wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	1EMW
Title	:	SOLUTION STRUCTURE OF THE RIBOSOMAL PROTEIN S16 FROM THERMUS THERMOPHILUS
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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:

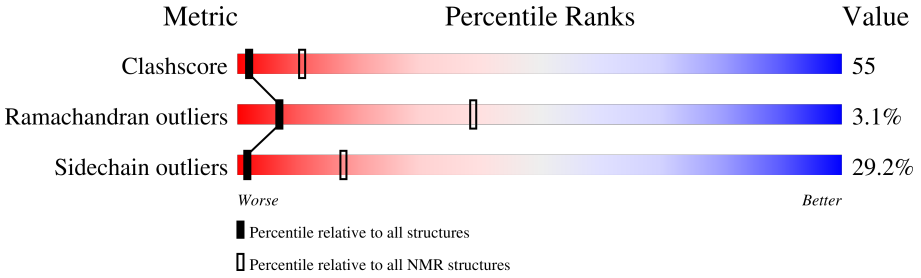
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.26
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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	88	7% 47% 6% • 40%

2 Ensemble composition and analysis (i)

This entry contains 47 models. Model 44 is the overall representative, medoid model (most similar to other models). The authors have identified model 11 as representative, based on the following criterion: *closest to the average*.

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:8, A:17-A:22, A:32-A:40, A:49-A:79 (53)	0.16	44

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 1 single-model cluster was found.

Cluster number	Models
1	1, 2, 3, 4, 5, 8, 9, 11, 12, 13, 14, 16, 17, 19, 21, 24, 25, 27, 28, 30, 31, 32, 33, 36, 37, 40, 41, 42, 43, 44, 45, 46, 47
2	6, 10, 18, 20, 26, 29, 35
3	15, 39
4	7, 22
5	34, 38
Single-model clusters	23

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 1487 atoms, of which 752 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called S16 RIBOSOMAL PROTEIN.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	88	1487	462	752	147	125	1	0

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	ARG	SER	conflict	UNP P80379
A	6	LEU	VAL	conflict	UNP P80379
A	7	ALA	SER	conflict	UNP P80379
A	24	ALA	VAL	conflict	UNP P80379
A	31	LYS	ALA	conflict	UNP P80379
A	68	ASP	-	SEE REMARK 999	UNP P80379
A	69	THR	-	SEE REMARK 999	UNP P80379
A	70	ALA	-	SEE REMARK 999	UNP P80379
A	71	ARG	-	SEE REMARK 999	UNP P80379
A	72	ARG	-	SEE REMARK 999	UNP P80379
A	73	LEU	-	SEE REMARK 999	UNP P80379
A	74	LEU	-	SEE REMARK 999	UNP P80379
A	75	ARG	-	SEE REMARK 999	UNP P80379
A	76	GLN	-	SEE REMARK 999	UNP P80379
A	77	ALA	-	SEE REMARK 999	UNP P80379
A	78	GLY	-	SEE REMARK 999	UNP P80379
A	79	VAL	-	SEE REMARK 999	UNP P80379
A	80	PHE	-	SEE REMARK 999	UNP P80379
A	81	ARG	-	SEE REMARK 999	UNP P80379
A	82	GLN	-	SEE REMARK 999	UNP P80379
A	83	GLU	-	SEE REMARK 999	UNP P80379
A	84	ALA	-	SEE REMARK 999	UNP P80379
A	85	ARG	-	SEE REMARK 999	UNP P80379
A	86	GLU	-	SEE REMARK 999	UNP P80379
A	87	GLY	-	SEE REMARK 999	UNP P80379
A	88	ALA	-	SEE REMARK 999	UNP P80379

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: S16 RIBOSOMAL PROTEIN

Chain A: 7% 47% 6% • 40%

M1 V2 K3 I4 R5 L6 A7 R8 F9 G10 S11 K12 H13 N14 P15 H16 Y17 R18 I19 V20 V21 T22 D23 A24 R25 R26 K27 R28 D29 G30 K31 Y32 I33 E34 K35 I36 G37 Y38 Y39 D40 P41 R42 K43 T44 T45 P46 D47 W48 L49 K50 V51 D52 V53 E54 R55 A56 R57 Y58 W59 L60

S61 V62 G63 A64 Q65 P66 T67 D68 T69 A70 R71 R72 L73 L74 R75 Q76 A77 G78 V79 F80 R81 Q82 E83 A84 R85 E86 G87 A88

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

The representative model is number 44. Colouring as in section 4.1 above.

- Molecule 1: S16 RIBOSOMAL PROTEIN

Chain A: 19% 34% 6% • 40%

M1 I4 R5 L6 A7 R8 F9 G10 S11 K12 H13 N14 P15 H16 Y17 R18 I19 V20 V21 T22 D23 A24 R25 R26 K27 R28 D29 G30 K31 Y32 I33 E34 K35 I36 G37 Y38 Y39 D40 P41 R42 K43 T44 T45 P46 D47 W48 L49 K50 V51 D52 V53 E54 R55 A56 R57 Y58 W59 L60 S61

V62 T67 A70 R71 R72 L73 L74 R75 F80 R81 Q82 E83 A84 R85 E86 G87 A88

5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *simulated annealing*.

Of the 50 calculated structures, 47 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
X-PLOR	structure solution	3.851
X-PLOR	refinement	3.851

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	0.99±0.00	0±0/449 (0.0± 0.0%)	1.34±0.01	0±0/609 (0.0± 0.0%)
All	All	0.99	0/21103 (0.0%)	1.34	4/28623 (0.0%)

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(^{o})	Ideal(^{o})	Models	
								Worst	Total
1	A	7	ALA	N-CA-CB	-5.12	102.93	110.10	15	3
1	A	59	TRP	CA-CB-CG	-5.02	104.17	113.70	36	1

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	441	467	467	50±6
All	All	20727	21949	21949	2363

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 55.

5 of 286 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:4:ILE:HD12	1:A:21:VAL:HG13	1.04	1.23	6	2
1:A:36:ILE:HG21	1:A:59:TRP:CD1	0.94	1.97	33	47
1:A:21:VAL:HG11	1:A:59:TRP:CD2	0.93	1.99	6	47
1:A:19:ILE:HD11	1:A:39:TYR:HB2	0.92	1.40	31	25
1:A:4:ILE:HD11	1:A:59:TRP:CB	0.91	1.94	39	42

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	53/88 (60%)	47±2 (88±3%)	5±1 (9±2%)	2±1 (3±1%)	7	39
All	All	2491/4136 (60%)	2195 (88%)	219 (9%)	77 (3%)	7	39

All 5 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	53	VAL	47
1	A	69	THR	17
1	A	67	THR	7
1	A	32	TYR	4
1	A	2	VAL	2

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	45/74 (61%)	32±2 (71±5%)	13±2 (29±5%)	1	17
All	All	2115/3478 (61%)	1498 (71%)	617 (29%)	1	17

5 of 32 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	20	VAL	47
1	A	53	VAL	45
1	A	21	VAL	43
1	A	19	ILE	38
1	A	75	ARG	35

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