



Full wwPDB NMR Structure Validation Report ⓘ

Apr 27, 2016 – 05:06 AM BST

PDB ID : 2RS8
Title : Solution structure of the N-terminal RNA recognition motif of NonO
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Deposited on : 2011-11-29

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
<http://wwpdb.org/validation/2016/NMRValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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

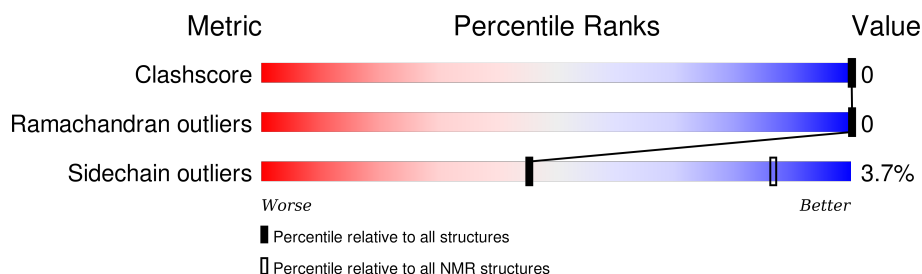
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR


The overall completeness of chemical shifts assignment is 88%.

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	99	 <div>72% 28%</div>

2 Ensemble composition and analysis

This entry contains 20 models. Model 9 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 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:76-A:146 (71)	0.16	9

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1539 atoms, of which 773 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Non-POU domain-containing octamer-binding protein.

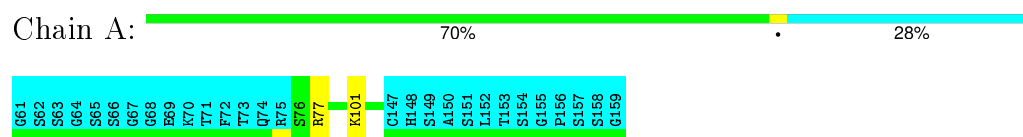
Mol	Chain	Residues	Atoms						Trace
1	A	99	Total	C	H	N	O	S	0
			1539	479	773	139	145	3	

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	61	GLY	-	EXPRESSION TAG	UNP Q99K48
A	62	SER	-	EXPRESSION TAG	UNP Q99K48
A	63	SER	-	EXPRESSION TAG	UNP Q99K48
A	64	GLY	-	EXPRESSION TAG	UNP Q99K48
A	65	SER	-	EXPRESSION TAG	UNP Q99K48
A	66	SER	-	EXPRESSION TAG	UNP Q99K48
A	67	GLY	-	EXPRESSION TAG	UNP Q99K48
A	154	SER	-	EXPRESSION TAG	UNP Q99K48
A	155	GLY	-	EXPRESSION TAG	UNP Q99K48
A	156	PRO	-	EXPRESSION TAG	UNP Q99K48
A	157	SER	-	EXPRESSION TAG	UNP Q99K48
A	158	SER	-	EXPRESSION TAG	UNP Q99K48
A	159	GLY	-	EXPRESSION TAG	UNP Q99K48

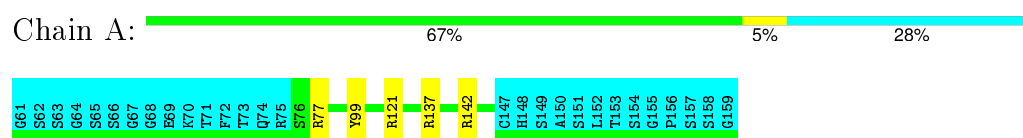
4.2.3 Score per residue for model 3

- Molecule 1: Non-POU domain-containing octamer-binding protein



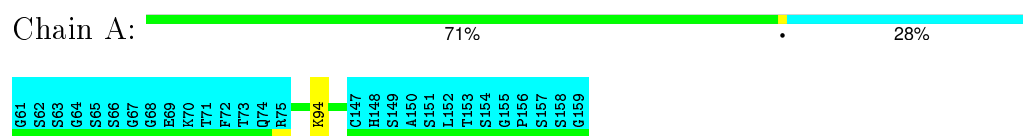
4.2.4 Score per residue for model 4

- Molecule 1: Non-POU domain-containing octamer-binding protein



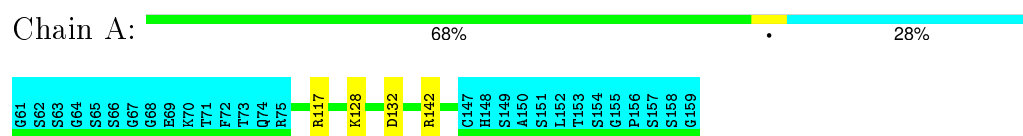
4.2.5 Score per residue for model 5

- Molecule 1: Non-POU domain-containing octamer-binding protein



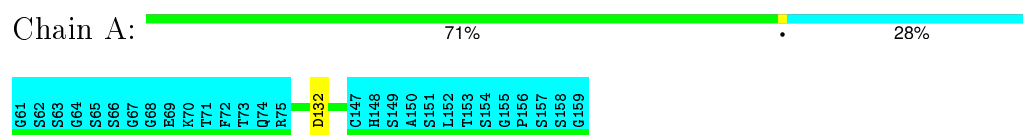
4.2.6 Score per residue for model 6

- Molecule 1: Non-POU domain-containing octamer-binding protein



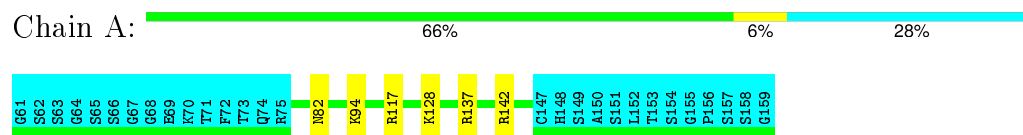
4.2.7 Score per residue for model 7

- Molecule 1: Non-POU domain-containing octamer-binding protein



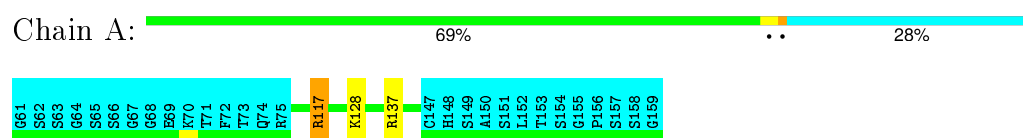
4.2.8 Score per residue for model 8

- Molecule 1: Non-POU domain-containing octamer-binding protein



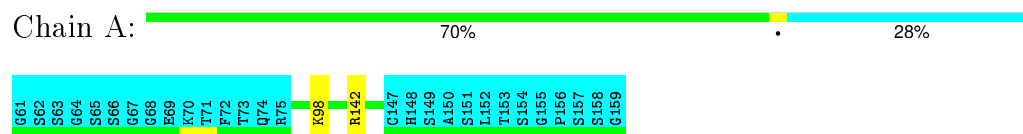
4.2.9 Score per residue for model 9 (medoid)

- Molecule 1: Non-POU domain-containing octamer-binding protein



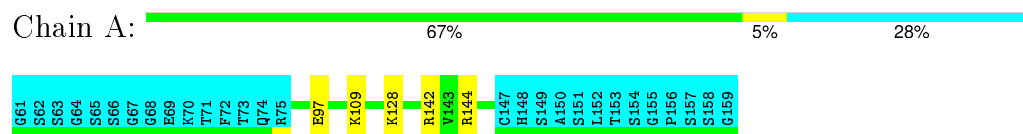
4.2.10 Score per residue for model 10

- Molecule 1: Non-POU domain-containing octamer-binding protein



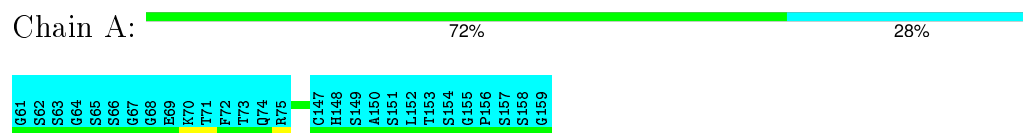
4.2.11 Score per residue for model 11

- Molecule 1: Non-POU domain-containing octamer-binding protein



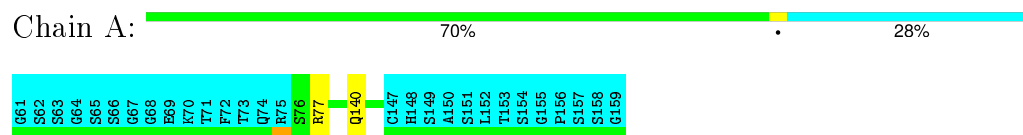
4.2.12 Score per residue for model 12

- Molecule 1: Non-POU domain-containing octamer-binding protein



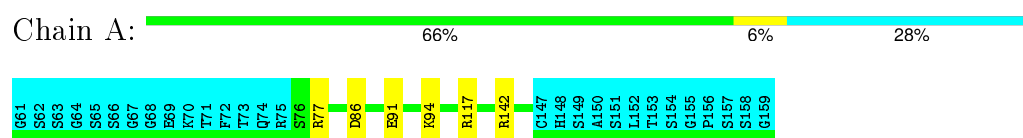
4.2.13 Score per residue for model 13

- Molecule 1: Non-POU domain-containing octamer-binding protein



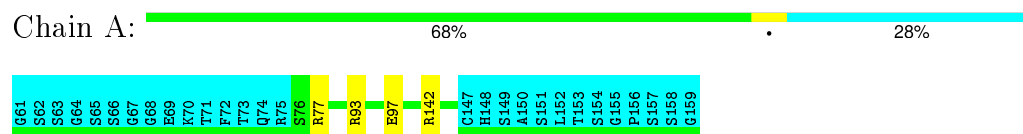
4.2.14 Score per residue for model 14

- Molecule 1: Non-POU domain-containing octamer-binding protein



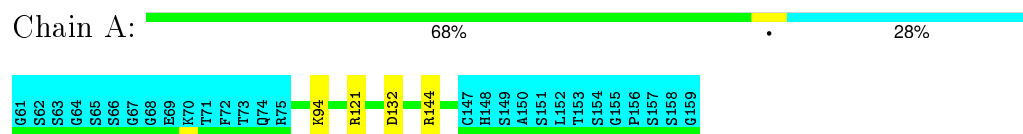
4.2.15 Score per residue for model 15

- Molecule 1: Non-POU domain-containing octamer-binding protein



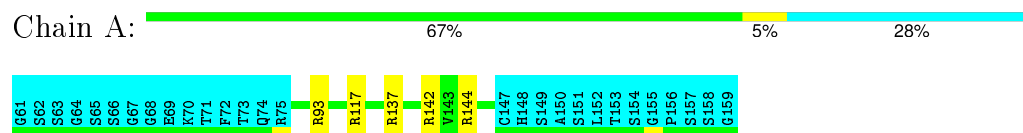
4.2.16 Score per residue for model 16

- Molecule 1: Non-POU domain-containing octamer-binding protein



4.2.17 Score per residue for model 17

- Molecule 1: Non-POU domain-containing octamer-binding protein





5 Refinement protocol and experimental data overview

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

Of the 200 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
AMBER	refinement	
CYANA	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)	2rs8_cs.str
Number of chemical shift lists	1
Total number of shifts	1091
Number of shifts mapped to atoms	1091
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	88%

No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality

6.1 Standard geometry

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	0.69±0.00	0±0/592 (0.0±0.0%)	0.92±0.03	1±2/791 (0.1±0.2%)
All	All	0.69	0/11840 (0.0%)	0.92	23/15820 (0.1%)

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	0.2±0.4
All	All	0	4

There are no bond-length outliers.

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(°)	Ideal(°)	Models	
								Worst	Total
1	A	117	ARG	NE-CZ-NH1	7.06	123.83	120.30	18	4
1	A	137	ARG	NE-CZ-NH1	6.50	123.55	120.30	2	4
1	A	77	ARG	NE-CZ-NH1	6.33	123.47	120.30	3	4
1	A	142	ARG	NE-CZ-NH1	6.17	123.39	120.30	4	1
1	A	93	ARG	NE-CZ-NH1	6.09	123.34	120.30	17	2
1	A	77	ARG	NE-CZ-NH2	-6.05	117.28	120.30	2	2
1	A	93	ARG	NE-CZ-NH2	-5.67	117.47	120.30	17	1
1	A	121	ARG	NE-CZ-NH1	5.66	123.13	120.30	2	3
1	A	144	ARG	NE-CZ-NH1	5.33	122.97	120.30	17	1
1	A	137	ARG	NE-CZ-NH2	-5.02	117.79	120.30	17	1

There are no chirality outliers.

All unique planar outliers are listed below.

Mol	Chain	Res	Type	Group	Models (Total)
1	A	99	TYR	Sidechain	4

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
All	All	11620	12060	12060	-

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 -.

There are no clashes.

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	71/99 (72%)	70±1 (99±1%)	1±1 (1±1%)	0±0 (0±0%)	100	100
All	All	1420/1980 (72%)	1407 (99%)	13 (1%)	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	61/82 (74%)	59±2 (96±2%)	2±2 (4±2%)	45	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1220/1640 (74%)	1175 (96%)	45 (4%)	45 87

All 19 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	142	ARG	8
1	A	128	LYS	5
1	A	117	ARG	4
1	A	94	LYS	4
1	A	97	GLU	3
1	A	101	LYS	3
1	A	132	ASP	3
1	A	98	LYS	2
1	A	77	ARG	2
1	A	144	ARG	2
1	A	91	GLU	1
1	A	109	LYS	1
1	A	140	GLN	1
1	A	133	ASN	1
1	A	121	ARG	1
1	A	137	ARG	1
1	A	82	ASN	1
1	A	111	LYS	1
1	A	86	ASP	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.6 Ligand geometry

There are no ligands in this entry.

6.7 Other polymers

There are no such molecules in this entry.

6.8 Polymer linkage issues

There are no chain breaks in this entry.

7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 88% for the well-defined parts and 80% for the entire structure.

7.1 Chemical shift list 1

File name: 2rs8_cs.str

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping

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	1091
Number of shifts mapped to atoms	1091
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	86	0.04 ± 0.10	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	77	0.19 ± 0.12	None needed (< 0.5 ppm)
$^{13}\text{C}'$	78	0.32 ± 0.14	None needed (< 0.5 ppm)
^{15}N	71	0.29 ± 0.26	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 88%, i.e. 832 atoms were assigned a chemical shift out of a possible 949. 0 out of 12 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	334/349 (96%)	138/139 (99%)	135/142 (95%)	61/68 (90%)
Sidechain	438/531 (82%)	275/313 (88%)	160/187 (86%)	3/31 (10%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	60/69 (87%)	32/38 (84%)	28/30 (93%)	0/1 (0%)
Overall	832/949 (88%)	445/490 (91%)	323/359 (90%)	64/100 (64%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 80%, i.e. 974 atoms were assigned a chemical shift out of a possible 1219. 0 out of 13 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	397/487 (82%)	162/194 (84%)	164/198 (83%)	71/95 (75%)
Sidechain	508/647 (79%)	320/384 (83%)	184/227 (81%)	4/36 (11%)
Aromatic	69/85 (81%)	37/47 (79%)	32/36 (89%)	0/2 (0%)
Overall	974/1219 (80%)	519/625 (83%)	380/461 (82%)	75/133 (56%)

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.

Random coil index (RCI) for chain A:

