



wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 2L2I
Title : NMR Structure of the complex between the Tfb1 subunit of TFIID and the activation domain of EKLF
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Deposited on : 2010-08-18

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 ⓘ 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
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : trunk28760
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

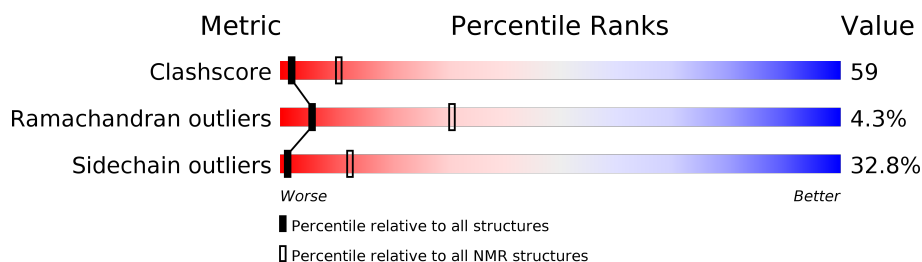
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment 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	125131	11601
Ramachandran outliers	121729	10391
Sidechain outliers	121581	10367

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	115	
2	B	42	

2 Ensemble composition and analysis

This entry contains 20 models. Model 14 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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:3-A:32, A:36-A:64, A:86-A:112, B:70-B:75 (92)	0.28	14

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 6 clusters and 7 single-model clusters were found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2207 atoms, of which 1097 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called RNA polymerase II transcription factor B subunit 1.

Mol	Chain	Residues	Atoms						Trace
1	A	115	Total	C	H	N	O	S	0
			1821	559	919	161	177	5	

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	PRO	-	EXPRESSION TAG	UNP P32776

- Molecule 2 is a protein called Krueppel-like factor 1.

Mol	Chain	Residues	Atoms					Trace
2	B	27	Total	C	H	N	O	0
			386	127	178	32	49	

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	49	GLY	-	EXPRESSION TAG	UNP Q13351
B	50	SER	-	EXPRESSION TAG	UNP Q13351

GLY	SER	LEU	HIS	VAL	LYS	SER	GLU	ASP	GLN	P59	G60	E61	E62	E63	D64	D65	E66	R67	G68	A69	D70	A71	T72	W73	D74	L75	D76	L77	L78	L79	T80	N81	F82	S83	G84	P85	GLU	PRO	GLY	GLY	ALA
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5 Refinement protocol and experimental data overview

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

Of the 100 calculated structures, 20 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
CNS	structure solution	
CNS	refinement	

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](#)

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

There are no bond-length outliers.

There are no bond-angle outliers.

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	678	708	705	77±7
2	B	50	41	41	20±3
All	All	14560	14980	14920	1725

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 59.

5 of 486 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:48:LEU:HD22	1:A:108:ILE:HD12	1.04	1.11	8	1
1:A:42:LEU:HD21	1:A:104:LEU:HD23	1.03	1.31	17	3
1:A:27:LEU:HD13	1:A:104:LEU:HD21	1.00	1.25	17	2
1:A:27:LEU:HD23	1:A:104:LEU:HD21	0.96	1.33	5	4
1:A:51:THR:HG22	2:B:72:THR:HG23	0.95	1.33	2	8

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	86/115 (75%)	77±2 (89±2%)	6±2 (8±2%)	3±1 (3±1%)	7	38
2	B	6/42 (14%)	2±1 (28±15%)	3±1 (54±16%)	1±1 (18±15%)	0	3
All	All	1840/3140 (59%)	1567 (85%)	194 (11%)	79 (4%)	6	30

5 of 16 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	25	ALA	20
1	A	53	ALA	19
2	B	74	ASP	8
2	B	71	ALA	7
1	A	57	LYS	6

6.3.2 Protein sidechains ⓘ

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	77/103 (75%)	53±3 (69±4%)	24±3 (31±4%)	1	14
2	B	5/33 (15%)	2±1 (46±18%)	3±1 (54±18%)	0	1
All	All	1640/2720 (60%)	1102 (67%)	538 (33%)	1	12

5 of 61 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	23	SER	20
1	A	22	VAL	20
1	A	107	ILE	20
1	A	89	PHE	20
1	A	58	MET	20

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

No chemical shift data were provided