wwPDB NMR Structure Validation Summary Report (i)

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PDB ID	:	1YUU
Title	:	Solution structure of Calcium-S100A13
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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
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:

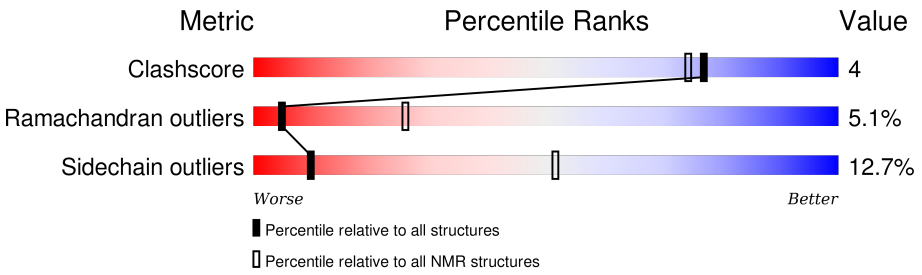
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	rb-20027457
Ideal 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 is 86%.

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	A	98	70% 17% 12%
1	B	98	64% 20% • 14%

2 Ensemble composition and analysis (i)

This entry contains 25 models. Model 7 is the overall representative, medoid model (most similar to other models). The authors have identified model 19 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:4-A:89, B:6-B:89 (170)	0.54	7

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 5 single-model clusters were found.

Cluster number	Models
1	1, 2, 3, 4, 5, 7, 8, 10, 11, 12, 13, 15, 16, 17, 20
2	14, 23, 24
3	19, 25
Single-model clusters	6; 9; 18; 21; 22

3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3284 atoms, of which 1666 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called S100 calcium-binding protein A13.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	98	1640	512	833	136	157	2	0
1	B	98	1640	512	833	136	157	2	0

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	
			Total	Ca
2	B	2	2	2
2	A	2	2	2

4 Residue-property plots (i)

4.1 Average score per residue in the NMR ensemble

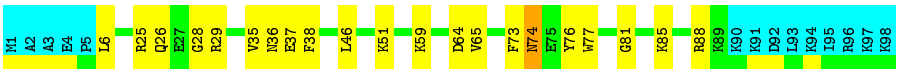
These plots are provided for all protein, RNA and DNA chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain 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: S100 calcium-binding protein A13

Chain A: 70% 17% 12%

- Molecule 1: S100 calcium-binding protein A13

Chain B: 64% 20% • 14%



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

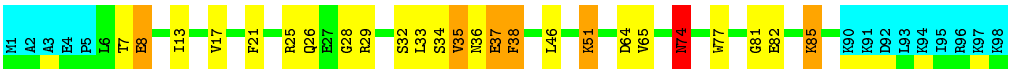
The representative model is number 7. Colouring as in section 4.1 above.

- Molecule 1: S100 calcium-binding protein A13

Chain A: 69% 16% • 12%

- Molecule 1: S100 calcium-binding protein A13

Chain B: 60% 18% 6% • 14%



5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *Restrained energy minimization.*

Of the 300 calculated structures, 25 were deposited, based on the following criterion: *target function.*

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

Software name	Classification	Version
DYANA	structure solution	1.5
AMBER	refinement	6

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)	BMRB entry 6531
Number of chemical shift lists	1
Total number of shifts	2274
Number of shifts mapped to atoms	2274
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	86%

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)

Bond lengths and bond angles in the following residue types are not validated in this section: CA

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	0.57±0.01	0±0/719 (0.0±0.0%)	1.06±0.04	2±1/968 (0.2±0.1%)
1	B	0.57±0.01	0±0/702 (0.0±0.0%)	1.10±0.04	2±1/944 (0.2±0.1%)
All	All	0.57	0/35525 (0.0%)	1.08	103/47800 (0.2%)

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	0.2±0.4	2.1±1.3
1	B	0.1±0.3	7.8±2.3
All	All	7	248

There are no bond-length outliers.

5 of 27 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	88	ARG	NE-CZ-NH2	-8.49	116.06	120.30	8	2
1	A	77	TRP	CZ3-CH2-CZ2	-7.96	112.05	121.60	15	5
1	B	35	VAL	C-N-CA	7.85	141.32	121.70	4	1
1	A	88	ARG	NE-CZ-NH2	-7.69	116.45	120.30	4	3
1	A	29	ARG	NE-CZ-NH2	-7.45	116.58	120.30	16	1

5 of 6 unique chiral outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Models (Total)
1	A	88	ARG	CA	2
1	A	49	LEU	CA	1
1	B	6	LEU	CA	1
1	A	86	GLU	CA	1
1	B	18	THR	CA	1

5 of 40 unique planar outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Group	Models (Total)
1	B	46	LEU	Peptide	24
1	B	64	ASP	Peptide	22
1	B	73	PHE	Peptide,Sidechain	22
1	B	28	GLY	Peptide	18
1	B	74	ASN	Peptide	16

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	708	708	708	8±5
1	B	692	695	694	9±5
All	All	35100	35075	35065	258

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 4.

5 of 94 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:49:LEU:HD23	1:A:83:LEU:HG	0.79	1.54	1	1
1:A:77:TRP:CZ2	1:B:77:TRP:CZ2	0.68	2.81	13	13
1:A:77:TRP:CZ2	1:B:77:TRP:CE2	0.68	2.82	22	13
1:A:77:TRP:CE2	1:B:77:TRP:CZ2	0.66	2.84	22	15
1:B:63:LEU:HB3	1:B:71:LEU:HD11	0.65	1.69	3	1

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	86/98 (88%)	69±2 (80±3%)	14±2 (16±3%)	4±1 (5±1%)	5 29
1	B	84/98 (86%)	62±2 (74±3%)	17±2 (21±3%)	5±2 (6±2%)	4 23
All	All	4250/4900 (87%)	3266 (77%)	769 (18%)	215 (5%)	5 26

5 of 37 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	B	65	VAL	25
1	A	28	GLY	21
1	A	29	ARG	20
1	B	29	ARG	20
1	B	35	VAL	15

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	81/91 (89%)	71±3 (88±3%)	10±3 (12±3%)	10 52
1	B	79/91 (87%)	69±3 (87±4%)	10±3 (13±4%)	9 51
All	All	4000/4550 (88%)	3492 (87%)	508 (13%)	10 51

5 of 108 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	59	LYS	23
1	A	26	GLN	20

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Mol	Chain	Res	Type	Models (Total)
1	B	51	LYS	20
1	B	74	ASN	17
1	A	51	LYS	17

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)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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)

The completeness of assignment taking into account all chemical shift lists is 86% for the well-defined parts and 79% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 6531

Chemical shift list name: *assigned_ chem_ shift_ list_ 1*

7.1.1 Bookkeeping (i)

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	2274
Number of shifts mapped to atoms	2274
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	6

7.1.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	182	2.34 ± 0.10	Should be applied
$^{13}C_{\beta}$	176	3.23 ± 0.05	Should be applied
$^{13}C'$	174	2.57 ± 0.07	Should be applied
^{15}N	176	0.81 ± 0.25	Should be applied

7.1.3 Completeness of resonance assignments (i)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 86%, i.e. 1891 atoms were assigned a chemical shift out of a possible 2202. 20 out of 36 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	829/844 (98%)	333/337 (99%)	331/340 (97%)	165/167 (99%)
Sidechain	1048/1214 (86%)	626/706 (89%)	410/454 (90%)	12/54 (22%)

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	Total	^{1}H	^{13}C	^{15}N
Aromatic	14/144 (10%)	12/78 (15%)	0/62 (0%)	2/4 (50%)
Overall	1891/2202 (86%)	971/1121 (87%)	741/856 (87%)	179/225 (80%)

7.1.4 Statistically unusual chemical shifts (i)

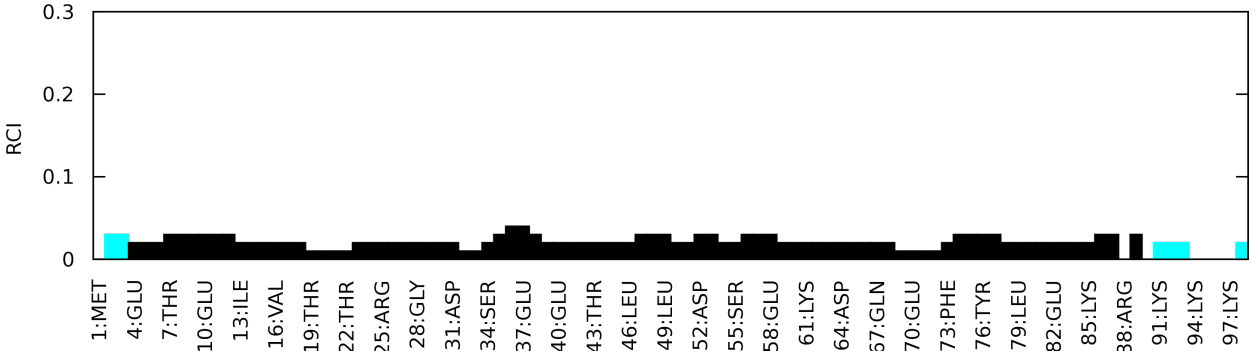
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, *ppm*	Expected range, *ppm*	Z-score
1	B	80	ILE	CG2	3.45	24.63 – 10.43	-9.9
1	A	80	ILE	CG2	3.45	24.63 – 10.43	-9.9
1	A	46	LEU	CG	35.20	32.55 – 21.05	7.3
1	B	46	LEU	CG	35.20	32.55 – 21.05	7.3
1	B	9	LEU	CD2	14.44	32.60 – 15.60	-5.7
1	A	9	LEU	CD2	14.44	32.60 – 15.60	-5.7

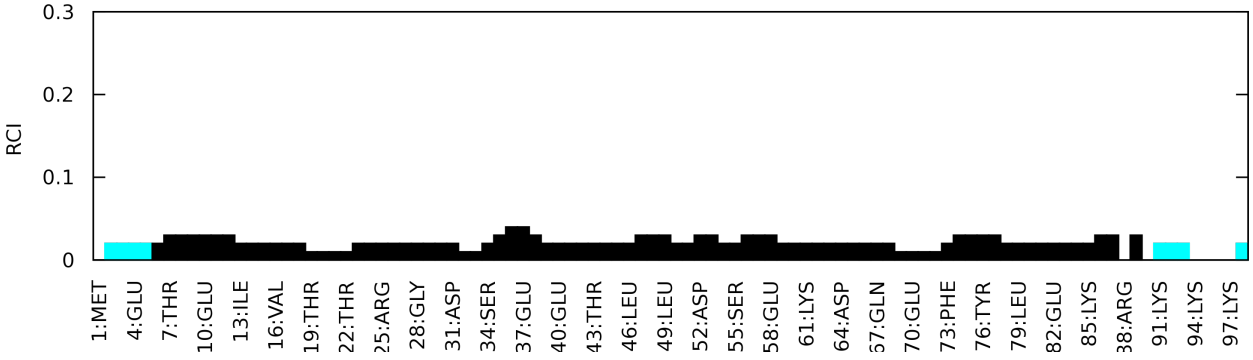
7.1.5 Random Coil Index (RCI) plots (i)

The images below report *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:



Random coil index (RCI) for chain B:


RCI
0.3
0.2
0.1
0
1:MET
4:GLU
7:THR
10:GLU
13:ILE
16:VAL
19:THR
22:THR
25:ARG
28:GLY
31:ASP
34:SER
37:GLU
40:GLU
43:THR
46:LEU
49:LEU
52:ASP
55:SER
58:GLU
61:LYS
64:ASP
57:GLN
70:GLU
73:PHE
76:TYR
79:LEU
82:GLU
85:LYS
38:ARG
91:LYS
94:LYS
97:LYS