wwPDB NMR Structure Validation Summary Report (i)

Apr 21, 2024 – 03:28 PM EDT

PDB ID	:	2KQF
Title	:	Solution structure of MAST205-PDZ complexed with the C-terminus of a rabies virus G protein
Authors	:	Terrien, E.; Wolff, N.; Cordier, F.; Simenel, C.; Bernard, A.; Lafon, M.; Delepierre, M.
Deposited on	:	2009-11-04

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
https://www.wwpdb.org/validation/2017/NMRValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

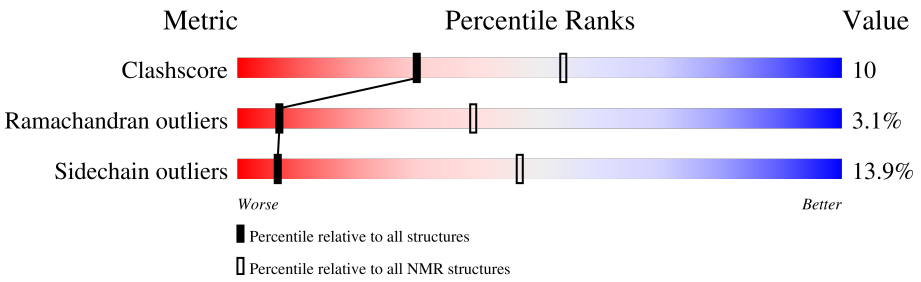
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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	96	55% 26% • 17%
2	B	13	100%

2 Ensemble composition and analysis (i)

This entry contains 10 models. Model 4 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:3-A:22, A:36-A:95 (80)	0.20	4

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. No single-model clusters were found.

Cluster number	Models
1	5, 6, 7, 10
2	1, 3, 4
3	2, 8, 9

3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1670 atoms, of which 831 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Microtubule-associated serine/threonine-protein kinase 2.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	96	1470	463	735	134	135	3	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP Q6P0Q8

- Molecule 2 is a protein called C-terminal motif from Glycoprotein.

Mol	Chain	Residues	Atoms					Trace
			Total	C	H	N	O	
2	B	13	200	62	96	20	22	0

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: Microtubule-associated serine/threonine-protein kinase 2

Chain A: 55% 26% • 17%

G1 G2 I8 I9 I10 H11 K16 Y17 G18 L21 R22 A23 I24 R25 V26 Y27 M28 G29 D30 S31 D32 V33 Y34 T35 H38 M39 V40 V43 E44 D45 P48 A49 S50 L54 R55 D58 H69 V72 V76 L79 I80 V87 A88 I89 S90 T91

L94 E95 N96

• Molecule 2: C-terminal motif from Glycoprotein

Chain B: 100%

S101 W102 E103 S104 H105 K106 S107 G108 G109 E110 T111 R112 L113

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

The representative model is number 4. Colouring as in section 4.1 above.

• Molecule 1: Microtubule-associated serine/threonine-protein kinase 2

Chain A: 49% 34% 17%

G1 G2 P6 P7 I8 I9 I10 H11 K16 Y17 G18 L21 R22 A23 I24 R25 V26 Y27 M28 G29 D30 S31 D32 V33 Y34 T35 H38 M39 V40 V43 E44 D45 P48 A49 S50 L54 R55 D58 L59 I60 N64 H69 V72 E75 V76 L79 I80

V87 A88 I89 S90 T91 L94 E95 N96

• Molecule 2: C-terminal motif from Glycoprotein

Chain B: 100%

S101
W102
E103
S104
H105
K106
S107
G108
G109
E110
T111
R112
L113

5 Refinement protocol and experimental data overview (i)

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

Of the 200 calculated structures, 10 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
ARIA	structure solution	
CNS	refinement	

No chemical shift data was provided.

6 Model quality (i)

6.1 Standard geometry (i)

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	615	624	624	12±2
2	B	0	0	0	0±0
All	All	6150	6240	6240	125

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 10.

5 of 32 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:40:VAL:HG11	1:A:54:LEU:HD21	0.72	1.60	9	7
1:A:80:ILE:HG23	1:A:87:VAL:HG11	0.57	1.74	6	9
1:A:10:ILE:HG21	1:A:17:TYR:HB3	0.56	1.78	7	8
1:A:38:HIS:CD2	1:A:76:VAL:HG11	0.55	2.37	1	9
1:A:45:ASP:HA	1:A:50:SER:CB	0.54	2.32	2	8

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	80/96 (83%)	66±1 (83±2%)	11±1 (14±2%)	2±1 (3±1%)	7 39
2	B	0	-	-	-	-
All	All	800/1090 (73%)	665 (83%)	110 (14%)	25 (3%)	7 39

5 of 6 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	48	PRO	10
1	A	69	HIS	10
1	A	18	GLY	2
1	A	15	LYS	1
1	A	93	PRO	1

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	67/79 (85%)	58±1 (86±2%)	9±1 (14±2%)	7 47
2	B	0	-	-	-
All	All	670/900 (74%)	577 (86%)	93 (14%)	7 47

5 of 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	21	LEU	10
1	A	55	ARG	10
1	A	79	LEU	10
1	A	94	LEU	10
1	A	95	GLU	10

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)

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