



Full wwPDB NMR Structure Validation Report ⓘ

Feb 20, 2022 – 02:17 AM EST

PDB ID : 1TNW
Title : NMR SOLUTION STRUCTURE OF CALCIUM SATURATED SKELETAL
MUSCLE TROPONIN C
Authors : Slupsky, C.M.; Sykes, B.D.
Deposited on : 1995-08-23

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 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)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : 2.26
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.26

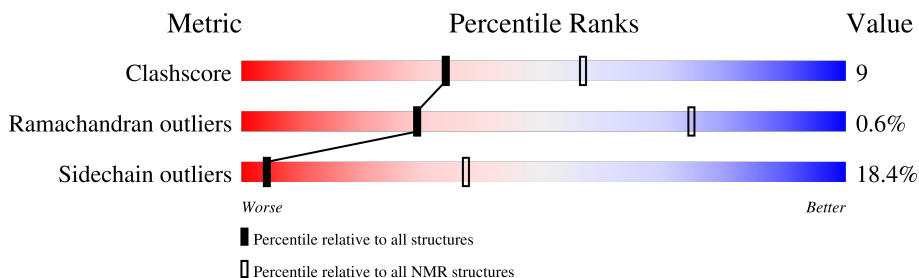
1 Overall quality at a glance

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	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	162	 63% 23% • 12%

2 Ensemble composition and analysis

This entry contains 23 models. Model 22 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:5-A:84 (80)	0.48	22
2	A:96-A:158 (63)	0.52	17

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 2 single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2471 atoms, of which 1199 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called TROPONIN C.

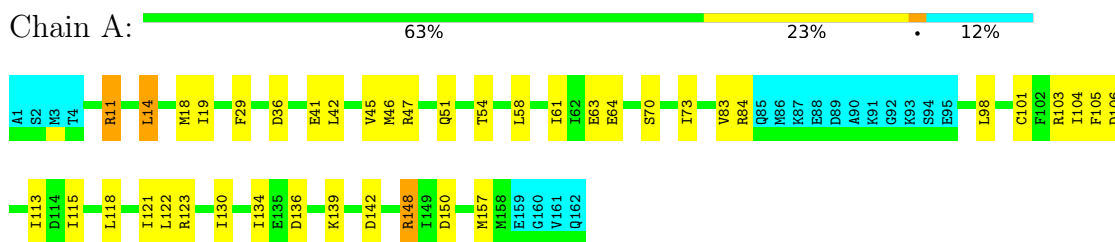
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	162	2471	783	1199	202	275	12	0

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: TROPONIN C

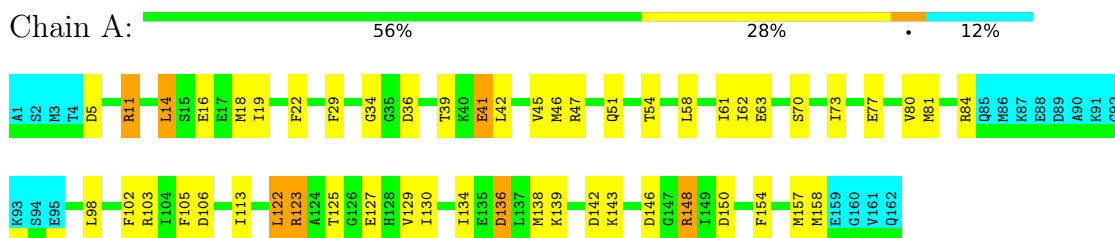


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

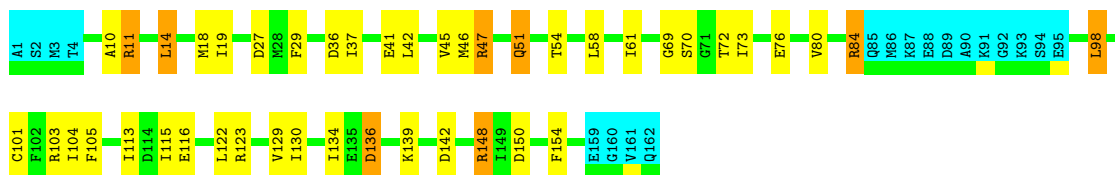
- Molecule 1: TROPONIN C



4.2.2 Score per residue for model 2

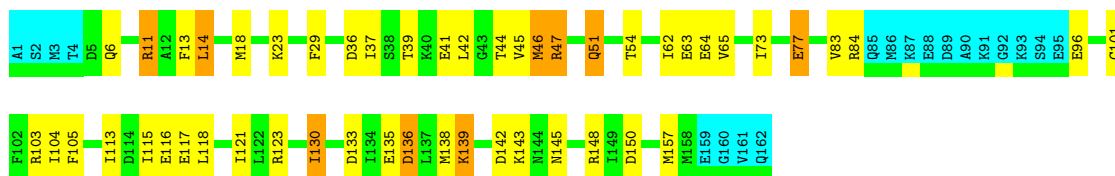
- Molecule 1: TROPONIN C





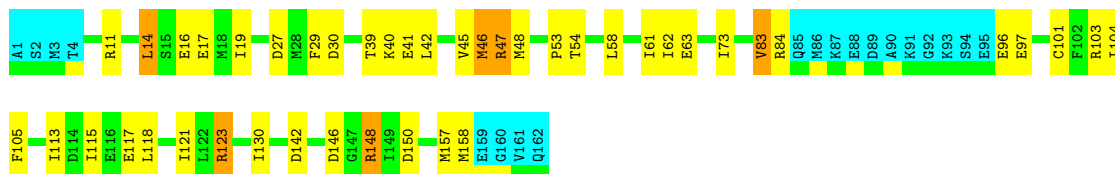
4.2.3 Score per residue for model 3

- Molecule 1: TROPONIN C



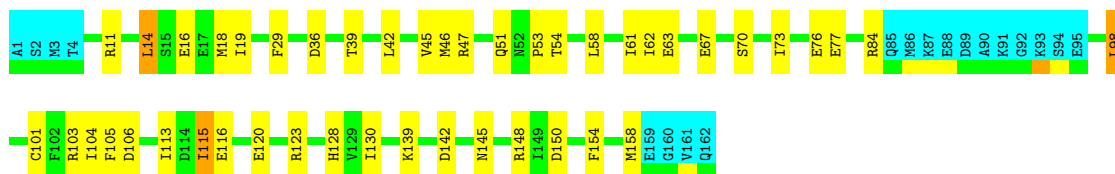
4.2.4 Score per residue for model 4

- Molecule 1: TROPONIN C



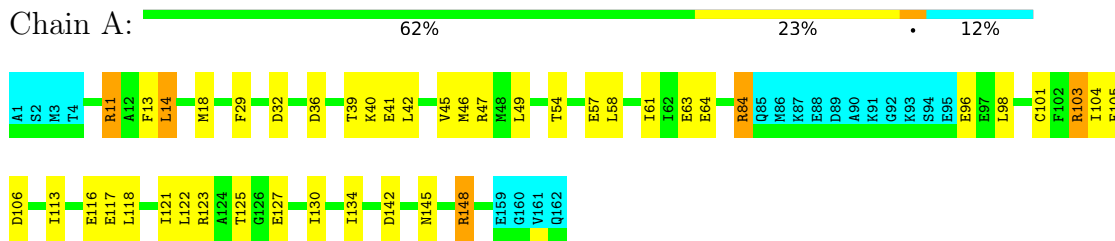
4.2.5 Score per residue for model 5

- Molecule 1: TROPONIN C



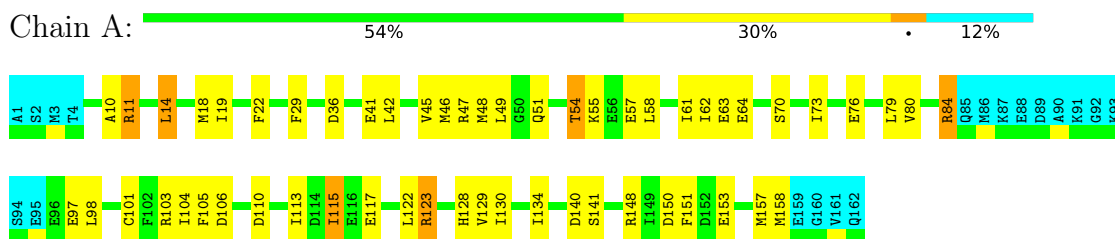
4.2.6 Score per residue for model 6

- Molecule 1: TROPONIN C



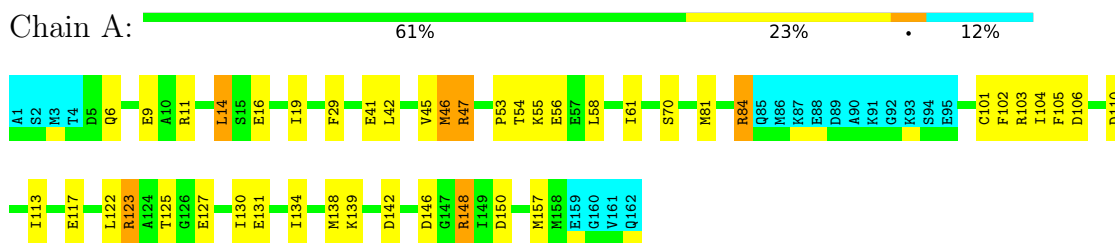
4.2.7 Score per residue for model 7

- Molecule 1: TROPONIN C



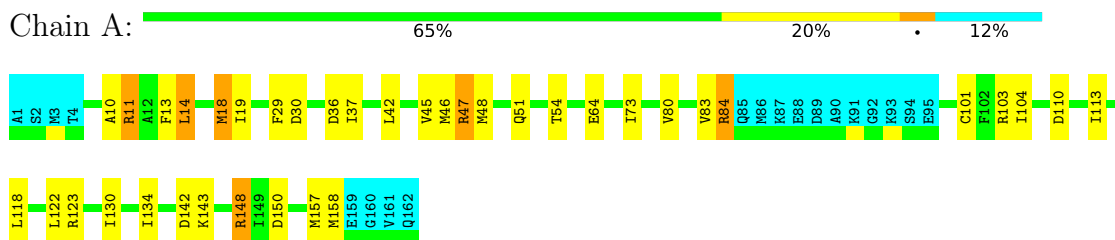
4.2.8 Score per residue for model 8

- Molecule 1: TROPONIN C



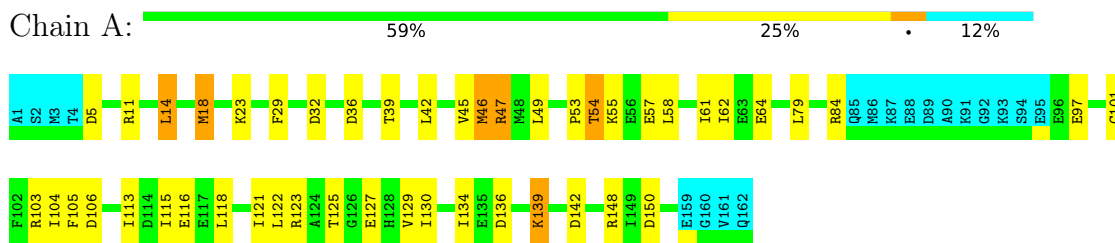
4.2.9 Score per residue for model 9

- Molecule 1: TROPONIN C



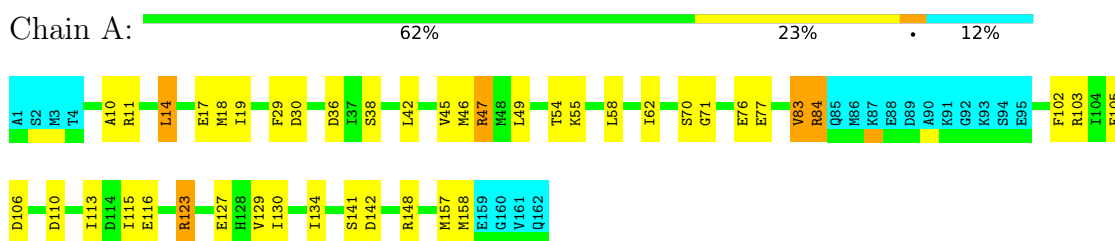
4.2.10 Score per residue for model 10

- Molecule 1: TROPONIN C



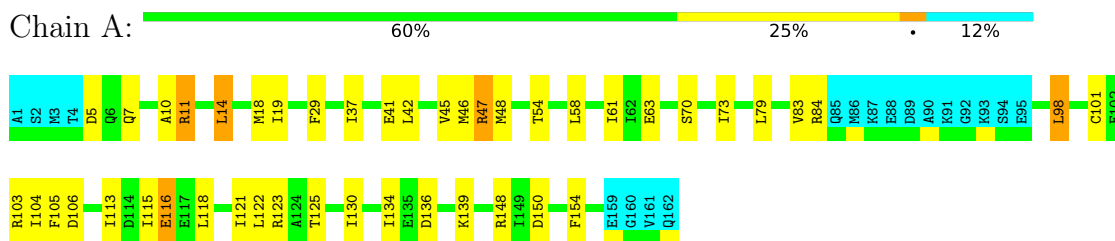
4.2.11 Score per residue for model 11

- Molecule 1: TROPONIN C



4.2.12 Score per residue for model 12

- Molecule 1: TROPONIN C



4.2.13 Score per residue for model 13

- Molecule 1: TROPONIN C



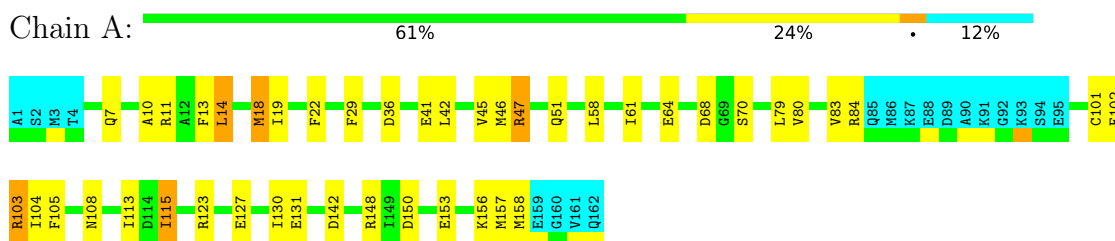
4.2.14 Score per residue for model 14

- Molecule 1: TROPONIN C



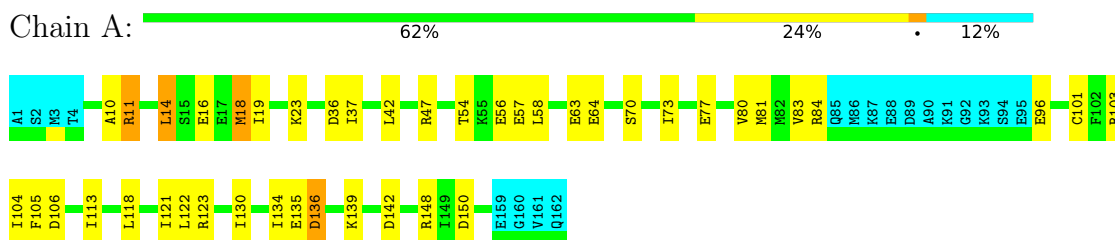
4.2.15 Score per residue for model 15

- Molecule 1: TROPONIN C



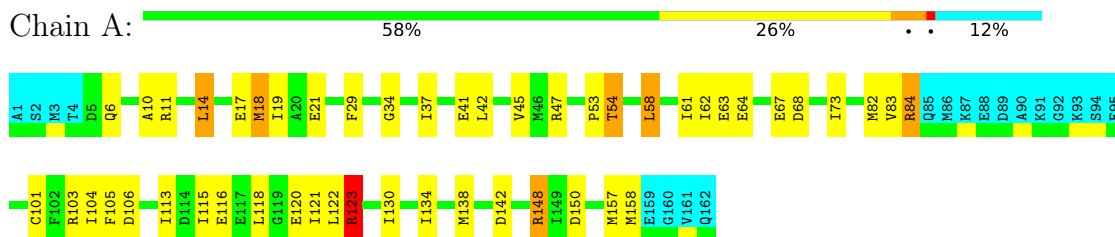
4.2.16 Score per residue for model 16

- Molecule 1: TROPONIN C



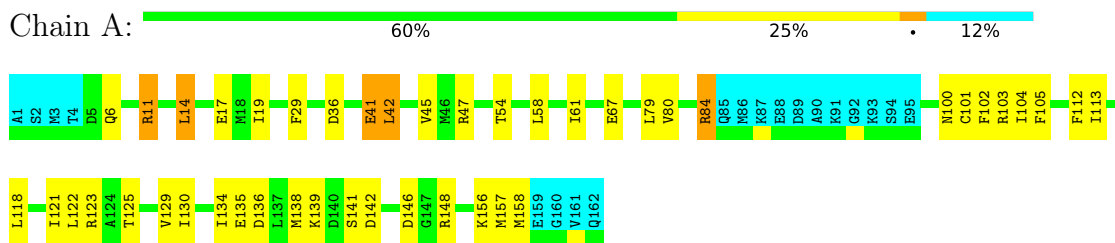
4.2.17 Score per residue for model 17

- Molecule 1: TROPONIN C



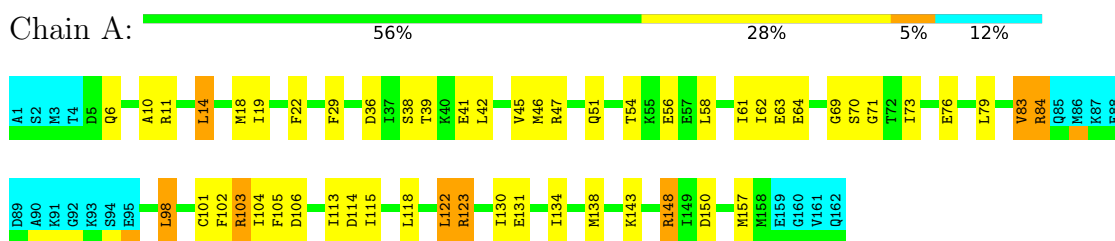
4.2.18 Score per residue for model 18

- Molecule 1: TROPONIN C



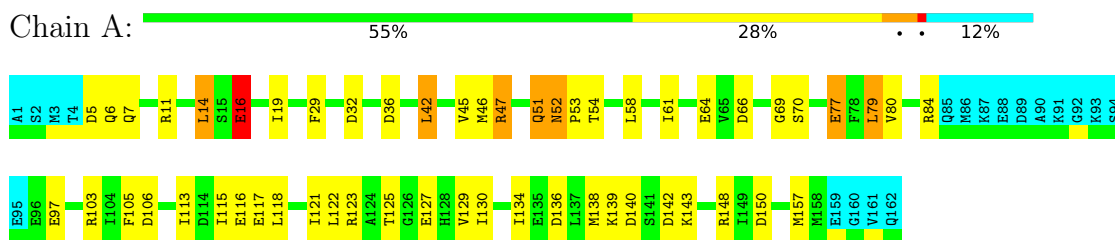
4.2.19 Score per residue for model 19

- Molecule 1: TROPONIN C



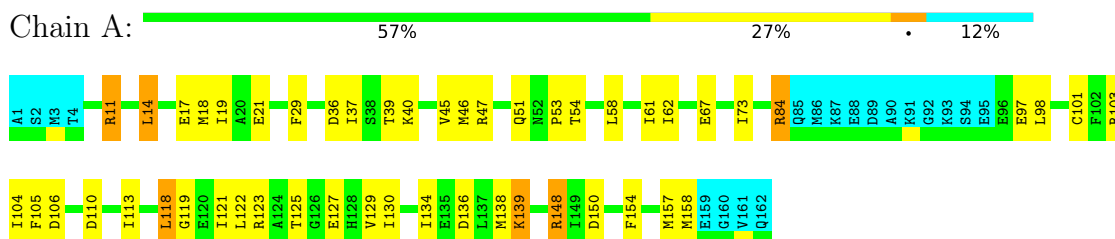
4.2.20 Score per residue for model 20

- Molecule 1: TROPONIN C



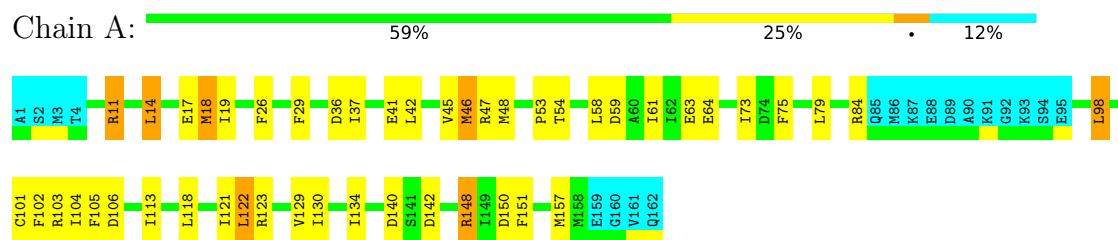
4.2.21 Score per residue for model 21

- Molecule 1: TROPONIN C



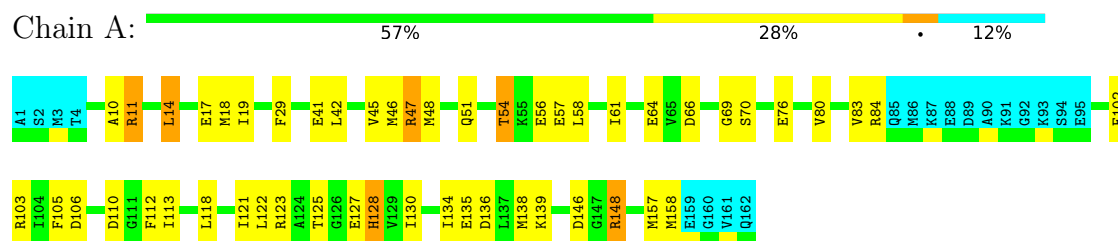
4.2.22 Score per residue for model 22 (medoid)

- Molecule 1: TROPONIN C



4.2.23 Score per residue for model 23

- Molecule 1: TROPONIN C



5 Refinement protocol and experimental data overview

Of the ? calculated structures, 23 were deposited, based on the following criterion: ?.

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

Software name	Classification	Version
X-PLOR	refinement	3.1

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

There are no covalent bond-length or bond-angle outliers.

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	6.0±0.0
All	All	0	138

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	A	11	ARG	Sidechain	23
1	A	47	ARG	Sidechain	23
1	A	84	ARG	Sidechain	23
1	A	103	ARG	Sidechain	23
1	A	123	ARG	Sidechain	23
1	A	148	ARG	Sidechain	23

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	1131	1060	1060	20±4
All	All	26013	24380	24380	453

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 9.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:42:LEU:HD23	1:A:58:LEU:HD22	0.99	1.34	14	9
1:A:122:LEU:HD13	1:A:134:ILE:HD12	0.94	1.37	2	3
1:A:14:LEU:HD13	1:A:19:ILE:HD11	0.86	1.46	8	1
1:A:42:LEU:HD11	1:A:73:ILE:HD11	0.84	1.48	19	4
1:A:42:LEU:HB3	1:A:58:LEU:HD13	0.81	1.53	10	2
1:A:129:VAL:HG13	1:A:134:ILE:HD11	0.78	1.55	20	4
1:A:101:CYS:HA	1:A:104:ILE:HD12	0.72	1.61	14	18
1:A:122:LEU:HD11	1:A:129:VAL:HG11	0.69	1.62	1	1
1:A:122:LEU:O	1:A:125:THR:HG22	0.69	1.86	13	11
1:A:16:GLU:HA	1:A:19:ILE:HD12	0.69	1.62	20	2
1:A:14:LEU:HD12	1:A:19:ILE:CG1	0.68	2.18	12	19
1:A:118:LEU:HA	1:A:121:ILE:HD12	0.66	1.68	20	12
1:A:39:THR:HA	1:A:42:LEU:HD12	0.65	1.67	19	1
1:A:14:LEU:HD13	1:A:19:ILE:CD1	0.65	2.18	8	1
1:A:14:LEU:HD12	1:A:19:ILE:HG13	0.64	1.68	15	18
1:A:42:LEU:HD23	1:A:58:LEU:CD2	0.63	2.22	11	7
1:A:129:VAL:CG1	1:A:134:ILE:HD11	0.62	2.24	20	6
1:A:29:PHE:HA	1:A:45:VAL:HG21	0.62	1.72	22	21
1:A:98:LEU:HD12	1:A:154:PHE:CE2	0.61	2.30	2	3
1:A:54:THR:HG23	1:A:57:GLU:HB2	0.61	1.71	10	3
1:A:58:LEU:HA	1:A:61:ILE:HD12	0.60	1.72	19	17
1:A:14:LEU:CD1	1:A:79:LEU:HD22	0.60	2.26	15	6
1:A:129:VAL:HG21	1:A:134:ILE:HD11	0.60	1.72	7	2
1:A:29:PHE:CD2	1:A:45:VAL:HG21	0.60	2.32	3	1
1:A:42:LEU:HD11	1:A:73:ILE:CD1	0.59	2.25	19	1
1:A:14:LEU:HD22	1:A:14:LEU:N	0.59	2.12	6	1
1:A:122:LEU:HD13	1:A:134:ILE:HG12	0.58	1.75	20	9
1:A:105:PHE:HB3	1:A:113:ILE:HG12	0.58	1.75	20	13
1:A:29:PHE:HB3	1:A:37:ILE:HD12	0.58	1.75	3	1
1:A:54:THR:HG23	1:A:57:GLU:CB	0.58	2.29	7	1
1:A:122:LEU:HD13	1:A:134:ILE:CD1	0.57	2.30	9	3
1:A:13:PHE:C	1:A:14:LEU:HD13	0.57	2.20	6	1
1:A:39:THR:HG23	1:A:62:ILE:HD12	0.57	1.76	21	1
1:A:122:LEU:HB3	1:A:129:VAL:HG11	0.57	1.76	22	1
1:A:42:LEU:HD12	1:A:62:ILE:HD11	0.57	1.75	3	2
1:A:14:LEU:HD22	1:A:14:LEU:H	0.57	1.58	6	1
1:A:14:LEU:HD13	1:A:79:LEU:HD12	0.57	1.76	20	1
1:A:41:GLU:O	1:A:45:VAL:HG23	0.56	2.01	2	11
1:A:77:GLU:O	1:A:80:VAL:HG12	0.56	2.01	1	3
1:A:122:LEU:HG	1:A:123:ARG:N	0.55	2.15	19	2
1:A:62:ILE:HG23	1:A:73:ILE:HG12	0.55	1.77	7	4

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:105:PHE:CB	1:A:113:ILE:HG12	0.54	2.32	21	17
1:A:14:LEU:HD12	1:A:19:ILE:HG12	0.54	1.79	14	11
1:A:53:PRO:HB2	1:A:58:LEU:HD22	0.54	1.78	17	1
1:A:14:LEU:HD11	1:A:79:LEU:HD22	0.53	1.78	10	4
1:A:62:ILE:HG21	1:A:71:GLY:O	0.53	2.04	19	2
1:A:102:PHE:CE1	1:A:113:ILE:HD13	0.53	2.39	23	3
1:A:46:MET:CB	1:A:53:PRO:HG3	0.53	2.34	20	5
1:A:125:THR:HG23	1:A:127:GLU:H	0.52	1.64	1	8
1:A:29:PHE:CD1	1:A:42:LEU:HD23	0.52	2.39	1	1
1:A:39:THR:CG2	1:A:62:ILE:HD12	0.52	2.35	21	4
1:A:6:GLN:HB3	1:A:80:VAL:HG21	0.52	1.81	18	1
1:A:98:LEU:HB3	1:A:154:PHE:CE2	0.51	2.40	21	1
1:A:122:LEU:HD21	1:A:134:ILE:HD13	0.51	1.82	19	2
1:A:39:THR:HG22	1:A:62:ILE:HD12	0.51	1.81	10	3
1:A:14:LEU:HB2	1:A:18:MET:HB3	0.51	1.81	10	8
1:A:113:ILE:HD12	1:A:113:ILE:N	0.51	2.21	1	19
1:A:130:ILE:CD1	1:A:133:ASP:HB2	0.51	2.36	3	1
1:A:14:LEU:N	1:A:14:LEU:HD13	0.51	2.19	6	1
1:A:42:LEU:HD22	1:A:62:ILE:HD11	0.51	1.82	11	1
1:A:39:THR:HA	1:A:62:ILE:HD11	0.51	1.81	4	2
1:A:14:LEU:HD22	1:A:79:LEU:HD12	0.50	1.83	20	1
1:A:118:LEU:HD21	1:A:122:LEU:HG	0.50	1.83	21	1
1:A:51:GLN:O	1:A:52:ASN:C	0.50	2.49	20	1
1:A:54:THR:O	1:A:58:LEU:HD23	0.50	2.05	17	1
1:A:115:ILE:O	1:A:115:ILE:HD13	0.50	2.06	13	3
1:A:10:ALA:CB	1:A:83:VAL:HG21	0.49	2.37	17	1
1:A:102:PHE:CE1	1:A:113:ILE:CD1	0.49	2.95	15	9
1:A:10:ALA:CB	1:A:80:VAL:HG23	0.49	2.37	7	3
1:A:10:ALA:HB1	1:A:83:VAL:HG21	0.49	1.85	17	2
1:A:58:LEU:HD23	1:A:61:ILE:HD12	0.49	1.84	8	2
1:A:13:PHE:CD2	1:A:83:VAL:HG22	0.48	2.42	15	3
1:A:54:THR:O	1:A:58:LEU:HB3	0.48	2.07	17	1
1:A:42:LEU:HD13	1:A:73:ILE:HD11	0.48	1.86	17	1
1:A:10:ALA:CB	1:A:83:VAL:HG11	0.48	2.38	23	1
1:A:42:LEU:CD1	1:A:62:ILE:HD11	0.48	2.39	19	3
1:A:14:LEU:CD1	1:A:79:LEU:HD12	0.48	2.39	20	1
1:A:37:ILE:HB	1:A:73:ILE:HB	0.48	1.86	22	9
1:A:135:GLU:CG	1:A:136:ASP:N	0.48	2.76	23	1
1:A:10:ALA:HB2	1:A:80:VAL:HG23	0.48	1.86	9	2
1:A:136:ASP:HA	1:A:139:LYS:HG2	0.47	1.84	16	8
1:A:136:ASP:HA	1:A:139:LYS:CG	0.47	2.39	3	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:14:LEU:CD2	1:A:83:VAL:HG23	0.47	2.39	16	3
1:A:14:LEU:CD1	1:A:19:ILE:HG12	0.47	2.39	20	1
1:A:98:LEU:HD23	1:A:154:PHE:CD2	0.47	2.44	1	1
1:A:49:LEU:CD1	1:A:49:LEU:N	0.47	2.77	10	2
1:A:10:ALA:HA	1:A:83:VAL:HG21	0.47	1.86	11	2
1:A:98:LEU:HD22	1:A:155:LEU:HD23	0.46	1.86	14	1
1:A:98:LEU:HD22	1:A:98:LEU:N	0.46	2.24	21	2
1:A:115:ILE:HD13	1:A:115:ILE:O	0.46	2.09	14	2
1:A:46:MET:HB2	1:A:53:PRO:HG3	0.46	1.87	20	1
1:A:65:VAL:HG12	1:A:77:GLU:HG2	0.46	1.88	3	1
1:A:76:GLU:HA	1:A:79:LEU:HD12	0.46	1.87	14	1
1:A:29:PHE:CE1	1:A:42:LEU:HD13	0.46	2.45	20	1
1:A:122:LEU:CD1	1:A:129:VAL:HG11	0.46	2.38	1	1
1:A:118:LEU:HD12	1:A:121:ILE:HD11	0.45	1.88	13	1
1:A:98:LEU:HD21	1:A:158:MET:HB3	0.45	1.86	1	1
1:A:14:LEU:HD13	1:A:22:PHE:CD2	0.45	2.47	19	5
1:A:130:ILE:CD1	1:A:133:ASP:CB	0.45	2.95	3	1
1:A:118:LEU:HD21	1:A:122:LEU:CG	0.45	2.42	21	1
1:A:118:LEU:HD21	1:A:122:LEU:CD1	0.45	2.42	21	1
1:A:62:ILE:HG23	1:A:73:ILE:CG1	0.45	2.40	7	1
1:A:14:LEU:HD12	1:A:14:LEU:O	0.44	2.11	3	1
1:A:105:PHE:HB2	1:A:113:ILE:HG12	0.44	1.89	4	2
1:A:14:LEU:HD13	1:A:79:LEU:HD22	0.44	1.90	18	2
1:A:105:PHE:CB	1:A:113:ILE:CG1	0.44	2.95	1	1
1:A:14:LEU:HD13	1:A:22:PHE:CE2	0.44	2.48	19	1
1:A:37:ILE:HG22	1:A:42:LEU:HG	0.44	1.87	22	1
1:A:14:LEU:HB3	1:A:18:MET:HB3	0.44	1.89	6	1
1:A:46:MET:CE	1:A:58:LEU:HD21	0.44	2.42	21	1
1:A:115:ILE:HG23	1:A:116:GLU:N	0.43	2.28	12	7
1:A:98:LEU:HD21	1:A:151:PHE:HZ	0.43	1.72	22	1
1:A:14:LEU:CD2	1:A:79:LEU:HD12	0.43	2.44	20	1
1:A:122:LEU:CD1	1:A:134:ILE:HG23	0.43	2.44	17	1
1:A:46:MET:SD	1:A:53:PRO:HG3	0.43	2.53	10	2
1:A:10:ALA:HB1	1:A:80:VAL:HA	0.43	1.90	15	1
1:A:42:LEU:HD22	1:A:62:ILE:CG1	0.42	2.43	17	1
1:A:112:PHE:CD1	1:A:112:PHE:N	0.42	2.87	23	1
1:A:98:LEU:HD12	1:A:154:PHE:CD2	0.42	2.49	2	1
1:A:79:LEU:N	1:A:79:LEU:CD2	0.42	2.83	20	1
1:A:45:VAL:HG13	1:A:48:MET:SD	0.42	2.55	12	1
1:A:14:LEU:CD1	1:A:19:ILE:HD11	0.42	2.33	8	1
1:A:42:LEU:HD13	1:A:73:ILE:CD1	0.42	2.45	17	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:105:PHE:CD2	1:A:121:ILE:HD13	0.41	2.49	18	1
1:A:37:ILE:HG21	1:A:42:LEU:CD1	0.41	2.46	14	1
1:A:122:LEU:HD13	1:A:134:ILE:HG23	0.41	1.91	2	1
1:A:37:ILE:O	1:A:72:THR:HB	0.41	2.16	2	1
1:A:118:LEU:HA	1:A:121:ILE:CG1	0.41	2.45	13	1
1:A:120:GLU:HA	1:A:123:ARG:CG	0.41	2.46	17	1
1:A:26:PHE:CE2	1:A:75:PHE:HA	0.41	2.51	22	1
1:A:114:ASP:HA	1:A:148:ARG:HG3	0.41	1.92	19	1
1:A:18:MET:SD	1:A:22:PHE:CZ	0.40	3.14	14	1
1:A:36:ASP:HB3	1:A:73:ILE:O	0.40	2.16	3	2
1:A:39:THR:HG22	1:A:59:ASP:HA	0.40	1.93	14	1
1:A:6:GLN:HB3	1:A:80:VAL:CG2	0.40	2.45	18	1
1:A:46:MET:HB3	1:A:51:GLN:HB3	0.40	1.92	3	2

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	143/162 (88%)	134±2 (94±2%)	8±2 (6±1%)	1±1 (1±1%)	29	74
All	All	3289/3726 (88%)	3084 (94%)	184 (6%)	21 (1%)	29	74

All 9 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	69	GLY	5
1	A	34	GLY	3
1	A	128	HIS	3
1	A	118	LEU	3
1	A	145	ASN	2
1	A	119	GLY	2
1	A	5	ASP	1
1	A	16	GLU	1
1	A	52	ASN	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	121/136 (89%)	99±3 (82±3%)	22±3 (18±3%)	4 37
All	All	2783/3128 (89%)	2270 (82%)	513 (18%)	4 37

All 81 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	14	LEU	23
1	A	130	ILE	23
1	A	54	THR	22
1	A	142	ASP	18
1	A	150	ASP	18
1	A	106	ASP	17
1	A	18	MET	16
1	A	157	MET	16
1	A	36	ASP	15
1	A	46	MET	15
1	A	11	ARG	13
1	A	51	GLN	13
1	A	63	GLU	13
1	A	70	SER	13
1	A	64	GLU	13
1	A	148	ARG	11
1	A	47	ARG	11
1	A	84	ARG	11
1	A	138	MET	10
1	A	158	MET	10
1	A	76	GLU	7
1	A	117	GLU	7
1	A	17	GLU	7
1	A	115	ILE	7
1	A	42	LEU	7
1	A	41	GLU	6
1	A	98	LEU	6
1	A	6	GLN	6
1	A	139	LYS	6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Models (Total)
1	A	48	MET	6
1	A	110	ASP	6
1	A	16	GLU	5
1	A	136	ASP	5
1	A	143	LYS	5
1	A	146	ASP	5
1	A	77	GLU	5
1	A	97	GLU	5
1	A	123	ARG	5
1	A	103	ARG	5
1	A	56	GLU	5
1	A	5	ASP	4
1	A	81	MET	4
1	A	122	LEU	4
1	A	96	GLU	4
1	A	30	ASP	4
1	A	40	LYS	4
1	A	67	GLU	4
1	A	116	GLU	4
1	A	55	LYS	4
1	A	66	ASP	4
1	A	23	LYS	3
1	A	135	GLU	3
1	A	83	VAL	3
1	A	32	ASP	3
1	A	140	ASP	3
1	A	141	SER	3
1	A	131	GLU	3
1	A	7	GLN	3
1	A	27	ASP	2
1	A	57	GLU	2
1	A	49	LEU	2
1	A	153	GLU	2
1	A	38	SER	2
1	A	127	GLU	2
1	A	68	ASP	2
1	A	156	LYS	2
1	A	21	GLU	2
1	A	44	THR	1
1	A	120	GLU	1
1	A	39	THR	1
1	A	145	ASN	1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Models (Total)
1	A	9	GLU	1
1	A	118	LEU	1
1	A	108	ASN	1
1	A	58	LEU	1
1	A	82	MET	1
1	A	100	ASN	1
1	A	112	PHE	1
1	A	79	LEU	1
1	A	59	ASP	1
1	A	128	HIS	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 monosaccharides 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