

# Full wwPDB X-ray Structure Validation Report (i)

#### Oct 22, 2023 – 12:58 PM EDT

PDB ID	:	3EXU
Title	:	A glycoside hydrolase family 11 xylanase with an extended thumb region
Authors	:	Vandermarliere, E.; Pollet, A.; Strelkov, S.V.; Delcour, J.A.; Courtin, C.M.
Deposited on		
Resolution	:	1.81 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* A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

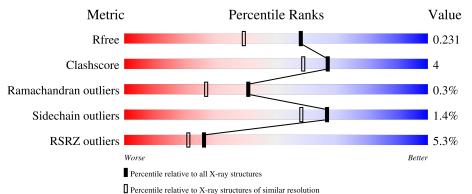
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber $(2001)$
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 1.81 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
$R_{free}$	130704	7484(1.84-1.80)
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)
RSRZ outliers	127900	7371 (1.84-1.80)

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 >=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 <=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	А	185	92%	8%	•					
1	В	185	5% 86%	6% • 6%	1					



#### 3EXU

# 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3096 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.

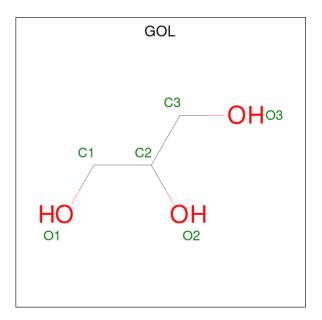
• Molecule 1 is a protein called Endo-1,4-beta-xylanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	А	185	Total 1451	-		0 286	${ m S} { m 2}$	0	1	0
1	В	173	Total 1358	-	N 230	O 264	$\begin{array}{c} \mathrm{S} \\ \mathrm{2} \end{array}$	0	1	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	11	PHE	ASP	engineered mutation	UNP Q59254
А	122	ASP	ARG	engineered mutation	UNP Q59254
В	11	PHE	ASP	engineered mutation	UNP Q59254
В	122	ASP	ARG	engineered mutation	UNP Q59254

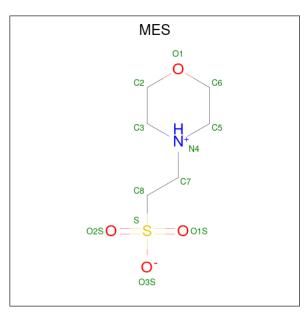
• Molecule 2 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
2	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0

• Molecule 3 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula:  $C_6H_{13}NO_4S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	В	1	Total 12	C 6	N 1	0 4	S 1	0	0

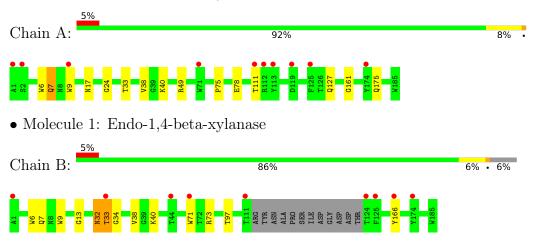
• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	136	Total O 136 136	0	0
4	В	127	Total O 127 127	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: Endo-1,4-beta-xylanase



# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	43.91Å 51.88Å 84.73Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $94.20^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	32.76 - 1.81	Depositor
	32.76 - 1.81	EDS
% Data completeness	87.9 (32.76-1.81)	Depositor
(in resolution range)	87.9 (32.76-1.81)	EDS
R <sub>merge</sub>	(Not available)	Depositor
R <sub>sym</sub>	0.08	Depositor
$< I/\sigma(I) > 1$	$1.96 (at 1.81 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
$R, R_{free}$	0.191 , $0.232$	Depositor
It, Itfree	0.190 , $0.231$	DCC
$R_{free}$ test set	1541 reflections $(5.03\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	18.2	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.37, $45.5$	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.48, < L^2>=0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3096	wwPDB-VP
Average B, all atoms $(Å^2)$	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 9.20% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<sup>&</sup>lt;sup>1</sup>Intensities 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, MES

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	А	0.43	0/1503	0.54	0/2060	
1	В	0.46	0/1407	0.64	1/1926~(0.1%)	
All	All	0.44	0/2910	0.59	1/3986~(0.0%)	

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
1	В	32	ASN	C-N-CA	6.48	137.91	121.70

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	А	1451	0	1323	12	0
1	В	1358	0	1242	9	0
2	А	6	0	8	2	0
2	В	6	0	8	1	0
3	В	12	0	12	1	0
4	А	136	0	0	0	0
4	В	127	0	0	2	0
All	All	3096	0	2593	21	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 4.

All $(21)$ close contacts	within the same	asymmetric unit	are listed bel	low, sorted by their clash
magnitude.				

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:B:9:TRP:CH2	1:B:166:TYR:OH	2.16	0.99
1:B:9:TRP:CZ2	1:B:166:TYR:OH	2.24	0.90
1:A:6:TRP:CH2	1:A:38:VAL:HG13	2.25	0.70
1:A:75:PRO:HD2	1:A:161:GLY:HA2	1.83	0.60
1:A:49:ARG:HA	2:A:186:GOL:C3	2.33	0.59
1:A:49:ARG:HA	2:A:186:GOL:H32	1.85	0.57
1:B:32:ASN:HB2	1:B:34:GLY:H	1.69	0.57
1:B:97:THR:H	2:B:187:GOL:H31	1.75	0.52
1:B:13:GLY:C	1:B:33:THR:OG1	2.48	0.52
1:B:6:TRP:CH2	1:B:38:VAL:HG13	2.46	0.51
1:B:9:TRP:HH2	1:B:166:TYR:OH	1.88	0.51
1:A:24:GLY:HA2	1:A:40:LYS:HD3	1.93	0.50
1:A:6:TRP:CH2	1:A:38:VAL:CG1	2.94	0.50
1:A:33:THR:O	1:A:175:GLN:HA	2.12	0.49
1:A:7:GLN:HG3	1:A:9:TRP:HD1	1.81	0.46
1:A:17:ASN:HB2	4:B:284:HOH:O	2.16	0.45
1:B:71:TRP:HZ3	1:B:73:ARG:CG	2.30	0.44
1:A:78:GLU:HB3	1:A:127:GLN:HG2	2.00	0.43
1:A:9:TRP:HH2	3:B:186:MES:H71	1.86	0.41
1:B:166:TYR:HB3	4:B:242:HOH:O	2.21	0.40
1:A:78:GLU:O	1:A:127:GLN:HA	2.20	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	ed Favoured A		Allowed Outliers			
1	А	184/185~(100%)	180 (98%)	4 (2%)	0	100 100	)	

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
1	В	170/185~(92%)	166 (98%)	3~(2%)	1 (1%)	25	12
All	All	354/370~(96%)	346~(98%)	7~(2%)	1 (0%)	41	27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	33	THR

#### 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	Percentile	es
1	А	152/151~(101%)	150~(99%)	2(1%)	69 61	
1	В	142/151~(94%)	140~(99%)	2(1%)	67 58	
All	All	294/302~(97%)	290~(99%)	4 (1%)	67 58	

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	7	GLN
1	А	111	THR
1	В	7	GLN
1	В	40	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	156	HIS
1	А	159	ASN

#### 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 monosaccharides in this entry.

### 5.6 Ligand geometry (i)

3 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).

Mal	Mol Type Chain Res		Chain Res Link		Bo	Bond lengths			Bond angles		
	туре	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z  > 2	
3	MES	В	186	-	12,12,12	2.21	1 (8%)	$14,\!16,\!16$	2.61	7 (50%)	
2	GOL	А	186	-	$5,\!5,\!5$	0.43	0	$5,\!5,\!5$	0.78	0	
2	GOL	В	187	-	$5,\!5,\!5$	0.32	0	$5,\!5,\!5$	0.45	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
3	MES	В	186	-	-	5/6/14/14	0/1/1/1
2	GOL	А	186	-	-	4/4/4/4	-
2	GOL	В	187	-	-	3/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	В	186	MES	C8-S	-7.35	1.67	1.77

All (7) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	В	186	MES	C5-N4-C3	5.07	120.24	108.83
3	В	186	MES	O1S-S-C8	4.02	111.76	106.92
3	В	186	MES	C7-N4-C5	3.63	120.51	111.23
3	В	186	MES	C7-N4-C3	3.13	119.25	111.23
3	В	186	MES	C2-C3-N4	-2.93	105.67	110.10
3	В	186	MES	C6-C5-N4	-2.75	105.93	110.10
3	В	186	MES	O3S-S-C8	2.00	109.01	105.77

There are no chirality outliers.

Mol	Chain	Res	Type	Atoms
3	В	186	MES	N4-C7-C8-S
3	В	186	MES	C7-C8-S-O2S
3	В	186	MES	C7-C8-S-O3S
2	А	186	GOL	O1-C1-C2-C3
2	В	187	GOL	C1-C2-C3-O3
2	А	186	GOL	O1-C1-C2-O2
2	В	187	GOL	O2-C2-C3-O3
2	А	186	GOL	O2-C2-C3-O3
2	В	187	GOL	O1-C1-C2-O2
2	А	186	GOL	C1-C2-C3-O3
3	В	186	MES	C7-C8-S-O1S
3	В	186	MES	C8-C7-N4-C3

All (12) torsion outliers are listed below:

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	В	186	MES	1	0
2	А	186	GOL	2	0
2	В	187	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 (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,  $95^{th}$  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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	$Q{<}0.9$
1	А	185/185~(100%)	0.26	10 (5%) 25 20	11, 17, 26, 29	0
1	В	173/185~(93%)	0.27	9 (5%) 27 22	10, 16, 22, 34	0
All	All	358/370~(96%)	0.26	19 (5%) 26 21	10, 17, 26, 34	0

All (19) RSRZ outliers are listed below:

Mol	Mol Chain		Type	RSRZ
1	В	124	THR	7.1
1	В	1	ALA	6.0
1	В	33	THR	5.0
1	В	125	PHE	4.9
1	А	9	TRP	4.7
1	В	71	TRP	4.4
1	А	174	TYR	4.1
1	А	1	ALA	3.4
1	В	174	TYR	3.3
1	В	111	THR	3.2
1	А	71	TRP	3.2
1	А	119	ASP	2.9
1	А	113	TYR	2.9
1	А	2	SER	2.9
1	В	44	THR	2.6
1	В	166	TYR	2.5
1	А	112	ARG	2.5
1	А	125	PHE	2.1
1	А	111	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)

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,  $95^{th}$  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	$B-factors(Å^2)$	Q < 0.9
2	GOL	В	187	6/6	0.53	0.32	40,41,42,43	0
2	GOL	А	186	6/6	0.65	0.26	28,29,30,31	0
3	MES	В	186	12/12	0.67	0.34	94,94,95,95	0

### 6.5 Other polymers (i)

There are no such residues in this entry.

