



# Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 4QPW  
Title : BiXyn10A CBM1 with Xylohexaose Bound  
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Deposited on : 2014-06-25  
Resolution : 1.14 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.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

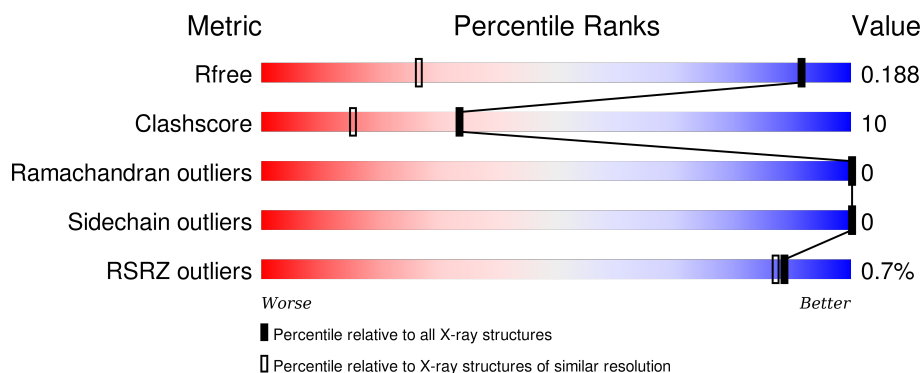
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.14 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	1495 (1.20-1.08)
Clashscore	102246	1593 (1.20-1.08)
Ramachandran outliers	100387	1521 (1.20-1.08)
Sidechain outliers	100360	1518 (1.20-1.08)
RSRZ outliers	91569	1498 (1.20-1.08)

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	142	<div> <div></div> <div>83%</div> <div>15%</div> <div></div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1466 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 glycosyl hydrolase family 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	142	Total	C	N	O	S	0	1	0
			1127	718	182	222	5			

- Molecule 2 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	3	Total	C	O	0	0
			28	15	13		


- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	311	Total	O	0	0
			311	311		

### 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 ( $\text{RSRZ} > 2$ ). 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: glycosyl hydrolase family 10

Chain A: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	36.87Å 57.05Å 85.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.44 – 1.14 47.44 – 1.14	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.44-1.14) 99.9 (47.44-1.14)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.33 (at 1.15Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.160 , 0.178 0.170 , 0.188	Depositor DCC
$R_{free}$ test set	3223 reflections (5.15%)	DCC
Wilson B-factor (Å <sup>2</sup> )	12.0	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 55.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 65782 reflections	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	1466	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.63% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: XYP

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	1.57	10/1156 (0.9%)	1.27	11/1559 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	211	PHE	CG-CD1	-6.68	1.28	1.38
1	A	221	PRO	C-O	6.61	1.36	1.23
1	A	178	GLU	CD-OE1	-6.45	1.18	1.25
1	A	169	SER	CB-OG	6.37	1.50	1.42
1	A	188	THR	CB-OG1	-6.37	1.30	1.43
1	A	250	SER	CB-OG	-5.75	1.34	1.42
1	A	199	GLU	CD-OE2	-5.67	1.19	1.25
1	A	166	ILE	N-CA	5.27	1.56	1.46
1	A	165	GLU	CD-OE1	-5.16	1.20	1.25
1	A	249	TRP	CZ2-CH2	-5.10	1.27	1.37

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	275	ASP	CB-CG-OD1	9.43	126.78	118.30
1	A	182	GLU	OE1-CD-OE2	-6.82	115.11	123.30
1	A	218	ASP	CB-CG-OD2	-6.58	112.37	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	178	GLU	OE1-CD-OE2	-6.54	115.46	123.30
1	A	235	ARG	NE-CZ-NH2	-6.23	117.18	120.30
1	A	232[A]	MET	CG-SD-CE	-6.15	90.35	100.20
1	A	232[B]	MET	CG-SD-CE	-6.15	90.35	100.20
1	A	163	ASP	CB-CG-OD2	5.51	123.26	118.30
1	A	275	ASP	CB-CG-OD2	-5.49	113.36	118.30
1	A	211	PHE	CB-CG-CD1	5.29	124.50	120.80
1	A	211	PHE	CB-CG-CD2	-5.02	117.29	120.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	221	PRO	Mainchain

## 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	1127	0	1078	23	0
2	A	28	0	24	0	0
3	A	311	0	0	16	3
All	All	1466	0	1102	23	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 10.

All (23) 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:232[A]:MET:CE	1:A:232[A]:MET:SD	2.03	1.46
1:A:170:THR:O	3:A:1291:HOH:O	1.94	0.85
1:A:165:GLU:HG3	3:A:1194:HOH:O	1.79	0.82
1:A:170:THR:C	3:A:1291:HOH:O	2.21	0.79
1:A:165:GLU:CG	3:A:1194:HOH:O	2.31	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:MET:HE3	3:A:1036:HOH:O	1.85	0.76
1:A:233:MET:CE	3:A:1036:HOH:O	2.42	0.67
1:A:232[A]:MET:CE	1:A:232[A]:MET:CG	2.73	0.67
1:A:161:LYS:HD3	3:A:1248:HOH:O	1.93	0.67
1:A:171:ALA:HA	3:A:1291:HOH:O	1.95	0.65
1:A:271:LYS:HE3	3:A:1133:HOH:O	2.02	0.59
1:A:165:GLU:CD	3:A:1194:HOH:O	2.38	0.59
1:A:232[A]:MET:CE	1:A:232[A]:MET:HB2	2.35	0.57
1:A:232[A]:MET:CE	1:A:232[A]:MET:CB	2.84	0.55
1:A:232[A]:MET:HE3	1:A:232[A]:MET:HB2	1.88	0.55
1:A:210:HIS:HD2	3:A:1037:HOH:O	1.90	0.54
1:A:173:TYR:HD2	3:A:1138:HOH:O	1.90	0.53
1:A:271:LYS:CE	3:A:1133:HOH:O	2.57	0.52
1:A:232[A]:MET:CB	1:A:232[A]:MET:HE3	2.45	0.45
1:A:161:LYS:CD	3:A:1245:HOH:O	2.66	0.42
1:A:171:ALA:N	3:A:1291:HOH:O	2.48	0.41
1:A:233:MET:CE	1:A:293:LYS:HG2	2.50	0.41
1:A:161:LYS:HD3	3:A:1245:HOH:O	2.19	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)
3:A:1170:HOH:O	3:A:1299:HOH:O[4_545]	1.86	0.34
3:A:1255:HOH:O	3:A:1262:HOH:O[4_545]	2.15	0.05
3:A:1071:HOH:O	3:A:1268:HOH:O[3_545]	2.18	0.02

## 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	141/142 (99%)	138 (98%)	3 (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	121/120 (101%)	121 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	210	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](#)

3 carbohydrates 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	XYP	A	901	2	9,9,10	1.38	2 (22%)	12,12,14	1.67	3 (25%)
2	XYP	A	902	2	9,9,10	1.40	2 (22%)	12,12,14	1.71	4 (33%)
2	XYP	A	903	2	10,10,10	2.02	2 (20%)	12,14,14	2.64	6 (50%)

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	XYP	A	901	2	-	0/0/14/17	0/1/1/1
2	XYP	A	902	2	-	0/0/14/17	0/1/1/1
2	XYP	A	903	2	-	0/0/17/17	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	902	XYP	C4B-C3B	-2.43	1.49	1.52
2	A	902	XYP	O5B-C1B	-2.22	1.38	1.42
2	A	901	XYP	C4B-C3B	2.21	1.55	1.52
2	A	901	XYP	O5B-C1B	2.79	1.48	1.42
2	A	903	XYP	O5B-C5B	4.08	1.50	1.43
2	A	903	XYP	O4B-C4B	4.48	1.53	1.43

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	903	XYP	O4B-C4B-C5B	-5.92	97.34	109.21
2	A	903	XYP	O5B-C5B-C4B	-3.71	104.84	110.86
2	A	902	XYP	C1B-C2B-C3B	-3.44	105.48	109.54
2	A	901	XYP	C4B-C3B-C2B	-3.03	108.24	111.24
2	A	901	XYP	C5B-O5B-C1B	-2.67	107.27	111.57
2	A	903	XYP	C1B-C2B-C3B	-2.63	106.52	110.43
2	A	903	XYP	O3B-C3B-C4B	-2.37	105.72	110.00
2	A	903	XYP	C4B-C3B-C2B	-2.36	107.03	111.04
2	A	902	XYP	O2B-C2B-C1B	2.05	113.32	109.21
2	A	901	XYP	O4B-C4B-C5B	2.13	113.47	109.21
2	A	902	XYP	O5B-C1B-C2B	2.65	115.26	110.31
2	A	902	XYP	C4B-C3B-C2B	2.69	113.89	111.24
2	A	903	XYP	O2B-C2B-C1B	3.44	117.39	109.82

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.6 Ligand geometry

There are no ligands in this entry.

## 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	142/142 (100%)	-0.07	1 (0%) 89 87	8, 13, 24, 28	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	170	THR	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](#)

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	XYP	A	901	9/10	0.95	0.06	-0.55	15,16,20,21	0
2	XYP	A	903	10/10	0.91	0.11	-	21,28,36,38	0
2	XYP	A	902	9/10	0.97	0.07	-	16,17,19,19	0

### 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.