



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:07 AM UTC

PDB ID : 8K2G / pdb_00008k2g
Title : Crystal structure of Group 1 Oligosaccharide-releasing beta-N-acetylgalactosaminidase NgaCa from Cohnella abietis, apo 2 form
Authors : Sumida, T.; Fushinobu, S.
Deposited on : 2023-07-12
Resolution : 1.75 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
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

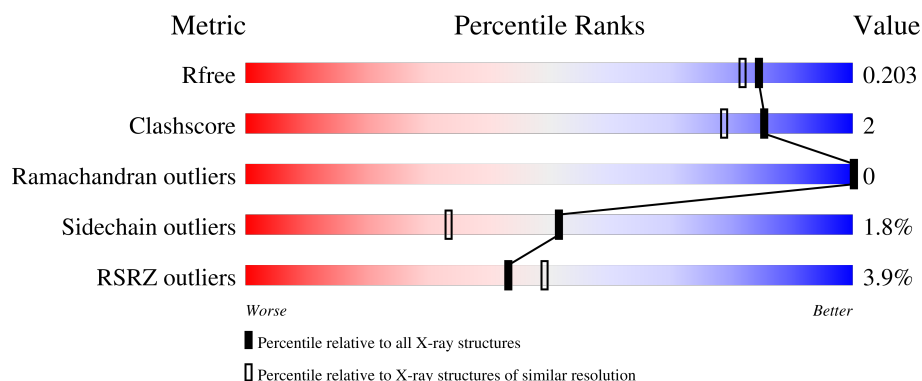
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	516	3% 90% 8%
1	B	516	4% 90% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8766 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 Oligosaccharide-releasing beta-N-acetylgalactosaminidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	505	Total	C	N	O	S	0	1	0
			4035	2578	671	775	11			
1	A	505	Total	C	N	O	S	0	1	0
			4035	2578	671	775	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	GLY	-	expression tag	UNP A0A3T1D3V8
B	-1	PRO	-	expression tag	UNP A0A3T1D3V8
B	0	GLY	-	expression tag	UNP A0A3T1D3V8
A	-2	GLY	-	expression tag	UNP A0A3T1D3V8
A	-1	PRO	-	expression tag	UNP A0A3T1D3V8
A	0	GLY	-	expression tag	UNP A0A3T1D3V8

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

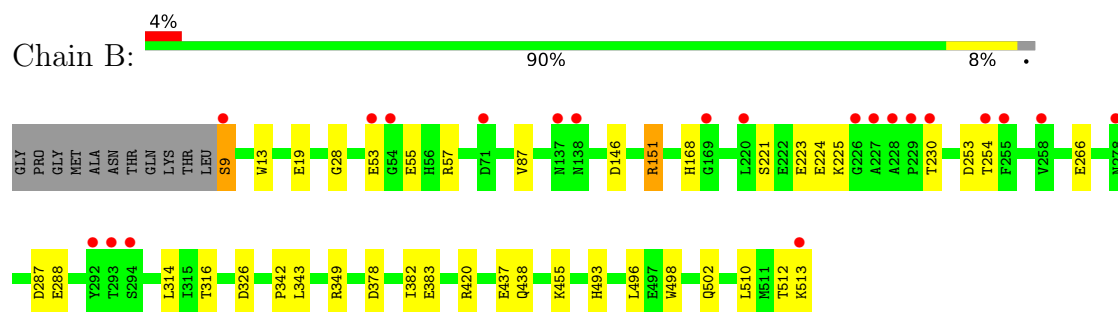
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	389	Total	O	0	0
			389	389		
3	A	305	Total	O	0	0
			305	305		

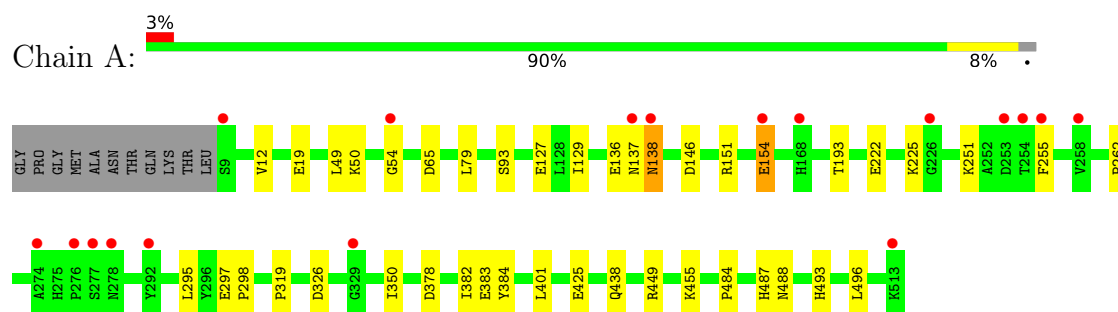
3 Residue-property plots [i](#)

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: Oligosaccharide-releasing beta-N-acetylgalactosaminidase



- Molecule 1: Oligosaccharide-releasing beta-N-acetylgalactosaminidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	53.21Å 69.68Å 71.11Å 88.41° 74.33° 87.93°	Depositor
Resolution (Å)	47.67 – 1.75 47.67 – 1.75	Depositor EDS
% Data completeness (in resolution range)	97.1 (47.67-1.75) 97.1 (47.67-1.75)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.12 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.8.0415	Depositor
R, R_{free}	0.156 , 0.193 0.169 , 0.203	Depositor DCC
R_{free} test set	4810 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	15.0	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.031 for -h,k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8766	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 5.80% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG

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	0.64	0/4156	1.02	7/5666 (0.1%)
1	B	0.69	0/4156	1.07	8/5666 (0.1%)
All	All	0.66	0/8312	1.05	15/11332 (0.1%)

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	2
1	B	0	3
All	All	0	5

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	65	ASP	CA-CB-CG	7.05	119.65	112.60
1	B	151	ARG	CG-CD-NE	-6.84	96.95	112.00
1	B	326	ASP	CA-CB-CG	6.68	119.28	112.60
1	B	266	GLU	CB-CG-CD	6.58	123.78	112.60
1	A	378	ASP	CA-CB-CG	6.46	119.06	112.60
1	A	326	ASP	CA-CB-CG	6.28	118.88	112.60
1	A	438	GLN	CB-CG-CD	6.15	123.05	112.60
1	B	438	GLN	CB-CG-CD	6.01	122.82	112.60
1	B	378	ASP	CA-CB-CG	5.83	118.43	112.60
1	A	262	PRO	N-CA-C	5.66	117.60	110.70
1	A	383	GLU	CB-CG-CD	5.38	121.74	112.60
1	B	420	ARG	CB-CA-C	5.37	118.74	110.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	383	GLU	CB-CG-CD	5.31	121.62	112.60
1	A	154	GLU	CB-CG-CD	5.10	121.27	112.60
1	B	146	ASP	CA-CB-CG	5.03	117.63	112.60

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	136	GLU	Peptide
1	A	151	ARG	Sidechain
1	B	151	ARG	Sidechain
1	B	349	ARG	Sidechain
1	B	57	ARG	Sidechain

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	4035	0	3857	16	0
1	B	4035	0	3857	18	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	305	0	0	1	0
3	B	389	0	0	4	0
All	All	8766	0	7714	34	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.

All (34) 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:B:87:VAL:HG13	3:B:751:HOH:O	1.90	0.70
1:B:437:GLU:HG3	3:B:881:HOH:O	1.95	0.66
1:B:512:THR:HG22	1:B:512:THR:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:484:PRO:HD2	1:A:488:ASN:O	2.08	0.54
1:A:382:ILE:HD11	1:A:496:LEU:HD22	1.92	0.51
1:A:222:GLU:HA	1:A:225:LYS:HD2	1.92	0.51
1:B:19:GLU:HG3	1:B:493:HIS:CE1	2.46	0.51
1:A:319:PRO:HG2	1:A:350:ILE:HD13	1.93	0.50
1:A:19:GLU:HG3	1:A:493:HIS:CE1	2.48	0.49
1:A:193:THR:CG2	1:A:401:LEU:HD12	2.43	0.48
1:B:382:ILE:HD11	1:B:496:LEU:HD22	1.94	0.48
1:B:254:THR:HG23	1:B:288:GLU:HB2	1.96	0.47
1:B:253:ASP:HA	1:B:287:ASP:O	2.16	0.46
1:A:295:LEU:C	1:A:298:PRO:HD2	2.41	0.46
1:B:221:SER:HB2	1:B:224:GLU:HG2	1.98	0.46
1:B:510:LEU:HD23	1:B:513:LYS:HE2	1.97	0.46
1:B:287:ASP:HA	1:B:316:THR:OG1	2.16	0.45
1:B:9:SER:HA	3:B:978:HOH:O	2.17	0.44
1:A:138:ASN:HD22	1:A:138:ASN:HA	1.60	0.44
1:B:498:TRP:O	1:B:502:GLN:HG2	2.18	0.43
1:B:87:VAL:CG1	3:B:751:HOH:O	2.58	0.43
1:A:54:GLY:O	1:A:137:ASN:OD1	2.36	0.43
1:A:93:SER:HA	1:A:425:GLU:O	2.19	0.43
1:A:12:VAL:HG22	1:A:49:LEU:HD12	2.01	0.42
1:A:487:HIS:HE1	3:A:976:HOH:O	2.01	0.42
1:B:221:SER:HB2	1:B:223:GLU:HG2	2.01	0.42
1:A:19:GLU:CG	1:A:493:HIS:NE2	2.83	0.41
1:B:254:THR:HG23	1:B:288:GLU:CB	2.50	0.41
1:B:13:TRP:CG	1:B:28:GLY:H	2.38	0.41
1:B:314:LEU:C	1:B:314:LEU:HD23	2.46	0.41
1:A:297:GLU:HB2	1:A:298:PRO:HD3	2.03	0.41
1:A:79:LEU:HA	1:A:146:ASP:O	2.21	0.40
1:B:343:LEU:HD12	1:B:343:LEU:C	2.46	0.40
1:A:384:TYR:CZ	1:A:449:ARG:HD2	2.57	0.40

There are no symmetry-related clashes.

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	504/516 (98%)	493 (98%)	11 (2%)	0	100	100
1	B	504/516 (98%)	493 (98%)	11 (2%)	0	100	100
All	All	1008/1032 (98%)	986 (98%)	22 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	436/443 (98%)	428 (98%)	8 (2%)	51	33
1	B	436/443 (98%)	428 (98%)	8 (2%)	51	33
All	All	872/886 (98%)	856 (98%)	16 (2%)	51	33

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	9	SER
1	B	53	GLU
1	B	55	GLU
1	B	168	HIS
1	B	225	LYS
1	B	230	THR
1	B	342	PRO
1	B	455	LYS
1	A	50	LYS
1	A	127	GLU
1	A	129	ILE
1	A	138	ASN
1	A	154	GLU
1	A	251	LYS
1	A	255	PHE

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Mol	Chain	Res	Type
1	A	455	LYS

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

Mol	Chain	Res	Type
1	B	23	GLN
1	B	291	HIS
1	A	23	GLN
1	A	291	HIS
1	A	487	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 ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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

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

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	505/516 (97%)	0.12	18 (3%)	46 52	7, 18, 38, 64	1 (0%)
1	B	505/516 (97%)	-0.08	21 (4%)	40 46	6, 14, 34, 72	1 (0%)
All	All	1010/1032 (97%)	0.02	39 (3%)	43 49	6, 16, 37, 72	2 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	ASN	4.5
1	B	254	THR	4.4
1	A	255	PHE	4.3
1	B	54	GLY	4.2
1	B	138	ASN	3.9
1	A	138	ASN	3.5
1	B	258	VAL	3.4
1	B	255	PHE	3.4
1	B	230	THR	3.4
1	A	513	LYS	3.3
1	B	226	GLY	3.3
1	A	54	GLY	3.2
1	B	293	THR	3.2
1	B	513	LYS	3.0
1	A	258	VAL	3.0
1	A	254	THR	2.9
1	B	9	SER	2.8
1	B	137	ASN	2.8
1	B	294	SER	2.8
1	A	276	PRO	2.7
1	B	71	ASP	2.6
1	A	278	ASN	2.6
1	A	9	SER	2.6
1	A	274	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	220	LEU	2.5
1	A	226	GLY	2.5
1	B	228	ALA	2.4
1	A	277	SER	2.4
1	B	227	ALA	2.3
1	A	168	HIS	2.3
1	B	229	PRO	2.2
1	A	292	TYR	2.2
1	A	329	GLY	2.1
1	B	169	GLY	2.1
1	B	278	ASN	2.1
1	A	154	GLU	2.1
1	A	253	ASP	2.1
1	B	292	TYR	2.1
1	B	53	GLU	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	MG	A	601	1/1	0.99	0.07	16,16,16,16	0
2	MG	B	601	1/1	1.00	0.05	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.