



wwPDB NMR Structure Validation Summary Report ⓘ

Mar 1, 2026 – 05:59 PM UTC

PDB ID : 2MFH / pdb_00002mfh
BMRB ID : 19549
Title : Csr/Rsm protein-RNA recognition - A molecular affinity ruler: RsmZ(36-44)
/RsmE(dimer) 2:1 complex
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Deposited on : 2013-10-11

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

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

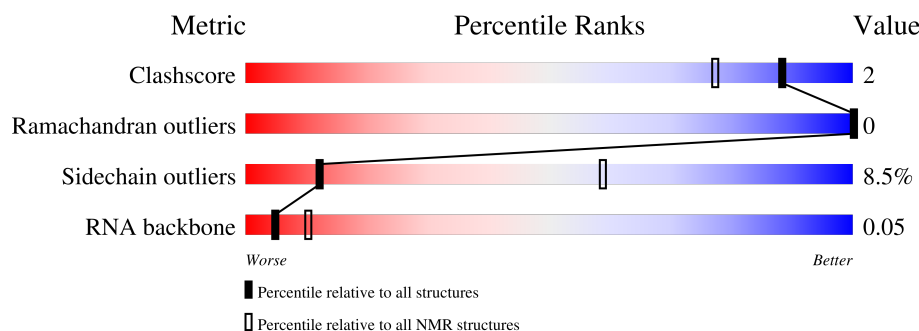
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 42%.

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	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823
RNA backbone	8273	777

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	70	
1	C	70	
2	B	9	
2	D	9	

2 Ensemble composition and analysis

This entry contains 20 models. Model 2 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:1-A:48, C:1-C:48 (96)	0.27	2

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 4 clusters and 3 single-model clusters were found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2416 atoms, of which 1142 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Carbon storage regulator homolog.

Mol	Chain	Residues	Atoms						Trace
1	A	59	Total	C	H	N	O	S	0
			920	280	472	81	86	1	
1	C	59	Total	C	H	N	O	S	0
			920	280	472	81	86	1	

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	60	LYS	-	expression tag	UNP Q5MXB2
A	61	ARG	-	expression tag	UNP Q5MXB2
A	62	GLU	-	expression tag	UNP Q5MXB2
A	63	THR	-	expression tag	UNP Q5MXB2
A	64	PRO	-	expression tag	UNP Q5MXB2
A	65	HIS	-	expression tag	UNP Q5MXB2
A	66	HIS	-	expression tag	UNP Q5MXB2
A	67	HIS	-	expression tag	UNP Q5MXB2
A	68	HIS	-	expression tag	UNP Q5MXB2
A	69	HIS	-	expression tag	UNP Q5MXB2
A	70	HIS	-	expression tag	UNP Q5MXB2
C	60	LYS	-	expression tag	UNP Q5MXB2
C	61	ARG	-	expression tag	UNP Q5MXB2
C	62	GLU	-	expression tag	UNP Q5MXB2
C	63	THR	-	expression tag	UNP Q5MXB2
C	64	PRO	-	expression tag	UNP Q5MXB2
C	65	HIS	-	expression tag	UNP Q5MXB2
C	66	HIS	-	expression tag	UNP Q5MXB2
C	67	HIS	-	expression tag	UNP Q5MXB2
C	68	HIS	-	expression tag	UNP Q5MXB2
C	69	HIS	-	expression tag	UNP Q5MXB2
C	70	HIS	-	expression tag	UNP Q5MXB2

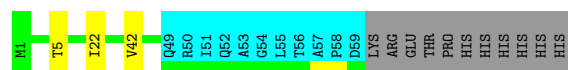
- Molecule 2 is a RNA chain called RsmZ(36-44) RNA.

Mol	Chain	Residues	Atoms						Trace
2	B	9	Total	C	H	N	O	P	0
			288	86	99	35	60	8	

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Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	P	
2	D	9	288	86	99	35	60	8	0



- | | | |
|----|--|-----|
| M1 | | I22 | V42 | Q49 | R50 | I51 | Q52 | A53 | G54 | L55 | T56 | A57 | P58 | D59 | LVS | ARG | GLU | THR | PRO | HIS | HIS | HIS | HIS | HIS |
|----|--|-----|

-

- | |
|-----|
| U36 |
| C37 |
| A38 |
| G39 |
| G40 |
| A41 |
| C42 |
| A43 |
| U44 |

5 Refinement protocol and experimental data overview

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

Of the 999 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
CYANA	structure solution	
Amber	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)	working_cs.cif
Number of chemical shift lists	2
Total number of shifts	878
Number of shifts mapped to atoms	824
Number of unparsed shifts	0
Number of shifts with mapping errors	54
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	42%

6 Model quality

6.1 Standard geometry

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	0.49±0.01	0±0/370 (0.0± 0.0%)	0.98±0.02	0±0/501 (0.0± 0.0%)
1	C	0.49±0.01	0±0/370 (0.0± 0.0%)	0.98±0.02	0±0/501 (0.0± 0.0%)
2	B	0.73±0.01	0±0/211 (0.0± 0.0%)	1.10±0.07	0±0/327 (0.1± 0.1%)
2	D	0.73±0.01	0±0/211 (0.0± 0.0%)	1.07±0.06	0±1/327 (0.1± 0.2%)
All	All	0.59	0/23240 (0.0%)	1.02	15/33120 (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	C	0.0±0.0	0.1±0.2
2	D	0.0±0.0	0.1±0.4
2	B	0.0±0.0	0.1±0.2
All	All	0	5

There are no bond-length outliers.

5 of 10 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(°)	Ideal(°)	Models	
								Worst	Total
2	D	39	G	C4'-C3'-C2'	-6.20	96.40	102.60	11	1
2	B	44	U	C5'-C4'-C3'	-6.04	106.93	116.00	6	2
2	D	44	U	C5'-C4'-C3'	-5.86	107.21	116.00	10	1
2	B	42	C	O4'-C4'-C3'	5.62	109.62	104.00	18	2
2	B	39	G	C4'-C3'-C2'	-5.61	96.99	102.60	19	1

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)
2	D	42	C	Sidechain	1
2	D	41	A	Sidechain	1
1	C	6	ARG	Sidechain	1
2	D	43	A	Sidechain	1
2	B	40	G	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	367	390	390	1±1
1	C	367	390	390	1±1
2	B	189	99	99	2±1
2	D	189	99	99	2±2
All	All	22240	19560	19560	78

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 2.

5 of 43 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:43:A:C5	2:B:44:U:H1'	0.61	2.31	13	2
2:D:43:A:O4'	2:D:44:U:C6	0.55	2.60	15	8
2:D:43:A:C5	2:D:44:U:H1'	0.53	2.37	10	1
2:D:42:C:C3'	2:D:43:A:H5''	0.52	2.35	8	4
1:A:23:LEU:HD21	2:D:42:C:C2	0.52	2.38	5	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	47/70 (67%)	44±1 (93±2%)	3±1 (7±2%)	0±0 (0±0%)	100	100
1	C	47/70 (67%)	44±1 (93±2%)	3±1 (7±2%)	0±0 (0±0%)	100	100
All	All	1880/2800 (67%)	1748 (93%)	132 (7%)	0 (0%)	100	100

There are no Ramachandran outliers.

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	41/60 (68%)	38±1 (92±3%)	3±1 (8±3%)	13	61
1	C	41/60 (68%)	37±1 (91±3%)	4±1 (9±3%)	11	57
All	All	1640/2400 (68%)	1501 (92%)	139 (8%)	12	59

5 of 25 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	42	VAL	20
1	C	42	VAL	19
1	C	22	ILE	17
1	A	22	ILE	17
1	C	5	THR	9

6.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers	Suiteness
2	B	8/9 (89%)	4±0 (50±0%)	1±1 (17±8%)	0.06±0.00
2	D	8/9 (89%)	4±0 (50±0%)	2±1 (23±15%)	0.10±0.00
All	All	332/360 (92%)	160 (48%)	64 (19%)	0.08

The overall RNA backbone suiteness is 0.05.

5 of 8 unique RNA backbone outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
2	B	37	C	20
2	B	40	G	20
2	B	42	C	20
2	B	43	A	20
2	D	37	C	20

5 of 14 unique RNA pucker outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
2	D	43	A	12
2	D	42	C	10
2	B	42	C	9
2	B	43	A	8
2	D	36	U	7

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 oligosaccharides 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

The completeness of assignment taking into account all chemical shift lists is 42% for the well-defined parts and 42% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chem_shift_list_1*

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	746
Number of shifts mapped to atoms	692
Number of unparsed shifts	0
Number of shifts with mapping errors	54
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

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

- No matching atom found in the structure. First 5 (of 54) occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	60	LYS	H	8.071	0.000	1
1	A	60	LYS	HA	4.241	0.000	1
1	A	60	LYS	HB2	1.786	0.000	.
1	A	60	LYS	HB3	1.671	0.000	.
1	A	60	LYS	HD2	1.568	0.000	.
1	A	60	LYS	HD3	1.568	0.000	.
1	A	60	LYS	HE2	2.896	0.000	.
1	A	60	LYS	HE3	2.896	0.000	.
1	A	60	LYS	HG2	1.339	0.000	.
1	A	60	LYS	HG3	1.289	0.000	.
1	A	60	LYS	CA	56.099	0.000	1
1	A	60	LYS	CB	32.623	0.017	1
1	A	60	LYS	CD	28.854	0.000	1
1	A	60	LYS	CE	42.23	0.000	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	60	LYS	CG	24.65	0.018	1
1	A	60	LYS	N	121.441	0.000	1
1	A	61	ARG	H	7.926	0.000	1
1	A	61	ARG	HA	4.052	0.000	1
1	A	61	ARG	HB2	1.676	0.000	.
1	A	61	ARG	HB3	1.652	0.000	.
1	A	61	ARG	HD2	2.996	0.000	.
1	A	61	ARG	HD3	2.996	0.000	.
1	A	61	ARG	HG2	1.476	0.000	.
1	A	61	ARG	HG3	1.404	0.000	.
1	A	61	ARG	CA	56.519	0.000	1
1	A	61	ARG	CB	30.568	0.022	1
1	A	61	ARG	CD	43.29	0.000	1
1	A	61	ARG	CG	27.02	0.011	1
1	A	61	ARG	N	121.098	0.000	1
1	A	62	GLU	H	8.371	0.000	1
1	A	62	GLU	HA	4.167	0.000	1
1	A	62	GLU	HB2	1.813	0.000	.
1	A	62	GLU	HB3	1.754	0.000	.
1	A	62	GLU	HG2	2.153	0.000	.
1	A	62	GLU	HG3	2.127	0.000	.
1	A	62	GLU	CA	56.835	0.000	1
1	A	62	GLU	CB	30.017	0.017	1
1	A	62	GLU	CG	36.286	0.001	1
1	A	62	GLU	N	121.228	0.000	1
1	A	63	THR	H	7.837	0.000	1
1	A	63	THR	HA	4.159	0.000	1
1	A	63	THR	HB	4.124	0.000	1
1	A	63	THR	HG21	1.051	0.000	1
1	A	63	THR	HG22	1.051	0.000	1
1	A	63	THR	HG23	1.051	0.000	1
1	A	63	THR	CA	62.064	0.000	1
1	A	63	THR	CB	69.578	0.000	1
1	A	63	THR	CG2	21.64	0.000	1
1	A	63	THR	N	122.477	0.000	1
1	A	65	HIS	HA	4.341	0.000	1
1	A	65	HIS	HB2	3.088	0.000	.
1	A	65	HIS	HB3	2.942	0.000	.
1	A	65	HIS	CA	57.503	0.000	1
1	A	65	HIS	CB	30.994	0.065	1

7.1.2 Chemical shift referencing ⓘ

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	64	-0.17 ± 0.11	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	57	0.05 ± 0.22	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	60	-0.98 ± 0.27	Should be applied

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 34%, i.e. 570 atoms were assigned a chemical shift out of a possible 1676. 0 out of 16 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	193/486 (40%)	99/200 (50%)	48/192 (25%)	46/94 (49%)
Sidechain	367/820 (45%)	250/538 (46%)	110/252 (44%)	7/30 (23%)
Aromatic	10/32 (31%)	6/16 (38%)	4/14 (29%)	0/2 (0%)
Sugar	0/198 (0%)	0/108 (0%)	0/90 (0%)	0/0 (—%)
Base	0/140 (0%)	0/86 (0%)	0/32 (0%)	0/22 (0%)
Overall	570/1676 (34%)	355/948 (37%)	162/580 (28%)	53/148 (36%)

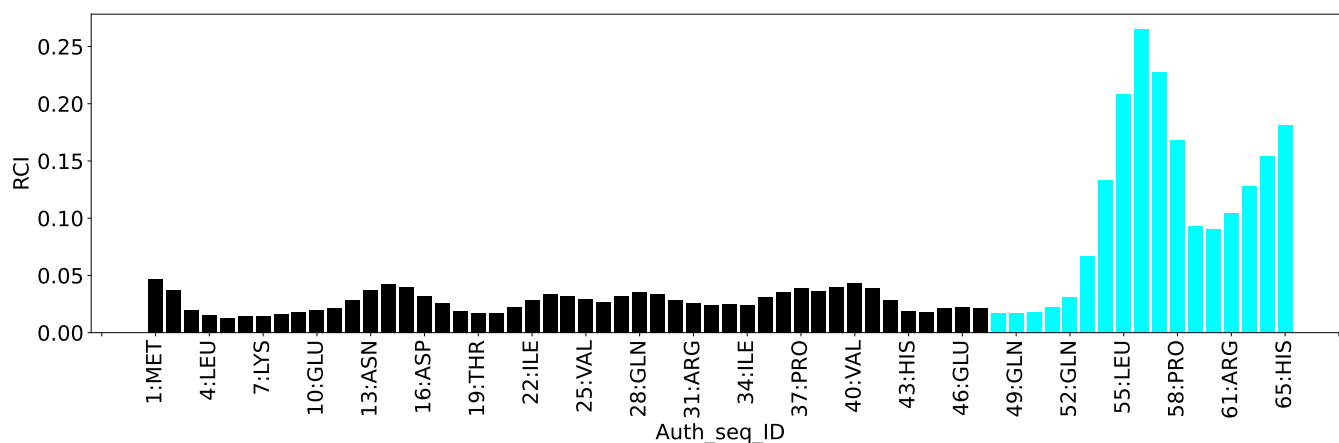
7.1.4 Statistically unusual chemical shifts ⓘ

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots ⓘ

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. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:



7.2 Chemical shift list 2

File name: working_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	132
Number of shifts mapped to atoms	132
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

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 8%, i.e. 128 atoms were assigned a chemical shift out of a possible 1676. 0 out of 16 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	0/486 (0%)	0/200 (0%)	0/192 (0%)	0/94 (0%)

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	Total	¹H	¹³C	¹⁵N
Sidechain	0/820 (0%)	0/538 (0%)	0/252 (0%)	0/30 (0%)
Aromatic	0/32 (0%)	0/16 (0%)	0/14 (0%)	0/2 (0%)
Sugar	98/198 (49%)	53/108 (49%)	45/90 (50%)	0/0 (—%)
Base	30/140 (21%)	15/86 (17%)	14/32 (44%)	1/22 (5%)
Overall	128/1676 (8%)	68/948 (7%)	59/580 (10%)	1/148 (1%)

7.2.4 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.

List Id	Chain	Res	Type	Atom	Shift, <i>ppm</i>	Expected range, <i>ppm</i>	Z-score
2	B	40	G	H5''	2.73	2.98 – 5.38	-6.0

7.2.5 Random Coil Index (RCI) plots [i](#)

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