wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	1IKC
Title	:	NMR Structure of alpha-Bungarotoxin
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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:

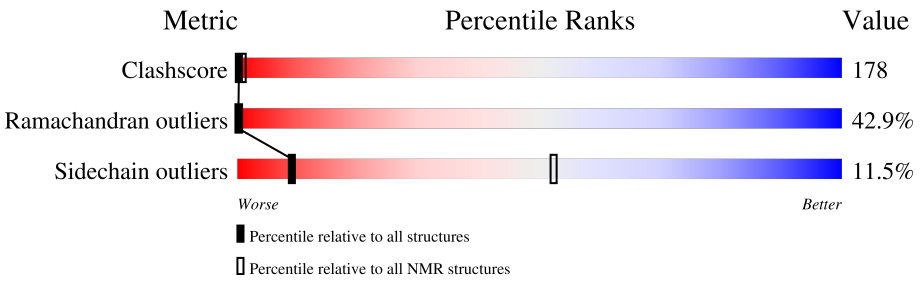
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.26
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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	74	11% 32% 34% 15% 8%

2 Ensemble composition and analysis (i)

This entry contains 30 models. The atoms present in the NMR models are not consistent. Some calculations may have failed as a result. All residues are included in the validation scores. Model 10 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	A:2-A:16, A:22-A:74 (68)	0.16	10

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 3 single-model clusters were found.

Cluster number	Models
1	1, 2, 3, 4, 5, 6, 7, 8, 9, 10, 11, 12, 13, 14, 15, 16, 18, 19, 22, 23, 24, 26, 29
2	20, 21
3	17, 25
Single-model clusters	27; 28; 30

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 694 atoms, of which 143 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called long neurotoxin 1.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	74	694	338	143	97	105	11	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: long neurotoxin 1

Chain A: 11% 32% 34% 15% 8%

I1 V2 C3 H4 T5 T6 A7 T8 S9 P10 I11 S12 A13 V14 T15 C16 P17 P18 G19 E20 N21 L22 C23 Y24 R25 K26 M27 W28 C29 D30 A31 S35 R36 G37 K38 V39 V40 E41 L42 G43 C44 A45 A46 T47 C48 K51 K52 P53 Y54 E55 E56 V57 T58 C59 C60 S61 T62 D63

K64 C65 N66 P67 H68 P69 K70 Q71 R72 P73 G74

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

The representative model is number 10. Colouring as in section 4.1 above.

- Molecule 1: long neurotoxin 1

Chain A: 9% 34% 32% 16% 8%

I1 V2 C3 H4 T5 T6 A7 T8 S9 P10 I11 S12 A13 V14 T15 C16 P17 P18 G19 E20 N21 L22 C23 Y24 R25 K26 M27 W28 C29 D30 A31 S35 R36 G37 K38 V39 V40 E41 L42 G43 C44 A45 A46 T47 C48 P49 S50 K51 K52 P53 Y54 E55 E56 V57 T58 C59 C60 S61 T62

D63 K64 C65 N66 P67 H68 P69 K70 Q71 R72 P73 G74

5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *distance geometry, simulated annealing, molecular dynamics.*

Of the 100 calculated structures, 30 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
DIANA	structure solution	1.5
Amber	refinement	4.1

No chemical shift data was provided.

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	1.01±0.12	2±1/520 (0.3± 0.3%)	2.12±0.18	25±4/705 (3.5± 0.6%)
All	All	1.02	50/15600 (0.3%)	2.13	736/21150 (3.5%)

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	2.9±0.9
All	All	0	86

5 of 21 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	68	HIS	CG-CD2	-14.58	1.10	1.35	15	2
1	A	71	GLN	CD-NE2	-10.70	1.06	1.32	15	1
1	A	28	TRP	CE2-CZ2	-9.60	1.23	1.39	21	2
1	A	68	HIS	CB-CG	9.24	1.66	1.50	15	2
1	A	28	TRP	CD2-CE2	-8.73	1.30	1.41	11	5

5 of 99 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(°)	Ideal(°)	Models	
								Worst	Total
1	A	5	THR	OG1-CB-CG2	-24.30	54.11	110.00	7	21
1	A	68	HIS	ND1-CG-CD2	-23.06	73.72	106.00	24	5
1	A	68	HIS	CG-CD2-NE2	20.50	148.14	109.20	15	5
1	A	6	THR	CA-CB-CG2	15.27	133.77	112.40	21	8
1	A	5	THR	CA-CB-CG2	14.77	133.07	112.40	25	12

There are no chirality outliers.

5 of 14 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	A	24	TYR	Sidechain,Mainchain	28
1	A	43	GLY	Mainchain,Peptide	12
1	A	68	HIS	Sidechain	10
1	A	37	GLY	Mainchain	9
1	A	6	THR	Mainchain	7

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	508	135	489	178±19
All	All	15240	4050	14687	5330

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 178.

5 of 676 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:68:HIS:HD2	1:A:71:GLN:NE2	1.14	1.37	24	11
1:A:5:THR:CB	1:A:64:LYS:HZ3	1.12	1.56	25	2
1:A:6:THR:HB	1:A:42:LEU:HD21	1.11	1.14	24	29
1:A:42:LEU:HD23	1:A:66:ASN:HB3	1.11	1.14	30	3
1:A:23:CYS:HA	1:A:44:CYS:SG	1.10	1.87	1	10

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	67/74 (91%)	13±1 (19±2%)	25±2 (38±3%)	29±2 (43±3%)	0 0
All	All	2010/2220 (91%)	390 (19%)	757 (38%)	863 (43%)	0 0

5 of 39 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	7	ALA	30
1	A	9	SER	30
1	A	11	ILE	30
1	A	12	SER	30
1	A	13	ALA	30

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	60/65 (92%)	53±4 (88±7%)	7±4 (12±7%)	9 52
All	All	1800/1950 (92%)	1593 (88%)	207 (12%)	9 52

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	24	TYR	30
1	A	42	LEU	30
1	A	55	GLU	26
1	A	16	CYS	24
1	A	44	CYS	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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	30-A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
30	A	15:THR	C	16:CYS	N	1.16

7 Chemical shift validation (i)

No chemical shift data were provided