Full wwPDB NMR Structure Validation Report (i)

Apr 26, 2016 – 04:18 PM BST

PDB ID	:	1PFD
Title	:	THE SOLUTION STRUCTURE OF HIGH PLANT PARSLEY [2FE-2S] FERREDOXIN, NMR, 18 STRUCTURES
Authors	:	Im, S.-C.; Liu, G.; Luchinat, C.; Sykes, A.G.; Bertini, I.
Deposited on	:	1998-05-05

This is a Full wwPDB NMR Structure Validation 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:

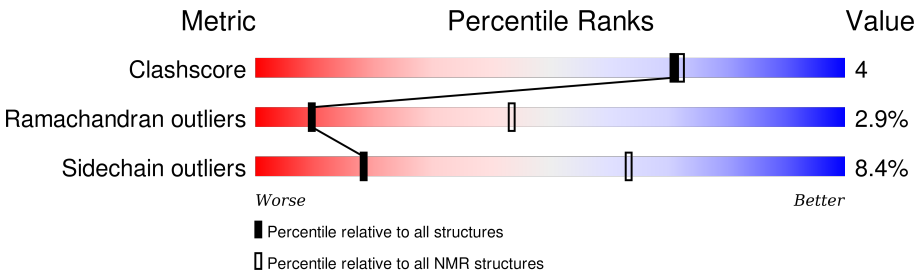
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	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 46%.

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	96	64% 15% 22%

2 Ensemble composition and analysis (i)

This entry contains 18 models. Model 5 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative.

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:35, A:48-A:59, A:64-A:74, A:79-A:96 (75)	0.21	5

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

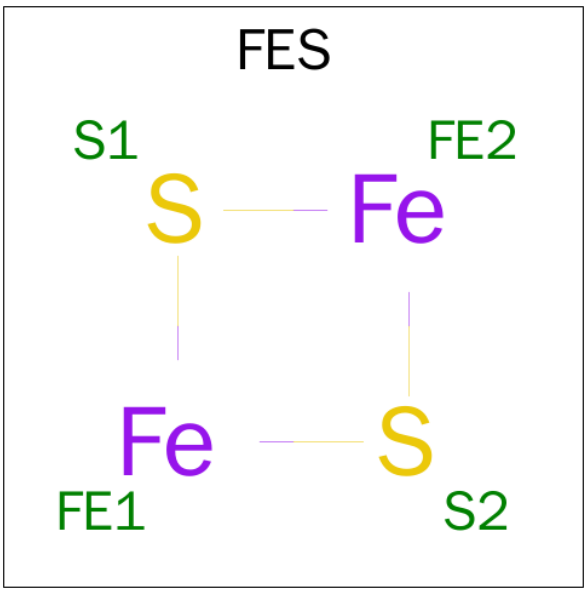
3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1404 atoms, of which 668 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called FERREDOXIN.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	96	1400	449	668	112	165	6	0

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms		
			Total	Fe	S
2	A	1	4	2	2

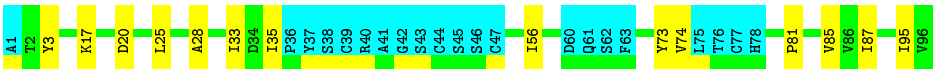
4 Residue-property plots (i)

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: FERREDOXIN

Chain A: 64% 15% 22%



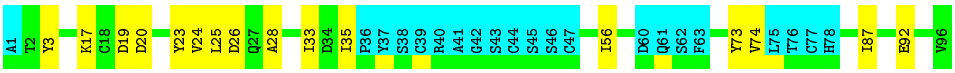
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: FERREDOXIN

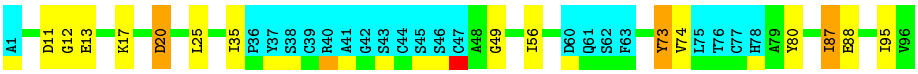
Chain A: 61% 17% 22%



4.2.2 Score per residue for model 2

- Molecule 1: FERREDOXIN

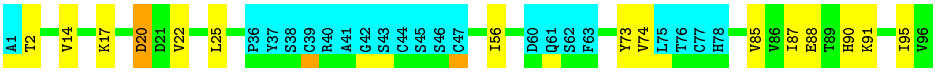
Chain A: 63% 13% • 22%



4.2.3 Score per residue for model 3

- Molecule 1: FERREDOXIN

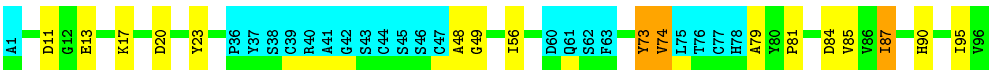
Chain A: 63% 15% • 22%



4.2.4 Score per residue for model 4

- Molecule 1: FERREDOXIN

Chain A: 60% 15% • 22%



4.2.5 Score per residue for model 5 (medoid)

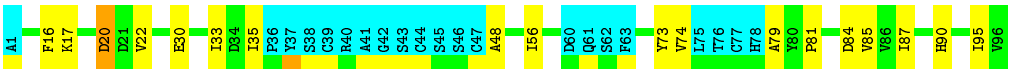
- Molecule 1: FERREDOXIN

Chain A: 64% 13% • 22%

4.2.6 Score per residue for model 6

- Molecule 1: FERREDOXIN

Chain A: 59% 18% • 22%



4.2.7 Score per residue for model 7

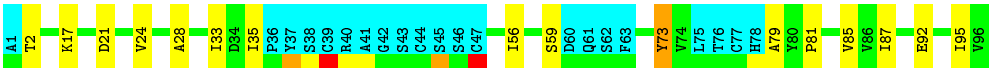
- Molecule 1: FERREDOXIN

Chain A: 59% 19% 22%

4.2.8 Score per residue for model 8

- Molecule 1: FERREDOXIN

Chain A: 61% 16% • 22%



4.2.9 Score per residue for model 9

- Molecule 1: FERREDOXIN

Chain A: 59% 16% • • 22%

4.2.10 Score per residue for model 10

- Molecule 1: FERREDOXIN

Chain A: 60% 17% • 22%

4.2.11 Score per residue for model 11

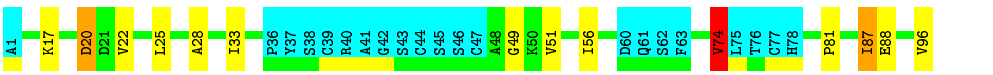
- Molecule 1: FERREDOXIN

Chain A: 60% 18% 22%

4.2.12 Score per residue for model 12

- Molecule 1: FERREDOXIN

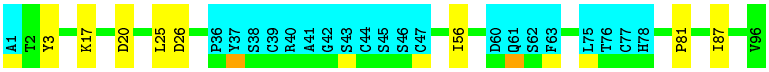
Chain A: 64% 11% • • 22%



4.2.13 Score per residue for model 13

- Molecule 1: FERREDOXIN

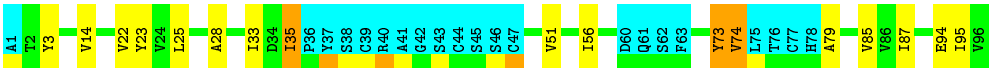
Chain A: 70% 8% 22%



4.2.14 Score per residue for model 14

- Molecule 1: FERREDOXIN

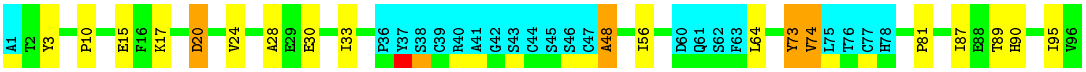
Chain A: 60% 15% • 22%



4.2.15 Score per residue for model 15

- Molecule 1: FERREDOXIN

Chain A: 58% 16% • 22%



4.2.16 Score per residue for model 16

- Molecule 1: FERREDOXIN

Chain A: 59% 17% • 22%

4.2.17 Score per residue for model 17

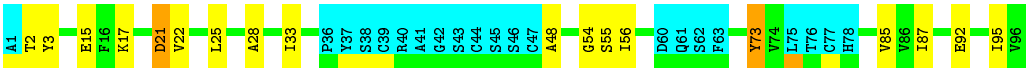
- Molecule 1: FERREDOXIN

Chain A: 64% 14% • 22%

4.2.18 Score per residue for model 18

- Molecule 1: FERREDOXIN

Chain A: 59% 17% • 22%



5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *DISTANCE GEOMETRY*.

Of the 50 calculated structures, 18 were deposited, based on the following criterion: *LEAST TARGET FUNCTION*.

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

Software name	Classification	Version
DYANA	refinement	1.4
DYANA	structure solution	
DYANAT1	structure solution	
AMBER	structure solution	

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 4181
Number of chemical shift lists	1
Total number of shifts	570
Number of shifts mapped to atoms	563
Number of unparsed shifts	0
Number of shifts with mapping errors	7
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	46%

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: FES

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	0.51±0.01	0±0/588 (0.0±0.0%)	0.98±0.06	0±0/799 (0.0±0.0%)
All	All	0.51	0/10584 (0.0%)	0.99	3/14382 (0.0%)

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	0.0±0.0	2.0±1.0
All	All	0	36

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	96	VAL	CA-C-O	-18.40	81.46	120.10	12	1
1	A	18	CYS	CA-CB-SG	-5.15	104.73	114.00	17	2

There are no chirality outliers.

All 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	A	73	TYR	Sidechain	14
1	A	3	TYR	Sidechain	9
1	A	23	TYR	Sidechain	7
1	A	20	ASP	Sidechain	4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group	Models (Total)
1	A	16	PHE	Sidechain	1
1	A	80	TYR	Sidechain	1

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	581	538	538	5±1
2	A	4	0	0	0±0
All	All	10530	9684	9684	89

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.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:56:ILE:HD11	1:A:74:VAL:HG21	0.98	1.31	7	11
1:A:56:ILE:HG22	1:A:81:PRO:HA	0.59	1.73	7	11
1:A:56:ILE:CD1	1:A:74:VAL:HG21	0.56	2.30	2	5
1:A:56:ILE:HG23	1:A:85:VAL:HG21	0.56	1.77	8	6
1:A:28:ALA:HB1	1:A:33:ILE:HB	0.55	1.78	16	9
1:A:25:LEU:HD12	1:A:26:ASP:N	0.55	2.17	13	1
1:A:49:GLY:HA3	1:A:87:ILE:HD11	0.51	1.81	12	3
1:A:10:PRO:HD3	1:A:90:HIS:CD2	0.49	2.43	15	2
1:A:56:ILE:CG2	1:A:85:VAL:HG21	0.49	2.38	18	2
1:A:56:ILE:HD11	1:A:74:VAL:CG2	0.49	2.35	5	2
1:A:25:LEU:HD13	1:A:26:ASP:N	0.48	2.24	5	1
1:A:73:TYR:CG	1:A:95:ILE:HG22	0.47	2.44	14	10
1:A:56:ILE:HD12	1:A:79:ALA:HB1	0.47	1.87	4	5
1:A:56:ILE:CG2	1:A:85:VAL:HG11	0.46	2.40	6	3
1:A:33:ILE:HG22	1:A:35:ILE:HG23	0.46	1.87	5	1
1:A:54:GLY:HA3	1:A:85:VAL:HG22	0.45	1.89	18	1
1:A:7:LEU:HD12	1:A:33:ILE:HD13	0.45	1.89	7	1
1:A:5:VAL:HG13	1:A:85:VAL:HG13	0.44	1.89	7	2
1:A:25:LEU:C	1:A:25:LEU:HD13	0.44	2.33	16	1
1:A:56:ILE:HD12	1:A:79:ALA:CB	0.43	2.44	11	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:51:VAL:HG23	1:A:74:VAL:HB	0.43	1.89	14	1
1:A:51:VAL:CG2	1:A:74:VAL:HB	0.43	2.43	12	1
1:A:87:ILE:HD13	1:A:88:GLU:N	0.42	2.30	12	2
1:A:14:VAL:HG23	1:A:15:GLU:H	0.42	1.74	5	1
1:A:16:PHE:HB3	1:A:33:ILE:HD11	0.41	1.91	6	1
1:A:20:ASP:N	1:A:20:ASP:OD1	0.41	2.53	9	1
1:A:25:LEU:HD23	1:A:25:LEU:C	0.41	2.36	9	1
1:A:20:ASP:OD1	1:A:20:ASP:N	0.41	2.54	16	1
1:A:88:GLU:OE1	1:A:91:LYS:NZ	0.41	2.52	3	1
1:A:56:ILE:HG21	1:A:85:VAL:HG11	0.41	1.93	6	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	74/96 (77%)	59±2 (80±2%)	13±1 (17±2%)	2±1 (3±2%)	9	43
All	All	1332/1728 (77%)	1066 (80%)	227 (17%)	39 (3%)	9	43

All 11 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	48	ALA	7
1	A	15	GLU	6
1	A	25	LEU	5
1	A	24	VAL	5
1	A	14	VAL	4
1	A	84	ASP	4
1	A	2	THR	3
1	A	13	GLU	2
1	A	92	GLU	1
1	A	74	VAL	1
1	A	12	GLY	1

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	66/84 (79%)	60±1 (92±2%)	6±1 (8±2%)	18	63
All	All	1188/1512 (79%)	1088 (92%)	100 (8%)	18	63

All 21 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	87	ILE	18
1	A	17	LYS	17
1	A	35	ILE	9
1	A	22	VAL	8
1	A	20	ASP	8
1	A	25	LEU	7
1	A	74	VAL	5
1	A	90	HIS	4
1	A	26	ASP	3
1	A	19	ASP	3
1	A	11	ASP	3
1	A	92	GLU	2
1	A	94	GLU	2
1	A	30	GLU	2
1	A	21	ASP	2
1	A	89	THR	2
1	A	27	GLN	1
1	A	31	GLU	1
1	A	59	SER	1
1	A	55	SER	1
1	A	4	ASN	1

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)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length is the number of standard deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
2	FES	A	97	1	0,4,4	0.00±0.00	-

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
2	FES	A	97	1	0,4,4	0.00±0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FES	A	97	1	-	0±0,0,4,4	0±0,1,1,1

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 46% for the well-defined parts and 43% for the entire structure.

7.1 Chemical shift list 1

File name: BMRB entry 4181

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	570
Number of shifts mapped to atoms	563
Number of unparsed shifts	0
Number of shifts with mapping errors	7
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- Residue not found in structure. All 7 occurences are reported below.

Chain	Res	Type	Atom	Shift Data		
				Value	Uncertainty	Ambiguity
A	64	LYS	HA	4.382	0.0	1
A	64	LYS	HB2	1.808	0.001	1
A	64	LYS	HD3	3.008	0.001	1
A	64	LYS	HB3	1.418	0.082	1
A	64	LYS	HD2	0.749	0.008	1
A	42	ASP	H	8.775	0.003	1
A	64	LYS	H	10.487	0.0	1

7.1.2 Chemical shift referencing (i)

No chemical shift referencing corrections were calculated (not enough data).

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 46%, i.e. 392 atoms were assigned a chemical shift out of a possible 846. 0 out of 13 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	145/371 (39%)	145/148 (98%)	0/150 (0%)	0/73 (0%)
Sidechain	225/427 (53%)	225/242 (93%)	0/177 (0%)	0/8 (0%)
Aromatic	22/48 (46%)	22/25 (88%)	0/22 (0%)	0/1 (0%)
Overall	392/846 (46%)	392/415 (94%)	0/349 (0%)	0/82 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 43%, i.e. 462 atoms were assigned a chemical shift out of a possible 1067. 0 out of 14 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	177/474 (37%)	177/189 (94%)	0/192 (0%)	0/93 (0%)
Sidechain	254/521 (49%)	254/300 (85%)	0/209 (0%)	0/12 (0%)
Aromatic	31/72 (43%)	31/38 (82%)	0/32 (0%)	0/2 (0%)
Overall	462/1067 (43%)	462/527 (88%)	0/433 (0%)	0/107 (0%)

7.1.4 Statistically unusual chemical shifts (i)

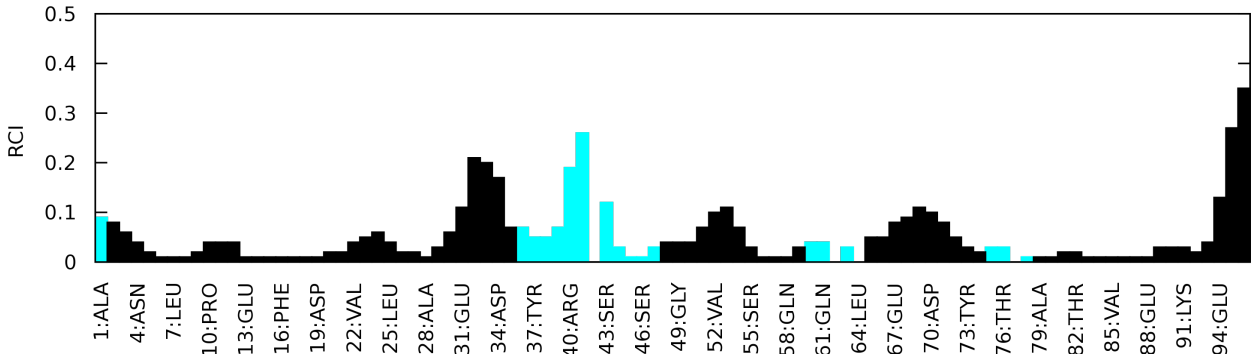
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, *ppm*	Expected range, *ppm*	Z-score
1	A	64	LYS	HD3	3.01	2.75 – 0.45	6.1

7.1.5 Random Coil Index (RCI) plots (i)

The image below reports *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:


RCI
0.5
0.4
0.3
0.2
0.1
0
1:ALA
4:ASN
7:LEU
10:PRO
13:GLU
16:PHE
19:ASP
22:VAL
25:LEU
28:ALA
31:GLU
34:ASP
37:TYR
40:ARG
43:SER
46:SER
49:GLY
52:VAL
55:SER
58:GLN
51:GLN
64:LEU
67:GLU
70:ASP
73:TYR
76:THR
79:ALA
82:THR
85:VAL
88:GLU
91:LYS
94:GLU