



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 02:53 am GMT

PDB ID : 4MFF
Title : Structure of human DNA polymerase beta complexed with O6MG in the template base paired with incoming non-hydrolyzable TTP
Authors : Koag, M.C.; Min, K.; Monzingo, A.F.; Lee, S.
Deposited on : 2013-08-27
Resolution : 2.55 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

<http://wwpdb.org/validation/2016/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
Mogul : 1.7.2 (RC1), CSD as538be (2017)
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

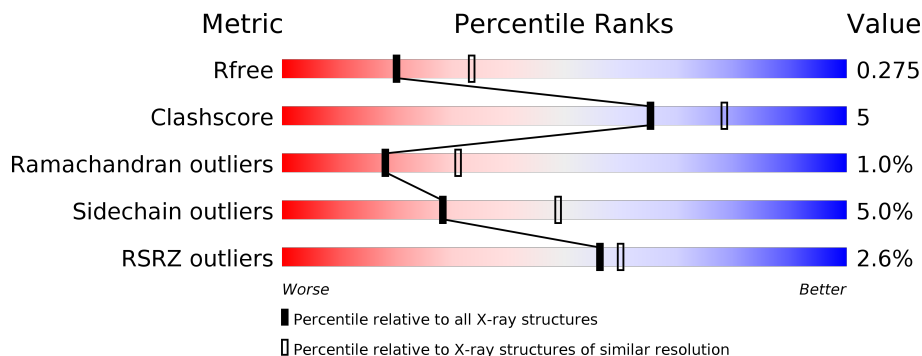
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	4993 (2.58-2.50)
Clashscore	112137	5755 (2.58-2.50)
Ramachandran outliers	110173	5652 (2.58-2.50)
Sidechain outliers	110143	5654 (2.58-2.50)
RSRZ outliers	101464	5026 (2.58-2.50)

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. 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	325	<div> <div>3%</div> <div> <div></div> <div>84%</div> <div>14%</div> <div>..</div> </div> </div>
2	T	16	<div> <div>75%</div> <div>25%</div> </div>
3	P	10	<div> <div>80%</div> <div>20%</div> </div>
4	D	5	<div> <div>80%</div> <div>20%</div> </div>

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 3276 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 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2559	1621	448	481	9			

- Molecule 2 is a DNA chain called template.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	16	Total	C	N	O	P	0	0	0
			322	154	60	93	15			

- Molecule 3 is a DNA chain called up primer.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	10	Total	C	N	O	P	0	0	0
			205	98	40	58	9			

- Molecule 4 is a DNA chain called dn primer.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	5	Total	C	N	O	P	0	0	0
			106	49	20	32	5			

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Mg	0	0
			2	2		

- Molecule 6 is SODIUM ION (three-letter code: NA) (formula: Na).

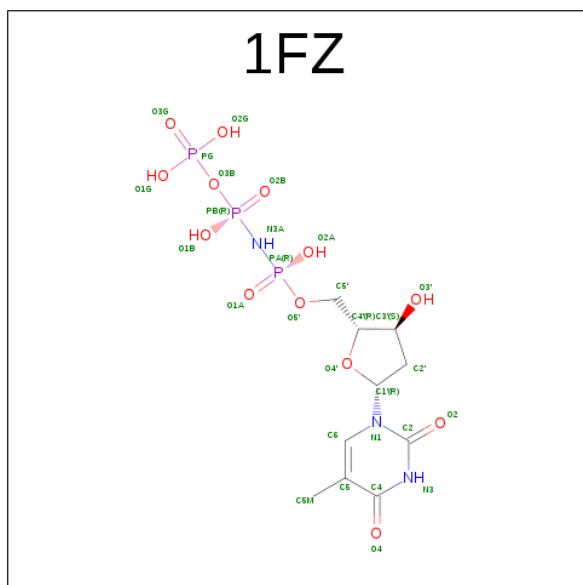
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Na	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	1	Total	Na	0	0
			1	1		

- Molecule 7 is 5'-O-[(R)-HYDROXY{[(R)-HYDROXY(PHOSPHONOOXY)PHOSPHORYL]AMINO}PHOSPHORYL]THYMIDINE (three-letter code: 1FZ) (formula: C₁₀H₁₈N₃O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	P	0	0
			29	10	3	13	3		

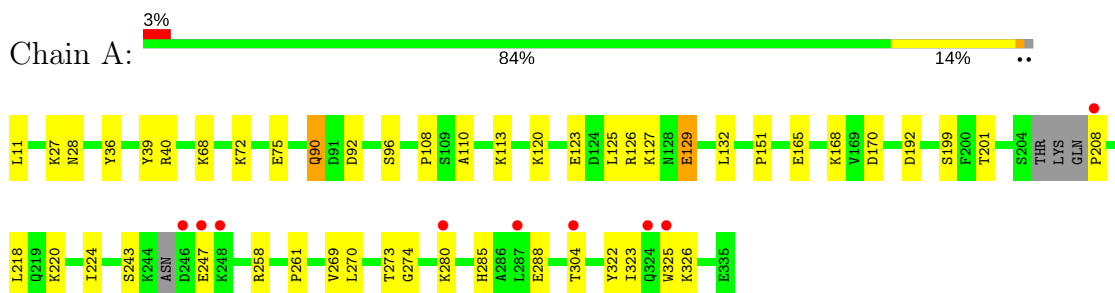
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	40	Total	O	0	0
			40	40		
8	T	5	Total	O	0	0
			5	5		
8	P	5	Total	O	0	0
			5	5		
8	D	1	Total	O	0	0
			1	1		

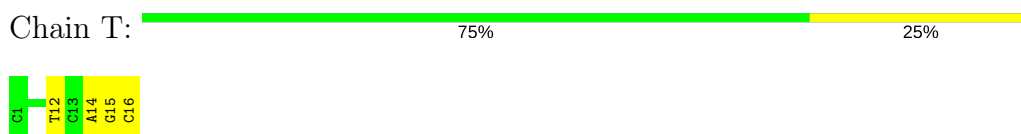
3 Residue-property plots [i](#)

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.

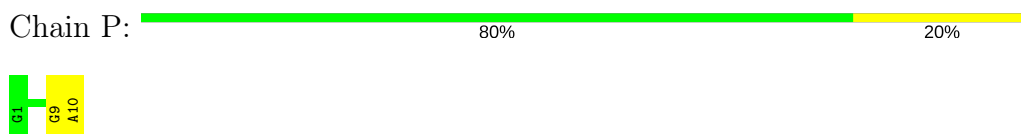
- Molecule 1: DNA polymerase beta



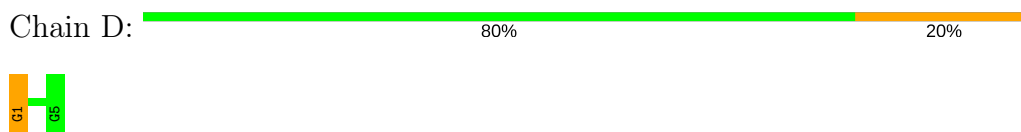
- Molecule 2: template



- Molecule 3: up primer



- Molecule 4: dn primer



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.55Å 78.84Å 54.75Å 90.00° 105.95° 90.00°	Depositor
Resolution (Å)	20.00 – 2.55 19.09 – 2.55	Depositor EDS
% Data completeness (in resolution range)	95.2 (20.00-2.55) 95.5 (19.09-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 2.56Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.222 , 0.294 0.211 , 0.275	Depositor DCC
R_{free} test set	695 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.183	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.034 for l,-k,h	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3276	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.29% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, MG, 6OG, 1FZ

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	0.63	0/2606	0.77	1/3499 (0.0%)
2	T	0.48	0/333	1.11	2/508 (0.4%)
3	P	0.48	0/230	0.95	0/354
4	D	1.00	1/118 (0.8%)	0.86	0/179
All	All	0.62	1/3287 (0.0%)	0.83	3/4540 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1	DG	OP3-P	-9.00	1.50	1.61

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	12	DT	O5'-P-OP1	-6.47	99.88	105.70
1	A	170	ASP	CB-CG-OD1	5.52	123.27	118.30
2	T	14	DA	P-O3'-C3'	5.42	126.20	119.70

There are no chirality outliers.

There are no planarity outliers.

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	2559	0	2572	27	0
2	T	322	0	182	1	0
3	P	205	0	114	1	0
4	D	106	0	57	1	0
5	A	2	0	0	0	1
6	A	1	0	0	0	0
6	D	1	0	0	0	0
7	A	29	0	14	1	0
8	A	40	0	0	10	1
8	D	1	0	0	0	0
8	P	5	0	0	1	0
8	T	5	0	0	0	0
All	All	3276	0	2939	29	1

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 5.

All (29) 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:27:LYS:HD2	8:A:504:HOH:O	1.28	1.32
1:A:199:SER:HA	8:A:513:HOH:O	1.62	1.00
1:A:192:ASP:HB2	8:A:538:HOH:O	1.61	0.99
1:A:208:PRO:HA	8:A:506:HOH:O	1.72	0.89
2:T:15:DG:H2''	2:T:16:DC:OP2	1.82	0.79
1:A:110:ALA:HB2	8:P:105:HOH:O	1.85	0.75
1:A:165:GLU:HA	8:A:524:HOH:O	1.85	0.74
1:A:92:ASP:OD2	1:A:120:LYS:HE2	1.89	0.73
1:A:123:GLU:HG3	1:A:126:ARG:NH2	2.13	0.64
1:A:285:HIS:HD2	1:A:323:ILE:HD12	1.66	0.59
1:A:90:GLN:O	1:A:90:GLN:HG2	2.02	0.58
1:A:168:LYS:CB	8:A:524:HOH:O	2.52	0.57
1:A:168:LYS:HB3	8:A:524:HOH:O	2.04	0.57
1:A:36:TYR:CE1	1:A:40:ARG:HD2	2.43	0.53
1:A:168:LYS:CE	8:A:524:HOH:O	2.57	0.52
1:A:192:ASP:OD1	1:A:258:ARG:NH2	2.33	0.52
1:A:125:LEU:HD22	1:A:132:LEU:HD11	1.93	0.50
1:A:168:LYS:HE3	8:A:524:HOH:O	2.12	0.49
1:A:218:LEU:HB2	1:A:224:ILE:HD12	1.95	0.49
1:A:68:LYS:HB2	4:D:1:DG:H3'	1.96	0.48
1:A:274:GLY:HA3	7:A:404:1FZ:H8	1.96	0.46
3:P:9:DG:C6	3:P:10:DA:C6	3.04	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:GLU:HA	8:A:535:HOH:O	2.17	0.44
1:A:39:TYR:OH	1:A:72:LYS:HE3	2.16	0.43
1:A:90:GLN:HE21	1:A:90:GLN:C	2.21	0.43
1:A:285:HIS:O	1:A:288:GLU:HG2	2.19	0.43
1:A:28:ASN:HA	1:A:108:PRO:HB3	2.02	0.42
1:A:269:VAL:O	1:A:273:THR:OG1	2.27	0.42
1:A:201:THR:C	1:A:261:PRO:HB3	2.42	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:402:MG:MG	8:A:510:HOH:O[1_655]	1.68	0.52

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	315/325 (97%)	292 (93%)	20 (6%)	3 (1%)	18	31

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	322	TYR
1	A	247	GLU
1	A	129	GLU

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	278/286 (97%)	264 (95%)	14 (5%)	28	49

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	LEU
1	A	90	GLN
1	A	96	SER
1	A	113	LYS
1	A	127	LYS
1	A	129	GLU
1	A	151	PRO
1	A	220	LYS
1	A	243	SER
1	A	270	LEU
1	A	280	LYS
1	A	304	THR
1	A	325	TRP
1	A	326	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	90	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	6OG	T	6	2	17,25,26	1.08	1 (5%)	20,36,39	1.99	5 (25%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	6OG	T	6	2	-	2/5/23/24	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	T	6	6OG	C5-C4	3.25	1.47	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	6	6OG	C4-C5-N7	-3.33	106.19	109.41
2	T	6	6OG	N3-C2-N1	-2.52	123.78	127.46
2	T	6	6OG	C2'-C3'-C4'	2.52	108.10	102.73
2	T	6	6OG	C2-N1-C6	4.36	122.48	116.18
2	T	6	6OG	C2-N3-C4	4.41	120.31	115.16

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	T	6	6OG	C-O6-C6-N1
2	T	6	6OG	C-O6-C6-C5

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
7	1FZ	A	404	5	25,30,30	2.38	8 (32%)	31,47,47	2.18	5 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	1FZ	A	404	5	-	0/14/34/34	0/2/2/2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	404	1FZ	C5M-C5	-5.77	1.40	1.51
7	A	404	1FZ	PB-O1B	2.27	1.63	1.56
7	A	404	1FZ	C2-N3	2.68	1.43	1.38
7	A	404	1FZ	PA-O2A	2.81	1.64	1.56
7	A	404	1FZ	O4'-C1'	3.14	1.49	1.42
7	A	404	1FZ	PA-N3A	3.57	1.72	1.63
7	A	404	1FZ	PA-O1A	4.27	1.51	1.46
7	A	404	1FZ	C4-N3	5.55	1.43	1.33

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	404	1FZ	C5-C4-N3	-5.89	118.75	125.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	404	1FZ	PG-O3B-PB	-3.80	118.96	132.38
7	A	404	1FZ	O4'-C1'-N1	2.38	111.79	107.78
7	A	404	1FZ	O1A-PA-N3A	4.03	117.81	111.79
7	A	404	1FZ	C4-N3-C2	7.56	121.77	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	404	1FZ	1	0

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	321/325 (98%)	-0.07	9 (2%) 53 57	7, 17, 46, 57	0
2	T	15/16 (93%)	-0.38	0 100 100	10, 16, 29, 32	0
3	P	10/10 (100%)	-0.49	0 100 100	12, 17, 26, 26	0
4	D	5/5 (100%)	-0.61	0 100 100	10, 11, 16, 20	0
All	All	351/356 (98%)	-0.11	9 (2%) 56 60	7, 17, 45, 57	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	208	PRO	3.6
1	A	304	THR	2.4
1	A	246	ASP	2.4
1	A	248	LYS	2.3
1	A	247	GLU	2.3
1	A	287	LEU	2.3
1	A	280	LYS	2.2
1	A	325	TRP	2.1
1	A	324	GLN	2.1

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	6OG	T	6	23/24	0.94	0.14	-	22,30,33,33	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
7	1FZ	A	404	29/29	0.95	0.16	0.50	24,31,36,37	0
6	NA	D	101	1/1	0.91	0.13	-0.15	16,16,16,16	0
6	NA	A	403	1/1	0.94	0.09	-1.85	12,12,12,12	0
5	MG	A	402	1/1	0.95	0.10	-	34,34,34,34	0
5	MG	A	401	1/1	0.93	0.06	-	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.