wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	1S4X
Title	:	NMR Structure of the integrin B3 cytoplasmic domain in DPC micelles
Authors	:	Vinogradova, O.; Vaynberg, J.; Kong, X.; Haas, T.A.; Plow, E.F.; Qin, J.
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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
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:

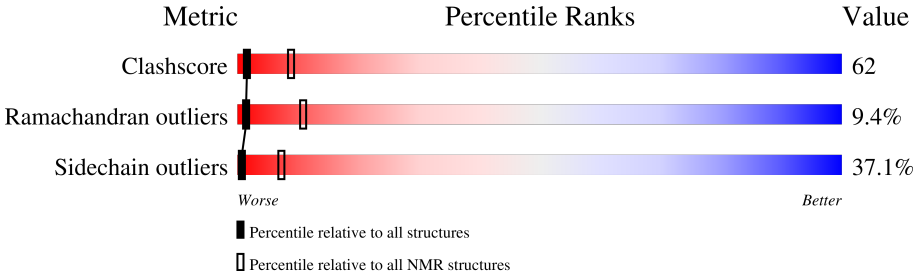
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.26
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.26

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	67	

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 15 is the overall representative, medoid model (most similar to other models). The authors have identified model 6 as representative, based on the following criterion: *fewest violations*.

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:21-A:34 (14)	0.20	16
2	A:35-A:64 (30)	0.40	15

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 4 clusters and 1 single-model cluster was found.

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

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 785 atoms, of which 390 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Integrin beta-3.

Mol	Chain	Residues	Atoms					Trace
			Total	C	H	N	O	
1	A	47	785	250	390	70	75	0

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	cloning artifact	UNP P05106
A	2	SER	-	cloning artifact	UNP P05106
A	3	SER	-	cloning artifact	UNP P05106
A	4	HIS	-	cloning artifact	UNP P05106
A	5	HIS	-	cloning artifact	UNP P05106
A	6	HIS	-	cloning artifact	UNP P05106
A	7	HIS	-	cloning artifact	UNP P05106
A	8	HIS	-	cloning artifact	UNP P05106
A	9	HIS	-	cloning artifact	UNP P05106
A	10	SER	-	cloning artifact	UNP P05106
A	11	SER	-	cloning artifact	UNP P05106
A	12	GLY	-	cloning artifact	UNP P05106
A	13	LEU	-	cloning artifact	UNP P05106
A	14	VAL	-	cloning artifact	UNP P05106
A	15	PRO	-	cloning artifact	UNP P05106
A	16	ARG	-	cloning artifact	UNP P05106
A	17	GLY	-	cloning artifact	UNP P05106
A	18	SER	-	cloning artifact	UNP P05106
A	19	HIS	-	cloning artifact	UNP P05106
A	20	MET	-	cloning artifact	UNP P05106

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: Integrin beta-3

Chain A: 12% 33% 19% • • 30%

GLY SER SER HIS HIS HIS HIS HIS HIS SER SER GLY LEU VAL PRO ARG GLY SER HIS MET K21 L22 L23 I24 T25 I26 H27 D28 R29 K30 A33 K34 F35 E36 E37 E38 R39 A40 R41 A42 K43 W44 D45 T46 A47 N48 N49 P50 L51 Y52 K53 E54 A55 T56 S57 T58 F59 T60 N61

I62 T63 Y64 R65 G66 T67

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

The representative model is number 15. Colouring as in section 4.1 above.

- Molecule 1: Integrin beta-3

Chain A: 10% 31% 24% • 30%

GLY SER SER HIS HIS HIS HIS HIS HIS SER SER GLY LEU VAL PRO ARG GLY SER HIS MET K21 L22 L23 I24 T25 D28 R29 K30 E31 F32 A33 K34 F35 E36 E37 E38 R39 A40 R41 A42 K43 W44 D45 T46 A47 N48 N49 P50 L51 Y52 K53 E54 A55 T56 S57 T58 F59 T60 N61

I62 T63 Y64 R65 G66 T67

5 Refinement protocol and experimental data overview (i)

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

Of the 88 calculated structures, 20 were deposited, based on the following criterion: *structures with the least restraint violations*.

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

Software name	Classification	Version
X-PLOR	structure solution	
X-PLOR	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	372	367	366	46±7
All	All	7440	7340	7320	913

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 62.

5 of 153 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:39:ARG:O	1:A:41:ARG:N	1.18	1.76	16	20
1:A:49:ASN:HD22	1:A:51:LEU:H	0.90	1.10	15	6
1:A:39:ARG:C	1:A:41:ARG:H	0.82	1.78	16	20
1:A:48:ASN:ND2	1:A:49:ASN:H	0.73	1.82	15	6
1:A:41:ARG:O	1:A:45:ASP:N	0.69	2.26	14	20

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	43/67 (64%)	29±2 (68±3%)	10±2 (23±4%)	4±1 (9±2%)	1 11
All	All	860/1340 (64%)	584 (68%)	195 (23%)	81 (9%)	1 11

5 of 7 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	40	ALA	20
1	A	45	ASP	20
1	A	48	ASN	18
1	A	50	PRO	15
1	A	64	TYR	6

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	39/58 (67%)	25±2 (63±4%)	14±2 (37±4%)	1 7
All	All	780/1160 (67%)	491 (63%)	289 (37%)	1 7

5 of 26 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	24	ILE	20
1	A	43	LYS	20
1	A	46	THR	20
1	A	49	ASN	20
1	A	51	LEU	20

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