



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 23, 2017 – 08:57 PM EST

PDB ID : 5TPV  
Title : X-ray structure of WlaRA (TDP-fucose-3,4-ketoisomerase) from *Campylobacter jejuni*  
Authors : Holden, H.M.; Thoden, J.B.; Li, Z.A.; Riegert, A.S.; Goneau, M.-F.; Cunningham, A.M.; Vinograd, E.; Schoenhofen, I.C.; Gilbert, M.; Li, J.  
Deposited on : 2016-10-21  
Resolution : 2.14 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.

---

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
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) : rb-20028442

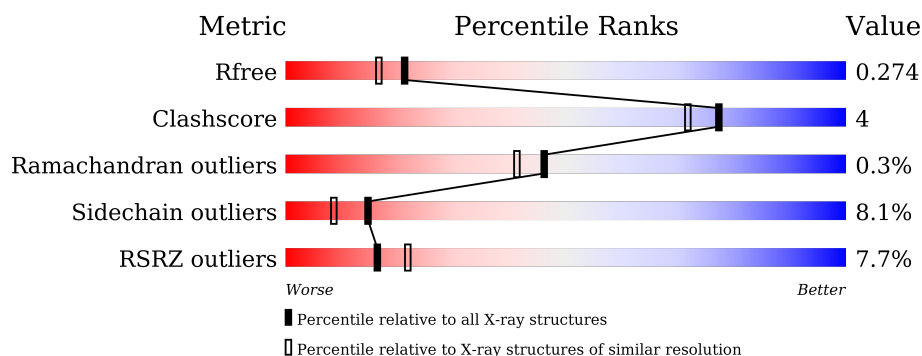
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.14 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	1693 (2.16-2.12)
Clashscore	102246	1824 (2.16-2.12)
Ramachandran outliers	100387	1798 (2.16-2.12)
Sidechain outliers	100360	1798 (2.16-2.12)
RSRZ outliers	91569	1699 (2.16-2.12)

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	153	<div> <div>9%</div> <div> <div></div> <div>70%</div> <div>16%</div> <div>•</div> <div>12%</div> </div> </div>
1	B	153	<div> <div>5%</div> <div> <div></div> <div>75%</div> <div>14%</div> <div>•</div> <div>10%</div> </div> </div>
1	C	153	<div> <div>7%</div> <div> <div></div> <div>71%</div> <div>14%</div> <div>•</div> <div>14%</div> </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
2	TYD	C	201	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3707 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 WlaRA, TDP-fucose-3,4-ketoisomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	135	Total	C	N	O	S	0	1	0
			1143	746	180	210	7			
1	B	138	Total	C	N	O	S	0	1	0
			1162	756	186	213	7			
1	C	131	Total	C	N	O	S	0	1	0
			1106	720	174	205	7			

There are 24 discrepancies between the modelled and reference sequences:

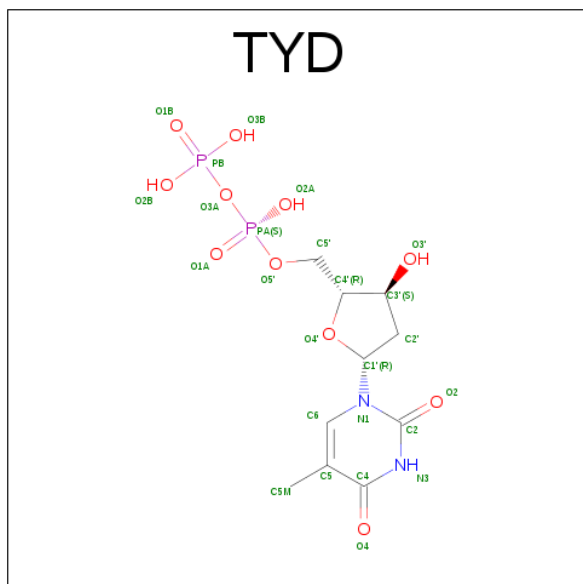
Chain	Residue	Modelled	Actual	Comment	Reference
A	146	LEU	-	expression tag	UNP Q9ALS3
A	147	GLU	-	expression tag	UNP Q9ALS3
A	148	HIS	-	expression tag	UNP Q9ALS3
A	149	HIS	-	expression tag	UNP Q9ALS3
A	150	HIS	-	expression tag	UNP Q9ALS3
A	151	HIS	-	expression tag	UNP Q9ALS3
A	152	HIS	-	expression tag	UNP Q9ALS3
A	153	HIS	-	expression tag	UNP Q9ALS3
B	146	LEU	-	expression tag	UNP Q9ALS3
B	147	GLU	-	expression tag	UNP Q9ALS3
B	148	HIS	-	expression tag	UNP Q9ALS3
B	149	HIS	-	expression tag	UNP Q9ALS3
B	150	HIS	-	expression tag	UNP Q9ALS3
B	151	HIS	-	expression tag	UNP Q9ALS3
B	152	HIS	-	expression tag	UNP Q9ALS3
B	153	HIS	-	expression tag	UNP Q9ALS3
C	146	LEU	-	expression tag	UNP Q9ALS3
C	147	GLU	-	expression tag	UNP Q9ALS3
C	148	HIS	-	expression tag	UNP Q9ALS3
C	149	HIS	-	expression tag	UNP Q9ALS3
C	150	HIS	-	expression tag	UNP Q9ALS3
C	151	HIS	-	expression tag	UNP Q9ALS3
C	152	HIS	-	expression tag	UNP Q9ALS3

*Continued on next page...*

*Continued from previous page...*

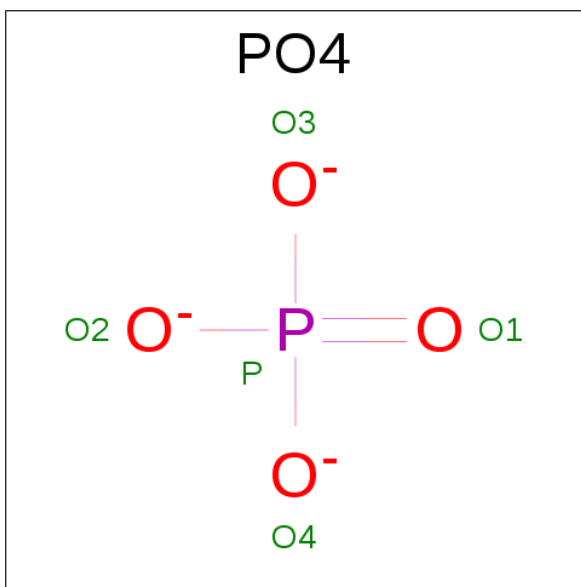
Chain	Residue	Modelled	Actual	Comment	Reference
C	153	HIS	-	expression tag	UNP Q9ALS3

- Molecule 2 is THYMIDINE-5'-DIPHOSPHATE (three-letter code: TYD) (formula:  $C_{10}H_{16}N_2O_{11}P_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
2	B	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
2	C	1	Total	C	N	O	P	0	0
			25	10	2	11	2		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula:  $O_4P$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	86	Total	O	0	0
			86	86		
4	B	84	Total	O	0	0
			84	84		
4	C	41	Total	O	0	0
			41	41		



- Molecule 1: WlaRA, TDP-fucose-3,4-ketoisomerase



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	177.95Å 177.95Å 88.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	125.83 – 2.14 26.52 – 2.14	Depositor EDS
% Data completeness (in resolution range)	99.1 (125.83-2.14) 99.2 (26.52-2.14)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.76 (at 2.13Å)	Xtriage
Refinement program	REFMAC 5.8.0124	Depositor
R, $R_{free}$	0.207 , 0.263 0.217 , 0.274	Depositor DCC
$R_{free}$ test set	1879 reflections (5.07%)	DCC
Wilson B-factor (Å <sup>2</sup> )	37.9	Xtriage
Anisotropy	0.055	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 44.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.53$ , $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3707	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: TYD, PO4

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	0.73	0/1173	1.10	7/1580 (0.4%)
1	B	0.68	0/1192	1.05	5/1604 (0.3%)
1	C	0.65	0/1134	1.03	3/1525 (0.2%)
All	All	0.69	0/3499	1.06	15/4709 (0.3%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	33	ARG	NE-CZ-NH1	-9.48	115.56	120.30
1	A	70	ASP	CB-CG-OD2	7.13	124.72	118.30
1	B	115	ASP	CB-CG-OD2	-7.07	111.94	118.30
1	A	90	ASP	CB-CG-OD2	7.07	124.66	118.30
1	C	90	ASP	CB-CG-OD2	6.77	124.40	118.30
1	C	103	ASP	CB-CG-OD2	6.54	124.19	118.30
1	B	115	ASP	CB-CG-OD1	5.90	123.61	118.30
1	C	33	ARG	NE-CZ-NH2	5.79	123.20	120.30
1	A	70	ASP	CB-CG-OD1	-5.34	113.49	118.30
1	B	33	ARG	NE-CZ-NH2	5.27	122.93	120.30
1	B	34	ILE	CA-CB-CG2	5.26	121.42	110.90
1	B	46	ARG	NE-CZ-NH2	5.23	122.91	120.30
1	A	39	ASP	CB-CG-OD1	5.20	122.98	118.30
1	A	34	ILE	CA-CB-CG2	5.17	121.25	110.90
1	A	33	ARG	NE-CZ-NH2	5.03	122.81	120.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts ⓘ

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	1143	0	1144	14	0
1	B	1162	0	1163	11	0
1	C	1106	0	1108	5	1
2	A	25	0	13	2	0
2	B	25	0	13	1	0
2	C	25	0	13	2	0
3	B	10	0	0	0	0
4	A	86	0	0	1	1
4	B	84	0	0	0	0
4	C	41	0	0	0	0
All	All	3707	0	3454	26	2

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 (26) 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:43:ASP:N	1:B:43:ASP:OD1	2.17	0.78
1:A:12:ASN:ND2	4:A:301:HOH:O	2.31	0.63
1:B:68:LEU:HD23	1:B:68:LEU:C	2.21	0.60
2:B:201:TYD:O5'	2:B:201:TYD:H6	2.01	0.59
1:A:113:TYR:OH	1:B:30:GLU:OE2	2.15	0.58
1:A:33:ARG:O	1:A:34:ILE:HD12	2.04	0.58
1:B:82:ASN:ND2	1:B:84:THR:OG1	2.36	0.56
2:A:201:TYD:O5'	2:A:201:TYD:H6	2.05	0.56
2:C:201:TYD:H6	2:C:201:TYD:O5'	2.07	0.55
2:C:201:TYD:H5'2	2:C:201:TYD:PB	2.47	0.54
1:A:30:GLU:OE2	1:B:113:TYR:OH	2.25	0.50
1:A:36:TYR:HA	1:A:106:LEU:O	2.15	0.46
1:A:33:ARG:HD3	1:B:22:GLN:OE1	2.17	0.45
1:A:22:GLN:OE1	1:B:33:ARG:HD3	2.17	0.45
1:C:45:ILE:H	1:C:45:ILE:HD12	1.81	0.45
1:A:117:LYS:HB3	1:C:11:ILE:HG23	1.99	0.44
1:A:10:THR:HG22	1:A:19:ASN:OD1	2.18	0.44

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:29:PHE:CD1	1:C:111:ASN:HB3	2.54	0.43
2:A:201:TYD:H1'	1:B:11:ILE:HD13	2.01	0.43
1:A:41:LEU:HB2	1:B:15:ASN:HB3	2.01	0.42
1:A:125:TYR:CE2	1:C:15:ASN:HA	2.54	0.42
1:B:36:TYR:HA	1:B:106:LEU:O	2.20	0.41
1:C:1:MET:HE3	1:C:71:ASP:HB3	2.01	0.41
1:A:1:MET:HE3	1:A:71:ASP:HB3	2.02	0.41
1:A:37:ILE:HA	1:B:17:ILE:O	2.21	0.41

All (2) 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)
4:A:353:HOH:O	4:A:353:HOH:O[7_555]	1.85	0.35
1:C:30:GLU:OE2	1:C:113:TYR:OH[5_455]	2.00	0.20

## 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	134/153 (88%)	128 (96%)	6 (4%)	0	100	100
1	B	137/153 (90%)	133 (97%)	4 (3%)	0	100	100
1	C	130/153 (85%)	126 (97%)	3 (2%)	1 (1%)	24	14
All	All	401/459 (87%)	387 (96%)	13 (3%)	1 (0%)	46	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	130	LYS

### 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	130/144 (90%)	120 (92%)	10 (8%)	16	9
1	B	131/144 (91%)	121 (92%)	10 (8%)	16	10
1	C	126/144 (88%)	115 (91%)	11 (9%)	13	7
All	All	387/432 (90%)	356 (92%)	31 (8%)	15	9

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

Mol	Chain	Res	Type
1	A	1	MET
1	A	32	LYS
1	A	41	LEU
1	A	42	ASP
1	A	85	LYS
1	A	91	LYS
1	A	95	LYS
1	A	116	GLU
1	A	133	ILE
1	A	135	TRP
1	B	9	LYS
1	B	32	LYS
1	B	34	ILE
1	B	42	ASP
1	B	68	LEU
1	B	81	ASN
1	B	85	LYS
1	B	95	LYS
1	B	102	LYS
1	B	118	GLU
1	C	1	MET
1	C	2	ASN
1	C	41	LEU
1	C	68	LEU
1	C	77	LYS
1	C	81	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	C	82	ASN
1	C	91	LYS
1	C	95	LYS
1	C	124	LYS
1	C	131	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	ASN
1	B	81	ASN
1	B	82	ASN
1	B	132	ASN
1	C	81	ASN
1	C	96	GLN
1	C	99	ASN
1	C	131	ASN

### 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](#)

5 ligands 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	$\# Z  > 2$
2	TYD	A	201	-	23,26,26	2.44	3 (13%)	26,40,40	2.07	5 (19%)
2	TYD	B	201	-	23,26,26	2.45	3 (13%)	26,40,40	2.07	5 (19%)
3	PO4	B	202	-	4,4,4	0.59	0	6,6,6	0.37	0
3	PO4	B	203	-	4,4,4	0.63	0	6,6,6	0.24	0
2	TYD	C	201	-	23,26,26	2.45	3 (13%)	26,40,40	2.07	5 (19%)

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	TYD	A	201	-	-	0/16/28/28	0/2/2/2
2	TYD	B	201	-	-	0/16/28/28	0/2/2/2
3	PO4	B	202	-	-	0/0/0/0	0/0/0/0
3	PO4	B	203	-	-	0/0/0/0	0/0/0/0
2	TYD	C	201	-	-	0/16/28/28	0/2/2/2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	201	TYD	C6-N1	-9.59	1.34	1.46
2	B	201	TYD	C6-N1	-9.56	1.34	1.46
2	A	201	TYD	C6-N1	-9.54	1.34	1.46
2	B	201	TYD	C6-C5	-3.97	1.39	1.51
2	C	201	TYD	C6-C5	-3.97	1.39	1.51
2	A	201	TYD	C6-C5	-3.95	1.39	1.51
2	B	201	TYD	C2-N1	3.99	1.41	1.35
2	A	201	TYD	C2-N1	4.00	1.41	1.35
2	C	201	TYD	C2-N1	4.03	1.41	1.35

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	201	TYD	C4-N3-C2	-4.78	120.64	126.81
2	C	201	TYD	C4-N3-C2	-4.76	120.65	126.81
2	B	201	TYD	C4-N3-C2	-4.72	120.70	126.81
2	B	201	TYD	C2'-C1'-N1	-4.70	109.72	115.65

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	201	TYD	C2'-C1'-N1	-4.69	109.74	115.65
2	C	201	TYD	C2'-C1'-N1	-4.68	109.75	115.65
2	C	201	TYD	O2-C2-N1	-2.93	119.33	123.17
2	A	201	TYD	O2-C2-N1	-2.92	119.34	123.17
2	B	201	TYD	O2-C2-N1	-2.90	119.37	123.17
2	A	201	TYD	C5M-C5-C6	3.67	120.66	112.52
2	C	201	TYD	C5M-C5-C6	3.69	120.70	112.52
2	B	201	TYD	C5M-C5-C6	3.70	120.72	112.52
2	B	201	TYD	N3-C2-N1	5.05	121.32	116.64
2	A	201	TYD	N3-C2-N1	5.06	121.33	116.64
2	C	201	TYD	N3-C2-N1	5.07	121.33	116.64

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	201	TYD	2	0
2	B	201	TYD	1	0
2	C	201	TYD	2	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	135/153 (88%)	0.10	14 (10%) 8 12	25, 38, 66, 100	0
1	B	138/153 (90%)	-0.10	7 (5%) 32 41	27, 40, 69, 98	0
1	C	131/153 (85%)	0.15	10 (7%) 17 22	34, 48, 74, 95	0
All	All	404/459 (88%)	0.05	31 (7%) 16 22	25, 42, 73, 100	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	133	ILE	6.7
1	A	134	VAL	4.0
1	C	125	TYR	3.7
1	B	108	VAL	3.7
1	B	107	LEU	3.4
1	A	135	TRP	3.3
1	A	107	LEU	3.3
1	B	135	TRP	3.2
1	A	108	VAL	3.2
1	C	131	ASN	3.2
1	A	109	LEU	3.2
1	A	57	ILE	3.1
1	A	42	ASP	2.9
1	B	42	ASP	2.9
1	B	138	GLY	2.9
1	A	106	LEU	2.9
1	C	127	CYS	2.8
1	C	95	LYS	2.7
1	A	43	ASP	2.7
1	C	107	LEU	2.4
1	A	35	PHE	2.4
1	C	102	LYS	2.4
1	C	128	GLU	2.4

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	C	129	LEU	2.3
1	A	58	PHE	2.2
1	B	61	LEU	2.1
1	C	35	PHE	2.1
1	A	59	ILE	2.0
1	A	41	LEU	2.0
1	B	34	ILE	2.0
1	C	43	ASP	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(Å <sup>2</sup> )	Q<0.9
2	TYD	C	201	25/25	0.77	0.22	2.35	40,69,123,134	0
2	TYD	A	201	25/25	0.90	0.12	-0.00	33,40,45,55	0
2	TYD	B	201	25/25	0.94	0.10	-0.26	30,34,37,39	0
3	PO4	B	203	5/5	0.83	0.29	-	75,80,86,102	0
3	PO4	B	202	5/5	0.88	0.19	-	58,67,72,96	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.