



Full wwPDB X-ray Structure Validation Report

Sep 23, 2014 – 06:41 PM EDT

PDB ID : 3M6L
Title : Crystal structure of transketolase in complex with thiamine diphosphate, ribose-5-phosphate and calcium ion
Authors : Nocek, B.; Makowska-Grzyska, M.; Maltseva, N.; Joachimiak, A.; Anderson, W.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2010-03-15
Resolution : 1.59 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

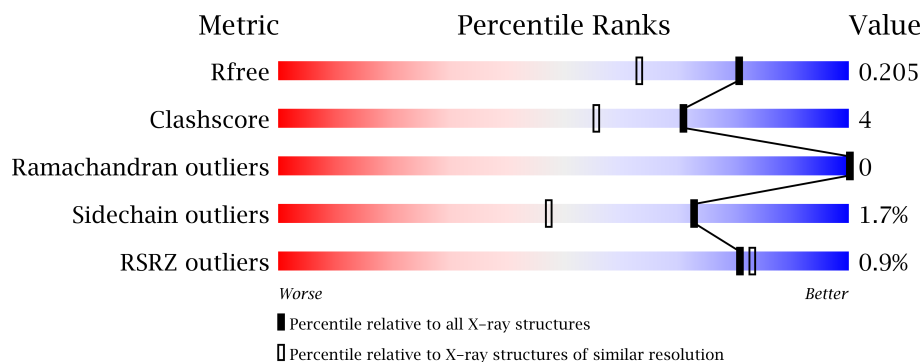
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable23489
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable23489

1 Overall quality at a glance

The reported resolution of this entry is 1.59 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}	66092	1872 (1.60-1.60)
Clashscore	79885	2199 (1.60-1.60)
Ramachandran outliers	78287	2126 (1.60-1.60)
Sidechain outliers	78261	2125 (1.60-1.60)
RSRZ outliers	66119	1872 (1.60-1.60)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	635	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
5	EDO	A	1002	-	X
6	R5P	A	701	-	X

2 Entry composition i

There are 8 unique types of molecules in this entry. The entry contains 5535 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

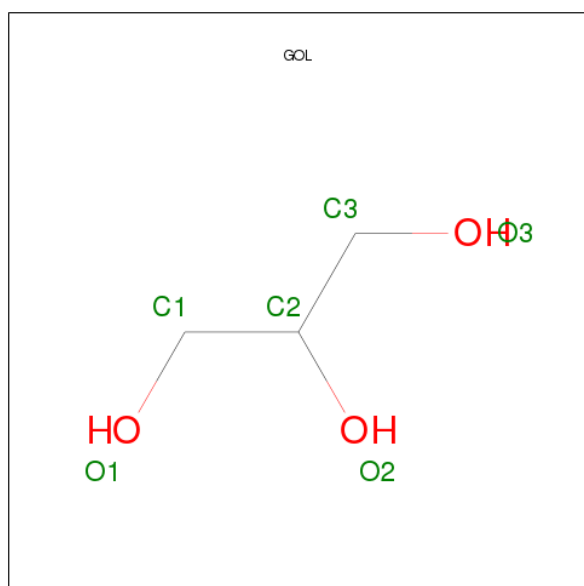
- Molecule 1 is a protein called Transketolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	632	4914	3142	831	923	8	10	3	3	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q0P7Y3
A	-1	ASN	-	expression tag	UNP Q0P7Y3
A	0	ALA	-	expression tag	UNP Q0P7Y3

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	6	3	3	0	0

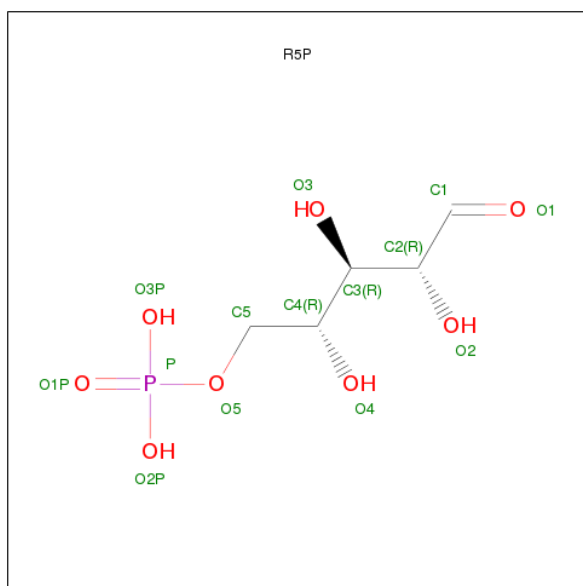
- TPP
-
- The chemical structure of TPP (Thiamine Pyrophosphate) is shown. It consists of a pyrophosphate group (two phosphorus atoms, P1 and P2, connected by an oxygen atom) and a thiazolium ring system. The pyrophosphate group is linked to the thiazolium ring via an oxygen atom (O7). The thiazolium ring is a five-membered ring containing a sulfur atom (S1) and a nitrogen atom (N3). The thiazolium ring is substituted with a methyl group (CM4) at C5, a methyl group (CM2) at C2', and an amino group (NH2) at C4'. The thiazolium ring is also linked to a pyrimidine ring via a methylene group (C7'). The pyrimidine ring is a six-membered ring containing two nitrogen atoms (N1' and N3'). The pyrimidine ring is substituted with a methyl group (CM2) at C2' and an amino group (NH2) at C4'.

EDO

The chemical structure of EDO (Ethane-1,2-diol) is shown. It consists of two carbon atoms (C1 and C2) connected by a single bond. Each carbon atom is bonded to a hydroxyl group (OH). The oxygen atoms are labeled O1 and O2, and the carbon atoms are labeled C1 and C2. The hydrogens are shown in red, and the oxygen and carbon atoms are shown in green.

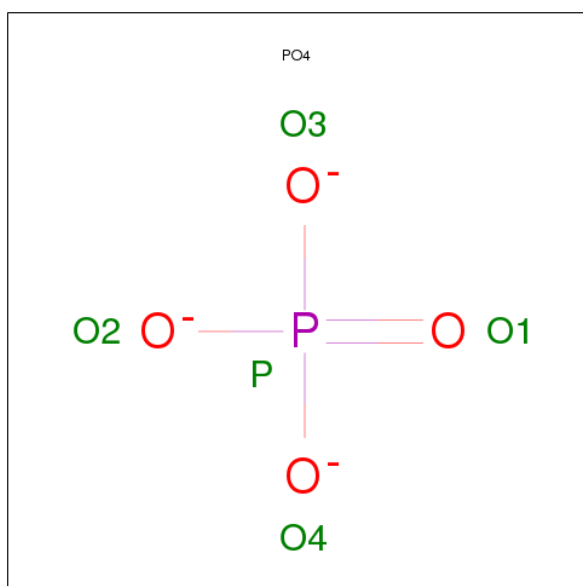
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is SUGAR (RIBOSE-5-PHOSPHATE) (three-letter code: R5P) (formula: $C_5H_{11}O_8P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	O	P	0	0
			14	5	8	1		

- Molecule 7 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	P	0	0
			5	4	1		

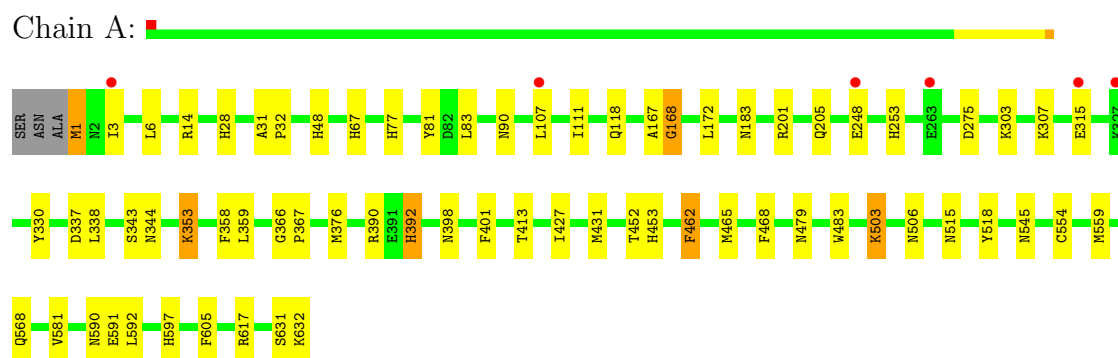
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	565	Total	O	0	0
			565	565		

3 Residue-property plots

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: Transketolase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	130.92Å 71.46Å 69.62Å 90.00° 109.48° 90.00°	Depositor
Resolution (Å)	40.00 – 1.59 35.66 – 1.59	Depositor EDS
% Data completeness (in resolution range)	94.3 (40.00-1.59) 94.3 (35.66-1.59)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 1.59Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.160 , 0.201 0.169 , 0.205	Depositor DCC
R_{free} test set	3821 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	17.4	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 38.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 76496 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5535	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: GOL, PO4, EDO, TPP, R5P, CA

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	1.29	8/5021 (0.2%)	1.10	14/6757 (0.2%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	554	CYS	CB-SG	-8.54	1.67	1.82
1	A	518	TYR	CD2-CE2	-5.92	1.30	1.39
1	A	462	PHE	CE1-CZ	5.68	1.48	1.37
1	A	591	GLU	CD-OE2	5.67	1.31	1.25
1	A	315	GLU	CB-CG	5.45	1.62	1.52
1	A	168	GLY	N-CA	5.05	1.53	1.46
1	A	343	SER	CB-OG	5.01	1.48	1.42
1	A	401	PHE	CE1-CZ	5.01	1.46	1.37

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	390	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	A	337	ASP	CB-CG-OD1	5.80	123.52	118.30
1	A	330	TYR	CB-CG-CD1	-5.71	117.57	121.00
1	A	275	ASP	CB-CG-OD1	5.64	123.38	118.30
1	A	605	PHE	CG-CD1-CE1	5.56	126.92	120.80
1	A	14	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	A	201	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	A	338	LEU	CB-CG-CD2	-5.41	101.81	111.00
1	A	617	ARG	NE-CZ-NH1	5.34	122.97	120.30
1	A	413	THR	CA-CB-CG2	-5.28	105.01	112.40
1	A	390	ARG	NE-CZ-NH2	-5.24	117.68	120.30
1	A	358	PHE	CB-CG-CD1	-5.16	117.19	120.80
1	A	605	PHE	CB-CG-CD2	5.07	124.34	120.80

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	359	LEU	CB-CG-CD2	5.03	119.56	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4914	0	4871	42	0
2	A	6	0	8	0	0
3	A	26	0	16	3	0
4	A	1	0	0	0	0
5	A	4	0	6	0	0
6	A	14	0	9	2	0
7	A	5	0	0	0	0
8	A	565	0	0	9	0
All	All	5535	0	4910	44	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (44) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:303:LYS:NZ	1:A:307:LYS:HZ1	1.63	0.97
1:A:1:MSE:HE2	1:A:6:LEU:HD11	1.47	0.95
1:A:581:VAL:H	1:A:597:HIS:HD2	1.15	0.89
1:A:303:LYS:NZ	1:A:307:LYS:NZ	2.27	0.83
1:A:48:HIS:HD2	1:A:81:TYR:OH	1.71	0.73
1:A:303:LYS:HZ2	1:A:307:LYS:NZ	1.89	0.69
1:A:90:ASN:ND2	8:A:941:HOH:O	2.24	0.68
1:A:118:GLN:HE22	1:A:392:HIS:CE1	2.13	0.66
1:A:303:LYS:HZ2	1:A:307:LYS:HZ1	1.38	0.64
1:A:303:LYS:HZ3	1:A:307:LYS:NZ	1.98	0.59
1:A:515:ASN:HD22	1:A:568:GLN:NE2	2.00	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:344:ASN:HD21	1:A:483:TRP:HE1	1.48	0.59
1:A:506:ASN:HB2	8:A:642:HOH:O	2.01	0.59
1:A:1:MSE:CE	1:A:6:LEU:HD11	2.30	0.57
1:A:344:ASN:ND2	1:A:483:TRP:HE1	2.03	0.57
1:A:353:LYS:HE2	8:A:1107:HOH:O	2.06	0.54
1:A:168:GLY:HA3	1:A:205:GLN:O	2.08	0.53
1:A:77:HIS:HD2	1:A:83:LEU:O	1.91	0.53
3:A:5001:TPP:H2	6:A:701:R5P:O1	2.09	0.52
1:A:167:ALA:HA	1:A:172:LEU:HD12	1.93	0.51
1:A:427:ILE:HG22	1:A:431:MSE:HE2	1.93	0.50
1:A:353:LYS:NZ	1:A:353:LYS:HB3	2.27	0.50
1:A:631:SER:O	1:A:632:LYS:CB	2.60	0.49
1:A:48:HIS:HE1	8:A:725:HOH:O	1.94	0.49
1:A:545:ASN:ND2	8:A:1023:HOH:O	2.47	0.47
1:A:67:HIS:HD2	8:A:857:HOH:O	1.96	0.47
1:A:67:HIS:HE1	3:A:5001:TPP:O1B	1.98	0.46
1:A:31:ALA:HB3	1:A:32:PRO:HD3	1.98	0.45
1:A:559:MSE:HE1	1:A:592:LEU:HD21	1.98	0.44
1:A:581:VAL:H	1:A:597:HIS:CD2	2.07	0.44
1:A:118:GLN:HE22	1:A:392:HIS:HE1	1.60	0.43
1:A:462:PHE:CE2	1:A:465:MSE:HE1	2.53	0.43
1:A:111:ILE:HG13	1:A:431:MSE:SE	2.69	0.42
3:A:5001:TPP:H2	6:A:701:R5P:C1	2.49	0.42
1:A:452:THR:OG1	1:A:453:HIS:HD2	2.03	0.42
1:A:28:HIS:CG	1:A:67:HIS:HB2	2.55	0.42
1:A:503:LYS:HB3	1:A:503:LYS:HE2	1.94	0.42
1:A:559:MSE:HE1	1:A:592:LEU:CD2	2.50	0.42
1:A:344:ASN:HB2	1:A:479:ASN:HD21	1.85	0.41
1:A:398:ASN:HA	1:A:398:ASN:HD22	1.68	0.41
1:A:253:HIS:HD2	8:A:988:HOH:O	2.03	0.41
1:A:366:GLY:N	1:A:367:PRO:CD	2.84	0.41
1:A:183[B]:ASN:ND2	8:A:948:HOH:O	2.54	0.40
1:A:590:ASN:ND2	8:A:1081:HOH:O	2.53	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	633/635 (100%)	617 (98%)	16 (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	521/514 (101%)	512 (98%)	9 (2%)	73	47

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	3	ILE
1	A	107	LEU
1	A	248	GLU
1	A	353	LYS
1	A	376	MSE
1	A	392	HIS
1	A	468	PHE
1	A	503	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	48	HIS
1	A	67	HIS
1	A	77	HIS
1	A	90	ASN
1	A	134	GLN
1	A	135	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	344	ASN
1	A	369	ASN
1	A	386	HIS
1	A	392	HIS
1	A	398	ASN
1	A	453	HIS
1	A	479	ASN
1	A	545	ASN
1	A	568	GLN
1	A	590	ASN
1	A	597	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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, 1 is monoatomic - leaving 5 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	EDO	A	1002	-	3,3,3	0.68	0	2,2,2	0.20	0
2	GOL	A	2001	-	5,5,5	0.50	0	5,5,5	0.62	0
3	TPP	A	5001	4	27,27,27	2.46	6 (22%)	40,40,40	2.39	13 (32%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	PO4	A	633	-	4,4,4	0.25	0	6,6,6	0.31	0
6	R5P	A	701	-	13,13,13	2.55	5 (38%)	18,18,18	3.05	8 (44%)

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	EDO	A	1002	-	-	0/1/1/1	0/0/0/0
2	GOL	A	2001	-	-	0/4/4/4	0/0/0/0
3	TPP	A	5001	4	-	0/17/17/17	0/2/2/2
7	PO4	A	633	-	-	0/0/0/0	0/0/0/0
6	R5P	A	701	-	-	0/15/16/16	0/0/0/0

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	5001	TPP	C5-C4	8.89	1.47	1.36
6	A	701	R5P	O1-C1	5.68	1.47	1.19
6	A	701	R5P	C5-C4	4.19	1.58	1.51
3	A	5001	TPP	C4-N3	-4.14	1.36	1.39
3	A	5001	TPP	C7'-N3	3.77	1.55	1.48
6	A	701	R5P	P-O1P	3.66	1.63	1.51
3	A	5001	TPP	C5'-C4'	3.26	1.50	1.42
3	A	5001	TPP	C7'-C5'	3.17	1.58	1.51
3	A	5001	TPP	C2-S1	2.84	1.77	1.68
6	A	701	R5P	C2-C1	2.37	1.54	1.50
6	A	701	R5P	P-O2P	2.19	1.62	1.54

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	5001	TPP	C2-S1-C5	-8.67	85.96	91.48
6	A	701	R5P	O3-C3-C2	-7.55	95.59	108.93
3	A	5001	TPP	S1-C2-N3	4.95	120.02	112.62
6	A	701	R5P	C3-C2-C1	4.85	118.49	111.81
6	A	701	R5P	O1-C1-C2	-4.66	111.79	125.33
3	A	5001	TPP	C5-C4-N3	4.60	116.77	107.57
3	A	5001	TPP	CM4-C4-C5	-4.38	117.94	129.06
6	A	701	R5P	C4-C3-C2	3.67	124.58	113.31
6	A	701	R5P	O3-C3-C4	-3.56	99.76	108.73

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	5001	TPP	C6-C5-C4	-3.46	123.35	127.33
6	A	701	R5P	O3P-P-O5	3.40	116.42	106.67
3	A	5001	TPP	O7-C7-C6	-2.96	96.95	109.09
6	A	701	R5P	O5-C5-C4	2.87	117.25	109.23
3	A	5001	TPP	C7'-N3-C2	2.47	130.60	125.20
3	A	5001	TPP	C6-C5-S1	2.42	125.67	122.28
3	A	5001	TPP	C5'-C7'-N3	-2.17	109.56	113.28
3	A	5001	TPP	C7-C6-C5	2.14	118.38	112.25
3	A	5001	TPP	CM4-C4-N3	2.14	125.26	122.53
3	A	5001	TPP	PA-O3A-PB	2.09	137.73	131.93
3	A	5001	TPP	O7-PA-O1A	-2.04	101.40	109.37
6	A	701	R5P	O4-C4-C5	2.01	114.25	110.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

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, 95th 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 ²)	Q<0.9
1	A	632/635 (99%)	-0.13	6 (0%) 81 83	9, 18, 31, 50	4 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	248	GLU	2.8
1	A	327	LYS	2.3
1	A	107	LEU	2.3
1	A	315	GLU	2.3
1	A	3	ILE	2.2
1	A	263	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 95th 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	EDO	A	1002	4/4	0.16	6.03	50,52,53,54	0
6	R5P	A	701	14/14	0.14	2.61	37,43,47,51	0
7	PO4	A	633	5/5	0.12	0.81	42,43,45,46	0
2	GOL	A	2001	6/6	0.10	0.69	22,36,40,46	0
3	TPP	A	5001	26/26	0.07	-0.06	11,21,34,36	0
4	CA	A	801	1/1	0.05	-1.53	16,16,16,16	0

6.5 Other polymers ⓘ

There are no such residues in this entry.