



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:38 AM GMT

PDB ID : 3PKY  
Title : Polymerase Domain from Mycobacterium tuberculosis Ligase D in complex with DNA, UTP and Manganese.  
Authors : Brissett, N.C.; Fox, G.C.; Pitcher, R.S.; Doherty, A.J.  
Deposited on : 2010-11-12  
Resolution : 3.10 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  
<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

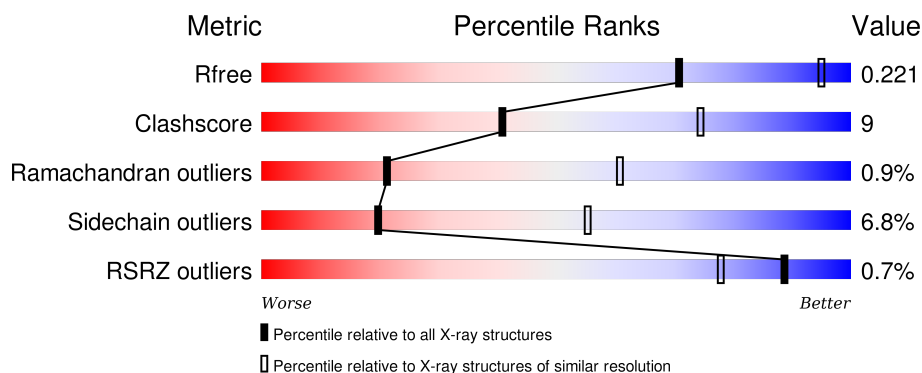
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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))
$R_{free}$	91344	1114 (3.14-3.06)
Clashscore	102246	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)
RSRZ outliers	91569	1119 (3.14-3.06)

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	303	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 78%, yellow 15%, orange 5%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>78%</span> <span>15%</span> <span>• 5%</span> </div> </div>
1	B	303	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, green 75%, yellow 17%, orange 6%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>%</span> <span>75%</span> <span>17%</span> <span>• 6%</span> </div> </div>
2	C	5	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 40%, yellow 60%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span></span> <span>40%</span> <span>60%</span> </div> </div>
3	D	13	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 8%, yellow 23%, orange 15%, grey 54%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>8%</span> <span>23%</span> <span>15%</span> <span>54%</span> </div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 4695 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 Putative DNA ligase-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	23	0	0
			2193	1385	390	413	5			
1	B	286	Total	C	N	O	S	25	0	0
			2184	1380	389	410	5			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP P71571
A	-1	SER	-	EXPRESSION TAG	UNP P71571
A	0	HIS	-	EXPRESSION TAG	UNP P71571
B	-2	GLY	-	EXPRESSION TAG	UNP P71571
B	-1	SER	-	EXPRESSION TAG	UNP P71571
B	0	HIS	-	EXPRESSION TAG	UNP P71571

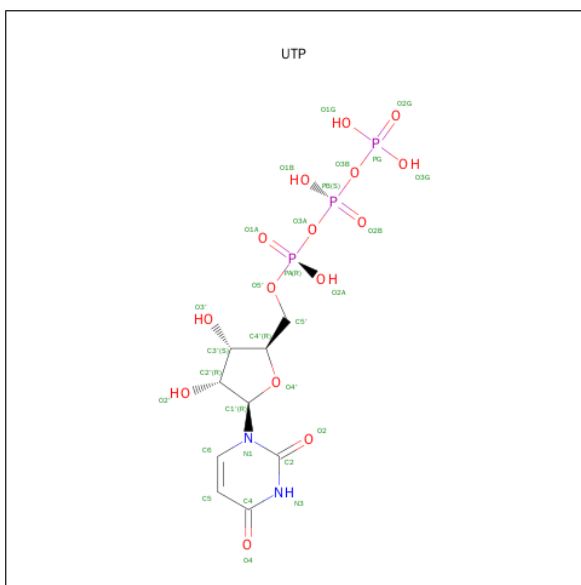
- Molecule 2 is a DNA chain called DNA 5'-D(P\*GP\*CP\*GP\*GP\*C)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	5	Total	C	N	O	P	0	0	0
			105	48	21	31	5			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	6	Total	C	N	O	P	0	0	1
			104	49	21	29	5			

- Molecule 4 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
4	A	1	Total	C	N	O	P	0	0
			29	9	2	15	3		
4	B	1	Total	C	N	O	P	0	0
			29	9	2	15	3		

- Molecule 5 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Mn	0	0
			2	2		
5	A	2	Total	Mn	0	0
			2	2		

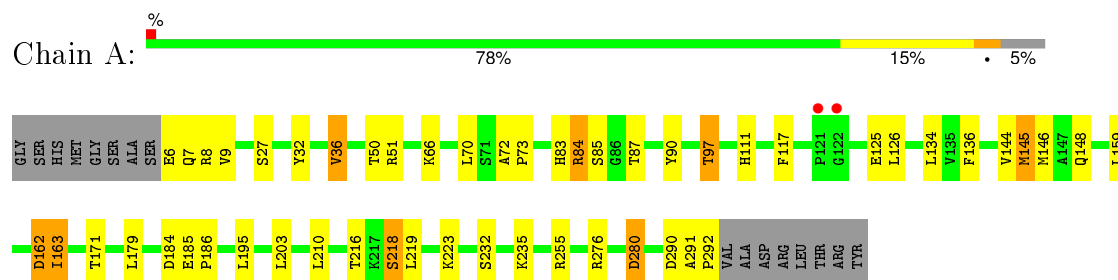
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	27	Total	O	0	0
			27	27		
6	B	17	Total	O	0	0
			17	17		
6	C	1	Total	O	0	0
			1	1		
6	D	2	Total	O	0	0
			2	2		

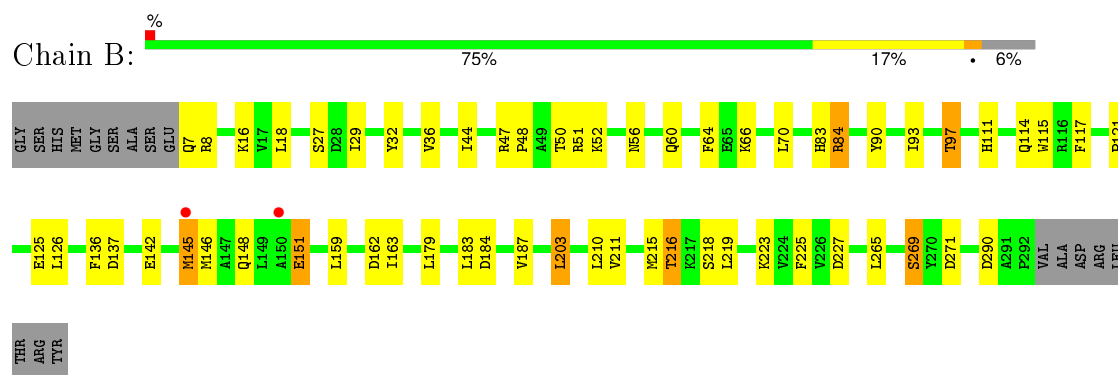
### 3 Residue-property plots [i](#)

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.

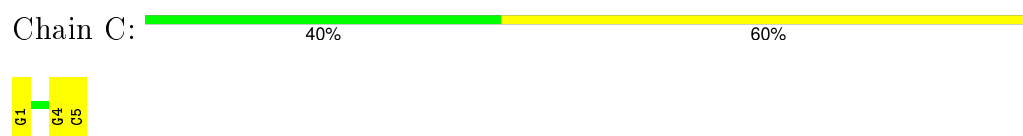
- Molecule 1: Putative DNA ligase-like protein



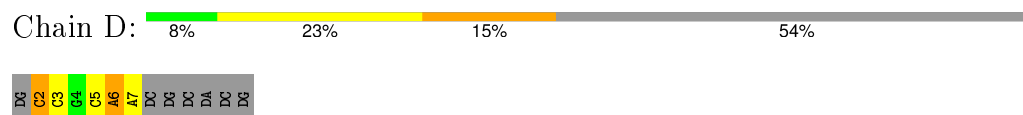
- Molecule 1: Putative DNA ligase-like protein



- Molecule 2: DNA 5'-D(P\*GP\*CP\*GP\*GP\*C)-3'



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



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	145.71Å 145.71Å 44.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.57 – 3.10 48.57 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.57-3.10) 99.3 (48.57-3.10)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	0.16	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.59 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, $R_{free}$	0.191 , 0.248 0.196 , 0.221	Depositor DCC
$R_{free}$ test set	894 reflections (5.37%)	DCC
Wilson B-factor (Å <sup>2</sup> )	78.5	Xtriage
Anisotropy	0.200	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 33.8	EDS
Estimated twinning fraction	0.458 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 17577 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4695	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% 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, MN

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	0.45	0/2244	0.61	0/3066
1	B	0.44	0/2235	0.61	0/3054
2	C	1.43	1/117 (0.9%)	1.76	3/177 (1.7%)
3	D	1.42	2/116 (1.7%)	1.78	2/177 (1.1%)
All	All	0.54	3/4712 (0.1%)	0.73	5/6474 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	DG	OP3-P	-11.14	1.47	1.61
3	D	2	DC	C3'-O3'	-5.83	1.36	1.44
3	D	3	DC	P-O5'	-5.32	1.54	1.59

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	DG	OP1-P-OP2	-9.79	104.92	119.60
2	C	1	DG	O4'-C1'-N9	-7.13	103.01	108.00
2	C	1	DG	O4'-C4'-C3'	-5.69	102.22	104.50
3	D	6	DA	P-O3'-C3'	5.61	126.43	119.70
3	D	5	DC	N3-C2-O2	-5.41	118.11	121.90

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	2193	0	2194	29	0
1	B	2184	0	2188	43	0
2	C	105	0	56	4	0
3	D	104	0	56	6	0
4	A	29	0	11	2	0
4	B	29	0	11	0	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	27	0	0	0	0
6	B	17	0	0	0	0
6	C	1	0	0	0	0
6	D	2	0	0	0	0
All	All	4695	0	4516	76	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 9.

All (76) 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:B:84:ARG:HB2	2:C:5:DC:H42	1.21	1.02
1:B:84:ARG:HB2	2:C:5:DC:N4	1.83	0.92
4:A:301:UTP:H5'1	4:A:301:UTP:H6	1.40	0.87
1:B:159:LEU:HD12	1:B:203:LEU:HD12	1.64	0.78
1:B:16:LYS:NZ	3:D:2:DC:C3'	2.48	0.77
1:A:66:LYS:HD3	1:A:90:TYR:OH	1.88	0.73
3:D:6:DA:H2''	3:D:7:DA:H5''	1.72	0.72
1:A:216:THR:HG22	1:A:218:SER:O	1.93	0.66
2:C:4:DG:H2''	2:C:5:DC:H4'	1.78	0.65
1:B:145:MET:HE2	1:B:145:MET:H	1.60	0.65
1:B:145:MET:HG2	1:B:148:GLN:HB3	1.77	0.65
1:B:16:LYS:HZ2	3:D:2:DC:C3'	2.10	0.64
1:B:148:GLN:HE21	1:B:210:LEU:HB3	1.62	0.64
1:B:16:LYS:HZ3	3:D:2:DC:C3'	2.11	0.63
1:B:183:LEU:HD13	1:B:187:VAL:HG12	1.81	0.62
1:A:144:VAL:HA	1:A:145:MET:HE1	1.83	0.60
1:B:203:LEU:HD23	1:B:211:VAL:HG21	1.83	0.60
1:B:162:ASP:O	1:B:163:ILE:HG13	2.03	0.58
1:B:216:THR:CG2	1:B:218:SER:O	2.51	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:ARG:CB	2:C:5:DC:N4	2.62	0.57
1:B:269:SER:HB2	1:B:271:ASP:OD1	2.05	0.57
1:A:117:PHE:HB3	1:A:126:LEU:HB3	1.87	0.56
1:A:145:MET:CE	1:A:148:GLN:HB3	2.36	0.56
1:B:145:MET:CE	1:B:145:MET:H	2.20	0.54
1:A:291:ALA:HB1	1:A:292:PRO:HD2	1.91	0.53
4:A:301:UTP:H3	3:D:6:DA:H61	1.56	0.52
1:A:159:LEU:HD12	1:A:203:LEU:HD12	1.91	0.52
1:A:145:MET:CE	1:A:145:MET:H	2.23	0.52
1:B:32:TYR:O	1:B:36:VAL:HG22	2.10	0.51
1:B:7:GLN:O	1:B:97:THR:HG21	2.10	0.51
1:B:216:THR:HG22	1:B:218:SER:O	2.11	0.50
1:B:163:ILE:HG22	1:B:163:ILE:O	2.11	0.50
1:B:50:THR:HB	1:B:111:HIS:HB2	1.94	0.50
1:B:162:ASP:C	1:B:163:ILE:HG13	2.32	0.50
1:B:44:ILE:HD12	1:B:93:ILE:HB	1.94	0.49
1:A:171:THR:HG23	1:A:255:ARG:HG3	1.95	0.48
1:A:185:GLU:OE2	1:A:186:PRO:HD2	2.13	0.48
1:B:52:LYS:HD2	1:B:64:PHE:CE2	2.49	0.48
1:B:52:LYS:HD2	1:B:64:PHE:CZ	2.48	0.48
1:A:32:TYR:O	1:A:36:VAL:HG22	2.13	0.48
1:A:145:MET:CE	1:A:145:MET:N	2.78	0.47
1:A:159:LEU:HD12	1:A:203:LEU:CD1	2.44	0.47
1:B:66:LYS:HD3	1:B:90:TYR:OH	2.15	0.47
1:A:8:ARG:HG3	1:A:9:VAL:N	2.30	0.47
1:A:162:ASP:C	1:A:163:ILE:HG13	2.34	0.47
1:A:276:ARG:O	1:A:280:ASP:HB2	2.15	0.46
1:A:50:THR:HB	1:A:111:HIS:HB2	1.98	0.46
1:B:48:PRO:HG3	1:B:115:TRP:CH2	2.50	0.45
1:A:83:HIS:O	1:A:84:ARG:C	2.54	0.45
1:B:148:GLN:HA	1:B:151:GLU:HB2	1.98	0.45
1:A:84:ARG:HB2	1:A:85:SER:H	1.57	0.45
1:B:146:MET:SD	1:B:265:LEU:HD23	2.57	0.45
1:B:136:PHE:HB2	1:B:179:LEU:HB2	1.99	0.45
1:B:216:THR:HG22	1:B:218:SER:C	2.37	0.45
1:B:137:ASP:HB3	1:B:227:ASP:HB3	1.98	0.45
1:B:16:LYS:NZ	3:D:2:DC:O3'	2.50	0.43
1:B:56:ASN:HB3	1:B:60:GLN:HG3	2.01	0.43
1:B:83:HIS:O	1:B:84:ARG:C	2.55	0.43
1:B:117:PHE:HB3	1:B:126:LEU:HB3	2.00	0.43
1:A:145:MET:HE3	1:A:148:GLN:HB3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:LEU:HD12	1:B:29:ILE:HD12	2.01	0.42
1:B:145:MET:HG2	1:B:148:GLN:CB	2.48	0.42
1:A:145:MET:HE2	1:A:145:MET:N	2.35	0.42
1:B:7:GLN:HB3	1:B:97:THR:HG21	2.00	0.42
1:A:134:LEU:HD11	1:A:195:LEU:HD23	2.02	0.42
1:B:47:ARG:HH12	1:B:114:GLN:C	2.22	0.42
1:A:216:THR:CG2	1:A:218:SER:O	2.65	0.41
1:A:7:GLN:O	1:A:97:THR:HG21	2.20	0.41
1:B:47:ARG:HG2	1:B:115:TRP:O	2.21	0.41
1:B:51:ARG:N	1:B:51:ARG:HD3	2.35	0.41
1:A:159:LEU:CD1	1:A:203:LEU:HD12	2.50	0.41
1:A:51:ARG:HD3	1:A:51:ARG:N	2.35	0.41
1:A:136:PHE:HB2	1:A:179:LEU:HB2	2.02	0.41
1:B:215:MET:HE2	1:B:225:PHE:HZ	1.86	0.41
1:A:232:SER:HB3	1:A:235:LYS:HG3	2.03	0.40
1:A:72:ALA:HB1	1:A:73:PRO:HD2	2.03	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	285/303 (94%)	266 (93%)	16 (6%)	3 (1%)	17	55
1	B	284/303 (94%)	266 (94%)	16 (6%)	2 (1%)	26	65
All	All	569/606 (94%)	532 (94%)	32 (6%)	5 (1%)	21	61

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	84	ARG
1	B	84	ARG

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Mol	Chain	Res	Type
1	A	146	MET
1	A	162	ASP
1	B	121	PRO

### 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	230/242 (95%)	214 (93%)	16 (7%)	19	54
1	B	229/242 (95%)	214 (93%)	15 (7%)	21	56
All	All	459/484 (95%)	428 (93%)	31 (7%)	20	55

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLU
1	A	27	SER
1	A	36	VAL
1	A	70	LEU
1	A	87	THR
1	A	97	THR
1	A	125	GLU
1	A	145	MET
1	A	163	ILE
1	A	184	ASP
1	A	210	LEU
1	A	218	SER
1	A	219	LEU
1	A	223	LYS
1	A	280	ASP
1	A	290	ASP
1	B	8	ARG
1	B	27	SER
1	B	70	LEU
1	B	97	THR

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Mol	Chain	Res	Type
1	B	125	GLU
1	B	142	GLU
1	B	145	MET
1	B	151	GLU
1	B	184	ASP
1	B	203	LEU
1	B	216	THR
1	B	219	LEU
1	B	223	LYS
1	B	269	SER
1	B	290	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	105	GLN
1	B	7	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 4 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	$\# Z  > 2$
4	UTP	A	301	5	20,30,30	0.73	0	30,47,47	1.87	4 (13%)
4	UTP	B	301	5	20,30,30	0.76	1 (5%)	30,47,47	1.78	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
4	UTP	A	301	5	-	0/18/38/38	0/2/2/2
4	UTP	B	301	5	-	0/18/38/38	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	301	UTP	O4'-C1'	2.35	1.44	1.41

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	301	UTP	PB-O3A-PA	-4.09	121.24	132.73
4	B	301	UTP	PB-O3B-PG	-3.79	119.95	132.67
4	A	301	UTP	PB-O3B-PG	-3.50	120.93	132.67
4	B	301	UTP	PB-O3A-PA	-2.44	125.87	132.73
4	B	301	UTP	O3G-PG-O1G	2.14	115.53	107.38
4	A	301	UTP	O4'-C1'-N1	2.21	112.74	108.08
4	B	301	UTP	O4'-C1'-N1	2.22	112.76	108.08
4	A	301	UTP	C4-N3-C2	6.95	121.02	114.14
4	B	301	UTP	C4-N3-C2	7.21	121.29	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	301	UTP	2	0

## 5.7 Other polymers

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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	287/303 (94%)	-0.00	2 (0%) 89 78	51, 70, 92, 115	17 (5%)
1	B	286/303 (94%)	-0.07	2 (0%) 89 78	51, 69, 93, 118	21 (7%)
2	C	5/5 (100%)	-0.20	0 100 100	65, 68, 91, 134	0
3	D	6/13 (46%)	0.20	0 100 100	66, 72, 90, 132	0
All	All	584/624 (93%)	-0.03	4 (0%) 89 78	51, 70, 93, 134	38 (6%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	122	GLY	2.7
1	B	145	MET	2.6
1	A	121	PRO	2.2
1	B	150	ALA	2.1

### 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	UTP	B	301	29/29	0.96	0.17	-1.03	55,63,74,74	0
4	UTP	A	301	29/29	0.97	0.14	-1.47	53,62,70,72	0
5	MN	A	303	1/1	0.93	0.15	-	64,64,64,64	0
5	MN	B	303	1/1	0.97	0.06	-	66,66,66,66	0
5	MN	A	302	1/1	0.96	0.16	-	57,57,57,57	0
5	MN	B	302	1/1	0.98	0.11	-	61,61,61,61	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.