

wwPDB X-ray Structure Validation Summary Report (i)

May 26, 2020 – 05:15 pm BST

PDB ID : 4ZE4

Title: Structure of Gan1D, a putative 6-phospho-beta-galactosidase from Geobacillus

stearothermophilus

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Deposited on : 2015-04-20

Resolution : 1.92 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.11

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac: 5.8.0158

 $\begin{array}{cccc} & CCP4 & : & 7.0.044 \; (Gargrove) \\ Ideal \; geometry \; (proteins) & : & Engh \; \& \; Huber \; (2001) \end{array}$

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

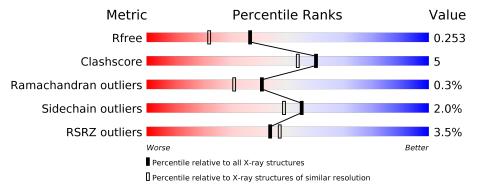
Validation Pipeline (wwPDB-VP) : 2.11

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 1.92 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
R_{free}	130704	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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		
1	A	485	85%	12%	
1	В	485	6% 85%	12%	



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8358 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 Putative 6-phospho-beta-galactobiosidase.

\mathbf{Mol}	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	474	Total	С	N	О	S	0	Q	0
1	Λ	474	3920	2518	674	717	11	0	0	0
1	B	474	Total	С	N	О	S	0	1	0
1	Ъ	474	3880	2490	664	715	11		1	

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MET	=	initiating methionine	UNP W8QF82
A	-5	ILE	-	expression tag	UNP W8QF82
A	-4	HIS	_	expression tag	UNP W8QF82
A	-3	HIS	ı	expression tag	UNP W8QF82
A	-2	HIS	-	expression tag	UNP W8QF82
A	-1	HIS	ı	expression tag	UNP W8QF82
A	0	HIS	-	expression tag	UNP W8QF82
A	1	HIS	-	expression tag	UNP W8QF82
В	-6	MET	_	initiating methionine	UNP W8QF82
В	-5	ILE	-	expression tag	UNP W8QF82
В	-4	HIS	ı	expression tag	UNP W8QF82
В	-3	HIS	-	expression tag	UNP W8QF82
В	-2	HIS	_	expression tag	UNP W8QF82
В	-1	HIS	=	expression tag	UNP W8QF82
В	0	HIS		expression tag	UNP W8QF82
В	1	HIS	-	expression tag	UNP W8QF82

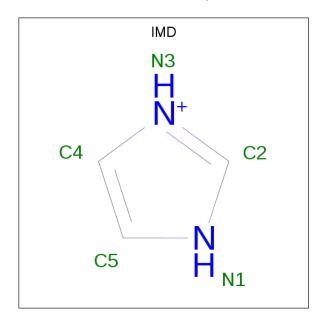
• Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).





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

 \bullet Molecule 3 is IMIDAZOLE (three-letter code: IMD) (formula: $\mathrm{C_3H_5N_2}).$



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N 5 3 2	0	0
3	A	1	Total C N 5 3 2	0	0



• Molecule 4 is water.

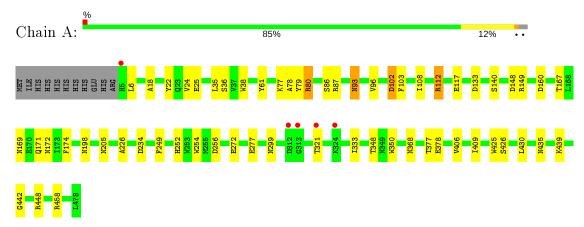
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	365	Total O 365 365	0	0
4	В	171	Total O 171 171	0	0



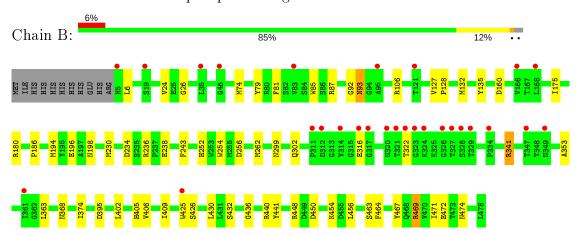
3 Residue-property plots (i)

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: Putative 6-phospho-beta-galactobiosidase



• Molecule 1: Putative 6-phospho-beta-galactobiosidase





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	108.01Å 68.88Å 152.88Å	Depositor
a, b, c, α , β , γ	90.00° 100.74° 90.00°	Depositor
Resolution (Å)	31.90 - 1.92	Depositor
Resolution (A)	31.88 - 1.92	EDS
% Data completeness	99.6 (31.90-1.92)	Depositor
(in resolution range)	99.7 (31.88-1.92)	EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	2.17 (at 1.92Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
P. P.	0.194 , 0.245	Depositor
R, R_{free}	0.203 , 0.253	DCC
R_{free} test set	4183 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.631	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.32 , 44.3	EDS
L-test for twinning ²	$ < L >=0.49, < L^2>=0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8358	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.21% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $< L^2>$ 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: GOL, IMD

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	Bo	nd lengths	Bond angles		
WIOI	Chain	RMSZ # Z > 5		RMSZ	# Z >5	
1	A	1.19	$17/4065 \ (0.4\%)$	1.06	$17/5528 \ (0.3\%)$	
1	В	0.76	0/4004	0.82	$4/5449 \ (0.1\%)$	
All	All	1.00	$17/8069 \ (0.2\%)$	0.95	21/10977~(0.2%)	

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(ext{\AA})$
1	A	87	ARG	CZ-NH2	13.12	1.50	1.33
1	A	36	SER	CB-OG	8.92	1.53	1.42
1	A	25	GLU	CG-CD	8.59	1.64	1.51
1	A	79	TYR	C-N	-6.58	1.19	1.34
1	A	38	TRP	C-O	6.14	1.35	1.23

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}(^{o})$
1	A	80	ARG	O-C-N	9.38	137.71	122.70
1	A	80	ARG	NE-CZ-NH2	-8.46	116.07	120.30
1	A	87	ARG	NE-CZ-NH1	8.39	124.49	120.30
1	A	87	ARG	NE-CZ-NH2	-8.08	116.26	120.30
1	A	80	ARG	CA-C-N	-7.93	99.75	117.20

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	$\mathbf{H}(\mathbf{model})$	H(added)	Clashes	Symm-Clashes
1	A	3920	0	3744	41	0
1	В	3880	0	3674	34	0
2	A	6	0	8	0	0
2	В	6	0	8	0	0
3	A	10	0	10	0	0
4	A	365	0	0	10	0
4	В	171	0	0	6	0
All	All	8358	0	7444	70	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 5.

The worst 5 of 70 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{c} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{array}$	Clash overlap (Å)
1:A:112[A]:ARG:HG3	1:A:112[A]:ARG:HH11	1.09	1.16
1:B:316:GLU:HG2	4:B:602:HOH:O	1.53	1.08
1:A:112[A]:ARG:CG	1:A:112[A]:ARG:HH11	1.66	1.05
1:A:112[A]:ARG:HG3	1:A:112[A]:ARG:NH1	1.80	0.90
1:A:102:ASP:HB2	4:A:601:HOH:O	1.72	0.90

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	Perce	ntiles
1	A	480/485 (99%)	462 (96%)	18 (4%)	0	100	100
1	В	473/485 (98%)	446 (94%)	24 (5%)	3 (1%)	25	14
All	All	953/970 (98%)	908 (95%)	42 (4%)	3 (0%)	41	31



All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	454	LYS
1	В	474	ASN
1	В	135	TYR

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	409/412 (99%)	403 (98%)	6 (2%)	65 61		
1	В	402/412 (98%)	391 (97%)	11 (3%)	44 36		
All	All	811/824 (98%)	794 (98%)	17 (2%)	55 46		

5 of 17 residues with a non-rotameric sidechain are listed below:

Mol	Chain	${f Res}$	\mathbf{Type}
1	В	24	VAL
1	В	93	ASN
1	В	322	THR
1	В	6	LEU
1	В	341	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	274	GLN
1	A	379	ASN
1	В	371	GLN
1	A	349	ASN
1	A	435	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 carbohydrates in this entry.

5.6 Ligand geometry (i)

4 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	Mol Type Chain Res		Res	Link	Bond lengths			Bond angles		
MIGI	Type	Chain	nes	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	IMD	A	503	-	3,5,5	0.58	0	$4,\!5,\!5$	0.28	0
2	GOL	В	501	-	5,5,5	0.39	0	5,5,5	0.44	0
3	IMD	A	502	_	3,5,5	0.30	0	$4,\!5,\!5$	0.48	0
2	GOL	A	501	_	5,5,5	0.35	0	5,5,5	0.29	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
3	IMD	A	503	-	-	-	0/1/1/1
2	GOL	В	501	-	-	0/4/4/4	-
3	IMD	A	502	-	-	-	0/1/1/1
2	GOL	A	501	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:



Mol	Chain	Res	Type	Atoms
2	A	501	GOL	C1-C2-C3-O3
2	A	501	GOL	O2-C2-C3-O3
2	A	501	GOL	O1-C1-C2-C3

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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	79:TYR	С	80:ARG	N	1.18



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^{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	<RSRZ $>$	$\#\mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q < 0.9
1	A	474/485 (97%)	-0.09	5 (1%) 80 82	11, 24, 48, 84	0
1	В	474/485 (97%)	0.40	28 (5%) 22 25	27, 39, 60, 91	0
All	All	948/970 (97%)	0.16	33 (3%) 44 47	11, 33, 56, 91	0

The worst 5 of 33 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	326	GLY	4.9
1	В	322	THR	4.4
1	В	95	ALA	4.4
1	A	5	HIS	4.3
1	В	321	THR	4.2

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)

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	$\operatorname{B-factors}({ m \AA}^2)$	Q<0.9
2	GOL	В	501	6/6	0.59	0.24	65,71,74,75	0
3	IMD	A	502	5/5	0.89	0.23	32,37,39,42	0
3	IMD	A	503	5/5	0.91	0.17	29,29,34,35	0
2	GOL	A	501	6/6	0.92	0.17	51,54,60,61	0

6.5 Other polymers (i)

There are no such residues in this entry.

