



wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 2L5X
Title : Solution structure of IL1A-S100A13 complex
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Deposited on : 2010-11-09

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
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

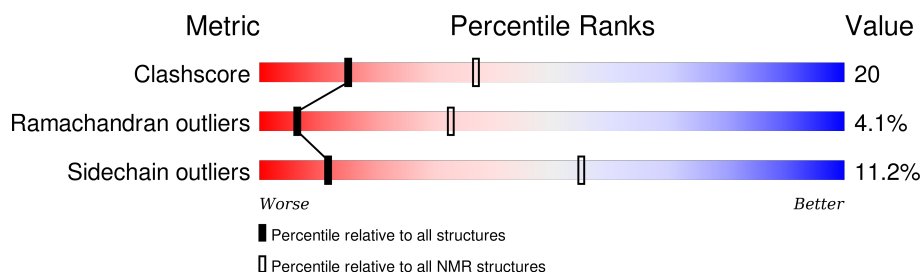
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	151	56% 40% .
1	D	151	58% 38% .
2	B	98	55% 36% 8% .
2	C	98	57% 35% 7% .

2 Ensemble composition and analysis ⓘ

This entry contains 20 models. Model 6 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:9-A:159, B:1-B:98, C:1-C:98, D:9-D:159 (498)	0.35	6

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 3 clusters and 1 single-model cluster was found.

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

3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 8128 atoms, of which 4084 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Interleukin-1 alpha.

Mol	Chain	Residues	Atoms						Trace
1	A	151	Total	C	H	N	O	S	0
			2423	782	1208	197	232	4	
1	D	151	Total	C	H	N	O	S	0
			2423	782	1208	197	232	4	

- Molecule 2 is a protein called Protein S100-A13.

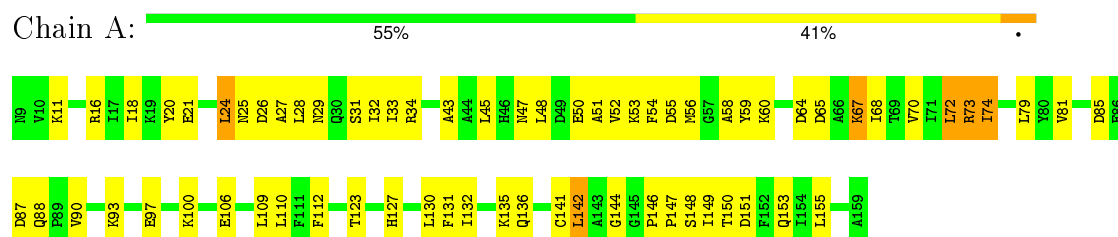
Mol	Chain	Residues	Atoms						Trace
2	B	98	Total	C	H	N	O	S	0
			1641	512	834	136	157	2	
2	C	98	Total	C	H	N	O	S	0
			1641	512	834	136	157	2	



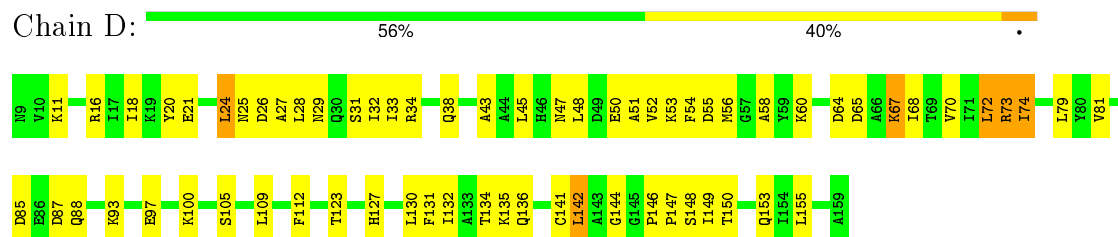
4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 6. Colouring as in section 4.1 above.

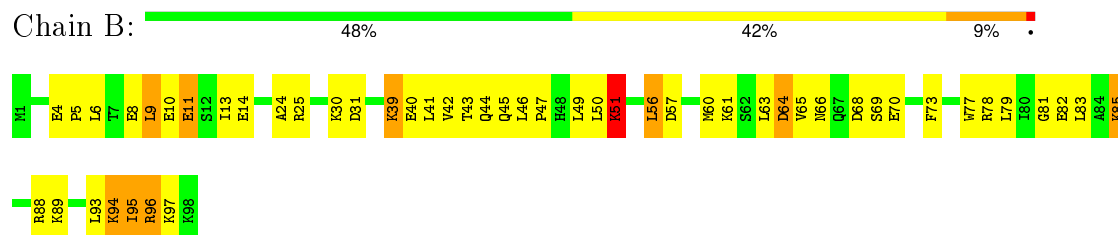
- Molecule 1: Interleukin-1 alpha



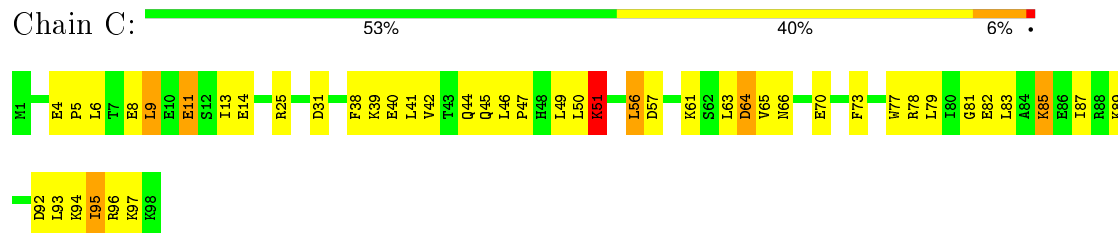
- Molecule 1: Interleukin-1 alpha



- Molecule 2: Protein S100-A13



- Molecule 2: Protein S100-A13



5 Refinement protocol and experimental data overview ⓘ

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

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
ARIA	structure solution	1.2 & 2.1
ARIA	refinement	1.2 & 2.1

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

6.1 Standard geometry

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

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	1215	1208	1202	49±2
1	D	1215	1208	1202	48±4
2	B	807	834	833	34±2
2	C	807	834	833	32±3
All	All	80880	81680	81400	3165

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 20.

5 of 270 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:63:LEU:HD21	2:B:79:LEU:HB2	0.84	1.49	14	20
2:C:63:LEU:HD21	2:C:79:LEU:HB2	0.84	1.50	18	20
1:D:65:ASP:HB3	1:D:112:PHE:CE2	0.81	2.10	16	20
1:D:60:LYS:HE2	1:D:68:ILE:HD12	0.81	1.53	8	20
1:A:60:LYS:HE2	1:A:68:ILE:HD12	0.80	1.53	10	20

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	149/151 (99%)	121±1 (81±1%)	23±1 (16±1%)	5±0 (3±0%)	9	41
1	D	149/151 (99%)	122±1 (82±1%)	23±1 (15±1%)	5±0 (3±0%)	9	41
2	B	96/98 (98%)	74±1 (77±1%)	17±1 (18±1%)	5±1 (5±1%)	4	24
2	C	96/98 (98%)	73±1 (76±1%)	17±1 (18±1%)	6±1 (6±1%)	4	21
All	All	9800/9960 (98%)	7796 (80%)	1598 (16%)	406 (4%)	6	33

5 of 33 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	147	PRO	20
1	D	18	ILE	20
1	A	64	ASP	20
2	C	73	PHE	20
1	D	147	PRO	20

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	133/133 (100%)	120±1 (90±1%)	13±1 (10±1%)	15	60
1	D	133/133 (100%)	120±1 (90±1%)	13±1 (10±1%)	15	59
2	B	91/91 (100%)	78±1 (86±1%)	13±1 (14±1%)	8	49
2	C	91/91 (100%)	79±1 (87±1%)	12±1 (13±1%)	9	50
All	All	8960/8960 (100%)	7952 (89%)	1008 (11%)	12	55

5 of 75 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	24	LEU	20
2	C	56	LEU	20
1	D	16	ARG	20
2	C	39	LYS	20
1	A	153	GLN	20

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

No chemical shift data were provided