



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 10, 2016 – 10:36 PM EDT

PDB ID : 4ZY9
Title : X-ray crystal structure of selenomethionine-labelled V110M mutant of chitosan-binding module 1 derived from chitosanase/glucanase from *Paenibacillus* sp. IK-5
Authors : Shinya, S.; Oi, H.; Kitaoku, Y.; Ohnuma, T.; Numata, T.; Fukamizo, T.
Deposited on : 2015-05-21
Resolution : 1.20 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027107
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-20027107

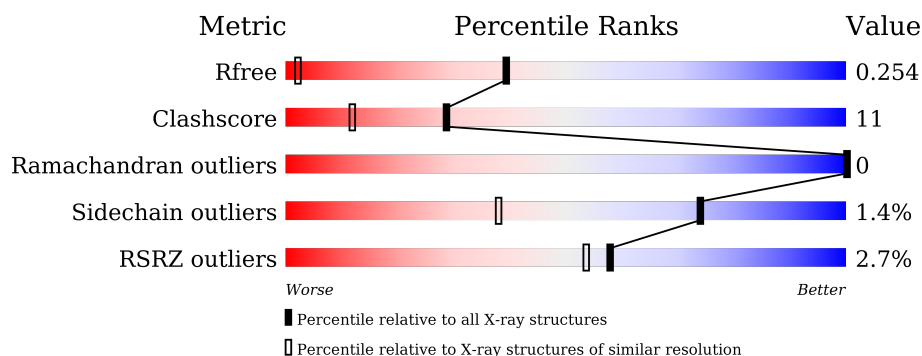
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.20 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	1495 (1.26-1.14)
Clashscore	102246	1607 (1.26-1.14)
Ramachandran outliers	100387	1540 (1.26-1.14)
Sidechain outliers	100360	1538 (1.26-1.14)
RSRZ outliers	91569	1500 (1.26-1.14)

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 $\leq 5\%$. 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	137	
1	B	137	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2748 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 Glucanase/chitosanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	133	Total	C	N	O	Se	0	13	0
			1092	690	185	216	1			
1	B	133	Total	C	N	O	Se	0	12	0
			1092	683	186	222	1			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	MSE	-	expression tag	UNP Q93IE7
A	-5	HIS	-	expression tag	UNP Q93IE7
A	-4	HIS	-	expression tag	UNP Q93IE7
A	-3	HIS	-	expression tag	UNP Q93IE7
A	-2	HIS	-	expression tag	UNP Q93IE7
A	-1	HIS	-	expression tag	UNP Q93IE7
A	0	HIS	-	expression tag	UNP Q93IE7
A	110	MSE	VAL	engineered mutation	UNP Q93IE7
B	-6	MSE	-	expression tag	UNP Q93IE7
B	-5	HIS	-	expression tag	UNP Q93IE7
B	-4	HIS	-	expression tag	UNP Q93IE7
B	-3	HIS	-	expression tag	UNP Q93IE7
B	-2	HIS	-	expression tag	UNP Q93IE7
B	-1	HIS	-	expression tag	UNP Q93IE7
B	0	HIS	-	expression tag	UNP Q93IE7
B	110	MSE	VAL	engineered mutation	UNP Q93IE7

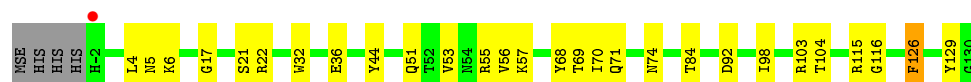
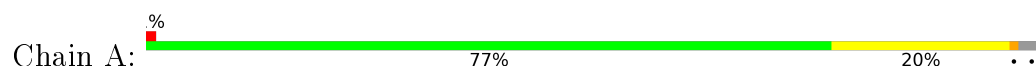
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	292	Total	O	0	0
			292	292		
2	B	272	Total	O	0	0
			272	272		

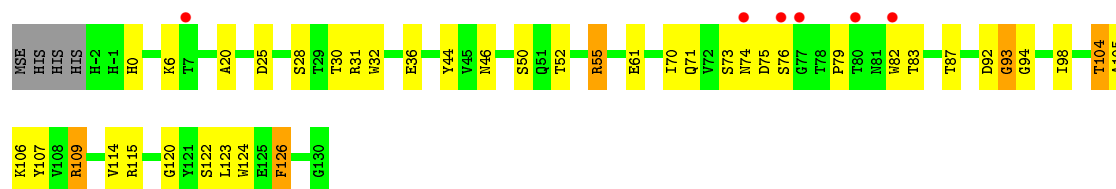
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: Glucanase/chitosanase



- Molecule 1: Glucanase/chitosanase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	28.23Å 65.33Å 69.89Å 90.00° 97.21° 90.00°	Depositor
Resolution (Å)	47.55 – 1.20 47.55 – 1.20	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.55-1.20) 99.2 (47.55-1.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.06 (at 1.20Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.208 , 0.249 0.217 , 0.254	Depositor DCC
R_{free} test set	1994 reflections (2.63%)	DCC
Wilson B-factor (Å ²)	8.3	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 32.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 77927 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2748	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.41% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	1.53	10/1154 (0.9%)	1.28	8/1573 (0.5%)
1	B	1.66	17/1139 (1.5%)	1.40	8/1552 (0.5%)
All	All	1.60	27/2293 (1.2%)	1.34	16/3125 (0.5%)

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	B	0	1

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	93	GLY	N-CA	-8.42	1.33	1.46
1	A	116	GLY	C-O	7.37	1.35	1.23
1	A	129	TYR	CE2-CZ	-6.67	1.29	1.38
1	B	124	TRP	CE3-CZ3	6.43	1.49	1.38
1	B	83	THR	N-CA	-6.26	1.33	1.46
1	A	36[A]	GLU	CD-OE2	5.71	1.31	1.25
1	A	36[B]	GLU	CD-OE2	5.71	1.31	1.25
1	A	21	SER	CA-CB	5.69	1.61	1.52
1	B	122[A]	SER	C-O	-5.61	1.12	1.23
1	B	122[B]	SER	C-O	-5.61	1.12	1.23
1	B	20	ALA	C-O	-5.59	1.12	1.23
1	A	32	TRP	CG-CD1	5.55	1.44	1.36
1	B	55	ARG	CZ-NH2	5.55	1.40	1.33
1	B	44	TYR	CE1-CZ	-5.46	1.31	1.38
1	A	44	TYR	CE1-CZ	-5.42	1.31	1.38
1	B	82	TRP	CG-CD1	5.42	1.44	1.36
1	B	123	LEU	C-O	-5.41	1.13	1.23
1	B	28	SER	CA-CB	-5.37	1.44	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	32	TRP	C-O	-5.36	1.13	1.23
1	B	87	THR	CB-OG1	-5.28	1.32	1.43
1	B	44	TYR	CG-CD2	-5.21	1.32	1.39
1	A	17	GLY	N-CA	-5.20	1.38	1.46
1	B	105	ALA	CA-C	-5.15	1.39	1.52
1	A	68	TYR	CE1-CZ	-5.13	1.31	1.38
1	B	28	SER	C-O	5.08	1.33	1.23
1	A	129	TYR	CE1-CZ	-5.06	1.31	1.38
1	B	83	THR	CB-CG2	-5.03	1.35	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	TYR	CB-CG-CD1	8.38	126.03	121.00
1	B	109	ARG	NE-CZ-NH2	-8.20	116.20	120.30
1	A	115	ARG	NE-CZ-NH2	-8.05	116.28	120.30
1	B	115	ARG	NE-CZ-NH2	-7.56	116.52	120.30
1	A	92	ASP	CB-CG-OD1	7.39	124.95	118.30
1	B	28	SER	CB-CA-C	-6.89	97.00	110.10
1	B	126	PHE	CB-CG-CD1	6.56	125.39	120.80
1	A	22	ARG	NE-CZ-NH2	-6.21	117.19	120.30
1	B	36	GLU	CG-CD-OE1	-6.17	105.95	118.30
1	B	31	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	A	126	PHE	CB-CG-CD1	5.77	124.84	120.80
1	A	55	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	B	31	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	A	57	LYS	CD-CE-NZ	5.28	123.83	111.70
1	A	68	TYR	CB-CG-CD1	-5.19	117.89	121.00
1	B	52	THR	CA-CB-CG2	-5.01	105.38	112.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	73	SER	Mainchain

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	1092	0	1057	21	0
1	B	1092	0	1028	28	0
2	A	292	0	0	7	5
2	B	272	0	0	14	4
All	All	2748	0	2085	49	6

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 11.

All (49) 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:92:ASP:OD2	2:B:202:HOH:O	1.69	1.10
1:B:93:GLY:O	2:B:201:HOH:O	1.67	1.09
1:A:70[B]:ILE:CD1	1:A:98:ILE:HD13	1.88	1.03
1:B:104[A]:THR:HG22	2:B:245:HOH:O	1.61	1.01
1:B:71[B]:GLN:NE2	2:B:203:HOH:O	1.98	0.96
1:A:70[B]:ILE:HD11	1:A:98:ILE:HD13	1.46	0.95
1:B:75[B]:ASP:O	1:B:106[B]:LYS:HD3	1.69	0.91
1:B:0:HIS:HA	2:B:237:HOH:O	1.72	0.89
1:A:69[B]:THR:HG23	1:A:84:THR:HG23	1.57	0.86
1:A:104[B]:THR:HG22	2:A:282:HOH:O	1.75	0.86
1:A:74:ASN:OD1	2:A:201:HOH:O	1.94	0.85
1:A:74:ASN:ND2	2:A:201:HOH:O	2.10	0.85
1:B:70[B]:ILE:HD11	1:B:98:ILE:HD13	1.59	0.85
1:B:75[A]:ASP:HA	2:B:230:HOH:O	1.78	0.82
1:B:70[B]:ILE:HD11	1:B:98:ILE:CD1	2.11	0.81
1:A:70[B]:ILE:HD11	1:A:98:ILE:CD1	2.11	0.80
1:A:56[B]:VAL:CG2	1:A:70[B]:ILE:HD13	2.11	0.80
1:A:56[B]:VAL:HG21	1:A:70[B]:ILE:HD13	1.69	0.75
1:B:70[B]:ILE:CD1	1:B:98:ILE:HD13	2.17	0.75
1:B:61:GLU:O	2:B:204:HOH:O	2.05	0.74
1:B:75[B]:ASP:OD2	1:B:79:PRO:HA	1.88	0.74
1:B:74[B]:ASN:O	2:B:205:HOH:O	2.09	0.71
1:A:53[B]:VAL:HG13	1:A:103[B]:ARG:HB2	1.73	0.71
1:A:70[B]:ILE:HD12	1:A:98:ILE:HD13	1.72	0.70
1:B:74[B]:ASN:OD1	2:B:205:HOH:O	2.07	0.70
1:B:120:GLY:O	2:B:206:HOH:O	2.11	0.68
1:B:114:VAL:HA	2:B:207:HOH:O	1.93	0.68
1:A:6[A]:LYS:NZ	2:A:205:HOH:O	2.31	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76[B]:SER:O	1:B:106[B]:LYS:NZ	2.28	0.62
1:A:70[B]:ILE:CD1	1:A:98:ILE:CD1	2.72	0.62
1:A:51[C]:GLN:NE2	2:A:206:HOH:O	2.32	0.59
1:B:0:HIS:CA	2:B:237:HOH:O	2.37	0.59
1:A:69[B]:THR:HG22	1:A:71:GLN:HG3	1.88	0.56
1:B:46[A]:ASN:HB2	1:B:107:TYR:CE1	2.45	0.52
1:B:109:ARG:NH1	2:B:213:HOH:O	2.44	0.51
1:A:69[B]:THR:HG22	1:A:71:GLN:CG	2.41	0.51
1:B:70[B]:ILE:CD1	1:B:98:ILE:CD1	2.86	0.48
1:A:69[B]:THR:CG2	1:A:84:THR:HG23	2.35	0.47
1:B:25[B]:ASP:OD2	1:B:30:THR:OG1	2.24	0.47
1:A:56[B]:VAL:HG21	1:A:70[B]:ILE:HG21	1.97	0.47
1:B:50:SER:HB3	1:B:104[B]:THR:CG2	2.45	0.46
1:A:51[A]:GLN:CG	2:A:202:HOH:O	2.64	0.44
1:B:55:ARG:HA	1:B:98:ILE:O	2.17	0.44
1:B:94:GLY:HA2	2:B:347:HOH:O	2.18	0.43
1:B:70[B]:ILE:HD11	1:B:98:ILE:HD12	1.95	0.42
1:A:51[C]:GLN:HG3	2:A:202:HOH:O	2.19	0.42
1:A:4:LEU:O	1:A:5[A]:ASN:HB2	2.22	0.40

All (6) 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)
2:A:258:HOH:O	2:A:421:HOH:O[2_555]	1.48	0.72
2:B:231:HOH:O	2:B:297:HOH:O[1_655]	1.93	0.27
2:A:368:HOH:O	2:A:425:HOH:O[2_545]	1.96	0.24
2:A:485:HOH:O	2:B:460:HOH:O[2_545]	2.08	0.12
2:A:477:HOH:O	2:B:465:HOH:O[2_545]	2.11	0.09
2:A:489:HOH:O	2:B:446:HOH:O[2_646]	2.12	0.08

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	143/137 (104%)	141 (99%)	2 (1%)	0	100	100
1	B	142/137 (104%)	138 (97%)	4 (3%)	0	100	100
All	All	285/274 (104%)	279 (98%)	6 (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	119/109 (109%)	118 (99%)	1 (1%)	86	62
1	B	117/109 (107%)	114 (97%)	3 (3%)	54	12
All	All	236/218 (108%)	232 (98%)	4 (2%)	74	29

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

Mol	Chain	Res	Type
1	A	126	PHE
1	B	104[A]	THR
1	B	104[B]	THR
1	B	126	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	81	ASN
1	B	41	GLN
1	B	51	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 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>2	OWAB(Å ²)	Q<0.9
1	A	132/137 (96%)	-0.01	1 (0%) 87 87	4, 7, 11, 21	0
1	B	132/137 (96%)	0.60	6 (4%) 37 33	5, 12, 22, 26	0
All	All	264/274 (96%)	0.29	7 (2%) 58 53	4, 9, 19, 26	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	77[A]	GLY	3.2
1	B	74[A]	ASN	2.3
1	B	7	THR	2.2
1	B	76[A]	SER	2.2
1	B	82	TRP	2.2
1	B	80	THR	2.1
1	A	-2	HIS	2.1

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.