



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 3GM8 / pdb_00003gm8
Title : Crystal structure of a beta-glycosidase from Bacteroides vulgatus
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Deposited on : 2009-03-13
Resolution : 2.40 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

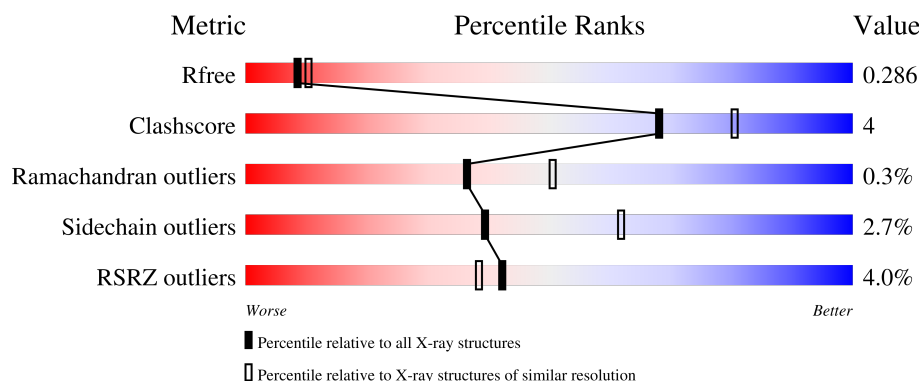
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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 $\leq 5\%$. 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	801	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6380 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 Glycoside hydrolase family 2, candidate beta-glycosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	770	6112	3926	1040	1121	25	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	17	SER	-	expression tag	UNP A6KXE5
A	18	LEU	-	expression tag	UNP A6KXE5
A	816	GLU	-	expression tag	UNP A6KXE5
A	817	GLY	-	expression tag	UNP A6KXE5

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	6	3	3	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		

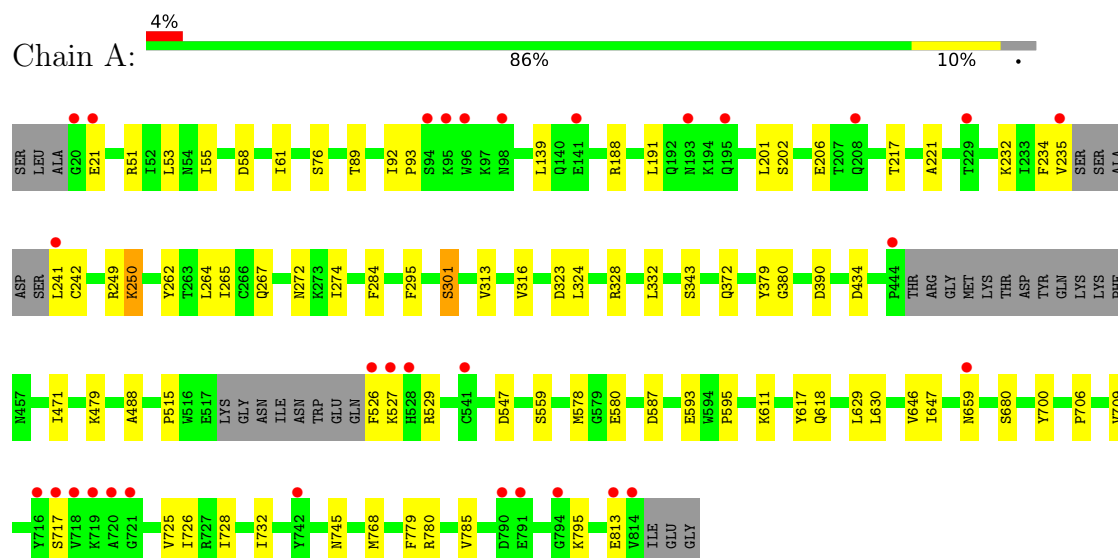
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	256	Total	O	0	0
			256	256		

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: Glycoside hydrolase family 2, candidate beta-glycosidase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	62.75Å 193.35Å 190.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 20.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.2 (20.00-2.40) 97.3 (20.00-2.40)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.83 (at 2.39Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.226 , 0.287 0.235 , 0.286	Depositor DCC
R_{free} test set	2281 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	31.8	Xtriage
Anisotropy	0.659	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 41.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6380	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

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

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	0.96	1/6295 (0.0%)	0.93	1/8571 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	785	VAL	CA-CB	6.01	1.62	1.54

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	21	GLU	N-CA-C	5.48	117.17	108.79

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	A	6112	0	5870	43	0
2	A	12	0	16	0	0
3	A	256	0	0	4	0
All	All	6380	0	5886	43	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 (43) 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:659:ASN:ND2	3:A:1005:HOH:O	2.03	0.91
1:A:201:LEU:HD21	1:A:264:LEU:HD21	1.56	0.87
1:A:301:SER:HB2	3:A:878:HOH:O	1.75	0.86
1:A:295:PHE:HB2	1:A:578:MET:HE3	1.69	0.73
1:A:267:GLN:NE2	1:A:274:ILE:HD13	2.05	0.70
1:A:55:ILE:C	1:A:55:ILE:HD12	2.20	0.67
1:A:526:PHE:N	3:A:856:HOH:O	2.31	0.62
1:A:726:ILE:HG22	1:A:728:ILE:HD11	1.86	0.56
1:A:629:LEU:HD22	1:A:700:TYR:HD2	1.72	0.55
1:A:313:VAL:HG13	1:A:328:ARG:CZ	2.38	0.54
1:A:201:LEU:HD21	1:A:264:LEU:CD2	2.35	0.52
1:A:191:LEU:HD21	1:A:284:PHE:HB3	1.93	0.51
1:A:313:VAL:HG12	1:A:313:VAL:O	2.11	0.50
1:A:234:PHE:CD1	1:A:242:CYS:HB2	2.48	0.49
1:A:324:LEU:HD13	1:A:725:VAL:HG11	1.94	0.48
1:A:58:ASP:O	1:A:61:ILE:HG22	2.13	0.48
1:A:717:SER:HA	1:A:813:GLU:O	2.15	0.47
1:A:51:ARG:CZ	1:A:53:LEU:HD23	2.45	0.47
1:A:221:ALA:HA	1:A:262:TYR:CD1	2.50	0.47
1:A:745:ASN:HD22	1:A:745:ASN:H	1.62	0.46
1:A:249:ARG:O	1:A:250:LYS:C	2.58	0.46
1:A:313:VAL:HG13	1:A:328:ARG:NE	2.30	0.46
1:A:515:PRO:HG3	1:A:593:GLU:CD	2.41	0.45
1:A:587:ASP:OD2	1:A:611:LYS:NZ	2.42	0.45
1:A:726:ILE:HG22	1:A:728:ILE:CD1	2.46	0.45
1:A:728:ILE:N	1:A:728:ILE:HD12	2.31	0.44
1:A:526:PHE:HD1	1:A:529:ARG:HD2	1.82	0.44
1:A:779:PHE:O	1:A:780:ARG:C	2.59	0.44
1:A:313:VAL:O	1:A:316:VAL:HG22	2.18	0.44
1:A:595:PRO:O	1:A:768:MET:HE3	2.17	0.44
1:A:217:THR:HA	1:A:265:ILE:O	2.19	0.43
1:A:92:ILE:HG22	1:A:93:PRO:O	2.20	0.41
1:A:202:SER:OG	1:A:241:LEU:HD22	2.21	0.41
1:A:488:ALA:HB3	1:A:580:GLU:HA	2.01	0.41
1:A:709:VAL:HG12	1:A:732:ILE:HD13	2.01	0.41
1:A:629:LEU:HD22	1:A:700:TYR:CD2	2.54	0.41
1:A:332:LEU:HD23	1:A:332:LEU:HA	1.74	0.41
1:A:188:ARG:NE	1:A:434:ASP:OD2	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:706:PRO:HG2	3:A:835:HOH:O	2.20	0.41
1:A:92:ILE:HD11	1:A:139:LEU:HB3	2.02	0.41
1:A:646:VAL:C	1:A:647:ILE:HD12	2.46	0.41
1:A:379:TYR:OH	1:A:390:ASP:OD2	2.27	0.40
1:A:617:TYR:O	1:A:618:GLN:C	2.61	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	Favoured	Allowed	Outliers	Percentiles
1	A	762/801 (95%)	729 (96%)	31 (4%)	2 (0%)	36 50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	380	GLY
1	A	527	LYS

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	655/701 (93%)	637 (97%)	18 (3%)	39 62

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	SER
1	A	89	THR
1	A	206	GLU
1	A	232	LYS
1	A	235	VAL
1	A	250	LYS
1	A	272	ASN
1	A	301	SER
1	A	323	ASP
1	A	343	SER
1	A	372	GLN
1	A	471	ILE
1	A	479	LYS
1	A	547	ASP
1	A	559	SER
1	A	630	LEU
1	A	680	SER
1	A	795	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	140	GLN
1	A	186	HIS
1	A	195	GLN
1	A	267	GLN
1	A	272	ASN
1	A	372	GLN
1	A	504	GLN
1	A	707	HIS
1	A	745	ASN
1	A	750	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GOL	A	1	-	5,5,5	0.52	0	5,5,5	0.48	0
2	GOL	A	2	-	5,5,5	0.40	0	5,5,5	0.53	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
2	GOL	A	1	-	-	1/4/4/4	-
2	GOL	A	2	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1	GOL	C1-C2-C3-O3

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

6.1 Protein, DNA and RNA chains

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, 95th 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 ²)	Q<0.9
1	A	770/801 (96%)	0.37	31 (4%) 42 38	36, 50, 69, 78	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	718	VAL	4.3
1	A	95	LYS	4.3
1	A	791	GLU	4.0
1	A	98	ASN	3.7
1	A	717	SER	3.6
1	A	720	ALA	3.5
1	A	20	GLY	3.5
1	A	790	ASP	3.4
1	A	235	VAL	3.4
1	A	526	PHE	3.3
1	A	193	ASN	3.3
1	A	141	GLU	3.1
1	A	195	GLN	2.9
1	A	814	VAL	2.8
1	A	444	PRO	2.6
1	A	96	TRP	2.5
1	A	229	THR	2.5
1	A	541	CYS	2.4
1	A	94	SER	2.3
1	A	241	LEU	2.3
1	A	716	TYR	2.3
1	A	208	GLN	2.3
1	A	528	HIS	2.3
1	A	719	LYS	2.2
1	A	813	GLU	2.1
1	A	659	ASN	2.1
1	A	721	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	21	GLU	2.0
1	A	794	GLY	2.0
1	A	527	LYS	2.0
1	A	742	TYR	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 oligosaccharides 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, 95th 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	GOL	A	2	6/6	0.90	0.13	66,67,69,70	0
2	GOL	A	1	6/6	0.95	0.10	42,45,51,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.