



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 18, 2021 – 08:03 AM EDT

PDB ID : 7RSU  
Title : TNA polymerase, n+2 product  
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Deposited on : 2021-08-11  
Resolution : 2.10 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  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.23.2  
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) : 2.23.2

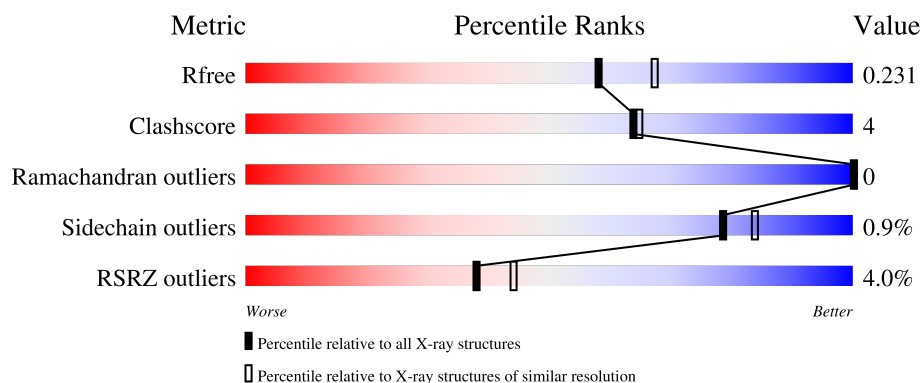
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	774	<div> <div>4%</div> <div>87%</div> <div>10%</div> <div>.</div> </div>
2	T	18	<div> <div>61%</div> <div>17%</div> <div>22%</div> </div>
3	P	13	<div> <div>8%</div> <div>46%</div> <div>38%</div> <div>15%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7043 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	756	Total	C	N	O	S	0	0	0
			6204	3996	1049	1143	16			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	141	ALA	ASP	conflict	UNP D0VWU9
A	143	ALA	GLU	conflict	UNP D0VWU9
A	147	HIS	GLU	conflict	UNP D0VWU9
A	485	ARG	ALA	conflict	UNP D0VWU9
A	491	SER	ASN	conflict	UNP D0VWU9
A	584	LYS	GLU	conflict	UNP D0VWU9
A	606	GLY	ARG	conflict	UNP D0VWU9
A	723	ALA	THR	conflict	UNP D0VWU9

- Molecule 2 is a DNA chain called Template.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	14	Total	C	N	O	P	0	0	0
			287	136	53	84	14			

- Molecule 3 is a DNA chain called Primer.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	13	Total	C	N	O	P	0	0	0
			265	124	51	77	13			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	262	Total	O	0	0
			262	262		

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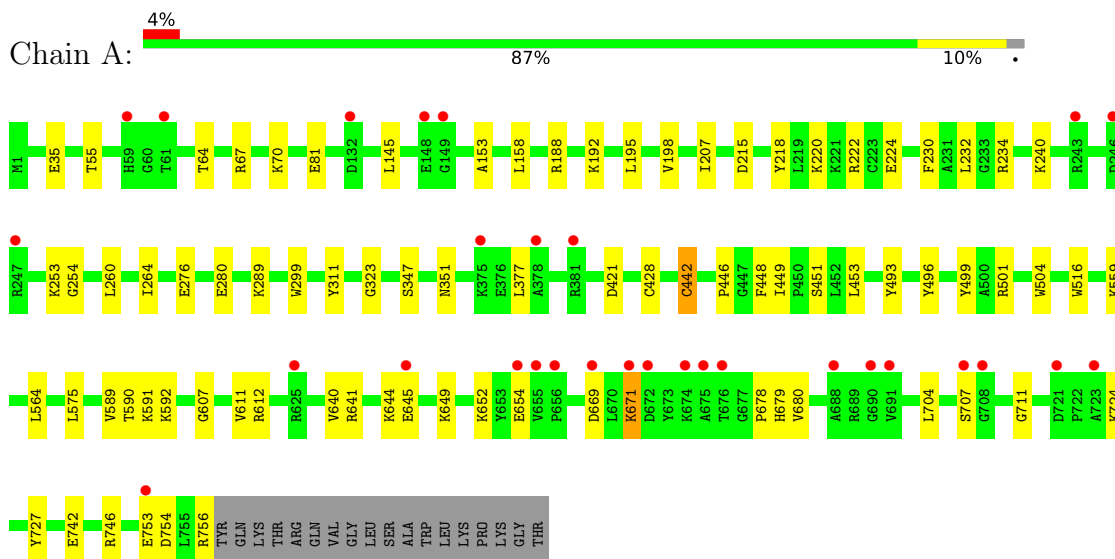
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	T	12	Total 12	O 12	0	0
4	P	13	Total 13	O 13	0	0

### 3 Residue-property plots

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.

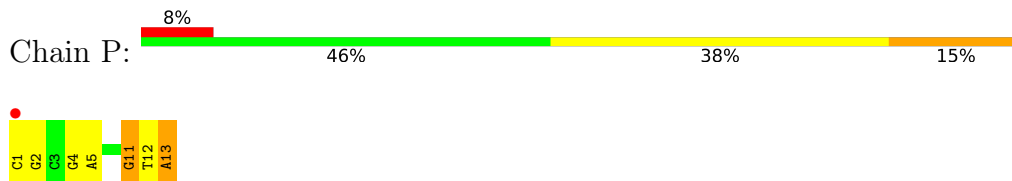
#### • Molecule 1: DNA polymerase



#### • Molecule 2: Template



#### • Molecule 3: Primer



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 <sub>1</sub> 2 <sub>1</sub> 2 <sub>1</sub>	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.54Å 111.69Å 148.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.30 – 2.10 45.29 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.9 (45.30-2.10) 97.9 (45.29-2.10)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.15 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, $R_{free}$	0.182 , 0.231 0.182 , 0.231	Depositor DCC
$R_{free}$ test set	1972 reflections (3.12%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.4	Xtriage
Anisotropy	0.384	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 51.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7043	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: TFT, FA2

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	0.39	0/6343	0.55	0/8556
2	T	0.87	0/321	0.97	0/492
3	P	0.81	0/253	1.12	3/388 (0.8%)
All	All	0.45	0/6917	0.61	3/9436 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	11	DG	O5'-P-OP1	8.74	121.19	110.70
3	P	11	DG	O5'-P-OP2	-5.76	100.52	105.70
3	P	11	DG	O4'-C4'-C3'	-5.52	102.29	104.50

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	6204	0	6250	52	0
2	T	287	0	158	2	0
3	P	265	0	141	5	0
4	A	262	0	0	1	0
4	P	13	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	T	12	0	0	0	0
All	All	7043	0	6549	56	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 4.

All (56) 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:669:ASP:HB3	1:A:671:LYS:HE2	1.72	0.72
1:A:753:GLU:OE2	1:A:753:GLU:N	2.25	0.69
1:A:591:LYS:HD3	2:T:9:DG:H5''	1.77	0.66
1:A:611:VAL:HG23	3:P:11:DG:H5''	1.76	0.66
1:A:449:ILE:HD12	1:A:449:ILE:N	2.12	0.65
1:A:276:GLU:O	1:A:280:GLU:HG3	1.99	0.62
1:A:240:LYS:HE3	1:A:253:LYS:HG2	1.88	0.56
1:A:742:GLU:OE1	1:A:746:ARG:HB3	2.08	0.54
1:A:207:ILE:HD13	1:A:323:GLY:HA3	1.90	0.53
1:A:612:ARG:HG2	3:P:11:DG:H5'	1.90	0.53
1:A:55:THR:HG22	1:A:64:THR:HG22	1.91	0.52
1:A:589:VAL:HG12	1:A:590:THR:HG23	1.92	0.51
1:A:35:GLU:OE2	1:A:67:ARG:NH2	2.44	0.51
1:A:641:ARG:O	1:A:645:GLU:HG2	2.11	0.50
1:A:153:ALA:HA	1:A:218:TYR:CZ	2.46	0.50
1:A:671:LYS:N	1:A:671:LYS:HD3	2.27	0.49
1:A:644:LYS:HE2	1:A:754:ASP:O	2.12	0.49
1:A:453:LEU:HD22	1:A:493:TYR:HD1	1.78	0.48
1:A:679:HIS:CE1	1:A:680:VAL:HG23	2.49	0.48
1:A:446:PRO:HB2	1:A:451:SER:HB2	1.95	0.48
1:A:449:ILE:N	1:A:449:ILE:CD1	2.78	0.47
1:A:640:VAL:HG12	1:A:644:LYS:HE3	1.97	0.46
1:A:652:LYS:HB2	1:A:654:GLU:HG3	1.98	0.46
2:T:15:DC:H2''	2:T:16:DG:C8	2.51	0.46
1:A:559:LYS:HA	1:A:559:LYS:HD3	1.72	0.45
3:P:4:DG:H2''	3:P:5:DA:C8	2.52	0.45
1:A:234:ARG:HD2	1:A:254:GLY:O	2.15	0.45
1:A:70:LYS:HE2	1:A:81:GLU:HB3	1.97	0.45
1:A:428:CYS:SG	1:A:442:CYS:HB2	2.57	0.45
3:P:1:DC:H1'	3:P:2:DG:N7	2.32	0.44
1:A:645:GLU:O	1:A:649:LYS:HG3	2.17	0.44
1:A:195:LEU:HD21	1:A:230:PHE:CD1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:LYS:HE2	1:A:224:GLU:OE1	2.18	0.44
1:A:198:VAL:HG11	1:A:232:LEU:HD22	2.00	0.44
1:A:289:LYS:HG3	4:A:964:HOH:O	2.17	0.44
1:A:428:CYS:HG	1:A:442:CYS:CB	2.31	0.44
1:A:421:ASP:OD1	1:A:421:ASP:N	2.52	0.42
1:A:158:LEU:HD22	1:A:299:TRP:CE2	2.52	0.42
1:A:756:ARG:HE	1:A:756:ARG:HB3	1.54	0.42
3:P:13:FA2:O3'	3:P:13:FA2:H8	2.19	0.42
1:A:289:LYS:NZ	1:A:311:TYR:OH	2.52	0.42
1:A:496:TYR:HE2	1:A:504:TRP:HB2	1.85	0.42
1:A:724:LYS:HB2	1:A:724:LYS:HE2	1.86	0.41
1:A:145:LEU:HB2	1:A:158:LEU:HD21	2.01	0.41
1:A:188:ARG:NH1	1:A:192:LYS:HD2	2.36	0.41
1:A:260:LEU:O	1:A:264:ILE:HG12	2.20	0.41
1:A:449:ILE:HD12	1:A:449:ILE:H	1.86	0.41
1:A:222:ARG:HD2	1:A:222:ARG:HA	1.76	0.41
1:A:448:PHE:CD2	1:A:449:ILE:HD12	2.56	0.41
1:A:499:TYR:CZ	1:A:501:ARG:HB2	2.56	0.41
1:A:564:LEU:HD23	1:A:564:LEU:HA	1.92	0.41
1:A:592:LYS:O	1:A:607:GLY:HA3	2.20	0.40
1:A:678:PRO:HA	1:A:711:GLY:HA2	2.02	0.40
1:A:377:LEU:HD12	1:A:377:LEU:HA	1.84	0.40
1:A:347:SER:OG	1:A:351:ASN:HB3	2.21	0.40
1:A:704:LEU:HA	1:A:727:TYR:HA	2.02	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	754/774 (97%)	740 (98%)	14 (2%)	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	656/671 (98%)	650 (99%)	6 (1%)	78 84

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	215	ASP
1	A	442	CYS
1	A	516	TRP
1	A	575	LEU
1	A	671	LYS
1	A	707	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are 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
3	TFT	P	12	2,3	13,20,21	4.96	11 (84%)	13,29,32	2.51	5 (38%)
3	FA2	P	13	2,3	16,22,23	6.03	12 (75%)	15,32,35	3.21	6 (40%)

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
3	TFT	P	12	2,3	-	2/3/21/22	0/2/2/2
3	FA2	P	13	2,3	-	1/1/21/22	0/3/3/3

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	13	FA2	C2-N3	13.29	1.53	1.32
3	P	13	FA2	C4-N3	9.93	1.49	1.35
3	P	13	FA2	O2'-C2'	-9.66	1.20	1.43
3	P	12	TFT	O2T-C2T	-8.63	1.22	1.43
3	P	12	TFT	O4T-C4T	6.77	1.58	1.43
3	P	13	FA2	C6-C5	-6.77	1.18	1.43
3	P	13	FA2	C8-N7	-6.09	1.23	1.34
3	P	12	TFT	C4-N3	5.94	1.43	1.33
3	P	12	TFT	C4-C5	5.84	1.54	1.41
3	P	12	TFT	O4T-C1T	5.58	1.50	1.41
3	P	12	TFT	C2T-C3T	-5.46	1.39	1.53
3	P	12	TFT	C2T-C1T	-5.37	1.45	1.53
3	P	13	FA2	C6-N1	-5.23	1.14	1.37
3	P	13	FA2	O4'-C1'	4.91	1.49	1.41
3	P	13	FA2	C5-C4	4.36	1.52	1.40
3	P	13	FA2	C2'-C1'	-4.24	1.47	1.53
3	P	13	FA2	O4'-C4'	3.94	1.52	1.43
3	P	13	FA2	C2'-C3'	-3.44	1.44	1.53
3	P	12	TFT	C5M-C5	3.33	1.57	1.51
3	P	13	FA2	C6-N6	3.08	1.45	1.34
3	P	12	TFT	O3T-C3T	-3.02	1.41	1.45
3	P	12	TFT	C2-N3	2.84	1.43	1.38
3	P	12	TFT	C4T-C3T	2.76	1.57	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	13	FA2	C4-C5-N7	-8.37	100.68	109.40
3	P	12	TFT	C4-N3-C2	5.96	120.18	115.14
3	P	13	FA2	C2-N1-C6	5.29	127.80	118.75
3	P	13	FA2	C1'-N9-C4	-4.77	118.27	126.64
3	P	12	TFT	O4T-C4T-C3T	-3.39	98.45	104.52
3	P	13	FA2	N3-C2-N1	-3.30	123.53	128.68
3	P	13	FA2	O4'-C4'-C3'	-3.05	99.07	104.52
3	P	12	TFT	C3T-C2T-C1T	3.03	106.64	99.92
3	P	12	TFT	C4T-C3T-C2T	2.63	106.87	102.28
3	P	13	FA2	C4'-C3'-C2'	2.40	106.47	102.28
3	P	12	TFT	O3T-C3T-C4T	-2.26	101.61	109.08

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	P	12	TFT	C2T-C1T-N1-C6
3	P	12	TFT	O4T-C1T-N1-C6
3	P	13	FA2	C2'-C3'-O3'-P

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	P	13	FA2	1	0

## 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 ⓘ

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	756/774 (97%)	0.39	30 (3%) 38 44	23, 40, 71, 104	0
2	T	14/18 (77%)	-0.04	0 100 100	33, 44, 74, 76	0
3	P	11/13 (84%)	0.44	1 (9%) 9 12	36, 56, 91, 104	0
All	All	781/805 (97%)	0.38	31 (3%) 38 44	23, 40, 73, 104	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	59	HIS	5.7
1	A	132	ASP	4.8
1	A	707	SER	4.2
1	A	671	LYS	3.8
1	A	721	ASP	3.7
1	A	690	GLY	3.3
1	A	708	GLY	3.1
3	P	1	DC	3.1
1	A	675	ALA	3.0
1	A	381	ARG	2.9
1	A	753	GLU	2.8
1	A	655	VAL	2.7
1	A	645	GLU	2.7
1	A	246	ASP	2.7
1	A	375	LYS	2.7
1	A	723	ALA	2.6
1	A	674	LYS	2.6
1	A	672	ASP	2.5
1	A	247	ARG	2.5
1	A	669	ASP	2.5
1	A	148	GLU	2.5
1	A	243	ARG	2.4
1	A	149	GLY	2.4

*Continued on next page...*

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Mol	Chain	Res	Type	RSRZ
1	A	691	VAL	2.4
1	A	688	ALA	2.3
1	A	61	THR	2.2
1	A	676	THR	2.1
1	A	654	GLU	2.1
1	A	656	PRO	2.0
1	A	378	ALA	2.0
1	A	625	ARG	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	FA2	P	13	20/21	0.95	0.12	24,37,48,50	0
3	TFT	P	12	19/20	0.97	0.12	26,32,35,37	0

## 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 [i](#)

There are no such residues in this entry.