



Full wwPDB NMR Structure Validation Report ⓘ

Apr 27, 2016 – 03:49 AM BST

PDB ID : 2MZW
Title : Staphylococcus aureus FusB:EF-GC3 complex
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Deposited on : 2015-02-25

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 ⓘ 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

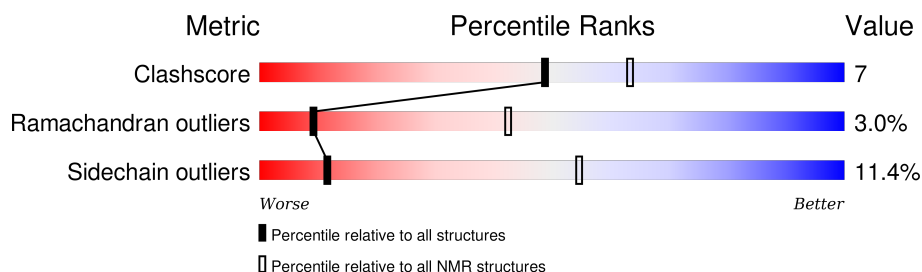
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 13%.

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	301	 66% 26% . . .
2	B	233	 76% 14% • 9%

2 Ensemble composition and analysis ⓘ

This entry contains 1 models. Identification of well-defined residues and clustering analysis are not possible.

3 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8050 atoms, of which 4018 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Elongation factor G.

Mol	Chain	Residues	Atoms						Trace
1	A	292	Total	C	H	N	O	S	0
			4471	1417	2214	378	445	17	

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	693	GLU	-	EXPRESSION TAG	UNP W8UT26
A	694	LEU	-	EXPRESSION TAG	UNP W8UT26
A	695	GLU	-	EXPRESSION TAG	UNP W8UT26
A	696	HIS	-	EXPRESSION TAG	UNP W8UT26
A	697	HIS	-	EXPRESSION TAG	UNP W8UT26
A	698	HIS	-	EXPRESSION TAG	UNP W8UT26
A	699	HIS	-	EXPRESSION TAG	UNP W8UT26
A	700	HIS	-	EXPRESSION TAG	UNP W8UT26
A	701	HIS	-	EXPRESSION TAG	UNP W8UT26

- Molecule 2 is a protein called Far1.

Mol	Chain	Residues	Atoms						Trace
2	B	213	Total	C	H	N	O	S	0
			3578	1143	1804	291	332	8	

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	EXPRESSION TAG	UNP Q8GNY5
B	-18	GLY	-	EXPRESSION TAG	UNP Q8GNY5
B	-17	SER	-	EXPRESSION TAG	UNP Q8GNY5
B	-16	SER	-	EXPRESSION TAG	UNP Q8GNY5
B	-15	HIS	-	EXPRESSION TAG	UNP Q8GNY5
B	-14	HIS	-	EXPRESSION TAG	UNP Q8GNY5
B	-13	HIS	-	EXPRESSION TAG	UNP Q8GNY5
B	-12	HIS	-	EXPRESSION TAG	UNP Q8GNY5
B	-11	HIS	-	EXPRESSION TAG	UNP Q8GNY5
B	-10	HIS	-	EXPRESSION TAG	UNP Q8GNY5
B	-9	SER	-	EXPRESSION TAG	UNP Q8GNY5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-8	SER	-	EXPRESSION TAG	UNP Q8GNY5
B	-7	GLY	-	EXPRESSION TAG	UNP Q8GNY5
B	-6	LEU	-	EXPRESSION TAG	UNP Q8GNY5
B	-5	VAL	-	EXPRESSION TAG	UNP Q8GNY5
B	-4	PRO	-	EXPRESSION TAG	UNP Q8GNY5
B	-3	ASN	-	EXPRESSION TAG	UNP Q8GNY5
B	-2	GLY	-	EXPRESSION TAG	UNP Q8GNY5
B	-1	SER	-	EXPRESSION TAG	UNP Q8GNY5
B	0	HIS	-	EXPRESSION TAG	UNP Q8GNY5

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	
3	B	1	Total	Zn
			1	1

5 Refinement protocol and experimental data overview

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

Of the 200 calculated structures, 1 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	2.33.0
X-PLOR NIH	refinement	2.33.0
HADDOCK	structure solution	2.1
HADDOCK	refinement	2.1

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)	2mzw_cs.str
Number of chemical shift lists	2
Total number of shifts	812
Number of shifts mapped to atoms	812
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	13%

No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality ⓘ

6.1 Standard geometry ⓘ

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

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	2257	2214	2204	51
2	B	1774	1804	1805	21
3	B	1	0	0	2
All	All	4032	4018	4009	60

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 7.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)
2:B:169:PHE:HB2	2:B:188:ILE:HG12	0.90	1.40
1:A:525:GLU:HG2	2:B:183:LYS:HB2	0.82	1.49
1:A:525:GLU:HA	2:B:183:LYS:HD3	0.73	1.58
2:B:195:CYS:SG	3:B:301:ZN:ZN	0.72	1.77
1:A:533:VAL:HB	1:A:534:PRO:HD3	0.70	1.63
2:B:195:CYS:HG	3:B:301:ZN:ZN	0.69	0.97
1:A:606:LYS:HA	1:A:645:TYR:HB3	0.68	1.64
1:A:522:PHE:HB3	2:B:168:LEU:HD13	0.67	1.67
2:B:157:CYS:SG	2:B:188:ILE:HA	0.65	2.31
1:A:467:LYS:HA	1:A:467:LYS:HE2	0.65	1.68
1:A:519:GLY:HA3	2:B:158:THR:HG23	0.65	1.67

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Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:546:LYS:O	1:A:550:GLU:HG2	0.64	1.91
1:A:639:ALA:HB3	2:B:35:VAL:HG21	0.60	1.72
1:A:602:GLU:HB2	1:A:678:VAL:CG1	0.58	2.28
1:A:606:LYS:HA	1:A:645:TYR:CB	0.55	2.30
2:B:8:HIS:CD2	2:B:201:ASP:HA	0.55	2.37
1:A:612:PRO:HA	1:A:639:ALA:HA	0.54	1.77
1:A:665:ARG:HD2	2:B:29:ASP:OD1	0.54	2.02
1:A:522:PHE:HB2	2:B:187:TYR:HE1	0.54	1.62
1:A:482:TYR:O	1:A:483:ARG:HB2	0.53	2.02
1:A:609:ILE:HB	1:A:642:VAL:HB	0.53	1.78
1:A:604:MET:HA	1:A:647:PRO:HA	0.52	1.80
2:B:23:VAL:HG21	2:B:36:ILE:HG13	0.52	1.81
1:A:485:THR:O	1:A:486:PHE:HB2	0.52	2.05
1:A:628:GLY:HA2	1:A:646:VAL:HG22	0.51	1.83
1:A:629:ARG:HB3	1:A:645:TYR:CE1	0.50	2.42
1:A:523:GLU:O	1:A:563:LYS:HA	0.49	2.06
1:A:486:PHE:CE2	1:A:515:PRO:HA	0.49	2.42
1:A:631:ASP:HB2	1:A:645:TYR:CE2	0.49	2.42
1:A:487:LYS:HG3	1:A:600:ILE:HD12	0.49	1.84
1:A:636:ARG:HB3	2:B:35:VAL:HG22	0.48	1.85
2:B:157:CYS:O	2:B:161:ASN:HA	0.48	2.08
2:B:90:SER:HB2	2:B:92:GLN:OE1	0.48	2.08
1:A:678:VAL:HB	1:A:679:PRO:HD2	0.47	1.85
1:A:535:ARG:O	1:A:538:ILE:HG12	0.47	2.09
1:A:510:HIS:HB2	1:A:568:ASP:HB3	0.47	1.86
1:A:647:PRO:O	1:A:679:PRO:HD3	0.47	2.09
1:A:538:ILE:HG13	1:A:539:PRO:HD3	0.46	1.87
1:A:603:PRO:HA	1:A:676:ALA:O	0.46	2.11
1:A:558:PRO:HA	2:B:156:PHE:N	0.46	2.25
1:A:484:GLU:HA	1:A:560:ILE:CG2	0.45	2.41
1:A:481:SER:HA	1:A:650:GLU:OE2	0.45	2.12
1:A:423:MET:O	1:A:427:LEU:HG	0.45	2.12
1:A:538:ILE:N	1:A:539:PRO:HD2	0.44	2.27
1:A:556:GLY:CA	2:B:153:VAL:HG21	0.44	2.43
1:A:413:GLU:HG3	1:A:447:GLN:HB3	0.44	1.88
1:A:602:GLU:HB2	1:A:678:VAL:HG11	0.44	1.89
1:A:406:PRO:O	1:A:453:MET:HA	0.44	2.12
1:A:684:GLU:HA	1:A:687:ILE:HG22	0.44	1.89
1:A:639:ALA:HB3	2:B:35:VAL:CG2	0.43	2.41
1:A:678:VAL:HB	1:A:679:PRO:CD	0.43	2.44
1:A:486:PHE:CE2	1:A:562:VAL:HA	0.43	2.49
1:A:602:GLU:HB2	1:A:678:VAL:HG13	0.42	1.91

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Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:600:ILE:HG23	1:A:676:ALA:O	0.42	2.15
1:A:539:PRO:HA	1:A:542:GLU:HG2	0.41	1.91
1:A:406:PRO:HB2	1:A:407:VAL:H	0.41	1.53
1:A:554:LEU:HD23	1:A:554:LEU:H	0.41	1.75
2:B:121:GLY:HA2	2:B:131:TYR:O	0.41	2.16
1:A:522:PHE:CA	1:A:561:ASP:HB3	0.40	2.46
1:A:520:ALA:HB3	2:B:187:TYR:HB2	0.40	1.92

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	290/301 (96%)	232 (80%)	44 (15%)	14 (5%)	5	28
2	B	211/233 (91%)	199 (94%)	11 (5%)	1 (0%)	38	79
All	All	501/534 (94%)	431 (86%)	55 (11%)	15 (3%)	9	42

All 15 Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	650	GLU
1	A	633	MET
1	A	522	PHE
1	A	406	PRO
1	A	597	ASP
1	A	572	HIS
2	B	159	ILE
1	A	520	ALA
1	A	483	ARG
1	A	558	PRO
1	A	561	ASP
1	A	501	GLY
1	A	486	PHE
1	A	533	VAL
1	A	442	ASP

6.3.2 Protein sidechains ⓘ

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	243/252 (96%)	210 (86%)	33 (14%)	9	49
2	B	204/221 (92%)	186 (91%)	18 (9%)	17	62
All	All	447/473 (95%)	396 (89%)	51 (11%)	11	54

All 51 residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
2	B	127	SER
1	A	681	SER
1	A	636	ARG
1	A	649	SER
1	A	678	VAL
2	B	177	SER
1	A	440	HIS
2	B	186	ASP
1	A	540	SER
1	A	445	THR
1	A	489	SER
1	A	436	THR
1	A	518	THR
2	B	163	GLU
1	A	498	ARG
1	A	522	PHE
1	A	472	GLU
2	B	71	GLU
1	A	458	LEU
2	B	175	THR
1	A	656	THR
2	B	144	LEU
1	A	431	GLN
1	A	486	PHE
2	B	158	THR
1	A	561	ASP
2	B	43	ASP
2	B	200	THR

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Mol	Chain	Res	Type
2	B	123	ASN
1	A	610	GLU
1	A	433	GLU
1	A	409	HIS
1	A	443	GLU
1	A	455	GLU
1	A	652	PHE
1	A	505	GLN
1	A	545	LEU
1	A	579	MET
1	A	411	SER
2	B	124	GLU
2	B	54	ASP
2	B	156	PHE
1	A	438	HIS
2	B	92	GLN
1	A	629	ARG
1	A	576	SER
2	B	161	ASN
2	B	52	ASP
1	A	550	GLU
2	B	53	ASP
1	A	499	GLN

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 1 ligands modelled in this entry, 1 is 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

The completeness of assignment taking into account all chemical shift lists is 13% for the well-defined parts and 13% for the entire structure.

7.1 Chemical shift list 1

File name: 2mzw_cs.str

Chemical shift list name: *C3_assigned_chem_shift_list*

7.1.1 Bookkeeping

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	560
Number of shifts mapped to atoms	560
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

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	129	0.34 ± 0.11	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	0	—	—
$^{13}\text{C}'$	117	0.11 ± 0.15	None needed (< 0.5 ppm)
^{15}N	157	-0.33 ± 0.22	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 9%, i.e. 560 atoms were assigned a chemical shift out of a possible 6300. 0 out of 72 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	560/2485 (23%)	157/990 (16%)	246/1010 (24%)	157/485 (32%)
Sidechain	0/3326 (0%)	0/1953 (0%)	0/1230 (0%)	0/143 (0%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	0/489 (0%)	0/255 (0%)	0/213 (0%)	0/21 (0%)
Overall	560/6300 (9%)	157/3198 (5%)	246/2453 (10%)	157/649 (24%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 9%, i.e. 560 atoms were assigned a chemical shift out of a possible 6300. 0 out of 72 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	560/2485 (23%)	157/990 (16%)	246/1010 (24%)	157/485 (32%)
Sidechain	0/3326 (0%)	0/1953 (0%)	0/1230 (0%)	0/143 (0%)
Aromatic	0/489 (0%)	0/255 (0%)	0/213 (0%)	0/21 (0%)
Overall	560/6300 (9%)	157/3198 (5%)	246/2453 (10%)	157/649 (24%)

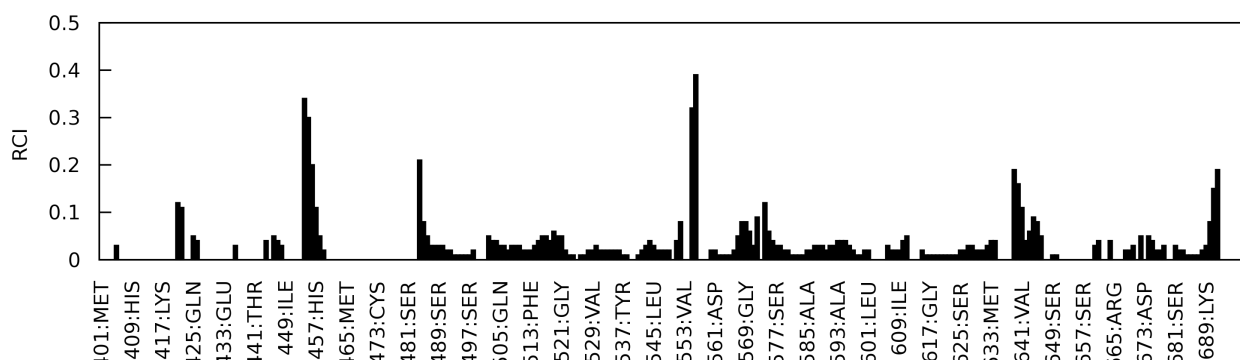
7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

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: 2mzw_cs.str

Chemical shift list name: *C3_assigned_chem_shift_list_dup*

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	252
Number of shifts mapped to atoms	252
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](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	0	—	—
$^{13}\text{C}_\beta$	0	—	—
$^{13}\text{C}'$	0	—	—
^{15}N	126	-0.24 ± 0.18	None needed (< 0.5 ppm)

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 4%, i.e. 248 atoms were assigned a chemical shift out of a possible 6300. 0 out of 72 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	248/2485 (10%)	124/990 (13%)	0/1010 (0%)	124/485 (26%)
Sidechain	0/3326 (0%)	0/1953 (0%)	0/1230 (0%)	0/143 (0%)
Aromatic	0/489 (0%)	0/255 (0%)	0/213 (0%)	0/21 (0%)
Overall	248/6300 (4%)	124/3198 (4%)	0/2453 (0%)	124/649 (19%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 4%, i.e. 248 atoms were assigned a chemical shift out of a possible 6300. 0 out of 72 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	248/2485 (10%)	124/990 (13%)	0/1010 (0%)	124/485 (26%)

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	Total	¹ H	¹³ C	¹⁵ N
Sidechain	0/3326 (0%)	0/1953 (0%)	0/1230 (0%)	0/143 (0%)
Aromatic	0/489 (0%)	0/255 (0%)	0/213 (0%)	0/21 (0%)
Overall	248/6300 (4%)	124/3198 (4%)	0/2453 (0%)	124/649 (19%)

7.2.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.2.5 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 B:

