Full wwPDB X-ray Structure Validation Report ⓘ

May 10, 2016 – 01:48 PM EDT

PDB ID	:	5FA0
Title	:	The structure of the beta-3-deoxy-D-manno-oct-2-ulosonic acid transferase domain from WbbB
Authors	:	Mallette, E.; Ovchinnikova, O.G.; Whitfield, C.; Kimber, M.S.
Deposited on	:	2015-12-10
Resolution	:	2.30 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see references ①) were used in the production of this report:

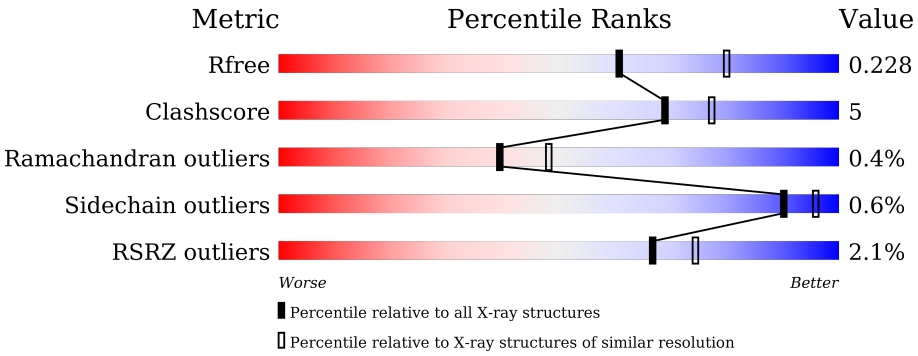
MolProbity	:	4.02b-467
Mogul	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027457
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20027457

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

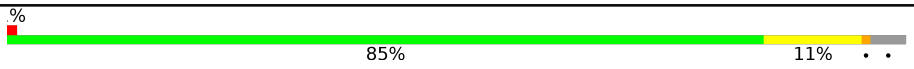
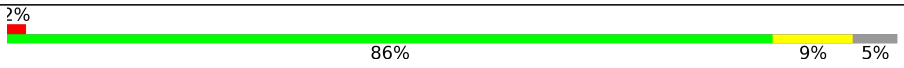
The reported resolution of this entry is 2.30 Å.

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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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. 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	410	
1	B	410	

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 6801 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 N-acetyl glucosaminyl transferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	394	Total	C	N	O	S	Se	0	1	0
			3140	2012	519	596	7	6			
1	B	388	Total	C	N	O	S	Se	0	5	0
			3098	1991	505	590	7	5			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MSE	-	initiating methionine	UNP Q6U8B0
A	1	GLY	-	expression tag	UNP Q6U8B0
A	402	LEU	-	expression tag	UNP Q6U8B0
A	403	GLU	-	expression tag	UNP Q6U8B0
A	404	HIS	-	expression tag	UNP Q6U8B0
A	405	HIS	-	expression tag	UNP Q6U8B0
A	406	HIS	-	expression tag	UNP Q6U8B0
A	407	HIS	-	expression tag	UNP Q6U8B0
A	408	HIS	-	expression tag	UNP Q6U8B0
A	409	HIS	-	expression tag	UNP Q6U8B0
B	0	MSE	-	initiating methionine	UNP Q6U8B0
B	1	GLY	-	expression tag	UNP Q6U8B0
B	402	LEU	-	expression tag	UNP Q6U8B0
B	403	GLU	-	expression tag	UNP Q6U8B0
B	404	HIS	-	expression tag	UNP Q6U8B0
B	405	HIS	-	expression tag	UNP Q6U8B0
B	406	HIS	-	expression tag	UNP Q6U8B0
B	407	HIS	-	expression tag	UNP Q6U8B0
B	408	HIS	-	expression tag	UNP Q6U8B0
B	409	HIS	-	expression tag	UNP Q6U8B0

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0

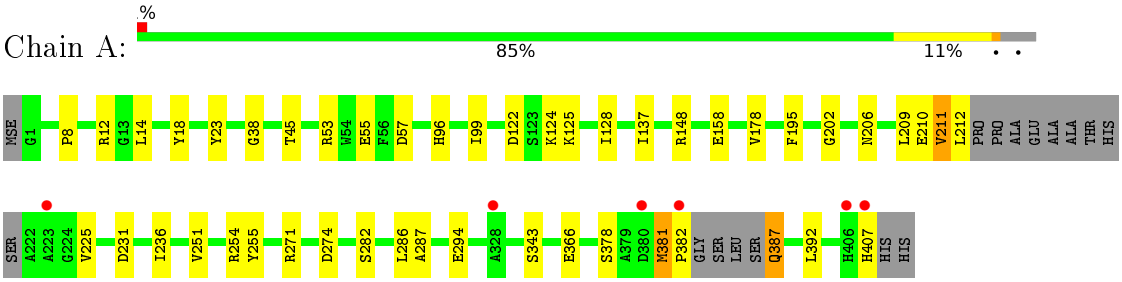
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	306	Total O 306 306	0	0
3	B	256	Total O 256 256	0	0

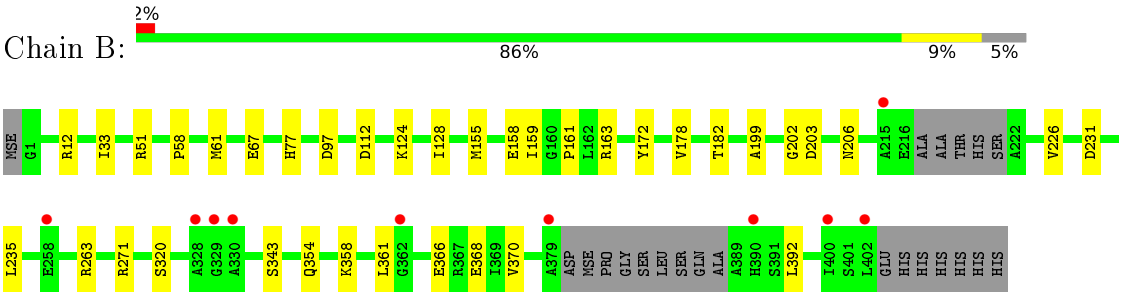
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 N-acetyl glucosaminyl transferase



- Molecule 1: Putative N-acetyl glucosaminyl transferase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	82.93Å 82.93Å 120.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.47 – 2.30 48.80 – 1.88	Depositor EDS
% Data completeness (in resolution range)	100.0 (41.47-2.30) 96.8 (48.80-1.88)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	1.32 (at 1.88Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.185 , 0.229 0.182 , 0.228	Depositor DCC
R_{free} test set	1141 reflections (3.03%)	DCC
Wilson B-factor (Å^2)	18.8	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 44.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	0.034 for k,h,-l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6801	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.49 % of the origin peak, indicating pseudo translational symmetry.The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 8.5948e-03.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CL

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	0.23	0/3215	0.39	0/4355
1	B	0.23	0/3183	0.38	0/4315
All	All	0.23	0/6398	0.38	0/8670

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

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	3140	0	3025	30	0
1	B	3098	0	3010	29	0
2	A	1	0	0	1	0
3	A	306	0	0	9	1
3	B	256	0	0	10	1
All	All	6801	0	6035	56	1

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.

All (56) 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:57:ASP:OD1	3:A:601:HOH:O	1.93	0.85
1:B:97:ASP:OD2	3:B:501:HOH:O	1.98	0.81
2:A:501:CL:CL	3:A:867:HOH:O	2.38	0.79
1:A:255:TYR:OH	1:A:294:GLU:OE2	2.01	0.76
1:B:368:GLU:OE1	3:B:502:HOH:O	2.10	0.68
1:B:67:GLU:OE2	3:B:503:HOH:O	2.13	0.66
1:A:125:LYS:NZ	3:A:607:HOH:O	2.26	0.65
1:B:182:THR:O	3:B:505:HOH:O	2.14	0.64
1:B:51:ARG:NH1	3:B:506:HOH:O	2.19	0.64
1:A:99:ILE:HG23	1:A:236:ILE:HD12	1.82	0.61
1:B:163:ARG:NH2	3:B:515:HOH:O	2.30	0.61
1:B:370:VAL:HG21	1:B:392:LEU:HD21	1.82	0.61
1:B:58:PRO:HA	1:B:61:MSE:HE2	1.83	0.58
1:A:96:HIS:ND1	3:A:603:HOH:O	2.17	0.57
1:B:202:GLY:O	1:B:206:ASN:ND2	2.37	0.56
1:A:148:ARG:NH2	3:A:618:HOH:O	2.38	0.56
1:A:382:PRO:HB3	1:A:387:GLN:HG2	1.88	0.56
1:A:210:GLU:OE1	3:A:605:HOH:O	2.18	0.55
1:A:202:GLY:O	1:A:206:ASN:ND2	2.32	0.53
1:A:271:ARG:HH11	1:B:271:ARG:NE	2.07	0.52
1:A:12:ARG:NH2	1:A:231:ASP:OD1	2.42	0.52
1:A:378:SER:OG	3:A:602:HOH:O	2.09	0.52
1:A:122:ASP:OD1	1:A:125:LYS:NZ	2.41	0.51
1:B:343:SER:OG	1:B:366:GLU:OE1	2.29	0.49
1:A:381:MSE:SE	1:A:382:PRO:HD2	2.63	0.49
1:A:407:HIS:N	3:A:622:HOH:O	2.45	0.48
1:B:263:ARG:NE	3:B:504:HOH:O	2.14	0.48
1:B:158:GLU:HB3	1:B:178:VAL:HG21	1.95	0.48
1:B:112:ASP:O	3:B:508:HOH:O	2.20	0.47
1:B:155:MSE:HE3	1:B:361:LEU:HD13	1.95	0.47
1:B:182:THR:HG21	1:B:320:SER:HB3	1.96	0.46
1:B:159:ILE:CG2	1:B:163:ARG:HD2	2.46	0.46
1:B:199:ALA:HB1	1:B:203:ASP:HB2	1.97	0.45
1:A:271:ARG:NH1	1:B:271:ARG:HH21	2.15	0.45
1:A:343:SER:OG	1:A:366:GLU:OE1	2.35	0.45
1:B:354[B]:GLN:HG2	1:B:358:LYS:HE2	1.98	0.45
1:A:124:LYS:O	1:A:128:ILE:HG12	2.16	0.44
1:A:23:TYR:HH	1:A:45:THR:HG1	1.59	0.44
1:A:18:TYR:CD1	1:A:137:ILE:HD13	2.53	0.44
1:B:12:ARG:NH2	1:B:231:ASP:OD1	2.51	0.44
1:A:282:SER:HB3	1:A:287:ALA:HB3	2.00	0.44
1:A:225:VAL:HG22	1:A:251:VAL:HG21	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:ARG:HH11	1:B:271:ARG:CZ	2.32	0.43
1:B:124:LYS:O	1:B:128:ILE:HG12	2.18	0.43
1:A:8:PRO:HB3	1:A:38:GLY:HA3	2.01	0.43
1:B:161:PRO:HD3	1:B:172:TYR:CZ	2.54	0.43
1:B:235:LEU:O	3:B:509:HOH:O	2.22	0.42
1:A:158:GLU:HB3	1:A:178:VAL:HG21	2.02	0.42
1:A:271:ARG:NH1	1:B:271:ARG:HE	2.17	0.42
1:A:209:LEU:HD21	1:A:286:LEU:HD11	2.01	0.42
1:B:33:ILE:O	1:B:77:HIS:HB3	2.20	0.41
1:A:254:ARG:NH1	3:A:631:HOH:O	2.53	0.41
1:B:226:VAL:O	3:B:510:HOH:O	2.22	0.41
1:A:195:PHE:CE1	1:A:392:LEU:HD21	2.56	0.41
1:A:53:ARG:HB3	1:A:55:GLU:OE1	2.21	0.41
1:B:159:ILE:HG23	1:B:163:ARG:HD2	2.02	0.40

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	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:855:HOH:O	3:B:749:HOH:O[1_655]	1.89	0.31

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	389/410 (95%)	375 (96%)	11 (3%)	3 (1%)	24 27
1	B	387/410 (94%)	376 (97%)	11 (3%)	0	100 100
All	All	776/820 (95%)	751 (97%)	22 (3%)	3 (0%)	39 48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	381	MSE
1	A	274	ASP
1	A	211	VAL

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	338/342 (99%)	334 (99%)	4 (1%)	78 89
1	B	336/342 (98%)	336 (100%)	0	100 100
All	All	674/684 (98%)	670 (99%)	4 (1%)	90 96

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LEU
1	A	211	VAL
1	A	212	LEU
1	A	387	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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

Of 1 ligands modelled in this entry, 1 is 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, 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>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	388/410 (94%)	-0.36	6 (1%)	76	81	9, 28, 58, 99	0
1	B	383/410 (93%)	-0.16	10 (2%)	59	68	10, 31, 69, 84	0
All	All	771/820 (94%)	-0.26	16 (2%)	67	74	9, 30, 65, 99	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	379	ALA	4.7
1	A	407	HIS	4.1
1	A	328	ALA	3.8
1	B	328	ALA	3.8
1	B	402	LEU	3.5
1	B	258	GLU	3.4
1	A	382	PRO	3.4
1	B	215	ALA	3.1
1	A	406	HIS	2.9
1	A	380	ASP	2.8
1	B	330	ALA	2.4
1	B	390	HIS	2.3
1	B	329	GLY	2.1
1	A	223	ALA	2.1
1	B	400	ILE	2.1
1	B	362	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å^2)	Q<0.9
2	CL	A	501	1/1	0.97	0.10	-	42,42,42,42	0

6.5 Other polymers (i)

There are no such residues in this entry.