099 (0.2%)	1.18	11/4768 (0.2%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	113	DA	C3'-O3'	-7.42	1.34	1.44
6	F	212	DT	C3'-O3'	6.85	1.52	1.44
6	F	212	DT	C1'-N1	6.57	1.57	1.49
1	A	104	DC	C3'-O3'	-6.06	1.36	1.44
4	D	108	DT	C1'-N1	5.89	1.56	1.49
7	G	311	DT	C1'-N1	5.41	1.56	1.49
1	A	119	DA	C3'-O3'	5.24	1.50	1.44

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	112	DC	O5'-P-OP2	-9.97	96.72	105.70
4	D	108	DT	O4'-C4'-C3'	-8.18	101.09	106.00
2	B	110	DA	O4'-C1'-N9	6.77	112.74	108.00
1	A	114	DG	O5'-P-OP1	6.73	118.78	110.70
6	F	206	DA	O4'-C1'-N9	6.32	112.42	108.00
3	C	110	DG	O4'-C1'-N9	5.58	111.91	108.00
1	A	102	DT	O4'-C1'-N1	5.28	111.69	108.00
5	E	103	DC	O4'-C1'-N1	5.26	111.68	108.00
3	C	108	DT	N3-C4-O4	5.15	122.99	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	DA	O4'-C1'-N9	5.08	111.56	108.00
1	A	111	DG	C8-N9-C4	-5.07	104.37	106.40

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	426	0	237	37	0
2	B	520	0	293	40	0
3	C	290	0	161	20	0
4	D	285	0	160	26	0
5	E	287	0	160	17	0
6	F	427	0	237	39	0
7	G	429	0	236	28	0
8	H	103	0	57	0	0
All	All	2767	0	1541	172	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 45.

All (172) 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:104:DC:N4	3:C:113:DG:O6	1.96	0.98
1:A:103:DC:N4	3:C:114:DG:O6	2.06	0.87
5:E:110:DG:N2	7:G:308:DC:N3	2.36	0.74
1:A:108:DC:O2	3:C:110:DG:N2	2.19	0.74
2:B:111:DC:H2'	2:B:112:DC:C2	2.27	0.70
6:F:209:DT:H3'	6:F:210:DG:C8	2.26	0.69
6:F:218:DT:H2''	6:F:219:DA:C8	2.28	0.69
6:F:210:DG:H2'	6:F:211:DA:C8	2.28	0.68
1:A:111:DG:OP2	1:A:111:DG:H8	1.77	0.68
5:E:112:DC:H2'	5:E:113:DG:C8	2.29	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:DG:H3'	1:A:111:DG:C8	2.30	0.66
3:C:112:DA:H2''	3:C:113:DG:H5'	1.77	0.66
2:B:109:DC:H2''	2:B:110:DA:C8	2.31	0.66
3:C:103:DT:H2''	3:C:104:DG:H8	1.61	0.66
2:B:120:DG:H2''	2:B:121:DT:OP2	1.97	0.65
1:A:113:DA:H1'	1:A:114:DG:C5	2.31	0.65
1:A:113:DA:H1'	1:A:114:DG:N7	2.13	0.64
2:B:106:DG:H2''	2:B:107:DA:C8	2.33	0.64
7:G:310:DG:C8	7:G:311:DT:H72	2.34	0.63
2:B:107:DA:H2'	2:B:108:DT:H71	1.81	0.62
5:E:103:DC:H1'	5:E:104:DG:C8	2.35	0.62
2:B:118:DC:H2'	2:B:119:DC:C4	2.35	0.61
5:E:105:DT:H1'	5:E:106:DA:C8	2.36	0.61
2:B:111:DC:H1'	4:D:108:DT:H2'	1.83	0.61
4:D:111:DT:H3	6:F:206:DA:H61	1.47	0.61
3:C:101:DT:H2''	3:C:102:DC:O5'	2.00	0.60
4:D:106:DC:H2''	4:D:107:DG:C8	2.36	0.60
5:E:108:DT:H4'	5:E:109:DG:OP2	2.03	0.59
6:F:202:DA:H8	6:F:202:DA:O5'	1.85	0.59
4:D:111:DT:H2''	4:D:112:DA:C8	2.38	0.59
3:C:111:DT:H6	3:C:111:DT:H5'	1.67	0.59
6:F:209:DT:H3'	6:F:210:DG:H8	1.68	0.59
7:G:312:DA:H2''	7:G:313:DC:O5'	2.03	0.59
7:G:312:DA:H2''	7:G:313:DC:C5'	2.33	0.59
6:F:202:DA:H1'	6:F:203:DG:OP2	2.03	0.58
3:C:103:DT:H2''	3:C:104:DG:C8	2.39	0.58
2:B:118:DC:H42	7:G:315:DG:H1	1.52	0.57
5:E:107:DG:H2''	5:E:108:DT:H5''	1.86	0.57
1:A:120:DC:H2''	1:A:121:DT:H5'	1.87	0.57
4:D:109:DG:C6	4:D:110:DG:C5	2.93	0.57
1:A:118:DG:C2	4:D:107:DG:N2	2.73	0.56
1:A:106:DA:H61	3:C:111:DT:H3	1.54	0.56
7:G:312:DA:H2''	7:G:313:DC:H5'	1.86	0.56
1:A:119:DA:H2''	1:A:120:DC:O5'	2.05	0.56
6:F:205:DT:H2'	6:F:206:DA:C8	2.41	0.56
4:D:108:DT:H4'	4:D:109:DG:OP2	2.06	0.55
4:D:110:DG:N2	6:F:208:DC:O2	2.38	0.55
6:F:206:DA:H1'	6:F:207:DC:H5'	1.88	0.55
4:D:111:DT:H3	6:F:206:DA:N6	2.06	0.54
1:A:109:DT:H3'	1:A:110:DG:H8	1.72	0.54
4:D:103:DA:C6	4:D:104:DG:C6	2.95	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:111:DT:H2''	3:C:112:DA:O5'	2.08	0.54
3:C:112:DA:H2'	3:C:113:DG:C8	2.43	0.54
3:C:108:DT:H4'	3:C:109:DG:OP2	2.08	0.54
2:B:118:DC:H4'	3:C:109:DG:OP1	2.08	0.53
7:G:317:DC:H1'	7:G:318:DA:C8	2.43	0.53
2:B:116:DC:H2'	2:B:117:DA:C8	2.44	0.53
2:B:105:DC:H2''	2:B:106:DG:N7	2.23	0.52
2:B:108:DT:H2''	2:B:109:DC:C5	2.45	0.52
4:D:112:DA:O5'	4:D:112:DA:H8	1.92	0.52
2:B:110:DA:H2''	4:D:109:DG:O4'	2.09	0.52
7:G:303:DG:H2''	7:G:304:DC:H5''	1.91	0.52
2:B:115:DC:H2''	2:B:116:DC:C6	2.45	0.51
7:G:304:DC:H2''	7:G:305:DG:H5'	1.92	0.51
1:A:113:DA:H1'	1:A:114:DG:C8	2.46	0.51
2:B:118:DC:N4	7:G:315:DG:H1	2.08	0.51
6:F:220:DC:H2''	6:F:221:DG:C8	2.46	0.51
1:A:108:DC:H6	1:A:108:DC:H5''	1.76	0.50
2:B:103:DA:C4	5:E:109:DG:C8	2.98	0.50
4:D:110:DG:N2	6:F:208:DC:C2	2.80	0.50
4:D:103:DA:H2''	4:D:104:DG:OP1	2.12	0.50
6:F:219:DA:H2''	6:F:220:DC:O5'	2.10	0.50
4:D:109:DG:C4	4:D:110:DG:C8	3.00	0.49
7:G:304:DC:C2'	7:G:305:DG:H5'	2.42	0.49
1:A:101:DG:H8	1:A:101:DG:HO5'	1.60	0.49
1:A:106:DA:N6	3:C:111:DT:H3	2.09	0.49
6:F:218:DT:H2''	6:F:219:DA:H8	1.77	0.49
6:F:202:DA:O5'	6:F:202:DA:C8	2.64	0.49
6:F:220:DC:H2''	6:F:221:DG:H8	1.78	0.49
1:A:106:DA:H1'	1:A:107:DC:H5'	1.95	0.49
1:A:114:DG:C5	1:A:115:DG:C6	3.01	0.49
6:F:215:DG:C5	6:F:216:DA:C6	3.01	0.48
2:B:110:DA:C6	4:D:109:DG:C5	3.01	0.48
1:A:113:DA:N6	2:B:112:DC:C5	2.81	0.48
2:B:120:DG:H21	7:G:314:DG:N2	2.10	0.48
3:C:107:DG:C2	7:G:318:DA:C2	3.00	0.48
5:E:111:DT:H73	5:E:112:DC:N4	2.28	0.48
6:F:201:DC:N4	6:F:202:DA:N6	2.62	0.48
7:G:319:DT:H2''	7:G:320:DC:O5'	2.14	0.48
2:B:111:DC:C1'	4:D:108:DT:H2'	2.43	0.47
2:B:110:DA:C5	4:D:109:DG:C8	3.02	0.47
4:D:103:DA:H2'	4:D:104:DG:C8	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:210:DG:H4'	6:F:210:DG:OP1	2.15	0.47
6:F:202:DA:H8	6:F:202:DA:P	2.38	0.47
2:B:103:DA:H1'	5:E:109:DG:C1'	2.44	0.47
1:A:101:DG:C8	1:A:102:DT:H72	2.50	0.47
4:D:103:DA:N6	4:D:104:DG:O6	2.47	0.47
6:F:205:DT:H2'	6:F:206:DA:H8	1.79	0.47
2:B:114:DG:C5	2:B:115:DC:C4	3.02	0.47
1:A:107:DC:H2''	1:A:108:DC:H6	1.80	0.47
1:A:113:DA:N1	2:B:113:DT:C4	2.83	0.46
1:A:118:DG:H2''	1:A:119:DA:OP2	2.15	0.46
5:E:104:DG:H2'	5:E:105:DT:H72	1.97	0.46
7:G:310:DG:C6	7:G:311:DT:C4	3.03	0.46
7:G:303:DG:H1'	7:G:304:DC:O4'	2.16	0.46
1:A:112:DC:H4'	1:A:113:DA:OP1	2.16	0.46
2:B:117:DA:H2''	3:C:109:DG:O4'	2.16	0.46
5:E:101:DT:H2'	5:E:102:DG:C8	2.51	0.46
5:E:106:DA:C6	5:E:107:DG:C6	3.04	0.46
2:B:108:DT:H2''	2:B:109:DC:C6	2.50	0.46
2:B:113:DT:H2''	2:B:114:DG:OP2	2.16	0.46
7:G:303:DG:H4'	7:G:304:DC:OP1	2.14	0.46
5:E:107:DG:C8	5:E:108:DT:H71	2.51	0.45
5:E:107:DG:N2	6:F:218:DT:O2	2.49	0.45
6:F:212:DT:H2''	6:F:213:DC:O5'	2.17	0.45
7:G:302:DA:H2''	7:G:303:DG:OP1	2.16	0.45
2:B:119:DC:H2''	2:B:120:DG:C5	2.51	0.45
2:B:101:DA:H5'	2:B:121:DT:H1'	1.99	0.45
6:F:204:DA:H5'	6:F:205:DT:OP2	2.17	0.45
1:A:120:DC:H2''	1:A:121:DT:C5'	2.47	0.45
6:F:209:DT:H6	6:F:209:DT:C5'	2.28	0.45
6:F:214:DG:H2''	6:F:215:DG:OP2	2.15	0.45
7:G:304:DC:N4	7:G:305:DG:O6	2.49	0.45
7:G:306:DA:H1'	7:G:307:DC:O5'	2.17	0.45
5:E:110:DG:N2	7:G:308:DC:C4	2.84	0.45
5:E:110:DG:H8	5:E:110:DG:O5'	2.00	0.44
4:D:111:DT:C2	4:D:112:DA:C5	3.05	0.44
6:F:206:DA:N3	6:F:207:DC:O4'	2.50	0.44
2:B:118:DC:C1'	3:C:108:DT:H2'	2.48	0.44
1:A:110:DG:N2	1:A:111:DG:C2	2.86	0.43
7:G:313:DC:H2''	7:G:314:DG:N7	2.33	0.43
1:A:107:DC:H2''	1:A:108:DC:C6	2.53	0.43
6:F:217:DC:H2''	6:F:218:DT:O5'	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:103:DA:OP2	2:B:103:DA:H2'	2.18	0.43
4:D:102:DC:H1'	4:D:103:DA:OP2	2.18	0.43
2:B:101:DA:H2	7:G:311:DT:H3	1.66	0.43
7:G:302:DA:C2	7:G:303:DG:C6	3.06	0.43
1:A:117:DC:H2''	1:A:118:DG:C8	2.54	0.43
6:F:214:DG:C8	6:F:214:DG:H5''	2.54	0.43
6:F:216:DA:C5	6:F:217:DC:N4	2.86	0.43
7:G:314:DG:H2''	7:G:315:DG:C8	2.54	0.43
1:A:113:DA:C2	2:B:114:DG:N2	2.87	0.42
1:A:119:DA:H8	1:A:119:DA:OP2	2.01	0.42
4:D:109:DG:C5	4:D:110:DG:N7	2.87	0.42
2:B:104:DC:N4	5:E:108:DT:C4	2.86	0.42
2:B:101:DA:N7	2:B:121:DT:C4	2.87	0.42
6:F:216:DA:C5	6:F:217:DC:C4	3.07	0.42
7:G:310:DG:C5	7:G:311:DT:C4	3.07	0.42
1:A:118:DG:C2	1:A:119:DA:C4	3.08	0.42
2:B:111:DC:C5	2:B:112:DC:N4	2.87	0.42
1:A:116:DA:N6	4:D:107:DG:O6	2.53	0.42
6:F:201:DC:C4	6:F:202:DA:N6	2.88	0.42
6:F:201:DC:H2'	6:F:202:DA:C8	2.54	0.42
4:D:101:DA:C8	4:D:101:DA:H5'	2.55	0.42
6:F:202:DA:H2''	6:F:203:DG:OP1	2.19	0.42
1:A:110:DG:C2	1:A:111:DG:C2	3.07	0.41
6:F:209:DT:H6	6:F:209:DT:H5''	1.85	0.41
2:B:118:DC:H2'	2:B:119:DC:C5	2.54	0.41
6:F:211:DA:H2''	6:F:212:DT:OP1	2.20	0.41
6:F:215:DG:C2	6:F:216:DA:C2	3.08	0.41
7:G:314:DG:H2''	7:G:315:DG:OP2	2.19	0.41
2:B:101:DA:N7	2:B:121:DT:C5	2.89	0.41
3:C:113:DG:H1'	3:C:114:DG:O5'	2.21	0.41
3:C:106:DT:C2	3:C:107:DG:C8	3.09	0.41
1:A:102:DT:H1'	1:A:103:DC:OP2	2.21	0.41
2:B:111:DC:H1'	4:D:108:DT:C2	2.55	0.41
6:F:216:DA:C6	6:F:217:DC:N4	2.89	0.41
7:G:318:DA:H2'	7:G:319:DT:H71	2.02	0.41
1:A:115:DG:H1	2:B:111:DC:H42	1.68	0.40
1:A:116:DA:H3'	1:A:116:DA:OP2	2.21	0.40
3:C:102:DC:H2''	3:C:103:DT:OP1	2.21	0.40
1:A:104:DC:C6	1:A:105:DT:H72	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein molecules in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein molecules in this entry.

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 carbohydrates 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, 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	21/21 (100%)	-1.38	0 100 100	235, 292, 367, 383	0
2	B	26/26 (100%)	-1.11	0 100 100	259, 287, 726, 835	0
3	C	14/14 (100%)	-1.42	0 100 100	283, 359, 421, 424	0
4	D	14/14 (100%)	-1.53	0 100 100	273, 340, 398, 399	0
5	E	14/14 (100%)	-1.52	0 100 100	294, 339, 388, 406	0
6	F	21/21 (100%)	-1.43	0 100 100	264, 333, 349, 357	0
7	G	21/21 (100%)	-1.40	0 100 100	307, 325, 380, 391	0
8	H	5/5 (100%)	-0.29	0 100 100	714, 843, 870, 877	0
All	All	136/136 (100%)	-1.33	0 100 100	235, 329, 714, 877	0

There are no RSRZ outliers to report.

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.