



Full wwPDB NMR Structure Validation Report ⓘ

Jun 3, 2023 – 06:04 PM EDT

PDB ID : 2MT5
BMRB ID : 25149
Title : Isolated Ring domain
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Deposited on : 2014-08-13

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

<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 : 4.02b-467
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
BMRB Restraints Analysis : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.33

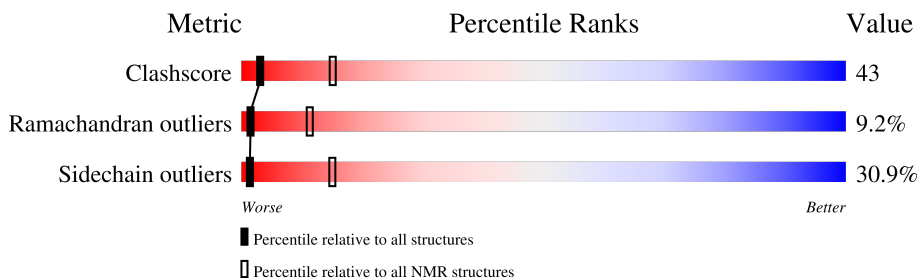
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 92%.

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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	70	

2 Ensemble composition and analysis

This entry contains 20 models. Model 1 is the overall representative, medoid model (most similar to other models).

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:7-A:21, A:30-A:70 (56)	0.56	1

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 2 clusters. No single-model clusters were found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1066 atoms, of which 511 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Anaphase-promoting complex subunit 11.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	70	1063	342	511	101	95	14	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP Q9NYG5
A	2	SER	-	expression tag	UNP Q9NYG5

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

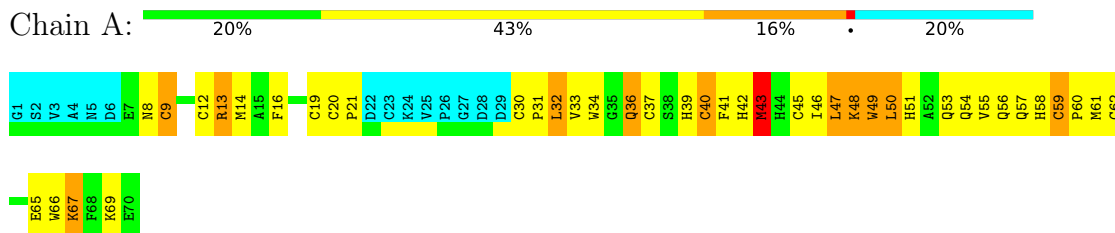
Mol	Chain	Residues	Atoms	
			Total	Zn
2	A	3	3	3

4 Residue-property plots [i](#)

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: Anaphase-promoting complex subunit 11

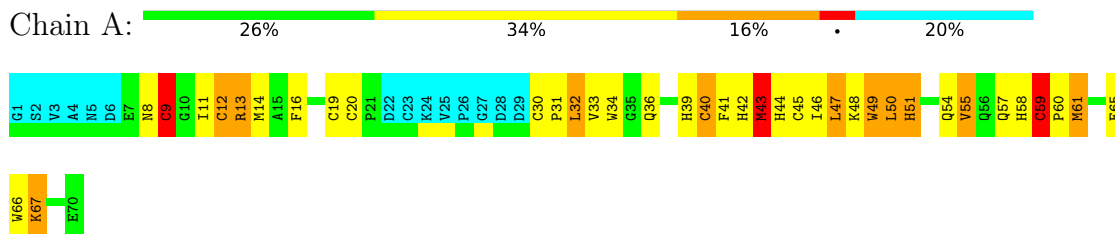


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

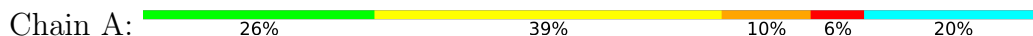
4.2.1 Score per residue for model 1 (medoid)

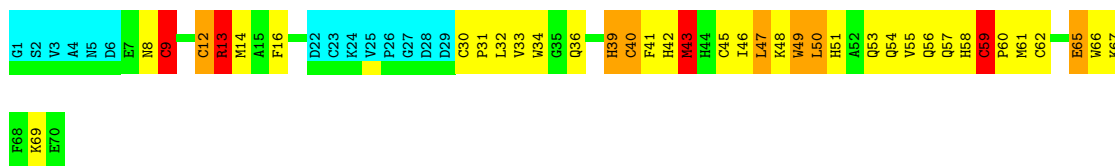
- Molecule 1: Anaphase-promoting complex subunit 11



4.2.2 Score per residue for model 2

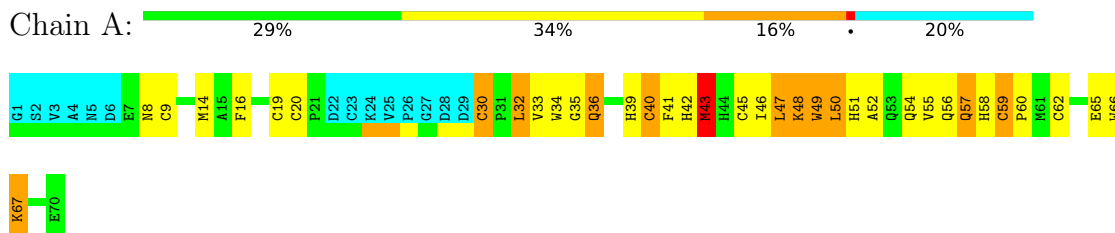
- Molecule 1: Anaphase-promoting complex subunit 11





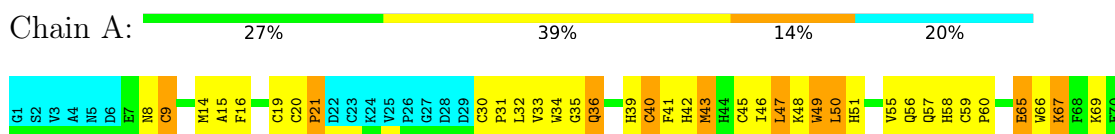
4.2.3 Score per residue for model 3

- Molecule 1: Anaphase-promoting complex subunit 11



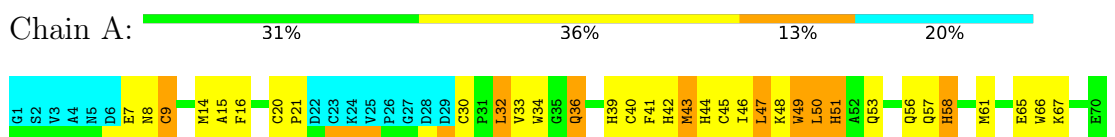
4.2.4 Score per residue for model 4

- Molecule 1: Anaphase-promoting complex subunit 11



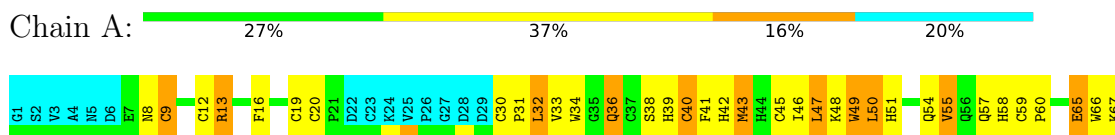
4.2.5 Score per residue for model 5

- Molecule 1: Anaphase-promoting complex subunit 11



4.2.6 Score per residue for model 6

- Molecule 1: Anaphase-promoting complex subunit 11



F68
K69
E70

4.2.7 Score per residue for model 7

- Molecule 1: Anaphase-promoting complex subunit 11

Chain A: 24% 37% 16% • 20%

G1 S2 V3 A4 N5 D6 E7 N8 C9 C12 R13 F16 C20 P21 D22 C23 K24 V25 P26 G27 D28 D29 L32 V33 W34 G35 C36 Q36 C37 S38 H39 C40 F41 F42 H43 H44 C45 I46 I47 K48 W49 L50 H51 A52 Q53 Q54 V55 Q56 Q57 H58 C59 P60 M61 E65 W66 K67

F68
K69
E70

4.2.8 Score per residue for model 8

- Molecule 1: Anaphase-promoting complex subunit 11

Chain A: 31% 33% 14% • 20%

G1 S2 V3 A4 M5 D6 C9 C12 R13 F16 C20 P21 D22 C23 K24 V25 P26 G27 D28 D29 L32 V33 W34 G35 Q36 C37 S38 H39 C40 F41 F42 H43 H44 C45 L47 K48 W49 L50 A52 V55 Q56 Q57 H58 C59 P60 M61 C62 W66 K67 K69

E70

4.2.9 Score per residue for model 9

- Molecule 1: Anaphase-promoting complex subunit 11

Chain A: 29% 34% 16% • 20%

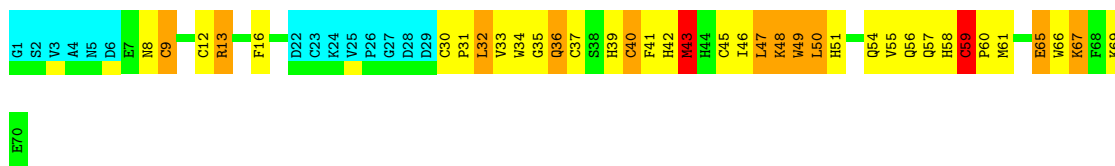
G1 S2 V3 A4 D6 E7 N8 C9 G10 I11 M14 A15 F16 C19 D22 C23 K24 V25 P26 G27 D28 D29 C30 P31 L32 V33 W34 G35 C36 S38 Q36 H39 C40 F41 F42 H43 M43 H44 C45 I46 I47 K48 W49 L50 H51 A52 Q53 Q54 V55 Q56 Q57 H58 C59 R63 Q64 E65 W66

K67
E70

4.2.10 Score per residue for model 10

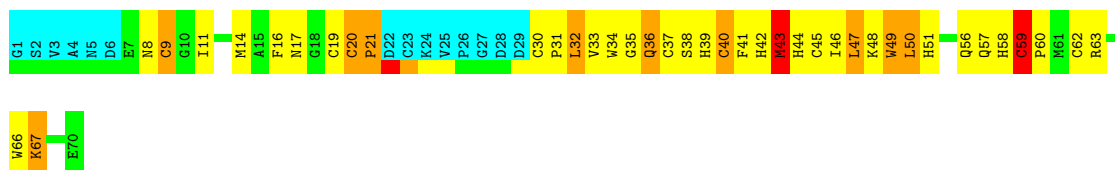
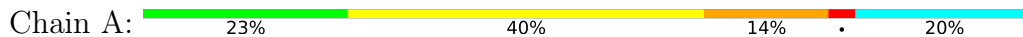
- Molecule 1: Anaphase-promoting complex subunit 11

Chain A: 27% 34% 16% • 20%



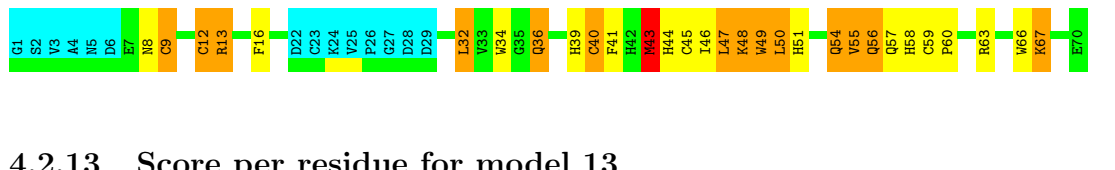
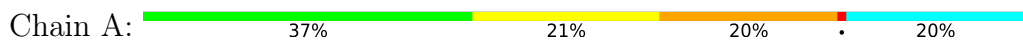
4.2.11 Score per residue for model 11

- Molecule 1: Anaphase-promoting complex subunit 11



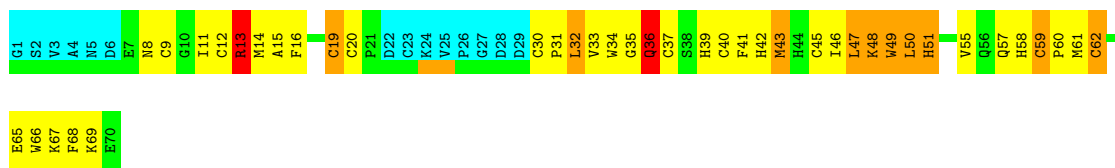
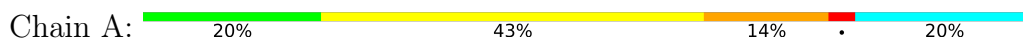
4.2.12 Score per residue for model 12

- Molecule 1: Anaphase-promoting complex subunit 11



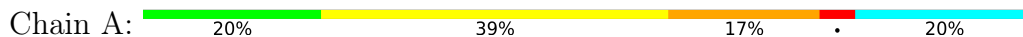
4.2.13 Score per residue for model 13

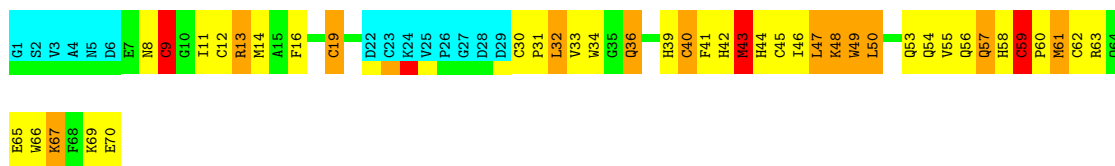
- Molecule 1: Anaphase-promoting complex subunit 11



4.2.14 Score per residue for model 14

- Molecule 1: Anaphase-promoting complex subunit 11

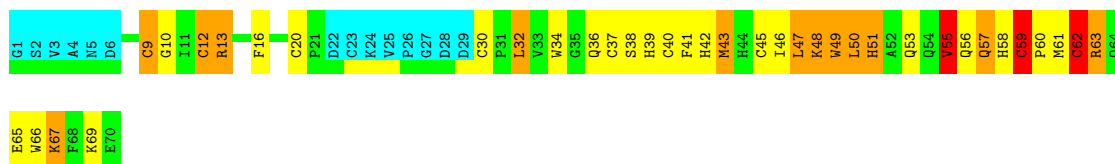




4.2.15 Score per residue for model 15

- Molecule 1: Anaphase-promoting complex subunit 11

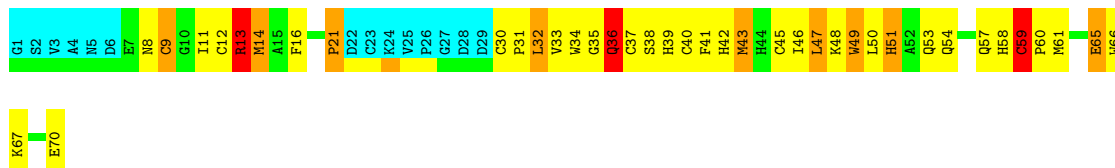
Chain A: 26% 31% 19% • 20%



4.2.16 Score per residue for model 16

- Molecule 1: Anaphase-promoting complex subunit 11

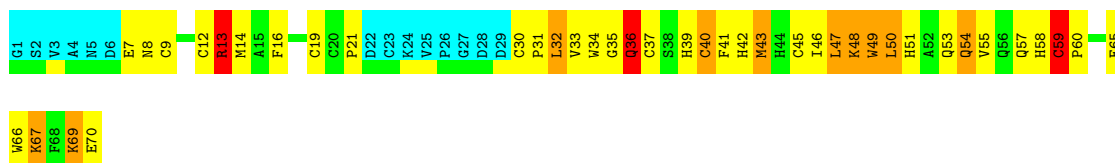
Chain A: 23% 40% 13% • 20%



4.2.17 Score per residue for model 17

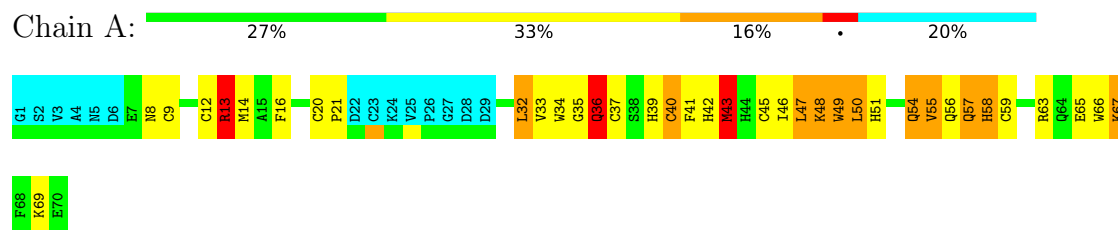
- Molecule 1: Anaphase-promoting complex subunit 11

Chain A: 21% 40% 14% • 20%



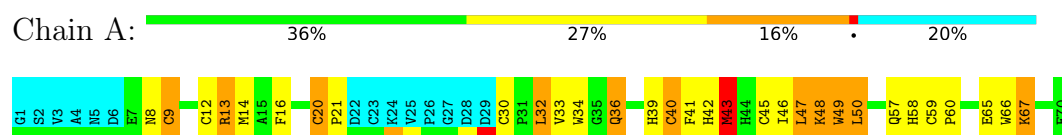
4.2.18 Score per residue for model 18

- Molecule 1: Anaphase-promoting complex subunit 11



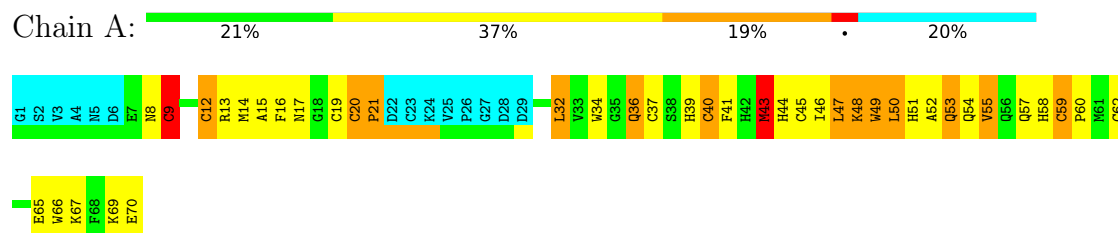
4.2.19 Score per residue for model 19

- Molecule 1: Anaphase-promoting complex subunit 11



4.2.20 Score per residue for model 20

- Molecule 1: Anaphase-promoting complex subunit 11



5 Refinement protocol and experimental data overview

The models were refined using the following method: *torsion angle dynamics*.

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *target function*.

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

Software name	Classification	Version
CYANA	structure solution	2.1
CYANA	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	1
Total number of shifts	837
Number of shifts mapped to atoms	837
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	92%

6 Model quality i

6.1 Standard geometry i

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 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	458	430	430	38±5
All	All	9220	8600	8600	763

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 43.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:32:LEU:HD11	1:A:34:TRP:CZ2	0.90	2.02	8	4
1:A:32:LEU:HD13	1:A:40:CYS:SG	0.84	2.13	16	1
1:A:32:LEU:HD11	1:A:34:TRP:CE2	0.79	2.11	8	7
1:A:39:HIS:CD2	1:A:41:PHE:CD1	0.76	2.73	16	18
1:A:32:LEU:HD21	1:A:34:TRP:CZ2	0.75	2.15	13	10
1:A:50:LEU:HD22	1:A:66:TRP:CD1	0.75	2.16	18	18
1:A:16:PHE:CD1	1:A:32:LEU:HD13	0.74	2.17	14	3
1:A:11:ILE:HD12	1:A:41:PHE:CD2	0.74	2.17	9	3
1:A:47:LEU:HD23	1:A:48:LYS:HD2	0.74	1.58	17	1
1:A:47:LEU:HD23	1:A:48:LYS:HE2	0.73	1.61	20	3
1:A:46:ILE:CD1	1:A:49:TRP:CE3	0.71	2.73	9	20
1:A:47:LEU:HD23	1:A:48:LYS:CE	0.70	2.16	14	10

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:47:LEU:HD23	1:A:48:LYS:HE3	0.70	1.63	14	6
1:A:32:LEU:CD1	1:A:34:TRP:CH2	0.70	2.75	15	4
1:A:47:LEU:HD23	1:A:48:LYS:CD	0.70	2.15	17	1
1:A:55:VAL:HG12	1:A:56:GLN:OE1	0.70	1.87	2	1
1:A:16:PHE:HB3	1:A:32:LEU:HD21	0.68	1.64	15	4
1:A:39:HIS:NE2	1:A:41:PHE:CD1	0.65	2.64	2	13
1:A:16:PHE:CD1	1:A:32:LEU:HD22	0.65	2.27	7	4
1:A:55:VAL:HG23	1:A:56:GLN:N	0.65	2.05	18	2
1:A:39:HIS:NE2	1:A:41:PHE:CD2	0.65	2.65	13	2
1:A:11:ILE:HD12	1:A:41:PHE:CE2	0.62	2.28	9	1
1:A:36:GLN:NE2	1:A:36:GLN:H	0.62	1.93	11	10
1:A:32:LEU:CD1	1:A:34:TRP:CZ2	0.61	2.81	8	5
1:A:16:PHE:HB3	1:A:32:LEU:HD22	0.61	1.71	13	4
1:A:49:TRP:CZ3	1:A:50:LEU:HB2	0.60	2.32	6	20
1:A:39:HIS:CD2	1:A:41:PHE:CE2	0.60	2.89	15	2
1:A:7:GLU:O	1:A:16:PHE:CE2	0.60	2.55	7	2
1:A:9:CYS:SG	1:A:41:PHE:CD2	0.60	2.95	11	12
1:A:16:PHE:CZ	1:A:40:CYS:SG	0.60	2.95	6	13
1:A:55:VAL:HG23	1:A:56:GLN:OE1	0.59	1.96	12	1
1:A:32:LEU:CD1	1:A:34:TRP:CE2	0.59	2.85	8	1
1:A:7:GLU:O	1:A:16:PHE:CD2	0.59	2.55	17	1
1:A:54:GLN:HE21	1:A:54:GLN:N	0.59	1.96	17	1
1:A:16:PHE:CE2	1:A:40:CYS:SG	0.58	2.96	11	12
1:A:46:ILE:HD12	1:A:49:TRP:CE3	0.58	2.32	16	20
1:A:46:ILE:HG23	1:A:47:LEU:N	0.58	2.13	2	20
1:A:16:PHE:CZ	1:A:40:CYS:O	0.58	2.57	2	3
1:A:32:LEU:HD12	1:A:33:VAL:N	0.58	2.14	10	8
1:A:50:LEU:HD22	1:A:66:TRP:HD1	0.58	1.54	18	4
1:A:66:TRP:CZ3	1:A:67:LYS:O	0.57	2.57	11	16
1:A:46:ILE:HD11	1:A:49:TRP:CZ3	0.57	2.35	9	20
1:A:34:TRP:CD1	1:A:40:CYS:SG	0.57	2.98	15	1
1:A:33:VAL:HG12	1:A:41:PHE:O	0.56	2.01	14	10
1:A:9:CYS:SG	1:A:42:HIS:CE1	0.56	2.98	14	2
1:A:30:CYS:HG	1:A:44:HIS:CE1	0.56	2.18	5	1
1:A:47:LEU:O	1:A:51:HIS:CG	0.55	2.60	9	4
1:A:54:GLN:NE2	1:A:55:VAL:HG13	0.55	2.16	6	2
1:A:30:CYS:N	1:A:31:PRO:CD	0.55	2.69	9	11
1:A:13:ARG:CD	1:A:13:ARG:N	0.55	2.70	6	2
1:A:20:CYS:SG	1:A:42:HIS:CG	0.55	3.00	5	1
1:A:14:MET:CG	1:A:42:HIS:NE2	0.54	2.71	17	3
1:A:39:HIS:CD2	1:A:61:MET:SD	0.54	3.00	14	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:40:CYS:O	1:A:41:PHE:CD1	0.54	2.60	11	2
1:A:30:CYS:SG	1:A:44:HIS:CE1	0.54	3.01	5	1
1:A:16:PHE:CB	1:A:32:LEU:HD21	0.53	2.33	15	1
1:A:16:PHE:CE1	1:A:40:CYS:O	0.52	2.62	16	6
1:A:46:ILE:HD11	1:A:49:TRP:CE3	0.52	2.39	9	20
1:A:41:PHE:CZ	1:A:61:MET:SD	0.52	3.03	13	1
1:A:32:LEU:HD12	1:A:34:TRP:CH2	0.52	2.40	16	2
1:A:57:GLN:HE22	1:A:66:TRP:N	0.52	2.02	9	2
1:A:32:LEU:HD12	1:A:33:VAL:O	0.51	2.05	11	4
1:A:13:ARG:CD	1:A:13:ARG:O	0.51	2.58	19	2
1:A:12:CYS:O	1:A:13:ARG:CD	0.51	2.59	17	5
1:A:10:GLY:N	1:A:41:PHE:CE1	0.51	2.79	15	1
1:A:52:ALA:HB3	1:A:53:GLN:NE2	0.51	2.19	20	1
1:A:47:LEU:O	1:A:51:HIS:CB	0.50	2.59	4	9
1:A:49:TRP:CZ2	1:A:60:PRO:HD3	0.50	2.42	6	16
1:A:34:TRP:CD1	1:A:40:CYS:HB3	0.50	2.42	13	12
1:A:65:GLU:O	1:A:65:GLU:CG	0.50	2.59	13	7
1:A:59:CYS:CB	1:A:62:CYS:SG	0.50	3.00	2	4
1:A:32:LEU:O	1:A:34:TRP:CZ3	0.50	2.65	9	2
1:A:12:CYS:O	1:A:13:ARG:C	0.50	2.49	10	15
1:A:43:MET:O	1:A:47:LEU:CB	0.49	2.60	13	4
1:A:58:HIS:O	1:A:59:CYS:C	0.49	2.51	10	17
1:A:36:GLN:NE2	1:A:36:GLN:N	0.49	2.60	5	8
1:A:16:PHE:CE1	1:A:32:LEU:HD13	0.49	2.42	14	1
1:A:53:GLN:CB	1:A:55:VAL:HG22	0.49	2.37	15	1
1:A:36:GLN:OE1	1:A:67:LYS:CB	0.49	2.61	5	4
1:A:47:LEU:HG	1:A:51:HIS:CD2	0.49	2.43	16	5
1:A:16:PHE:CD1	1:A:32:LEU:CD2	0.49	2.95	16	1
1:A:32:LEU:HG	1:A:34:TRP:CH2	0.49	2.43	14	11
1:A:46:ILE:HG23	1:A:47:LEU:H	0.48	1.68	14	19
1:A:54:GLN:N	1:A:54:GLN:NE2	0.48	2.61	17	1
1:A:34:TRP:CD1	1:A:40:CYS:HB2	0.48	2.43	15	2
1:A:20:CYS:SG	1:A:42:HIS:CD2	0.48	3.06	5	1
1:A:16:PHE:CE2	1:A:40:CYS:HB2	0.48	2.44	9	1
1:A:14:MET:CE	1:A:19:CYS:O	0.47	2.62	4	2
1:A:55:VAL:HG12	1:A:56:GLN:H	0.47	1.68	3	2
1:A:32:LEU:HG	1:A:34:TRP:CZ3	0.47	2.45	17	2
1:A:16:PHE:CZ	1:A:40:CYS:HB2	0.47	2.45	9	2
1:A:11:ILE:HD12	1:A:41:PHE:HD2	0.47	1.64	9	3
1:A:12:CYS:O	1:A:14:MET:N	0.46	2.49	19	5
1:A:33:VAL:HG23	1:A:43:MET:CA	0.46	2.40	8	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:43:MET:O	1:A:44:HIS:C	0.46	2.53	8	6
1:A:48:LYS:O	1:A:52:ALA:HB3	0.46	2.10	8	1
1:A:36:GLN:CD	1:A:36:GLN:N	0.46	2.69	16	4
1:A:32:LEU:HD13	1:A:34:TRP:CZ2	0.46	2.46	15	1
1:A:33:VAL:CG1	1:A:41:PHE:O	0.46	2.63	2	2
1:A:49:TRP:CZ2	1:A:57:GLN:O	0.46	2.69	8	1
1:A:53:GLN:HB3	1:A:55:VAL:HG22	0.46	1.88	7	1
1:A:16:PHE:CE1	1:A:40:CYS:HB2	0.46	2.46	11	1
1:A:12:CYS:O	1:A:13:ARG:NE	0.46	2.49	15	2
1:A:54:GLN:C	1:A:54:GLN:NE2	0.46	2.70	18	1
1:A:56:GLN:O	1:A:56:GLN:CG	0.45	2.64	7	1
1:A:47:LEU:O	1:A:51:HIS:ND1	0.45	2.49	9	4
1:A:11:ILE:O	1:A:13:ARG:NH2	0.45	2.49	1	1
1:A:49:TRP:CH2	1:A:57:GLN:O	0.45	2.69	8	1
1:A:39:HIS:CD2	1:A:40:CYS:N	0.45	2.84	13	1
1:A:7:GLU:CG	1:A:7:GLU:O	0.45	2.64	5	1
1:A:66:TRP:CE3	1:A:67:LYS:N	0.45	2.85	11	20
1:A:14:MET:SD	1:A:15:ALA:N	0.45	2.90	20	5
1:A:51:HIS:ND1	1:A:51:HIS:O	0.45	2.50	15	7
1:A:36:GLN:N	1:A:36:GLN:HE21	0.45	2.10	13	1
1:A:19:CYS:SG	1:A:20:CYS:N	0.45	2.90	13	1
1:A:56:GLN:NE2	1:A:56:GLN:O	0.45	2.50	5	1
1:A:53:GLN:CA	1:A:53:GLN:OE1	0.45	2.65	9	1
1:A:35:GLY:O	1:A:37:CYS:N	0.44	2.50	13	8
1:A:53:GLN:O	1:A:54:GLN:NE2	0.44	2.50	16	1
1:A:46:ILE:CG2	1:A:47:LEU:N	0.44	2.79	10	20
1:A:12:CYS:C	1:A:13:ARG:CD	0.44	2.85	13	1
1:A:47:LEU:CD2	1:A:48:LYS:HE2	0.44	2.42	13	1
1:A:49:TRP:O	1:A:53:GLN:NE2	0.44	2.50	20	1
1:A:12:CYS:SG	1:A:14:MET:CB	0.44	3.06	18	3
1:A:42:HIS:O	1:A:43:MET:C	0.44	2.56	5	17
1:A:34:TRP:NE1	1:A:40:CYS:SG	0.44	2.90	15	2
1:A:49:TRP:CH2	1:A:60:PRO:HD3	0.44	2.48	13	2
1:A:9:CYS:SG	1:A:41:PHE:CG	0.43	3.11	2	1
1:A:48:LYS:HA	1:A:48:LYS:CE	0.43	2.43	17	1
1:A:20:CYS:SG	1:A:30:CYS:CB	0.43	3.06	19	1
1:A:46:ILE:O	1:A:49:TRP:N	0.43	2.51	11	3
1:A:11:ILE:HG22	1:A:12:CYS:SG	0.43	2.54	13	1
1:A:52:ALA:O	1:A:54:GLN:NE2	0.43	2.51	3	1
1:A:32:LEU:CD1	1:A:40:CYS:HB2	0.43	2.44	2	1
1:A:39:HIS:NE2	1:A:41:PHE:CE2	0.43	2.86	15	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:54:GLN:NE2	1:A:54:GLN:N	0.43	2.67	3	1
1:A:19:CYS:O	1:A:20:CYS:C	0.43	2.57	11	3
1:A:43:MET:CG	1:A:44:HIS:N	0.43	2.81	20	1
1:A:11:ILE:O	1:A:13:ARG:CZ	0.42	2.67	1	1
1:A:53:GLN:N	1:A:53:GLN:CD	0.42	2.71	14	1
1:A:16:PHE:CZ	1:A:40:CYS:HB3	0.42	2.49	16	1
1:A:69:LYS:O	1:A:70:GLU:C	0.42	2.58	14	3
1:A:33:VAL:HG11	1:A:68:PHE:HD1	0.42	1.74	13	1
1:A:47:LEU:CD2	1:A:48:LYS:CE	0.42	2.95	14	1
1:A:32:LEU:CD2	1:A:34:TRP:CZ2	0.42	3.01	3	1
1:A:35:GLY:O	1:A:36:GLN:C	0.42	2.58	17	8
1:A:61:MET:O	1:A:63:ARG:N	0.42	2.52	15	1
1:A:49:TRP:CZ2	1:A:58:HIS:O	0.42	2.73	18	3
1:A:65:GLU:O	1:A:66:TRP:C	0.42	2.59	3	13
1:A:34:TRP:CE2	1:A:40:CYS:SG	0.41	3.07	15	1
1:A:54:GLN:NE2	1:A:54:GLN:H	0.41	2.13	12	1
1:A:38:SER:O	1:A:39:HIS:C	0.41	2.59	16	4
1:A:39:HIS:CD2	1:A:41:PHE:CE1	0.41	3.08	2	1
1:A:36:GLN:NE2	1:A:67:LYS:N	0.41	2.69	9	1
1:A:49:TRP:CH2	1:A:58:HIS:O	0.41	2.74	18	3
1:A:65:GLU:O	1:A:65:GLU:CD	0.41	2.59	16	2
1:A:57:GLN:CD	1:A:57:GLN:C	0.41	2.79	14	2
1:A:65:GLU:OE1	1:A:65:GLU:N	0.41	2.54	16	1
1:A:57:GLN:O	1:A:58:HIS:O	0.41	2.39	18	2
1:A:14:MET:SD	1:A:19:CYS:O	0.40	2.79	1	2
1:A:35:GLY:C	1:A:69:LYS:CB	0.40	2.89	10	1
1:A:30:CYS:N	1:A:31:PRO:HD3	0.40	2.31	9	2
1:A:13:ARG:N	1:A:13:ARG:HD2	0.40	2.32	6	1
1:A:33:VAL:CG2	1:A:43:MET:HB2	0.40	2.46	13	1
1:A:59:CYS:SG	1:A:62:CYS:SG	0.40	3.20	2	1
1:A:16:PHE:HD1	1:A:32:LEU:HD22	0.40	1.74	8	1
1:A:58:HIS:O	1:A:59:CYS:O	0.40	2.39	10	1
1:A:32:LEU:HD21	1:A:34:TRP:CH2	0.40	2.51	11	1
1:A:33:VAL:HG22	1:A:34:TRP:N	0.40	2.32	11	1
1:A:59:CYS:HG	1:A:62:CYS:C	0.40	2.20	13	1
1:A:56:GLN:O	1:A:56:GLN:OE1	0.40	2.40	7	1
1:A:37:CYS:SG	1:A:62:CYS:SG	0.40	3.19	15	1
1:A:57:GLN:O	1:A:58:HIS:C	0.40	2.60	3	1
1:A:51:HIS:O	1:A:51:HIS:ND1	0.40	2.54	10	1
1:A:37:CYS:O	1:A:39:HIS:N	0.40	2.54	20	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	55/70 (79%)	37±2 (67±4%)	13±2 (24±4%)	5±1 (9±2%)	1	11
All	All	1100/1400 (79%)	735 (67%)	264 (24%)	101 (9%)	1	11

All 15 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	45	CYS	20
1	A	9	CYS	17
1	A	43	MET	12
1	A	59	CYS	11
1	A	21	PRO	9
1	A	13	ARG	7
1	A	55	VAL	6
1	A	56	GLN	4
1	A	36	GLN	4
1	A	58	HIS	3
1	A	69	LYS	3
1	A	19	CYS	2
1	A	30	CYS	1
1	A	53	GLN	1
1	A	62	CYS	1

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	51/62 (82%)	35±3 (69±5%)	16±3 (31±5%)	1	15
All	All	1020/1240 (82%)	705 (69%)	315 (31%)	1	15

All 34 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	36	GLN	20
1	A	43	MET	20
1	A	47	LEU	20
1	A	48	LYS	20
1	A	49	TRP	20
1	A	57	GLN	20
1	A	50	LEU	19
1	A	8	ASN	18
1	A	32	LEU	15
1	A	40	CYS	15
1	A	13	ARG	12
1	A	67	LYS	12
1	A	59	CYS	11
1	A	55	VAL	10
1	A	54	GLN	8
1	A	51	HIS	7
1	A	61	MET	7
1	A	53	GLN	6
1	A	69	LYS	6
1	A	20	CYS	6
1	A	63	ARG	6
1	A	9	CYS	5
1	A	12	CYS	5
1	A	65	GLU	5
1	A	56	GLN	4
1	A	62	CYS	4
1	A	70	GLU	3
1	A	19	CYS	3
1	A	14	MET	2
1	A	17	ASN	2
1	A	39	HIS	1
1	A	7	GLU	1
1	A	37	CYS	1
1	A	11	ILE	1

6.3.3 RNA

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](#)

Of 3 ligands modelled in this entry, 3 are 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 [i](#)

The completeness of assignment taking into account all chemical shift lists is 92% for the well-defined parts and 92% 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 [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	837
Number of shifts mapped to atoms	837
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	6

7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	70	-0.27 ± 0.39	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	64	-0.19 ± 0.20	None needed (< 0.5 ppm)
$^{13}\text{C}'$	64	-0.02 ± 0.34	None needed (< 0.5 ppm)
^{15}N	64	0.54 ± 0.43	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 92%, i.e. 706 atoms were assigned a chemical shift out of a possible 764. 0 out of 5 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	274/277 (99%)	112/112 (100%)	109/112 (97%)	53/53 (100%)
Sidechain	352/386 (91%)	240/250 (96%)	104/119 (87%)	8/17 (47%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	80/101 (79%)	43/53 (81%)	34/40 (85%)	3/8 (38%)
Overall	706/764 (92%)	395/415 (95%)	247/271 (91%)	64/78 (82%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 92%, i.e. 837 atoms were assigned a chemical shift out of a possible 909. 0 out of 7 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	337/347 (97%)	139/141 (99%)	134/140 (96%)	64/66 (97%)
Sidechain	420/461 (91%)	287/297 (97%)	124/145 (86%)	9/19 (47%)
Aromatic	80/101 (79%)	43/53 (81%)	34/40 (85%)	3/8 (38%)
Overall	837/909 (92%)	469/491 (96%)	292/325 (90%)	76/93 (82%)

7.1.4 Statistically unusual chemical shifts [i](#)

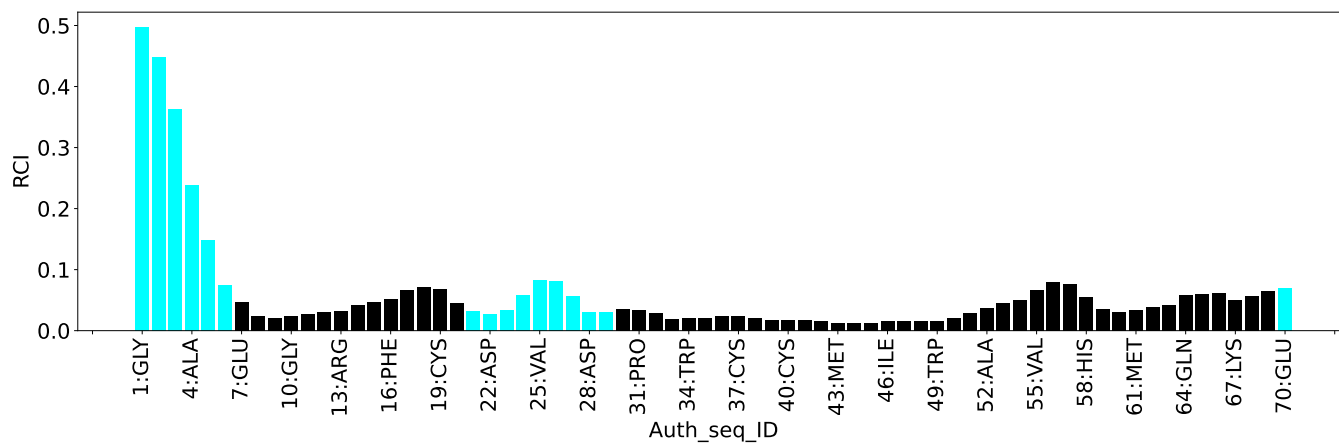
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	60	PRO	HG3	-0.37	0.33 – 3.48	-7.2
1	A	66	TRP	HE1	5.83	6.88 – 13.28	-6.7
1	A	60	PRO	HB2	0.24	0.37 – 3.78	-5.4
1	A	46	ILE	HD11	-0.81	-0.72 – 2.09	-5.3
1	A	46	ILE	HD12	-0.81	-0.72 – 2.09	-5.3
1	A	46	ILE	HD13	-0.81	-0.72 – 2.09	-5.3

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:



8 NMR restraints analysis

8.1 Conformationally restricting restraints

The following table provides the summary of experimentally observed NMR restraints in different categories. Restraints are classified into different categories based on the sequence separation of the atoms involved.

Description	Value
Total distance restraints	998
Intra-residue ($ i-j =0$)	392
Sequential ($ i-j =1$)	250
Medium range ($ i-j >1$ and $ i-j <5$)	115
Long range ($ i-j \geq 5$)	235
Inter-chain	0
Hydrogen bond restraints	0
Disulfide bond restraints	6
Total dihedral-angle restraints	0
Number of unmapped restraints	9
Number of restraints per residue	14.3
Number of long range restraints per residue ¹	3.4

¹Long range hydrogen bonds and disulfide bonds are counted as long range restraints while calculating the number of long range restraints per residue

8.2 Residual restraint violations

This section provides the overview of the restraint violations analysis. The violations are binned as small, medium and large violations based on its absolute value. Average number of violations per model is calculated by dividing the total number of violations in each bin by the size of the ensemble.

8.2.1 Average number of distance violations per model

Distance violations less than 0.1 Å are not included in the calculation.

Bins (Å)	Average number of violations per model	Max (Å)
0.1-0.2 (Small)	0.6	0.18
0.2-0.5 (Medium)	None	None
>0.5 (Large)	None	None

8.2.2 Average number of dihedral-angle violations per model [i](#)

Dihedral-angle violations less than 1° are not included in the calculation. There are no dihedral-angle violations

9 Distance violation analysis [i](#)

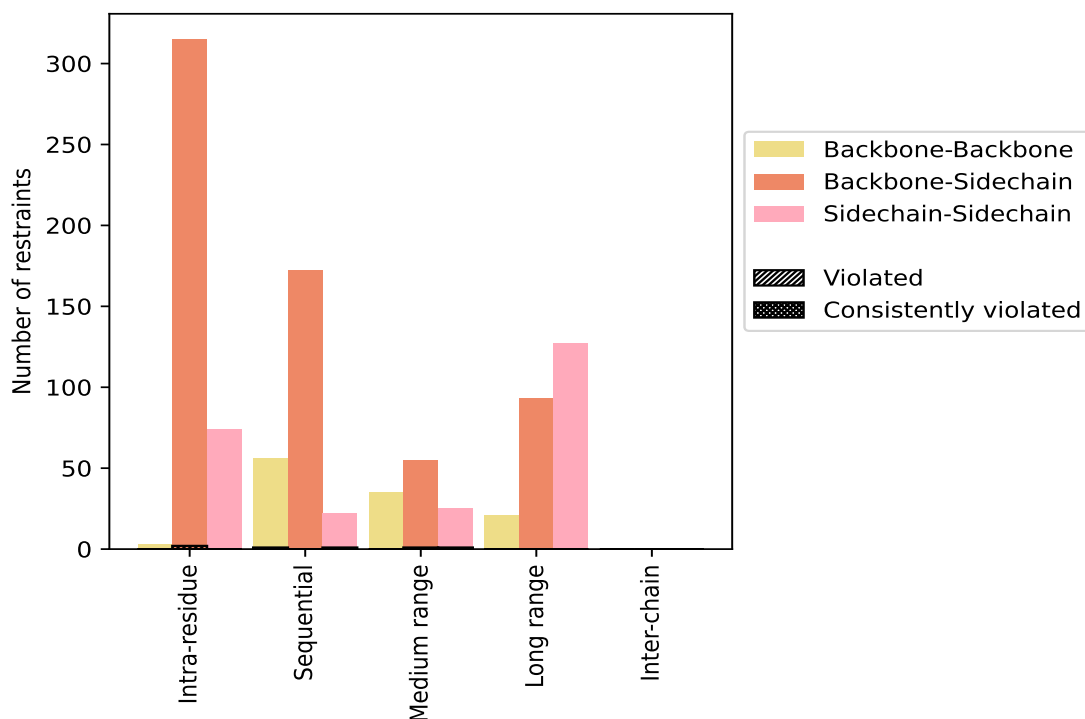
9.1 Summary of distance violations [i](#)

The following table shows the summary of distance violations in different restraint categories based on the sequence separation of the atoms involved. Each category is further sub-divided into three sub-categories based on the atoms involved. Violations less than 0.1 Å are not included in the statistics.

Restrains type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue ($i-j =0$)	392	39.3	2	0.5	0.2	0	0.0	0.0
Backbone-Backbone	3	0.3	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	315	31.6	2	0.6	0.2	0	0.0	0.0
Sidechain-Sidechain	74	7.4	0	0.0	0.0	0	0.0	0.0
Sequential ($i-j =1$)	250	25.1	2	0.8	0.2	0	0.0	0.0
Backbone-Backbone	56	5.6	1	1.8	0.1	0	0.0	0.0
Backbone-Sidechain	172	17.2	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	22	2.2	1	4.5	0.1	0	0.0	0.0
Medium range ($i-j >1$ & $i-j <5$)	115	11.5	2	1.7	0.2	0	0.0	0.0
Backbone-Backbone	35	3.5	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	55	5.5	1	1.8	0.1	0	0.0	0.0
Sidechain-Sidechain	25	2.5	1	4.0	0.1	0	0.0	0.0
Long range ($i-j \geq 5$)	235	23.5	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	21	2.1	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	93	9.3	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	121	12.1	0	0.0	0.0	0	0.0	0.0
Inter-chain	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Hydrogen bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Disulfide bond	6	0.6	0	0.0	0.0	0	0.0	0.0
Total	998	100.0	6	0.6	0.6	0	0.0	0.0
Backbone-Backbone	115	11.5	1	0.9	0.1	0	0.0	0.0
Backbone-Sidechain	635	63.6	3	0.5	0.3	0	0.0	0.0
Sidechain-Sidechain	248	24.8	2	0.8	0.2	0	0.0	0.0

¹ percentage calculated with respect to the total number of distance restraints, ² percentage calculated with respect to the number of restraints in a particular restraint category, ³ violated in at least one model, ⁴ violated in all the models

9.1.1 Bar chart : Distribution of distance restraints and violations [i](#)



Violated and consistently violated restraints are shown using different hatch patterns in their respective categories. The hydrogen bonds and disulfied bonds are counted in their appropriate category on the x-axis

9.2 Distance violation statistics for each model [i](#)

The following table provides the distance violation statistics for each model in the ensemble. Violations less than 0.1 Å are not included in the statistics.

Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
1	0	0	0	0	0	0	0.0	0.0	0.0	0.0
2	0	0	0	0	0	0	0.0	0.0	0.0	0.0
3	0	0	0	0	0	0	0.0	0.0	0.0	0.0
4	0	0	0	0	0	0	0.0	0.0	0.0	0.0
5	0	0	0	0	0	0	0.0	0.0	0.0	0.0
6	1	0	0	0	0	1	0.14	0.14	0.0	0.14
7	1	0	0	0	0	1	0.14	0.14	0.0	0.14
8	0	0	0	0	0	0	0.0	0.0	0.0	0.0
9	1	0	1	0	0	2	0.14	0.18	0.03	0.14
10	1	0	0	0	0	1	0.13	0.13	0.0	0.13
11	0	0	0	0	0	0	0.0	0.0	0.0	0.0

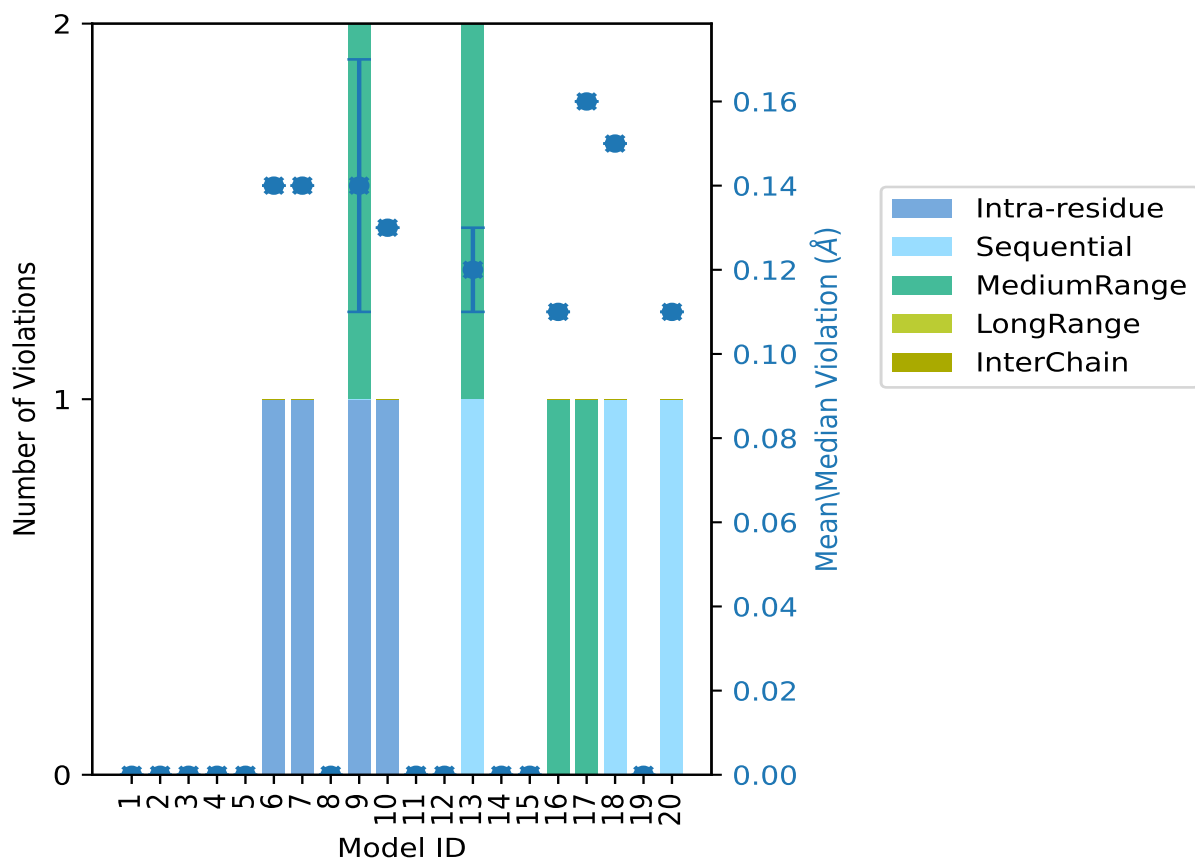
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Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
12	0	0	0	0	0	0	0.0	0.0	0.0	0.0
13	0	1	1	0	0	2	0.12	0.13	0.01	0.12
14	0	0	0	0	0	0	0.0	0.0	0.0	0.0
15	0	0	0	0	0	0	0.0	0.0	0.0	0.0
16	0	0	1	0	0	1	0.11	0.11	0.0	0.11
17	0	0	1	0	0	1	0.16	0.16	0.0	0.16
18	0	1	0	0	0	1	0.15	0.15	0.0	0.15
19	0	0	0	0	0	0	0.0	0.0	0.0	0.0
20	0	1	0	0	0	1	0.11	0.11	0.0	0.11

¹Intra-residue restraints, ²Sequential restraints, ³Medium range restraints, ⁴Long range restraints, ⁵Inter-chain restraints, ⁶Standard deviation

9.2.1 Bar graph : Distance Violation statistics for each model [\(i\)](#)



The mean(dot),median(x) and the standard deviation are shown in blue with respect to the y axis on the right

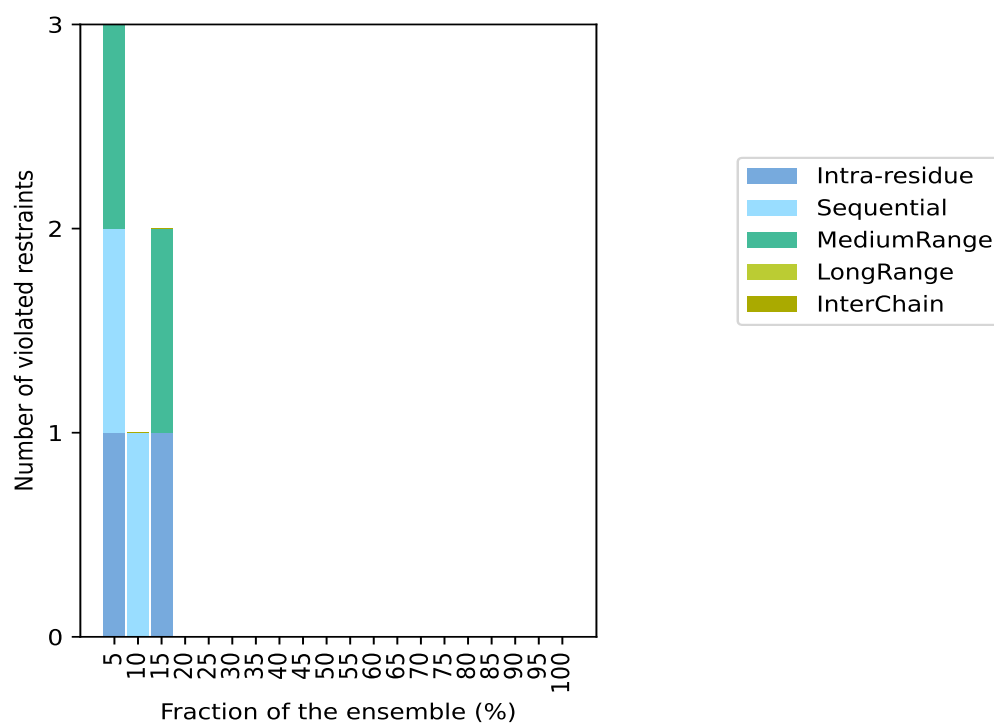
9.3 Distance violation statistics for the ensemble

Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 986(IR:390, SQ:248, MR:113, LR:235, IC:0) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total	Count ⁶	%
1	1	1	0	0	3	1	5.0
0	1	0	0	0	1	2	10.0
1	0	1	0	0	2	3	15.0
0	0	0	0	0	0	4	20.0
0	0	0	0	0	0	5	25.0
0	0	0	0	0	0	6	30.0
0	0	0	0	0	0	7	35.0
0	0	0	0	0	0	8	40.0
0	0	0	0	0	0	9	45.0
0	0	0	0	0	0	10	50.0
0	0	0	0	0	0	11	55.0
0	0	0	0	0	0	12	60.0
0	0	0	0	0	0	13	65.0
0	0	0	0	0	0	14	70.0
0	0	0	0	0	0	15	75.0
0	0	0	0	0	0	16	80.0
0	0	0	0	0	0	17	85.0
0	0	0	0	0	0	18	90.0
0	0	0	0	0	0	19	95.0
0	0	0	0	0	0	20	100.0

¹Intra-residue restraints, ²Sequential restraints, ³Medium range restraints, ⁴Long range restraints, ⁵Inter-chain restraints, ⁶ Number of models with violations

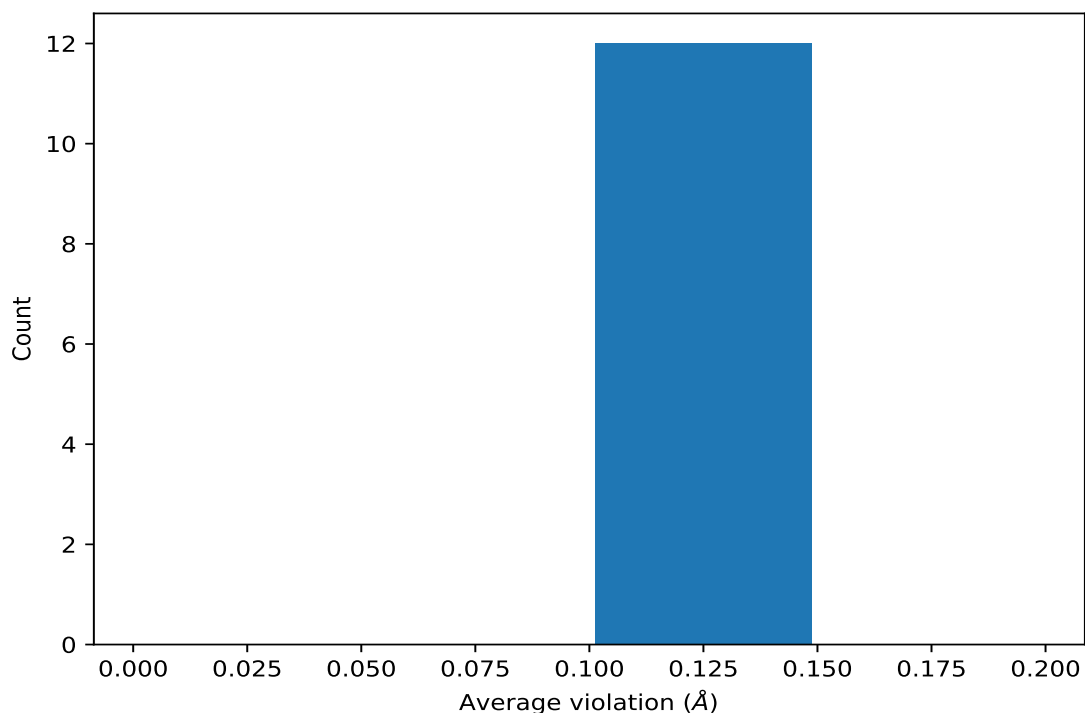
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



9.4 Most violated distance restraints in the ensemble [i](#)

9.4.1 Histogram : Distribution of mean distance violations [i](#)

The following histogram shows the distribution of the average value of the violation. The average is calculated for each restraint that is violated in more than one model over all the violated models in the ensemble



9.4.2 Table: Most violated distance restraints [i](#)

The following table provides the mean and the standard deviation of the violation for each restraint sorted by number of violated models and the mean value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

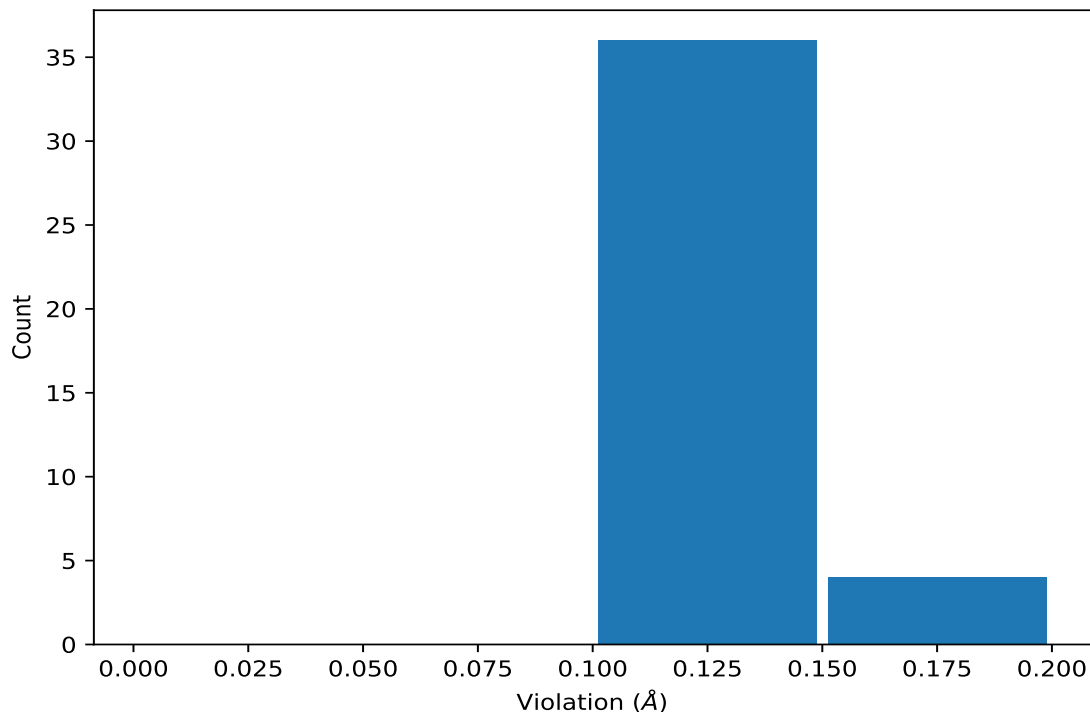
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,827)	1:A:14:MET:H	1:A:14:MET:HG2	3	0.14	0.0	0.14
(1,827)	1:A:14:MET:H	1:A:14:MET:HG3	3	0.14	0.0	0.14
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD21	3	0.11	0.0	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD22	3	0.11	0.0	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD23	3	0.11	0.0	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD21	3	0.11	0.0	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD22	3	0.11	0.0	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD23	3	0.11	0.0	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD21	3	0.11	0.0	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD22	3	0.11	0.0	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD23	3	0.11	0.0	0.11
(1,247)	1:A:49:TRP:HE1	1:A:50:LEU:HB2	2	0.14	0.01	0.14

¹Number of violated models, ²Standard deviation

9.5 All violated distance restraints [i](#)

9.5.1 Histogram : Distribution of distance violations [i](#)

The following histogram shows the distribution of the absolute value of the violation for all violated restraints in the ensemble.



9.5.2 Table : All distance violations [i](#)

The following table lists the absolute value of the violation for each restraint in the ensemble sorted by its value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,358)	1:A:7:GLU:H	1:A:7:GLU:HB3	9	0.18
(1,611)	1:A:47:LEU:HD21	1:A:50:LEU:H	17	0.16
(1,611)	1:A:47:LEU:HD22	1:A:50:LEU:H	17	0.16
(1,611)	1:A:47:LEU:HD23	1:A:50:LEU:H	17	0.16
(1,247)	1:A:49:TRP:HE1	1:A:50:LEU:HB2	18	0.15
(1,827)	1:A:14:MET:H	1:A:14:MET:HG2	6	0.14
(1,827)	1:A:14:MET:H	1:A:14:MET:HG3	6	0.14
(1,827)	1:A:14:MET:H	1:A:14:MET:HG2	7	0.14
(1,827)	1:A:14:MET:H	1:A:14:MET:HG3	7	0.14
(1,827)	1:A:14:MET:H	1:A:14:MET:HG2	10	0.13

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,827)	1:A:14:MET:H	1:A:14:MET:HG3	10	0.13
(1,247)	1:A:49:TRP:HE1	1:A:50:LEU:HB2	13	0.13
(1,664)	1:A:23:CYS:HA	1:A:24:LYS:H	20	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD21	9	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD22	9	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD23	9	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD21	9	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD22	9	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD23	9	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD21	9	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD22	9	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD23	9	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD21	13	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD22	13	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD23	13	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD21	13	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD22	13	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD23	13	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD21	13	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD22	13	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD23	13	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD21	16	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD22	16	0.11
(1,552)	1:A:46:ILE:HD11	1:A:50:LEU:HD23	16	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD21	16	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD22	16	0.11
(1,552)	1:A:46:ILE:HD12	1:A:50:LEU:HD23	16	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD21	16	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD22	16	0.11
(1,552)	1:A:46:ILE:HD13	1:A:50:LEU:HD23	16	0.11

10 Dihedral-angle violation analysis

Dihedral angle analysis failed due to data error in the dihedral angle restraints, possibly missing target value