

Full wwPDB X-ray Structure Validation Report (i)

Dec 8, 2023 - 07:53 am GMT

PDB ID	:	8BRY
Title	:	Room-temperature structure of Pedobacter heparinus N-acetylglucosamine 2-
		epimerase at atmospheric pressure
Authors	:	Lieske, J.; Saouane, S.; Assmann, M.; Zaun, H.; Kuballa, J.; Meents, A.
Deposited on		
Resolution	:	1.50 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:

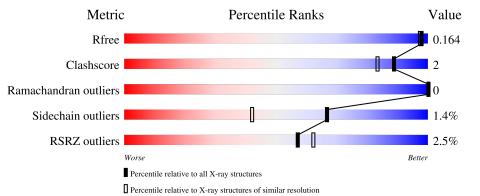
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as 541 be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 1.50 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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	418	^{2%} 90%	• 7%			
1	В	418	^{2%} 87%	5% 8%			



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2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 13517 atoms, of which 6356 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 N-acylglucosamine 2-epimerase.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace			
1	Λ	388	Total	С	Η	Ν	0	\mathbf{S}	0	22	0
1	Л		6645	2178	3241	583	627	16			
1	В	386	Total	С	Η	Ν	Ο	S	0	0	0
1	D	500	6388	2101	3115	557	599	16	0	9	U

Chain	Residue	Modelled	Actual	Comment	Reference
А	-14	MET	-	initiating methionine	UNP C6Y403
А	-13	GLY	-	expression tag	UNP C6Y403
А	-12	SER	-	expression tag	UNP C6Y403
А	-11	SER	-	expression tag	UNP C6Y403
А	-10	HIS	-	expression tag	UNP C6Y403
А	-9	HIS	-	expression tag	UNP C6Y403
А	-8	HIS	-	expression tag	UNP C6Y403
А	-7	HIS	-	expression tag	UNP C6Y403
А	-6	HIS	-	expression tag	UNP C6Y403
А	-5	HIS	-	expression tag	UNP C6Y403
А	-4	SER	-	expression tag	UNP C6Y403
А	-3	GLN	-	expression tag	UNP C6Y403
А	-2	ASP	-	expression tag	UNP C6Y403
А	-1	PRO	-	expression tag	UNP C6Y403
А	0	ASN	-	expression tag	UNP C6Y403
А	1	SER	-	expression tag	UNP C6Y403
В	-14	MET	-	initiating methionine	UNP C6Y403
В	-13	GLY	-	expression tag	UNP C6Y403
В	-12	SER	-	expression tag	UNP C6Y403
В	-11	SER	-	expression tag	UNP C6Y403
В	-10	HIS	-	expression tag	UNP C6Y403
В	-9	HIS	-	expression tag	UNP C6Y403
В	-8	HIS	-	expression tag	UNP C6Y403
В	-7	HIS	-	expression tag	UNP C6Y403
В	-6	HIS	-	expression tag	UNP C6Y403

There are 32 discrepancies between the modelled and reference sequences:

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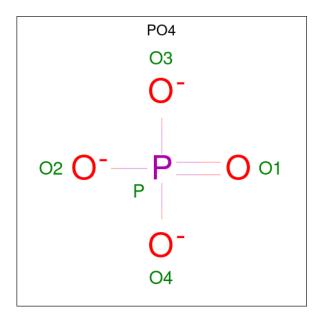
Chain	Residue	Modelled	Actual	Comment	Reference
В	-5	HIS	-	expression tag	UNP C6Y403
В	-4	SER	-	expression tag	UNP C6Y403
В	-3	GLN	-	expression tag	UNP C6Y403
В	-2	ASP	-	expression tag	UNP C6Y403
В	-1	PRO	-	expression tag	UNP C6Y403
В	0	ASN	-	expression tag	UNP C6Y403
B	1	SER	-	expression tag	UNP C6Y403

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• Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Μ	ol	Chain	Residues	Atoms	ZeroOcc	AltConf
2 2	2	А	1	Total Cl 1 1	0	0
2	2	В	1	Total Cl 1 1	0	0

• Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0
3	В	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{P} \\ 5 & 4 & 1 \end{array}$	0	0

• Molecule 4 is water.

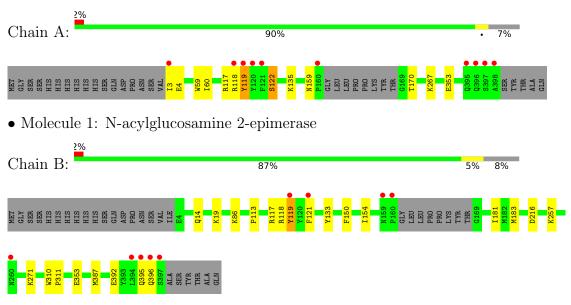


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	262	Total O 262 262	0	0
4	В	210	Total O 210 210	0	0



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: N-acylglucosamine 2-epimerase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	56.95Å 94.63Å 82.47Å	Depositor
a, b, c, α , β , γ	90.00° 102.53° 90.00°	Depositor
Resolution (Å)	17.62 - 1.50	Depositor
Resolution (A)	47.93 - 1.30	EDS
% Data completeness	99.8 (17.62-1.50)	Depositor
(in resolution range)	76.0(47.93-1.30)	EDS
R _{merge}	0.18	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	-0.06 (at 1.30Å)	Xtriage
Refinement program	PHENIX 1.18-3855_9999	Depositor
D D	0.154 , 0.164	Depositor
R, R_{free}	0.154 , 0.164	DCC
R_{free} test set	1306 reflections (0.63%)	wwPDB-VP
Wilson B-factor $(Å^2)$	13.8	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.44 , 53.6	EDS
L-test for twinning ²	$ \langle L \rangle = 0.51, \langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	13517	wwPDB-VP
Average B, all atoms $(Å^2)$	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.15% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities 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: CL, $\rm PO4$

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		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.46	0/3500	0.67	0/4739	
1	В	0.43	0/3369	0.66	0/4563	
All	All	0.45	0/6869	0.67	0/9302	

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 (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	3404	3241	3220	14	0
1	В	3273	3115	3103	12	0
2	А	1	0	0	0	0
2	В	1	0	0	0	0
3	А	5	0	0	0	0
3	В	5	0	0	0	0
4	А	262	0	0	3	0
4	В	210	0	0	0	0
All	All	7161	6356	6323	25	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 2.



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:ARG:HG3	1:A:117:ARG:HH11	1.51	0.75
1:A:117:ARG:HD3	1:A:119:TYR:CZ	2.31	0.64
1:A:170:THR:HG23	1:B:113:PRO:O	1.99	0.63
1:B:121:PHE:HE1	1:B:181:ILE:HG21	1.66	0.59
1:B:121:PHE:CE1	1:B:181:ILE:HG21	2.38	0.58
1:A:60[B]:ILE:HD12	4:A:691:HOH:O	2.05	0.56
1:A:117:ARG:HD3	1:A:119:TYR:CE1	2.42	0.54
1:B:86:LYS:HD2	1:B:133:TYR:OH	2.10	0.52
1:B:216:ASP:OD1	1:B:271[A]:LYS:HE2	2.11	0.50
1:A:117:ARG:HH11	1:A:117:ARG:CG	2.22	0.47
1:B:396:GLN:HA	1:B:396:GLN:OE1	2.14	0.47
1:A:119:TYR:N	1:A:119:TYR:CD2	2.82	0.47
1:B:19:LYS:HE2	1:B:19:LYS:HB3	1.85	0.46
1:A:159:ASN:ND2	1:A:159:ASN:O	2.51	0.44
1:A:135[B]:LYS:CE	4:A:738:HOH:O	2.66	0.44
1:B:14:GLN:NE2	1:B:387:MET:HG3	2.34	0.43
1:B:310:TRP:CG	1:B:311:PRO:HD3	2.54	0.42
1:A:267:LYS:HB2	1:A:267:LYS:HE2	1.78	0.42
1:B:257:LYS:NZ	1:B:392:GLU:OE2	2.45	0.42
1:A:59:TRP:HE1	1:A:122:SER:HG	1.65	0.42
1:A:59:TRP:CD1	1:A:122:SER:OG	2.68	0.41
1:A:135[B]:LYS:HE3	4:A:738:HOH:O	2.21	0.41
1:A:3:ILE:CG2	1:A:4:GLU:N	2.83	0.41
1:B:117:ARG:NE	1:B:119:TYR:CE2	2.89	0.41
1:B:150:PHE:CE2	1:B:154:ILE:HD11	2.56	0.40

All (25) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

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	406/418~(97%)	397~(98%)	9~(2%)	0	100	100
1	В	391/418 (94%)	382 (98%)	9~(2%)	0	100	100
All	All	797/836~(95%)	779~(98%)	18 (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 side chain 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 side chain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	А	352/357~(99%)	348~(99%)	4 (1%)	73 53
1	В	338/357~(95%)	333~(98%)	5(2%)	65 39
All	All	690/714~(97%)	681~(99%)	9~(1%)	67 44

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	118	ARG
1	А	119	TYR
1	А	122	SER
1	А	353	GLU
1	В	118	ARG
1	В	119	TYR
1	В	183	MET
1	В	353	GLU
1	В	395	GLN

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such side chains are listed below:

Mol	Chain	Res	Type
1	А	159	ASN
1	В	14	GLN
1	В	79	GLN
1	В	102	GLN

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Mol	Chain	\mathbf{Res}	Type
1	В	296	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)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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 Li	Link	Bond lengths			Bond angles		
IVIOI	Mol Type Chain	LIIIK		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2	
3	PO4	А	502	-	4,4,4	0.80	0	$6,\!6,\!6$	0.71	0
3	PO4	В	502	-	4,4,4	0.83	0	$6,\!6,\!6$	0.42	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 (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



$6 \quad \text{Fit of model and data} \quad (i)$

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	388/418~(92%)	-0.42	10 (2%) 56 61	14, 21, 48, 100	0
1	В	386/418~(92%)	-0.36	9 (2%) 60 65	15, 27, 56, 97	0
All	All	774/836~(92%)	-0.39	19 (2%) 57 62	14, 24, 54, 100	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	159	ASN	10.6
1	А	160	PRO	6.8
1	В	119	TYR	6.0
1	А	120	TYR	5.7
1	В	397	SER	5.2
1	А	3	ILE	4.9
1	А	398	ALA	4.6
1	А	118	ARG	4.0
1	А	396	GLN	3.7
1	В	394	LEU	3.7
1	А	119	TYR	3.7
1	В	396	GLN	3.5
1	А	397	SER	3.2
1	В	395	GLN	2.7
1	В	121	PHE	2.7
1	В	260	ASN	2.6
1	В	160	PRO	2.4
1	А	121	PHE	2.4
1	А	395[A]	GLN	2.2

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
3	PO4	В	502	5/5	0.91	0.20	$31,\!37,\!52,\!52$	5
3	PO4	А	502	5/5	0.97	0.10	34,47,69,81	0
2	CL	А	501	1/1	1.00	0.07	14,14,14,14	0
2	CL	В	501	1/1	1.00	0.04	$17,\!17,\!17,\!17$	0

6.5 Other polymers (i)

There are no such residues in this entry.

