Full wwPDB NMR Structure Validation Report (i)

May 28, 2020 – 11:38 pm BST

PDB ID	:	2ME6
Title	:	NMR Structure of the homeodomain transcription factor Gbx1 from Homo sapiens in complex with the DNA sequence CGACTAATTAGTCG
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Deposited on	:	2013-09-23

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
https://www.wwpdb.org/validation/2017/NMRValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

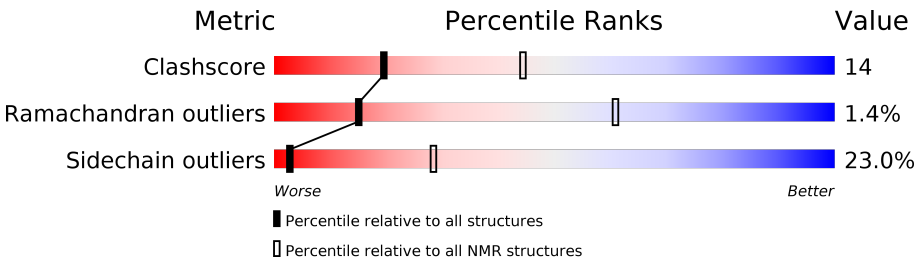
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	2.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 59%.

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	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

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	71	42% 32% 25%
2	B	14	29% 71%
2	C	14	29% 71%

2 Ensemble composition and analysis (i)

This entry contains 4 models. Model 4 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *closest to the average*.

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:11-A:63 (53)	0.65	4

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 1 clusters. No single-model clusters were found.

Cluster number	Models
1	1, 2, 3, 4

3 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2096 atoms, of which 937 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Homeobox protein GBX-1.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	71	1207	370	619	116	101	1	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	EXPRESSION TAG	UNP Q14549

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

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	P	
2	B	14	443	137	159	52	82	13	0
			Total	C	H	N	O	P	
2	C	14	446	137	159	52	84	14	0

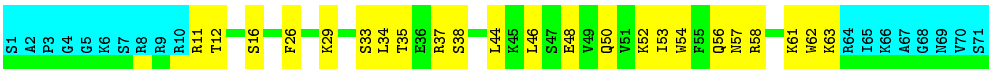
4 Residue-property plots (i)

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.

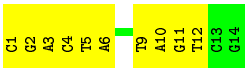
- Molecule 1: Homeobox protein GBX-1

Chain A:



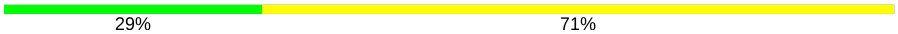
- Molecule 2: DNA (5'-D(*CP*GP*AP*CP*TP*AP*AP*TP*TP*AP*GP*TP*CP*G)-3')

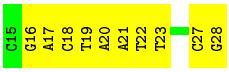
Chain B:



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

Chain C:





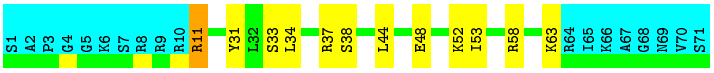
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: Homeobox protein GBX-1

Chain A:



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

Chain B:

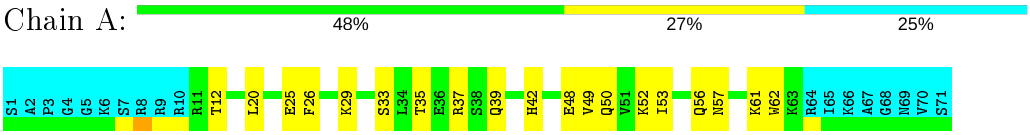
• Molecule 2: DNA (5'-D(*CP*GP*AP*CP*TP*AP*AP*TP*TP*AP*GP*TP*CP*G)-3')

Chain C:

4.2.2 Score per residue for model 2

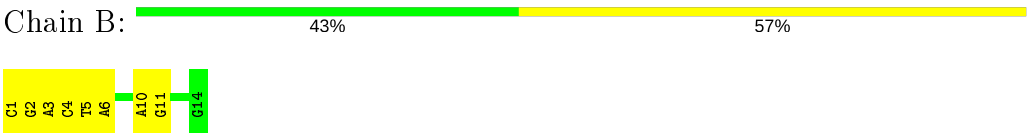
• Molecule 1: Homeobox protein GBX-1

Chain A:



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

Chain B:



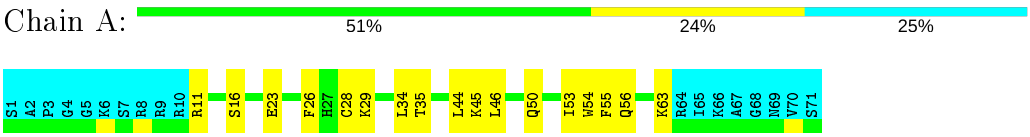
• Molecule 2: DNA (5'-D(*CP*GP*AP*CP*TP*AP*AP*TP*TP*AP*GP*TP*CP*G)-3')

Chain C:

4.2.3 Score per residue for model 3

• Molecule 1: Homeobox protein GBX-1

Chain A:



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

Chain B: 14% 86%

C1 G2 A3 C4 T5 A6 A7 T8 T9 A10 G11 T12 C13 G14

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

Chain C: 36% 64%

C15 G16 A17 C18 T19 A20 A21 T22 T23 G28

4.2.4 Score per residue for model 4 (medoid)

- Molecule 1: Homeobox protein GBX-1

Chain A: 49% 20% 6% 25%

S1 A2 P3 G4 G5 K6 S7 R8 R9 R10 R11 T12 S16 L34 T35 S38 L44 K45 L46 S47 Q50 I53 W54 F55 Q56 N57 R58 K61 W62 K63 R64 I65 K66 A67 G68 N69 V70 S71

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

Chain B: 29% 71%

C1 G2 A3 C4 T5 A6 T9 A10 G11 T12 C13 G14

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

Chain C: 43% 57%

C15 G16 A17 C18 T19 A20 A21 T22 G28

5 Refinement protocol and experimental data overview (i)

The models were refined using the following method: *molecular dynamics*.

Of the 200 calculated structures, 4 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
CNS	structure solution	
CNS	structure solution	
CNS	structure solution	
CNS	structure solution	
UNIO	structure solution	
ARIA	refinement	
HADDOCK	refinement	
3D DART	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)	input_cs.cif
Number of chemical shift lists	2
Total number of shifts	997
Number of shifts mapped to atoms	997
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	59%

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	455	472	470	10±3
2	B	284	159	160	10±3
2	C	287	159	159	16±4
All	All	4104	3160	3156	102

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 14.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:53:ILE:HG21	2:C:21:DA:H2'	0.84	1.49	2	1
1:A:53:ILE:HG21	2:C:22:DT:H72	0.80	1.53	3	2
1:A:53:ILE:HD13	2:C:22:DT:C7	0.78	2.07	3	1
2:B:1:DC:O2	2:C:28:DG:N2	0.76	2.18	2	4
1:A:53:ILE:HG23	2:C:22:DT:C7	0.76	2.11	2	2
2:C:22:DT:C6	2:C:23:DT:H72	0.71	2.20	1	1
2:C:17:DA:H2"	2:C:18:DC:C5	0.70	2.21	3	2
1:A:53:ILE:HD13	2:C:22:DT:H73	0.69	1.62	3	1
1:A:53:ILE:HG23	2:C:22:DT:H71	0.69	1.63	2	2
1:A:53:ILE:CG2	2:C:22:DT:H71	0.68	2.18	2	1
2:B:2:DG:H2"	2:B:3:DA:C8	0.68	2.24	1	4
2:B:5:DT:H2"	2:B:6:DA:C8	0.63	2.28	4	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:10:DA:H2"	2:B:11:DG:H5'	0.62	1.69	3	1
2:C:20:DA:H2"	2:C:21:DA:C8	0.62	2.30	4	1
1:A:34:LEU:O	1:A:35:THR:HB	0.60	1.94	4	1
2:C:17:DA:H2"	2:C:18:DC:C6	0.60	2.32	3	2
1:A:11:ARG:HA	2:C:20:DA:H5"	0.60	1.72	3	1
2:B:10:DA:H2"	2:B:11:DG:C8	0.59	2.32	2	1
2:C:18:DC:H2"	2:C:19:DT:C7	0.59	2.27	4	2
1:A:61:LYS:NZ	2:C:20:DA:H3'	0.59	2.12	2	1
2:C:18:DC:H2"	2:C:19:DT:C5	0.58	2.33	2	2
2:C:15:DC:H2"	2:C:16:DG:C8	0.58	2.33	4	1
1:A:63:LYS:NZ	2:B:6:DA:H3'	0.58	2.13	3	2
2:B:10:DA:C2'	2:B:11:DG:H5'	0.58	2.28	3	1
1:A:53:ILE:HD13	2:C:22:DT:H72	0.57	1.76	4	1
1:A:53:ILE:HG21	2:C:22:DT:C7	0.57	2.29	3	1
2:C:16:DG:H2"	2:C:17:DA:C8	0.56	2.35	3	2
2:C:24:DA:H2"	2:C:25:DG:C8	0.55	2.37	2	1
1:A:12:THR:H	2:C:20:DA:H5"	0.54	1.61	4	1
2:C:22:DT:C7	2:C:23:DT:H73	0.53	2.33	3	1
2:C:21:DA:H2"	2:C:22:DT:C7	0.53	2.33	3	1
2:B:2:DG:OP2	2:B:2:DG:H2'	0.52	2.04	3	2
2:C:18:DC:H2"	2:C:19:DT:H71	0.52	1.82	4	2
1:A:34:LEU:HD13	2:B:4:DC:H5'	0.52	1.80	4	1
2:B:4:DC:H2"	2:B:5:DT:C5	0.51	2.40	2	1
2:C:27:DC:H2"	2:C:28:DG:C8	0.51	2.40	2	1
1:A:53:ILE:CD1	2:C:22:DT:H73	0.51	2.36	3	1
2:B:3:DA:H2"	2:B:4:DC:C6	0.50	2.42	1	2
1:A:53:ILE:HD13	2:C:21:DA:C2'	0.49	2.37	2	1
2:B:9:DT:H2"	2:B:10:DA:C8	0.49	2.42	3	3
1:A:53:ILE:CG2	2:C:22:DT:C7	0.49	2.89	4	2
1:A:38:SER:OG	2:B:3:DA:H5"	0.49	2.08	1	1
2:B:10:DA:H2"	2:B:11:DG:C5'	0.49	2.38	3	1
2:C:21:DA:OP2	2:C:21:DA:H8	0.49	1.91	3	1
2:C:22:DT:H73	2:C:23:DT:H73	0.49	1.83	3	1
2:C:21:DA:C2'	2:C:22:DT:H72	0.49	2.38	3	1
1:A:53:ILE:CG2	2:C:20:DA:H2"	0.48	2.38	1	1
1:A:37:ARG:NH1	2:B:4:DC:H3'	0.48	2.23	1	1
1:A:56:GLN:HE22	2:C:22:DT:H73	0.48	1.69	2	1
2:B:11:DG:H2"	2:B:12:DT:C7	0.47	2.39	4	1
2:B:2:DG:H8	2:B:2:DG:OP2	0.47	1.92	4	1
1:A:53:ILE:CD1	2:C:22:DT:C7	0.47	2.88	3	1
1:A:48:GLU:O	1:A:52:LYS:HG3	0.47	2.10	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:10:DA:C1'	2:B:11:DG:H5'	0.46	2.41	3	1
2:B:3:DA:H8	2:B:3:DA:OP2	0.46	1.94	1	1
2:B:12:DT:H2"	2:B:13:DC:C6	0.46	2.46	3	1
2:B:10:DA:H1'	2:B:11:DG:H5'	0.46	1.86	3	1
1:A:53:ILE:HD13	2:C:21:DA:H2'	0.46	1.88	4	1
1:A:49:VAL:O	1:A:53:ILE:HG13	0.46	2.11	2	1
2:C:20:DA:H1'	2:C:21:DA:C8	0.46	2.46	2	1
1:A:53:ILE:CG2	2:C:22:DT:H72	0.45	2.41	4	1
2:B:4:DC:H2"	2:B:5:DT:C6	0.44	2.48	2	1
1:A:12:THR:N	2:C:20:DA:H5"	0.43	2.27	4	1
2:C:28:DG:OP2	2:C:28:DG:H8	0.43	1.96	2	1
1:A:11:ARG:N	2:C:20:DA:H5'	0.43	2.28	4	1
2:B:11:DG:H8	2:B:11:DG:OP2	0.43	1.95	4	1
2:C:21:DA:H2"	2:C:22:DT:C6	0.43	2.49	4	1
1:A:56:GLN:NE2	2:C:22:DT:H73	0.43	2.28	2	1
2:C:26:DT:H2"	2:C:27:DC:C5	0.42	2.48	1	1
1:A:26:PHE:HA	1:A:29:LYS:O	0.42	2.14	2	1
1:A:25:GLU:H	1:A:25:GLU:CD	0.42	2.17	2	1
2:B:3:DA:H2"	2:B:4:DC:C5	0.41	2.51	4	1
1:A:61:LYS:HZ2	2:C:20:DA:H3'	0.41	1.75	2	1
2:C:19:DT:H2'	2:C:19:DT:OP2	0.41	2.15	4	1
1:A:57:ASN:OD1	2:C:20:DA:H8	0.41	1.98	4	1
1:A:37:ARG:NH1	2:B:3:DA:H3'	0.41	2.31	2	1
2:B:8:DT:H2"	2:B:9:DT:C7	0.41	2.46	3	1
1:A:53:ILE:CD1	2:C:22:DT:H72	0.40	2.46	1	1
1:A:53:ILE:CD1	2:C:21:DA:H2'	0.40	2.46	1	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	53/71 (75%)	49±1 (92±2%)	4±1 (7±2%)	1±0 (1±1%)	15	61
All	All	212/284 (75%)	194 (92%)	15 (7%)	3 (1%)	15	61

All 3 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	34	LEU	1
1	A	11	ARG	1
1	A	35	THR	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	49/62 (79%)	38±3 (77±5%)	11±3 (23±5%)	3 28
All	All	196/248 (79%)	151 (77%)	45 (23%)	3 28

All 31 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	50	GLN	3
1	A	44	LEU	3
1	A	16	SER	2
1	A	35	THR	2
1	A	62	TRP	2
1	A	58	ARG	2
1	A	54	TRP	2
1	A	56	GLN	2
1	A	33	SER	2
1	A	11	ARG	2
1	A	63	LYS	2
1	A	46	LEU	2
1	A	29	LYS	1
1	A	31	TYR	1
1	A	12	THR	1
1	A	28	CYS	1
1	A	39	GLN	1
1	A	61	LYS	1
1	A	38	SER	1
1	A	55	PHE	1
1	A	47	SER	1

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Mol	Chain	Res	Type	Models (Total)
1	A	42	HIS	1
1	A	53	ILE	1
1	A	23	GLU	1
1	A	45	LYS	1
1	A	57	ASN	1
1	A	20	LEU	1
1	A	34	LEU	1
1	A	52	LYS	1
1	A	48	GLU	1
1	A	26	PHE	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)

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)

The completeness of assignment taking into account all chemical shift lists is 59% for the well-defined parts and 60% for the entire structure.

7.1 Chemical shift list 1

File name: input_cs.cif

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	816
Number of shifts mapped to atoms	816
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	15

7.1.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	68	1.90 ± 0.18	Should be applied
$^{13}C_{\beta}$	65	2.48 ± 0.15	Should be applied
$^{13}C'$	0	—	None (insufficient data)
^{15}N	67	0.35 ± 0.41	None needed (< 0.5 ppm)

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. 589 atoms were assigned a chemical shift out of a possible 1294. 9 out of 9 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	210/265 (79%)	105/106 (99%)	52/106 (49%)	53/53 (100%)
Sidechain	316/404 (78%)	196/239 (82%)	115/141 (82%)	5/24 (21%)

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	Total	^{1}H	^{13}C	^{15}N
Aromatic	63/75 (84%)	32/39 (82%)	29/30 (97%)	2/6 (33%)
Overall	589/1294 (46%)	333/710 (47%)	196/467 (42%)	60/117 (51%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 48%, i.e. 738 atoms were assigned a chemical shift out of a possible 1524. 10 out of 10 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	269/353 (76%)	134/141 (95%)	68/142 (48%)	67/70 (96%)
Sidechain	406/546 (74%)	253/325 (78%)	147/182 (81%)	6/39 (15%)
Aromatic	63/75 (84%)	32/39 (82%)	29/30 (97%)	2/6 (33%)
Overall	738/1524 (48%)	419/831 (50%)	244/544 (45%)	75/149 (50%)

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.

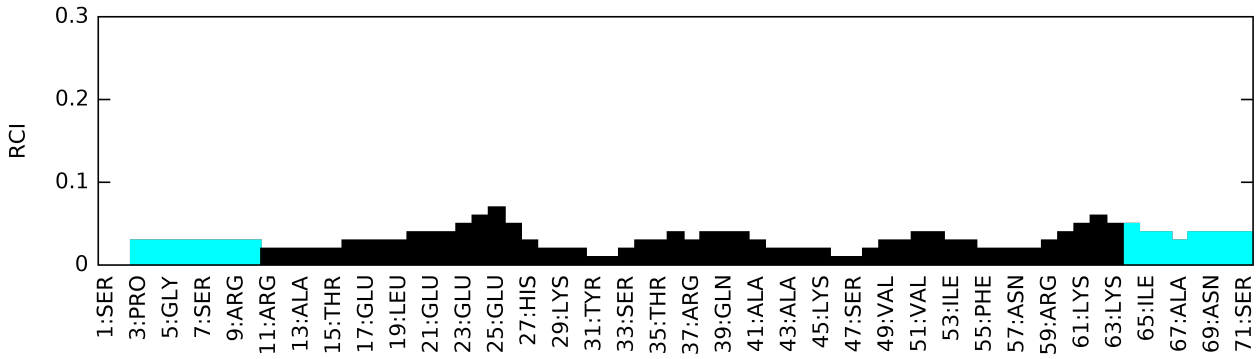
Mol	Chain	Res	Type	Atom	Shift, *ppm*	Expected range, *ppm*	Z-score
1	A	58	ARG	HG2	-0.69	2.92 – 0.22	-8.4
1	A	58	ARG	HB2	-0.46	3.15 – 0.45	-8.4
1	A	22	LEU	HB2	-1.02	3.32 – -0.08	-7.8
1	A	30	LYS	HE2	1.58	3.87 – 1.97	-7.0
1	A	33	SER	HB3	1.92	5.25 – 2.45	-6.9
1	A	38	SER	HB2	2.12	5.18 – 2.58	-6.8
1	A	33	SER	HB2	2.29	5.18 – 2.58	-6.1
1	A	58	ARG	HG3	-0.20	3.00 – 0.10	-6.0
1	A	30	LYS	HE3	1.82	3.86 – 1.96	-5.7
1	A	30	LYS	HD2	0.30	2.76 – 0.46	-5.7
1	A	22	LEU	HD23	-0.78	2.14 – -0.66	-5.4
1	A	22	LEU	HD22	-0.78	2.14 – -0.66	-5.4
1	A	22	LEU	HD21	-0.78	2.14 – -0.66	-5.4
1	A	37	ARG	HG2	0.12	2.92 – 0.22	-5.4
1	A	58	ARG	CG	21.10	33.23 – 21.23	-5.1

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:



7.2 Chemical shift list 2

File name: input_cs.cif

Chemical shift list name: *assigned_chem_shift_list_2*

7.2.1 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	181
Number of shifts mapped to atoms	181
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.2.2 Chemical shift referencing (i)

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

7.2.3 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 13%, i.e. 173 atoms were assigned a chemical shift out of a possible 1294. 0 out of 9 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	0/265 (0%)	0/106 (0%)	0/106 (0%)	0/53 (0%)
Sidechain	0/404 (0%)	0/239 (0%)	0/141 (0%)	0/24 (0%)
Aromatic	0/75 (0%)	0/39 (0%)	0/30 (0%)	0/6 (0%)
Overall	173/1294 (13%)	173/710 (24%)	0/467 (0%)	0/117 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 11%, i.e. 173 atoms were assigned a chemical shift out of a possible 1524. 0 out of 10 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	0/353 (0%)	0/141 (0%)	0/142 (0%)	0/70 (0%)
Sidechain	0/546 (0%)	0/325 (0%)	0/182 (0%)	0/39 (0%)
Aromatic	0/75 (0%)	0/39 (0%)	0/30 (0%)	0/6 (0%)
Overall	173/1524 (11%)	173/831 (21%)	0/544 (0%)	0/149 (0%)

7.2.4 Statistically unusual chemical shifts (i)

There are no statistically unusual chemical shifts.

7.2.5 Random Coil Index (RCI) plots (i)

No *random coil index* (RCI) plot could be generated from the current chemical shift list (assigned_chem_shift_list_2). RCI is only applicable to proteins.