



# Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 04:49 PM GMT

PDB ID : 4GEY  
Title : High pH structure of Pseudomonas putida OprB  
Authors : van den Berg, Bert  
Deposited on : 2012-08-02  
Resolution : 2.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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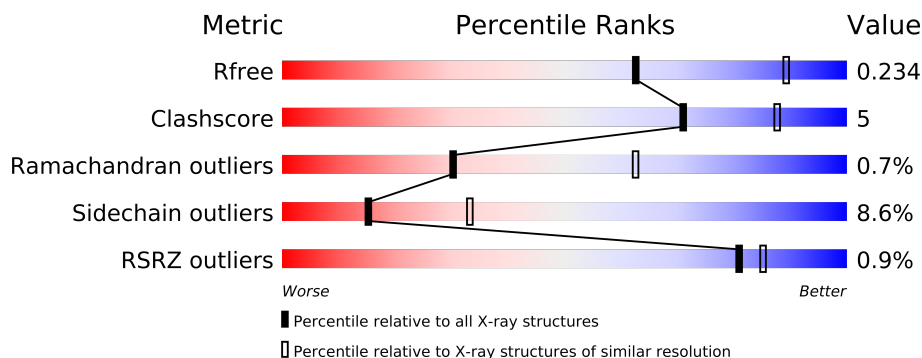
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
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) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.70 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	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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. 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	436	

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

Mol	Type	Chain	Res	Geometry	Electron density
2	BGC	A	501	-	X
3	C8E	A	502	-	X
3	C8E	A	503	-	X
3	C8E	A	504	-	X
3	C8E	A	505	-	X
3	C8E	A	506	-	X
3	C8E	A	507	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
3	C8E	A	508	-	X
3	C8E	A	509	-	X
4	DMU	A	510	-	X
5	PO4	A	514	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 3597 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.

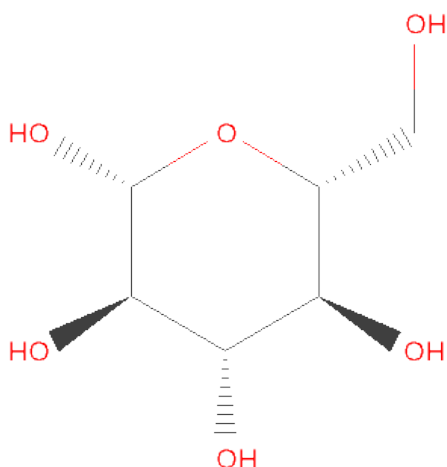
- Molecule 1 is a protein called Porin B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	420	Total	C	N	O	S	Se	0	0	0
			3309	2104	570	629	2	4			

There are 17 discrepancies between the modelled and reference sequences:

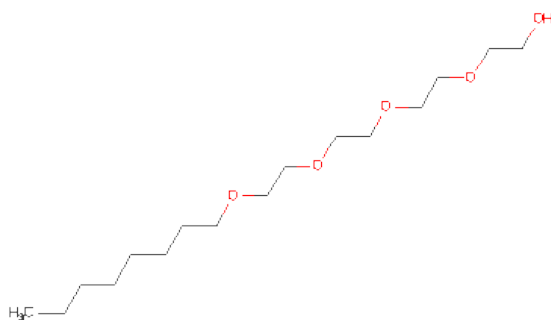
Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	ALA	-	EXPRESSION TAG	UNP E4R6F8
A	-13	ASN	-	EXPRESSION TAG	UNP E4R6F8
A	-12	VAL	-	EXPRESSION TAG	UNP E4R6F8
A	-11	ARG	-	EXPRESSION TAG	UNP E4R6F8
A	-10	LEU	-	EXPRESSION TAG	UNP E4R6F8
A	-9	GLN	-	EXPRESSION TAG	UNP E4R6F8
A	-8	HIS	-	EXPRESSION TAG	UNP E4R6F8
A	-7	HIS	-	EXPRESSION TAG	UNP E4R6F8
A	-6	HIS	-	EXPRESSION TAG	UNP E4R6F8
A	-5	HIS	-	EXPRESSION TAG	UNP E4R6F8
A	-4	HIS	-	EXPRESSION TAG	UNP E4R6F8
A	-3	HIS	-	EXPRESSION TAG	UNP E4R6F8
A	-2	HIS	-	EXPRESSION TAG	UNP E4R6F8
A	-1	LEU	-	EXPRESSION TAG	UNP E4R6F8
A	0	GLU	-	EXPRESSION TAG	UNP E4R6F8
A	220	MSE	LEU	ENGINEERED MUTATION	UNP E4R6F8
A	222	MSE	VAL	ENGINEERED MUTATION	UNP E4R6F8

- Molecule 2 is SUGAR (BETA-D-GLUCOSE) (three-letter code: BGC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is (HYDROXYETHYLOXY)TRI(ETHYLOXY)OCTANE (three-letter code: C8E) (formula:  $C_{16}H_{34}O_5$ ).



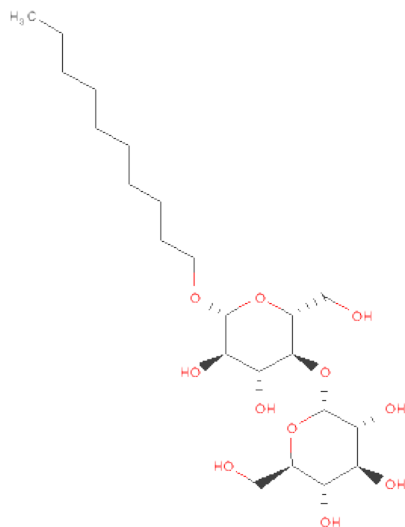
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	10	2		
3	A	1	Total	C	O	0	0
			21	16	5		

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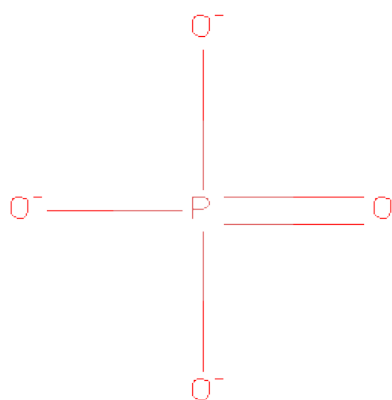
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	8	4		
3	A	1	Total	C	O	0	0
			15	10	5		
3	A	1	Total	C	O	0	0
			12	10	2		
3	A	1	Total	C	O	0	0
			10	9	1		
3	A	1	Total	C	O	0	0
			13	11	2		
3	A	1	Total	C	O	0	0
			9	6	3		

- Molecule 4 is SUGAR (DECYL-BETA-D-MALTOPYRANOSIDE) (three-letter code: DMU) (formula: C<sub>22</sub>H<sub>42</sub>O<sub>11</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			30	19	11		
4	A	1	Total	C	O	0	0
			31	20	11		
4	A	1	Total	C	O	0	0
			30	19	11		
4	A	1	Total	C	O	0	0
			29	18	11		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



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

- Molecule 6 is water.

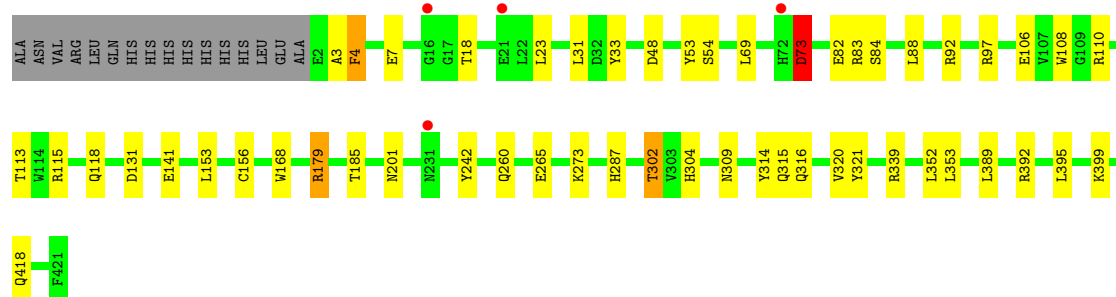
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	47	Total	O	0	0
			47	47		

### 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: Porin B

Chain A: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.34Å 88.34Å 177.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.98 – 2.70 42.87 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.2 (19.98-2.70) 96.8 (42.87-2.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.91 (at 2.69Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8_1069)	Depositor
R, $R_{free}$	0.188 , 0.239 0.190 , 0.234	Depositor DCC
$R_{free}$ test set	1950 reflections (8.88%)	DCC
Wilson B-factor (Å <sup>2</sup> )	47.1	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 21.3	EDS
Estimated twinning fraction	0.038 for -h,-k,l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 22670 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3597	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, BGC, C8E, DMU

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	0.41	0/3398	0.58	0/4609

There are no bond length outliers.

There are no bond angle outliers.

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	3309	0	0	16	0
2	A	12	0	0	3	0
3	A	104	0	0	0	0
4	A	120	0	0	0	0
5	A	5	0	0	0	0
6	A	47	0	0	0	0
All	All	3597	0	0	16	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 5.

All (16) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:106:GLU:O	2:A:501:BGC:O2	1.99	0.80
1:A:33:TYR:OH	1:A:53:TYR:OH	2.02	0.78
1:A:115:ARG:NH2	1:A:201:ASN:O	2.19	0.75
1:A:110:ARG:O	1:A:115:ARG:NH1	2.29	0.65
1:A:48:ASP:OD1	1:A:92:ARG:NH1	2.39	0.56
1:A:108:TRP:O	2:A:501:BGC:O2	2.23	0.55
1:A:287:HIS:ND1	1:A:321:TYR:OH	2.44	0.51
1:A:118:GLN:NE2	1:A:141:GLU:OE1	2.44	0.51
1:A:131:ASP:OD1	1:A:179:ARG:NH2	2.45	0.49
1:A:242:TYR:OH	1:A:304:HIS:ND1	2.52	0.43
1:A:83:ARG:NH2	2:A:501:BGC:O3	2.51	0.42
1:A:273:LYS:NZ	1:A:309:ASN:OD1	2.52	0.42
1:A:302:THR:CG2	1:A:304:HIS:NE2	2.83	0.42
1:A:315:GLN:OE1	1:A:339:ARG:NH1	2.53	0.42
1:A:392:ARG:NH1	1:A:418:GLN:OE1	2.53	0.42
1:A:73:ASP:OD1	1:A:73:ASP:N	2.54	0.41

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	418/436 (96%)	398 (95%)	17 (4%)	3 (1%)	30 62

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	ALA
1	A	4	PHE
1	A	73	ASP

### 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	337/352 (96%)	308 (91%)	29 (9%)	15	33

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	PHE
1	A	7	GLU
1	A	18	THR
1	A	23	LEU
1	A	31	LEU
1	A	54	SER
1	A	69	LEU
1	A	73	ASP
1	A	82	GLU
1	A	84	SER
1	A	88	LEU
1	A	97	ARG
1	A	113	THR
1	A	153	LEU
1	A	156	CYS
1	A	168	TRP
1	A	179	ARG
1	A	185	THR
1	A	260	GLN
1	A	265	GLU
1	A	302	THR
1	A	314	TYR
1	A	316	GLN
1	A	320	VAL
1	A	352	LEU
1	A	353	LEU
1	A	389	LEU
1	A	395	LEU
1	A	399	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 ⓘ

14 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	BGC	A	501	-	12,12,12	2.71	5 (41%)	17,17,17	1.15	2 (11%)
3	C8E	A	502	-	11,11,20	0.46	0	10,10,19	0.63	0
3	C8E	A	503	-	20,20,20	0.47	0	19,19,19	0.41	0
3	C8E	A	504	-	11,11,20	1.59	1 (9%)	9,10,19	0.25	0
3	C8E	A	505	-	14,14,20	1.53	1 (7%)	12,13,19	0.37	0
3	C8E	A	506	-	11,11,20	0.43	0	10,10,19	0.69	0
3	C8E	A	507	-	8,9,20	0.51	0	7,8,19	0.94	1 (14%)
3	C8E	A	508	-	12,12,20	1.86	2 (16%)	10,11,19	0.55	0
3	C8E	A	509	-	8,8,20	1.50	1 (12%)	6,7,19	0.43	0
4	DMU	A	510	-	31,31,34	2.78	10 (32%)	42,42,45	1.61	8 (19%)
4	DMU	A	511	-	32,32,34	3.09	10 (31%)	43,43,45	2.04	9 (20%)
4	DMU	A	512	-	31,31,34	2.83	10 (32%)	42,42,45	1.73	9 (21%)
4	DMU	A	513	-	30,30,34	2.67	12 (40%)	41,41,45	1.70	7 (17%)
5	PO4	A	514	-	4,4,4	0.24	0	6,6,6	0.31	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
2	BGC	A	501	-	-	0/2/22/22	0/1/1/1
3	C8E	A	502	-	-	0/9/9/18	0/0/0/0
3	C8E	A	503	-	-	0/18/18/18	0/0/0/0
3	C8E	A	504	-	-	0/9/9/18	0/0/0/0
3	C8E	A	505	-	-	0/12/12/18	0/0/0/0
3	C8E	A	506	-	-	0/9/9/18	0/0/0/0
3	C8E	A	507	-	-	0/7/7/18	0/0/0/0
3	C8E	A	508	-	-	0/10/10/18	0/0/0/0
3	C8E	A	509	-	-	0/6/6/18	0/0/0/0
4	DMU	A	510	-	-	0/16/56/59	0/2/2/2
4	DMU	A	511	-	-	0/17/57/59	0/2/2/2
4	DMU	A	512	-	-	0/16/56/59	0/2/2/2
4	DMU	A	513	-	-	0/15/55/59	0/2/2/2
5	PO4	A	514	-	-	0/0/0/0	0/0/0/0

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	511	DMU	C37-C34	-12.58	1.50	1.55
4	A	512	DMU	C34-C31	-9.82	1.51	1.55
4	A	510	DMU	C34-C31	-8.83	1.51	1.55
4	A	513	DMU	C31-C28	-7.58	1.52	1.55
4	A	511	DMU	C11-C9	-5.56	1.32	1.52
3	A	505	C8E	C7-C8	-5.51	1.53	1.55
4	A	510	DMU	C11-C9	-5.43	1.32	1.52
4	A	513	DMU	C11-C9	-5.39	1.33	1.52
4	A	512	DMU	C11-C9	-5.38	1.33	1.52
3	A	504	C8E	C10-C11	-5.07	1.53	1.55
3	A	508	C8E	C2-C3	-4.96	1.53	1.55
4	A	513	DMU	O1-C9	4.60	1.55	1.44
4	A	513	DMU	C57-C4	-4.60	1.35	1.52
4	A	510	DMU	C57-C4	-4.58	1.35	1.52
4	A	511	DMU	C57-C4	-4.55	1.35	1.52
2	A	501	BGC	C3-C2	-4.54	1.40	1.52
4	A	510	DMU	O5-C4	4.52	1.55	1.44
4	A	512	DMU	C57-C4	-4.50	1.36	1.52
2	A	501	BGC	C6-C5	-4.48	1.36	1.52
2	A	501	BGC	C1-C2	-4.38	1.44	1.52
4	A	512	DMU	O5-C4	4.35	1.55	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	511	DMU	O5-C4	4.22	1.54	1.44
4	A	510	DMU	O1-C9	4.18	1.54	1.44
3	A	509	C8E	C13-C14	-4.13	1.53	1.55
4	A	512	DMU	O1-C9	4.08	1.54	1.44
4	A	513	DMU	O5-C4	4.05	1.54	1.44
4	A	511	DMU	O1-C9	3.89	1.54	1.44
3	A	508	C8E	C14-C13	-3.88	1.53	1.55
4	A	510	DMU	C5-C7	-3.64	1.42	1.52
2	A	501	BGC	O5-C5	3.48	1.53	1.44
4	A	511	DMU	C5-C7	-3.41	1.43	1.52
4	A	510	DMU	C2-C1	-3.38	1.43	1.52
4	A	513	DMU	C5-C7	-3.37	1.43	1.52
4	A	512	DMU	C5-C7	-3.30	1.43	1.52
4	A	512	DMU	C2-C1	-3.11	1.44	1.52
4	A	513	DMU	C2-C1	-2.89	1.44	1.52
4	A	511	DMU	C2-C1	-2.73	1.45	1.52
4	A	511	DMU	O49-C1	2.67	1.49	1.43
4	A	510	DMU	C6-C1	-2.60	1.44	1.52
4	A	513	DMU	O49-C1	2.57	1.49	1.43
4	A	513	DMU	C6-C1	-2.56	1.44	1.52
4	A	512	DMU	C6-C1	-2.37	1.45	1.52
4	A	512	DMU	O49-C1	2.35	1.48	1.43
4	A	513	DMU	C8-C9	2.31	1.58	1.53
4	A	510	DMU	O49-C1	2.24	1.48	1.43
4	A	510	DMU	O3-C5	2.22	1.48	1.43
2	A	501	BGC	C4-C3	-2.19	1.46	1.52
4	A	511	DMU	O16-C6	2.13	1.44	1.40
4	A	512	DMU	O3-C5	2.12	1.48	1.43
4	A	513	DMU	O3-C5	2.11	1.48	1.43
4	A	513	DMU	C10-C5	-2.04	1.46	1.52
4	A	511	DMU	O55-C2	2.01	1.47	1.43

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	511	DMU	C10-O1-C9	5.51	124.44	113.73
4	A	511	DMU	C1-C2-C3	5.38	121.31	109.61
4	A	510	DMU	C10-O1-C9	5.36	124.14	113.73
4	A	512	DMU	C10-O1-C9	4.93	123.31	113.73
4	A	513	DMU	C10-O1-C9	4.73	122.93	113.73
4	A	512	DMU	C18-O16-C6	4.67	122.36	113.96
4	A	511	DMU	C6-C1-C2	4.40	118.55	110.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	511	DMU	C10-C5-C7	4.03	117.84	110.00
4	A	511	DMU	O16-C6-C1	3.91	113.15	108.18
4	A	513	DMU	C7-C8-C9	3.91	117.18	110.20
4	A	513	DMU	C18-O16-C6	3.90	120.97	113.96
4	A	513	DMU	C1-C2-C3	3.55	117.33	109.61
4	A	510	DMU	O1-C10-C5	3.16	116.79	110.31
4	A	510	DMU	C18-O16-C6	3.15	119.62	113.96
4	A	512	DMU	C1-C2-C3	2.87	115.85	109.61
4	A	511	DMU	O1-C10-C5	2.85	116.16	110.31
4	A	512	DMU	C7-C8-C9	2.67	114.96	110.20
4	A	511	DMU	C18-O16-C6	2.56	118.56	113.96
4	A	510	DMU	O16-C6-C1	2.46	111.32	108.18
4	A	512	DMU	O16-C6-C1	2.45	111.31	108.18
2	A	501	BGC	O6-C6-C5	2.44	119.75	111.36
4	A	513	DMU	C6-O5-C4	2.34	118.27	113.73
4	A	513	DMU	C6-C1-C2	2.30	114.47	110.00
4	A	510	DMU	C34-C31-C28	2.25	119.52	114.46
4	A	511	DMU	O7-C10-C5	2.23	113.47	108.12
3	A	507	C8E	O9-C8-C7	2.22	111.70	107.35
4	A	512	DMU	O7-C10-C5	2.22	113.44	108.12
4	A	512	DMU	C34-C31-C28	2.21	119.43	114.46
4	A	512	DMU	O7-C10-O1	-2.19	105.24	110.69
4	A	510	DMU	O16-C18-C19	2.18	118.36	109.87
4	A	510	DMU	C6-O5-C4	2.17	117.94	113.73
4	A	513	DMU	C8-C7-C5	2.17	114.82	110.82
4	A	512	DMU	O1-C10-C5	2.14	114.70	110.31
4	A	510	DMU	C1-C2-C3	2.08	114.12	109.61
4	A	511	DMU	O16-C18-C19	2.04	117.83	109.87
2	A	501	BGC	O5-C5-C4	-2.02	106.02	109.76

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, 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	420/436 (96%)	-0.35	4 (0%) 79 83	22, 35, 55, 65	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	231	ASN	2.3
1	A	72	HIS	2.2
1	A	16	GLY	2.1
1	A	21	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, 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	BGC	A	501	12/12	0.34	10.86	48,55,61,71	12

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	C8E	A	504	12/21	0.36	7.73	66,76,87,88	0
4	DMU	A	510	30/33	0.28	7.20	43,57,85,92	0
3	C8E	A	506	12/21	0.32	5.87	54,60,67,72	0
3	C8E	A	503	21/21	0.29	5.85	50,67,79,80	0
3	C8E	A	507	10/21	0.32	5.12	44,51,59,63	0
3	C8E	A	505	15/21	0.26	4.42	50,62,89,90	0
5	PO4	A	514	5/5	0.26	4.04	106,108,110,111	0
3	C8E	A	508	13/21	0.26	3.73	40,70,73,73	0
3	C8E	A	509	9/21	0.27	3.10	47,50,69,70	0
3	C8E	A	502	12/21	0.18	2.05	34,43,52,55	0
4	DMU	A	512	30/33	0.28	1.96	38,66,91,95	0
4	DMU	A	513	29/33	0.23	1.84	43,69,110,113	0
4	DMU	A	511	31/33	0.15	0.47	31,53,64,69	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.