wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	1A66
Title	:	SOLUTION NMR STRUCTURE OF THE CORE NFATC1/DNA COMPLEX, 18 STRUCTURES
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Deposited on	:	1998-03-06

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
http://wwpdb.org/validation/2016/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:

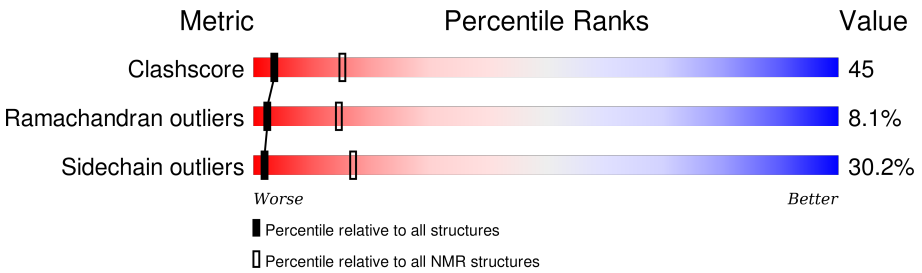
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	rb-20027457
Ideal 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 (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	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 <=5%

Mol	Chain	Length	Quality of chain
1	B	12	8% 92%
2	C	12	33% 67%
3	A	178	25% 58% 9% 8%

2 Ensemble composition and analysis (i)

This entry contains 18 models. Model 3 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:10-A:40, A:45-A:177 (164)	0.77	3

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 2 clusters and 3 single-model clusters were found.

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

3 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3595 atoms, of which 1704 are hydrogens and 0 are deuteriums.

- Molecule 1 is a DNA chain called DNA (5'-D(*CP*GP*AP*GP*GP*AP*AP*AP*AP*TP*TP*G)-3').

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	P	
1	B	12	385	119	136	52	67	11	0

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*AP*AP*TP*TP*TP*TP*CP*CP*TP*CP*G)-3').

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	P	
2	C	12	376	116	139	37	73	11	0

- Molecule 3 is a protein called CORE NFATC1.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
3	A	178	2834	875	1429	267	258	5	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	ALA	ENGINEERED	UNP O95644
A	2	LYS	LEU	ENGINEERED	UNP O95644
A	28	ARG	HIS	ENGINEERED	UNP O95644

4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

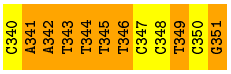
These plots are provided for all protein, RNA and DNA chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain 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: DNA (5'-D(*CP*GP*AP*GP*GP*AP*AP*AP*AP*TP*TP*G)-3')

Chain B:

- Molecule 2: DNA (5'-D(*CP*AP*AP*TP*TP*TP*TP*CP*CP*TP*CP*G)-3')

Chain C:



- Molecule 3: CORE NFATC1

Chain A:

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

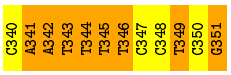
The representative model is number 3. Colouring as in section 4.1 above.

- Molecule 1: DNA (5'-D(*CP*GP*AP*GP*GP*AP*AP*AP*AP*TP*TP*G)-3')

Chain B:

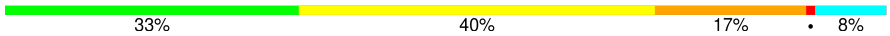
- Molecule 2: DNA (5'-D(*CP*AP*AP*TP*TP*TP*TP*CP*CP*TP*CP*G)-3')

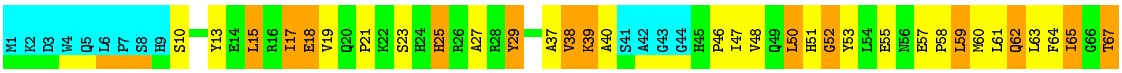
Chain C:

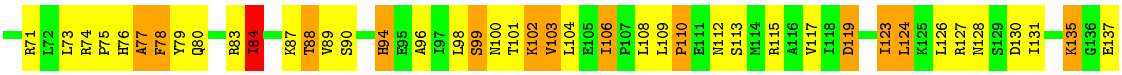


- Molecule 3: CORE NFATC1

Chain A:







5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *TORSION ANGLE DYNAMICS AND SIMULATED ANNEALING.*

Of the 18 calculated structures, 18 were deposited, based on the following criterion: *NOE VIOLATION <= 0.4 ANGSTROM, DIHEDRAL ANGLE VIOLATION <= 5 DEGREE.*

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

Software name	Classification	Version
X-PLOR	refinement	3.1
DYANA-1.4	structure solution	
X-PLOR	structure solution	3.1

No chemical shift data was provided. No validations of the models with respect to experimental NMR restraints is performed at this time.

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	B	1.28±0.02	2±0/281 (0.7±0.1%)	2.45±0.01	25±1/433 (5.8±0.3%)
2	C	1.23±0.01	1±0/263 (0.3±0.2%)	2.26±0.01	17±1/403 (4.3±0.3%)
3	A	1.02±0.01	0±0/1325 (0.0±0.0%)	0.80±0.01	0±0/1794 (0.0±0.0%)
All	All	1.10	51/33642 (0.2%)	1.49	765/47340 (1.6%)

All 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	B	325	DT	C5-C7	5.54	1.53	1.50	12	17
2	C	345	DT	C5-C7	5.49	1.53	1.50	1	15
1	B	324	DT	C5-C7	5.48	1.53	1.50	3	18
2	C	343	DT	C5-C7	5.11	1.53	1.50	13	1

5 of 59 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	B	326	DG	N7-C8-N9	9.71	117.95	113.10	8	18
1	B	316	DG	N7-C8-N9	9.38	117.79	113.10	16	18
2	C	351	DG	N7-C8-N9	9.23	117.71	113.10	5	18
1	B	319	DG	N7-C8-N9	9.15	117.68	113.10	17	18
1	B	318	DG	N7-C8-N9	8.94	117.57	113.10	11	18

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	B	249	136	136	15±3
2	C	237	139	139	21±3
3	A	1300	1332	1332	127±13
All	All	32148	28926	28926	2773

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 45.

5 of 1007 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
3:A:61:LEU:HD13	3:A:153:VAL:HG22	1.11	1.11	14	18
3:A:15:LEU:HD13	3:A:153:VAL:HG21	1.09	1.23	10	18
3:A:37:ALA:HB2	3:A:123:ILE:HG22	1.05	1.17	14	6
3:A:79:TYR:CE2	3:A:126:LEU:HD21	1.04	1.87	13	7
3:A:47:ILE:HD12	3:A:117:VAL:HG13	1.00	1.34	17	12

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
3	A	164/178 (92%)	114±5 (70±3%)	37±5 (22±3%)	13±3 (8±2%)	2 14
All	All	2952/3204 (92%)	2054 (70%)	659 (22%)	239 (8%)	2 14

5 of 49 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
3	A	99	SER	18

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Mol	Chain	Res	Type	Models (Total)
3	A	110	PRO	18
3	A	77	ALA	16
3	A	18	GLU	15
3	A	52	GLY	14

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
3	A	145/156 (93%)	101±7 (70±5%)	44±7 (30±5%)	2 17
All	All	2610/2808 (93%)	1822 (70%)	788 (30%)	2 17

5 of 113 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)
3	A	65	ILE	18
3	A	15	LEU	16
3	A	161	ARG	16
3	A	62	GLN	15
3	A	67	THR	15

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 carbohydrates 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