

Full wwPDB X-ray Structure Validation Report (i)

May 21, 2020 - 03:03 am BST

PDB ID	:	2WMF
Title	:	Crystal structure of the catalytic module of a family 98 glycoside hydrolase
		from Streptococcus pneumoniae TIGR4 (Sp4GH98) in its native form.
Authors	:	Higgins, M.A.; Whitworth, G.E.; El Warry, N.; Randriantsoa, M.; Samain, E.;
		Burke, R.D.; Vocadlo, D.J.; Boraston, A.B.
Deposited on	:	2009-06-30
$\operatorname{Resolution}$:	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 following versions of software and data (see references (1)) were used in the production of this report:

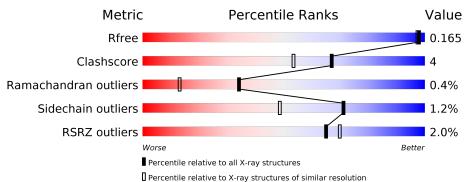
MolProbity Xtriage (Phenix) EDS		
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
$\operatorname{CCP4}$:	$7.0.044 (\mathrm{Gargrove})$
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 1.50 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
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 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% 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		
			2%		
1	А	581	87%	8%	•



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5103 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 FUCOLECTIN-RELATED PROTEIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Δ	555	Total	С	Ν	Ο	S	0	0	0
	А	000	4503	2873	756	853	21	0	U	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	544	SER	LYS	$\operatorname{conflict}$	UNP Q97N96

• Molecule 2 is water.

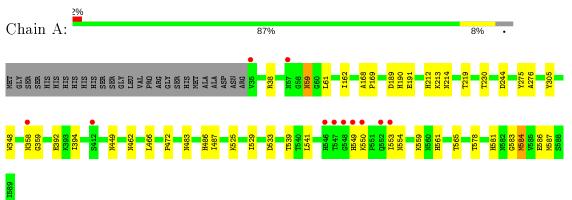
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	600	Total O 600 600	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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: FUCOLECTIN-RELATED PROTEIN





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	50.84Å 90.49 Å 116.78 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.35 - 1.50	Depositor
	38.34 - 1.50	EDS
% Data completeness	84.3(38.35-1.50)	Depositor
(in resolution range)	84.3(38.34-1.50)	EDS
R _{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$3.42 \; ({\rm at} \; 1.50 {\rm \AA})$	Xtriage
Refinement program	REFMAC $5.2.0019$	Depositor
R, R_{free}	0.124 , 0.160	Depositor
III, IIIfree	0.133 , 0.165	DCC
R_{free} test set	3675 reflections $(5.01%)$	wwPDB-VP
Wilson B-factor (Å ²)	12.3	Xtriage
Anisotropy	0.249	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.37, 49.7	EDS
L-test for twinning ²	$ \langle L \rangle = 0.48, \langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5103	wwPDB-VP
Average B, all atoms $(Å^2)$	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.40% 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)

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		
	Ullalli	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.77	0/4636	0.78	5/6317~(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	А	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	А	584	MET	CG-SD-CE	7.62	112.38	100.20
1	А	359	GLY	N-CA-C	7.00	130.61	113.10
1	А	38	ARG	NE-CZ-NH1	5.66	123.13	120.30
1	А	587	MET	CG-SD-CE	-5.61	91.22	100.20
1	А	213	LYS	CD-CE-NZ	5.15	123.55	111.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	358	ASN	Peptide

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



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	А	4503	0	4304	34	0
2	А	600	0	0	18	1
All	All	5103	0	4304	34	1

the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

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 (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:A:584:MET:HG2	2:A:2597:HOH:O	$\frac{1.29}{1.29}$	1.26
1:A:584:MET:CE	2:A:2337:HOH:O 2:A:2327:HOH:O	1.25	1.15
1:A:584:MET:HE1	2:A:2327:HOH:O 2:A:2327:HOH:O	1.54	0.96
1:A:584:MET:SD	2:A:2571:HOH:O	2.32	0.86
1:A:529:ILE:HD11	2:A:2535:HOH:O	1.74	0.86
1:A:529:ILE:CD1	2:A:2535:HOH:O 2:A:2535:HOH:O	2.32	0.00
1:A:584:MET:HE3	2:A:2327:HOH:O	1.71	0.76
1:A:584:MET:CG	2:A:2597:HOH:O	2.06	0.74
1:A:392:GLU:OE2	2:A:2402:HOH:O	2.09	0.69
1:A:486:HIS:CD2	2:A:2500:HOH:O	2.48	0.67
1:A:212:HIS:HD2	1:A:244:ASP:OD1	1.80	0.63
1:A:584:MET:HE3	2:A:2491:HOH:O	2.00	0.62
1:A:533:ASP:OD1	1:A:561:HIS:HE1	1.85	0.60
1:A:586:GLU:HG3	2:A:2597:HOH:O	2.02	0.59
1:A:539:THR:OG1	1:A:581:HIS:HE1	1.86	0.58
1:A:212:HIS:HE1	2:A:2114:HOH:O	1.86	0.57
1:A:559:LYS:O	1:A:561:HIS:HD2	1.90	0.54
1:A:486:HIS:HD2	2:A:2492:HOH:O	1.97	0.48
1:A:219:THR:HG22	1:A:230:THR:HG22	1.96	0.47
1:A:462:ASN:HD21	1:A:483:ASN:HA	1.78	0.47
1:A:190:HIS:CD2	1:A:230:THR:HG21	2.49	0.47
1:A:59:ASN:ND2	1:A:61:LEU:H	2.13	0.46
1:A:487:ILE:HD13	1:A:541:LEU:HD21	1.98	0.46
1:A:168:ALA:HB3	1:A:169:PRO:HD3	1.99	0.45
1:A:565:THR:HG23	2:A:2595:HOH:O	2.16	0.44
1:A:466:LEU:HD21	1:A:541:LEU:CD1	2.47	0.44
1:A:348:TRP:CD1	1:A:394:ILE:HD11	2.52	0.43
1:A:553:ILE:HG22	2:A:2583:HOH:O	2.18	0.42
1:A:578:THR:HG23	2:A:2561:HOH:O	2.20	0.42
1:A:525:LYS:HA	1:A:529:ILE:HD11	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:581:HIS:HD2	1:A:583:GLY:O	2.03	0.41
1:A:449:ASN:HA	1:A:472:PRO:HG2	2.03	0.41
1:A:214:ASN:ND2	2:A:2236:HOH:O	2.55	0.40
1:A:276:ALA:HA	1:A:305:TYR:CG	2.56	0.40

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All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1			Clash overlap (Å)
2:A:2181:HOH:O	2:A:2499:HOH:O[4_556]	0.99	1.21

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	553/581~(95%)	526~(95%)	25~(4%)	2(0%)	34 13

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	191	GLU
1	А	162	ILE

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	493/515~(96%)	487~(99%)	6 (1%)	71 48	

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

Mol	Chain	Res	Type
1	А	59	ASN
1	А	189	ASP
1	А	275	TYR
1	А	549	HIS
1	А	550	LYS
1	А	554	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	59	ASN
1	А	190	HIS
1	А	200	ASN
1	А	212	HIS
1	А	214	ASN
1	А	375	ASN
1	А	433	GLN
1	А	462	ASN
1	А	481	ASN
1	А	486	HIS
1	А	549	HIS
1	А	561	HIS
1	А	571	ASN
1	А	580	ASN
1	А	581	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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, 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 $>$	$\cdot 2$	$\mathbf{OWAB}(\mathbf{A}^2)$	Q<0.9
1	А	555/581~(95%)	-0.26	11 (1%) 65	70	7, 11, 21, 48	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	549	HIS	6.5
1	А	547	THR	5.7
1	А	358	ASN	3.7
1	А	548	GLY	3.2
1	А	553	ILE	3.1
1	А	546	HIS	3.0
1	А	57	ASN	3.0
1	А	35	VAL	3.0
1	А	550	LYS	2.4
1	А	552	GLN	2.3
1	А	412	SER	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 carbohydrates in this entry.

6.4 Ligands (i)

There are no ligands in this entry.



6.5 Other polymers (i)

There are no such residues in this entry.

