



Full wwPDB NMR Structure Validation Report ⓘ

Mar 5, 2026 – 01:49 PM UTC

PDB ID : 1N3G / pdb_00001n3g
Title : Solution structure of the ribosome-associated cold shock response protein Yfia of Escherichia coli
Authors : Rak, A.; Kalinin, A.; Shcherbakov, D.; Bayer, P.
Deposited on : 2002-10-28

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	:	4-5-2 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

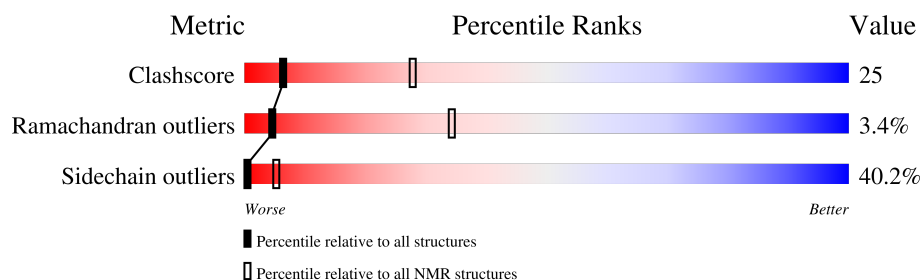
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

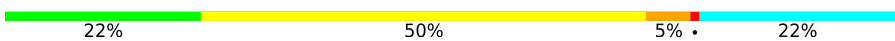
The overall completeness of chemical shifts assignment was not calculated.

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

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	113	

2 Ensemble composition and analysis

This entry contains 29 models. Model 14 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:88 (88)	0.77	14

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 5 clusters and 1 single-model cluster was found.

Cluster number	Models
1	2, 3, 4, 5, 10, 11, 12, 13, 14, 18, 19, 20, 21, 24, 25, 29
2	7, 8, 16, 17, 23, 26
3	9, 28
4	22, 27
5	1, 6
Single-model clusters	15

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 1810 atoms, of which 913 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Protein yfiA.

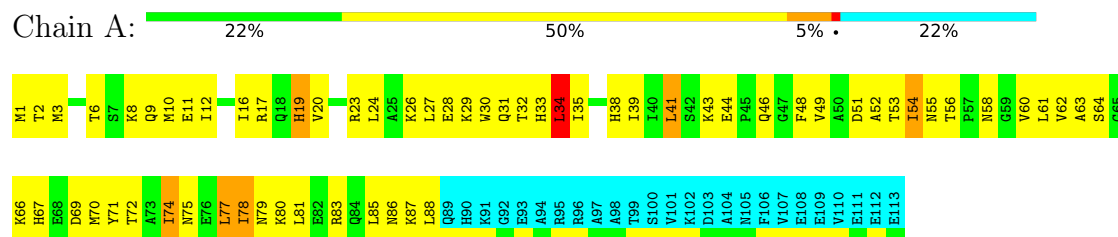
Mol	Chain	Residues	Atoms						Trace
1	A	113	Total	C	H	N	O	S	0
			1810	558	913	162	173	4	

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: Protein yfIA

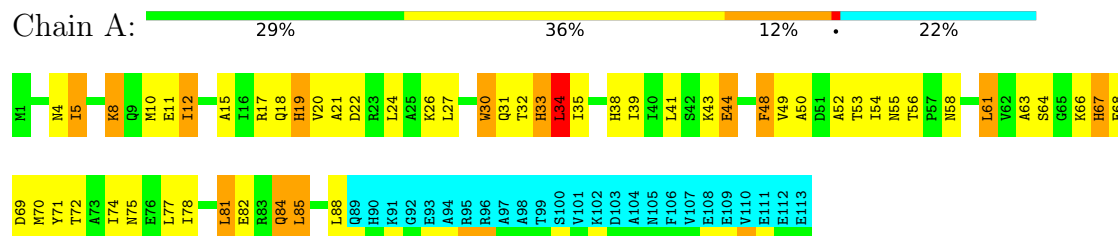


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

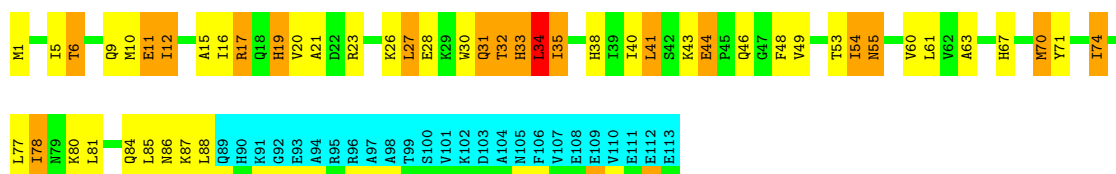
- Molecule 1: Protein yfIA



4.2.2 Score per residue for model 2

- Molecule 1: Protein yfIA





4.2.3 Score per residue for model 3

- Molecule 1: Protein yfiA

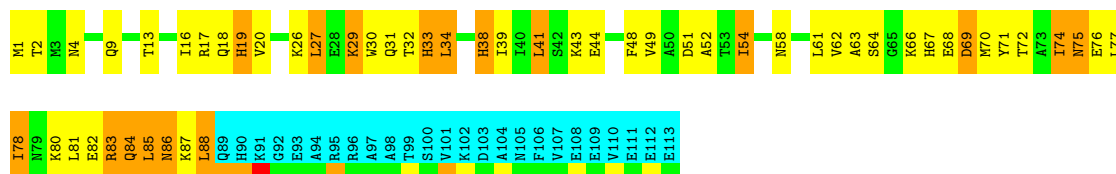
Chain A: 33% 30% 15% 22%



4.2.4 Score per residue for model 4

- Molecule 1: Protein yfiA

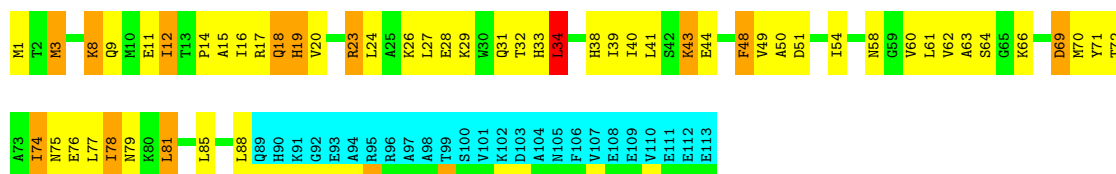
Chain A: 30% 33% 15% 22%



4.2.5 Score per residue for model 5

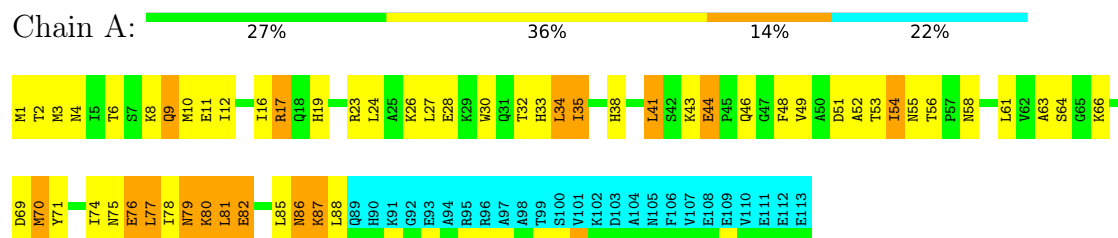
- Molecule 1: Protein yfiA

Chain A: 30% 36% 11% 22%



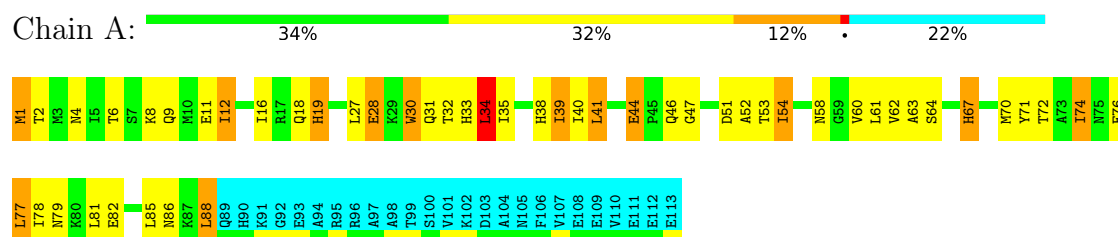
4.2.6 Score per residue for model 6

- Molecule 1: Protein yfiA



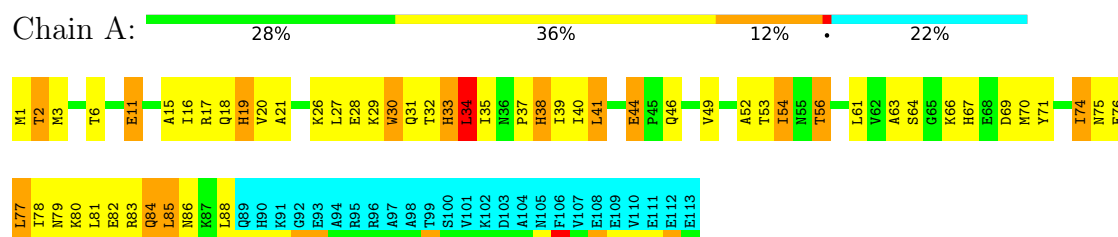
4.2.7 Score per residue for model 7

- Molecule 1: Protein yfiA



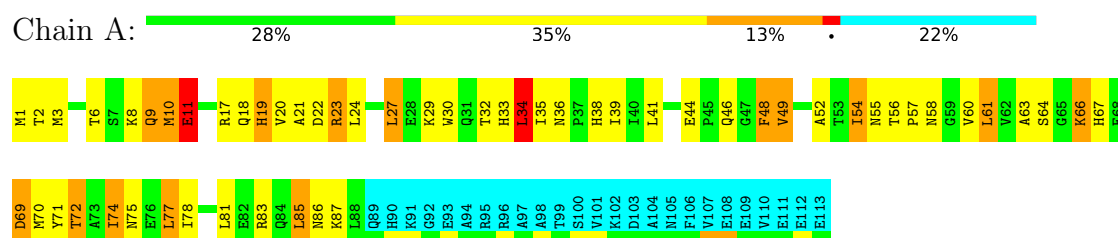
4.2.8 Score per residue for model 8

- Molecule 1: Protein yfiA



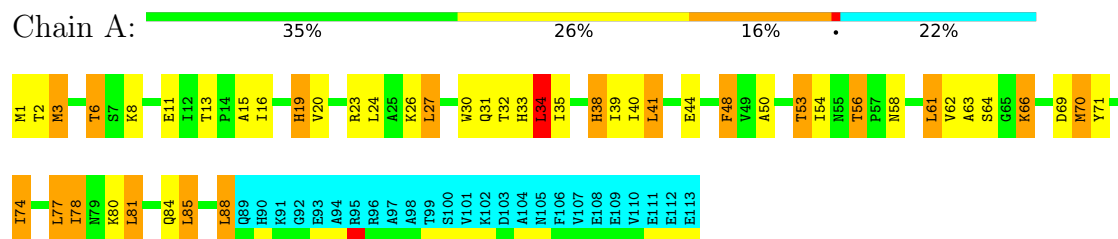
4.2.9 Score per residue for model 9

- Molecule 1: Protein yfiA



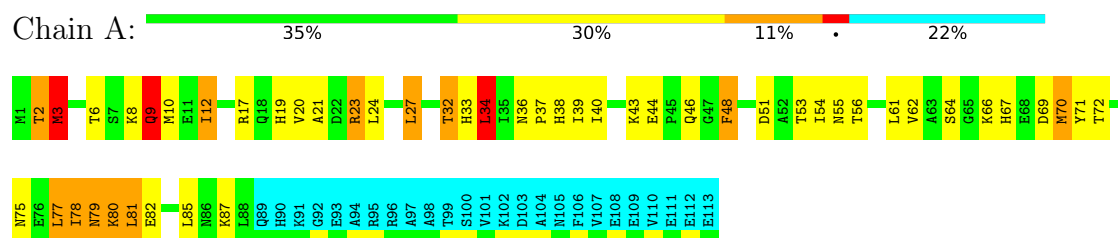
4.2.10 Score per residue for model 10

- Molecule 1: Protein yfiA



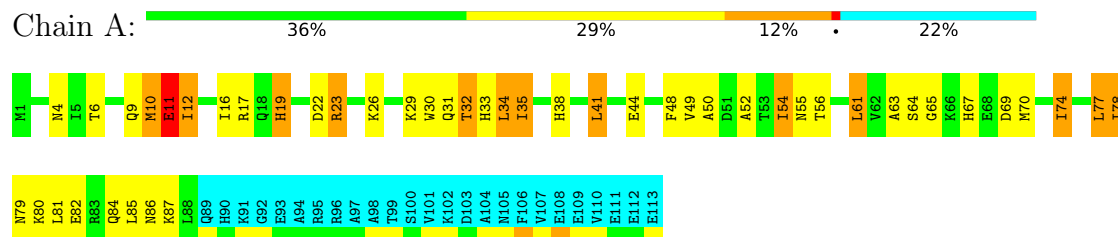
4.2.11 Score per residue for model 11

- Molecule 1: Protein yfiA



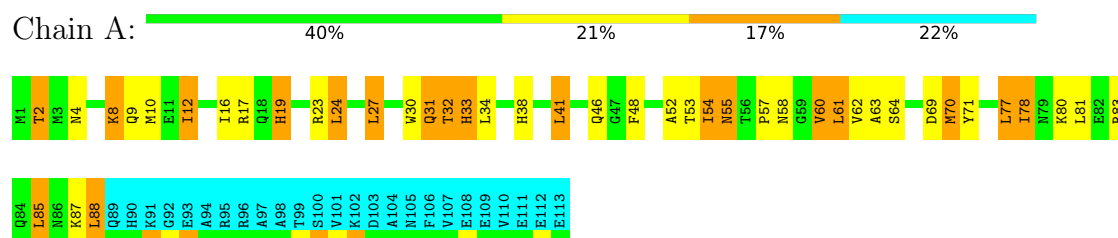
4.2.12 Score per residue for model 12

- Molecule 1: Protein yfiA



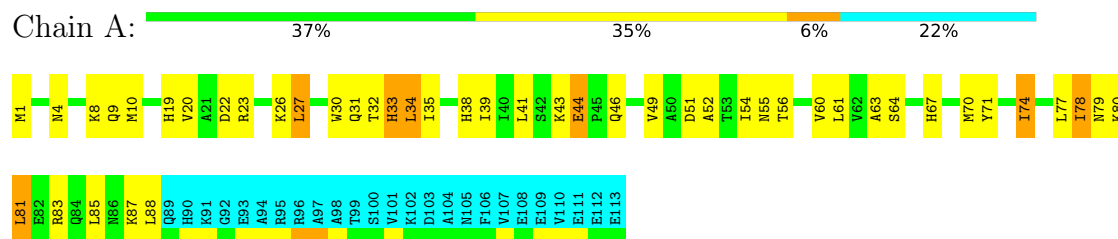
4.2.13 Score per residue for model 13

- Molecule 1: Protein yfiA



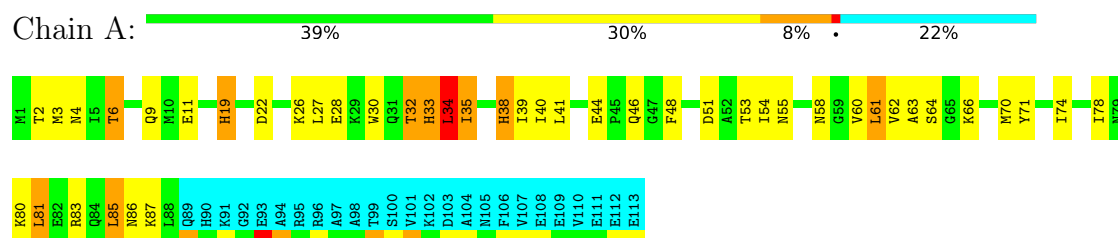
4.2.14 Score per residue for model 14 (medoid)

- Molecule 1: Protein yfiA



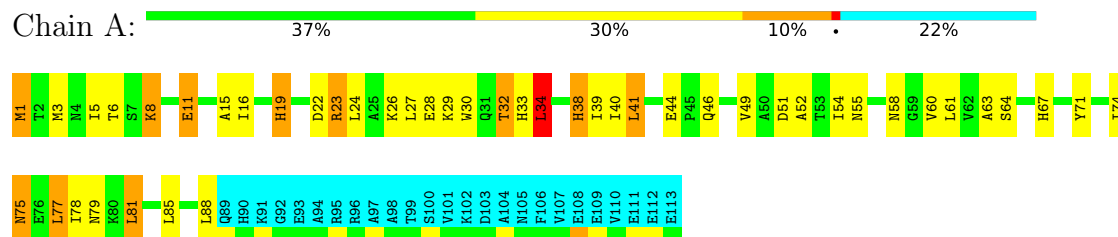
4.2.15 Score per residue for model 15

- Molecule 1: Protein yfiA



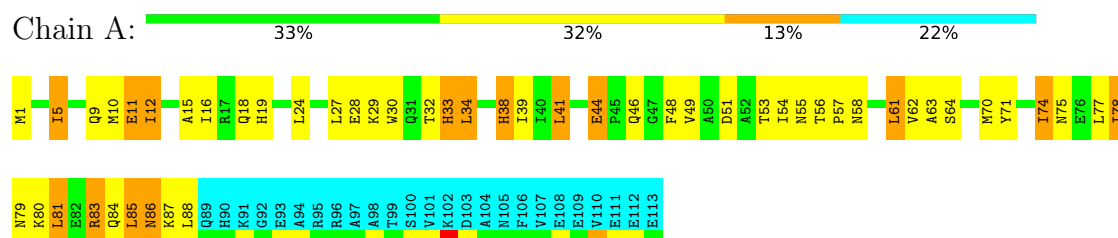
4.2.16 Score per residue for model 16

- Molecule 1: Protein yfiA



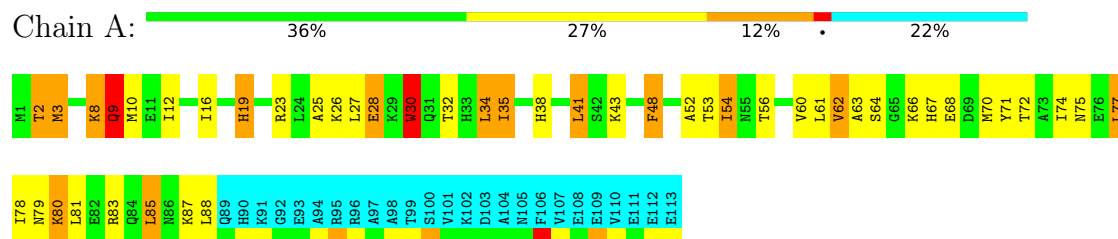
4.2.17 Score per residue for model 17

- Molecule 1: Protein yfiA



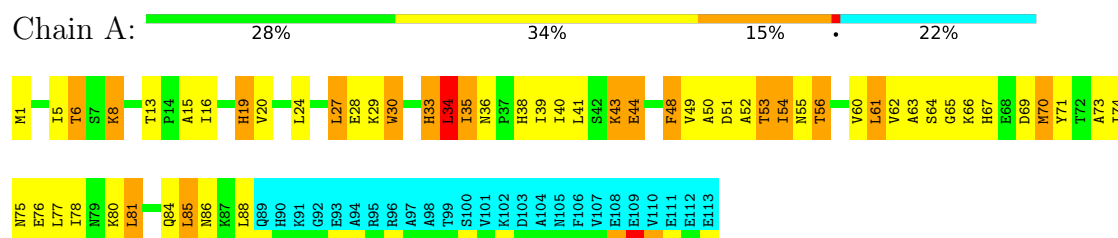
4.2.18 Score per residue for model 18

- Molecule 1: Protein yfiA



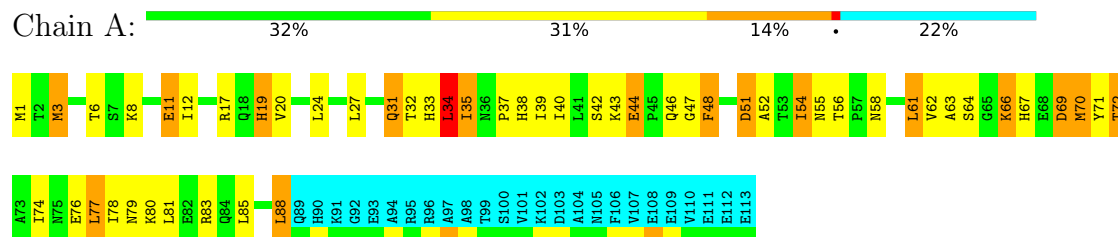
4.2.19 Score per residue for model 19

- Molecule 1: Protein yfiA



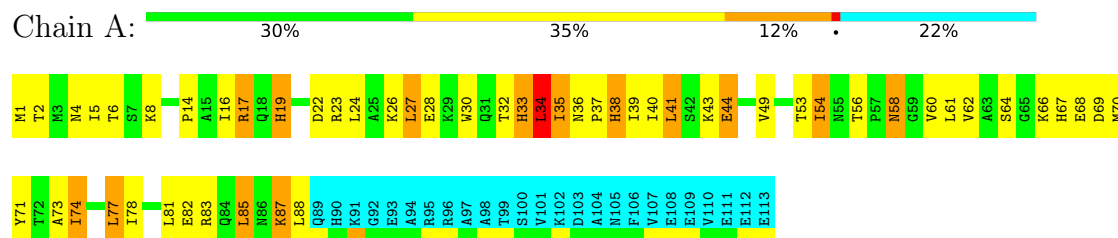
4.2.20 Score per residue for model 20

- Molecule 1: Protein yfiA



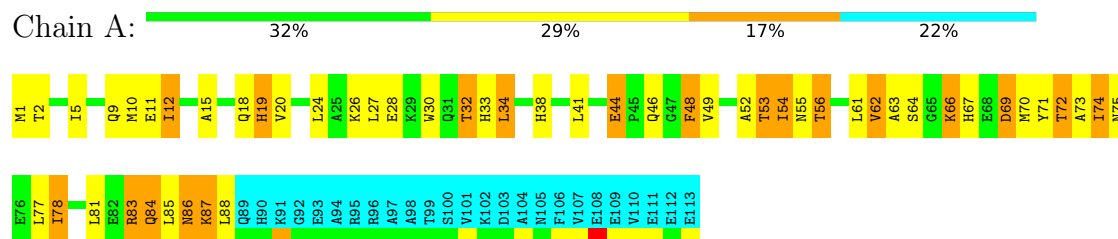
4.2.21 Score per residue for model 21

- Molecule 1: Protein yfiA



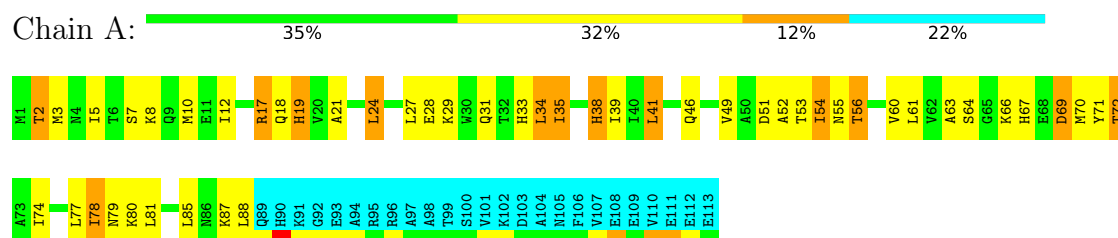
4.2.22 Score per residue for model 22

- Molecule 1: Protein yfiA



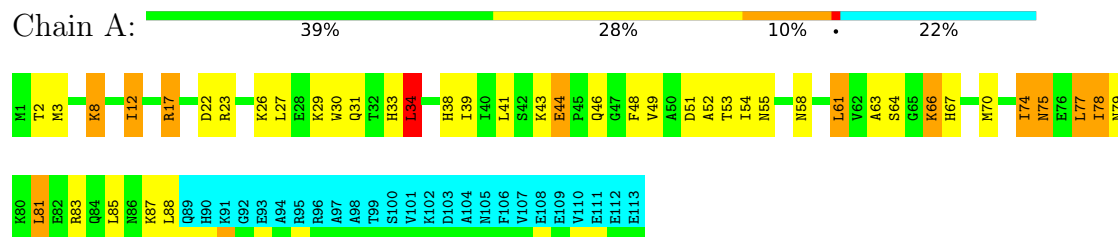
4.2.23 Score per residue for model 23

- Molecule 1: Protein yfiA



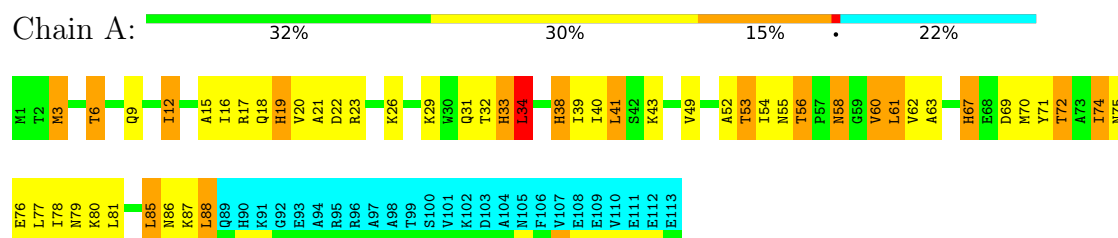
4.2.24 Score per residue for model 24

- Molecule 1: Protein yfiA



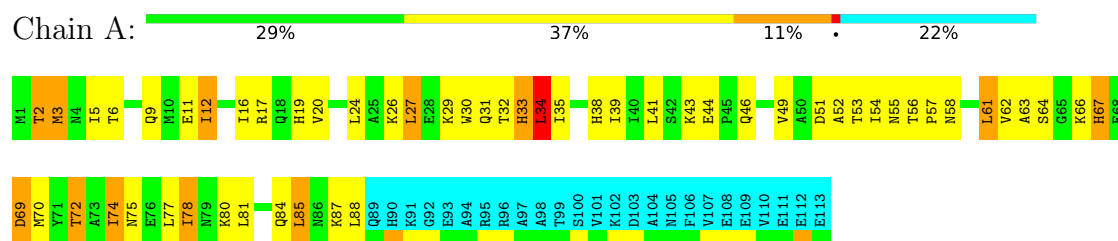
4.2.25 Score per residue for model 25

- Molecule 1: Protein yfiA



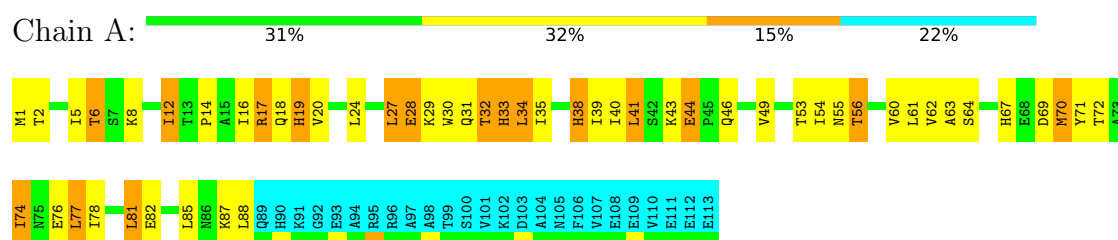
4.2.26 Score per residue for model 26

- Molecule 1: Protein yfiA



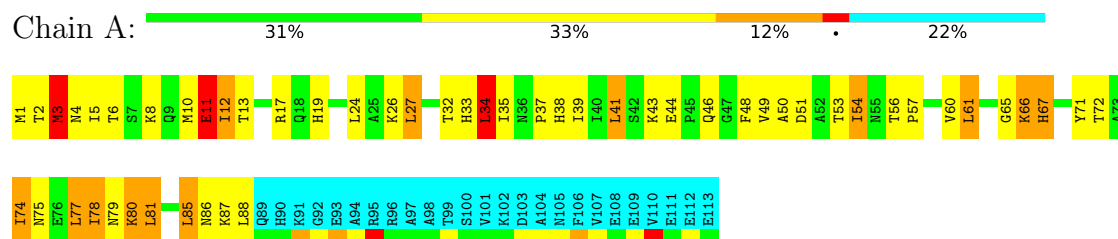
4.2.27 Score per residue for model 27

- Molecule 1: Protein yfiA



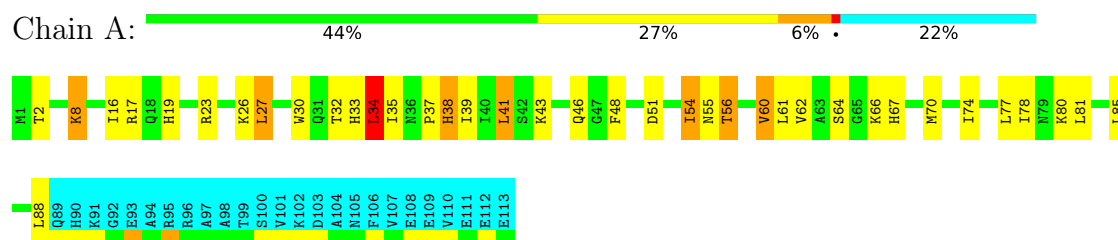
4.2.28 Score per residue for model 28

- Molecule 1: Protein yfiA



4.2.29 Score per residue for model 29

- Molecule 1: Protein yfiA



5 Refinement protocol and experimental data overview

The models were refined using the following method: *Torsion angle dynamics and simulated annealing*.

Of the 50 calculated structures, 29 were deposited, based on the following criterion: *structures with acceptable covalent geometry, structures with the least restraint violations*.

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

Software name	Classification	Version
CNS	refinement	1.0

No chemical shift data was provided.

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	698	728	724	36±8
All	All	20242	21112	20996	1049

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 25.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:63:ALA:HB2	1:A:81:LEU:HD12	1.12	1.15	16	6
1:A:34:LEU:HD13	1:A:56:THR:HG23	1.01	1.29	22	1
1:A:61:LEU:HD11	1:A:85:LEU:HD21	1.00	1.33	27	1
1:A:63:ALA:HB1	1:A:77:LEU:HD21	0.93	1.39	4	1
1:A:41:LEU:HD21	1:A:70:MET:HE1	0.92	1.39	24	1
1:A:52:ALA:HB3	1:A:63:ALA:HB3	0.91	1.41	22	17
1:A:54:ILE:HD11	1:A:81:LEU:HD21	0.90	1.40	2	3
1:A:63:ALA:HB2	1:A:81:LEU:CD1	0.89	1.97	6	12
1:A:77:LEU:HD13	1:A:78:ILE:N	0.85	1.85	7	1
1:A:61:LEU:CD2	1:A:88:LEU:HD21	0.85	2.00	19	1
1:A:15:ALA:HB1	1:A:71:TYR:CE1	0.84	2.06	10	5
1:A:27:LEU:HD12	1:A:78:ILE:HD11	0.84	1.49	19	1
1:A:63:ALA:CB	1:A:81:LEU:HD12	0.84	2.01	16	2
1:A:56:THR:HG21	1:A:61:LEU:HD12	0.84	1.49	9	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:5:ILE:HG22	1:A:39:ILE:HB	0.83	1.49	17	1
1:A:33:HIS:C	1:A:34:LEU:HD22	0.83	1.98	2	10
1:A:34:LEU:HD12	1:A:54:ILE:HG21	0.82	1.52	26	8
1:A:34:LEU:CD1	1:A:56:THR:HG22	0.81	2.06	27	4
1:A:34:LEU:HD12	1:A:54:ILE:CG2	0.81	2.06	22	13
1:A:75:ASN:O	1:A:78:ILE:HG22	0.80	1.76	9	12
1:A:61:LEU:HD12	1:A:62:VAL:N	0.79	1.92	4	1
1:A:63:ALA:HB1	1:A:77:LEU:CD2	0.79	2.08	4	1
1:A:63:ALA:HB2	1:A:81:LEU:HD11	0.78	1.56	20	4
1:A:61:LEU:HD22	1:A:85:LEU:HD12	0.78	1.56	24	1
1:A:6:THR:O	1:A:40:ILE:HG23	0.77	1.79	10	11
1:A:48:PHE:CE2	1:A:70:MET:HE2	0.77	2.13	24	1
1:A:77:LEU:HD12	1:A:78:ILE:N	0.77	1.95	26	3
1:A:23:ARG:NE	1:A:78:ILE:HD12	0.77	1.95	11	1
1:A:61:LEU:HD21	1:A:88:LEU:HD21	0.77	1.57	19	1
1:A:34:LEU:HD12	1:A:54:ILE:HG22	0.76	1.55	14	10
1:A:58:ASN:CB	1:A:88:LEU:HD21	0.76	2.10	5	2
1:A:53:THR:HG22	1:A:60:VAL:HG12	0.76	1.56	7	3
1:A:32:THR:HG23	1:A:34:LEU:HD21	0.76	1.55	10	3
1:A:30:TRP:CE3	1:A:85:LEU:HD21	0.76	2.15	21	2
1:A:27:LEU:HD12	1:A:34:LEU:HD21	0.76	1.56	14	1
1:A:61:LEU:HD23	1:A:81:LEU:CD2	0.76	2.10	28	1
1:A:58:ASN:O	1:A:88:LEU:HD11	0.76	1.81	5	1
1:A:16:ILE:O	1:A:20:VAL:HG22	0.75	1.80	19	3
1:A:34:LEU:HD22	1:A:54:ILE:HB	0.75	1.57	19	3
1:A:77:LEU:HD23	1:A:78:ILE:N	0.75	1.96	3	11
1:A:16:ILE:HD11	1:A:70:MET:HE2	0.75	1.57	13	1
1:A:38:HIS:C	1:A:39:ILE:HD12	0.75	2.06	21	7
1:A:63:ALA:HB1	1:A:77:LEU:HG	0.75	1.58	20	11
1:A:61:LEU:HD11	1:A:85:LEU:CD2	0.75	2.11	27	1
1:A:34:LEU:HD23	1:A:54:ILE:HG21	0.74	1.57	1	2
1:A:77:LEU:HD22	1:A:77:LEU:O	0.74	1.82	7	1
1:A:61:LEU:HD11	1:A:85:LEU:HD12	0.74	1.59	13	1
1:A:27:LEU:O	1:A:27:LEU:HD22	0.74	1.81	27	1
1:A:74:ILE:O	1:A:78:ILE:HG23	0.74	1.83	24	2
1:A:88:LEU:O	1:A:88:LEU:HD13	0.73	1.83	8	1
1:A:32:THR:OG1	1:A:85:LEU:HD21	0.73	1.83	1	2
1:A:33:HIS:O	1:A:35:ILE:HD13	0.73	1.82	21	1
1:A:16:ILE:HG21	1:A:41:LEU:HD13	0.73	1.59	7	7
1:A:23:ARG:HD3	1:A:78:ILE:HG22	0.73	1.60	3	1
1:A:61:LEU:HD21	1:A:81:LEU:CB	0.73	2.13	4	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:63:ALA:HB1	1:A:77:LEU:HD13	0.73	1.61	9	2
1:A:30:TRP:CE3	1:A:85:LEU:HD22	0.72	2.18	1	10
1:A:58:ASN:CB	1:A:88:LEU:HD13	0.72	2.15	20	1
1:A:10:MET:HE3	1:A:42:SER:C	0.72	2.09	3	1
1:A:30:TRP:CE3	1:A:85:LEU:HD13	0.71	2.20	15	2
1:A:16:ILE:HG23	1:A:74:ILE:CD1	0.71	2.15	7	1
1:A:67:HIS:CD2	1:A:72:THR:HG21	0.71	2.20	1	3
1:A:19:HIS:HB3	1:A:74:ILE:HG21	0.71	1.63	4	7
1:A:58:ASN:HB2	1:A:88:LEU:HD21	0.70	1.63	5	1
1:A:44:GLU:HB2	1:A:49:VAL:HG23	0.70	1.63	8	13
1:A:23:ARG:NH2	1:A:74:ILE:HG22	0.70	2.02	5	1
1:A:27:LEU:HD23	1:A:27:LEU:O	0.70	1.87	18	3
1:A:34:LEU:HD23	1:A:56:THR:HG22	0.69	1.61	8	4
1:A:44:GLU:OE2	1:A:49:VAL:HG21	0.69	1.87	19	2
1:A:5:ILE:HG22	1:A:39:ILE:CB	0.69	2.16	17	1
1:A:81:LEU:HD22	1:A:85:LEU:CD1	0.69	2.17	24	1
1:A:61:LEU:HD21	1:A:88:LEU:HD12	0.69	1.63	18	1
1:A:41:LEU:HD23	1:A:41:LEU:O	0.68	1.88	9	4
1:A:74:ILE:HA	1:A:77:LEU:HD23	0.68	1.65	9	3
1:A:48:PHE:CZ	1:A:70:MET:HE2	0.68	2.23	24	1
1:A:58:ASN:HB3	1:A:88:LEU:HD21	0.68	1.63	21	3
1:A:34:LEU:N	1:A:34:LEU:HD13	0.68	2.04	2	2
1:A:1:MET:HE3	1:A:28:GLU:HB2	0.68	1.66	7	1
1:A:63:ALA:HB2	1:A:81:LEU:CG	0.67	2.19	3	6
1:A:53:THR:HG23	1:A:62:VAL:HG22	0.67	1.66	25	4
1:A:61:LEU:CD1	1:A:85:LEU:HD12	0.67	2.19	13	2
1:A:35:ILE:HD12	1:A:55:ASN:O	0.67	1.89	15	1
1:A:27:LEU:HD22	1:A:34:LEU:HD21	0.67	1.65	22	2
1:A:41:LEU:HD11	1:A:70:MET:SD	0.67	2.30	24	1
1:A:34:LEU:CD2	1:A:56:THR:HG22	0.66	2.20	8	2
1:A:50:ALA:HB1	1:A:77:LEU:HD12	0.66	1.67	1	2
1:A:78:ILE:O	1:A:78:ILE:HD13	0.66	1.89	4	7
1:A:16:ILE:CG2	1:A:41:LEU:HD13	0.66	2.21	5	3
1:A:63:ALA:HB1	1:A:77:LEU:CD1	0.66	2.21	9	2
1:A:30:TRP:CD2	1:A:85:LEU:HD21	0.65	2.25	21	1
1:A:10:MET:HE2	1:A:48:PHE:CE1	0.65	2.26	3	2
1:A:77:LEU:HD22	1:A:77:LEU:C	0.65	2.16	7	1
1:A:63:ALA:CB	1:A:77:LEU:HD21	0.65	2.19	4	1
1:A:61:LEU:HD11	1:A:84:GLN:HB3	0.65	1.66	8	1
1:A:63:ALA:HA	1:A:81:LEU:HD11	0.65	1.67	4	1
1:A:63:ALA:O	1:A:77:LEU:HD12	0.65	1.91	3	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:52:ALA:HB3	1:A:77:LEU:HD23	0.65	1.66	7	1
1:A:32:THR:CG2	1:A:85:LEU:HD21	0.65	2.22	13	2
1:A:15:ALA:HB1	1:A:71:TYR:OH	0.65	1.92	17	4
1:A:56:THR:HG21	1:A:61:LEU:HD13	0.65	1.68	17	1
1:A:12:ILE:HD13	1:A:12:ILE:N	0.64	2.07	22	10
1:A:32:THR:HG23	1:A:85:LEU:HD21	0.64	1.67	13	2
1:A:4:ASN:C	1:A:5:ILE:HD12	0.64	2.16	21	1
1:A:50:ALA:HB1	1:A:77:LEU:HD13	0.64	1.67	10	2
1:A:16:ILE:HD13	1:A:41:LEU:HD21	0.64	1.66	4	8
1:A:34:LEU:HD11	1:A:56:THR:HG22	0.64	1.70	12	2
1:A:16:ILE:HG21	1:A:41:LEU:HD22	0.64	1.68	13	2
1:A:61:LEU:HD13	1:A:81:LEU:HD22	0.64	1.70	26	2
1:A:88:LEU:HD12	1:A:88:LEU:O	0.64	1.92	20	1
1:A:61:LEU:HD12	1:A:81:LEU:HD22	0.64	1.69	23	1
1:A:20:VAL:HG22	1:A:74:ILE:HG23	0.64	1.71	8	1
1:A:27:LEU:HD13	1:A:30:TRP:HB2	0.63	1.70	4	3
1:A:54:ILE:O	1:A:60:VAL:HG13	0.63	1.92	9	4
1:A:32:THR:HG23	1:A:34:LEU:CD2	0.63	2.23	10	1
1:A:34:LEU:CD1	1:A:54:ILE:HG22	0.63	2.24	20	2
1:A:61:LEU:HD13	1:A:84:GLN:NE2	0.63	2.08	4	1
1:A:16:ILE:CD1	1:A:70:MET:HE2	0.62	2.22	13	1
1:A:40:ILE:HD12	1:A:51:ASP:O	0.62	1.94	16	2
1:A:61:LEU:HD22	1:A:84:GLN:OE1	0.62	1.94	22	1
1:A:23:ARG:HG3	1:A:78:ILE:HD12	0.62	1.71	16	2
1:A:32:THR:O	1:A:33:HIS:HB2	0.62	1.94	21	2
1:A:58:ASN:OD1	1:A:88:LEU:HD13	0.62	1.93	10	1
1:A:61:LEU:CD2	1:A:81:LEU:HD21	0.62	2.24	29	1
1:A:53:THR:HG22	1:A:60:VAL:CG1	0.62	2.25	19	3
1:A:61:LEU:HD22	1:A:81:LEU:HD21	0.61	1.71	29	1
1:A:3:MET:HE3	1:A:24:LEU:CD2	0.61	2.25	5	1
1:A:27:LEU:CD1	1:A:78:ILE:HD11	0.61	2.25	28	2
1:A:61:LEU:HD13	1:A:81:LEU:HD11	0.61	1.71	5	2
1:A:5:ILE:CG2	1:A:39:ILE:HB	0.61	2.26	17	1
1:A:61:LEU:CB	1:A:81:LEU:HD21	0.61	2.26	16	1
1:A:62:VAL:O	1:A:62:VAL:HG13	0.61	1.96	15	9
1:A:34:LEU:C	1:A:35:ILE:HD13	0.60	2.20	20	1
1:A:39:ILE:HD13	1:A:52:ALA:HB2	0.60	1.73	8	3
1:A:34:LEU:CD2	1:A:34:LEU:N	0.60	2.64	26	2
1:A:85:LEU:HD23	1:A:85:LEU:O	0.60	1.96	29	1
1:A:16:ILE:HD13	1:A:41:LEU:CD2	0.60	2.26	4	4
1:A:15:ALA:HB3	1:A:70:MET:HE1	0.59	1.74	19	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:41:LEU:HD21	1:A:70:MET:CE	0.59	2.24	24	1
1:A:2:THR:O	1:A:3:MET:CB	0.59	2.50	11	1
1:A:61:LEU:HD11	1:A:85:LEU:CD1	0.59	2.26	13	1
1:A:27:LEU:C	1:A:27:LEU:HD13	0.59	2.23	27	2
1:A:34:LEU:HD22	1:A:34:LEU:N	0.59	2.12	15	12
1:A:53:THR:HG23	1:A:62:VAL:HG23	0.59	1.74	22	1
1:A:1:MET:HB2	1:A:24:LEU:HD11	0.59	1.72	16	1
1:A:63:ALA:HB2	1:A:81:LEU:HG	0.59	1.73	17	6
1:A:61:LEU:HD23	1:A:84:GLN:CD	0.59	2.23	10	2
1:A:12:ILE:HG22	1:A:17:ARG:HB3	0.59	1.74	20	4
1:A:44:GLU:HB3	1:A:49:VAL:HG23	0.59	1.75	4	2
1:A:16:ILE:HD13	1:A:41:LEU:HD11	0.59	1.73	25	3
1:A:61:LEU:HD21	1:A:81:LEU:HB3	0.58	1.75	4	1
1:A:63:ALA:CB	1:A:81:LEU:HD11	0.58	2.29	3	4
1:A:61:LEU:HD13	1:A:84:GLN:CD	0.58	2.23	4	2
1:A:39:ILE:HD12	1:A:39:ILE:N	0.58	2.13	4	7
1:A:50:ALA:HB3	1:A:65:GLY:O	0.58	1.98	19	2
1:A:30:TRP:HB3	1:A:85:LEU:HD13	0.58	1.75	27	1
1:A:56:THR:HG21	1:A:61:LEU:HG	0.58	1.75	27	1
1:A:33:HIS:O	1:A:34:LEU:C	0.58	2.47	28	13
1:A:27:LEU:HD12	1:A:27:LEU:O	0.58	1.99	10	1
1:A:81:LEU:HD22	1:A:85:LEU:HD12	0.58	1.74	10	2
1:A:49:VAL:HG12	1:A:51:ASP:OD2	0.58	1.99	6	1
1:A:16:ILE:CG1	1:A:70:MET:HE2	0.58	2.27	13	2
1:A:10:MET:HE2	1:A:48:PHE:CZ	0.58	2.34	28	1
1:A:19:HIS:CD2	1:A:71:TYR:CD1	0.57	2.92	4	3
1:A:61:LEU:HD22	1:A:81:LEU:CD2	0.57	2.29	6	2
1:A:15:ALA:CB	1:A:70:MET:HE1	0.57	2.28	19	1
1:A:30:TRP:CG	1:A:85:LEU:HD23	0.57	2.34	24	1
1:A:58:ASN:O	1:A:88:LEU:HD22	0.57	2.00	7	1
1:A:35:ILE:HD11	1:A:55:ASN:OD1	0.57	2.00	6	2
1:A:16:ILE:HG23	1:A:74:ILE:HD12	0.57	1.75	7	3
1:A:39:ILE:HD13	1:A:52:ALA:CB	0.57	2.30	26	2
1:A:3:MET:HG3	1:A:24:LEU:HD22	0.57	1.77	28	1
1:A:39:ILE:HD13	1:A:52:ALA:HB1	0.57	1.77	20	1
1:A:61:LEU:HB3	1:A:81:LEU:HD22	0.57	1.76	20	1
1:A:32:THR:O	1:A:34:LEU:N	0.57	2.37	13	1
1:A:40:ILE:HD12	1:A:40:ILE:N	0.57	2.15	21	2
1:A:78:ILE:HD13	1:A:78:ILE:C	0.57	2.25	4	4
1:A:77:LEU:HD12	1:A:77:LEU:C	0.57	2.25	23	6
1:A:32:THR:O	1:A:33:HIS:CB	0.56	2.52	21	6

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:61:LEU:HB3	1:A:81:LEU:HD21	0.56	1.77	28	8
1:A:15:ALA:HB1	1:A:71:TYR:CZ	0.56	2.35	22	3
1:A:16:ILE:O	1:A:20:VAL:HG23	0.56	2.00	26	3
1:A:54:ILE:CD1	1:A:81:LEU:HD11	0.56	2.30	19	1
1:A:24:LEU:HA	1:A:27:LEU:HD12	0.56	1.78	20	1
1:A:34:LEU:HD13	1:A:34:LEU:H	0.55	1.62	2	2
1:A:88:LEU:O	1:A:88:LEU:HD23	0.55	2.00	21	3
1:A:52:ALA:HB3	1:A:77:LEU:CD2	0.55	2.30	7	1
1:A:61:LEU:HD12	1:A:62:VAL:H	0.55	1.61	4	1
1:A:34:LEU:H	1:A:34:LEU:HD13	0.55	1.59	21	1
1:A:32:THR:O	1:A:33:HIS:HB3	0.55	2.02	26	1
1:A:61:LEU:HD22	1:A:81:LEU:CD1	0.55	2.32	1	1
1:A:3:MET:HE2	1:A:24:LEU:HD23	0.55	1.77	10	1
1:A:12:ILE:N	1:A:12:ILE:HD12	0.55	2.17	20	1
1:A:63:ALA:HB2	1:A:81:LEU:HD23	0.55	1.77	25	4
1:A:61:LEU:HD13	1:A:85:LEU:HD12	0.55	1.79	25	1
1:A:32:THR:HG23	1:A:34:LEU:CD1	0.55	2.32	28	1
1:A:74:ILE:HA	1:A:77:LEU:HD22	0.55	1.79	3	1
1:A:56:THR:CG2	1:A:61:LEU:HD12	0.55	2.31	10	1
1:A:63:ALA:HB2	1:A:81:LEU:CD2	0.55	2.32	27	3
1:A:5:ILE:HD12	1:A:5:ILE:N	0.54	2.16	21	1
1:A:27:LEU:HD13	1:A:27:LEU:O	0.54	2.01	7	1
1:A:87:LYS:N	1:A:87:LYS:HD2	0.54	2.15	14	2
1:A:2:THR:HG22	1:A:35:ILE:O	0.54	2.02	18	2
1:A:56:THR:HG21	1:A:61:LEU:CD1	0.54	2.33	10	3
1:A:34:LEU:CD2	1:A:54:ILE:HG21	0.54	2.31	1	2
1:A:61:LEU:HD21	1:A:81:LEU:CG	0.54	2.31	4	1
1:A:77:LEU:HD13	1:A:77:LEU:C	0.54	2.27	7	1
1:A:67:HIS:CD2	1:A:72:THR:HG22	0.54	2.38	26	1
1:A:81:LEU:HD23	1:A:81:LEU:O	0.54	2.03	18	4
1:A:77:LEU:C	1:A:77:LEU:HD23	0.54	2.28	25	4
1:A:58:ASN:O	1:A:88:LEU:HD12	0.54	2.03	13	2
1:A:63:ALA:HB1	1:A:77:LEU:CG	0.53	2.33	4	4
1:A:77:LEU:HD23	1:A:78:ILE:H	0.53	1.63	3	1
1:A:3:MET:SD	1:A:24:LEU:HD22	0.53	2.43	23	1
1:A:61:LEU:HB2	1:A:81:LEU:HD21	0.53	1.80	16	2
1:A:61:LEU:HD22	1:A:88:LEU:HD21	0.53	1.78	19	1
1:A:3:MET:SD	1:A:5:ILE:HD11	0.53	2.43	28	1
1:A:81:LEU:O	1:A:81:LEU:HD13	0.53	2.04	29	1
1:A:23:ARG:NH2	1:A:77:LEU:HD22	0.53	2.19	5	1
1:A:55:ASN:HD22	1:A:60:VAL:HG22	0.53	1.64	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:10:MET:O	1:A:11:GLU:O	0.53	2.26	28	2
1:A:16:ILE:HG21	1:A:41:LEU:CD2	0.53	2.33	13	1
1:A:48:PHE:CE2	1:A:70:MET:CB	0.53	2.92	11	3
1:A:16:ILE:HD11	1:A:70:MET:SD	0.53	2.44	29	1
1:A:5:ILE:HG23	1:A:39:ILE:HB	0.53	1.79	1	1
1:A:61:LEU:HD22	1:A:81:LEU:HD13	0.53	1.79	1	1
1:A:63:ALA:CA	1:A:81:LEU:HD11	0.53	2.34	4	1
1:A:10:MET:CE	1:A:41:LEU:HD21	0.52	2.34	6	1
1:A:77:LEU:HD23	1:A:77:LEU:C	0.52	2.29	6	2
1:A:78:ILE:C	1:A:78:ILE:HD13	0.52	2.29	13	1
1:A:30:TRP:CD2	1:A:85:LEU:HD23	0.52	2.39	24	1
1:A:9:GLN:HG2	1:A:10:MET:N	0.52	2.19	6	2
1:A:53:THR:HG23	1:A:62:VAL:CG2	0.52	2.34	22	1
1:A:13:THR:H	1:A:16:ILE:HD12	0.52	1.64	19	1
1:A:70:MET:HG2	1:A:71:TYR:N	0.52	2.20	13	9
1:A:54:ILE:HD11	1:A:81:LEU:CD2	0.52	2.26	2	1
1:A:56:THR:HG21	1:A:61:LEU:HD11	0.52	1.80	25	1
1:A:27:LEU:HD13	1:A:30:TRP:CB	0.52	2.34	4	3
1:A:54:ILE:HD11	1:A:81:LEU:HD11	0.52	1.82	21	1
1:A:28:GLU:HA	1:A:34:LEU:HD23	0.52	1.82	15	1
1:A:16:ILE:HG21	1:A:41:LEU:CD1	0.52	2.35	27	4
1:A:48:PHE:O	1:A:66:LYS:HA	0.52	2.04	22	6
1:A:3:MET:CE	1:A:24:LEU:HD23	0.52	2.35	10	1
1:A:10:MET:C	1:A:11:GLU:HG3	0.52	2.29	28	1
1:A:19:HIS:CE1	1:A:71:TYR:CE2	0.52	2.98	16	6
1:A:2:THR:HG21	1:A:36:ASN:OD1	0.52	2.05	21	1
1:A:19:HIS:CE1	1:A:71:TYR:CZ	0.52	2.98	21	1
1:A:16:ILE:HG13	1:A:70:MET:HE2	0.51	1.82	19	2
1:A:81:LEU:HD13	1:A:81:LEU:C	0.51	2.30	29	2
1:A:38:HIS:HB2	1:A:53:THR:HG23	0.51	1.83	8	1
1:A:50:ALA:HB1	1:A:77:LEU:CD1	0.51	2.35	1	2
1:A:61:LEU:HD22	1:A:84:GLN:HE22	0.51	1.65	4	1
1:A:19:HIS:CD2	1:A:71:TYR:CE1	0.51	2.99	4	6
1:A:33:HIS:O	1:A:34:LEU:O	0.51	2.28	16	7
1:A:62:VAL:HG12	1:A:62:VAL:O	0.51	2.05	3	2
1:A:61:LEU:CG	1:A:81:LEU:HG	0.51	2.36	4	1
1:A:8:LYS:HG3	1:A:9:GLN:N	0.51	2.21	9	1
1:A:34:LEU:HD12	1:A:54:ILE:CB	0.51	2.36	22	1
1:A:19:HIS:CB	1:A:71:TYR:CE1	0.51	2.93	6	1
1:A:6:THR:O	1:A:6:THR:HG23	0.51	2.06	8	2
1:A:88:LEU:C	1:A:88:LEU:HD12	0.51	2.30	19	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:58:ASN:HD21	1:A:88:LEU:HD13	0.51	1.66	25	1
1:A:19:HIS:CE1	1:A:71:TYR:CE1	0.50	2.99	3	4
1:A:34:LEU:HD13	1:A:56:THR:HG22	0.50	1.82	18	1
1:A:53:THR:HG23	1:A:53:THR:O	0.50	2.07	28	3
1:A:20:VAL:HG21	1:A:39:ILE:HG21	0.50	1.82	9	3
1:A:55:ASN:ND2	1:A:60:VAL:HG23	0.50	2.22	13	1
1:A:19:HIS:CG	1:A:71:TYR:CE1	0.50	2.99	18	3
1:A:56:THR:CG2	1:A:61:LEU:HD23	0.50	2.37	11	1
1:A:12:ILE:HG22	1:A:17:ARG:CB	0.50	2.37	24	1
1:A:61:LEU:HD22	1:A:85:LEU:CD1	0.50	2.33	24	1
1:A:37:PRO:HB3	1:A:54:ILE:HG23	0.50	1.84	28	1
1:A:34:LEU:HD23	1:A:56:THR:HG23	0.50	1.83	29	1
1:A:49:VAL:HG12	1:A:49:VAL:O	0.49	2.06	8	1
1:A:61:LEU:HB2	1:A:81:LEU:HD23	0.49	1.84	15	1
1:A:33:HIS:CD2	1:A:33:HIS:N	0.49	2.79	27	7
1:A:12:ILE:HG22	1:A:16:ILE:CG2	0.49	2.36	5	1
1:A:3:MET:HE1	1:A:21:ALA:CB	0.49	2.38	11	1
1:A:32:THR:HG1	1:A:34:LEU:HD13	0.49	1.67	15	1
1:A:27:LEU:HD12	1:A:32:THR:CG2	0.49	2.37	4	1
1:A:20:VAL:HG21	1:A:39:ILE:HG13	0.49	1.84	14	1
1:A:33:HIS:CD2	1:A:33:HIS:C	0.49	2.91	26	1
1:A:61:LEU:HD23	1:A:81:LEU:HD23	0.49	1.82	28	1
1:A:10:MET:HE2	1:A:48:PHE:CD1	0.49	2.42	3	1
1:A:38:HIS:O	1:A:39:ILE:HD13	0.49	2.07	25	3
1:A:23:ARG:HE	1:A:77:LEU:HD11	0.48	1.68	9	1
1:A:30:TRP:CD2	1:A:85:LEU:HD13	0.48	2.43	9	2
1:A:61:LEU:HD11	1:A:85:LEU:CG	0.48	2.38	13	1
1:A:33:HIS:O	1:A:33:HIS:CG	0.48	2.66	26	1
1:A:16:ILE:HD13	1:A:41:LEU:CD1	0.48	2.38	8	2
1:A:16:ILE:CD1	1:A:41:LEU:HD21	0.48	2.38	4	1
1:A:19:HIS:CE1	1:A:71:TYR:CD1	0.48	3.02	3	3
1:A:19:HIS:ND1	1:A:71:TYR:CE1	0.48	2.82	19	8
1:A:5:ILE:HA	1:A:39:ILE:HB	0.48	1.85	17	1
1:A:5:ILE:CG2	1:A:12:ILE:HD12	0.48	2.38	22	1
1:A:1:MET:N	1:A:34:LEU:O	0.48	2.46	10	1
1:A:2:THR:HG22	1:A:4:ASN:OD1	0.48	2.09	6	2
1:A:5:ILE:HG21	1:A:12:ILE:HD12	0.48	1.86	22	1
1:A:76:GLU:HA	1:A:79:ASN:OD1	0.48	2.09	6	2
1:A:69:ASP:CB	1:A:72:THR:OG1	0.48	2.62	23	6
1:A:35:ILE:N	1:A:35:ILE:HD13	0.47	2.24	2	2
1:A:70:MET:CG	1:A:71:TYR:N	0.47	2.77	22	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:5:ILE:HG23	1:A:39:ILE:O	0.47	2.08	19	1
1:A:62:VAL:C	1:A:81:LEU:HD21	0.47	2.33	20	1
1:A:23:ARG:NH2	1:A:24:LEU:HD12	0.47	2.23	13	1
1:A:61:LEU:HD22	1:A:81:LEU:HD23	0.47	1.86	17	1
1:A:5:ILE:HG21	1:A:12:ILE:CD1	0.47	2.39	27	1
1:A:7:SER:OG	1:A:12:ILE:HD11	0.47	2.09	23	1
1:A:55:ASN:HA	1:A:60:VAL:HG22	0.47	1.86	2	1
1:A:34:LEU:HD22	1:A:54:ILE:CB	0.47	2.39	9	1
1:A:32:THR:O	1:A:33:HIS:CG	0.47	2.68	1	2
1:A:70:MET:O	1:A:74:ILE:CD1	0.47	2.62	3	8
1:A:43:LYS:HG2	1:A:48:PHE:CE1	0.47	2.44	19	2
1:A:61:LEU:HD23	1:A:88:LEU:CD1	0.47	2.39	25	1
1:A:23:ARG:HD3	1:A:78:ILE:HD11	0.47	1.86	12	1
1:A:81:LEU:C	1:A:81:LEU:HD13	0.47	2.34	12	1
1:A:61:LEU:HD23	1:A:61:LEU:N	0.47	2.25	15	1
1:A:14:PRO:HA	1:A:17:ARG:HG2	0.47	1.85	21	2
1:A:27:LEU:HD21	1:A:85:LEU:HD11	0.47	1.87	22	1
1:A:61:LEU:HD11	1:A:85:LEU:HG	0.47	1.87	6	1
1:A:3:MET:HE1	1:A:21:ALA:HA	0.47	1.87	8	1
1:A:35:ILE:HD12	1:A:36:ASN:H	0.47	1.69	19	1
1:A:54:ILE:O	1:A:56:THR:HG22	0.47	2.10	23	1
1:A:61:LEU:HD13	1:A:84:GLN:OE1	0.46	2.10	4	1
1:A:24:LEU:HD21	1:A:37:PRO:HB2	0.46	1.87	11	1
1:A:55:ASN:ND2	1:A:60:VAL:HG22	0.46	2.25	14	1
1:A:61:LEU:HD21	1:A:88:LEU:CD1	0.46	2.39	18	1
1:A:67:HIS:CD2	1:A:72:THR:CG2	0.46	2.98	26	4
1:A:32:THR:HG21	1:A:56:THR:CG2	0.46	2.40	11	1
1:A:37:PRO:HA	1:A:54:ILE:HG23	0.46	1.86	11	5
1:A:61:LEU:HD12	1:A:81:LEU:HD23	0.46	1.88	11	3
1:A:30:TRP:CE3	1:A:85:LEU:CD2	0.46	2.98	26	1
1:A:34:LEU:HD12	1:A:56:THR:HG22	0.46	1.86	27	1
1:A:82:GLU:O	1:A:86:ASN:ND2	0.46	2.48	4	2
1:A:11:GLU:OE1	1:A:12:ILE:HD12	0.46	2.11	3	1
1:A:81:LEU:HD12	1:A:81:LEU:H	0.46	1.69	4	1
1:A:56:THR:CG2	1:A:61:LEU:HD11	0.46	2.40	25	1
1:A:48:PHE:CE2	1:A:70:MET:HB3	0.46	2.46	11	2
1:A:48:PHE:CE2	1:A:70:MET:HB2	0.45	2.47	1	3
1:A:30:TRP:CH2	1:A:86:ASN:OD1	0.45	2.69	19	2
1:A:17:ARG:O	1:A:21:ALA:HB2	0.45	2.11	23	2
1:A:10:MET:HE1	1:A:48:PHE:CE2	0.45	2.47	6	1
1:A:32:THR:O	1:A:57:PRO:CG	0.45	2.64	26	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:27:LEU:HD22	1:A:27:LEU:C	0.45	2.33	27	1
1:A:63:ALA:HB2	1:A:81:LEU:HD21	0.45	1.89	3	2
1:A:79:ASN:CG	1:A:80:LYS:N	0.45	2.74	11	1
1:A:32:THR:OG1	1:A:56:THR:HG22	0.45	2.12	26	1
1:A:43:LYS:CG	1:A:48:PHE:CE1	0.45	3.00	19	2
1:A:70:MET:O	1:A:73:ALA:HB3	0.45	2.11	19	1
1:A:20:VAL:HG11	1:A:39:ILE:HG21	0.45	1.87	11	1
1:A:34:LEU:CD2	1:A:34:LEU:O	0.45	2.65	28	2
1:A:39:ILE:N	1:A:39:ILE:CD1	0.45	2.80	7	5
1:A:44:GLU:CD	1:A:49:VAL:HG21	0.45	2.36	6	1
1:A:30:TRP:HE3	1:A:85:LEU:HD22	0.45	1.71	8	2
1:A:34:LEU:HD13	1:A:34:LEU:N	0.45	2.26	21	1
1:A:24:LEU:C	1:A:24:LEU:HD23	0.45	2.37	9	1
1:A:39:ILE:C	1:A:40:ILE:HD12	0.45	2.36	5	1
1:A:36:ASN:O	1:A:36:ASN:OD1	0.45	2.35	9	1
1:A:32:THR:O	1:A:57:PRO:CD	0.45	2.64	13	1
1:A:34:LEU:HD12	1:A:54:ILE:HB	0.45	1.89	22	1
1:A:75:ASN:O	1:A:78:ILE:HG12	0.45	2.12	24	2
1:A:62:VAL:O	1:A:62:VAL:CG2	0.45	2.64	18	1
1:A:88:LEU:HD12	1:A:88:LEU:C	0.45	2.36	20	1
1:A:23:ARG:CG	1:A:78:ILE:HG21	0.45	2.42	25	1
1:A:33:HIS:O	1:A:34:LEU:HD22	0.44	2.11	2	1
1:A:50:ALA:CB	1:A:77:LEU:CD1	0.44	2.95	12	1
1:A:62:VAL:O	1:A:62:VAL:CG1	0.44	2.65	29	3
1:A:20:VAL:HG13	1:A:21:ALA:N	0.44	2.27	25	3
1:A:5:ILE:N	1:A:5:ILE:CD1	0.44	2.80	21	1
1:A:67:HIS:HB3	1:A:72:THR:CG2	0.44	2.43	25	1
1:A:44:GLU:CB	1:A:49:VAL:CG2	0.44	2.95	26	2
1:A:61:LEU:HD22	1:A:84:GLN:NE2	0.44	2.26	4	1
1:A:10:MET:HE3	1:A:48:PHE:CZ	0.44	2.47	12	1
1:A:5:ILE:HD13	1:A:12:ILE:HG21	0.44	1.88	22	1
1:A:13:THR:HG23	1:A:16:ILE:HD12	0.44	1.90	10	1
1:A:74:ILE:O	1:A:77:LEU:HG	0.44	2.13	24	1
1:A:67:HIS:CD2	1:A:72:THR:HB	0.44	2.48	28	1
1:A:67:HIS:CD2	1:A:72:THR:CB	0.44	3.00	27	1
1:A:6:THR:HG23	1:A:40:ILE:HG12	0.44	1.89	8	2
1:A:52:ALA:HB2	1:A:77:LEU:HD11	0.44	1.89	1	1
1:A:23:ARG:CD	1:A:78:ILE:HD11	0.44	2.42	12	1
1:A:23:ARG:HG3	1:A:78:ILE:CG2	0.44	2.42	25	1
1:A:34:LEU:HB2	1:A:54:ILE:HG22	0.44	1.89	26	1
1:A:62:VAL:N	1:A:81:LEU:HD21	0.44	2.27	26	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:56:THR:HB	1:A:57:PRO:HD2	0.44	1.90	9	2
1:A:44:GLU:HG3	1:A:49:VAL:HG21	0.44	1.90	3	2
1:A:80:LYS:HG3	1:A:81:LEU:N	0.44	2.27	6	1
1:A:34:LEU:HD23	1:A:56:THR:CG2	0.44	2.39	8	1
1:A:49:VAL:HG22	1:A:66:LYS:CE	0.44	2.43	23	1
1:A:32:THR:O	1:A:57:PRO:HG3	0.44	2.13	26	1
1:A:27:LEU:HD21	1:A:85:LEU:CD1	0.43	2.43	22	2
1:A:54:ILE:HD11	1:A:81:LEU:HD13	0.43	1.89	8	1
1:A:70:MET:SD	1:A:71:TYR:CD1	0.43	3.11	19	2
1:A:44:GLU:O	1:A:47:GLY:O	0.43	2.36	20	1
1:A:30:TRP:C	1:A:32:THR:N	0.43	2.76	22	9
1:A:10:MET:SD	1:A:41:LEU:HD23	0.43	2.53	3	1
1:A:24:LEU:HD11	1:A:37:PRO:HB2	0.43	1.89	3	1
1:A:20:VAL:HA	1:A:23:ARG:CG	0.43	2.42	5	1
1:A:34:LEU:N	1:A:34:LEU:CD2	0.43	2.80	22	3
1:A:32:THR:O	1:A:33:HIS:CD2	0.43	2.72	1	3
1:A:12:ILE:HG22	1:A:16:ILE:HG22	0.43	1.89	5	1
1:A:40:ILE:N	1:A:40:ILE:CD1	0.43	2.81	5	1
1:A:3:MET:SD	1:A:24:LEU:HD13	0.43	2.53	9	1
1:A:27:LEU:CD1	1:A:81:LEU:HD13	0.43	2.44	9	2
1:A:83:ARG:HG2	1:A:84:GLN:N	0.43	2.28	17	1
1:A:83:ARG:O	1:A:86:ASN:ND2	0.43	2.51	22	2
1:A:10:MET:SD	1:A:48:PHE:CZ	0.43	3.11	6	1
1:A:34:LEU:HD13	1:A:56:THR:CG2	0.43	2.22	22	1
1:A:34:LEU:O	1:A:34:LEU:HD23	0.43	2.14	26	2
1:A:32:THR:CG2	1:A:34:LEU:CD1	0.43	2.96	28	1
1:A:50:ALA:CB	1:A:77:LEU:HD12	0.43	2.40	1	1
1:A:23:ARG:CD	1:A:78:ILE:HD12	0.43	2.43	6	1
1:A:61:LEU:HB2	1:A:81:LEU:HD22	0.43	1.89	7	1
1:A:61:LEU:CD1	1:A:81:LEU:HD22	0.43	2.42	23	1
1:A:63:ALA:CB	1:A:81:LEU:HD23	0.43	2.44	25	1
1:A:34:LEU:N	1:A:34:LEU:CD1	0.43	2.79	28	1
1:A:30:TRP:CZ3	1:A:86:ASN:OD1	0.43	2.72	2	1
1:A:85:LEU:O	1:A:86:ASN:C	0.43	2.62	2	2
1:A:81:LEU:O	1:A:85:LEU:N	0.43	2.51	4	1
1:A:24:LEU:HD22	1:A:39:ILE:HD11	0.43	1.90	11	1
1:A:32:THR:O	1:A:57:PRO:HD2	0.43	2.14	13	1
1:A:48:PHE:CZ	1:A:70:MET:CE	0.43	3.01	24	1
1:A:88:LEU:HD23	1:A:88:LEU:C	0.43	2.38	26	1
1:A:30:TRP:CZ3	1:A:85:LEU:HD22	0.43	2.47	18	2
1:A:81:LEU:HD12	1:A:81:LEU:N	0.43	2.28	4	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:54:ILE:HD12	1:A:81:LEU:HD11	0.43	1.91	14	1
1:A:34:LEU:CD1	1:A:54:ILE:HG21	0.43	2.43	15	1
1:A:87:LYS:HG3	1:A:88:LEU:N	0.43	2.29	21	1
1:A:19:HIS:CD2	1:A:74:ILE:HG13	0.43	2.49	3	1
1:A:19:HIS:CE1	1:A:71:TYR:CD2	0.42	3.07	8	3
1:A:1:MET:HE1	1:A:28:GLU:HB2	0.42	1.91	16	1
1:A:69:ASP:O	1:A:73:ALA:HB2	0.42	2.14	22	1
1:A:23:ARG:HG3	1:A:78:ILE:HG21	0.42	1.91	25	1
1:A:12:ILE:N	1:A:12:ILE:CD1	0.42	2.74	22	2
1:A:2:THR:O	1:A:37:PRO:HG3	0.42	2.14	8	1
1:A:73:ALA:O	1:A:77:LEU:CD2	0.42	2.67	21	1
1:A:44:GLU:HB2	1:A:49:VAL:CG2	0.42	2.45	27	1
1:A:32:THR:OG1	1:A:34:LEU:HD22	0.42	2.15	20	1
1:A:4:ASN:O	1:A:39:ILE:HD12	0.42	2.13	28	1
1:A:14:PRO:O	1:A:18:GLN:HB2	0.42	2.15	5	1
1:A:44:GLU:HB3	1:A:49:VAL:HG22	0.42	1.90	9	1
1:A:54:ILE:HD11	1:A:81:LEU:CD1	0.42	2.45	23	1
1:A:63:ALA:HB2	1:A:81:LEU:HB2	0.42	1.90	14	2
1:A:74:ILE:O	1:A:78:ILE:HG22	0.42	2.15	7	1
1:A:27:LEU:HD23	1:A:30:TRP:HB2	0.42	1.91	26	2
1:A:61:LEU:CD2	1:A:81:LEU:CG	0.42	2.97	4	1
1:A:78:ILE:HG23	1:A:79:ASN:N	0.42	2.30	7	1
1:A:32:THR:OG1	1:A:57:PRO:HD2	0.42	2.14	17	1
1:A:19:HIS:ND1	1:A:71:TYR:CZ	0.42	2.87	20	1
1:A:87:LYS:C	1:A:87:LYS:HD2	0.42	2.40	22	1
1:A:63:ALA:CA	1:A:81:LEU:HD23	0.42	2.44	25	1
1:A:12:ILE:H	1:A:12:ILE:HD12	0.42	1.75	28	1
1:A:88:LEU:CD2	1:A:88:LEU:O	0.42	2.68	4	1
1:A:78:ILE:HD13	1:A:78:ILE:O	0.42	2.15	13	1
1:A:2:THR:OG1	1:A:3:MET:N	0.42	2.52	18	1
1:A:61:LEU:HD12	1:A:61:LEU:C	0.42	2.39	4	1
1:A:13:THR:OG1	1:A:16:ILE:HD12	0.42	2.15	10	1
1:A:61:LEU:HD13	1:A:84:GLN:HB3	0.42	1.90	12	1
1:A:60:VAL:C	1:A:61:LEU:HD23	0.42	2.39	15	1
1:A:1:MET:HE2	1:A:28:GLU:OE1	0.42	2.15	17	1
1:A:53:THR:HG22	1:A:60:VAL:HG11	0.42	1.92	19	1
1:A:32:THR:CG2	1:A:56:THR:CG2	0.42	2.97	21	1
1:A:63:ALA:N	1:A:81:LEU:HD23	0.41	2.30	1	1
1:A:76:GLU:O	1:A:79:ASN:OD1	0.41	2.38	7	1
1:A:30:TRP:CD2	1:A:85:LEU:CD1	0.41	3.03	15	1
1:A:20:VAL:HG21	1:A:39:ILE:HD12	0.41	1.90	25	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:27:LEU:CD1	1:A:30:TRP:HB2	0.41	2.46	10	1
1:A:48:PHE:HB2	1:A:67:HIS:O	0.41	2.15	18	1
1:A:27:LEU:C	1:A:27:LEU:CD1	0.41	2.92	27	1
1:A:79:ASN:ND2	1:A:80:LYS:N	0.41	2.68	28	1
1:A:20:VAL:HG23	1:A:21:ALA:N	0.41	2.31	3	1
1:A:78:ILE:C	1:A:78:ILE:CD1	0.41	2.93	14	1
1:A:8:LYS:O	1:A:9:GLN:C	0.41	2.63	18	1
1:A:34:LEU:N	1:A:34:LEU:HD22	0.41	2.30	21	1
1:A:56:THR:HG21	1:A:61:LEU:CG	0.41	2.44	27	1
1:A:78:ILE:CG2	1:A:79:ASN:N	0.41	2.84	28	1
1:A:87:LYS:CD	1:A:88:LEU:N	0.41	2.83	6	1
1:A:75:ASN:O	1:A:78:ILE:CG2	0.41	2.68	8	1
1:A:54:ILE:CG1	1:A:61:LEU:HB2	0.41	2.44	19	1
1:A:44:GLU:CB	1:A:49:VAL:HG22	0.41	2.44	9	1
1:A:27:LEU:HD12	1:A:32:THR:HG23	0.41	1.92	4	1
1:A:50:ALA:O	1:A:77:LEU:HD12	0.41	2.15	5	1
1:A:30:TRP:O	1:A:32:THR:N	0.41	2.52	16	2
1:A:8:LYS:C	1:A:9:GLN:HG3	0.41	2.41	18	1
1:A:7:SER:OG	1:A:41:LEU:O	0.41	2.37	23	1
1:A:27:LEU:HD13	1:A:28:GLU:N	0.41	2.29	27	1
1:A:54:ILE:O	1:A:60:VAL:HA	0.41	2.16	2	1
1:A:30:TRP:CH2	1:A:86:ASN:CG	0.41	2.98	2	1
1:A:44:GLU:CB	1:A:49:VAL:HG23	0.41	2.45	4	1
1:A:8:LYS:O	1:A:10:MET:N	0.41	2.54	13	1
1:A:6:THR:O	1:A:6:THR:CG2	0.41	2.68	26	1
1:A:34:LEU:HD11	1:A:56:THR:HG23	0.41	1.92	3	1
1:A:44:GLU:CB	1:A:47:GLY:O	0.41	2.69	7	1
1:A:12:ILE:HG22	1:A:17:ARG:CA	0.41	2.46	11	1
1:A:33:HIS:O	1:A:34:LEU:HB2	0.41	2.14	13	1
1:A:61:LEU:N	1:A:61:LEU:CD2	0.41	2.84	16	1
1:A:38:HIS:O	1:A:53:THR:N	0.41	2.52	23	1
1:A:32:THR:HG23	1:A:33:HIS:N	0.41	2.31	26	1
1:A:12:ILE:HD12	1:A:12:ILE:H	0.41	1.74	20	2
1:A:61:LEU:HB3	1:A:81:LEU:HD11	0.41	1.93	5	1
1:A:16:ILE:HA	1:A:71:TYR:OH	0.41	2.16	6	1
1:A:65:GLY:N	1:A:77:LEU:CD2	0.41	2.84	12	1
1:A:42:SER:O	1:A:48:PHE:CD1	0.41	2.74	20	1
1:A:27:LEU:HD12	1:A:34:LEU:HD12	0.41	1.93	21	1
1:A:81:LEU:HD13	1:A:81:LEU:O	0.41	2.15	25	1
1:A:70:MET:HG3	1:A:71:TYR:N	0.40	2.31	4	1
1:A:6:THR:HG23	1:A:6:THR:O	0.40	2.16	9	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:23:ARG:NE	1:A:77:LEU:HD11	0.40	2.30	9	1
1:A:33:HIS:H	1:A:34:LEU:HD22	0.40	1.76	14	1
1:A:10:MET:HE2	1:A:43:LYS:HB2	0.40	1.93	18	1
1:A:79:ASN:OD1	1:A:80:LYS:N	0.40	2.54	18	1
1:A:78:ILE:CG1	1:A:79:ASN:N	0.40	2.84	24	1
1:A:33:HIS:C	1:A:34:LEU:CD2	0.40	2.93	13	1
1:A:78:ILE:HG13	1:A:79:ASN:N	0.40	2.31	25	1
1:A:38:HIS:CB	1:A:53:THR:HG23	0.40	2.46	15	1
1:A:83:ARG:O	1:A:86:ASN:OD1	0.40	2.39	4	1
1:A:19:HIS:HB2	1:A:71:TYR:CE1	0.40	2.52	6	1
1:A:11:GLU:C	1:A:12:ILE:HD13	0.40	2.41	12	1
1:A:63:ALA:O	1:A:77:LEU:HD21	0.40	2.16	12	1
1:A:25:ALA:O	1:A:28:GLU:CG	0.40	2.69	18	1
1:A:63:ALA:CB	1:A:81:LEU:CD1	0.40	2.98	3	1
1:A:9:GLN:O	1:A:11:GLU:CG	0.40	2.70	9	1
1:A:32:THR:OG1	1:A:33:HIS:N	0.40	2.55	28	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	87/113 (77%)	71±2 (81±3%)	13±3 (15±3%)	3±1 (3±1%)	4	34
All	All	2523/3277 (77%)	2053 (81%)	383 (15%)	87 (3%)	4	34

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	34	LEU	27
1	A	11	GLU	11
1	A	8	LYS	10
1	A	3	MET	9
1	A	33	HIS	6
1	A	2	THR	6

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Mol	Chain	Res	Type	Models (Total)
1	A	30	TRP	4
1	A	31	GLN	4
1	A	60	VAL	3
1	A	9	GLN	2
1	A	87	LYS	1
1	A	82	GLU	1
1	A	10	MET	1
1	A	69	ASP	1
1	A	12	ILE	1

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	78/98 (80%)	47±3 (60±4%)	31±3 (40±4%)	0 5
All	All	2262/2842 (80%)	1353 (60%)	909 (40%)	0 5

All 70 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	38	HIS	29
1	A	74	ILE	27
1	A	19	HIS	26
1	A	64	SER	25
1	A	41	LEU	23
1	A	27	LEU	22
1	A	85	LEU	21
1	A	46	GLN	20
1	A	80	LYS	20
1	A	26	LYS	19
1	A	34	LEU	19
1	A	66	LYS	19
1	A	67	HIS	19
1	A	54	ILE	19
1	A	35	ILE	18
1	A	69	ASP	18

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Mol	Chain	Res	Type	Models (Total)
1	A	78	ILE	18
1	A	17	ARG	17
1	A	88	LEU	17
1	A	77	LEU	17
1	A	43	LYS	16
1	A	44	GLU	16
1	A	55	ASN	16
1	A	9	GLN	16
1	A	87	LYS	16
1	A	8	LYS	15
1	A	48	PHE	15
1	A	1	MET	15
1	A	70	MET	15
1	A	12	ILE	14
1	A	31	GLN	14
1	A	53	THR	14
1	A	32	THR	14
1	A	29	LYS	14
1	A	51	ASP	14
1	A	61	LEU	13
1	A	81	LEU	13
1	A	11	GLU	13
1	A	23	ARG	13
1	A	83	ARG	13
1	A	56	THR	12
1	A	28	GLU	12
1	A	2	THR	12
1	A	18	GLN	11
1	A	24	LEU	11
1	A	79	ASN	11
1	A	58	ASN	10
1	A	6	THR	10
1	A	3	MET	10
1	A	86	ASN	10
1	A	22	ASP	9
1	A	33	HIS	9
1	A	10	MET	7
1	A	82	GLU	7
1	A	60	VAL	7
1	A	76	GLU	7
1	A	72	THR	7
1	A	4	ASN	6

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Mol	Chain	Res	Type	Models (Total)
1	A	5	ILE	6
1	A	84	GLN	6
1	A	68	GLU	5
1	A	75	ASN	5
1	A	30	TRP	4
1	A	13	THR	3
1	A	62	VAL	3
1	A	20	VAL	2
1	A	49	VAL	2
1	A	16	ILE	1
1	A	39	ILE	1
1	A	36	ASN	1

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 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

No chemical shift data were provided