

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

#### Aug 6, 2020 – 11:35 PM BST

PDB ID	:	1REE
$\operatorname{Title}$	:	ENDO-1,4-BETA-XYLANASE II COMPLEX WITH 3,4-EPOXYBUTYL-B
		ETA-D-XYLOSIDE
Authors	:	Rouvinen, J.; Havukainen, R.; Torronen, A.
Deposited on	:	1995-12-21
Resolution	:	1.60  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 following versions of software and data (see references (1)) were used in the production of this report:

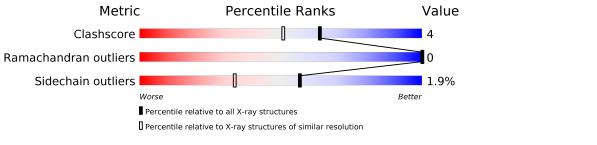
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

## 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
Clashscore	141614	3665(1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563(1.60-1.60)

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 >=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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain					
1	А	190	87%	12% •				
1	В	190	91%	9%				



## 2 Entry composition (i)

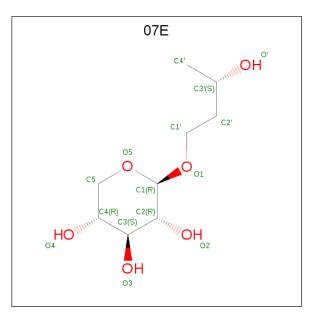
There are 4 unique types of molecules in this entry. The entry contains 3222 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 II.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Δ	190	Total	С	Ν	Ο	S	0	0	0
		190	1480	934	253	292	1	0	0	
1	р	190	Total	С	Ν	Ο	S	0	0	0
	D	190	1480	934	253	292	1	0	0	0

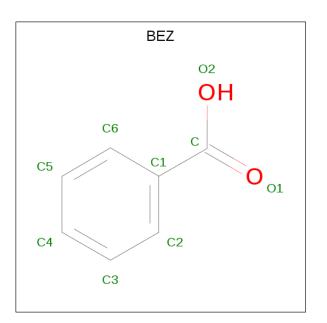
• Molecule 2 is (3S)-3-hydroxy butyl beta-D-xylopyranoside (three-letter code: 07E) (formula:  $\rm C_9H_{18}O_6).$ 



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
2	В	1	Total 15	С 9	O 6	0	0

• Molecule 3 is BENZOIC ACID (three-letter code: BEZ) (formula:  $C_7H_6O_2$ ).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	В	1	Total 9	$\begin{array}{c} \mathrm{C} \\ 7 \end{array}$	O 2	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	118	Total O 118 118	0	0
4	В	120	Total         O           120         120	0	0

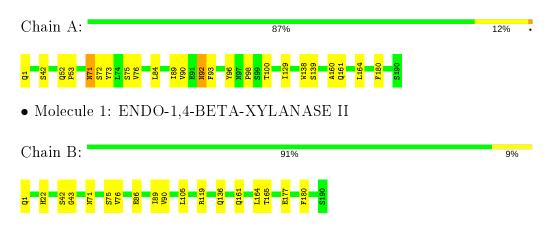


## 3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

• Molecule 1: ENDO-1,4-BETA-XYLANASE II





## 4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source	
Space group	P 1 21 1	Depositor	
Cell constants	81.84Å $60.81$ Å $38.09$ Å	Depositor	
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $94.20^{\circ}$ $90.00^{\circ}$	Depositor	
Resolution (Å)	8.00 - 1.60	Depositor	
% Data completeness	(Not available) (8.00-1.60)	Depositor	
(in resolution range)	(1007 available) (0.00-1.00)	Depositor	
$R_{merge}$	0.09	Depositor	
$R_{sym}$	(Not available)	Depositor	
Refinement program	X-PLOR	Depositor	
$R, R_{free}$	0.181 , $0.223$	Depositor	
Estimated twinning fraction	No twinning to report.	Xtriage	
Total number of atoms	3222	wwPDB-VP	
Average B, all atoms $(Å^2)$	21.0	wwPDB-VP	



# 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: BEZ, PCA,  $07\mathrm{E}$ 

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	Boi	nd lengths	Bond angles		
	Unam	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.36	0/1519	0.65	0/2073	
1	В	0.55	1/1519~(0.1%)	0.64	0/2073	
All	All	0.46	1/3038~(0.0%)	0.65	0/4146	

All (1) bond length outliers are listed below:

Mol	Chain	$\mathbf{Res}$	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	В	177	$\operatorname{GLU}$	CD-OE2	16.32	1.43	1.25

There are no bond angle outliers.

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	1480	0	1350	14	0
1	В	1480	0	1350	10	0
2	В	15	0	0	0	0
3	В	9	0	5	0	0
4	А	118	0	0	0	0
4	В	120	0	0	0	0
All	All	3222	0	2705	24	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 (24) 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:75:SER:HB3	1:A:90:VAL:HA	1.62	0.80
1:A:73:TYR:CZ	1:A:98:PRO:HB3	2.35	0.62
1:B:75:SER:HB3	1:B:90:VAL:HA	1.85	0.59
1:B:105:LEU:HD11	1:B:119:ARG:HB2	1.86	0.56
1:A:92:ASN:HD22	1:A:93:PHE:H	1.54	0.56
1:A:98:PRO:HB2	1:A:138:TRP:CZ3	2.43	0.54
1:B:76:VAL:HB	1:B:89:ILE:HB	1.91	0.53
1:A:73:TYR:CE1	1:A:98:PRO:HB3	2.46	0.51
1:A:160:ALA:HA	1:A:164:LEU:O	2.11	0.51
1:A:96:TYR:OH	1:A:100:THR:HB	2.12	0.50
1:A:42:SER:O	1:A:180:PHE:HA	2.12	0.49
1:A:84:LEU:HD22	1:A:129:ILE:HD11	1.96	0.47
1:B:164:LEU:HD12	1:B:165:THR:H	1.79	0.46
1:B:42:SER:O	1:B:180:PHE:HA	2.15	0.46
1:A:76:VAL:HB	1:A:89:ILE:HB	1.98	0.45
1:B:165:THR:O	1:B:165:THR:HG23	2.17	0.45
1:B:22:HIS:HB3	1:B:43:GLY:HA3	2.00	0.44
1:B:86:GLU:O	1:B:136:GLN:HA	2.17	0.44
1:A:92:ASN:HD22	1:A:93:PHE:N	2.16	0.43
1:A:71:ASN:HD22	1:A:72:SER:N	2.17	0.43
1:A:52:GLN:HA	1:A:53:PRO:HA	1.94	0.41
1:B:164:LEU:HD12	1:B:165:THR:N	2.36	0.41
1:A:71:ASN:C	1:A:71:ASN:HD22	2.23	0.41
1:B:75:SER:CB	1:B:90:VAL:HA	2.49	0.41

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	А	188/190~(99%)	181 (96%)	7 (4%)	0	100	100
1	В	188/190~(99%)	180~(96%)	8 (4%)	0	100	100
All	All	376/380~(99%)	361~(96%)	15~(4%)	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	А	155/155~(100%)	151~(97%)	4 (3%)	46 21		
1	В	155/155~(100%)	153~(99%)	2(1%)	69 50		
All	All	310/310~(100%)	304~(98%)	6~(2%)	57 34		

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	71	ASN
1	А	92	ASN
1	А	139	SER
1	А	161	GLN
1	В	71	ASN
1	В	161	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	41	ASN
1	А	71	ASN
1	А	92	ASN
1	А	144	HIS
1	В	38	ASN
1	В	71	ASN
1	В	121	GLN

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Mol	Chain	Res	Type
1	В	124	ASN
1	В	157	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)

2 non-standard protein/DNA/RNA residues 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	Link	Bond lengths			Bond angles		
	туре	Chain	nes		Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z >2
1	PCA	А	1	1	7,8,9	2.55	3 (42%)	$9,\!10,\!12$	1.78	<mark>3 (33%)</mark>
1	PCA	В	1	1	7,8,9	2.59	3 (42%)	$9,\!10,\!12$	1.78	<mark>3 (33%)</mark>

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
1	PCA	А	1	1	-	0/0/11/13	0/1/1/1
1	PCA	В	1	1	-	0/0/11/13	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
1	В	1	PCA	CD-N	5.59	1.49	1.34
1	А	1	PCA	CD-N	5.41	1.48	1.34
1	В	1	PCA	CA-N	3.14	1.50	1.46
1	А	1	PCA	CA-N	3.09	1.50	1.46
1	А	1	PCA	CB-CG	-2.34	1.47	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	В	1	PCA	CB-CG	-2.19	1.48	1.53

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$\mathbf{Ideal}(^{o})$
1	А	1	PCA	OE-CD-CG	-2.53	122.35	126.76
1	В	1	PCA	OE-CD-CG	-2.51	122.38	126.76
1	А	1	PCA	CB-CA-C	-2.45	109.33	112.70
1	В	1	PCA	CB-CA-C	-2.35	109.47	112.70
1	В	1	PCA	CB-CG-CD	2.09	107.78	104.40
1	А	1	PCA	CA-N-CD	-2.05	106.57	113.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

#### 5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry (i)

2 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 Ch	Chain	Dec	Res Link	Bo	Bond lengths			Bond angles		
	Type	Chain	nes		Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2	
2	07E	В	401	1	15, 15, 15	1.18	1(6%)	$20,\!20,\!20$	1.44	3(15%)	
3	BEZ	В	402	-	7,9,9	2.90	1 (14%)	8,11,11	0.30	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 a	re reported in the	Torsion and	Rings columns.
'-' means no outliers of that kind	were identified.			

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	07E	В	401	1	-	2/6/23/23	0/1/1/1
3	BEZ	В	402	-	-	0/0/4/4	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	$\mathbf{Res}$	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
3	В	402	BEZ	С1-С	-7.55	1.40	1.47
2	В	401	07E	O1-C1	3.00	1.45	1.40

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
2	В	401	$07\mathrm{E}$	O1-C1-C2	3.56	113.86	108.30
2	В	401	07E	C1'-C2'-C3'	3.27	118.96	113.92
2	В	401	07E	O1-C1'-C2'	2.64	114.30	108.77

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	В	401	07E	O5-C1-O1-C1'
2	В	401	07E	C2-C1-O1-C1'

There are no ring outliers.

No monomer is involved in short contacts.

### 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)

EDS was not executed - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

#### 6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

### 6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

#### 6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

