



# Full wwPDB NMR Structure Validation Report ⓘ

Feb 13, 2017 – 02:35 am GMT

PDB ID : 5AGQ  
Title : Solution structure of the TAM domain of human TIP5 BAZ2A involved in epigenetic regulation of rRNA genes  
Authors : Anosova, I.; Melnik, S.; Tripsianes, K.; Kateb, F.; Grummt, I.; Sattler, M.  
Deposited on : 2015-02-03

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
RCI : v\_1n\_11\_5\_13\_A (Berjanski et al., 2005)  
PANAV : Wang et al. (2010)  
ShiftChecker : trunk28760  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

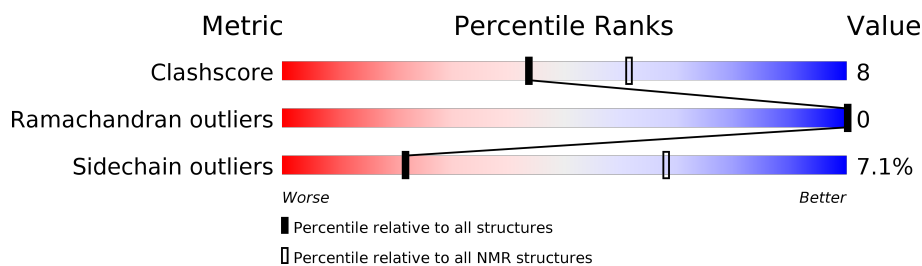
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 80%.

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	125131	11601
Ramachandran outliers	121729	10391
Sidechain outliers	121581	10367

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  $\geq 3$ , 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	113	<div>68% 22% 6% •</div>

## 2 Ensemble composition and analysis

This entry contains 20 models. Model 16 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:520-A:621 (102)	0.16	16

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 4 single-model clusters were found.

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

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1829 atoms, of which 909 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A.

Mol	Chain	Residues	Atoms						Trace
1	A	109	Total	C	H	N	O	S	0
			1829	581	909	178	158	3	

There are 5 discrepancies between the modelled and reference sequences:

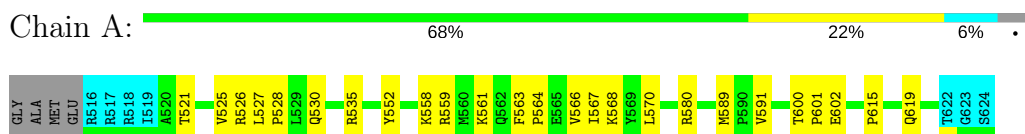
Chain	Residue	Modelled	Actual	Comment	Reference
A	512	GLY	-	EXPRESSION TAG	UNP Q9UIF9
A	513	ALA	-	EXPRESSION TAG	UNP Q9UIF9
A	514	MET	-	EXPRESSION TAG	UNP Q9UIF9
A	515	GLU	-	EXPRESSION TAG	UNP Q9UIF9
A	624	SER	-	CLONING ARTIFACT	UNP Q9UIF9

## 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: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A

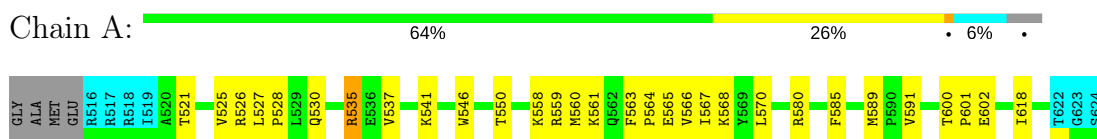


### 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

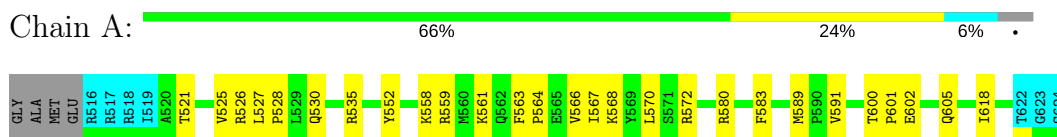
#### 4.2.1 Score per residue for model 1

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



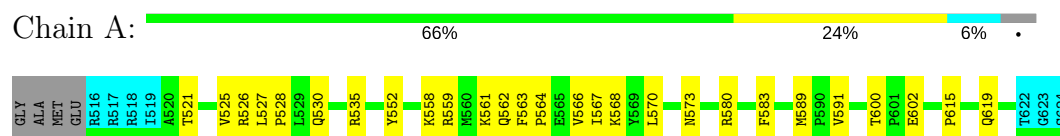
#### 4.2.2 Score per residue for model 2

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



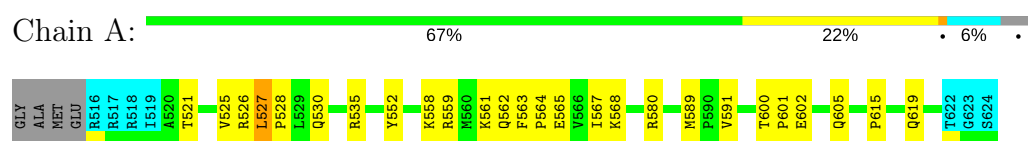
### 4.2.3 Score per residue for model 3

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



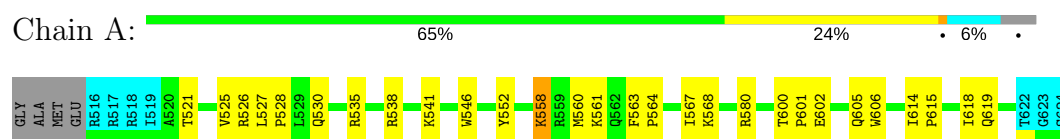
### 4.2.4 Score per residue for model 4

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



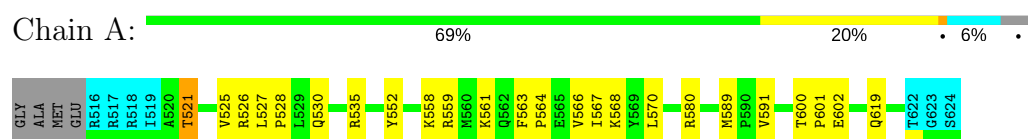
### 4.2.5 Score per residue for model 5

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



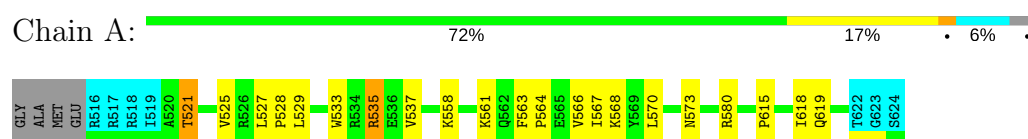
### 4.2.6 Score per residue for model 6

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



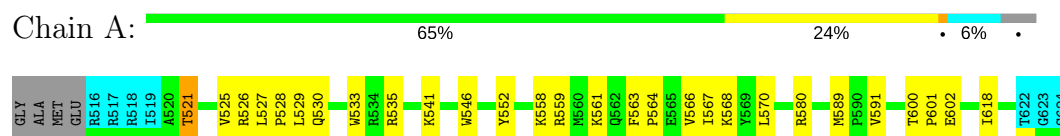
### 4.2.7 Score per residue for model 7

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



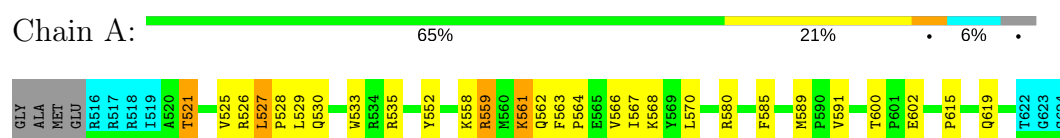
### 4.2.8 Score per residue for model 8

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



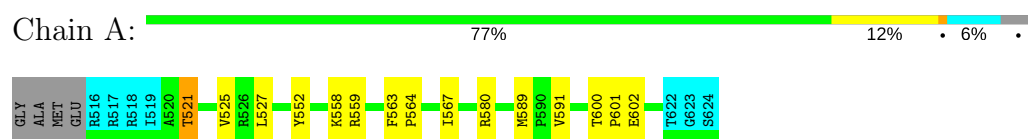
### 4.2.9 Score per residue for model 9

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



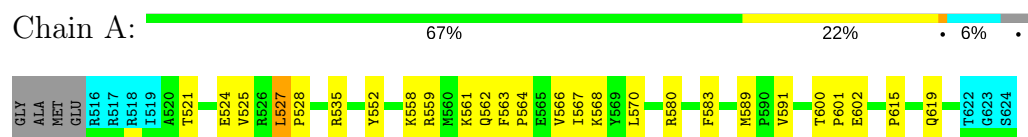
### 4.2.10 Score per residue for model 10

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



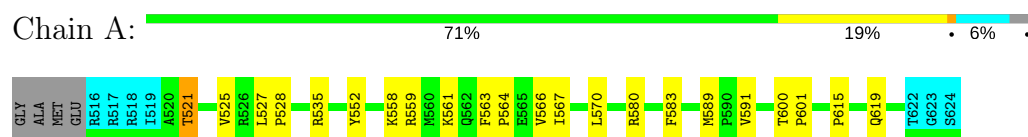
### 4.2.11 Score per residue for model 11

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



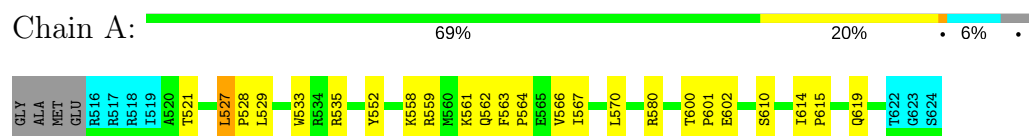
### 4.2.12 Score per residue for model 12

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



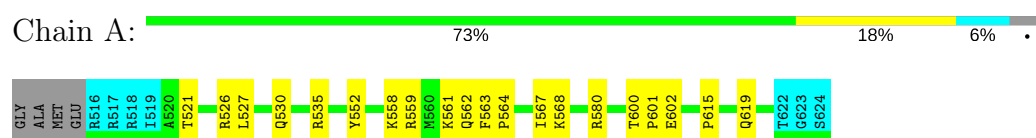
### 4.2.13 Score per residue for model 13

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



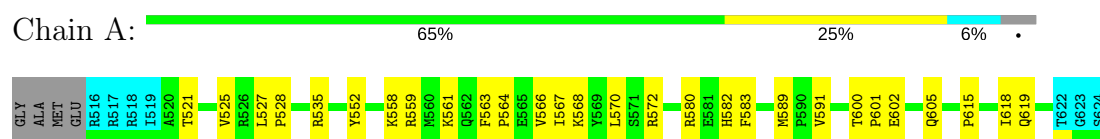
### 4.2.14 Score per residue for model 14

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



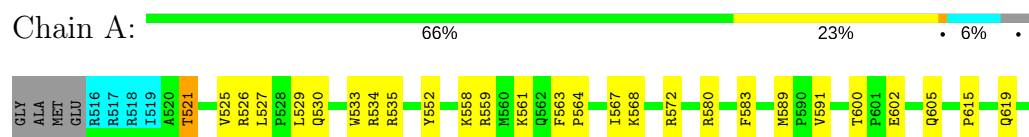
### 4.2.15 Score per residue for model 15

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



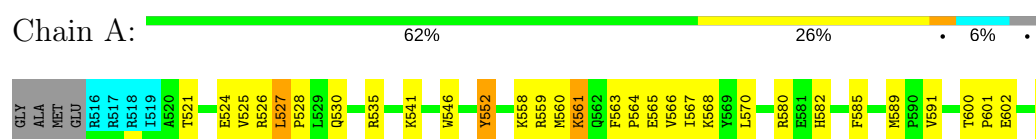
### 4.2.16 Score per residue for model 16 (medoid)

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



### 4.2.17 Score per residue for model 17

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A



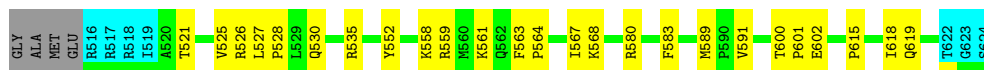




#### 4.2.18 Score per residue for model 18

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A

Chain A: 68% 22% 6% .



#### 4.2.19 Score per residue for model 19

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A

Chain A: 66% 23% 6% .



#### 4.2.20 Score per residue for model 20

- Molecule 1: BROMODOMAIN ADJACENT TO ZINC FINGER DOMAIN PROTEIN 2A

Chain A: 65% 25% 6% .



## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *ARIA2*, *CNS*.

Of the 200 calculated structures, 20 were deposited, based on the following criterion: *BEST ENERGY*.

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

Software name	Classification	Version
CNS	refinement	
SPARKY	structure solution	
CYANA	structure solution	3.0

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)	5agq_cs.str
Number of chemical shift lists	1
Total number of shifts	1287
Number of shifts mapped to atoms	1287
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	80%

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

## 6 Model quality

### 6.1 Standard geometry

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	861	842	842	14±3
All	All	17220	16840	16840	274

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 8.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:563:PHE:CE1	1:A:567:ILE:HD11	0.65	2.26	3	20
1:A:535:ARG:HG3	1:A:552:TYR:CZ	0.64	2.28	20	13
1:A:589:MET:HG3	1:A:591:VAL:HG23	0.63	1.68	17	16
1:A:535:ARG:HG3	1:A:552:TYR:CE1	0.59	2.33	9	13
1:A:525:VAL:HB	1:A:591:VAL:HG21	0.57	1.75	17	16
1:A:615:PRO:O	1:A:619:GLN:HG2	0.54	2.03	3	14
1:A:563:PHE:N	1:A:564:PRO:HD2	0.53	2.18	15	20
1:A:600:THR:OG1	1:A:602:GLU:HG2	0.53	2.03	14	18
1:A:526:ARG:O	1:A:530:GLN:HG3	0.52	2.05	1	13
1:A:566:VAL:O	1:A:570:LEU:HG	0.52	2.04	12	14
1:A:535:ARG:HG3	1:A:552:TYR:CE2	0.51	2.41	17	2
1:A:564:PRO:O	1:A:568:LYS:HE3	0.51	2.04	3	11
1:A:529:LEU:HA	1:A:533:TRP:O	0.51	2.06	7	5
1:A:525:VAL:O	1:A:528:PRO:HD2	0.49	2.08	12	13

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:568:LYS:O	1:A:572:ARG:HG3	0.49	2.08	20	4
1:A:535:ARG:CZ	1:A:537:VAL:HG23	0.48	2.38	7	2
1:A:563:PHE:O	1:A:567:ILE:HG12	0.46	2.11	15	7
1:A:558:LYS:HG2	1:A:560:MET:SD	0.46	2.50	5	1
1:A:563:PHE:CZ	1:A:567:ILE:HD11	0.45	2.46	6	6
1:A:521:THR:O	1:A:525:VAL:HG23	0.45	2.11	16	8
1:A:541:LYS:HD3	1:A:546:TRP:CE2	0.45	2.47	17	4
1:A:560:MET:CG	1:A:565:GLU:HB3	0.44	2.42	17	2
1:A:552:TYR:CE2	1:A:583:PHE:HD1	0.44	2.30	19	7
1:A:610:SER:O	1:A:614:ILE:HG13	0.44	2.12	19	3
1:A:533:TRP:O	1:A:534:ARG:HD2	0.44	2.12	16	1
1:A:524:GLU:O	1:A:527:LEU:HB2	0.43	2.13	17	2
1:A:590:PRO:HB3	1:A:614:ILE:HG21	0.43	1.88	20	1
1:A:552:TYR:CE1	1:A:583:PHE:HD1	0.43	2.30	11	1
1:A:600:THR:HB	1:A:601:PRO:HD2	0.43	1.90	6	15
1:A:600:THR:HG21	1:A:605:GLN:CD	0.43	2.35	15	5
1:A:527:LEU:N	1:A:528:PRO:HD2	0.42	2.29	13	4
1:A:561:LYS:HA	1:A:585:PHE:CE1	0.42	2.49	9	2
1:A:564:PRO:O	1:A:568:LYS:HE2	0.42	2.14	8	1
1:A:538:ARG:HB3	1:A:606:TRP:CH2	0.42	2.50	5	1
1:A:525:VAL:HA	1:A:582:HIS:O	0.41	2.15	17	2
1:A:614:ILE:O	1:A:618:ILE:HG13	0.41	2.16	5	1
1:A:559:ARG:H	1:A:559:ARG:NE	0.41	2.12	9	1
1:A:550:THR:HB	1:A:585:PHE:HA	0.40	1.93	1	1
1:A:589:MET:HG3	1:A:591:VAL:CG2	0.40	2.45	1	1
1:A:593:ASP:HB3	1:A:595:PHE:CZ	0.40	2.51	20	1
1:A:563:PHE:N	1:A:564:PRO:CD	0.40	2.85	15	1
1:A:562:GLN:HB2	1:A:565:GLU:CG	0.40	2.46	4	1

## 6.3 Torsion angles ⓘ

### 6.3.1 Protein backbone ⓘ

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	102/113 (90%)	97±2 (95±1%)	5±2 (5±1%)	0±0 (0±0%)	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	2040/2260 (90%)	1945 (95%)	95 (5%)	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	92/100 (92%)	86±1 (93±1%)	7±1 (7±1%)	22	68
All	All	1840/2000 (92%)	1710 (93%)	130 (7%)	22	68

All 11 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	527	LEU	20
1	A	558	LYS	20
1	A	521	THR	20
1	A	580	ARG	20
1	A	561	LYS	19
1	A	559	ARG	18
1	A	562	GLN	6
1	A	535	ARG	2
1	A	552	TYR	2
1	A	573	ASN	2
1	A	568	LYS	1

### 6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 6.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

The completeness of assignment taking into account all chemical shift lists is 80% for the well-defined parts and 78% for the entire structure.

### 7.1 Chemical shift list 1

File name: 5agq\_cs.str

Chemical shift list name: *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	1287
Number of shifts mapped to atoms	1287
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	12

#### 7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	98	$-0.08 \pm 0.22$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}_\beta$	99	$-0.08 \pm 0.21$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
$^{15}\text{N}$	96	$-0.11 \pm 0.34$	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 80%, i.e. 1113 atoms were assigned a chemical shift out of a possible 1388. 1 out of 13 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	371/494 (75%)	191/196 (97%)	91/204 (45%)	89/94 (95%)
Sidechain	609/749 (81%)	386/449 (86%)	216/252 (86%)	7/48 (15%)

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	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Aromatic	133/145 (92%)	69/77 (90%)	60/60 (100%)	4/8 (50%)
Overall	1113/1388 (80%)	646/722 (89%)	367/516 (71%)	100/150 (67%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 78%, i.e. 1171 atoms were assigned a chemical shift out of a possible 1493. 1 out of 13 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	398/529 (75%)	204/210 (97%)	98/218 (45%)	96/101 (95%)
Sidechain	640/819 (78%)	408/491 (83%)	225/271 (83%)	7/57 (12%)
Aromatic	133/145 (92%)	69/77 (90%)	60/60 (100%)	4/8 (50%)
Overall	1171/1493 (78%)	681/778 (88%)	383/549 (70%)	107/166 (64%)

#### 7.1.4 Statistically unusual chemical shifts ⓘ

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	552	TYR	CD1	118.61	139.11 – 126.41	-11.1
1	A	552	TYR	CD2	118.61	140.11 – 125.31	-9.5
1	A	538	ARG	HD2	1.08	4.27 – 1.97	-8.9
1	A	538	ARG	HD3	0.97	4.36 – 1.86	-8.6
1	A	597	GLU	HG3	0.70	3.31 – 1.21	-7.4
1	A	537	VAL	HB	-0.14	3.59 – 0.39	-6.6
1	A	561	LYS	HE3	1.71	3.86 – 1.96	-6.3
1	A	555	PRO	HD3	1.28	5.52 – 1.72	-6.2
1	A	597	GLU	HG2	1.15	3.33 – 1.23	-5.4
1	A	535	ARG	HG3	0.01	3.00 – 0.10	-5.3
1	A	538	ARG	HG2	0.18	2.92 – 0.22	-5.1
1	A	541	LYS	HB3	0.37	3.10 – 0.40	-5.1

#### 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.



Random coil index (RCI) for chain A:

