wwPDB NMR Structure Validation Summary Report

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Title	:	NMR structure of the gamma subunit of cGMP phosphodiesterase
Authors	:	Song, J.; Guo, L.W.; Ruoho, A.E.; Markley, J.L.; Center for Eukaryotic Structural Genomics (CESG)
Deposited on	:	2007-08-14

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 (i) 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 (i)) were used in the production of this report:

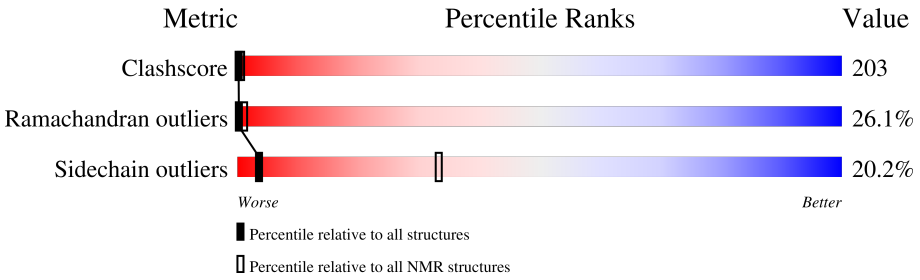
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 72%.

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	87	21% · 77%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA and RNA chains that are outliers for geometric criteria:

Mol	Chain	Compound	Res	Total models with violations	
				Chirality	Geometry
2	A	RCY	110	47	-
2	A	RCY	121	45	-
2	A	RCY	130	61	1
2	A	RCY	138	52	-
2	A	RCY	150	53	-
2	A	RCY	160	70	-
2	A	RCY	168	69	-

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Mol	Chain	Compound	Res	Total models with violations	
				Chirality	Geometry
2	A	RCY	173	50	-
2	A	RCY	176	56	-
2	A	RCY	187	57	-

2 Ensemble composition and analysis (i)

This entry contains 100 models. Model 6 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *fewest violations*.

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:59-A:78 (20)	3.41	6

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 6 clusters and 2 single-model clusters were found.

Cluster number	Models
1	11, 12, 16, 17, 19, 20, 21, 22, 26, 27, 28, 29, 32, 33, 36, 37, 38, 39, 40, 41, 44, 45, 48, 49, 52, 53, 56, 58, 67, 70, 71, 74, 82, 83, 84, 86, 88, 89, 91, 92, 93, 94, 96, 97, 99
2	1, 2, 3, 7, 8, 10, 13, 14, 23, 24, 25, 30, 31, 35, 42, 43, 46, 50, 55, 57, 59, 60, 61, 62, 64, 65, 66, 72, 73, 75, 79, 80, 87, 98, 100
3	4, 5, 6, 15, 18, 47, 54, 69, 77, 78, 85
4	51, 63, 68
5	34, 76
6	9, 95
Single-model clusters	81; 90

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1478 atoms, of which 659 are hydrogens and 0 are deuteriums.

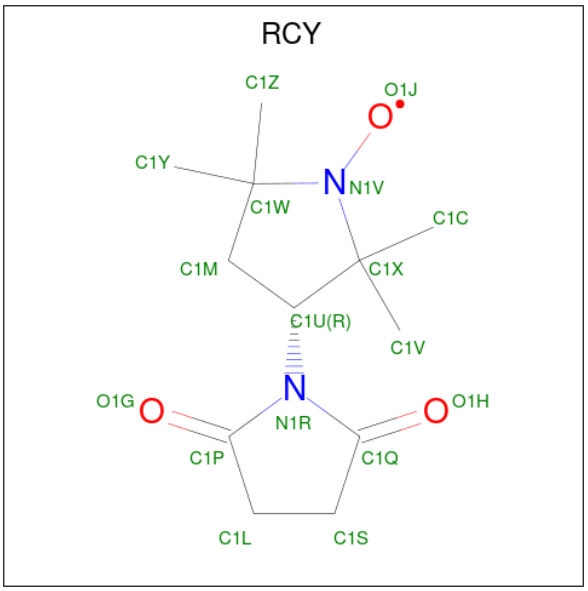
- Molecule 1 is a protein called Retinal rod rhodopsin-sensitive cGMP 3',5'-cyclic phosphodiesterase subunit gamma.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	87	1288	395	639	122	119	13	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	CYS	PHE	engineered mutation	UNP P04972
A	21	CYS	VAL	engineered mutation	UNP P04972
A	30	CYS	PHE	engineered mutation	UNP P04972
A	38	CYS	PHE	engineered mutation	UNP P04972
A	50	CYS	PHE	engineered mutation	UNP P04972
A	60	CYS	LEU	engineered mutation	UNP P04972
A	73	CYS	PHE	engineered mutation	UNP P04972
A	76	CYS	LEU	engineered mutation	UNP P04972
A	87	CYS	ILE	engineered mutation	UNP P04972

- Molecule 2 is (3'R)-1'-oxyl-2',2',5',5'-tetramethyl-1,3'-bipyrrolidine-2,5-dione (three-letter code: RCY) (formula: $C_{12}H_{19}N_2O_3$).



Mol	Chain	Residues	Atoms				
2	A	1	Total 19	C 12	H 2	N 2	O 3
2	A	1	Total 19	C 12	H 2	N 2	O 3
2	A	1	Total 19	C 12	H 2	N 2	O 3
2	A	1	Total 19	C 12	H 2	N 2	O 3
2	A	1	Total 19	C 12	H 2	N 2	O 3
2	A	1	Total 19	C 12	H 2	N 2	O 3
2	A	1	Total 19	C 12	H 2	N 2	O 3
2	A	1	Total 19	C 12	H 2	N 2	O 3
2	A	1	Total 19	C 12	H 2	N 2	O 3
2	A	1	Total 19	C 12	H 2	N 2	O 3

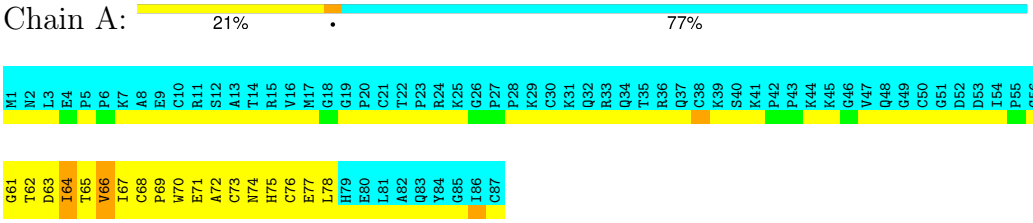
4 Residue-property plots

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide 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: Retinal rod rhodopsin-sensitive cGMP 3',5'-cyclic phosphodiesterase subunit gamma

Chain A:

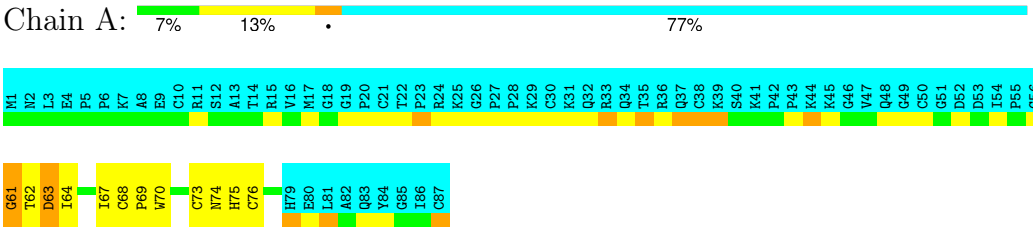


4.2 Residue scores for the representative (medoid) model from the NMR ensemble

The representative model is number 6. Colouring as in section 4.1 above.

• Molecule 1: Retinal rod rhodopsin-sensitive cGMP 3',5'-cyclic phosphodiesterase subunit gamma

Chain A:



5 Refinement protocol and experimental data overview

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

Of the 200 calculated structures, 100 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
XPLOR-NIH	refinement	2.11.2

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	1
Total number of shifts	845
Number of shifts mapped to atoms	727
Number of unparsed shifts	0
Number of shifts with mapping errors	118
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	72%

6 Model quality

6.1 Standard geometry

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

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	147	130	128	57±23
2	A	170	20	152	105±32
All	All	31700	15000	27846	12091

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 203.

5 of 7677 unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:A:160:RCY:C1M	2:A:168:RCY:H1Z	1.67	1.18	9	1
1:A:70:TRP:CZ2	2:A:168:RCY:H1MA	1.66	1.25	59	1
1:A:70:TRP:CD1	2:A:176:RCY:H1CA	1.66	1.15	62	1
1:A:64:ILE:CD1	2:A:168:RCY:C1P	1.66	1.74	12	2
1:A:70:TRP:CB	2:A:121:RCY:H1V	1.65	1.17	30	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	20/87 (23%)	7±2 (36±10%)	8±2 (38±10%)	5±2 (26±9%)	0 1
All	All	2000/8700 (23%)	723 (36%)	754 (38%)	523 (26%)	0 1

5 of 18 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	66	VAL	57
1	A	64	ILE	48
1	A	65	THR	41
1	A	77	GLU	39
1	A	69	PRO	33

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	17/73 (23%)	14±1 (80±9%)	3±1 (20±9%)	3 33
All	All	1700/7300 (23%)	1356 (80%)	344 (20%)	3 33

5 of 15 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	68	CYS	48
1	A	70	TRP	41
1	A	73	CYS	32
1	A	62	THR	31
1	A	76	CYS	26

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 monosaccharides in this entry.

6.6 Ligand geometry (i)

10 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length is the number of standard deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
2	RCY	A	110	1	17,18,18	2.86±0.33	6±0 (35±0%)
2	RCY	A	138	1	17,18,18	2.88±0.30	6±0 (35±2%)
2	RCY	A	187	1	17,18,18	2.88±0.42	6±0 (35±0%)
2	RCY	A	168	1	17,18,18	2.94±1.01	6±1 (36±5%)
2	RCY	A	173	1	17,18,18	2.86±0.24	6±0 (35±1%)
2	RCY	A	160	1	17,18,18	2.88±0.37	6±0 (35±1%)
2	RCY	A	121	1	17,18,18	2.88±0.31	6±0 (35±1%)
2	RCY	A	150	1	17,18,18	2.87±0.36	6±0 (35±1%)
2	RCY	A	130	1	17,18,18	3.26±4.12	6±1 (35±5%)
2	RCY	A	176	1	17,18,18	2.91±0.60	6±0 (35±2%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond angle is the number of standard

deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
2	RCY	A	110	1	22,30,30	4.32±0.30	8±1 (36±2%)
2	RCY	A	138	1	22,30,30	7.44±10.41	9±2 (38±7%)
2	RCY	A	187	1	22,30,30	4.73±3.94	8±1 (36±3%)
2	RCY	A	168	1	22,30,30	4.87±3.94	8±1 (37±4%)
2	RCY	A	173	1	22,30,30	5.49±6.41	8±1 (37±3%)
2	RCY	A	160	1	22,30,30	5.93±7.62	8±1 (37±5%)
2	RCY	A	121	1	22,30,30	6.64±9.13	8±1 (37±5%)
2	RCY	A	150	1	22,30,30	5.08±5.31	8±1 (36±3%)
2	RCY	A	130	1	22,30,30	6.14±7.98	8±2 (37±7%)
2	RCY	A	176	1	22,30,30	6.74±9.35	8±2 (38±8%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	RCY	A	121	1	-	0±0,4,40,40	0±0,2,2,2
2	RCY	A	187	1	-	0±0,4,40,40	0±0,2,2,2
2	RCY	A	176	1	-	1±0,4,40,40	0±0,2,2,2
2	RCY	A	110	1	-	0±0,4,40,40	0±0,2,2,2
2	RCY	A	168	1	1±0,1,7,7	0±0,4,40,40	0±0,2,2,2
2	RCY	A	130	1	1±0,1,7,7	0±0,4,40,40	0±0,2,2,2
2	RCY	A	173	1	1±0,1,7,7	0±0,4,40,40	0±0,2,2,2
2	RCY	A	150	1	1±0,1,7,7	0±0,4,40,40	0±0,2,2,2
2	RCY	A	160	1	1±0,1,7,7	0±0,4,40,40	0±0,2,2,2
2	RCY	A	138	1	1±0,1,7,7	0±0,4,40,40	0±0,2,2,2

5 of 116 unique bond 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(Å)	Ideal(Å)	Models	
								Worst	Total
2	A	130	RCY	C1Q-N1R	105.07	3.31	1.39	79	100
2	A	130	RCY	C1U-N1R	104.36	5.00	1.46	79	98
2	A	130	RCY	C1P-N1R	84.86	2.94	1.39	79	100
2	A	130	RCY	C1M-C1U	42.38	2.86	1.53	79	99

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
2	A	168	RCY	C1S-C1Q	35.22	2.51	1.51	79	2

5 of 179 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
2	A	138	RCY	C1Q-N1R-C1P	141.30	23.53	112.48	7	100
2	A	121	RCY	C1Q-N1R-C1P	140.81	23.83	112.48	52	100
2	A	173	RCY	C1Q-N1R-C1P	140.81	23.84	112.48	84	100
2	A	176	RCY	C1Q-N1R-C1P	140.79	23.85	112.48	91	100
2	A	130	RCY	C1Q-N1R-C1P	140.78	23.86	112.48	15	100

5 of 10 unique chiral outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Models (Total)
2	A	160	RCY	C1U	70
2	A	168	RCY	C1U	69
2	A	130	RCY	C1U	61
2	A	187	RCY	C1U	57
2	A	176	RCY	C1U	56

5 of 20 unique torsion outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Models (Total)
2	A	110	RCY	C1X-C1U-N1R-C1Q	15
2	A	168	RCY	C1X-C1U-N1R-C1Q	14
2	A	160	RCY	C1X-C1U-N1R-C1P	13
2	A	160	RCY	C1X-C1U-N1R-C1Q	12
2	A	121	RCY	C1X-C1U-N1R-C1Q	11

All unique ring outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Models (Total)
2	A	130	RCY	C1L-C1P-C1Q-C1S-N1R	1
2	A	130	RCY	C1M-C1U-C1W-C1X-N1V	1

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 72% for the well-defined parts and 66% 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	845
Number of shifts mapped to atoms	727
Number of unparsed shifts	0
Number of shifts with mapping errors	118
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 118) occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	10	ILE	H	8.289	0.05	.
1	A	10	ILE	HA	4.166	0.05	.
1	A	10	ILE	HB	1.88	0.05	.
1	A	10	ILE	HD11	0.859	0.05	.
1	A	10	ILE	HD12	0.859	0.05	.
1	A	10	ILE	HD13	0.859	0.05	.
1	A	10	ILE	HG12	1.214	0.05	.
1	A	10	ILE	HG13	1.501	0.05	.
1	A	10	ILE	HG21	0.913	0.05	.
1	A	10	ILE	HG22	0.913	0.05	.
1	A	10	ILE	HG23	0.913	0.05	.
1	A	10	ILE	C	176.308	0.15	.
1	A	10	ILE	CA	61.279	0.15	.
1	A	10	ILE	CB	38.666	0.15	.

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	10	ILE	CD1	12.714	0.15	.
1	A	10	ILE	CG1	27.278	0.15	.
1	A	10	ILE	CG2	17.53	0.15	.
1	A	10	ILE	N	123.217	0.15	.
1	A	21	VAL	H	8.376	0.05	.
1	A	21	VAL	HA	4.167	0.05	.
1	A	21	VAL	HB	1.962	0.05	.
1	A	21	VAL	HG21	0.956	0.05	.
1	A	21	VAL	HG22	0.956	0.05	.
1	A	21	VAL	HG23	0.956	0.05	.
1	A	21	VAL	C	176.368	0.15	.
1	A	21	VAL	CA	62.522	0.15	.
1	A	21	VAL	CB	32.722	0.15	.
1	A	21	VAL	CG2	20.766	0.15	.
1	A	21	VAL	N	120.972	0.15	.
1	A	30	PHE	H	8.168	0.05	.
1	A	30	PHE	HA	4.651	0.05	.
1	A	30	PHE	HB2	3.016	0.05	.
1	A	30	PHE	HB3	3.112	0.05	.
1	A	30	PHE	HD2	7.222	0.05	.
1	A	30	PHE	C	175.376	0.15	.
1	A	30	PHE	CA	57.554	0.15	.
1	A	30	PHE	CB	39.988	0.15	.
1	A	30	PHE	N	121.014	0.15	.
1	A	38	PHE	H	8.373	0.05	.
1	A	38	PHE	HA	4.659	0.05	.
1	A	38	PHE	HB2	3.011	0.05	.
1	A	38	PHE	HB3	3.101	0.05	.
1	A	38	PHE	C	175.417	0.15	.
1	A	38	PHE	CA	57.615	0.15	.
1	A	38	PHE	CB	40.008	0.15	.
1	A	38	PHE	N	122.108	0.15	.
1	A	50	PHE	H	8.222	0.05	.
1	A	50	PHE	HA	4.632	0.05	.
1	A	50	PHE	HB2	3.015	0.05	.
1	A	50	PHE	HB3	3.185	0.05	.
1	A	50	PHE	C	176.416	0.15	.
1	A	50	PHE	CA	58.008	0.15	.
1	A	50	PHE	CB	39.865	0.15	.
1	A	50	PHE	N	119.885	0.15	.
1	A	60	LEU	H	8.124	0.05	.

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	60	LEU	HA	4.388	0.05	.
1	A	60	LEU	HB2	1.648	0.05	.
1	A	60	LEU	HB3	1.663	0.05	.
1	A	60	LEU	HD21	0.907	0.05	.
1	A	60	LEU	HD22	0.907	0.05	.
1	A	60	LEU	HD23	0.907	0.05	.
1	A	60	LEU	C	177.67	0.15	.
1	A	60	LEU	CA	55.332	0.15	.
1	A	60	LEU	CB	42.514	0.15	.
1	A	60	LEU	CD2	23.502	0.15	.
1	A	60	LEU	CG	25.049	0.15	.
1	A	60	LEU	N	121.371	0.15	.
1	A	68	ALA	H	8.239	0.05	.
1	A	68	ALA	HA	4.361	0.05	.
1	A	68	ALA	HB1	1.002	0.05	.
1	A	68	ALA	HB2	1.002	0.05	.
1	A	68	ALA	HB3	1.002	0.05	.
1	A	68	ALA	CA	50.582	0.15	.
1	A	68	ALA	CB	18.099	0.15	.
1	A	68	ALA	N	128.782	0.15	.
1	A	73	PHE	H	7.935	0.05	.
1	A	73	PHE	HA	4.539	0.05	.
1	A	73	PHE	HB2	3.129	0.05	.
1	A	73	PHE	HB3	3.023	0.05	.
1	A	73	PHE	HD1	7.231	0.05	.
1	A	73	PHE	HD2	7.231	0.05	.
1	A	73	PHE	C	175.668	0.15	.
1	A	73	PHE	CA	57.985	0.15	.
1	A	73	PHE	CB	39.356	0.15	.
1	A	73	PHE	N	117.759	0.15	.
1	A	76	LEU	H	8.166	0.05	.
1	A	76	LEU	HA	4.287	0.05	.
1	A	76	LEU	HB2	1.568	0.05	.
1	A	76	LEU	HB3	1.642	0.05	.
1	A	76	LEU	HD11	0.863	0.05	.
1	A	76	LEU	HD12	0.863	0.05	.
1	A	76	LEU	HD13	0.863	0.05	.
1	A	76	LEU	HD21	0.9	0.05	.
1	A	76	LEU	HD22	0.9	0.05	.
1	A	76	LEU	HD23	0.9	0.05	.
1	A	76	LEU	C	177.284	0.15	.

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	76	LEU	CA	55.995	0.15	.
1	A	76	LEU	CB	42.241	0.15	.
1	A	76	LEU	CD1	23.706	0.15	.
1	A	76	LEU	CD2	24.084	0.15	.
1	A	76	LEU	N	122.091	0.15	.
1	A	87	ILE	H	7.855	0.05	.
1	A	87	ILE	HA	4.426	0.05	.
1	A	87	ILE	HB	2.018	0.05	.
1	A	87	ILE	HD11	0.848	0.05	.
1	A	87	ILE	HD12	0.848	0.05	.
1	A	87	ILE	HD13	0.848	0.05	.
1	A	87	ILE	HG12	1.433	0.05	.
1	A	87	ILE	HG13	1.188	0.05	.
1	A	87	ILE	HG21	0.895	0.05	.
1	A	87	ILE	HG22	0.895	0.05	.
1	A	87	ILE	HG23	0.895	0.05	.
1	A	87	ILE	CA	66.943	0.15	.
1	A	87	ILE	CB	39.016	0.15	.
1	A	87	ILE	CD1	12.969	0.15	.
1	A	87	ILE	CG1	27.193	0.15	.
1	A	87	ILE	CG2	17.895	0.15	.
1	A	87	ILE	N	127.372	0.15	.

7.1.2 Chemical shift referencing (i)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction ± precision, *ppm*	Suggested action
$^{13}C_{\alpha}$	84	-0.04 ± 0.09	None needed (< 0.5 ppm)
$^{13}C_{\beta}$	74	0.26 ± 0.13	None needed (< 0.5 ppm)
$^{13}C'$	76	-0.07 ± 0.07	None needed (< 0.5 ppm)
^{15}N	76	-0.98 ± 0.36	Should be applied

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 72%, i.e. 171 atoms were assigned a chemical shift out of a possible 237. 0 out of 2 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^{1}H	^{13}C	^{15}N
Backbone	78/100 (78%)	31/41 (76%)	32/40 (80%)	15/19 (79%)
Sidechain	91/117 (78%)	61/77 (79%)	30/39 (77%)	0/1 (0%)
Aromatic	2/20 (10%)	2/10 (20%)	0/7 (0%)	0/3 (0%)
Overall	171/237 (72%)	94/128 (73%)	62/86 (72%)	15/23 (65%)

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.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: