



wwPDB NMR Structure Validation Summary Report ⓘ

Feb 12, 2017 – 06:36 pm GMT

PDB ID : 1K45
Title : The Solution Structure of the CBM4-2 Carbohydrate Binding Module from a
Thermostable Rhodothermus marinus Xylanase.
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Deposited on : 2001-10-05

This is a wwPDB NMR 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

<http://wwpdb.org/validation/2016/NMRValidationReportHelp>

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:

Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	trunk28760
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

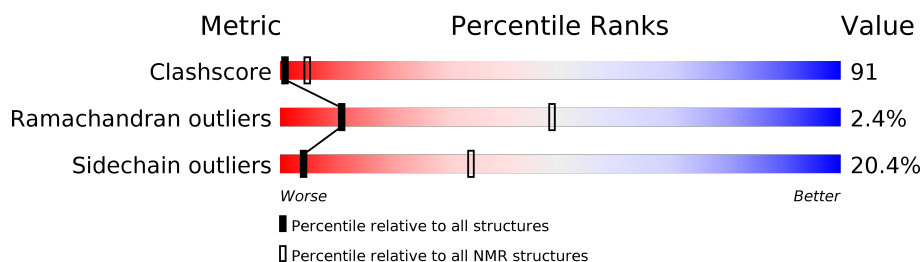
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 92%.

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)	NMR archive (#Entries)
Clashscore	125131	11601
Ramachandran outliers	121729	10391
Sidechain outliers	121581	10367

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	168	<div> <div></div> <div>23%</div> <div>60%</div> <div>17%</div> <div>.</div> </div>

2 Ensemble composition and analysis ⓘ

This entry contains 1 models. Identification of well-defined residues and clustering analysis are not possible.

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2490 atoms, of which 1208 are hydrogens and 0 are deuteriums.

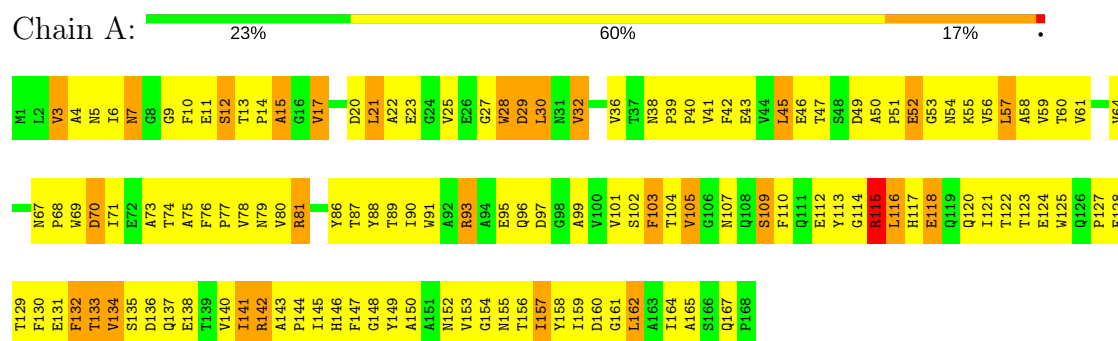
- Molecule 1 is a protein called Xylanase.

Mol	Chain	Residues	Atoms						Trace
1	A	168	Total	C	H	N	O	S	0
			2490	814	1208	209	258	1	

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	Initiating Methionine	UNP P96988
A	165	ALA	-	CLONING ARTIFACT	UNP P96988
A	166	SER	-	CLONING ARTIFACT	UNP P96988
A	167	GLN	-	CLONING ARTIFACT	UNP P96988
A	168	PRO	-	CLONING ARTIFACT	UNP P96988

- Molecule 1: Xylanase



5 Refinement protocol and experimental data overview

The models were refined using the following method: *hybrid distance geometry/simulated annealing using XPLOR*.

Of the ? calculated structures, 1 were deposited, based on the following criterion: ?.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
X-PLOR	structure solution	3.1
X-PLOR	refinement	3.1

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	BMRB entry 5181
Number of chemical shift lists	1
Total number of shifts	2023
Number of shifts mapped to atoms	2023
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	92%

No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

There are no covalent bond-length or bond-angle outliers.

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	Planarity
1	A	0	4
All	All	0	4

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

All planar outliers are listed below.

Mol	Chain	Res	Type	Group
1	A	93	ARG	Sidechain
1	A	81	ARG	Sidechain
1	A	115	ARG	Sidechain
1	A	142	ARG	Sidechain

6.2 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 each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	1282	1208	1208	226
All	All	1282	1208	1208	226

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 91.

5 of 226 clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:45:LEU:HD21	1:A:47:THR:HG23	0.92	1.40
1:A:45:LEU:HD21	1:A:47:THR:CG2	0.90	1.97
1:A:21:LEU:HD13	1:A:30:LEU:O	0.88	1.67
1:A:3:VAL:CG2	1:A:165:ALA:HB2	0.84	2.02
1:A:99:ALA:HB1	1:A:152:ASN:OD1	0.82	1.74

6.3 Torsion angles [i](#)

6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	166/168 (99%)	136 (82%)	26 (16%)	4 (2%)	11	48
All	All	166/168 (99%)	136 (82%)	26 (16%)	4 (2%)	11	48

All 4 Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	3	VAL
1	A	15	ALA
1	A	134	VAL
1	A	118	GLU

6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	137/137 (100%)	109 (80%)	28 (20%)	4	34
All	All	137/137 (100%)	109 (80%)	28 (20%)	4	34

5 of 28 residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	133	THR
1	A	30	LEU
1	A	135	SER
1	A	70	ASP
1	A	12	SER

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 92% for the well-defined parts and 92% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 5181

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	2023
Number of shifts mapped to atoms	2023
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	14

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	167	-0.06 ± 0.14	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	151	-0.16 ± 0.13	None needed (< 0.5 ppm)
$^{13}\text{C}'$	165	0.40 ± 0.07	None needed (< 0.5 ppm)
^{15}N	156	-0.10 ± 0.40	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 92%, i.e. 1757 atoms were assigned a chemical shift out of a possible 1917. 28 out of 28 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	811/820 (99%)	323/326 (99%)	332/336 (99%)	156/158 (99%)
Sidechain	798/914 (87%)	483/525 (92%)	292/356 (82%)	23/33 (70%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	148/183 (81%)	77/97 (79%)	67/80 (84%)	4/6 (67%)
Overall	1757/1917 (92%)	883/948 (93%)	691/772 (90%)	183/197 (93%)

7.1.4 Statistically unusual chemical shifts ⓘ

The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	67	ASN	HB3	-0.33	4.41 – 1.11	-9.4
1	A	118	GLU	HG3	0.70	3.31 – 1.21	-7.4
1	A	93	ARG	HD3	1.29	4.36 – 1.86	-7.3
1	A	67	ASN	HB2	0.58	4.36 – 1.26	-7.2
1	A	57	LEU	HB3	-0.89	3.34 – -0.26	-6.8
1	A	93	ARG	HG3	-0.38	3.00 – 0.10	-6.7
1	A	51	PRO	HG3	-0.04	3.56 – 0.26	-5.9
1	A	118	GLU	HG2	1.10	3.33 – 1.23	-5.6
1	A	93	ARG	HD2	1.84	4.27 – 1.97	-5.6
1	A	127	PRO	HB3	0.04	3.81 – 0.21	-5.5
1	A	51	PRO	HB3	0.05	3.81 – 0.21	-5.4
1	A	51	PRO	HB2	0.19	3.82 – 0.32	-5.4
1	A	51	PRO	HD3	1.69	5.52 – 1.72	-5.1
1	A	92	ALA	HA	6.49	6.46 – 2.06	5.1

7.1.5 Random Coil Index (RCI) plots ⓘ

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition.

Random coil index (RCI) for chain A:

