



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:11 PM GMT

PDB ID : 4BPW  
Title : Crystal structure of human primase bound to UTP  
Authors : Kilkenny, M.L.; Perera, R.L.; Pellegrini, L.  
Deposited on : 2013-05-28  
Resolution : 3.00 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

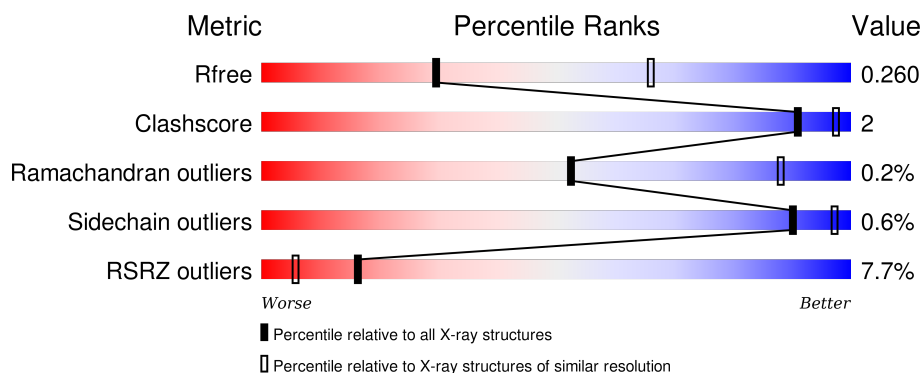
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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	423	<div> <div>6%</div> <div>82% 9% 10%</div> </div>
1	C	423	<div> <div>6%</div> <div>83% • 13%</div> </div>
2	B	253	<div> <div>13%</div> <div>77% 9% 13%</div> </div>
2	D	253	<div> <div>11%</div> <div>77% 5% 18%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	MG	A	1413	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 19731 atoms, of which 9873 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 PRIMASE SMALL SUBUNIT.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	382	Total	C	H	N	O	S	0	2	0
			6406	2066	3195	558	574	13			
1	C	367	Total	C	H	N	O	S	0	0	0
			6128	1982	3057	529	547	13			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P49642
A	-1	THR	-	EXPRESSION TAG	UNP P49642
A	0	SER	-	EXPRESSION TAG	UNP P49642
A	72	ALA	LYS	ENGINEERED MUTATION	UNP P49642
A	73	ALA	MET	ENGINEERED MUTATION	UNP P49642
C	-2	GLY	-	EXPRESSION TAG	UNP P49642
C	-1	THR	-	EXPRESSION TAG	UNP P49642
C	0	SER	-	EXPRESSION TAG	UNP P49642
C	72	ALA	LYS	ENGINEERED MUTATION	UNP P49642
C	73	ALA	MET	ENGINEERED MUTATION	UNP P49642

- Molecule 2 is a protein called DNA PRIMASE LARGE SUBUNIT.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	219	Total	C	H	N	O	S	0	0	0
			3646	1156	1849	310	329	2			
2	D	208	Total	C	H	N	O	S	0	0	0
			3465	1101	1750	295	317	2			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

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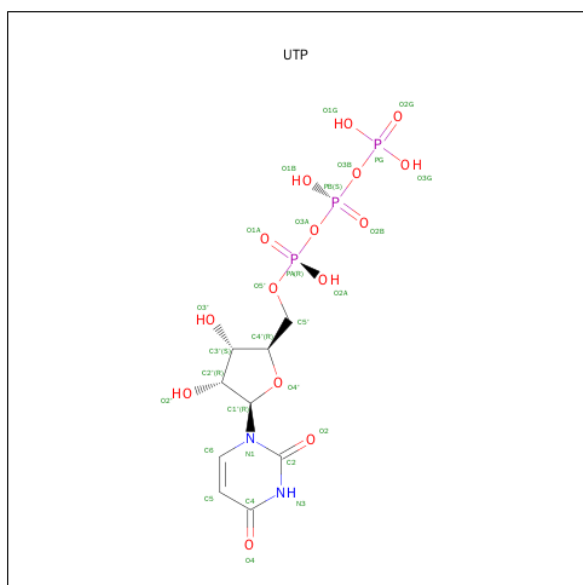
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total	Zn	0	0
			1	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		
4	C	2	Total	Mg	0	0
			2	2		

- Molecule 5 is URIDINE 5'-TRIPHOSPHATE (three-letter code: UTP) (formula: C<sub>9</sub>H<sub>15</sub>N<sub>2</sub>O<sub>15</sub>P<sub>3</sub>).

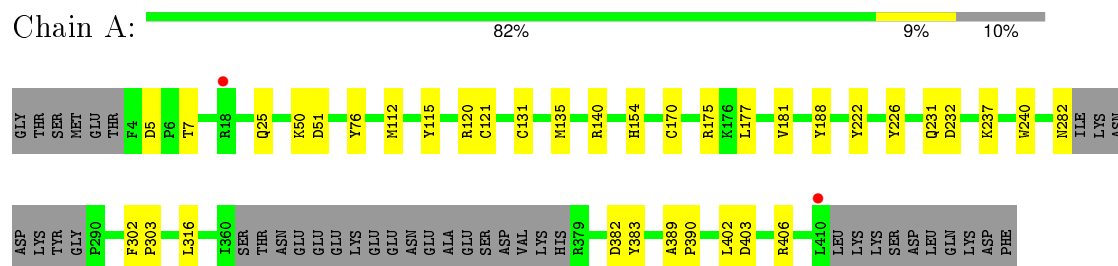


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	P	0	0
			40	9	11	2	15	3		
5	C	1	Total	C	H	N	O	P	0	0
			40	9	11	2	15	3		

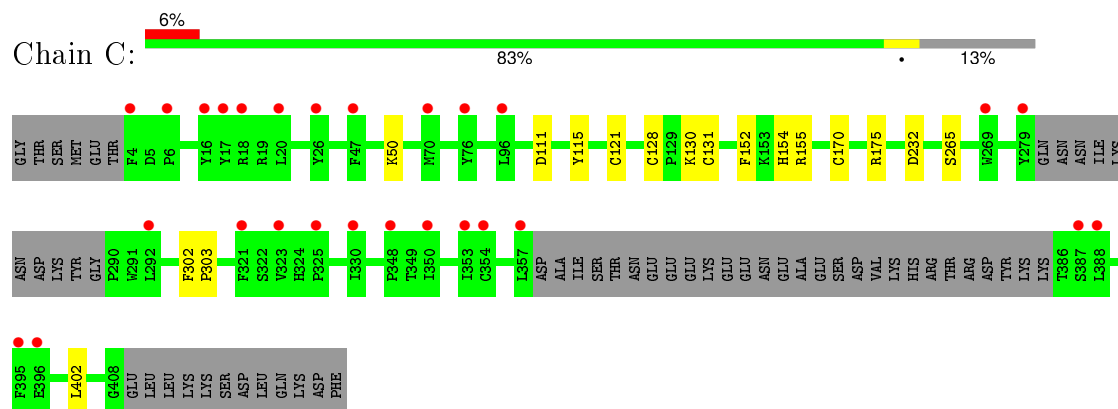
### 3 Residue-property plots [i](#)

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.

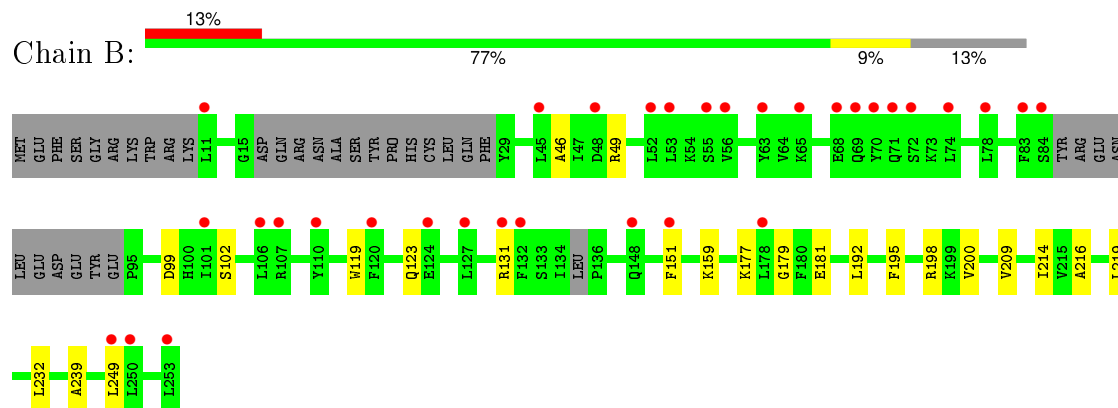
#### • Molecule 1: DNA PRIMASE SMALL SUBUNIT



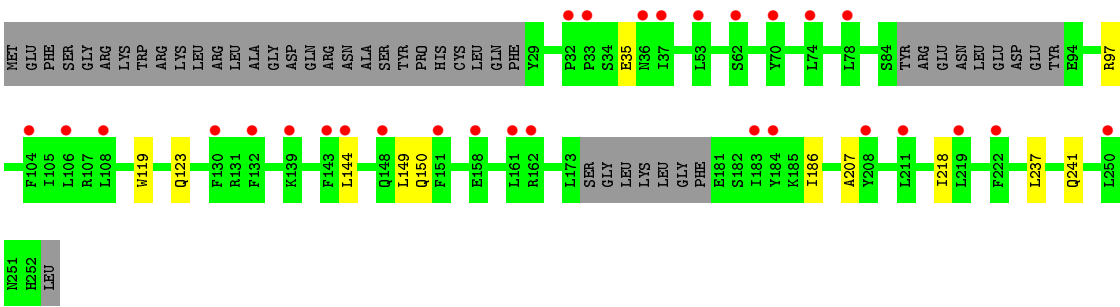
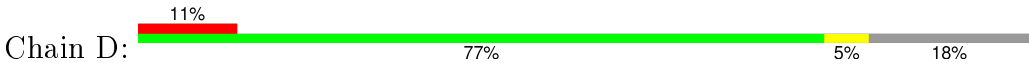
#### • Molecule 1: DNA PRIMASE SMALL SUBUNIT



#### • Molecule 2: DNA PRIMASE LARGE SUBUNIT



#### • Molecule 2: DNA PRIMASE LARGE SUBUNIT



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	116.39Å 69.44Å 127.09Å 90.00° 104.90° 90.00°	Depositor
Resolution (Å)	46.78 – 3.00 48.87 – 3.00	Depositor EDS
% Data completeness (in resolution range)	98.3 (46.78-3.00) 99.1 (48.87-3.00)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.76 (at 3.01Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.215 , 0.253 0.227 , 0.260	Depositor DCC
$R_{free}$ test set	1995 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	106.5	Xtriage
Anisotropy	0.363	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 65.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 39279 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	19731	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	103.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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: UTP, MG, ZN

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	0.29	0/3306	0.50	0/4461
1	C	0.25	0/3152	0.45	0/4255
2	B	0.25	0/1827	0.45	0/2450
2	D	0.24	0/1745	0.44	0/2347
All	All	0.26	0/10030	0.47	0/13513

There are no bond length outliers.

There are no bond angle outliers.

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	3211	3195	3171	21	0
1	C	3071	3057	3044	8	0
2	B	1797	1849	1846	10	0
2	D	1715	1750	1741	6	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	A	2	0	0	0	0
4	C	2	0	0	0	0
5	A	29	11	11	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	29	11	11	0	0
All	All	9858	9873	9824	44	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 (44) 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:C:170:CYS:O	1:C:175:ARG:NH1	2.20	0.73
1:A:403:ASP:OD1	1:A:406:ARG:NH2	2.22	0.71
1:A:140:ARG:NH1	1:A:222:TYR:OH	2.26	0.69
1:A:120:ARG:NH1	1:A:232:ASP:O	2.28	0.66
2:B:159:LYS:NZ	2:B:181:GLU:O	2.32	0.62
1:A:170:CYS:O	1:A:175:ARG:NH1	2.33	0.61
1:A:188:TYR:O	2:B:198:ARG:NH1	2.36	0.58
2:B:49:ARG:NH1	2:B:102:SER:OG	2.38	0.56
1:A:316:LEU:O	5:A:1414:UTP:H1'	2.06	0.56
1:C:232:ASP:OD1	1:C:265:SER:OG	2.24	0.55
1:A:237:LYS:HA	1:A:240:TRP:CE2	2.44	0.53
2:D:186:ILE:O	2:D:207:ALA:N	2.40	0.51
1:A:50:LYS:HG2	1:A:76:TYR:HE2	1.76	0.51
1:A:25:GLN:NE2	1:A:383:TYR:OH	2.44	0.50
1:A:112:MET:HE1	1:A:131:CYS:HB2	1.93	0.50
1:A:112:MET:HG2	1:A:135:MET:HE3	1.95	0.49
2:B:46:ALA:HB1	2:B:249:LEU:HD13	1.95	0.48
1:A:302:PHE:CG	1:A:303:PRO:HD2	2.49	0.48
1:A:5:ASP:OD2	1:A:7:THR:OG1	2.31	0.48
1:C:121:CYS:SG	1:C:131:CYS:HB3	2.54	0.47
2:D:119:TRP:O	2:D:123:GLN:NE2	2.49	0.46
1:A:115:TYR:OH	1:A:303:PRO:HA	2.15	0.46
1:A:121:CYS:SG	1:A:131:CYS:HB3	2.56	0.46
2:D:35:GLU:OE2	2:D:97:ARG:NE	2.39	0.45
1:A:154:HIS:CG	1:A:402:LEU:HD23	2.52	0.44
2:D:237:LEU:HG	2:D:241:GLN:HG3	1.99	0.44
2:B:99:ASP:OD1	2:B:131:ARG:NH2	2.43	0.43
1:C:115:TYR:OH	1:C:303:PRO:HA	2.18	0.43
1:C:154:HIS:CG	1:C:402:LEU:HD23	2.54	0.43
2:D:149:LEU:HG	2:D:150:GLN:H	1.82	0.43
2:B:119:TRP:O	2:B:123:GLN:NE2	2.51	0.43
1:A:302:PHE:CD1	1:A:303:PRO:HD2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:LEU:HD22	1:A:181:VAL:HG11	2.01	0.43
1:C:152:PHE:O	1:C:155:ARG:NH1	2.52	0.42
2:B:151:PHE:CE1	2:B:214:ILE:HD11	2.55	0.42
2:B:216:ALA:HA	2:B:219:LEU:HB3	1.99	0.42
1:C:302:PHE:CG	1:C:303:PRO:HD2	2.53	0.42
1:A:389:ALA:N	1:A:390:PRO:HD2	2.35	0.42
1:A:50:LYS:O	1:A:51:ASP:HB2	2.19	0.42
1:A:226:TYR:CZ	1:A:231:GLN:HG3	2.56	0.41
2:B:192:LEU:HA	2:B:195:PHE:CE2	2.56	0.41
1:C:128:CYS:SG	1:C:130:LYS:HB2	2.61	0.40
2:B:200:VAL:HG21	2:B:209:VAL:HG22	2.03	0.40
2:D:144:LEU:HD23	2:D:218:ILE:HD11	2.02	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	378/423 (89%)	367 (97%)	11 (3%)	0	100	100
1	C	361/423 (85%)	350 (97%)	11 (3%)	0	100	100
2	B	211/253 (83%)	201 (95%)	8 (4%)	2 (1%)	21	64
2	D	202/253 (80%)	195 (96%)	7 (4%)	0	100	100
All	All	1152/1352 (85%)	1113 (97%)	37 (3%)	2 (0%)	52	88

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	239	ALA
2	B	179	GLY

### 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	357/393 (91%)	355 (99%)	2 (1%)	90	97
1	C	341/393 (87%)	339 (99%)	2 (1%)	90	97
2	B	198/230 (86%)	196 (99%)	2 (1%)	82	95
2	D	189/230 (82%)	189 (100%)	0	100	100
All	All	1085/1246 (87%)	1079 (99%)	6 (1%)	90	97

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

Mol	Chain	Res	Type
1	A	282	ASN
1	A	382	ASP
2	B	177	LYS
2	B	232	LEU
1	C	50	LYS
1	C	111	ASP

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	154	HIS
1	C	154	HIS

### 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](#)

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 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	$\# Z  > 2$
5	UTP	A	1414	4	20,30,30	1.26	2 (10%)	27,47,47	1.25	4 (14%)
5	UTP	C	1412	4	20,30,30	1.14	2 (10%)	27,47,47	1.24	4 (14%)

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
5	UTP	A	1414	4	-	0/18/38/38	0/2/2/2
5	UTP	C	1412	4	-	0/18/38/38	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1414	UTP	O4'-C1'	-3.50	1.36	1.41
5	C	1412	UTP	O4'-C1'	-2.63	1.37	1.41
5	A	1414	UTP	O4'-C4'	-2.49	1.39	1.45
5	C	1412	UTP	O4'-C4'	-2.26	1.39	1.45

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1414	UTP	PB-O3A-PA	-3.66	122.47	132.73
5	C	1412	UTP	PB-O3B-PG	-2.70	123.61	132.67
5	A	1414	UTP	PB-O3B-PG	-2.37	124.73	132.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	1412	UTP	PB-O3A-PA	-2.22	126.50	132.73
5	A	1414	UTP	O3A-PA-O5'	2.04	108.34	102.94
5	A	1414	UTP	C4'-O4'-C1'	2.64	112.61	109.72
5	C	1412	UTP	O4'-C1'-N1	2.88	114.16	108.08
5	C	1412	UTP	O3A-PA-O5'	2.95	110.77	102.94

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
5	A	1414	UTP	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, 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	382/423 (90%)	0.16	2 (0%) 91 76	22, 53, 96, 123	0
1	C	367/423 (86%)	0.45	27 (7%) 17 6	50, 97, 154, 178	0
2	B	219/253 (86%)	0.79	33 (15%) 3 1	45, 128, 174, 190	0
2	D	208/253 (82%)	0.72	29 (13%) 4 1	72, 118, 164, 186	0
All	All	1176/1352 (86%)	0.46	91 (7%) 16 6	22, 94, 159, 190	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	36	ASN	7.7
2	B	11	LEU	6.0
2	D	151	PHE	5.3
2	B	110	TYR	5.2
2	D	32	PRO	5.2
2	B	132	PHE	5.1
2	B	249	LEU	5.0
2	D	144	LEU	5.0
2	B	83	PHE	4.9
1	C	354	CYS	4.8
2	D	184	TYR	4.8
1	C	388	LEU	4.7
2	B	45	LEU	4.7
2	B	74	LEU	4.6
1	C	70	MET	4.1
2	B	52	LEU	4.1
2	B	48	ASP	3.9
2	B	53	LEU	3.9
1	C	330	ILE	3.8
2	B	68	GLU	3.8
1	C	350	ILE	3.7

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Mol	Chain	Res	Type	RSRZ
2	D	74	LEU	3.7
2	D	37	ILE	3.5
1	C	20	LEU	3.5
2	B	84	SER	3.4
1	C	325	PRO	3.4
2	D	222	PHE	3.4
1	C	323	VAL	3.3
1	C	321	PHE	3.3
2	D	250	LEU	3.3
2	D	183	ILE	3.3
1	C	76	TYR	3.2
2	D	143	PHE	3.2
2	D	53	LEU	3.2
2	B	65	LYS	3.2
2	B	151	PHE	3.1
2	D	70	TYR	3.1
1	C	6	PRO	3.1
2	B	148	GLN	3.1
1	C	4	PHE	3.0
1	C	353	ILE	2.9
2	D	162	ARG	2.9
2	B	178	LEU	2.9
2	B	69	GLN	2.9
2	B	70	TYR	2.9
2	B	253	LEU	2.8
2	B	101	ILE	2.8
1	A	410	LEU	2.8
2	D	161	LEU	2.7
2	D	62	SER	2.7
2	B	78	LEU	2.7
2	B	72	SER	2.7
2	D	108	LEU	2.6
1	C	357	LEU	2.6
1	C	279	TYR	2.6
2	B	127	LEU	2.5
2	B	131	ARG	2.5
1	C	96	LEU	2.5
1	C	396	GLU	2.5
2	B	124	GLU	2.5
2	D	130	PHE	2.5
1	C	18	ARG	2.5
1	A	18	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	120	PHE	2.4
2	B	107	ARG	2.4
1	C	387	SER	2.4
2	D	132	PHE	2.4
2	D	208	TYR	2.4
1	C	348	PRO	2.4
1	C	17	TYR	2.4
2	D	139	LYS	2.3
1	C	26	TYR	2.3
2	D	158	GLU	2.3
2	B	106	LEU	2.3
2	D	211	LEU	2.2
2	B	56	VAL	2.2
1	C	16	TYR	2.2
2	D	148	GLN	2.2
2	D	33	PRO	2.2
2	D	104	PHE	2.1
2	B	63	TYR	2.1
1	C	269	TRP	2.1
2	D	106	LEU	2.1
2	D	219	LEU	2.1
2	B	55	SER	2.1
1	C	292	LEU	2.1
2	B	71	GLN	2.1
2	B	250	LEU	2.0
1	C	47	PHE	2.0
2	D	78	LEU	2.0
1	C	395	PHE	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](#)

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( $\text{\AA}^2$ )	Q<0.9
4	MG	A	1413	1/1	0.95	0.37	2.07	39,39,39,39	0
5	UTP	A	1414	29/29	0.89	0.36	1.28	28,52,71,71	40
5	UTP	C	1412	29/29	0.84	0.34	1.18	88,124,151,158	40
3	ZN	A	1411	1/1	0.99	0.32	1.00	40,40,40,40	0
4	MG	C	1411	1/1	0.95	0.24	-0.10	95,95,95,95	0
3	ZN	C	1409	1/1	0.98	0.22	-0.47	100,100,100,100	0
4	MG	C	1410	1/1	0.69	0.74	-	90,90,90,90	0
4	MG	A	1412	1/1	0.79	0.55	-	42,42,42,42	0

## 6.5 Other polymers

There are no such residues in this entry.