Full wwPDB NMR Structure Validation Report (i)

Jun 6, 2023 – 12:35 AM EDT

PDB ID	:	2MBO
BMRB ID	:	19406
Title	:	K11-linked Diubiquitin average solution structure at pH 6.8, 0 mM NaCl
Authors	:	Castaneda, C.A.; Fushman, D.
Deposited on	:	2013-08-02

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

The following versions of software and data (see references (1)) were used in the production of this report:

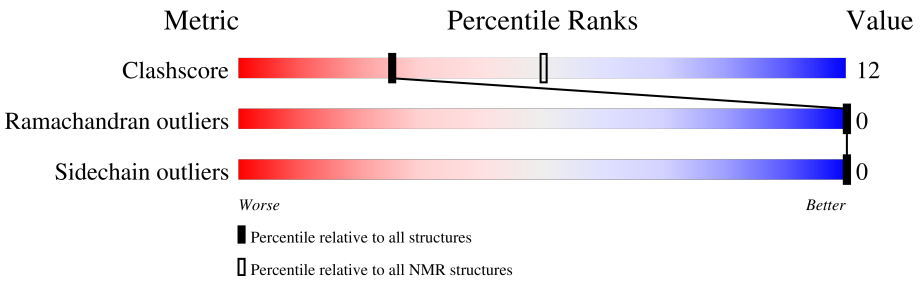
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
BMRB Restraints Analysis	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.33

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 6%.

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	76	87% 7% • 5%
1	B	76	86% 9% 5%

2 Ensemble composition and analysis (i)

This entry contains 1 models. Identification of well-defined residues and clustering analysis are not possible.

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2346 atoms, of which 1198 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Ubiquitin.

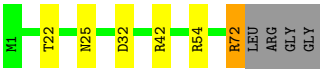
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	72	1173	362	599	98	113	1	0
1	B	72	1173	362	599	98	113	1	0

4 Residue-property plots (i)

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.

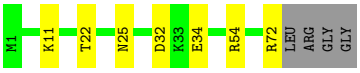
- Molecule 1: Ubiquitin

Chain A:



- Molecule 1: Ubiquitin

Chain B:



5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *in-house RDC (orientation) and MTSL (distance).*

Of the 1 calculated structures, 1 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
in-house	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)	working_cs.cif
Number of chemical shift lists	2
Total number of shifts	280
Number of shifts mapped to atoms	264
Number of unparsed shifts	0
Number of shifts with mapping errors	16
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	6%

6 Model quality (i)

6.1 Standard geometry (i)

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

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	2
1	B	0	2
All	All	0	4

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

All planar outliers are listed below.

Mol	Chain	Res	Type	Group
1	A	54	ARG	Sidechain
1	A	72	ARG	Sidechain
1	B	54	ARG	Sidechain
1	B	72	ARG	Sidechain

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	574	599	599	24
1	B	574	599	599	24
All	All	1148	1198	1198	27

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 12.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:72:ARG:C	1:B:34:GLU:O	1.30	1.66
1:A:72:ARG:CG	1:B:34:GLU:HA	1.19	1.66
1:A:72:ARG:HG3	1:B:34:GLU:OE1	1.17	0.99
1:A:72:ARG:CG	1:B:34:GLU:OE1	1.17	1.91
1:A:72:ARG:HG2	1:B:34:GLU:CA	1.10	1.76
1:A:72:ARG:HG2	1:B:34:GLU:HA	1.02	1.12
1:A:72:ARG:O	1:B:34:GLU:O	1.00	1.78
1:A:72:ARG:HG3	1:B:34:GLU:CD	0.94	1.82
1:A:72:ARG:CG	1:B:34:GLU:CA	0.84	2.46
1:A:72:ARG:HG2	1:B:34:GLU:C	0.77	1.98
1:A:72:ARG:NE	1:B:34:GLU:HA	0.73	1.99
1:A:72:ARG:HG2	1:B:34:GLU:O	0.68	1.89
1:A:42:ARG:NH1	1:B:11:LYS:HE2	0.67	2.04
1:A:72:ARG:CD	1:B:34:GLU:HA	0.66	2.18
1:A:42:ARG:HH12	1:B:11:LYS:CE	0.64	2.05
1:A:42:ARG:HH12	1:B:11:LYS:HE2	0.60	1.54
1:A:42:ARG:HH12	1:B:11:LYS:NZ	0.56	1.96
1:A:42:ARG:NH1	1:B:11:LYS:CE	0.55	2.70
1:A:72:ARG:CG	1:B:34:GLU:CD	0.46	2.64
1:A:72:ARG:NE	1:B:34:GLU:CA	0.45	2.77
1:B:32:ASP:OD1	1:B:32:ASP:O	0.43	2.37
1:A:32:ASP:O	1:A:32:ASP:OD1	0.43	2.37
1:A:42:ARG:NH1	1:B:11:LYS:NZ	0.41	2.65
1:A:22:THR:OG1	1:A:25:ASN:OD1	0.41	2.39
1:B:22:THR:OG1	1:B:25:ASN:OD1	0.41	2.39
1:B:32:ASP:OD1	1:B:32:ASP:C	0.40	2.60
1:A:32:ASP:OD1	1:A:32:ASP:C	0.40	2.60

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	70/76 (92%)	70 (100%)	0 (0%)	0 (0%)	100	100
1	B	70/76 (92%)	70 (100%)	0 (0%)	0 (0%)	100	100
All	All	140/152 (92%)	140 (100%)	0 (0%)	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	66/68 (97%)	66 (100%)	0 (0%)	100	100
1	B	66/68 (97%)	66 (100%)	0 (0%)	100	100
All	All	132/136 (97%)	132 (100%)	0 (0%)	100	100

There are no protein residues with a non-rotameric sidechain to report.

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)

The completeness of assignment taking into account all chemical shift lists is 6% for the well-defined parts and 6% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

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	140
Number of shifts mapped to atoms	132
Number of unparsed shifts	0
Number of shifts with mapping errors	8
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- No matching atom found in the structure. All 8 occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	73	LEU	H	8.357	.	.
1	A	73	LEU	N	124.023	.	.
1	A	74	ARG	H	8.518	.	.
1	A	74	ARG	N	121.993	.	.
1	A	75	GLY	H	8.546	.	.
1	A	75	GLY	N	110.52	.	.
1	A	76	GLY	H	8.38	.	.
1	A	76	GLY	N	108.997	.	.

7.1.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	0	—	None (insufficient data)
$^{13}C_{\beta}$	0	—	None (insufficient data)
$^{13}C'$	0	—	None (insufficient data)
^{15}N	70	1.11 ± 0.42	Should be applied

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 6%, i.e. 132 atoms were assigned a chemical shift out of a possible 2036. 0 out of 24 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	132/716 (18%)	66/290 (23%)	0/288 (0%)	66/138 (48%)
Sidechain	0/1248 (0%)	0/806 (0%)	0/394 (0%)	0/48 (0%)
Aromatic	0/72 (0%)	0/36 (0%)	0/34 (0%)	0/2 (0%)
Overall	132/2036 (6%)	66/1132 (6%)	0/716 (0%)	66/188 (35%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 6%, i.e. 132 atoms were assigned a chemical shift out of a possible 2036. 0 out of 24 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	132/716 (18%)	66/290 (23%)	0/288 (0%)	66/138 (48%)
Sidechain	0/1248 (0%)	0/806 (0%)	0/394 (0%)	0/48 (0%)
Aromatic	0/72 (0%)	0/36 (0%)	0/34 (0%)	0/2 (0%)
Overall	132/2036 (6%)	66/1132 (6%)	0/716 (0%)	66/188 (35%)

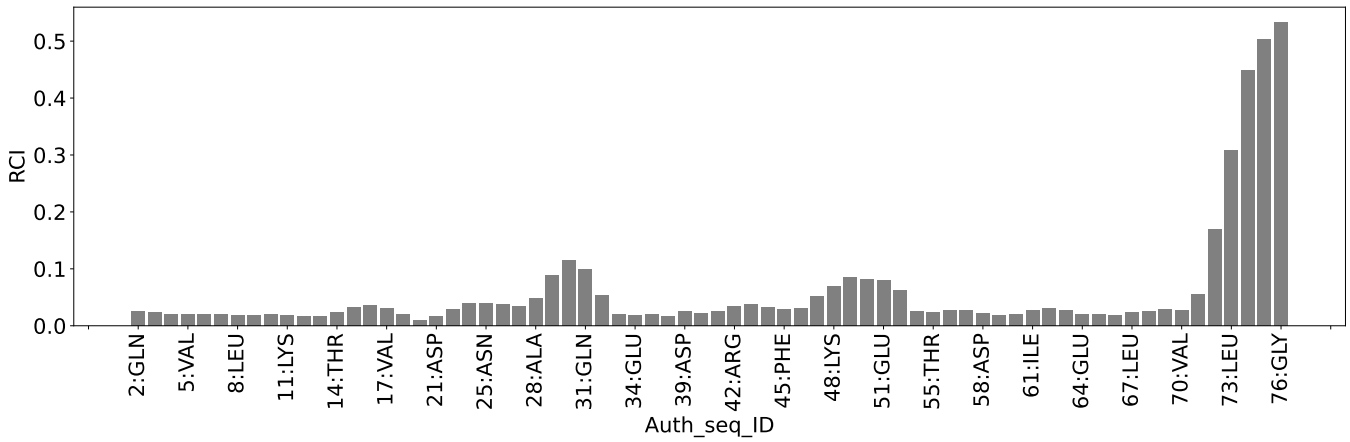
7.1.4 Statistically unusual chemical shifts (i)

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots (i)

The image below reports *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.If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:



7.2 Chemical shift list 2

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_2*

7.2.1 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	140
Number of shifts mapped to atoms	132
Number of unparsed shifts	0
Number of shifts with mapping errors	8
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- No matching atom found in the structure. All 8 occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
2	A	73	LEU	H	8.389	.	.
2	A	73	LEU	N	124.45	.	.
2	A	74	ARG	H	8.442	.	.
2	A	74	ARG	N	121.767	.	.
2	A	75	GLY	H	8.555	.	.
2	A	75	GLY	N	110.768	.	.
2	A	76	GLY	H	8.192	.	.

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
2	A	76	GLY	N	109.395	.	.

7.2.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	0	—	None (insufficient data)
$^{13}C_{\beta}$	0	—	None (insufficient data)
$^{13}C'$	0	—	None (insufficient data)
^{15}N	70	1.07 ± 0.50	Should be applied

7.2.3 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 6%, i.e. 132 atoms were assigned a chemical shift out of a possible 2036. 0 out of 24 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	132/716 (18%)	66/290 (23%)	0/288 (0%)	66/138 (48%)
Sidechain	0/1248 (0%)	0/806 (0%)	0/394 (0%)	0/48 (0%)
Aromatic	0/72 (0%)	0/36 (0%)	0/34 (0%)	0/2 (0%)
Overall	132/2036 (6%)	66/1132 (6%)	0/716 (0%)	66/188 (35%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 6%, i.e. 132 atoms were assigned a chemical shift out of a possible 2036. 0 out of 24 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	132/716 (18%)	66/290 (23%)	0/288 (0%)	66/138 (48%)
Sidechain	0/1248 (0%)	0/806 (0%)	0/394 (0%)	0/48 (0%)
Aromatic	0/72 (0%)	0/36 (0%)	0/34 (0%)	0/2 (0%)
Overall	132/2036 (6%)	66/1132 (6%)	0/716 (0%)	66/188 (35%)

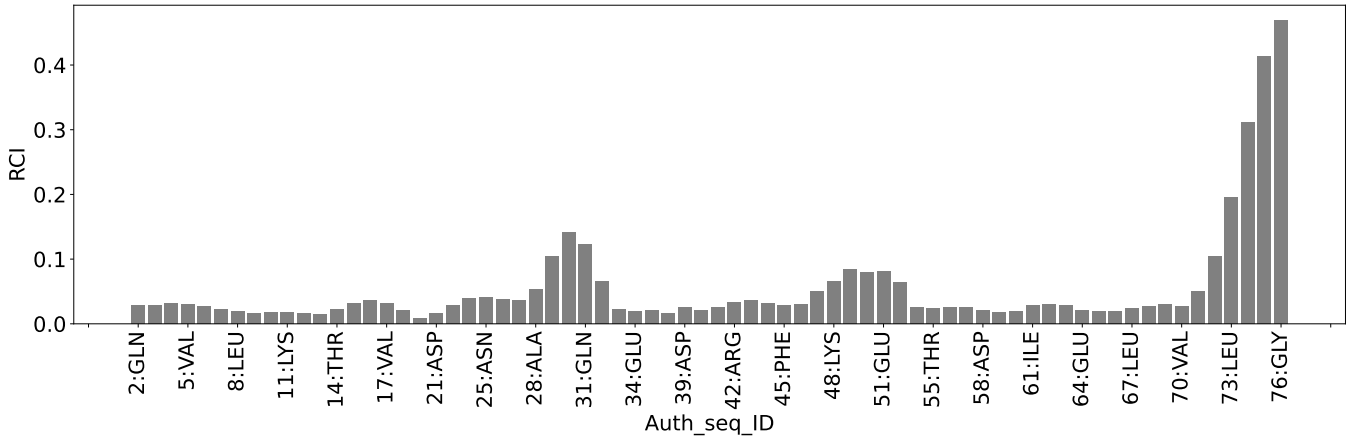
7.2.4 Statistically unusual chemical shifts (i)

There are no statistically unusual chemical shifts.

7.2.5 Random Coil Index (RCI) plots (i)

The image below reports *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.If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:



8 NMR restraints analysis (i)

No restraints data found

9 Distance violation analysis (i)

No distance restraints data found

10 Dihedral-angle violation analysis (i)

No dihedral-angle restraints found