wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	1HEJ
Title	:	C-TERMINAL XYLAN BINDING DOMAIN FROM CELLULOMONAS FIMI XYLANASE 11A
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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 (i) symbol.

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

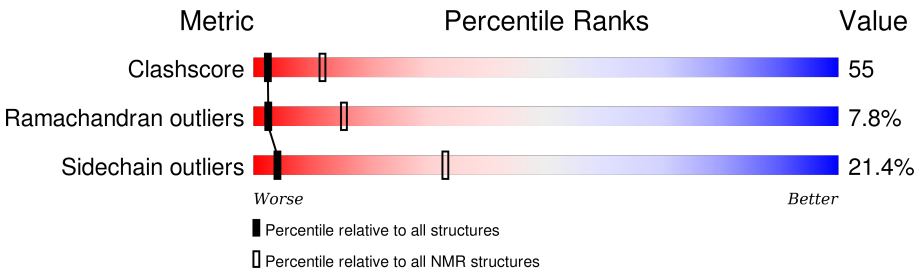
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	rb-20027457
Ideal 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:
SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

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	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

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 <=5%

Mol	Chain	Length	Quality of chain
1	C	88	20% 49% 11% • 18%

2 Ensemble composition and analysis (i)

This entry contains 5 models. Model 1 is the overall representative, medoid model (most similar to other models).

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	**Residue range (total)**	**Backbone RMSD (Å)**	**Medoid model**
1	C:562-C:581, C:585-C:630, C:636-C:641 (72)	0.47	1

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 1 clusters and 1 single-model cluster was found.

Cluster number	Models
1	1, 3, 4, 5
Single-model clusters	2

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 1185 atoms, of which 570 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called ENDO-1,4-BETA-XYLANASE D.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	C	88	1185	372	570	109	132	2	0

4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA and DNA chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: ENDO-1,4-BETA-XYLANASE D

Chain C: 20% 49% 11% • 18%

T557 G558 S559 C560 S561 V562 S563 A564 V565 R566 G567 E568 E569 W570 A571 D572 R573 F574 N575 V576 T577 Y578 S579 V580 S581 G582 S583 S584 S585 W586 V587 V588 T589 L590 G591 L592 N593 Q596 S597 V598 Q599 S600 S601 W602 N603 A604 A605 L606 S609 S610 G611 T612 V613 T614 A615 R616 S620

G621 N622 S623 F624 G625 V626 T627 F628 Y629 K630 N631 G632 S633 S634 A635 T636 P637 G638 A639 T640 C641 A642 T643 G644

4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 1. Colouring as in section 4.1 above.

- Molecule 1: ENDO-1,4-BETA-XYLANASE D

Chain C: 27% 36% 17% • 18%

T557 G558 S559 C560 S561 V562 S563 A564 V565 R566 E569 D572 R573 F574 N575 V576 T577 Y578 S579 V580 S581 G582 S583 S584 S585 W586 V587 V588 T589 L590 G591 L592 N593 G594 G595 Q596 Q599 S600 S601 W602 N603 A604 A605 L606 S609 S610 G611 T612 V613 T614 A615 R616 S620 G621 N622

S623 F624 G625 V626 T627 F628 Y629 K630 N631 G632 S633 S634 A635 T636 P637 G638 A639 T640 C641 A642 T643 G644

5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *HYBRID DISTANCE GEOMETRY/ SIMULATED ANNEALING.*

Of the 50 calculated structures, 5 were deposited, based on the following criterion: *RANDOM SELECTION OF 5 FROM THE 23 LOWEST ENERGY STRUCTURES.*

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

Software name	Classification	Version
X-PLOR 3.1	refinement	
MSI FELIX 97.0	structure solution	

No chemical shift data was provided. 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	C	0.0±0.0	2.8±0.4
All	All	0	14

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

All unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
1	C	566	ARG	Sidechain	5
1	C	616	ARG	Sidechain	5
1	C	573	ARG	Sidechain	4

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	C	524	492	492	56±9
All	All	2620	2460	2460	281

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 55.

5 of 182 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:C:589:THR:HG23	1:C:612:THR:HG23	1.00	1.27	1	2
1:C:588:VAL:HG23	1:C:640:THR:O	1.00	1.56	1	3
1:C:580:VAL:HG21	1:C:586:TRP:CZ2	0.97	1.94	5	4
1:C:574:PHE:CE1	1:C:636:THR:HG22	0.94	1.97	4	1
1:C:580:VAL:HG21	1:C:586:TRP:CH2	0.93	1.99	1	4

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	C	72/88 (82%)	53±3 (74±4%)	13±2 (18±2%)	6±1 (8±2%)	2	15
All	All	360/440 (82%)	267 (74%)	65 (18%)	28 (8%)	2	15

5 of 11 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	C	604	ALA	5
1	C	587	VAL	4
1	C	567	GLY	3
1	C	610	SER	3
1	C	620	SER	3

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	C	56/66 (85%)	44±1 (79±3%)	12±1 (21±3%)	4	33
All	All	280/330 (85%)	220 (79%)	60 (21%)	4	33

5 of 27 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	C	573	ARG	4
1	C	566	ARG	4
1	C	630	LYS	4
1	C	603	ASN	4
1	C	640	THR	4

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 (i)

No chemical shift data were provided