



# Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 2PGV / pdb\_00002pgv  
Title : GTB C209A  
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Deposited on : 2007-04-10  
Resolution : 1.79 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Xtrriage (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

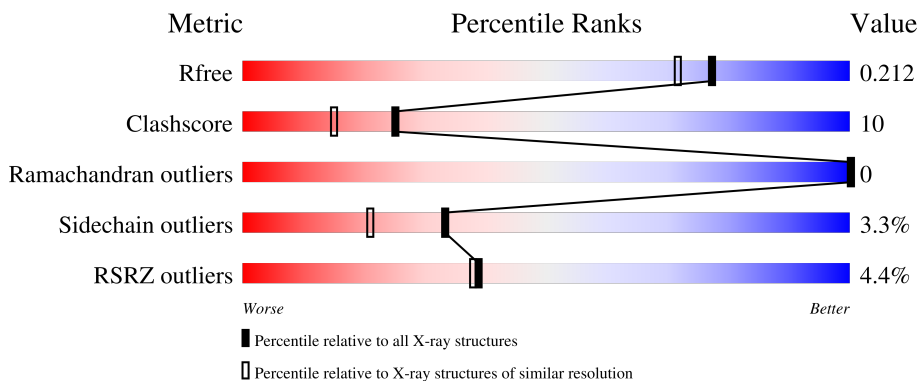
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.79 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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  $\geq 3$ , 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	292	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2421 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 Glycoprotein-fucosylgalactoside alpha-galactosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	271	2227	1444	380	391	12	0	2	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	63	MET	-	initiating methionine	UNP P16442
A	176	GLY	ARG	variant	UNP P16442
A	209	ALA	CYS	engineered mutation	UNP P16442
A	235	SER	GLY	variant	UNP P16442
A	266	MET	LEU	variant	UNP P16442
A	268	ALA	GLY	variant	UNP P16442

- Molecule 2 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Hg	0	0
			3	3		

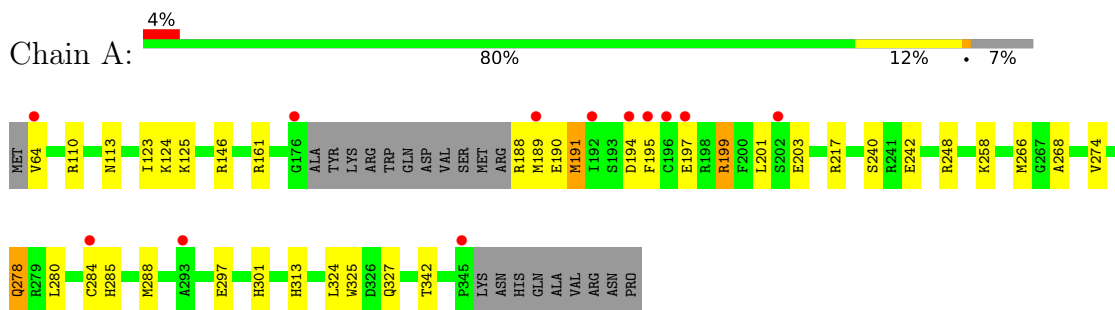
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	191	Total	O	0	0
			191	191		

### 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: Glycoprotein-fucosylgalactoside alpha-galactosyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.50Å 149.26Å 78.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.76 – 1.79 19.76 – 1.79	Depositor EDS
% Data completeness (in resolution range)	93.5 (19.76-1.79) 93.3 (19.76-1.79)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.26 (at 1.79Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.186 , 0.213 0.183 , 0.212	Depositor DCC
$R_{free}$ test set	1409 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.3	Xtrriage
Anisotropy	0.234	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 40.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	2421	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.68% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<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.

## 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: HG

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	0.83	0/2289	0.88	0/3110

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	A	2227	0	2202	43	0
2	A	3	0	0	0	0
3	A	191	0	0	2	0
All	All	2421	0	2202	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 10.

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:284[A]:CYS:SG	1:A:288:MET:HE2	1.82	1.18

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:ARG:HH11	1:A:199:ARG:HG3	1.39	0.86
1:A:284[B]:CYS:O	1:A:288:MET:HG3	1.78	0.84
1:A:240:SER:HB3	1:A:242:GLU:OE2	1.79	0.81
1:A:217:ARG:HH22	1:A:327:GLN:HE21	1.29	0.80
1:A:217:ARG:HH12	1:A:327:GLN:HE22	1.31	0.78
1:A:113[B]:ASN:HD21	1:A:146:ARG:HH21	1.31	0.77
1:A:217:ARG:HH12	1:A:327:GLN:NE2	1.84	0.76
1:A:199:ARG:HH11	1:A:199:ARG:CG	2.05	0.70
1:A:189:MET:HG3	1:A:288:MET:HE1	1.77	0.66
1:A:191:MET:HE3	1:A:195:PHE:HB2	1.77	0.65
1:A:284[A]:CYS:SG	1:A:288:MET:CE	2.74	0.65
1:A:189:MET:HG3	1:A:288:MET:CE	2.28	0.63
1:A:297:GLU:OE1	1:A:301:HIS:HD2	1.80	0.62
1:A:113[B]:ASN:HD21	1:A:146:ARG:NH2	2.01	0.58
1:A:284[A]:CYS:O	1:A:288:MET:HG3	2.03	0.58
1:A:123:ILE:HG22	1:A:124:LYS:HG3	1.86	0.58
1:A:189:MET:CG	1:A:288:MET:HE1	2.35	0.56
1:A:274:VAL:O	1:A:278:GLN:HG2	2.06	0.56
1:A:188:ARG:N	3:A:511:HOH:O	2.40	0.55
1:A:199:ARG:HG3	1:A:199:ARG:NH1	2.16	0.50
1:A:242:GLU:H	1:A:242:GLU:CD	2.20	0.49
1:A:280:LEU:C	1:A:280:LEU:HD23	2.38	0.49
1:A:189:MET:HB2	1:A:288:MET:HE1	1.96	0.48
1:A:110:ARG:HD2	1:A:110:ARG:HA	1.75	0.47
1:A:113[B]:ASN:ND2	1:A:146:ARG:HH21	2.07	0.46
1:A:217:ARG:HH22	1:A:327:GLN:NE2	2.04	0.46
1:A:189:MET:HG3	1:A:284[A]:CYS:SG	2.57	0.44
1:A:217:ARG:NH1	1:A:327:GLN:NE2	2.60	0.43
1:A:266:MET:HB3	1:A:324:LEU:HD23	2.00	0.43
1:A:313:HIS:HE1	3:A:579:HOH:O	2.02	0.43
1:A:189:MET:HB2	1:A:288:MET:CE	2.49	0.43
1:A:197:GLU:HA	1:A:278:GLN:OE1	2.19	0.42
1:A:194:ASP:HB3	1:A:195:PHE:HD2	1.84	0.42
1:A:266:MET:HB3	1:A:324:LEU:CD2	2.49	0.42
1:A:189:MET:HG3	1:A:288:MET:HE2	2.02	0.41
1:A:266:MET:HG2	1:A:268:ALA:H	1.85	0.41
1:A:297:GLU:OE1	1:A:301:HIS:CD2	2.69	0.41
1:A:199:ARG:O	1:A:203:GLU:HG3	2.21	0.41
1:A:248:ARG:CZ	1:A:258:LYS:HD2	2.51	0.41
1:A:285:HIS:HA	1:A:288:MET:HE3	2.03	0.40
1:A:189:MET:CB	1:A:288:MET:HE1	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:325:TRP:HB3	1:A:342:THR:HG22	2.04	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	269/292 (92%)	263 (98%)	6 (2%)	0	<b>100</b> <b>100</b>

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	A	242/259 (93%)	234 (97%)	8 (3%)	<b>33</b> <b>21</b>

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	VAL
1	A	125	LYS
1	A	161	ARG
1	A	190	GLU
1	A	191	MET

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Mol	Chain	Res	Type
1	A	199	ARG
1	A	201	LEU
1	A	278	GLN

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

Mol	Chain	Res	Type
1	A	73	GLN
1	A	155	GLN
1	A	301	HIS
1	A	313	HIS
1	A	327	GLN

### 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 oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 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 [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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	271/292 (92%)	-0.00	12 (4%) 39 38	13, 24, 41, 62	2 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	195	PHE	4.4
1	A	284[A]	CYS	3.6
1	A	176	GLY	3.3
1	A	345	PRO	3.2
1	A	293	ALA	2.5
1	A	194	ASP	2.4
1	A	192	ILE	2.4
1	A	202	SER	2.3
1	A	196	CYS	2.3
1	A	64	VAL	2.3
1	A	189	MET	2.3
1	A	197	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	HG	A	403	1/1	0.94	0.06	41,41,41,41	1
2	HG	A	401	1/1	0.98	0.06	27,27,27,27	1
2	HG	A	402	1/1	0.99	0.03	30,30,30,30	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.