



# Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 12:35 am BST

PDB ID : 4JOM  
Title : Structure of E. coli Pol III 3mPHP mutant  
Authors : Barros, T.; Guenther, J.; Kelch, B.; Anaya, J.; Prabhakar, A.; O'Donnell, M.; Kuriyan, J.; Lamers, M.H.  
Deposited on : 2013-03-18  
Resolution : 2.90 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

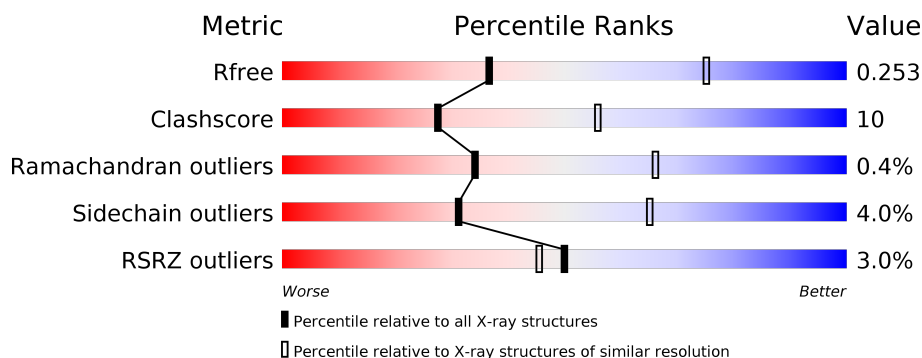
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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))
$R_{free}$	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 respectively. 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	918	<div> <div>3%</div> <div>77%</div> <div>20%</div> <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	PO4	A	1003	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 14296 atoms, of which 7107 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 III subunit alpha.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	909	Total	C	H	N	O	S	0	2	0
			14246	4547	7099	1221	1339	40			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	EXPRESSION TAG	UNP P10443
A	10	HIS	ARG	ENGINEERED MUTATION	UNP P10443
A	44	HIS	PHE	ENGINEERED MUTATION	UNP P10443
A	203	HIS	ARG	ENGINEERED MUTATION	UNP P10443

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



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

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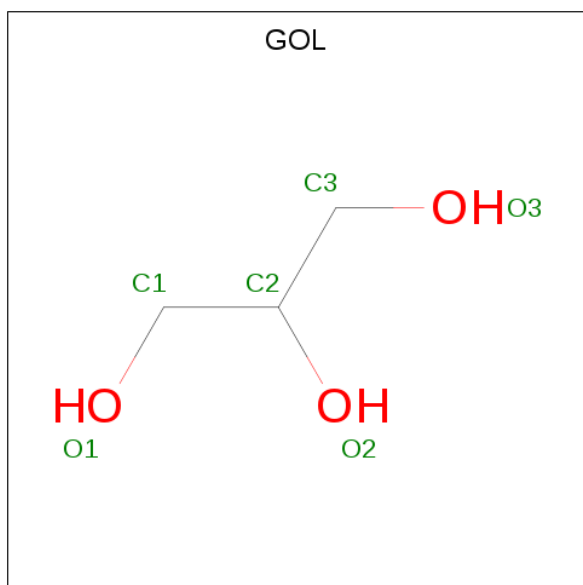
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	19	Total	O	0	0
			19	19		

### 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 the various outlier classes 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 III subunit alpha



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	82.43Å 98.94Å 139.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.95 – 2.90 46.94 – 2.89	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.95-2.90) 99.8 (46.94-2.89)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.89 (at 2.91Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, $R_{free}$	0.194 , 0.245 0.200 , 0.253	Depositor DCC
$R_{free}$ test set	1350 reflections (5.15%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	61.5	Xtriage
Anisotropy	0.549	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 48.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	14296	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ZN, 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	0.23	0/7303	0.43	1/9875 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	118	ARG	NE-CZ-NH1	5.51	123.06	120.30

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	7147	7099	7080	143	0
2	A	15	0	0	0	0
3	A	2	0	0	0	0
4	A	6	8	8	1	0
5	A	19	0	0	1	0
All	All	7189	7107	7088	143	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 10.

All (143) 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:A:3:GLU:N	1:A:3:GLU:OE1	2.04	0.89
1:A:225:LEU:O	1:A:230:ARG:NH1	2.11	0.83
1:A:444:ASP:OD2	1:A:506:ARG:NH1	2.20	0.74
1:A:216:ARG:NH2	1:A:505:THR:O	2.23	0.72
1:A:509:GLY:HA2	1:A:510:LYS:HB2	1.76	0.68
1:A:674:LYS:CG	1:A:675:PRO:HD3	2.26	0.66
1:A:706:ASP:OD1	1:A:709:ARG:NH1	2.28	0.65
1:A:872:LYS:HE2	1:A:872:LYS:HA	1.77	0.65
1:A:740:LEU:HD12	1:A:741:ALA:N	2.11	0.64
1:A:867:ALA:HA	1:A:909:ALA:HA	1.80	0.64
1:A:868:ARG:C	1:A:868:ARG:HD3	2.18	0.64
1:A:857:GLY:O	1:A:858:TYR:HB2	1.98	0.63
1:A:314:ARG:O	1:A:318:ARG:N	2.32	0.62
1:A:318:ARG:N	1:A:319:PRO:CD	2.64	0.61
1:A:509:GLY:CA	1:A:510:LYS:HB2	2.33	0.59
1:A:337:PHE:N	1:A:338:PRO:CD	2.65	0.58
1:A:310:ASP:O	1:A:312:GLU:N	2.36	0.58
1:A:287:SER:O	1:A:290:ASP:N	2.37	0.57
1:A:118:ARG:NH1	1:A:119:ASP:OD1	2.37	0.57
1:A:81:LEU:HD22	1:A:115:ILE:HD11	1.86	0.57
1:A:4:PRO:HG3	1:A:37:PRO:HG3	1.85	0.57
1:A:12:HIS:CE1	1:A:69:ASP:OD2	2.57	0.56
1:A:872:LYS:HA	1:A:872:LYS:CE	2.35	0.56
1:A:616:MET:CE	1:A:634:LEU:HB3	2.36	0.56
1:A:746:ASP:N	1:A:746:ASP:OD1	2.38	0.56
1:A:702:LEU:HA	1:A:705:ALA:HB3	1.88	0.56
1:A:858:TYR:CZ	1:A:868:ARG:HG3	2.40	0.56
1:A:674:LYS:HG3	1:A:675:PRO:HD3	1.88	0.55
1:A:581:GLU:HB2	1:A:582:PRO:HD2	1.88	0.55
1:A:118:ARG:HH11	1:A:118:ARG:CG	2.21	0.54
1:A:714:LYS:C	1:A:715:LYS:HG2	2.28	0.54
1:A:389:GLU:HG3	1:A:702:LEU:HB3	1.90	0.54
1:A:311:GLU:O	1:A:313:GLU:N	2.41	0.54
1:A:674:LYS:CB	1:A:675:PRO:HD3	2.38	0.53
1:A:661:ILE:HD12	1:A:663:TYR:CE1	2.45	0.52
1:A:143:LEU:HD12	1:A:151:VAL:HG22	1.92	0.52
1:A:606:THR:O	1:A:607:ALA:HB3	2.09	0.52
1:A:175:ARG:NH1	1:A:233:ASN:O	2.44	0.51
1:A:283:THR:HG21	1:A:286:MET:O	2.10	0.51
1:A:337:PHE:N	1:A:338:PRO:HD2	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:ALA:HB3	1:A:26:PRO:HD3	1.92	0.51
1:A:325:LEU:O	1:A:329:LEU:HB2	2.10	0.51
1:A:10:HIS:HB3	1:A:199:THR:HG23	1.92	0.51
1:A:9:LEU:HA	1:A:40:ALA:HB3	1.91	0.51
1:A:911:GLN:N	1:A:911:GLN:OE1	2.44	0.51
1:A:290:ASP:O	1:A:293:VAL:HG22	2.10	0.51
1:A:741:ALA:O	1:A:745:PHE:HB2	2.10	0.51
1:A:255:GLU:N	1:A:255:GLU:OE1	2.42	0.50
1:A:310:ASP:O	1:A:311:GLU:C	2.50	0.49
1:A:506:ARG:HD3	1:A:507:ASN:OD1	2.13	0.49
1:A:481:LEU:HB2	1:A:482:PRO:HD3	1.94	0.49
1:A:50:LEU:HD11	1:A:66:VAL:HG22	1.94	0.49
1:A:215:ILE:HD12	1:A:497:MET:HG2	1.94	0.49
1:A:216:ARG:C	1:A:216:ARG:HD3	2.33	0.49
1:A:446:GLY:HA2	1:A:494:LEU:HD21	1.94	0.48
1:A:595:LYS:H	1:A:595:LYS:CD	2.25	0.48
1:A:687:GLN:HB3	1:A:709:ARG:HD2	1.96	0.48
1:A:73:GLN:HB3	1:A:115:ILE:HD11	1.94	0.48
1:A:328:GLU:HG2	1:A:388:PHE:CG	2.48	0.48
1:A:241:ARG:HB2	1:A:246:MET:CE	2.43	0.47
1:A:693:ILE:HG13	1:A:694:ALA:N	2.29	0.47
1:A:814:PRO:HG3	1:A:833:TYR:CD2	2.49	0.47
1:A:6:PHE:CE2	1:A:246:MET:HG2	2.49	0.47
1:A:745:PHE:HA	1:A:748:VAL:HG12	1.96	0.47
1:A:609:PHE:CE2	1:A:610:GLN:HG3	2.50	0.47
1:A:392:LEU:HD12	1:A:400:PRO:HG3	1.97	0.47
1:A:309:PRO:O	1:A:310:ASP:C	2.54	0.46
1:A:871:THR:HG22	1:A:872:LYS:HE3	1.96	0.46
1:A:360:PRO:HD3	1:A:559:LEU:HD23	1.98	0.46
1:A:103:LYS:HD3	1:A:120:TRP:CD2	2.50	0.46
1:A:576:ARG:NH2	1:A:582:PRO:O	2.48	0.46
1:A:578:LYS:NZ	1:A:827:ASP:OD1	2.49	0.46
1:A:863:PHE:CE2	1:A:905:ALA:HA	2.51	0.46
1:A:318:ARG:N	1:A:319:PRO:HD2	2.30	0.46
1:A:792:ASP:N	1:A:792:ASP:OD1	2.48	0.46
1:A:609:PHE:O	1:A:610:GLN:HB2	2.16	0.46
1:A:688:GLU:N	1:A:688:GLU:OE2	2.47	0.46
1:A:6:PHE:CZ	1:A:246:MET:HG2	2.51	0.46
1:A:86:VAL:HG22	1:A:130:LEU:HD22	1.98	0.45
1:A:826:ASN:OD1	1:A:826:ASN:C	2.54	0.45
1:A:726:VAL:HA	1:A:729:GLU:OE1	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:875:ASN:OD1	1:A:876:ARG:N	2.49	0.45
1:A:542:ASP:HB2	4:A:1006:GOL:H2	1.97	0.45
1:A:245:GLU:O	1:A:249:LEU:HB2	2.17	0.45
1:A:314:ARG:HG3	1:A:315:LEU:HD22	1.98	0.45
1:A:572:ILE:O	1:A:576:ARG:HG2	2.16	0.45
1:A:162:PHE:N	1:A:163:PRO:HD3	2.32	0.45
1:A:576:ARG:HD2	1:A:581:GLU:CG	2.47	0.45
1:A:283:THR:CG2	1:A:286:MET:HB3	2.48	0.44
1:A:666:VAL:HG13	1:A:667:GLN:N	2.33	0.44
1:A:859:PHE:O	1:A:891:LEU:HD13	2.18	0.43
1:A:336:GLY:C	1:A:338:PRO:HD2	2.38	0.43
1:A:722:LYS:O	1:A:726:VAL:HG23	2.17	0.43
1:A:678:GLU:N	1:A:679:PRO:HD2	2.33	0.43
1:A:872:LYS:CE	1:A:872:LYS:CA	2.96	0.43
1:A:241:ARG:HB2	1:A:246:MET:HE2	2.01	0.43
1:A:143:LEU:HD23	1:A:184:ALA:HB1	1.99	0.43
1:A:360:PRO:HB3	1:A:767:VAL:HG21	2.01	0.43
1:A:871:THR:O	1:A:872:LYS:HB2	2.18	0.43
1:A:630:ASP:O	1:A:634:LEU:HG	2.19	0.43
1:A:708:LEU:HD13	1:A:723:GLN:HG2	2.01	0.43
1:A:684:ILE:HB	1:A:751:PHE:CD2	2.54	0.43
1:A:845:PRO:HB3	1:A:874:LEU:HD12	2.01	0.43
1:A:118:ARG:NH1	1:A:157:PHE:CD1	2.87	0.43
1:A:169:GLU:OE1	1:A:199:THR:HG21	2.19	0.43
1:A:139:VAL:O	1:A:143:LEU:HD13	2.19	0.43
1:A:686:TYR:CD2	1:A:758:LYS:HD2	2.54	0.43
1:A:841:VAL:HG13	1:A:842:GLY:N	2.34	0.43
1:A:254:PRO:O	1:A:257:LEU:N	2.52	0.42
1:A:611:LEU:HD11	1:A:631:MET:CE	2.48	0.42
1:A:717:PRO:O	1:A:718:GLU:C	2.57	0.42
1:A:4:PRO:CG	1:A:37:PRO:HG3	2.49	0.42
1:A:509:GLY:CA	1:A:510:LYS:CB	2.96	0.42
1:A:602:ARG:HB2	1:A:604:GLU:HG3	2.02	0.42
1:A:650:ASN:O	1:A:654:ARG:HG3	2.19	0.42
1:A:835:ILE:HG23	1:A:838:ILE:HD12	2.01	0.42
1:A:54:TYR:CZ	1:A:271:VAL:HG23	2.55	0.42
1:A:794:THR:O	1:A:798:VAL:HG23	2.19	0.42
1:A:100:LEU:HD23	1:A:124:LEU:HD12	2.02	0.42
1:A:20:GLY:O	1:A:216:ARG:NH1	2.50	0.42
1:A:639:ARG:HB2	1:A:642:PRO:HD2	2.02	0.42
1:A:745:PHE:HA	1:A:748:VAL:CG1	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:858:TYR:CE2	1:A:868:ARG:HG3	2.55	0.41
1:A:81:LEU:CD2	1:A:115:ILE:HD11	2.50	0.41
1:A:436:MET:HE2	1:A:436:MET:HB3	1.88	0.41
1:A:639:ARG:HB2	1:A:642:PRO:CD	2.50	0.41
1:A:797:VAL:O	1:A:801:VAL:HG23	2.21	0.41
1:A:559:LEU:HD12	5:A:1113:HOH:O	2.21	0.41
1:A:642:PRO:HB2	1:A:648:VAL:HG23	2.02	0.41
1:A:801:VAL:HG22	1:A:811:ILE:HD13	2.03	0.41
1:A:509:GLY:HA2	1:A:510:LYS:CB	2.48	0.41
1:A:118:ARG:CG	1:A:118:ARG:NH1	2.83	0.41
1:A:360:PRO:HD3	1:A:559:LEU:CD2	2.51	0.41
1:A:390:ARG:HG3	1:A:691:MET:HE3	2.03	0.41
1:A:247:CYS:HA	1:A:257:LEU:CD1	2.50	0.41
1:A:579:ASN:HB2	1:A:581:GLU:OE1	2.21	0.41
1:A:314:ARG:HG3	1:A:315:LEU:N	2.36	0.41
1:A:600:LEU:HB2	1:A:772:LEU:HD21	2.03	0.41
1:A:668:TRP:HB3	1:A:744:ILE:HG22	2.04	0.41
1:A:314:ARG:HG2	1:A:315:LEU:HD23	2.04	0.40
1:A:12:HIS:CE1	1:A:44[A]:HIS:NE2	2.90	0.40
1:A:227:ASP:OD1	1:A:227:ASP:C	2.60	0.40
1:A:446:GLY:CA	1:A:494:LEU:HD21	2.51	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	908/918 (99%)	863 (95%)	41 (4%)	4 (0%)	34 66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	311	GLU
1	A	858	TYR
1	A	312	GLU
1	A	337	PHE

### 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	753/758 (99%)	723 (96%)	30 (4%)	31	65

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	MET
1	A	118	ARG
1	A	148	SER
1	A	175	ARG
1	A	204	PHE
1	A	216	ARG
1	A	266	ARG
1	A	315	LEU
1	A	390	ARG
1	A	407	CYS
1	A	421	MET
1	A	492	LYS
1	A	500	LYS
1	A	507	ASN
1	A	548	TYR
1	A	614	ARG
1	A	630	ASP
1	A	669	GLN
1	A	674	LYS
1	A	685	LEU
1	A	723	GLN
1	A	729	GLU
1	A	739	GLU

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Mol	Chain	Res	Type
1	A	740	LEU
1	A	749	GLU
1	A	751	PHE
1	A	754	TYR
1	A	877	ARG
1	A	884	MET
1	A	911	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	822	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 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\# Z  > 2$
2	PO4	A	1003	-	4,4,4	0.89	0	6,6,6	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	GOL	A	1006	-	5,5,5	0.35	0	5,5,5	0.33	0
2	PO4	A	1001	-	4,4,4	0.85	0	6,6,6	0.48	0
2	PO4	A	1002	3	4,4,4	0.82	0	6,6,6	0.41	0

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	GOL	A	1006	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1006	GOL	C1-C2-C3-O3
4	A	1006	GOL	O1-C1-C2-O2
4	A	1006	GOL	O1-C1-C2-C3
4	A	1006	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1006	GOL	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	909/918 (99%)	0.22	27 (2%)	50 45	33, 70, 137, 173	1 (0%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	856	GLY	5.7
1	A	310	ASP	3.4
1	A	684	ILE	3.3
1	A	321	TYR	3.2
1	A	645	SER	3.1
1	A	666	VAL	3.1
1	A	668	TRP	3.1
1	A	664	PRO	2.8
1	A	859	PHE	2.8
1	A	651	PHE	2.7
1	A	667	GLN	2.7
1	A	683	ILE	2.6
1	A	727	PHE	2.6
1	A	693	ILE	2.5
1	A	896	ALA	2.4
1	A	895	ARG	2.4
1	A	682	GLY	2.3
1	A	888	PHE	2.3
1	A	640	PRO	2.2
1	A	869	THR	2.2
1	A	578	LYS	2.2
1	A	736	ILE	2.1
1	A	846	ILE	2.1
1	A	857	GLY	2.1
1	A	644	GLN	2.1
1	A	866	CYS	2.1
1	A	643	LEU	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. 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	B-factors( $\text{\AA}^2$ )	Q<0.9
2	PO4	A	1003	5/5	0.70	0.44	113,122,124,127	0
4	GOL	A	1006	6/6	0.71	0.29	52,72,86,86	0
2	PO4	A	1002	5/5	0.94	0.21	66,92,108,111	0
3	ZN	A	1004	1/1	0.95	0.14	78,78,78,78	0
2	PO4	A	1001	5/5	0.97	0.14	67,73,85,87	0
3	ZN	A	1005	1/1	0.98	0.11	87,87,87,87	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.