



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 23, 2024 – 10:50 AM EDT

PDB ID : 3W6L  
Title : Contribution of disulfide bond toward thermostability in hyperthermostable endocellulase  
Authors : Kim, H.-W.; Ishikawa, K.  
Deposited on : 2013-02-15  
Resolution : 1.75 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

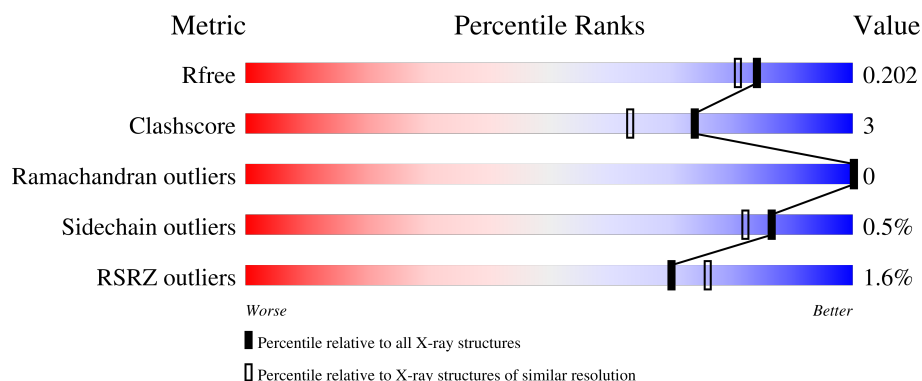
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

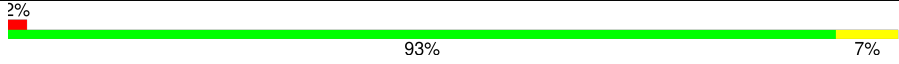
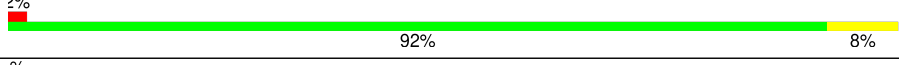
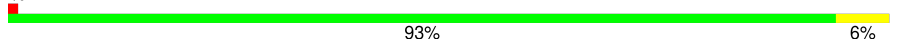
The reported resolution of this entry is 1.75 Å.

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))
$R_{free}$	164625	2888 (1.76-1.76)
Clashscore	180529	3097 (1.76-1.76)
Ramachandran outliers	177936	3072 (1.76-1.76)
Sidechain outliers	177891	3072 (1.76-1.76)
RSRZ outliers	164620	2887 (1.76-1.76)

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 of 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	377	
1	B	377	
1	C	377	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10022 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.

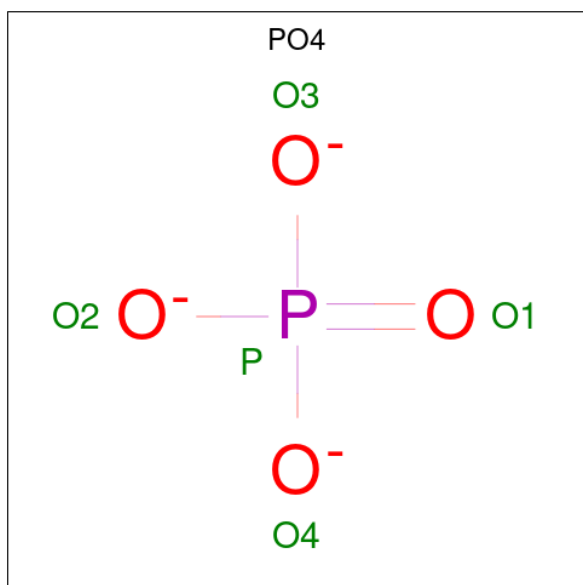
- Molecule 1 is a protein called 458aa long hypothetical endo-1,4-beta-glucanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	377	Total	C	N	O	S	0	0	0
			3069	2002	500	557	10			
1	B	377	Total	C	N	O	S	0	0	0
			3069	2002	500	557	10			
1	C	377	Total	C	N	O	S	0	0	0
			3069	2002	500	557	10			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	289	LYS	ARG	conflict	UNP O58925
B	289	LYS	ARG	conflict	UNP O58925
C	289	LYS	ARG	conflict	UNP O58925

- 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 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0
2	C	1	Total O P 5 4 1	0	0

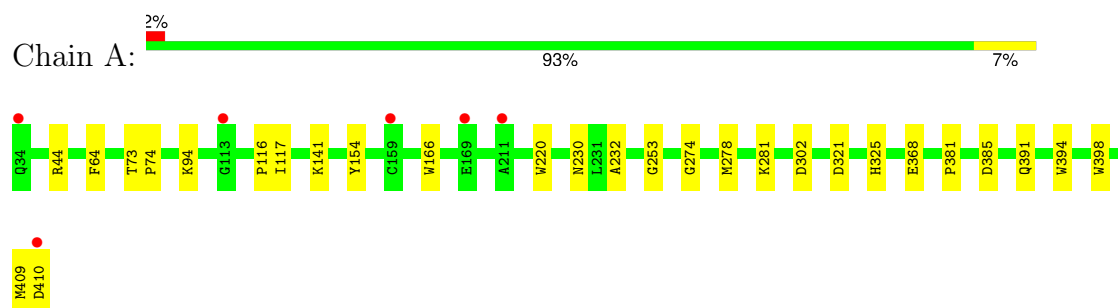
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	254	Total O 254 254	0	0
3	B	258	Total O 258 258	0	0
3	C	228	Total O 228 228	0	0

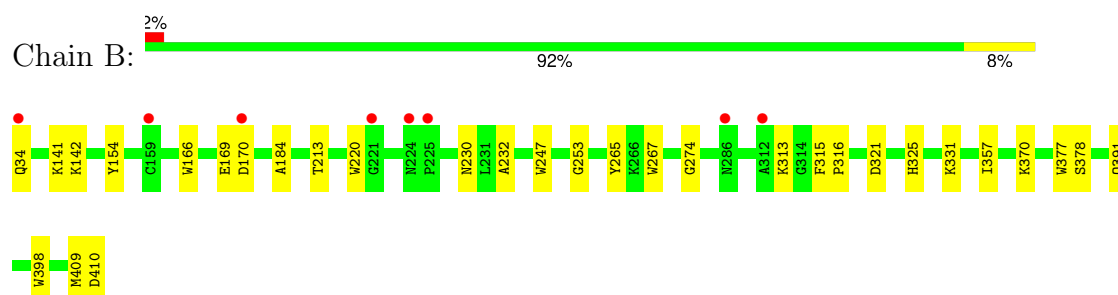
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

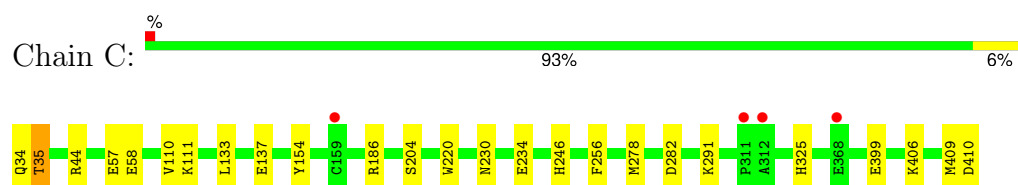
- Molecule 1: 458aa long hypothetical endo-1,4-beta-glucanase



- Molecule 1: 458aa long hypothetical endo-1,4-beta-glucanase



- Molecule 1: 458aa long hypothetical endo-1,4-beta-glucanase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	163.10Å 58.34Å 137.91Å 90.00° 109.65° 90.00°	Depositor
Resolution (Å)	37.53 – 1.75 37.53 – 1.75	Depositor EDS
% Data completeness (in resolution range)	98.2 (37.53-1.75) 98.2 (37.53-1.75)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.74 (at 1.75Å)	Xtriage
Refinement program	PHENIX 1.7.1 _743	Depositor
R, $R_{free}$	0.176 , 0.205 0.174 , 0.202	Depositor DCC
$R_{free}$ test set	6097 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.8	Xtriage
Anisotropy	0.060	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.41 , 45.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	10022	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.00% 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: 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	0.39	0/3180	0.54	0/4341
1	B	0.38	0/3180	0.54	0/4341
1	C	0.33	0/3180	0.49	0/4341
All	All	0.37	0/9540	0.52	0/13023

There are no bond length outliers.

There are no bond angle outliers.

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	3069	0	2908	19	0
1	B	3069	0	2908	17	2
1	C	3069	0	2908	18	0
2	A	20	0	0	0	0
2	B	40	0	0	0	0
2	C	15	0	0	0	0
3	A	254	0	0	6	1
3	B	258	0	0	3	2
3	C	228	0	0	4	0
All	All	10022	0	8724	54	3

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 3.

All (54) 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:281:LYS:NZ	3:A:804:HOH:O	2.02	0.91
1:A:44:ARG:NH1	3:A:819:HOH:O	2.03	0.89
1:B:315:PHE:CE2	1:B:357:ILE:HD11	2.07	0.89
1:C:282:ASP:OD2	3:C:812:HOH:O	1.93	0.86
1:A:44:ARG:NH2	3:A:829:HOH:O	2.17	0.78
1:A:321:ASP:OD2	3:A:740:HOH:O	2.09	0.69
1:A:368:GLU:OE2	3:A:752:HOH:O	2.13	0.67
1:C:34:GLN:N	3:C:808:HOH:O	2.27	0.67
1:C:409:MET:O	1:C:410:ASP:HB2	1.95	0.66
1:B:213:THR:OG1	3:B:761:HOH:O	2.17	0.60
1:A:409:MET:O	1:A:410:ASP:CG	2.45	0.55
1:A:409:MET:O	1:A:410:ASP:CB	2.54	0.55
1:A:166:TRP:CZ2	1:A:232:ALA:HB2	2.43	0.53
1:B:409:MET:O	1:B:410:ASP:HB2	2.08	0.53
1:B:220:TRP:CE2	1:B:230:ASN:HB3	2.43	0.53
1:A:141:LYS:HD2	3:A:661:HOH:O	2.09	0.53
1:B:253:GLY:O	1:B:274:GLY:HA2	2.08	0.52
1:C:406:LYS:HA	1:C:409:MET:CE	2.40	0.52
1:C:57:GLU:HG2	3:C:736:HOH:O	2.12	0.50
1:C:399:GLU:CG	3:C:789:HOH:O	2.59	0.50
1:C:230:ASN:O	1:C:234:GLU:HG3	2.12	0.50
1:C:204:SER:HG	1:C:256:PHE:HD2	1.60	0.49
1:C:291:LYS:HA	1:C:291:LYS:HE2	1.94	0.49
1:C:406:LYS:O	1:C:410:ASP:N	2.46	0.48
1:A:381:PRO:HD2	1:A:394:TRP:CH2	2.48	0.48
1:A:220:TRP:CE2	1:A:230:ASN:HB3	2.50	0.47
1:C:406:LYS:HA	1:C:409:MET:HE2	1.96	0.47
1:A:73:THR:HB	1:A:74:PRO:HD2	1.97	0.46
1:A:391:GLN:HG2	1:A:398:TRP:CE2	2.51	0.46
1:B:331:LYS:HE3	1:B:370:LYS:O	2.16	0.46
1:A:381:PRO:HD2	1:A:394:TRP:CZ2	2.51	0.46
1:B:391:GLN:HG2	1:B:398:TRP:CE2	2.51	0.45
1:A:116:PRO:O	1:A:117:ILE:HD13	2.17	0.44
1:A:253:GLY:O	1:A:274:GLY:HA2	2.18	0.44
1:B:166:TRP:CZ2	1:B:232:ALA:HB2	2.53	0.43
1:B:169:GLU:HG2	1:B:170:ASP:OD2	2.18	0.43
1:C:110:VAL:HA	1:C:133:LEU:HD13	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:PHE:CD2	1:A:409:MET:HG2	2.54	0.43
1:C:35:THR:HB	1:C:246:HIS:HB3	2.00	0.43
1:B:34:GLN:NE2	3:B:811:HOH:O	2.45	0.42
1:A:278:MET:HG2	1:A:325:HIS:O	2.19	0.42
1:B:321:ASP:OD1	1:B:325:HIS:HE1	2.02	0.42
1:C:44:ARG:NH2	1:C:58:GLU:OE2	2.47	0.42
1:C:278:MET:HG2	1:C:325:HIS:O	2.20	0.42
1:B:315:PHE:CZ	1:B:357:ILE:HD11	2.53	0.42
1:B:265:TYR:CE2	1:B:267:TRP:HB2	2.55	0.41
1:B:316:PRO:HG3	1:B:357:ILE:CD1	2.50	0.41
1:C:111:LYS:HA	1:C:111:LYS:HD3	1.90	0.41
1:C:137:GLU:OE2	1:C:186:ARG:HD3	2.21	0.41
1:C:220:TRP:CE2	1:C:230:ASN:HB3	2.55	0.41
1:A:302:ASP:HB3	1:A:385:ASP:O	2.20	0.41
1:B:184:ALA:HA	1:B:247:TRP:CH2	2.56	0.41
1:B:313:LYS:O	3:B:736:HOH:O	2.21	0.40
1:B:377:TRP:HA	1:B:378:SER:HA	1.82	0.40

All (3) 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)
1:B:141:LYS:NZ	3:B:796:HOH:O[4_555]	1.86	0.34
1:B:142:LYS:NZ	3:B:766:HOH:O[4_555]	1.91	0.29
3:A:713:HOH:O	3:A:797:HOH:O[4_556]	1.95	0.25

## 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	375/377 (100%)	366 (98%)	9 (2%)	0	<b>100</b> <b>100</b>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	375/377 (100%)	367 (98%)	8 (2%)	0	100	100
1	C	375/377 (100%)	368 (98%)	7 (2%)	0	100	100
All	All	1125/1131 (100%)	1101 (98%)	24 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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	324/324 (100%)	322 (99%)	2 (1%)	84	78
1	B	324/324 (100%)	323 (100%)	1 (0%)	91	88
1	C	324/324 (100%)	322 (99%)	2 (1%)	84	78
All	All	972/972 (100%)	967 (100%)	5 (0%)	86	82

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	94	LYS
1	A	154	TYR
1	B	154	TYR
1	C	35	THR
1	C	154	TYR

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

### 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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

15 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	PO4	A	502	-	4,4,4	0.92	0	6,6,6	0.43	0
2	PO4	B	502	-	4,4,4	0.85	0	6,6,6	0.36	0
2	PO4	C	501	-	4,4,4	0.99	0	6,6,6	0.53	0
2	PO4	B	508	-	4,4,4	0.85	0	6,6,6	0.70	0
2	PO4	B	503	-	4,4,4	0.93	0	6,6,6	0.55	0
2	PO4	A	503	-	4,4,4	0.80	0	6,6,6	0.53	0
2	PO4	B	506	-	4,4,4	0.92	0	6,6,6	0.43	0
2	PO4	A	501	-	4,4,4	0.95	0	6,6,6	0.42	0
2	PO4	B	507	-	4,4,4	0.92	0	6,6,6	0.35	0
2	PO4	C	503	-	4,4,4	1.00	0	6,6,6	0.49	0
2	PO4	B	505	-	4,4,4	0.81	0	6,6,6	0.64	0
2	PO4	B	504	-	4,4,4	0.74	0	6,6,6	0.54	0
2	PO4	B	501	-	4,4,4	0.84	0	6,6,6	0.33	0
2	PO4	A	504	-	4,4,4	0.94	0	6,6,6	0.43	0
2	PO4	C	502	-	4,4,4	0.89	0	6,6,6	0.64	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	377/377 (100%)	-0.17	6 (1%) 70 76	12, 20, 33, 46	6 (1%)
1	B	377/377 (100%)	-0.04	8 (2%) 63 69	12, 21, 36, 46	8 (2%)
1	C	377/377 (100%)	-0.03	4 (1%) 77 83	13, 25, 38, 50	8 (2%)
All	All	1131/1131 (100%)	-0.08	18 (1%) 70 76	12, 22, 36, 50	22 (1%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	312	ALA	4.9
1	A	410	ASP	4.1
1	C	312	ALA	3.7
1	C	311	PRO	3.3
1	A	34	GLN	2.9
1	A	159	CYS	2.7
1	B	34	GLN	2.6
1	C	159	CYS	2.5
1	B	170	ASP	2.4
1	B	286	ASN	2.2
1	A	211	ALA	2.2
1	A	169	GLU	2.1
1	B	225	PRO	2.1
1	A	113	GLY	2.1
1	B	159	CYS	2.1
1	B	224	ASN	2.1
1	C	368	GLU	2.1
1	B	221	GLY	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 monosaccharides 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	B	505	5/5	0.67	0.15	42,49,54,60	0
2	PO4	A	504	5/5	0.74	0.12	48,50,64,65	0
2	PO4	B	507	5/5	0.80	0.11	32,36,50,50	0
2	PO4	A	501	5/5	0.83	0.12	36,40,46,51	0
2	PO4	C	503	5/5	0.83	0.11	39,47,53,55	0
2	PO4	B	506	5/5	0.84	0.11	37,38,49,50	0
2	PO4	A	502	5/5	0.86	0.11	34,38,43,51	0
2	PO4	B	508	5/5	0.93	0.14	30,30,38,41	0
2	PO4	B	504	5/5	0.94	0.15	26,28,34,36	0
2	PO4	A	503	5/5	0.94	0.21	22,31,37,37	0
2	PO4	B	503	5/5	0.94	0.14	31,32,38,39	0
2	PO4	C	502	5/5	0.95	0.17	26,28,34,36	0
2	PO4	C	501	5/5	0.97	0.07	19,23,31,35	0
2	PO4	B	502	5/5	0.98	0.09	23,23,26,28	0
2	PO4	B	501	5/5	0.98	0.06	22,24,25,27	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.