wwPDB NMR Structure Validation Summary Report (i)

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Title	:	MAGI-1 PDZ1 / E6CT
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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:

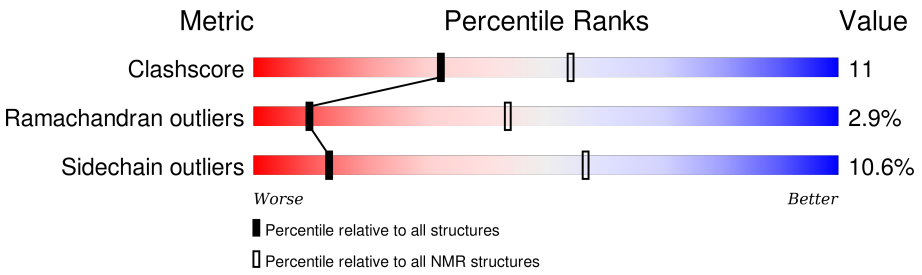
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	rb-20027457
Ideal 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 is 90%.

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	A	129	51% 22% • 25%
2	B	11	36% 64%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 12 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:5-A:101, B:158-B:161 (101)	0.30	12

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 5 clusters and 3 single-model clusters were found.

Cluster number	Models
1	3, 4, 9, 11, 12, 19
2	8, 15, 20
3	2, 6, 17
4	1, 10, 16
5	5, 18
Single-model clusters	7; 13; 14

3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2152 atoms, of which 1081 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Membrane-associated guanylate kinase, WW and PDZ domain-containing protein 1.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	129	1955	621	980	162	188	4	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	EXPRESSION TAG	UNP Q96QZ7
A	-2	ALA	-	EXPRESSION TAG	UNP Q96QZ7
A	-1	MET	-	EXPRESSION TAG	UNP Q96QZ7

- Molecule 2 is a protein called Protein E6.

Mol	Chain	Residues	Atoms					Trace
			Total	C	H	N	O	
2	B	11	197	53	101	24	19	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	161	VAL	-	EXPRESSION TAG	UNP P03126

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: Membrane-associated guanylate kinase, WW and PDZ domain-containing protein 1

Chain A: 51% 22% • 25%

G-3 A-2 M-1 G0 K1 P2 F3 F4 T5 R6 N7 E10 K14 F15 I16 H17 T18 K19 L20 R21 K22 F27 G28 V31 V32 G33 G34 D38 L41 K44 G50 P51 L54 D62 V63 N68 L73 G74 H75 I87 V93 D94 L97 C98 R99

P102 L103 P104 F105 D106 P107 D108 D109 P110 N111 T112 S113 L114 V115 T116 S117 V118 A119 I120 L121 D122 K123 E124 P125

• Molecule 2: Protein E6

Chain B: 36% 64%

R151 S152 S153 R154 T155 R156 R157 V161

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

The representative model is number 12. Colouring as in section 4.1 above.

• Molecule 1: Membrane-associated guanylate kinase, WW and PDZ domain-containing protein 1

Chain A: 48% 23% • 25%

G-3 A-2 M-1 G0 K1 P2 F3 F4 T5 R6 N7 P8 S9 E10 G13 K14 F15 I16 K19 L20 R21 K22 S23 S24 F27 G28 F29 T30 V31 G34 P37 L41 S45 L46 V47 G50 P51 L54 M58 V63 N68 G74 H75 V80 V81 K82

I87 V93 D94 L97 C98 R99 P102 L103 P104 F105 D106 P107 D108 D109 P110 N111 T112 S113 L114 V115 T116 S117 V118 A119 I120 L121 D122 K123 E124 P125

• Molecule 2: Protein E6

Chain B:

5 Refinement protocol and experimental data overview (i)

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

Of the 64 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
ATNOS-CANDID	structure solution	
X-PLOR NIH	refinement	

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 16559
Number of chemical shift lists	1
Total number of shifts	1421
Number of shifts mapped to atoms	1421
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	90%

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)

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 (average) 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	1.09±0.02	1±1/747 (0.1±0.1%)	0.90±0.02	0±0/1010 (0.0±0.0%)
2	B	1.00±0.10	0±0/32 (0.0±0.0%)	0.76±0.08	0±0/41 (0.0±0.0%)
All	All	1.09	12/15580 (0.1%)	0.89	2/21020 (0.0%)

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	0.0±0.0	0.1±0.4
All	All	0	3

All unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	9	SER	N-CA	-6.30	1.33	1.46	6	10
1	A	6	ARG	C-N	-5.10	1.22	1.34	11	2

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})	Models	
								Worst	Total
1	A	99	ARG	NE-CZ-NH2	-5.25	117.67	120.30	9	1
1	A	9	SER	N-CA-CB	-5.12	102.82	110.50	20	1

There are no chirality outliers.

All unique planar outliers are listed below.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	99	ARG	Sidechain	3

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	734	744	744	17±4
2	B	33	30	30	1±1
All	All	15340	15480	15480	348

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.

5 of 116 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:68:ASN:ND2	1:A:93:VAL:HB	0.79	1.93	12	1
1:A:41:LEU:HD21	1:A:80:VAL:HG21	0.71	1.62	18	8
1:A:7:ASN:O	1:A:10:GLU:HG2	0.70	1.87	12	20
1:A:22:LYS:HG3	1:A:87:ILE:CG2	0.64	2.21	2	9
1:A:31:VAL:O	2:B:158:GLU:HA	0.63	1.93	20	6

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	97/129 (75%)	84±2 (86±2%)	10±2 (11±2%)	3±1 (3±1%)	9 42
2	B	3/11 (27%)	3±0 (97±10%)	0±0 (3±10%)	0±0 (0±0%)	100 100
All	All	2000/2800 (71%)	1731 (87%)	211 (11%)	58 (3%)	9 43

5 of 10 unique Ramachandran outliers are listed below. They are sorted by the frequency of

occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	5	THR	13
1	A	68	ASN	9
1	A	24	SER	8
1	A	38	ASP	8
1	A	22	LYS	7

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	82/110 (75%)	73±2 (89±2%)	9±2 (11±2%)	13	56
2	B	4/11 (36%)	4±0 (92±11%)	0±0 (8±11%)	21	67
All	All	1720/2420 (71%)	1538 (89%)	182 (11%)	13	57

5 of 30 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	19	LYS	20
1	A	16	ILE	20
1	A	75	HIS	17
1	A	14	LYS	16
1	A	22	LYS	16

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)

The completeness of assignment taking into account all chemical shift lists is 90% for the well-defined parts and 75% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 16559

Chemical shift list name: *assigned_ chem_ shift_ list_ 1*

7.1.1 Bookkeeping (i)

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	1421
Number of shifts mapped to atoms	1421
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	5

7.1.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	109	0.03 ± 0.09	None needed (< 0.5 ppm)
$^{13}C_{\beta}$	96	0.20 ± 0.16	None needed (< 0.5 ppm)
$^{13}C'$	98	0.03 ± 0.27	None needed (< 0.5 ppm)
^{15}N	97	-0.32 ± 0.33	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments (i)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 90%, i.e. 1078 atoms were assigned a chemical shift out of a possible 1193. 19 out of 20 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	473/497 (95%)	196/198 (99%)	187/202 (93%)	90/97 (93%)
Sidechain	542/622 (87%)	332/361 (92%)	202/236 (86%)	8/25 (32%)

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	Total	^{1}H	^{13}C	^{15}N
Aromatic	63/74 (85%)	34/41 (83%)	29/30 (97%)	0/3 (0%)
Overall	1078/1193 (90%)	562/600 (94%)	418/468 (89%)	98/125 (78%)

7.1.4 Statistically unusual chemical shifts (i)

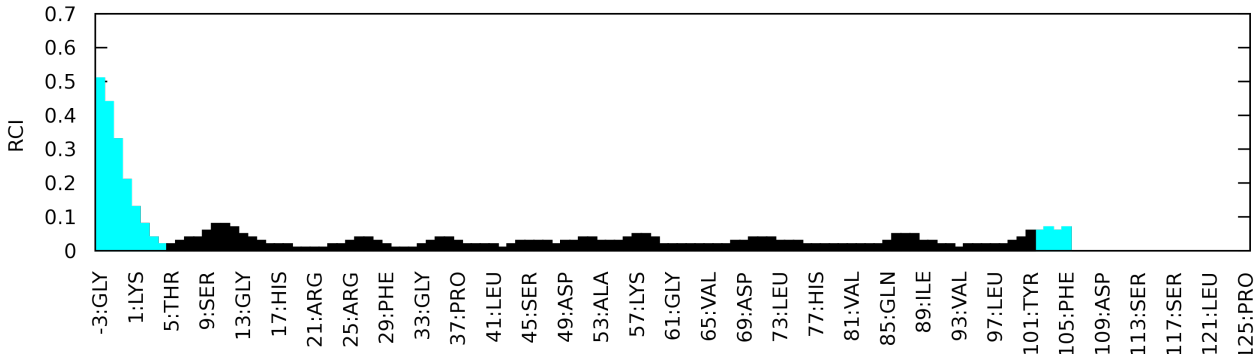
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, *ppm*	Expected range, *ppm*	Z-score
1	A	99	ARG	NE	115.29	92.63 – 76.73	19.3
1	A	6	ARG	NE	115.12	92.63 – 76.73	19.1
1	A	99	ARG	HH11	9.78	9.21 – 4.61	6.2
1	A	22	LYS	HG3	-0.35	2.76 – -0.04	-6.1
1	A	38	ASP	HA	2.73	6.15 – 3.05	-6.0

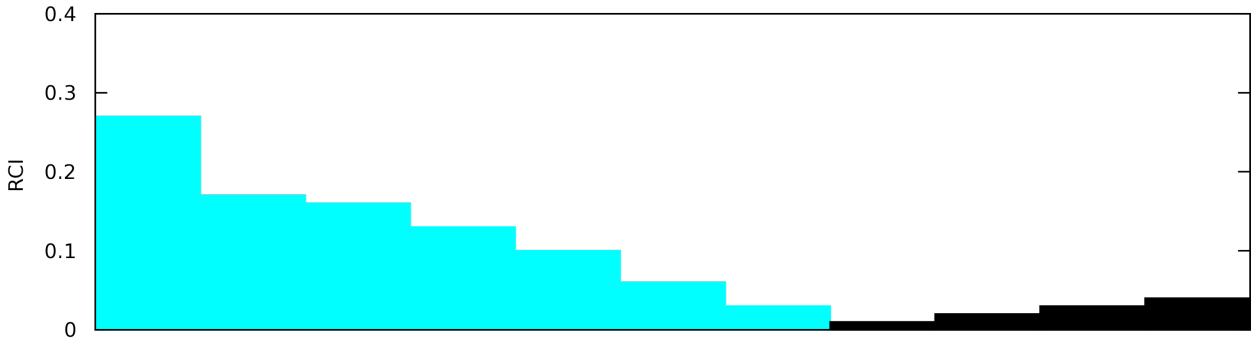
7.1.5 Random Coil Index (RCI) plots (i)

The images below report *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:



Random coil index (RCI) for chain B:


0.4
0.3
0.2
0.1
0
RCI