



Full wwPDB NMR Structure Validation Report ⓘ

Mar 5, 2026 – 02:24 PM UTC

PDB ID : 6ES6 / pdb_00006es6
BMRB ID : 34188
Title : Structure and dynamics conspire in the evolution of affinity between intrinsically disordered proteins
Authors : Chi, N.C.
Deposited on : 2017-10-19

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

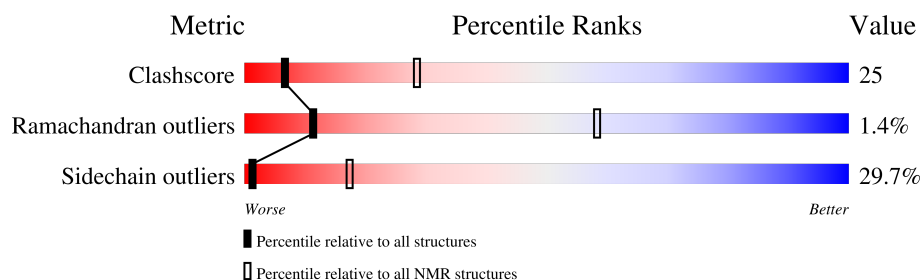
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 71%.

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	45	
2	B	50	

2 Ensemble composition and analysis

This entry contains 20 models. Model 13 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *target function*.

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:1044-A:1074, B:2062-B:2075, B:2081-B:2109 (74)	0.60	13

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 4 clusters and 3 single-model clusters were found.

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

3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 1463 atoms, of which 736 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called CID.

Mol	Chain	Residues	Atoms						Trace
1	A	45	Total	C	H	N	O	S	0
			658	200	323	55	79	1	

- Molecule 2 is a protein called NCBD.

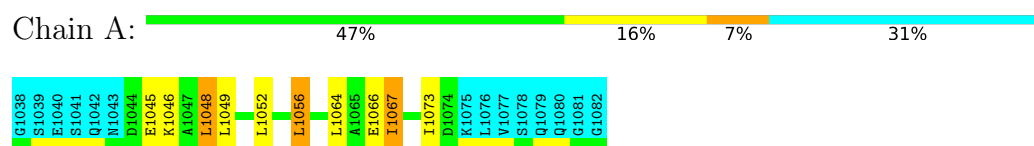
Mol	Chain	Residues	Atoms						Trace
2	B	50	Total	C	H	N	O	S	0
			805	246	413	74	71	1	

4 Residue-property plots [i](#)

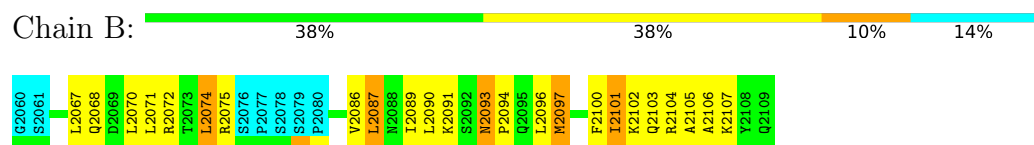
4.1 Average score per residue in the NMR ensemble

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: CID



• Molecule 2: NCBD

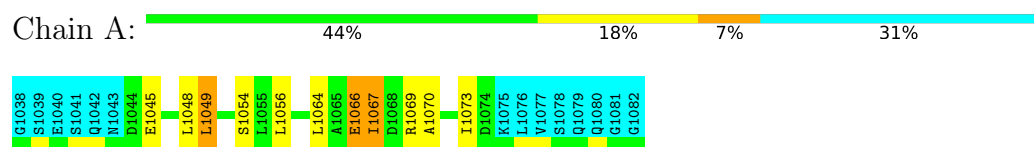


4.2 Scores per residue for each member of the ensemble

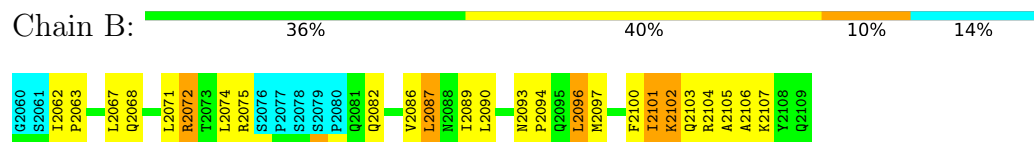
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

• Molecule 1: CID

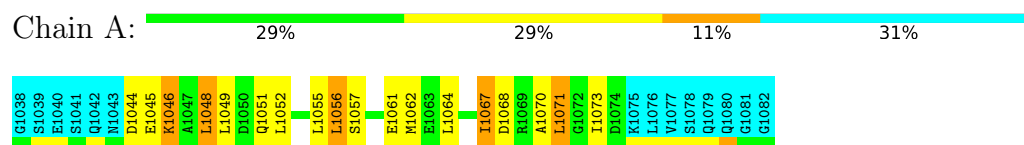


• Molecule 2: NCBD

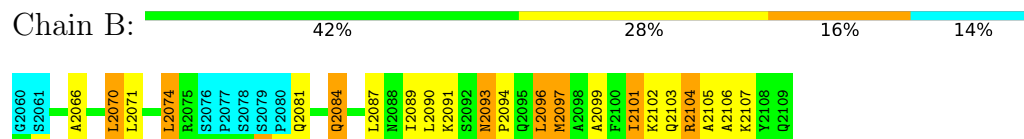


4.2.2 Score per residue for model 2

- Molecule 1: CID

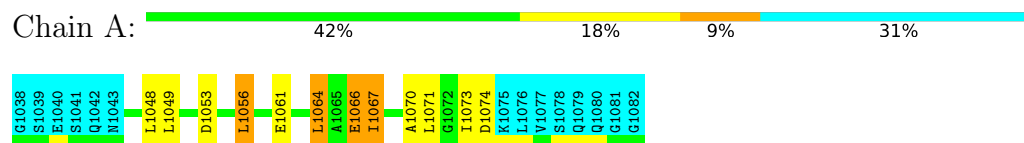


- Molecule 2: NCBD

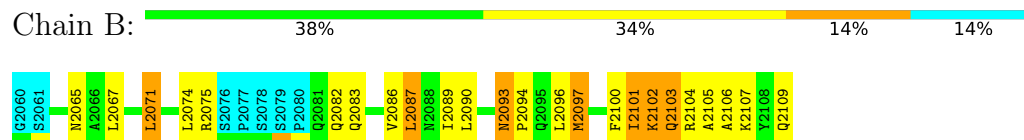


4.2.3 Score per residue for model 3

- Molecule 1: CID

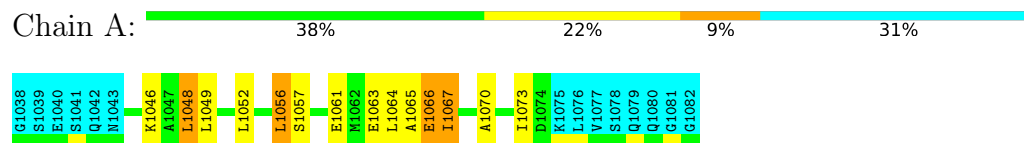


- Molecule 2: NCBD

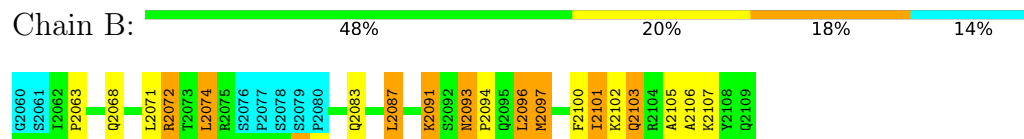


4.2.4 Score per residue for model 4

- Molecule 1: CID

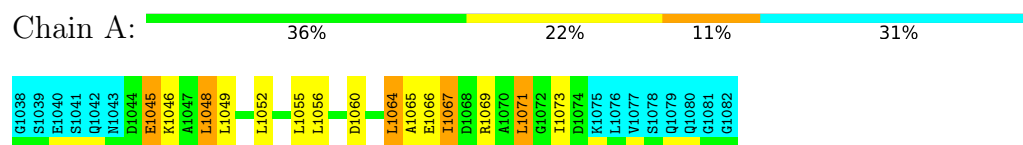


- Molecule 2: NCBD

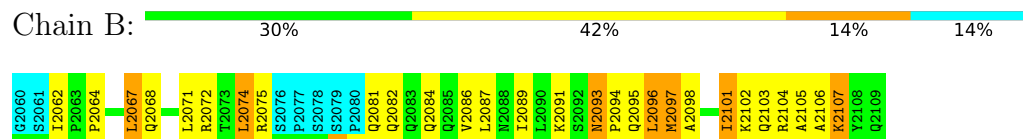


4.2.5 Score per residue for model 5

- Molecule 1: CID

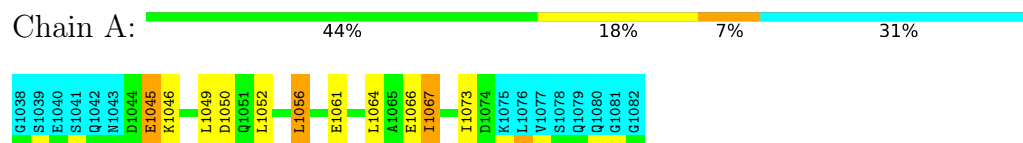


- Molecule 2: NCBD

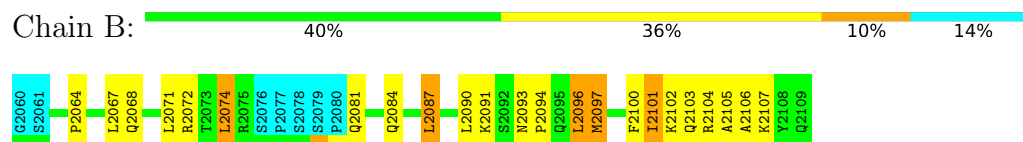


4.2.6 Score per residue for model 6

- Molecule 1: CID

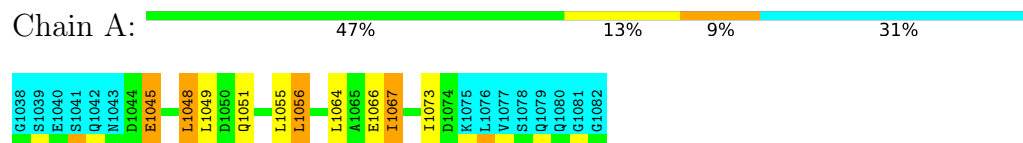


- Molecule 2: NCBD

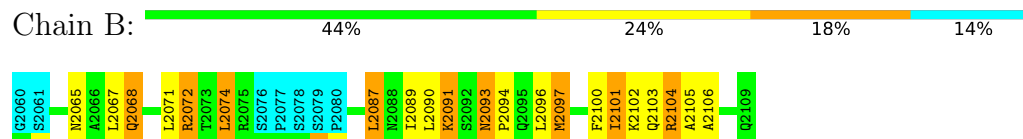


4.2.7 Score per residue for model 7

- Molecule 1: CID

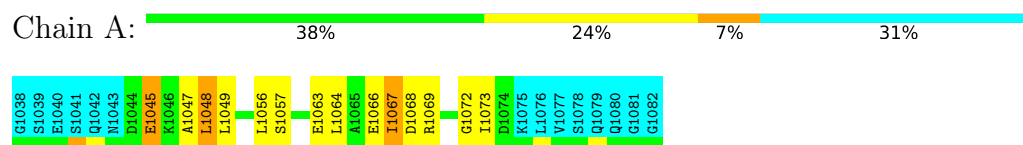


- Molecule 2: NCBD

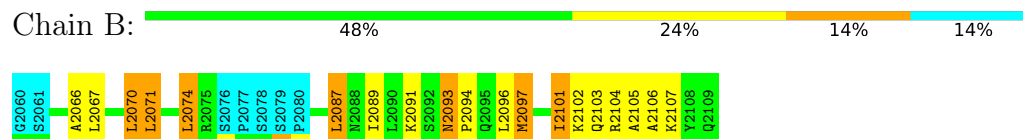


4.2.8 Score per residue for model 8

- Molecule 1: CID

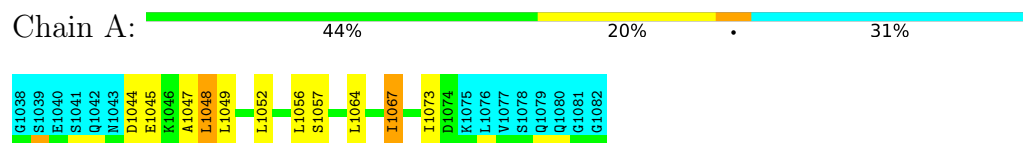


- Molecule 2: NCBD

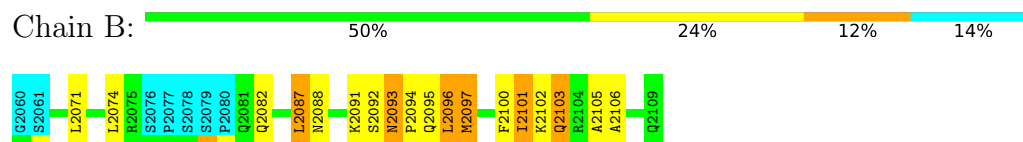


4.2.9 Score per residue for model 9

- Molecule 1: CID

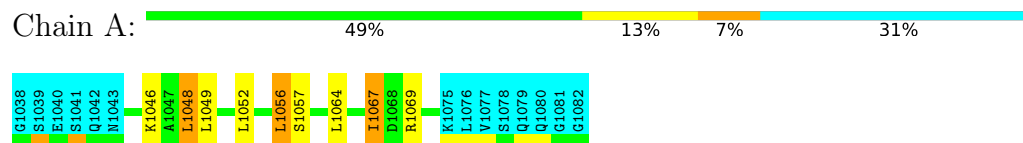


- Molecule 2: NCBD

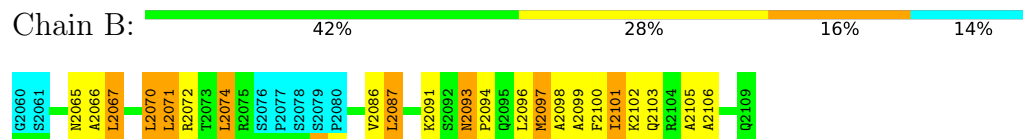


4.2.10 Score per residue for model 10

- Molecule 1: CID

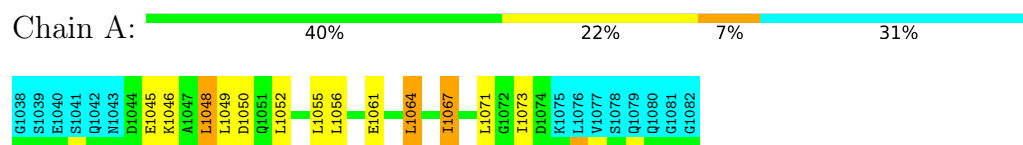


- Molecule 2: NCBD

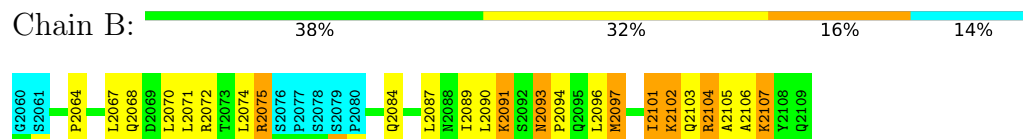


4.2.11 Score per residue for model 11

- Molecule 1: CID

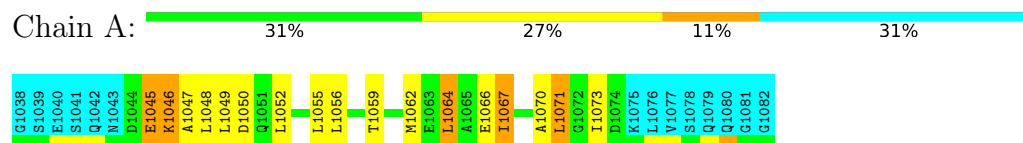


- Molecule 2: NCBD

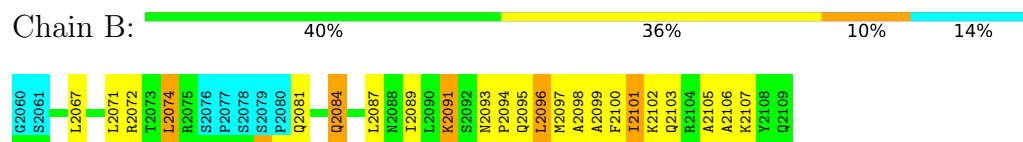


4.2.12 Score per residue for model 12

- Molecule 1: CID

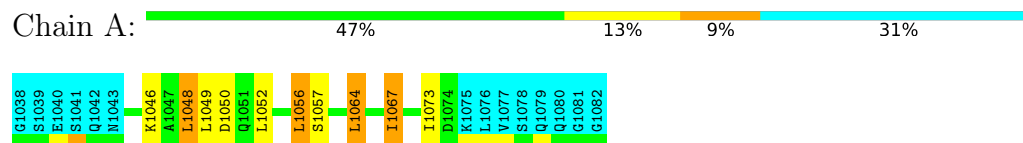


- Molecule 2: NCBD

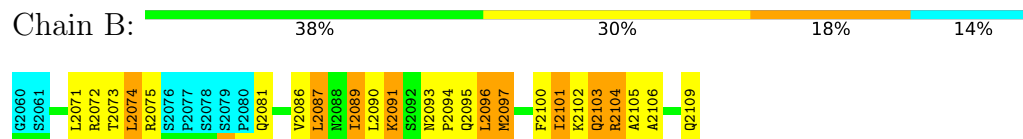


4.2.13 Score per residue for model 13 (medoid)

- Molecule 1: CID

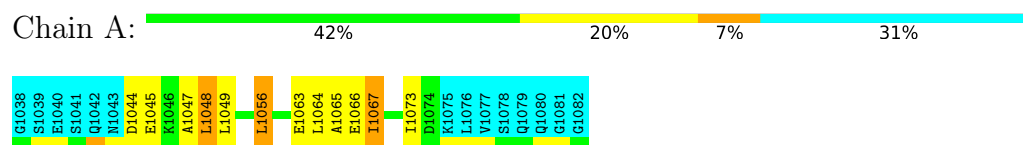


- Molecule 2: NCBD

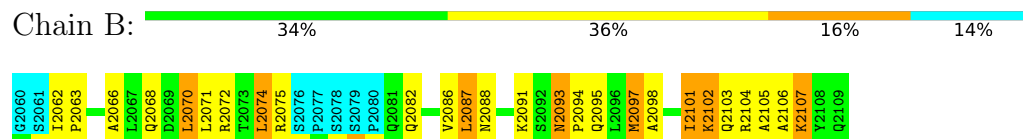


4.2.14 Score per residue for model 14

- Molecule 1: CID

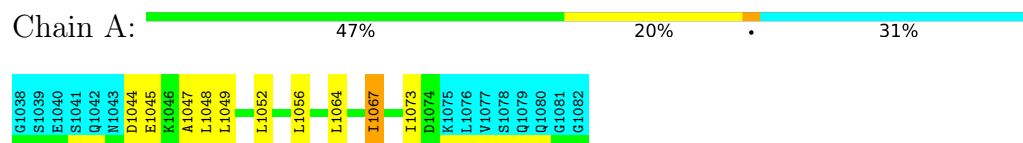


- Molecule 2: NCBD

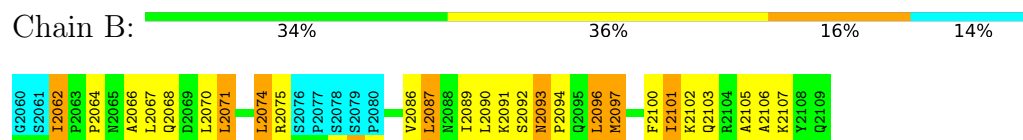


4.2.15 Score per residue for model 15

- Molecule 1: CID

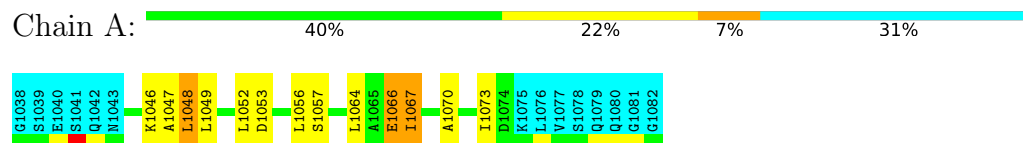


- Molecule 2: NCBD

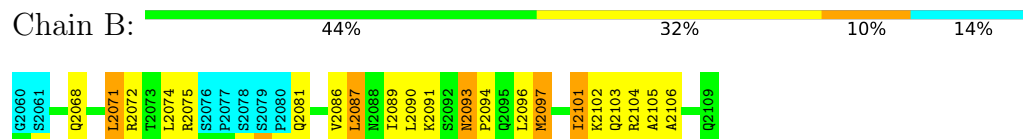


4.2.16 Score per residue for model 16

- Molecule 1: CID

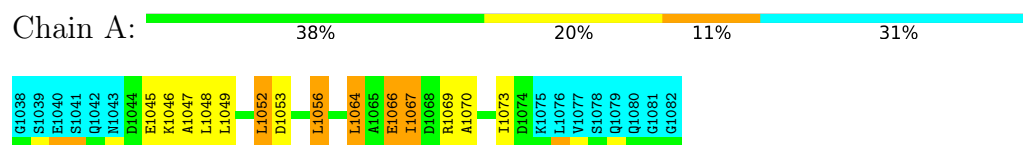


- Molecule 2: NCBD

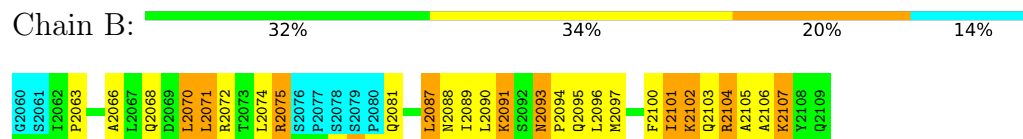


4.2.17 Score per residue for model 17

• Molecule 1: CID

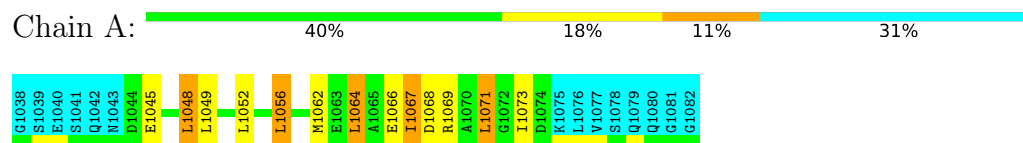


• Molecule 2: NCBD

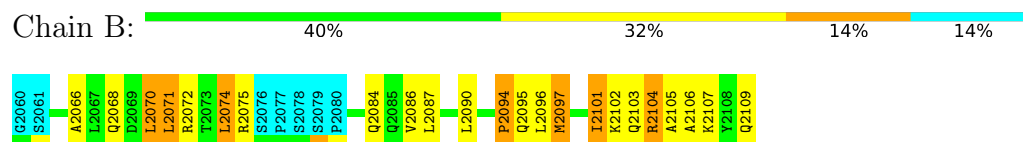


4.2.18 Score per residue for model 18

• Molecule 1: CID

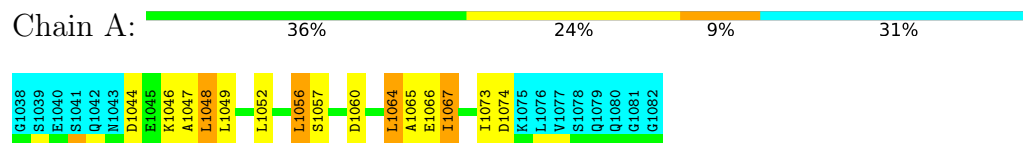


• Molecule 2: NCBD

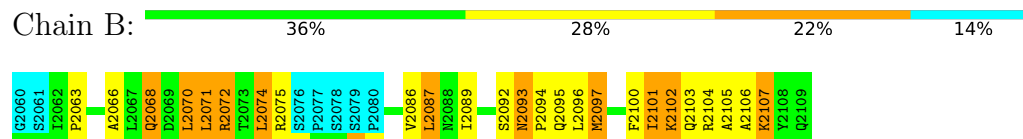


4.2.19 Score per residue for model 19

• Molecule 1: CID

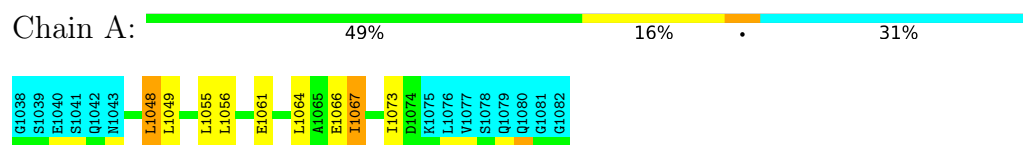


• Molecule 2: NCBD

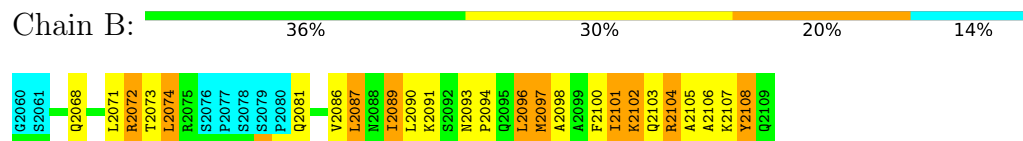


4.2.20 Score per residue for model 20

• Molecule 1: CID



• Molecule 2: NCBD



5 Refinement protocol and experimental data overview

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

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	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	896
Number of shifts mapped to atoms	896
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	71%

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	237	230	230	18±3
2	B	350	376	376	25±4
All	All	11740	12120	12120	592

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:1056:LEU:HD11	1:A:1064:LEU:HD21	1.03	1.28	12	5
2:B:2071:LEU:HD23	2:B:2096:LEU:HD11	0.98	1.33	6	1
1:A:1049:LEU:HD21	2:B:2074:LEU:HD21	0.98	1.34	3	3
2:B:2101:ILE:O	2:B:2105:ALA:HB3	0.79	1.76	18	20
1:A:1049:LEU:HD13	2:B:2071:LEU:HD13	0.74	1.59	1	6
1:A:1049:LEU:HD22	2:B:2096:LEU:HD11	0.74	1.59	12	1
1:A:1073:ILE:HG23	2:B:2097:MET:HE2	0.73	1.60	18	2
1:A:1056:LEU:HD21	1:A:1064:LEU:HD21	0.72	1.60	16	4
1:A:1049:LEU:HD11	2:B:2074:LEU:HD12	0.69	1.64	13	2
1:A:1049:LEU:HD11	2:B:2074:LEU:HB3	0.68	1.63	8	9
1:A:1049:LEU:HD11	2:B:2074:LEU:CB	0.68	2.18	10	4
1:A:1049:LEU:CD2	2:B:2096:LEU:HD11	0.68	2.18	12	1
2:B:2102:LYS:O	2:B:2106:ALA:HB2	0.67	1.89	5	20

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:2090:LEU:HD11	2:B:2104:ARG:HD2	0.67	1.65	2	1
2:B:2084:GLN:O	2:B:2087:LEU:HD12	0.67	1.90	18	5
1:A:1049:LEU:HD11	2:B:2074:LEU:HD11	0.66	1.67	3	2
2:B:2067:LEU:HD13	2:B:2093:ASN:ND2	0.66	2.05	10	2
2:B:2066:ALA:O	2:B:2070:LEU:HD22	0.65	1.91	10	8
1:A:1056:LEU:HD13	1:A:1056:LEU:O	0.64	1.91	16	9
2:B:2093:ASN:ND2	2:B:2096:LEU:HD12	0.64	2.07	12	5
2:B:2074:LEU:HD13	2:B:2074:LEU:O	0.64	1.92	17	1
1:A:1065:ALA:HB2	2:B:2107:LYS:HE3	0.64	1.70	5	1
1:A:1048:LEU:HD23	2:B:2071:LEU:HD21	0.64	1.68	16	14
1:A:1049:LEU:HD21	2:B:2074:LEU:HB2	0.64	1.69	7	4
1:A:1049:LEU:HD21	2:B:2074:LEU:HD23	0.64	1.70	6	1
1:A:1045:GLU:CD	1:A:1048:LEU:HD22	0.63	2.18	1	1
1:A:1071:LEU:HD21	2:B:2087:LEU:HB3	0.63	1.70	5	4
1:A:1068:ASP:OD2	1:A:1073:ILE:HD13	0.62	1.94	18	1
1:A:1048:LEU:HD22	2:B:2063:PRO:HG3	0.62	1.70	14	4
2:B:2074:LEU:HD12	2:B:2086:VAL:HG11	0.62	1.71	14	2
1:A:1049:LEU:HD21	2:B:2074:LEU:CD2	0.61	2.20	3	3
2:B:2064:PRO:CG	2:B:2067:LEU:HD22	0.61	2.25	5	2
1:A:1056:LEU:O	1:A:1059:THR:HG22	0.61	1.95	12	1
1:A:1049:LEU:HB3	2:B:2096:LEU:HD21	0.61	1.71	20	6
2:B:2103:GLN:O	2:B:2106:ALA:HB3	0.60	1.96	3	20
1:A:1049:LEU:HD11	2:B:2074:LEU:HD22	0.60	1.71	19	3
1:A:1056:LEU:O	1:A:1056:LEU:HD13	0.60	1.96	12	1
1:A:1056:LEU:CD1	1:A:1064:LEU:HD11	0.60	2.26	19	7
1:A:1052:LEU:HD13	2:B:2075:ARG:HB2	0.60	1.72	18	6
1:A:1049:LEU:CD2	2:B:2074:LEU:HD21	0.60	2.21	3	2
2:B:2074:LEU:HD22	2:B:2100:PHE:CE2	0.60	2.32	6	1
1:A:1044:ASP:OD2	1:A:1047:ALA:HB2	0.60	1.96	14	1
2:B:2074:LEU:C	2:B:2074:LEU:HD13	0.59	2.22	13	2
1:A:1056:LEU:CD1	1:A:1064:LEU:HD21	0.59	2.27	9	3
2:B:2074:LEU:HD23	2:B:2086:VAL:HG11	0.59	1.73	13	1
1:A:1067:ILE:HD11	2:B:2090:LEU:CD1	0.59	2.28	11	7
1:A:1045:GLU:HG2	2:B:2067:LEU:HD23	0.59	1.75	5	2
1:A:1073:ILE:HG21	2:B:2101:ILE:CD1	0.58	2.28	19	1
1:A:1067:ILE:HA	2:B:2087:LEU:HD21	0.58	1.73	11	4
1:A:1073:ILE:HD12	2:B:2097:MET:HE1	0.58	1.75	8	2
2:B:2108:TYR:C	2:B:2108:TYR:CD2	0.58	2.80	20	1
1:A:1049:LEU:CD1	2:B:2074:LEU:HD12	0.58	2.28	13	1
2:B:2090:LEU:HD13	2:B:2104:ARG:NH1	0.57	2.14	13	4
1:A:1048:LEU:HD22	2:B:2071:LEU:HD11	0.57	1.76	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:1048:LEU:HD22	2:B:2068:GLN:OE1	0.57	1.99	11	1
1:A:1045:GLU:OE2	2:B:2071:LEU:HD21	0.57	2.00	1	1
2:B:2093:ASN:ND2	2:B:2096:LEU:HD13	0.57	2.15	6	1
1:A:1064:LEU:HD22	2:B:2104:ARG:HG2	0.56	1.77	14	4
1:A:1065:ALA:HB2	2:B:2107:LYS:HD3	0.56	1.77	19	3
1:A:1049:LEU:HD13	2:B:2071:LEU:CD1	0.56	2.30	20	9
1:A:1056:LEU:HD11	1:A:1064:LEU:HD13	0.56	1.77	18	1
2:B:2074:LEU:HD12	2:B:2086:VAL:CG1	0.56	2.29	19	2
1:A:1049:LEU:HD13	2:B:2071:LEU:HB2	0.56	1.77	3	12
1:A:1056:LEU:CD1	1:A:1064:LEU:HD13	0.56	2.31	18	1
2:B:2074:LEU:HD22	2:B:2086:VAL:HG13	0.55	1.78	20	2
1:A:1073:ILE:HG22	2:B:2101:ILE:HD11	0.55	1.77	20	3
2:B:2093:ASN:O	2:B:2097:MET:HB2	0.55	2.01	6	13
2:B:2070:LEU:O	2:B:2089:ILE:HD12	0.55	2.00	11	1
1:A:1067:ILE:HD11	2:B:2090:LEU:HD11	0.55	1.78	18	2
1:A:1049:LEU:HD11	2:B:2074:LEU:CD1	0.55	2.31	3	2
1:A:1052:LEU:HD22	2:B:2075:ARG:CD	0.55	2.31	18	2
2:B:2090:LEU:HD13	2:B:2104:ARG:HE	0.54	1.63	18	1
1:A:1067:ILE:HD11	2:B:2090:LEU:HD12	0.54	1.79	11	3
1:A:1064:LEU:HD21	2:B:2104:ARG:HG2	0.54	1.80	11	2
1:A:1071:LEU:HD23	2:B:2104:ARG:HH22	0.53	1.63	3	1
1:A:1067:ILE:HB	2:B:2087:LEU:HD13	0.53	1.80	13	13
1:A:1073:ILE:CD1	2:B:2097:MET:HE1	0.53	2.33	8	1
2:B:2071:LEU:HD12	2:B:2072:ARG:N	0.53	2.19	12	4
1:A:1064:LEU:HD13	2:B:2107:LYS:HD3	0.53	1.79	5	1
1:A:1048:LEU:HD22	2:B:2063:PRO:CG	0.53	2.34	14	1
2:B:2074:LEU:CD2	2:B:2086:VAL:HG13	0.53	2.34	15	2
1:A:1049:LEU:HD22	2:B:2071:LEU:HB2	0.53	1.81	15	12
1:A:1067:ILE:HB	2:B:2087:LEU:HD22	0.52	1.82	16	7
1:A:1045:GLU:OE2	2:B:2067:LEU:HD12	0.52	2.04	12	1
1:A:1052:LEU:HD22	2:B:2075:ARG:HD3	0.52	1.80	18	1
1:A:1046:LYS:CG	2:B:2099:ALA:HB2	0.52	2.35	12	1
1:A:1049:LEU:HD11	2:B:2074:LEU:CD2	0.52	2.35	19	2
2:B:2074:LEU:HD22	2:B:2074:LEU:O	0.52	2.05	13	2
2:B:2071:LEU:HD12	2:B:2071:LEU:C	0.51	2.30	16	16
1:A:1049:LEU:HD21	2:B:2074:LEU:CB	0.51	2.35	9	6
2:B:2074:LEU:HD22	2:B:2090:LEU:CD2	0.51	2.34	11	1
2:B:2093:ASN:CG	2:B:2096:LEU:HD12	0.51	2.30	3	5
1:A:1049:LEU:HD11	2:B:2074:LEU:HD23	0.51	1.81	6	1
1:A:1073:ILE:CG2	2:B:2101:ILE:HD11	0.51	2.36	13	2
2:B:2064:PRO:CG	2:B:2067:LEU:HD13	0.51	2.35	15	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:2074:LEU:HB3	2:B:2086:VAL:HG13	0.51	1.82	3	3
1:A:1073:ILE:HG22	1:A:1073:ILE:O	0.51	2.05	13	6
1:A:1049:LEU:CD1	2:B:2074:LEU:HD22	0.51	2.35	12	2
1:A:1067:ILE:HD13	2:B:2104:ARG:CZ	0.50	2.36	5	1
1:A:1071:LEU:HD23	2:B:2104:ARG:NH2	0.50	2.21	3	1
1:A:1056:LEU:HD13	1:A:1056:LEU:C	0.50	2.30	12	5
2:B:2074:LEU:HD22	2:B:2100:PHE:CZ	0.50	2.41	6	1
1:A:1056:LEU:HD11	1:A:1064:LEU:HD11	0.50	1.82	1	5
2:B:2093:ASN:OD1	2:B:2096:LEU:HD12	0.50	2.06	3	4
2:B:2101:ILE:HG12	2:B:2105:ALA:HB2	0.50	1.83	5	5
2:B:2071:LEU:CD2	2:B:2096:LEU:HD11	0.50	2.37	11	2
1:A:1049:LEU:CD1	2:B:2074:LEU:HD11	0.50	2.37	3	2
2:B:2074:LEU:HA	2:B:2086:VAL:HG22	0.49	1.84	16	3
2:B:2067:LEU:HD13	2:B:2093:ASN:HD21	0.49	1.66	10	1
1:A:1056:LEU:CD2	1:A:1064:LEU:HD21	0.49	2.36	20	1
2:B:2071:LEU:HD23	2:B:2096:LEU:CD1	0.49	2.23	6	1
1:A:1056:LEU:HD23	2:B:2100:PHE:CZ	0.49	2.42	13	6
1:A:1073:ILE:HG13	2:B:2097:MET:HE1	0.48	1.85	13	2
1:A:1064:LEU:HD23	2:B:2107:LYS:HB3	0.48	1.84	15	1
2:B:2074:LEU:HD22	2:B:2086:VAL:HG22	0.48	1.83	18	1
1:A:1071:LEU:CD2	2:B:2087:LEU:HD23	0.48	2.38	11	1
2:B:2097:MET:HE3	2:B:2104:ARG:CZ	0.48	2.38	18	1
1:A:1045:GLU:OE1	1:A:1048:LEU:HD22	0.48	2.08	1	1
2:B:2086:VAL:HG11	2:B:2104:ARG:NH2	0.48	2.23	5	1
1:A:1045:GLU:CD	2:B:2067:LEU:HD12	0.48	2.33	8	2
2:B:2074:LEU:HD13	2:B:2074:LEU:C	0.48	2.33	17	1
1:A:1066:GLU:O	1:A:1070:ALA:HB2	0.48	2.09	3	5
1:A:1056:LEU:HD12	1:A:1064:LEU:HD21	0.47	1.86	4	3
1:A:1056:LEU:HD11	1:A:1064:LEU:CD1	0.47	2.39	1	2
1:A:1067:ILE:HD12	2:B:2087:LEU:CD2	0.47	2.38	7	7
2:B:2068:GLN:O	2:B:2072:ARG:CB	0.47	2.63	17	9
2:B:2064:PRO:HG3	2:B:2067:LEU:HD13	0.47	1.85	15	1
1:A:1067:ILE:O	1:A:1070:ALA:HB3	0.47	2.10	12	2
1:A:1049:LEU:HD22	2:B:2071:LEU:HA	0.47	1.86	8	4
1:A:1049:LEU:HD21	2:B:2074:LEU:HD11	0.47	1.86	11	1
2:B:2094:PRO:O	2:B:2097:MET:HB3	0.47	2.09	18	1
1:A:1073:ILE:O	1:A:1073:ILE:HG22	0.47	2.10	7	7
1:A:1073:ILE:HD12	2:B:2097:MET:SD	0.46	2.50	19	5
2:B:2093:ASN:CG	2:B:2096:LEU:HB2	0.46	2.34	12	1
2:B:2086:VAL:HG11	2:B:2104:ARG:HH21	0.46	1.70	5	1
2:B:2097:MET:HE3	2:B:2097:MET:HA	0.46	1.86	20	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:2101:ILE:O	2:B:2105:ALA:CB	0.46	2.59	18	8
2:B:2074:LEU:HD23	2:B:2074:LEU:C	0.46	2.36	14	3
1:A:1049:LEU:HD22	2:B:2071:LEU:CB	0.45	2.41	6	4
2:B:2090:LEU:HD21	2:B:2100:PHE:CE2	0.45	2.46	17	1
1:A:1073:ILE:HG21	2:B:2101:ILE:HG13	0.45	1.88	1	6
2:B:2062:ILE:HD12	2:B:2062:ILE:O	0.45	2.11	15	3
1:A:1067:ILE:O	1:A:1071:LEU:HD22	0.45	2.11	12	3
1:A:1056:LEU:HD21	2:B:2104:ARG:NE	0.45	2.26	2	1
2:B:2090:LEU:HD21	2:B:2100:PHE:CE1	0.45	2.47	13	4
1:A:1052:LEU:HD22	2:B:2075:ARG:CB	0.45	2.42	11	1
2:B:2093:ASN:HB3	2:B:2096:LEU:HB3	0.45	1.87	1	1
1:A:1067:ILE:HD12	2:B:2087:LEU:HD22	0.45	1.88	7	6
1:A:1052:LEU:HD13	2:B:2075:ARG:HB3	0.45	1.87	13	1
2:B:2073:THR:HG21	2:B:2089:ILE:HG13	0.45	1.88	20	2
1:A:1064:LEU:HD12	1:A:1064:LEU:O	0.45	2.12	18	1
1:A:1048:LEU:HD11	2:B:2063:PRO:HG3	0.45	1.88	1	1
2:B:2074:LEU:O	2:B:2074:LEU:HD22	0.44	2.11	17	1
1:A:1049:LEU:CG	2:B:2074:LEU:HD22	0.44	2.41	12	1
1:A:1049:LEU:HD22	2:B:2096:LEU:HD13	0.44	1.89	3	2
1:A:1073:ILE:HG12	2:B:2097:MET:HE1	0.44	1.89	15	1
2:B:2070:LEU:HA	2:B:2089:ILE:HD12	0.44	1.90	2	1
2:B:2090:LEU:HB3	2:B:2097:MET:HE3	0.44	1.89	15	1
2:B:2074:LEU:HD13	2:B:2100:PHE:CE2	0.44	2.48	19	1
1:A:1056:LEU:HD11	2:B:2104:ARG:CZ	0.44	2.43	2	1
1:A:1064:LEU:HD13	2:B:2107:LYS:HB3	0.44	1.89	11	2
1:A:1071:LEU:HD21	2:B:2087:LEU:CB	0.43	2.42	18	2
1:A:1056:LEU:HD12	1:A:1064:LEU:HD11	0.43	1.89	17	2
1:A:1073:ILE:HG21	2:B:2101:ILE:CG1	0.43	2.43	17	3
1:A:1056:LEU:HD11	1:A:1064:LEU:CD2	0.43	2.36	9	1
1:A:1073:ILE:HD12	1:A:1073:ILE:N	0.43	2.28	11	1
1:A:1067:ILE:CB	2:B:2087:LEU:HD22	0.43	2.44	16	2
2:B:2074:LEU:CB	2:B:2086:VAL:HG22	0.43	2.43	3	1
1:A:1056:LEU:HD12	2:B:2103:GLN:HB3	0.43	1.90	12	2
1:A:1049:LEU:HD21	2:B:2074:LEU:HB3	0.43	1.91	13	2
1:A:1071:LEU:HD21	2:B:2087:LEU:HD23	0.43	1.90	11	1
1:A:1068:ASP:O	1:A:1072:GLY:N	0.42	2.52	8	1
1:A:1064:LEU:HD22	2:B:2107:LYS:HB3	0.42	1.91	17	2
2:B:2096:LEU:O	2:B:2100:PHE:HB2	0.42	2.14	1	4
1:A:1049:LEU:HD11	2:B:2075:ARG:HB2	0.42	1.92	1	1
1:A:1073:ILE:CG2	2:B:2101:ILE:CG1	0.42	2.97	8	3
1:A:1046:LYS:CE	2:B:2099:ALA:HB1	0.42	2.44	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:1048:LEU:HB3	2:B:2071:LEU:HD22	0.42	1.91	10	1
2:B:2108:TYR:O	2:B:2108:TYR:CG	0.42	2.72	20	1
2:B:2074:LEU:C	2:B:2074:LEU:CD1	0.42	2.93	13	1
2:B:2101:ILE:HD12	2:B:2101:ILE:HA	0.42	1.68	19	1
1:A:1049:LEU:CB	2:B:2096:LEU:HD21	0.42	2.44	2	1
1:A:1056:LEU:HD21	1:A:1064:LEU:CD2	0.42	2.43	20	1
1:A:1048:LEU:HD13	2:B:2071:LEU:HD21	0.42	1.91	10	1
1:A:1045:GLU:HG3	2:B:2067:LEU:HD22	0.42	1.92	15	1
1:A:1046:LYS:HG2	2:B:2099:ALA:HB2	0.41	1.91	2	1
2:B:2074:LEU:CA	2:B:2086:VAL:HG22	0.41	2.45	16	1
1:A:1064:LEU:HD13	2:B:2107:LYS:CB	0.41	2.45	11	1
1:A:1073:ILE:O	1:A:1073:ILE:CG2	0.41	2.68	18	2
1:A:1056:LEU:HD13	1:A:1064:LEU:HD11	0.41	1.92	19	1
2:B:2083:GLN:OE1	2:B:2087:LEU:HD21	0.41	2.15	3	1
1:A:1073:ILE:CG2	2:B:2105:ALA:HB2	0.41	2.46	2	1
2:B:2064:PRO:HG2	2:B:2067:LEU:HD13	0.41	1.92	11	1
2:B:2074:LEU:CB	2:B:2086:VAL:HG13	0.41	2.45	16	1
2:B:2093:ASN:HD22	2:B:2096:LEU:HD12	0.40	1.75	12	1
1:A:1065:ALA:HB2	2:B:2107:LYS:CE	0.40	2.44	5	1
1:A:1071:LEU:HD22	1:A:1071:LEU:N	0.40	2.32	12	2
1:A:1073:ILE:HD12	2:B:2097:MET:CE	0.40	2.47	2	1
1:A:1052:LEU:HD22	2:B:2075:ARG:HB2	0.40	1.93	11	1

6.3 Torsion angles

6.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the backbone conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	31/45 (69%)	29±1 (92±3%)	2±1 (8±3%)	0±0 (0±0%)	100	100
2	B	42/50 (84%)	36±1 (86±2%)	5±1 (12±2%)	1±0 (2±0%)	7	44
All	All	1460/1900 (77%)	1295 (89%)	145 (10%)	20 (1%)	11	58

All 1 unique Ramachandran outliers are listed below.

Mol	Chain	Res	Type	Models (Total)
2	B	2094	PRO	20

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	27/38 (71%)	19±3 (71±10%)	8±3 (29±10%)	1	18
2	B	38/44 (86%)	26±2 (70±6%)	12±2 (30±6%)	1	16
All	All	1300/1640 (79%)	914 (70%)	386 (30%)	1	17

All 53 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	1067	ILE	20
2	B	2101	ILE	20
2	B	2097	MET	19
1	A	1048	LEU	16
2	B	2087	LEU	15
2	B	2093	ASN	15
1	A	1066	GLU	14
2	B	2074	LEU	14
2	B	2091	LYS	14
2	B	2107	LYS	13
2	B	2089	ILE	12
2	B	2096	LEU	12
1	A	1056	LEU	12
1	A	1045	GLU	11
1	A	1046	LYS	10
1	A	1064	LEU	9
2	B	2104	ARG	9
2	B	2072	ARG	8
1	A	1052	LEU	8
1	A	1057	SER	8
2	B	2081	GLN	8
2	B	2071	LEU	8
2	B	2095	GLN	8
2	B	2102	LYS	7

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Mol	Chain	Res	Type	Models (Total)
2	B	2070	LEU	7
1	A	1069	ARG	6
1	A	1055	LEU	6
1	A	1061	GLU	6
2	B	2068	GLN	5
2	B	2082	GLN	5
2	B	2075	ARG	5
1	A	1044	ASP	4
1	A	1071	LEU	4
2	B	2103	GLN	4
1	A	1050	ASP	4
2	B	2067	LEU	3
1	A	1062	MET	3
2	B	2084	GLN	3
1	A	1053	ASP	3
2	B	2065	ASN	3
2	B	2109	GLN	3
1	A	1063	GLU	3
2	B	2088	ASN	3
2	B	2092	SER	3
1	A	1051	GLN	2
1	A	1074	ASP	2
1	A	1060	ASP	2
2	B	2062	ILE	2
1	A	1049	LEU	1
1	A	1054	SER	1
1	A	1068	ASP	1
2	B	2083	GLN	1
2	B	2108	TYR	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.6 Ligand geometry

There are no ligands in this entry.

6.7 Other polymers

There are no such molecules in this entry.

6.8 Polymer linkage issues

There are no chain breaks in this entry.

7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 71% for the well-defined parts and 70% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *2R_1R.prot*

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	896
Number of shifts mapped to atoms	896
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	89	-0.56 ± 0.09	Should be checked
$^{13}\text{C}_\beta$	87	0.24 ± 0.07	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	None (insufficient data)
^{15}N	86	-0.41 ± 0.15	None needed (< 0.5 ppm)

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 71%, i.e. 743 atoms were assigned a chemical shift out of a possible 1051. 0 out of 15 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	285/365 (78%)	145/146 (99%)	69/148 (47%)	71/71 (100%)
Sidechain	456/667 (68%)	344/432 (80%)	111/206 (54%)	1/29 (3%)

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	Total	¹H	¹³C	¹⁵N
Aromatic	2/19 (11%)	2/9 (22%)	0/10 (0%)	0/0 (—%)
Overall	743/1051 (71%)	491/587 (84%)	180/364 (49%)	72/100 (72%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 70%, i.e. 893 atoms were assigned a chemical shift out of a possible 1275. 0 out of 17 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	356/470 (76%)	181/190 (95%)	89/190 (47%)	86/90 (96%)
Sidechain	535/786 (68%)	402/508 (79%)	132/244 (54%)	1/34 (3%)
Aromatic	2/19 (11%)	2/9 (22%)	0/10 (0%)	0/0 (—%)
Overall	893/1275 (70%)	585/707 (83%)	221/444 (50%)	87/124 (70%)

7.1.4 Statistically unusual chemical shifts ⓘ

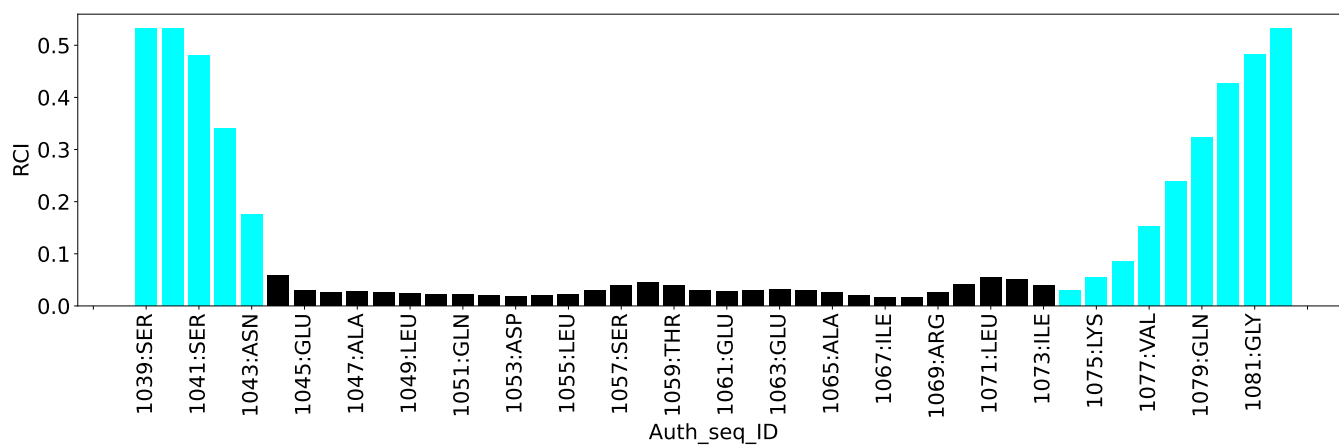
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	1059	THR	HG1	6.76	0.08 – 2.19	26.7

7.1.5 Random Coil Index (RCI) plots ⓘ

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:



Random coil index (RCI) for chain B:

