



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:59 AM GMT

PDB ID : 3QNO  
Title : RB69 DNA Polymerase (Y567A) Ternary Complex with dATP Opposite 3tCo  
Authors : Xia, S.; Wang, M.; Wang, J.; Konigsberg, W.H.  
Deposited on : 2011-02-08  
Resolution : 1.88 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  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

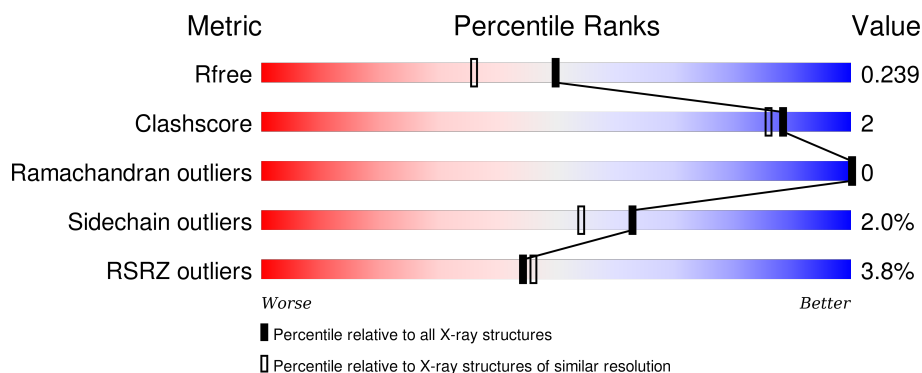
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.88 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))
$R_{free}$	91344	6965 (1.90-1.86)
Clashscore	102246	7778 (1.90-1.86)
Ramachandran outliers	100387	7691 (1.90-1.86)
Sidechain outliers	100360	7692 (1.90-1.86)
RSRZ outliers	91569	6979 (1.90-1.86)

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	901	<div> <div>4%</div> <div>93%</div> <div>7%</div> </div>
2	T	18	<div> <div>11%</div> <div>61%</div> <div>33%</div> <div>6%</div> </div>
3	P	13	<div> <div>69%</div> <div>23%</div> <div>8%</div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9142 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	901	Total	C	N	O	S	0	16	0
			7437	4784	1238	1381	34			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	222	ALA	ASP	CONFLICT	UNP Q38087
A	327	ALA	ASP	CONFLICT	UNP Q38087
A	567	ALA	TYR	ENGINEERED MUTATION	UNP Q38087

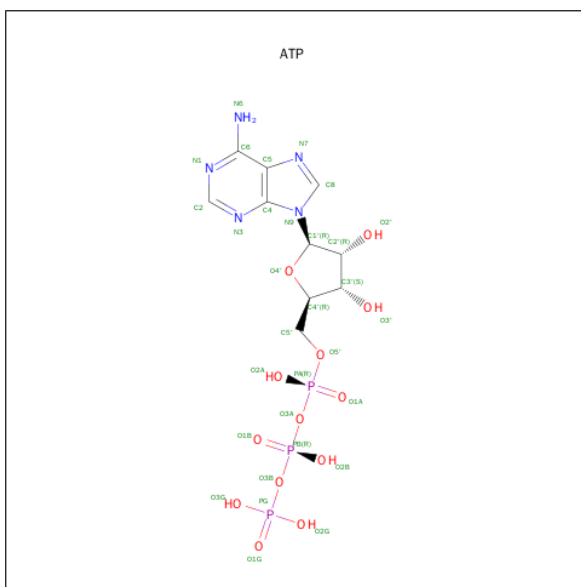
- Molecule 2 is a DNA chain called DNA Tempalte.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	17	Total	C	N	O	P	0	2	0
			395	189	75	112	19			

- Molecule 3 is a DNA chain called DNA Primer.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	13	Total	C	N	O	P	0	0	0
			262	126	48	76	12			

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>13</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			30	10	5	12	3		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total	Ca	0	0
			4	4		

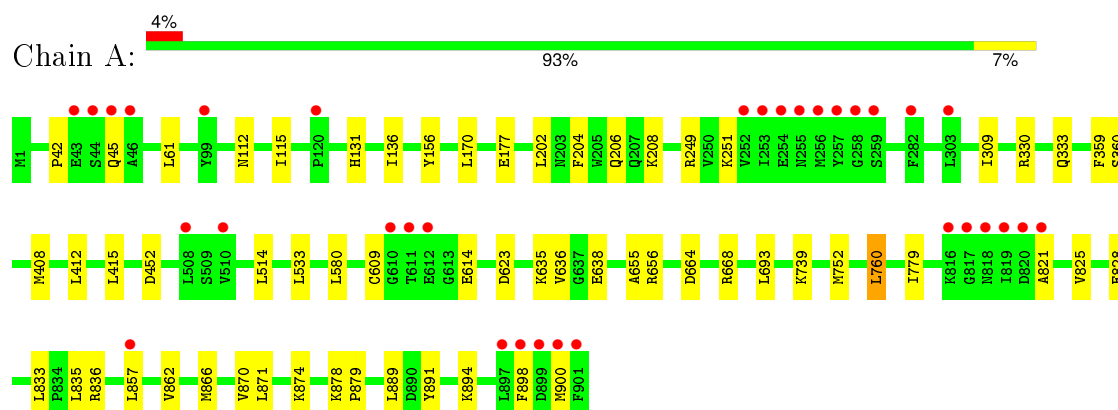
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	885	Total	O	0	12
			897	897		
6	T	80	Total	O	0	0
			80	80		
6	P	36	Total	O	0	1
			37	37		

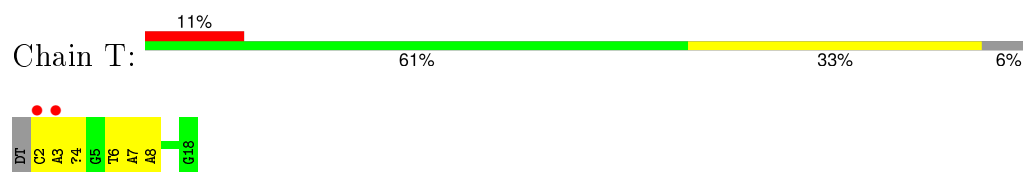
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

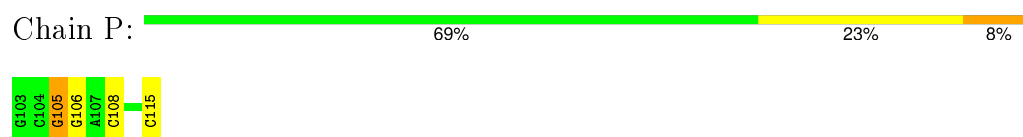
#### • Molecule 1: DNA Polymerase



#### • Molecule 2: DNA Tempalte



#### • Molecule 3: DNA Primer



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.88Å 120.50Å 130.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.88 45.60 – 1.88	Depositor EDS
% Data completeness (in resolution range)	99.0 (50.00-1.88) 99.0 (45.60-1.88)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.17 (at 1.88Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.174 , 0.212 0.200 , 0.239	Depositor DCC
$R_{free}$ test set	4778 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.7	Xtriage
Anisotropy	0.287	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 33.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 95488 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9142	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: DOC, CA, YCO, ATP

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	0.38	0/7665	0.49	0/10356
2	T	0.67	0/412	1.20	3/628 (0.5%)
3	P	0.69	0/273	1.43	3/420 (0.7%)
All	All	0.41	0/8350	0.61	6/11404 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	105	DG	P-O3'-C3'	8.26	129.61	119.70
2	T	8	DA	O4'-C1'-N9	7.25	113.08	108.00
3	P	108	DC	O4'-C1'-N1	6.85	112.80	108.00
3	P	108	DC	C1'-O4'-C4'	-5.92	104.18	110.10
2	T	7	DA	O4'-C1'-N9	-5.58	104.10	108.00
2	T	6	DT	P-O3'-C3'	5.45	126.24	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	7437	0	7396	30	0
2	T	395	0	214	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	P	262	0	148	2	0
4	A	30	0	10	0	0
5	A	4	0	0	0	0
6	A	897	0	0	5	0
6	P	37	0	0	0	0
6	T	80	0	0	0	0
All	All	9142	0	7768	39	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 2.

All (39) 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)
2:T:2[B]:DC:H4'	2:T:3[B]:DA:H2'	1.57	0.86
1:A:330[B]:ARG:HH11	1:A:333:GLN:HE22	1.25	0.81
2:T:2[B]:DC:H4'	2:T:3[B]:DA:H5'	1.67	0.77
2:T:2[B]:DC:H4'	2:T:3[B]:DA:C2'	2.18	0.74
6:A:1236:HOH:O	3:P:115:DOC:H5'	1.98	0.63
2:T:3[A]:DA:O5'	2:T:3[A]:DA:H2'	2.00	0.60
1:A:112[A]:ASN:ND2	6:A:1026:HOH:O	2.34	0.60
1:A:825:VAL:HB	1:A:828:GLU:HG3	1.87	0.56
1:A:870:VAL:HG13	1:A:874:LYS:HD3	1.88	0.56
2:T:2[B]:DC:C4'	2:T:3[B]:DA:H5'	2.37	0.55
1:A:202:LEU:O	1:A:206:GLN:HG3	2.07	0.55
1:A:170:LEU:HA	1:A:177:GLU:HG3	1.90	0.53
2:T:2[B]:DC:C4'	2:T:3[B]:DA:H2'	2.37	0.51
1:A:836:ARG:NH2	6:A:1430:HOH:O	2.42	0.51
1:A:514:LEU:HD13	1:A:533:LEU:HD11	1.92	0.51
1:A:131:HIS:HD2	1:A:156:TYR:OH	1.94	0.51
2:T:2[B]:DC:H4'	2:T:3[B]:DA:C3'	2.41	0.51
1:A:330[B]:ARG:NH1	1:A:333:GLN:HE22	2.03	0.48
1:A:415:LEU:HD22	1:A:623:ASP:HB3	1.96	0.47
1:A:412:LEU:HD13	1:A:415:LEU:HD13	1.97	0.47
1:A:408:MET:SD	1:A:655:ALA:HB2	2.54	0.47
1:A:760:LEU:HD13	1:A:891:TYR:HA	1.95	0.47
1:A:42:PRO:HD2	1:A:45:GLN:HE21	1.81	0.46
3:P:105:DG:H2''	3:P:106:DG:C8	2.51	0.46
1:A:862:VAL:O	1:A:866:MET:HG3	2.16	0.46
1:A:609:CYS:HA	1:A:635:LYS:HE2	1.98	0.45
1:A:249:ARG:HH12	1:A:251:LYS:HG3	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:833:LEU:HD13	1:A:866:MET:HG2	2.00	0.44
1:A:360:SER:OG	2:T:4:YCO:OP1	2.25	0.43
1:A:359:PHE:HB3	2:T:3[A]:DA:C4	2.55	0.42
1:A:664:ASP:O	1:A:668:ARG:HG3	2.19	0.42
1:A:739:LYS:HE3	6:A:1267:HOH:O	2.19	0.42
1:A:330[B]:ARG:HH11	1:A:333:GLN:NE2	2.03	0.41
1:A:752:MET:HG3	1:A:760:LEU:HG	2.01	0.41
1:A:779:ILE:HB	1:A:871:LEU:HD21	2.03	0.41
1:A:204:PHE:CE1	1:A:208:LYS:HD2	2.56	0.41
1:A:878:LYS:HB3	1:A:879:PRO:HD3	2.02	0.41
1:A:115:ILE:HG22	1:A:136:ILE:HG12	2.02	0.40
1:A:821:ALA:HA	6:A:1296:HOH:O	2.21	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	914/901 (101%)	898 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	813/797 (102%)	796 (98%)	17 (2%)	61	52

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	LEU
1	A	309	ILE
1	A	452	ASP
1	A	580	LEU
1	A	614	GLU
1	A	636	VAL
1	A	638	GLU
1	A	656	ARG
1	A	693	LEU
1	A	760	LEU
1	A	835	LEU
1	A	857	LEU
1	A	889[A]	LEU
1	A	889[B]	LEU
1	A	894	LYS
1	A	898	PHE
1	A	900	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	GLN
1	A	131	HIS
1	A	203	ASN
1	A	333	GLN
1	A	339	GLN
1	A	354	GLN
1	A	539	ASN
1	A	761	GLN
1	A	773	GLN
1	A	823	GLN

### 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	DOC	P	115	3,2	11,19,20	0.58	0	14,26,29	1.77	4 (28%)
2	YCO	T	4	2	21,29,30	1.61	2 (9%)	26,42,45	2.43	8 (30%)

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	DOC	P	115	3,2	-	0/3/18/19	0/2/2/2
2	YCO	T	4	2	-	0/3/21/22	0/3/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	T	4	YCO	C5-C4	4.12	1.48	1.41
2	T	4	YCO	C6-N1	5.02	1.42	1.35

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	4	YCO	C13-C8-N7	-4.95	120.96	126.00
2	T	4	YCO	C5-C4-N3	-4.38	118.76	121.28
2	T	4	YCO	C5-C4-N7	-3.23	119.43	121.28
2	T	4	YCO	C12-C13-C8	-2.37	120.14	122.92
3	P	115	DOC	C3'-C2'-C1'	2.11	105.07	102.71
2	T	4	YCO	O14-C13-C12	2.24	118.88	116.10
3	P	115	DOC	O4'-C1'-N1	2.72	112.43	107.72
3	P	115	DOC	O4'-C4'-C5'	3.44	114.61	109.54
3	P	115	DOC	C2-N3-C4	3.57	120.65	115.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	4	YCO	C9-C8-C13	4.10	119.76	115.93
2	T	4	YCO	N7-C4-N3	4.53	121.81	116.03
2	T	4	YCO	C2-N3-C4	6.73	121.43	114.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	P	115	DOC	1	0
2	T	4	YCO	1	0

## 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	$\# Z  > 2$
4	ATP	A	902	5	24,32,33	0.96	1 (4%)	32,50,52	1.85	4 (12%)

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
4	ATP	A	902	5	-	0/18/34/38	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	902	ATP	C5-C4	3.19	1.47	1.40

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	902	ATP	N3-C2-N1	-7.38	123.24	128.89
4	A	902	ATP	C2'-C1'-N9	-3.41	105.86	114.16
4	A	902	ATP	C4-C5-N7	-2.87	106.84	109.48
4	A	902	ATP	C2-N1-C6	2.27	122.82	118.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	901/901 (100%)	0.22	33 (3%) 45 47	13, 22, 45, 75	0
2	T	16/18 (88%)	0.14	2 (12%) 5 5	18, 24, 39, 40	0
3	P	12/13 (92%)	0.02	0 100 100	18, 32, 48, 48	0
All	All	929/932 (99%)	0.22	35 (3%) 44 46	13, 22, 45, 75	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	256	MET	8.8
1	A	819	ILE	6.4
1	A	257	TYR	6.1
1	A	900	MET	6.0
1	A	44	SER	5.4
1	A	252	VAL	5.3
1	A	255	ASN	5.3
2	T	2[A]	DC	5.0
1	A	254	GLU	5.0
1	A	901	PHE	4.6
1	A	46	ALA	4.6
1	A	857	LEU	4.5
1	A	510	VAL	4.0
1	A	253	ILE	3.9
1	A	610	GLY	3.7
1	A	508	LEU	3.7
1	A	817	GLY	3.6
1	A	303	LEU	3.5
1	A	820	ASP	3.0
1	A	897	LEU	2.9
1	A	45	GLN	2.9
1	A	259	SER	2.8
1	A	43	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	898	PHE	2.7
1	A	120	PRO	2.6
1	A	258	GLY	2.6
1	A	99	TYR	2.5
2	T	3[A]	DA	2.4
1	A	282	PHE	2.4
1	A	821	ALA	2.4
1	A	612	GLU	2.3
1	A	816	LYS	2.3
1	A	899	ASP	2.1
1	A	611	THR	2.1
1	A	818	ASN	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. 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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	DOC	P	115	18/19	0.97	0.17	-	16,17,25,26	0
2	YCO	T	4	26/27	0.96	0.14	-	16,19,23,25	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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	ATP	A	902	30/31	0.98	0.15	-0.28	16,16,21,23	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	CA	A	903	1/1	0.98	0.12	-1.33	19,19,19,19	0
5	CA	A	905	1/1	0.96	0.05	-2.34	38,38,38,38	0
5	CA	A	906	1/1	0.92	0.07	-	43,43,43,43	0
5	CA	A	904	1/1	0.99	0.06	-	34,34,34,34	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.