Full wwPDB NMR Structure Validation Report (i)

Mar 1, 2022 – 03:24 PM EST

PDB ID	:	2FFW
Title	:	Solution structure of the RBCC/TRIM B-box1 domain of human MID1: B-box with a RING
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Deposited on	:	2005-12-20

This is a Full wwPDB NMR Structure Validation 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/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:

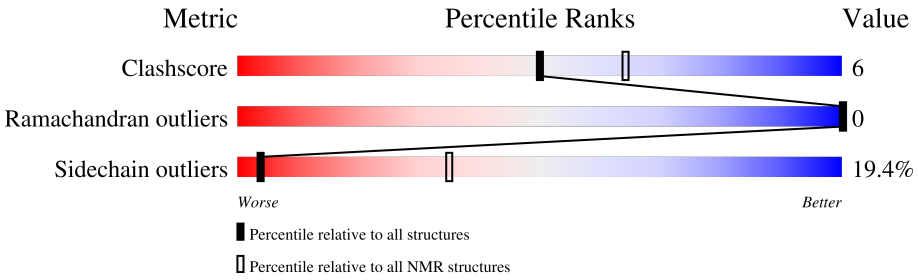
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.27
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.27

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	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

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	A	78	31% 13% 56%

2 Ensemble composition and analysis (i)

This entry contains 13 models. Model 5 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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	A:118-A:151 (34)	0.13	5

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 3 clusters and 1 single-model cluster was found.

Cluster number	Models
1	1, 2, 3, 5, 12, 13
2	4, 7, 9, 10
3	6, 11
Single-model clusters	8

3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1168 atoms, of which 568 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Midline-1.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	78	1166	363	568	107	121	7	0

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	
			Total	Zn
2	A	2	2	2

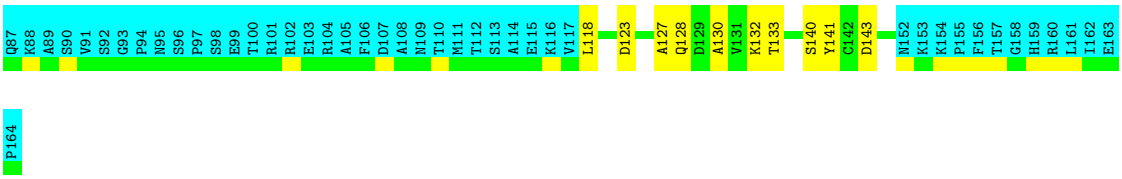
4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide 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: Midline-1

Chain A: 31% 13% 56%



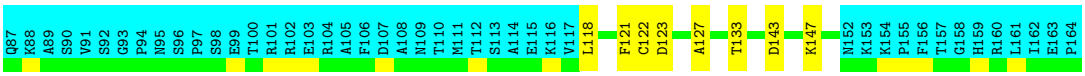
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

• Molecule 1: Midline-1

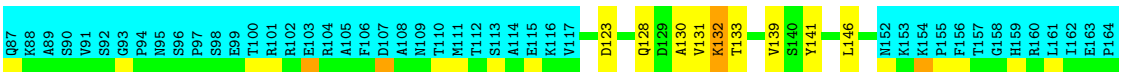
Chain A: 33% 10% 56%



4.2.2 Score per residue for model 2

• Molecule 1: Midline-1

Chain A: 32% 10% • 56%



4.2.3 Score per residue for model 3

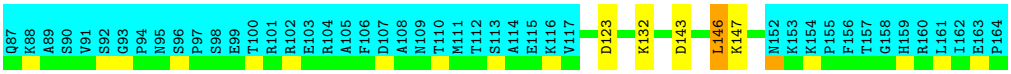
- Molecule 1: Midline-1

Chain A: 35% 9% 56%

4.2.4 Score per residue for model 4

- Molecule 1: Midline-1

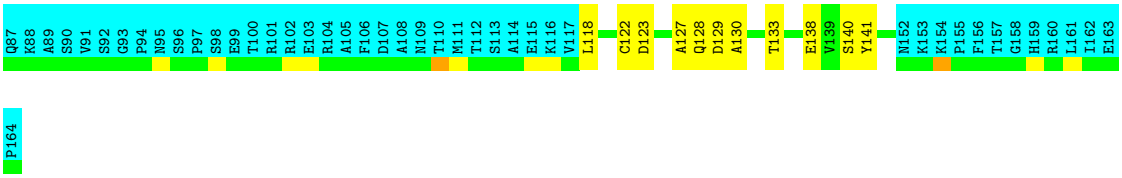
Chain A: 37% 5% • 56%



4.2.5 Score per residue for model 5 (medoid)

- Molecule 1: Midline-1

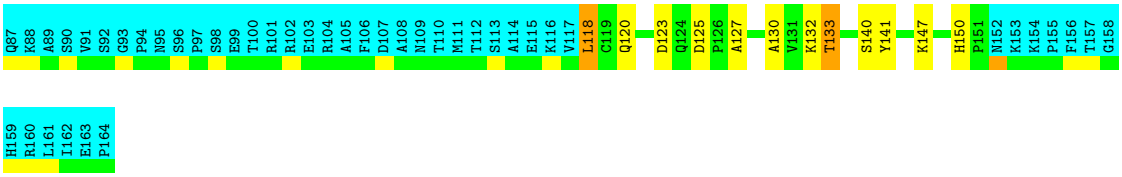
Chain A: 29% 14% 56%



4.2.6 Score per residue for model 6

- Molecule 1: Midline-1

Chain A: 28% 13% • 56%



4.2.7 Score per residue for model 7

- Molecule 1: Midline-1

Chain A: 29% 12% • 56%

Q87 K88 A89 S90 V91 S92 G93 P94 N95 S96 P97 S98 E99 T100 R101 R102 E103 R104 A105 F106 D107 A108 N109 T110 M111 T112 S113 A114 E115 K116 V117 F121 C122 D123 A127 Q128 D129 A130 V131 K132 T133 S140 Y141 C142 D143 N152 K153 K154 P155 F156 T157 G158 H159 R160 L161 I162

E163 P164

4.2.8 Score per residue for model 8

- Molecule 1: Midline-1

Chain A: 28% 10% 5% 56%

Q87 K88 A89 S90 V91 S92 G93 P94 N95 S96 P97 S98 E99 T100 R101 R102 E103 R104 A105 F106 D107 A108 N109 T110 M111 T112 S113 A114 E115 K116 V117 L118 D123 A127 Q128 D129 A130 V131 K132 T133 S140 D143 K147 N152 K153 K154 P155 F156 T157 G158 H159 R160 L161 I162

E163 P164

4.2.9 Score per residue for model 9

- Molecule 1: Midline-1

Chain A: 31% 12% • 56%

Q87 K88 A89 S90 V91 S92 G93 P94 N95 S96 P97 S98 E99 T100 R101 R102 E103 R104 A105 F106 D107 A108 N109 T110 M111 T112 S113 A114 E115 K116 V117 L118 C122 D123 A127 V131 K132 T133 S140 E144 C145 L146 N152 K153 K154 P155 F156 T157 G158 H159 R160 L161 I162 E163

P164

4.2.10 Score per residue for model 10

- Molecule 1: Midline-1

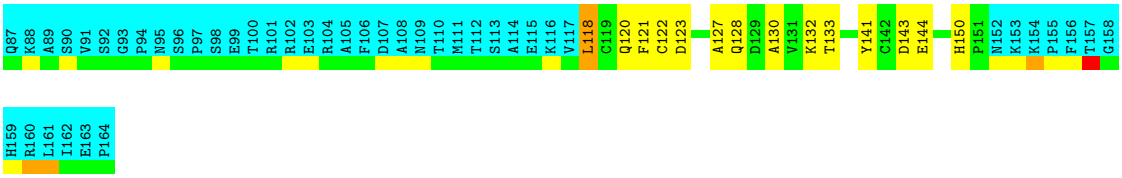
Chain A: 35% 8% • 56%

Q87 K88 A89 S90 V91 S92 G93 P94 N95 S96 P97 S98 E99 T100 R101 R102 E103 R104 A105 F106 D107 A108 N109 T110 M111 T112 S113 A114 E115 K116 V117 Q120 D123 A130 T133 S140 Y141 C142 D143 N152 K153 K154 P155 F156 T157 G158 H159 R160 L161 I162 E163 P164

4.2.11 Score per residue for model 11

- Molecule 1: Midline-1

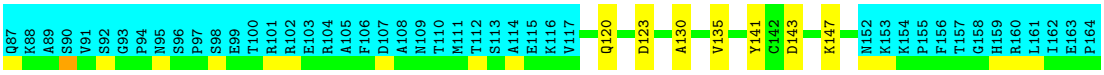
Chain A: 26% 17% • 56%



4.2.12 Score per residue for model 12

- Molecule 1: Midline-1

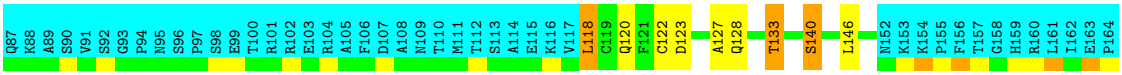
Chain A: 35% 9% 56%



4.2.13 Score per residue for model 13

- Molecule 1: Midline-1

Chain A: 32% 8% • 56%



5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *torsion angle dynamics*.

Of the 100 calculated structures, 13 were deposited, based on the following criterion: *target function*.

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

Software name	Classification	Version
CYANA	structure solution	2.1
CYANA	refinement	2.1

No chemical shift data was provided.

6 Model quality (i)

6.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	257	230	231	3±1
All	All	3367	2990	3003	38

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 6.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:118:LEU:HD22	1:A:127:ALA:HB3	0.70	1.64	5	5
1:A:131:VAL:HG23	1:A:132:LYS:HD2	0.61	1.71	8	2
1:A:133:THR:HG23	1:A:140:SER:CB	0.60	2.27	6	1
1:A:130:ALA:HB1	1:A:141:TYR:O	0.59	1.98	6	8
1:A:132:LYS:CD	1:A:146:LEU:HD22	0.57	2.30	4	1
1:A:133:THR:HG23	1:A:140:SER:HB3	0.54	1.79	13	2
1:A:133:THR:HG23	1:A:140:SER:CA	0.53	2.34	7	2
1:A:133:THR:HG23	1:A:140:SER:HA	0.50	1.82	7	2
1:A:132:LYS:HD3	1:A:146:LEU:HD22	0.49	1.84	4	1
1:A:131:VAL:HG23	1:A:132:LYS:CD	0.49	2.38	8	1
1:A:133:THR:HG23	1:A:140:SER:HB2	0.48	1.84	6	1
1:A:122:CYS:CB	1:A:127:ALA:HB1	0.45	2.41	7	6

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:133:THR:HG23	1:A:139:VAL:C	0.45	2.32	2	1
1:A:132:LYS:HE2	1:A:146:LEU:HD22	0.43	1.91	2	1
1:A:132:LYS:CE	1:A:146:LEU:HD22	0.42	2.45	2	1
1:A:118:LEU:CD2	1:A:127:ALA:HB3	0.41	2.42	8	1
1:A:131:VAL:HG23	1:A:132:LYS:HE3	0.41	1.91	9	1
1:A:118:LEU:HD23	1:A:128:GLN:O	0.41	2.15	8	1

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	34/78 (44%)	28±1 (81±3%)	6±1 (19±3%)	0±0 (0±0%)	100	100
All	All	442/1014 (44%)	358 (81%)	84 (19%)	0 (0%)	100	100

There are no Ramachandran outliers.

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	31/69 (45%)	25±2 (81±6%)	6±2 (19±6%)	4	35
All	All	403/897 (45%)	325 (81%)	78 (19%)	4	35

All 17 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	A	123	ASP	13
1	A	133	THR	10

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Mol	Chain	Res	Type	Models (Total)
1	A	118	LEU	7
1	A	143	ASP	7
1	A	128	GLN	6
1	A	132	LYS	6
1	A	120	GLN	6
1	A	147	LYS	5
1	A	140	SER	4
1	A	121	PHE	3
1	A	146	LEU	2
1	A	129	ASP	2
1	A	150	HIS	2
1	A	144	GLU	2
1	A	138	GLU	1
1	A	125	ASP	1
1	A	135	VAL	1

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 monosaccharides in this entry.

6.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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