Full wwPDB NMR Structure Validation Report (i)

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PDB ID	:	2MPB
Title	:	NMR structure of BA42 protein from the psychrophilic bacteria Bizionia argentinensis sp. nov
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Deposited on	:	2014-05-14

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:

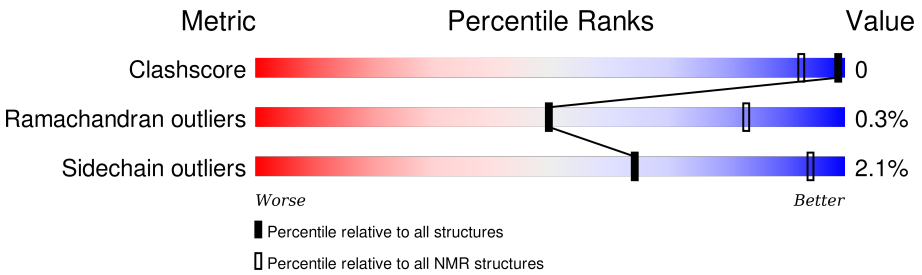
Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	rb-20027457
Ideal 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 90%.

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	145	90% · 7%

2 Ensemble composition and analysis (i)

This entry contains 20 models. Model 11 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:8-A:135, A:139-A:145 (135)	0.21	11

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 3 single-model clusters were found.

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

3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 2304 atoms, of which 1144 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called BA42.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	145	2304	734	1144	191	231	4	0

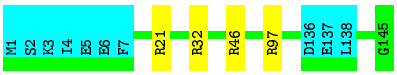
4 Residue-property plots (i)

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: BA42

Chain A:



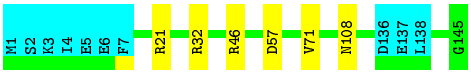
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: BA42

Chain A:



4.2.2 Score per residue for model 2

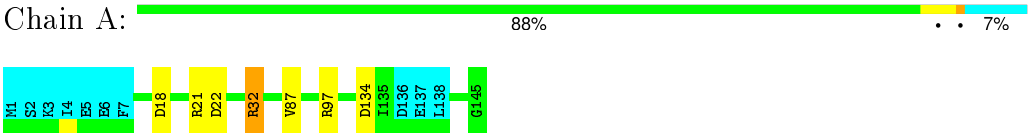
- Molecule 1: BA42

Chain A:

4.2.3 Score per residue for model 3

- Molecule 1: BA42

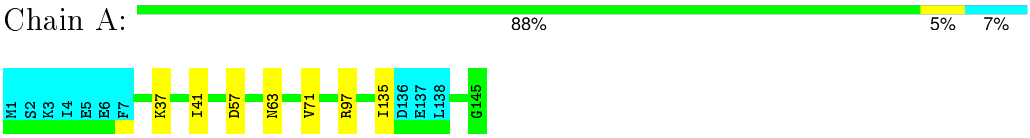
Chain A:



4.2.4 Score per residue for model 4

- Molecule 1: BA42

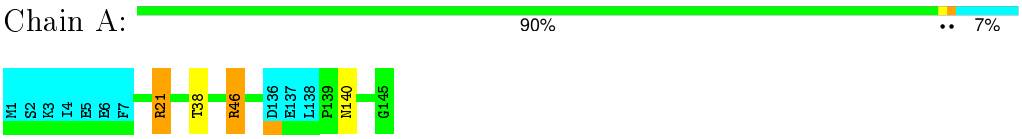
Chain A:



4.2.5 Score per residue for model 5

- Molecule 1: BA42

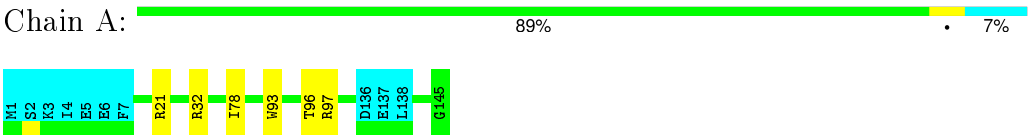
Chain A:



4.2.6 Score per residue for model 6

- Molecule 1: BA42

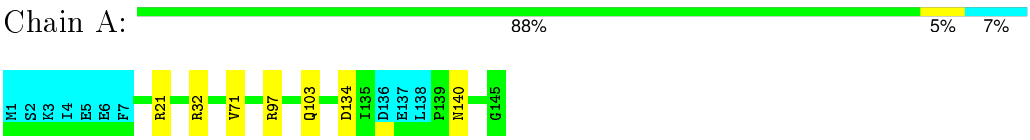
Chain A:



4.2.7 Score per residue for model 7

- Molecule 1: BA42

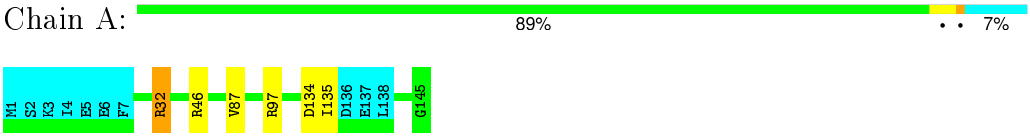
Chain A:



4.2.8 Score per residue for model 8

- Molecule 1: BA42

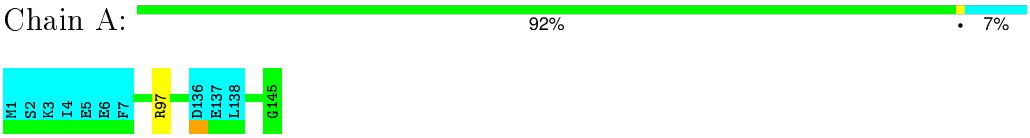
Chain A:



4.2.9 Score per residue for model 9

- Molecule 1: BA42

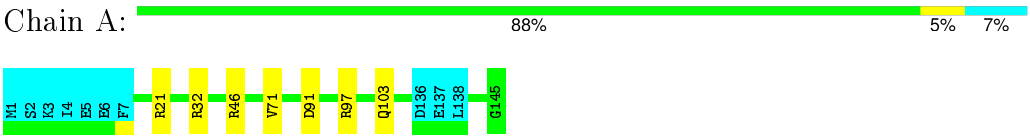
Chain A:



4.2.10 Score per residue for model 10

- Molecule 1: BA42

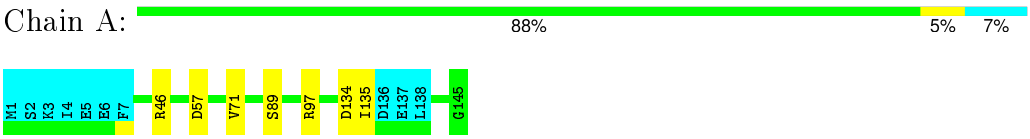
Chain A:



4.2.11 Score per residue for model 11 (medoid)

- Molecule 1: BA42

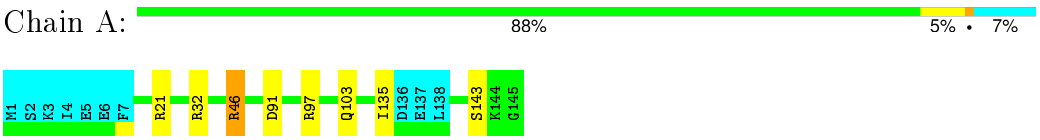
Chain A:



4.2.12 Score per residue for model 12

- Molecule 1: BA42

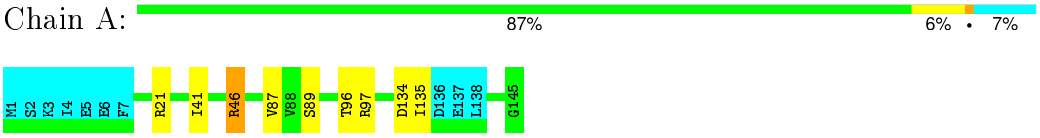
Chain A:



4.2.13 Score per residue for model 13

- Molecule 1: BA42

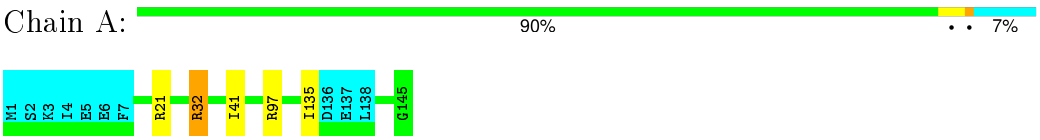
Chain A:



4.2.14 Score per residue for model 14

- Molecule 1: BA42

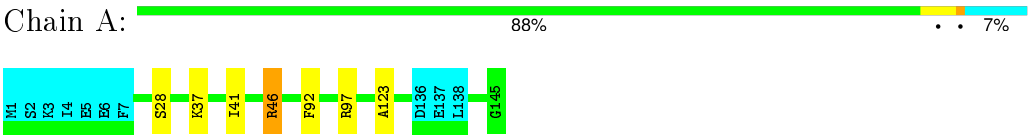
Chain A:



4.2.15 Score per residue for model 15

- Molecule 1: BA42

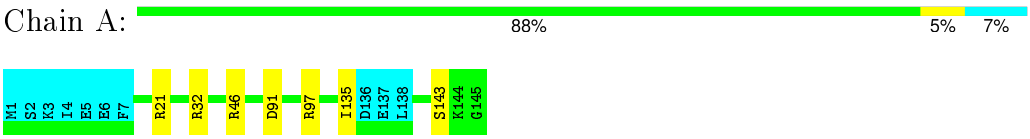
Chain A:



4.2.16 Score per residue for model 16

- Molecule 1: BA42

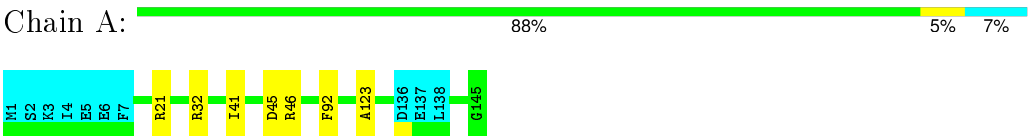
Chain A:



4.2.17 Score per residue for model 17

- Molecule 1: BA42

Chain A:



4.2.18 Score per residue for model 18

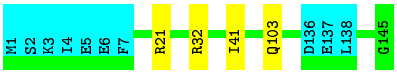
- Molecule 1: BA42

Chain A:

4.2.19 Score per residue for model 19

- Molecule 1: BA42

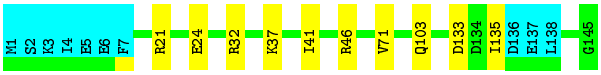
Chain A:



4.2.20 Score per residue for model 20

- Molecule 1: BA42

Chain A:



5 Refinement protocol and experimental data overview (i)

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

Of the 1000 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
X-PLOR NIH	structure solution	
X-PLOR NIH	refinement	

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)	2mpb_cs.str
Number of chemical shift lists	1
Total number of shifts	1770
Number of shifts mapped to atoms	1770
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	90%

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)

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	0.66±0.01	0±0/1092 (0.0±0.0%)	0.95±0.02	3±1/1469 (0.2±0.1%)
All	All	0.66	0/21840 (0.0%)	0.95	61/29380 (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	0.0±0.0	0.7±0.7
All	All	0	13

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	46	ARG	NE-CZ-NH1	9.83	125.21	120.30	11	12
1	A	97	ARG	NE-CZ-NH1	9.32	124.96	120.30	8	14
1	A	46	ARG	NE-CZ-NH2	-9.01	115.79	120.30	16	2
1	A	21	ARG	NE-CZ-NH1	8.37	124.48	120.30	10	14
1	A	32	ARG	NE-CZ-NH1	8.32	124.46	120.30	7	10
1	A	97	ARG	NE-CZ-NH2	-6.47	117.06	120.30	7	3
1	A	21	ARG	NE-CZ-NH2	-6.04	117.28	120.30	2	2
1	A	32	ARG	NE-CZ-NH2	-5.50	117.55	120.30	14	4

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	32	ARG	Sidechain	5
1	A	46	ARG	Sidechain	4
1	A	21	ARG	Sidechain	2
1	A	97	ARG	Sidechain	2

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	1075	1062	1062	0±1
All	All	21500	21240	21240	6

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 0.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:92:PHE:CE1	1:A:123:ALA:HB1	0.50	2.42	17	1
1:A:78:ILE:HG22	1:A:93:TRP:CZ3	0.46	2.44	6	1
1:A:92:PHE:CE2	1:A:123:ALA:HB1	0.43	2.48	2	2
1:A:37:LYS:HA	1:A:71:VAL:HG22	0.41	1.90	4	1
1:A:34:HIS:HB2	1:A:66:LEU:HD11	0.41	1.91	2	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	134/145 (92%)	129±1 (96±1%)	5±2 (4±1%)	0±1 (0±0%)	50	83
All	All	2680/2900 (92%)	2577 (96%)	94 (4%)	9 (0%)	50	83

All 2 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	41	ILE	7
1	A	37	LYS	2

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	116/126 (92%)	114±1 (98±1%)	2±1 (2±1%)	64 94
All	All	2320/2520 (92%)	2272 (98%)	48 (2%)	64 94

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	135	ILE	9
1	A	103	GLN	6
1	A	134	ASP	5
1	A	57	ASP	3
1	A	87	VAL	3
1	A	91	ASP	3
1	A	96	THR	2
1	A	140	ASN	2
1	A	89	SER	2
1	A	143	SER	2
1	A	24	GLU	1
1	A	63	ASN	1
1	A	133	ASP	1
1	A	22	ASP	1
1	A	28	SER	1
1	A	71	VAL	1
1	A	46	ARG	1
1	A	18	ASP	1
1	A	45	ASP	1
1	A	38	THR	1
1	A	108	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)

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 90% for the well-defined parts and 89% for the entire structure.

7.1 Chemical shift list 1

File name: 2mpb_cs.str

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	1770
Number of shifts mapped to atoms	1770
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	11

7.1.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	143	0.03 ± 0.14	None needed (< 0.5 ppm)
$^{13}C_{\beta}$	134	0.29 ± 0.17	None needed (< 0.5 ppm)
$^{13}C'$	143	0.24 ± 0.19	None needed (< 0.5 ppm)
^{15}N	140	0.65 ± 0.34	None needed (imprecise)

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 90%, i.e. 1496 atoms were assigned a chemical shift out of a possible 1664. 12 out of 19 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	671/671 (100%)	268/268 (100%)	270/270 (100%)	133/133 (100%)
Sidechain	723/868 (83%)	444/501 (89%)	276/328 (84%)	3/39 (8%)

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	Total	^{1}H	^{13}C	^{15}N
Aromatic	102/125 (82%)	54/67 (81%)	46/54 (85%)	2/4 (50%)
Overall	1496/1664 (90%)	766/836 (92%)	592/652 (91%)	138/176 (78%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 89%, i.e. 1593 atoms were assigned a chemical shift out of a possible 1793. 12 out of 20 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	709/721 (98%)	283/288 (98%)	286/290 (99%)	140/143 (98%)
Sidechain	776/938 (83%)	477/542 (88%)	296/356 (83%)	3/40 (8%)
Aromatic	108/134 (81%)	58/72 (81%)	48/58 (83%)	2/4 (50%)
Overall	1593/1793 (89%)	818/902 (91%)	630/704 (89%)	145/187 (78%)

7.1.4 Statistically unusual chemical shifts (i)

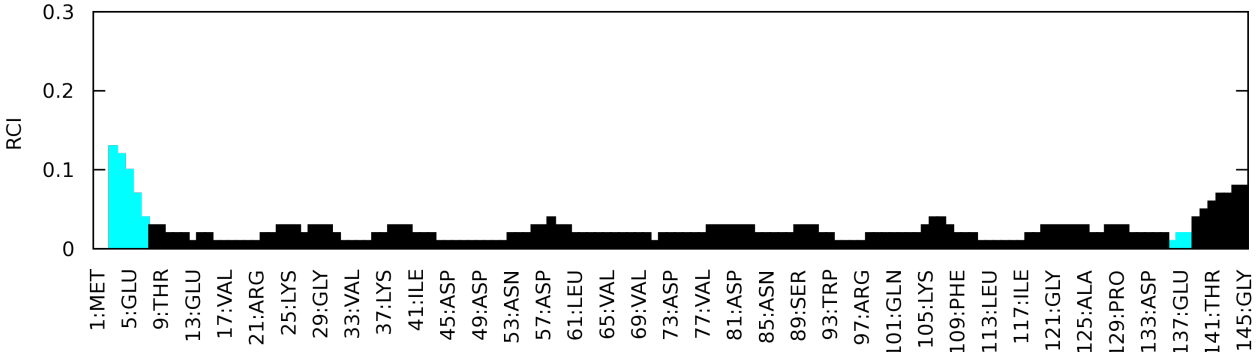
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

Mol	Chain	Res	Type	Atom	Shift, *ppm*	Expected range, *ppm*	Z-score
1	A	100	ILE	HG23	-1.33	2.13 – -0.57	-7.8
1	A	100	ILE	HG22	-1.33	2.13 – -0.57	-7.8
1	A	100	ILE	HG21	-1.33	2.13 – -0.57	-7.8
1	A	56	MET	HB2	-0.50	3.73 – 0.33	-7.4
1	A	88	VAL	HG21	-0.88	2.20 – -0.60	-6.0
1	A	88	VAL	HG22	-0.88	2.20 – -0.60	-6.0
1	A	88	VAL	HG23	-0.88	2.20 – -0.60	-6.0
1	A	63	ASN	HB2	0.96	4.36 – 1.26	-6.0
1	A	78	ILE	HG22	-0.73	2.13 – -0.57	-5.6
1	A	78	ILE	HG23	-0.73	2.13 – -0.57	-5.6
1	A	78	ILE	HG21	-0.73	2.13 – -0.57	-5.6

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
0.3
0.2
0.1
0
1:MET
5:GLU
9:THR
13:GLU
17:VAL
21:ARG
25:LYS
29:GLY
33:VAL
37:LYS
41:ILE
45:ASP
49:ASP
53:ASN
57:ASP
61:LEU
65:VAL
69:VAL
73:ASP
77:VAL
81:ASP
85:ASN
89:SER
93:TRP
97:ARG
101:GLN
105:LYS
109:PHE
113:LEU
117:ILE
121:GLY
125:ALA
129:PRO
133:ASP
137:GLU
141:THR
145:GLY