



Full wwPDB NMR Structure Validation Report ⓘ

Mar 7, 2026 – 01:43 AM UTC

PDB ID : 1BLD / pdb_00001bld
Title : BASIC FIBROBLAST GROWTH FACTOR (FGF-2) MUTANT WITH CYS
78 REPLACED BY SER AND CYS 96 REPLACED BY SER, NMR
Authors : Powers, R.; Seddon, A.P.; Bohlen, P.; Moy, F.J.
Deposited on : 1996-05-20

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

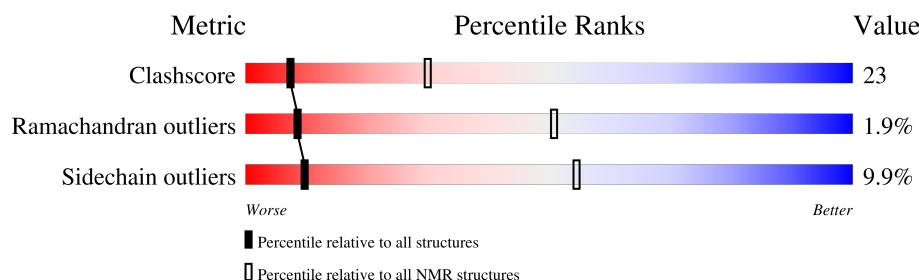
1 Overall quality at a glance

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	155	50% 32% 19%

2 Ensemble composition and analysis

This entry contains 30 models. Model 24 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:28-A:153 (126)	0.51	24

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 4 single-model clusters were found.

Cluster number	Models
1	1, 2, 4, 5, 6, 8, 12, 13, 18, 20, 21, 22, 24, 25, 26, 27, 28, 30
2	7, 14, 23, 29
3	10, 17
4	3, 9
Single-model clusters	11; 15; 16; 19

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2443 atoms, of which 1222 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called BASIC FIBROBLAST GROWTH FACTOR.

Mol	Chain	Residues	Atoms						Trace
1	A	155	Total	C	H	N	O	S	0
			2443	773	1222	218	225	5	

There are 4 discrepancies between the modelled and reference sequences:

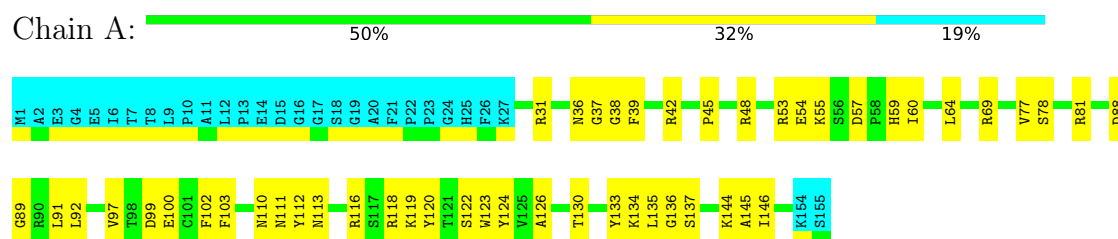
Chain	Residue	Modelled	Actual	Comment	Reference
A	3	GLU	ALA	conflict	UNP P09038
A	5	GLU	SER	conflict	UNP P09038
A	78	SER	CYS	engineered mutation	UNP P09038
A	96	SER	CYS	engineered mutation	UNP P09038

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: BASIC FIBROBLAST GROWTH FACTOR

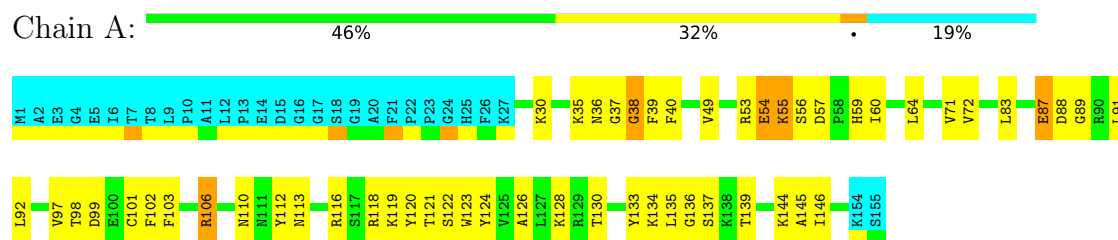


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

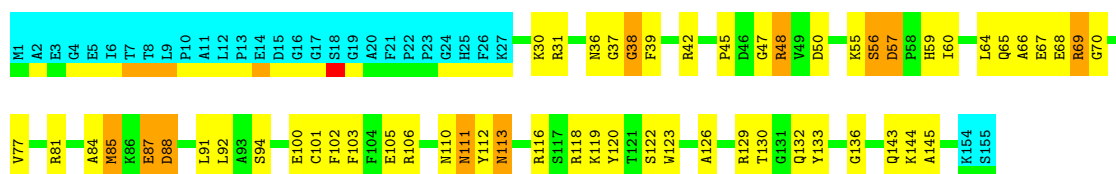
• Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



4.2.2 Score per residue for model 2

• Molecule 1: BASIC FIBROBLAST GROWTH FACTOR

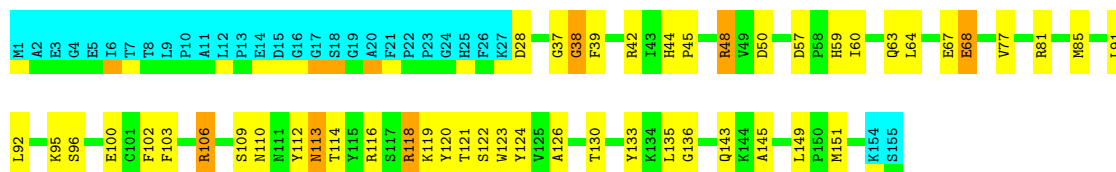




4.2.3 Score per residue for model 3

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR

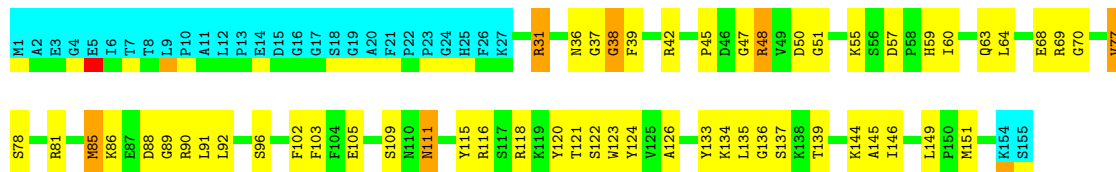
Chain A: 50% 28% 19%



4.2.4 Score per residue for model 4

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR

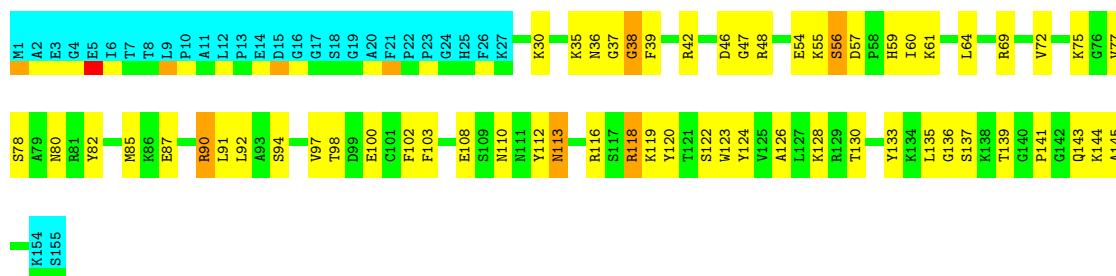
Chain A: 45% 32% 19%



4.2.5 Score per residue for model 5

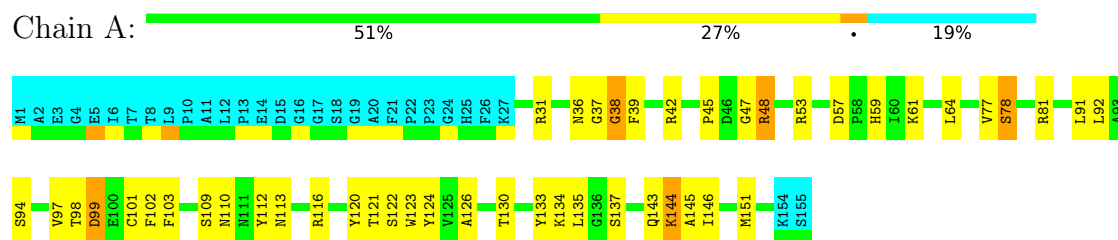
- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR

Chain A: 43% 35% 19%



4.2.6 Score per residue for model 6

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



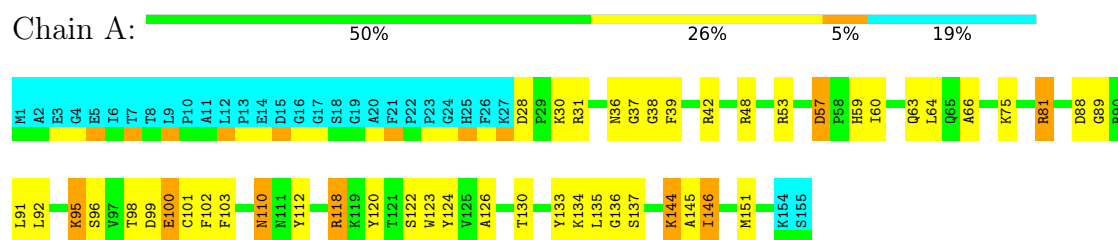
4.2.7 Score per residue for model 7

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



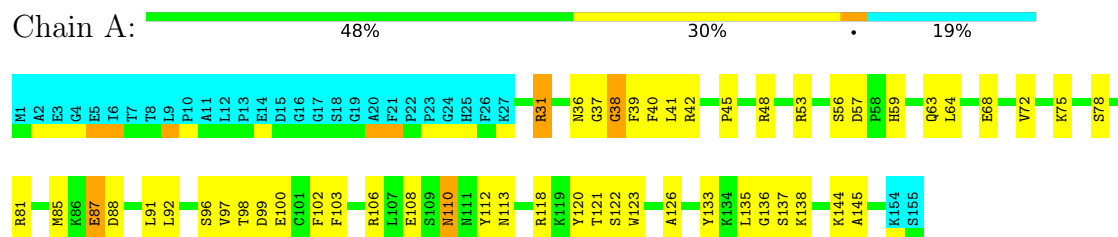
4.2.8 Score per residue for model 8

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



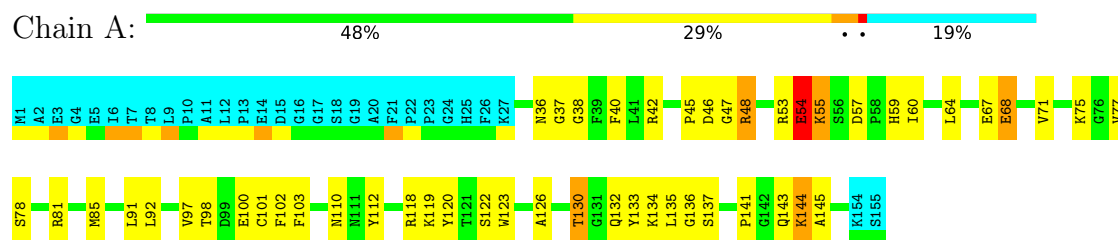
4.2.9 Score per residue for model 9

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



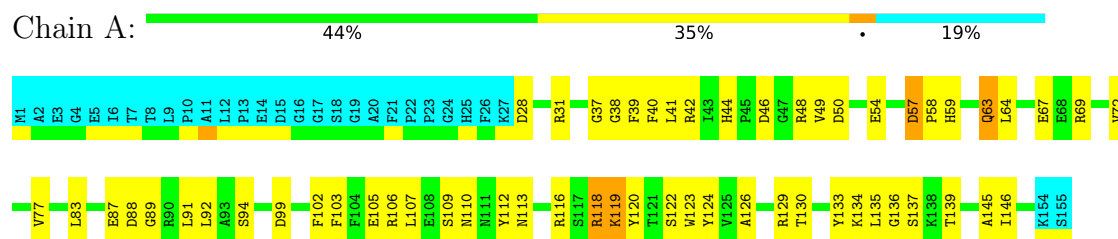
4.2.10 Score per residue for model 10

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



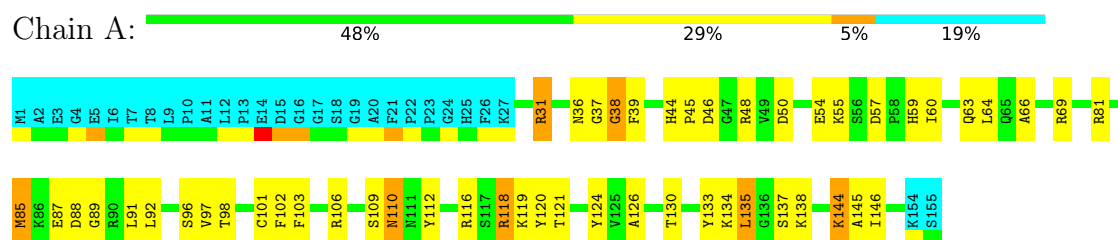
4.2.11 Score per residue for model 11

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



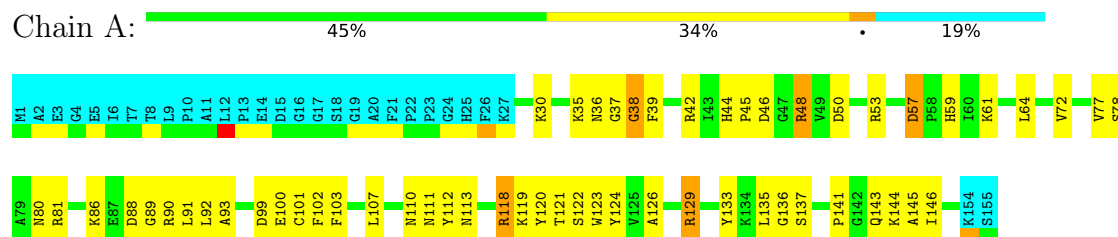
4.2.12 Score per residue for model 12

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



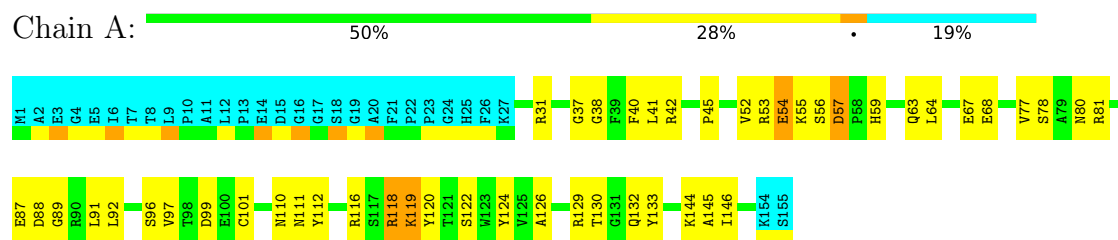
4.2.13 Score per residue for model 13

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



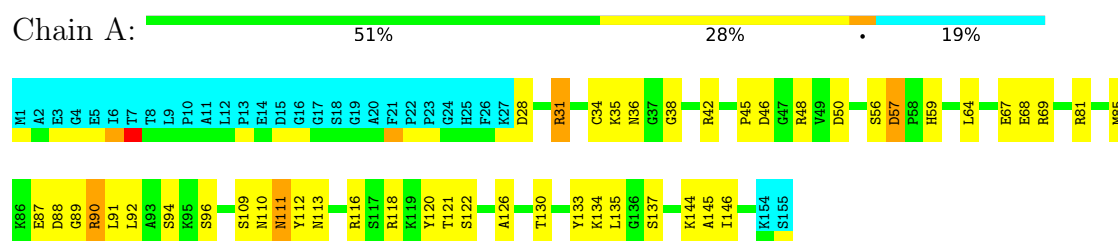
4.2.14 Score per residue for model 14

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



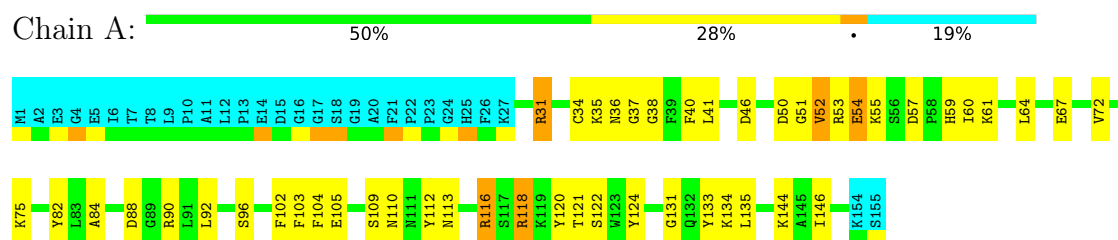
4.2.15 Score per residue for model 15

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



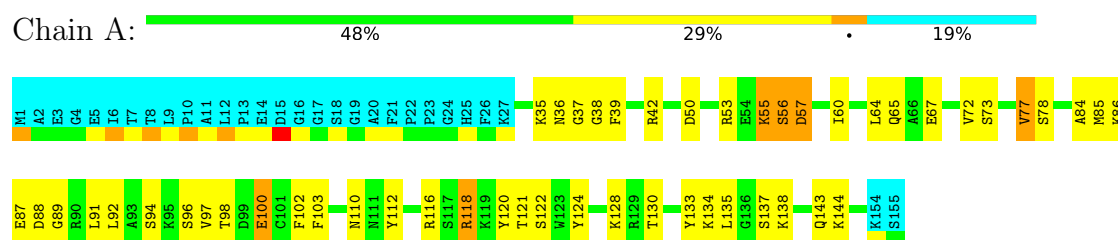
4.2.16 Score per residue for model 16

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



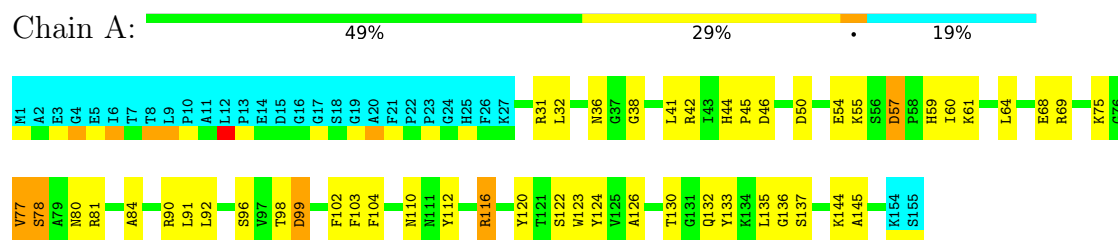
4.2.17 Score per residue for model 17

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



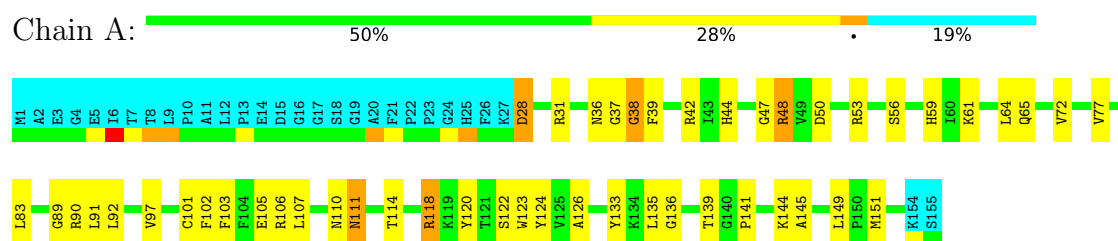
4.2.18 Score per residue for model 18

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



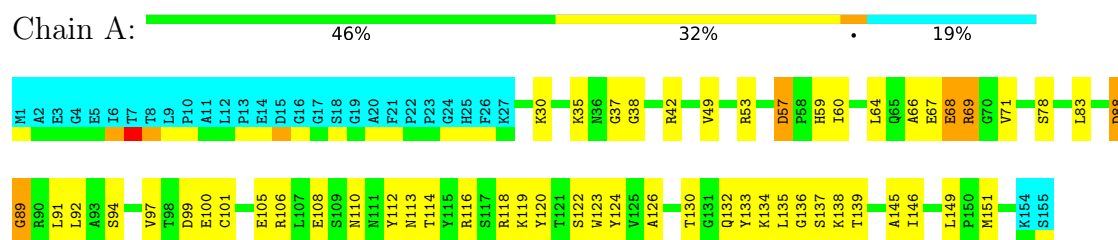
4.2.19 Score per residue for model 19

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



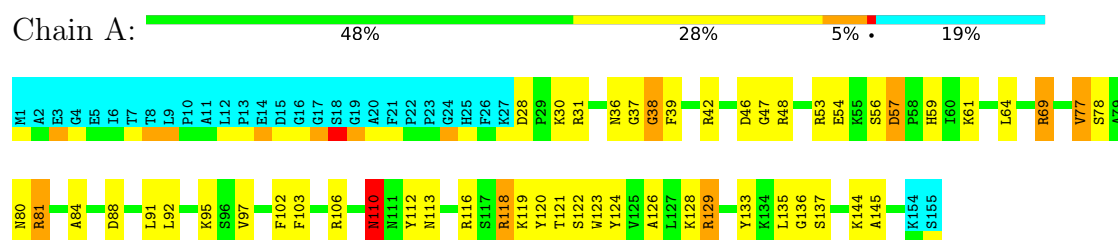
4.2.20 Score per residue for model 20

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



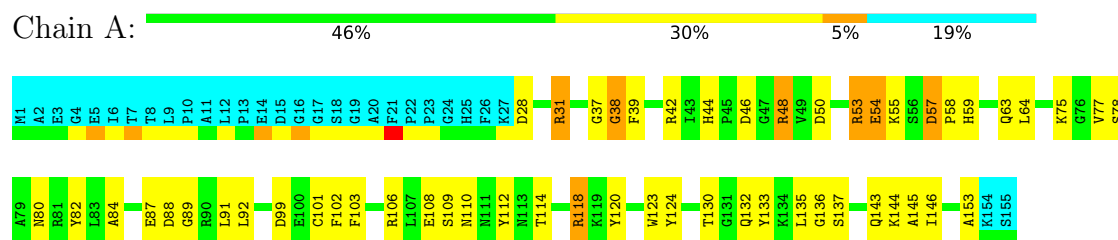
4.2.21 Score per residue for model 21

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



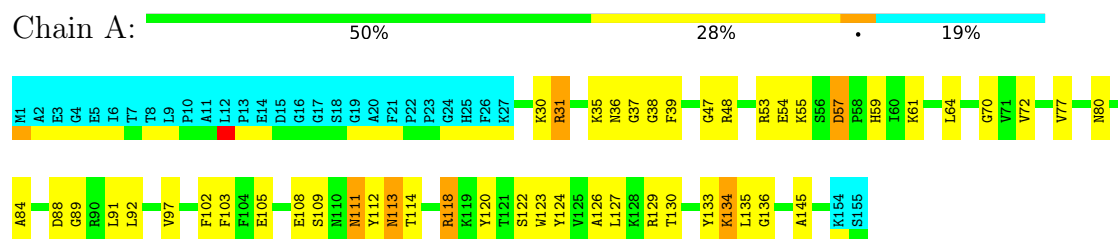
4.2.22 Score per residue for model 22

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



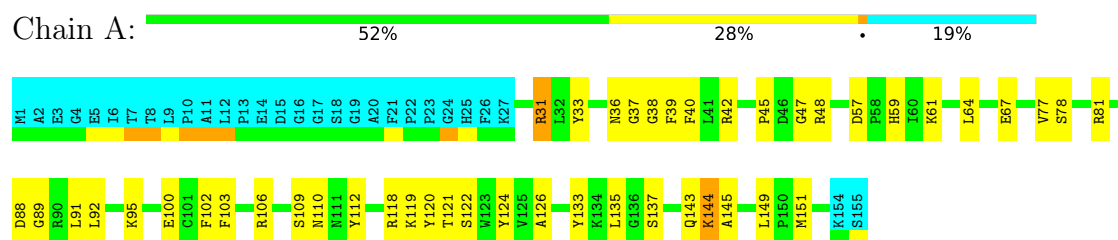
4.2.23 Score per residue for model 23

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



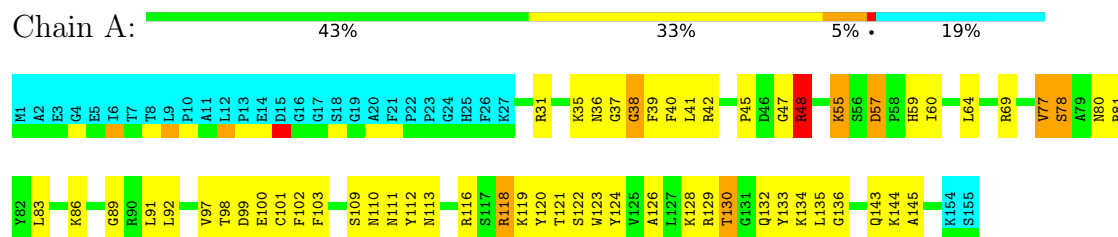
4.2.24 Score per residue for model 24 (medoid)

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



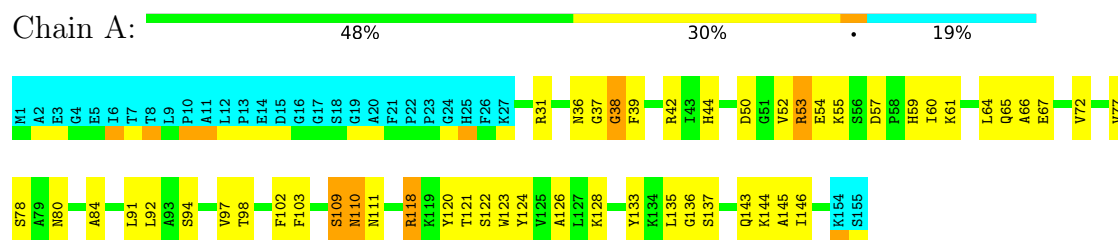
4.2.25 Score per residue for model 25

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



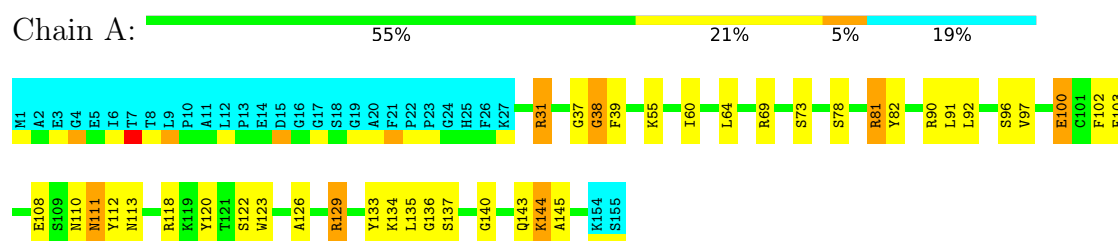
4.2.26 Score per residue for model 26

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



4.2.27 Score per residue for model 27

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



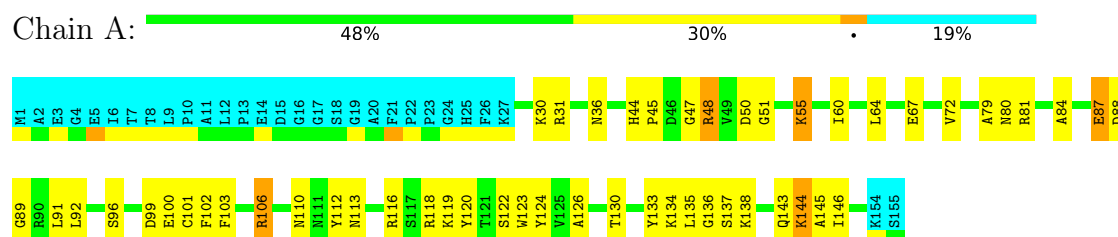
4.2.28 Score per residue for model 28

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



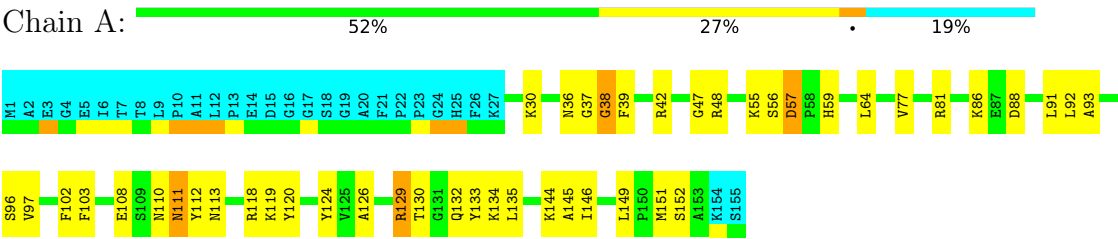
4.2.29 Score per residue for model 29

- Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



4.2.30 Score per residue for model 30

● Molecule 1: BASIC FIBROBLAST GROWTH FACTOR



5 Refinement protocol and experimental data overview ⓘ

Of the ? calculated structures, 30 were deposited, based on the following criterion: ?.

The authors did not provide any information on software used for structure solution, optimization or refinement.

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	1.41±0.01	0±0/1036 (0.0± 0.0%)	1.25±0.01	0±0/1390 (0.0± 0.0%)
All	All	1.41	0/31080 (0.0%)	1.25	2/41700 (0.0%)

There are no bond-length outliers.

All unique angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	68	GLU	CB-CA-C	-6.12	109.51	116.54	18	2

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	1014	1023	1023	48±6
All	All	30420	30690	30690	1426

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 23.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:44:HIS:HE2	1:A:50:ASP:CG	0.70	1.94	18	2

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:91:LEU:HD12	1:A:91:LEU:N	0.70	2.02	25	25
1:A:92:LEU:C	1:A:92:LEU:HD12	0.68	2.15	11	30
1:A:130:THR:HG23	1:A:132:GLN:OE1	0.66	1.90	25	2
1:A:126:ALA:HB1	1:A:145:ALA:HB3	0.65	1.68	15	6
1:A:97:VAL:O	1:A:97:VAL:HG13	0.65	1.90	14	10
1:A:126:ALA:HB1	1:A:145:ALA:CB	0.62	2.25	7	21
1:A:75:LYS:NZ	1:A:82:TYR:CE1	0.62	2.68	16	1
1:A:42:ARG:CZ	1:A:52:VAL:HG21	0.61	2.25	26	1
1:A:31:ARG:NH1	1:A:61:LYS:NZ	0.61	2.48	16	1
1:A:134:LYS:NZ	1:A:143:GLN:NE2	0.61	2.48	10	3
1:A:39:PHE:CZ	1:A:53:ARG:NH1	0.61	2.68	26	2
1:A:135:LEU:HD23	1:A:137:SER:OG	0.60	1.97	13	18
1:A:149:LEU:HD23	1:A:151:MET:HE2	0.60	1.72	24	6
1:A:55:LYS:H	1:A:55:LYS:CD	0.60	2.09	25	1
1:A:113:ASN:HD22	1:A:113:ASN:N	0.60	1.94	7	4
1:A:36:ASN:ND2	1:A:145:ALA:N	0.59	2.50	28	4
1:A:134:LYS:NZ	1:A:143:GLN:HE21	0.59	1.94	17	2
1:A:64:LEU:HD12	1:A:64:LEU:N	0.58	2.13	14	29
1:A:132:GLN:N	1:A:132:GLN:NE2	0.58	2.51	20	1
1:A:85:MET:SD	1:A:91:LEU:HD11	0.58	2.38	10	8
1:A:84:ALA:HB3	1:A:92:LEU:CD1	0.58	2.29	16	8
1:A:39:PHE:CE1	1:A:53:ARG:NH1	0.58	2.72	26	2
1:A:126:ALA:HB1	1:A:145:ALA:HB1	0.58	1.76	26	18
1:A:118:ARG:O	1:A:121:THR:HG23	0.57	1.98	13	5
1:A:108:GLU:OE2	1:A:114:THR:HG23	0.57	1.99	23	1
1:A:116:ARG:CZ	1:A:124:TYR:CZ	0.56	2.89	3	1
1:A:98:THR:N	1:A:101:CYS:SG	0.56	2.79	10	2
1:A:91:LEU:N	1:A:91:LEU:CD1	0.56	2.69	25	25
1:A:113:ASN:N	1:A:113:ASN:ND2	0.56	2.53	7	6
1:A:87:GLU:H	1:A:87:GLU:CD	0.56	2.09	9	4
1:A:106:ARG:CZ	1:A:116:ARG:NH2	0.55	2.70	12	1
1:A:75:LYS:NZ	1:A:82:TYR:CZ	0.55	2.73	5	2
1:A:111:ASN:N	1:A:111:ASN:ND2	0.55	2.54	23	2
1:A:106:ARG:CG	1:A:106:ARG:HH11	0.54	2.16	3	3
1:A:90:ARG:CG	1:A:90:ARG:HH11	0.54	2.16	15	3
1:A:48:ARG:CG	1:A:48:ARG:HH11	0.54	2.16	23	10
1:A:118:ARG:HH11	1:A:118:ARG:CG	0.54	2.16	21	8
1:A:53:ARG:CG	1:A:53:ARG:HH11	0.54	2.16	26	2
1:A:69:ARG:N	1:A:69:ARG:CD	0.54	2.70	18	2
1:A:48:ARG:NH1	1:A:48:ARG:CG	0.54	2.71	9	3
1:A:31:ARG:NH1	1:A:31:ARG:CG	0.54	2.71	22	5

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:106:ARG:HE	1:A:116:ARG:HH22	0.54	1.44	29	1
1:A:69:ARG:HH11	1:A:69:ARG:CG	0.54	2.16	2	3
1:A:31:ARG:CG	1:A:31:ARG:HH11	0.54	2.16	12	9
1:A:31:ARG:CG	1:A:31:ARG:NH1	0.54	2.71	27	6
1:A:81:ARG:CG	1:A:81:ARG:HH11	0.54	2.16	8	3
1:A:133:TYR:O	1:A:133:TYR:CG	0.53	2.61	2	30
1:A:118:ARG:CG	1:A:118:ARG:NH1	0.53	2.71	16	8
1:A:135:LEU:O	1:A:139:THR:HG23	0.53	2.03	1	4
1:A:90:ARG:CG	1:A:90:ARG:NH1	0.53	2.71	19	3
1:A:118:ARG:NH1	1:A:118:ARG:CG	0.53	2.71	19	4
1:A:53:ARG:O	1:A:55:LYS:N	0.53	2.42	16	6
1:A:45:PRO:O	1:A:81:ARG:NH1	0.53	2.42	10	16
1:A:44:HIS:NE2	1:A:50:ASP:OD2	0.53	2.42	13	6
1:A:36:ASN:ND2	1:A:144:LYS:O	0.53	2.42	4	18
1:A:72:VAL:O	1:A:72:VAL:HG23	0.53	2.03	9	8
1:A:119:LYS:O	1:A:120:TYR:CD1	0.53	2.62	20	12
1:A:123:TRP:CG	1:A:136:GLY:O	0.53	2.62	1	10
1:A:48:ARG:CG	1:A:48:ARG:NH1	0.53	2.71	23	8
1:A:42:ARG:NH1	1:A:57:ASP:OD1	0.53	2.42	17	3
1:A:116:ARG:NH1	1:A:124:TYR:OH	0.53	2.42	3	1
1:A:118:ARG:CG	1:A:118:ARG:HH11	0.53	2.16	11	4
1:A:132:GLN:OE1	1:A:132:GLN:N	0.53	2.42	10	1
1:A:91:LEU:N	1:A:91:LEU:HD12	0.53	2.19	12	2
1:A:91:LEU:CD1	1:A:91:LEU:N	0.53	2.72	12	1
1:A:67:GLU:O	1:A:69:ARG:N	0.53	2.42	28	1
1:A:37:GLY:O	1:A:39:PHE:CE2	0.52	2.62	30	21
1:A:102:PHE:C	1:A:103:PHE:CD1	0.52	2.88	11	26
1:A:123:TRP:CD1	1:A:136:GLY:O	0.52	2.62	11	17
1:A:133:TYR:O	1:A:133:TYR:CD2	0.52	2.62	4	24
1:A:87:GLU:OE1	1:A:88:ASP:N	0.52	2.42	29	3
1:A:116:ARG:NE	1:A:121:THR:O	0.52	2.42	25	6
1:A:75:LYS:NZ	1:A:82:TYR:OH	0.52	2.43	5	1
1:A:88:ASP:OD1	1:A:88:ASP:N	0.52	2.42	21	4
1:A:37:GLY:O	1:A:53:ARG:NH2	0.52	2.42	17	2
1:A:37:GLY:O	1:A:53:ARG:NH1	0.52	2.42	20	1
1:A:88:ASP:N	1:A:88:ASP:OD1	0.52	2.42	29	2
1:A:100:GLU:CD	1:A:100:GLU:N	0.52	2.68	27	3
1:A:118:ARG:C	1:A:120:TYR:N	0.52	2.67	12	20
1:A:124:TYR:CD1	1:A:124:TYR:N	0.52	2.78	18	21
1:A:42:ARG:NH1	1:A:57:ASP:CG	0.52	2.68	17	13
1:A:90:ARG:NH1	1:A:133:TYR:OH	0.52	2.42	16	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:63:GLN:OE1	1:A:64:LEU:N	0.52	2.42	22	1
1:A:64:LEU:N	1:A:64:LEU:CD1	0.52	2.73	14	30
1:A:113:ASN:N	1:A:113:ASN:OD1	0.52	2.42	6	8
1:A:42:ARG:NH1	1:A:44:HIS:CE1	0.52	2.77	18	1
1:A:42:ARG:NH2	1:A:57:ASP:CG	0.52	2.68	24	1
1:A:122:SER:O	1:A:123:TRP:CD1	0.52	2.63	26	4
1:A:68:GLU:O	1:A:70:GLY:N	0.52	2.42	2	2
1:A:132:GLN:N	1:A:132:GLN:CD	0.52	2.68	14	6
1:A:42:ARG:NH1	1:A:57:ASP:OD2	0.52	2.42	22	2
1:A:36:ASN:ND2	1:A:144:LYS:C	0.52	2.68	16	5
1:A:37:GLY:O	1:A:39:PHE:CD2	0.52	2.62	22	20
1:A:120:TYR:C	1:A:122:SER:H	0.52	2.13	26	25
1:A:116:ARG:NH2	1:A:124:TYR:CE1	0.52	2.78	3	1
1:A:77:VAL:O	1:A:80:ASN:ND2	0.52	2.42	21	5
1:A:39:PHE:CZ	1:A:129:ARG:O	0.52	2.63	13	3
1:A:42:ARG:HE	1:A:59:HIS:CB	0.52	2.18	18	2
1:A:110:ASN:ND2	1:A:112:TYR:H	0.52	2.03	21	2
1:A:69:ARG:CG	1:A:69:ARG:NH1	0.52	2.71	2	3
1:A:116:ARG:NH2	1:A:121:THR:O	0.52	2.43	3	2
1:A:50:ASP:N	1:A:50:ASP:OD1	0.52	2.42	13	3
1:A:67:GLU:OE1	1:A:102:PHE:CE1	0.52	2.62	17	1
1:A:106:ARG:CG	1:A:106:ARG:NH1	0.51	2.71	29	3
1:A:64:LEU:N	1:A:64:LEU:HD12	0.51	2.20	22	1
1:A:42:ARG:NH2	1:A:52:VAL:HG21	0.51	2.21	26	1
1:A:104:PHE:CD2	1:A:116:ARG:NH1	0.51	2.78	16	1
1:A:53:ARG:NH1	1:A:53:ARG:CG	0.51	2.71	26	2
1:A:137:SER:C	1:A:139:THR:N	0.51	2.68	20	2
1:A:81:ARG:CG	1:A:81:ARG:NH1	0.51	2.71	27	3
1:A:116:ARG:NE	1:A:124:TYR:OH	0.51	2.44	12	2
1:A:44:HIS:NE2	1:A:50:ASP:OD1	0.51	2.43	18	2
1:A:42:ARG:NH2	1:A:57:ASP:OD1	0.51	2.43	24	4
1:A:111:ASN:N	1:A:111:ASN:OD1	0.51	2.43	30	5
1:A:132:GLN:N	1:A:132:GLN:OE1	0.51	2.44	2	1
1:A:88:ASP:CG	1:A:89:GLY:N	0.51	2.69	7	8
1:A:123:TRP:CD2	1:A:136:GLY:O	0.51	2.64	9	1
1:A:108:GLU:CD	1:A:114:THR:HG23	0.51	2.30	23	1
1:A:135:LEU:C	1:A:137:SER:N	0.51	2.68	26	5
1:A:106:ARG:CZ	1:A:116:ARG:HH21	0.51	2.18	12	1
1:A:88:ASP:OD1	1:A:89:GLY:N	0.51	2.43	4	6
1:A:44:HIS:CD2	1:A:50:ASP:OD1	0.51	2.64	3	2
1:A:143:GLN:CD	1:A:143:GLN:N	0.51	2.68	6	2

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:140:GLY:N	1:A:143:GLN:OE1	0.51	2.44	27	1
1:A:28:ASP:N	1:A:28:ASP:OD1	0.51	2.43	19	1
1:A:73:SER:OG	1:A:102:PHE:CD2	0.51	2.63	27	1
1:A:106:ARG:HE	1:A:116:ARG:NH2	0.51	2.04	29	1
1:A:106:ARG:NH2	1:A:108:GLU:CD	0.51	2.69	9	1
1:A:134:LYS:HZ2	1:A:143:GLN:NE2	0.50	2.04	10	1
1:A:135:LEU:O	1:A:137:SER:N	0.50	2.44	26	1
1:A:118:ARG:O	1:A:120:TYR:N	0.50	2.44	14	17
1:A:86:LYS:NZ	1:A:90:ARG:NH2	0.50	2.60	4	1
1:A:87:GLU:CD	1:A:87:GLU:N	0.50	2.68	9	4
1:A:113:ASN:N	1:A:113:ASN:HD22	0.50	2.05	23	1
1:A:112:TYR:C	1:A:113:ASN:ND2	0.50	2.69	13	1
1:A:55:LYS:CD	1:A:55:LYS:N	0.50	2.74	25	1
1:A:42:ARG:HH11	1:A:57:ASP:CG	0.50	2.15	15	3
1:A:137:SER:O	1:A:139:THR:N	0.49	2.45	20	1
1:A:86:LYS:N	1:A:86:LYS:CD	0.49	2.75	25	1
1:A:81:ARG:NE	1:A:93:ALA:O	0.49	2.45	30	1
1:A:83:LEU:O	1:A:101:CYS:SG	0.49	2.70	19	2
1:A:36:ASN:HD21	1:A:144:LYS:CB	0.49	2.20	17	4
1:A:57:ASP:C	1:A:59:HIS:H	0.49	2.16	11	23
1:A:97:VAL:HG12	1:A:98:THR:N	0.49	2.23	5	2
1:A:110:ASN:ND2	1:A:110:ASN:N	0.49	2.60	28	3
1:A:106:ARG:NH1	1:A:116:ARG:NH2	0.49	2.60	12	1
1:A:92:LEU:C	1:A:92:LEU:CD1	0.49	2.83	2	25
1:A:44:HIS:CD2	1:A:50:ASP:OD2	0.49	2.65	12	3
1:A:129:ARG:HH11	1:A:129:ARG:CG	0.49	2.20	30	2
1:A:50:ASP:OD1	1:A:51:GLY:N	0.49	2.46	29	1
1:A:77:VAL:HG12	1:A:78:SER:N	0.49	2.22	17	3
1:A:143:GLN:HE21	1:A:143:GLN:N	0.49	2.06	25	1
1:A:31:ARG:NH1	1:A:61:LYS:HZ2	0.49	2.05	16	1
1:A:129:ARG:CG	1:A:129:ARG:NH1	0.49	2.73	30	2
1:A:106:ARG:NH2	1:A:108:GLU:OE1	0.49	2.45	22	2
1:A:36:ASN:HD21	1:A:144:LYS:C	0.49	2.16	28	2
1:A:120:TYR:C	1:A:122:SER:N	0.48	2.69	8	27
1:A:63:GLN:NE2	1:A:64:LEU:O	0.48	2.45	3	2
1:A:46:ASP:OD2	1:A:48:ARG:NH1	0.48	2.46	11	2
1:A:134:LYS:NZ	1:A:143:GLN:OE1	0.48	2.43	28	1
1:A:110:ASN:N	1:A:110:ASN:HD22	0.48	2.05	28	3
1:A:30:LYS:HG3	1:A:64:LEU:HD22	0.48	1.86	21	1
1:A:46:ASP:OD2	1:A:48:ARG:CD	0.48	2.62	21	1
1:A:46:ASP:OD1	1:A:48:ARG:CG	0.47	2.62	10	2

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:112:TYR:O	1:A:113:ASN:ND2	0.47	2.47	30	1
1:A:84:ALA:HB3	1:A:92:LEU:O	0.47	2.10	2	1
1:A:86:LYS:HZ3	1:A:90:ARG:NH2	0.47	2.06	4	1
1:A:61:LYS:O	1:A:77:VAL:CG2	0.47	2.63	19	4
1:A:129:ARG:CD	1:A:129:ARG:C	0.47	2.87	27	1
1:A:50:ASP:CG	1:A:51:GLY:N	0.47	2.69	4	2
1:A:40:PHE:CD1	1:A:40:PHE:N	0.47	2.83	24	3
1:A:110:ASN:C	1:A:112:TYR:H	0.47	2.18	21	23
1:A:116:ARG:CZ	1:A:124:TYR:OH	0.47	2.63	3	4
1:A:84:ALA:HB2	1:A:101:CYS:SG	0.47	2.50	22	1
1:A:97:VAL:O	1:A:97:VAL:CG1	0.47	2.63	23	10
1:A:110:ASN:N	1:A:110:ASN:OD1	0.47	2.48	18	5
1:A:118:ARG:C	1:A:120:TYR:H	0.47	2.18	29	17
1:A:31:ARG:HH22	1:A:153:ALA:CB	0.47	2.23	22	1
1:A:137:SER:OG	1:A:138:LYS:N	0.47	2.48	20	2
1:A:89:GLY:O	1:A:91:LEU:CD1	0.46	2.63	13	13
1:A:120:TYR:O	1:A:122:SER:N	0.46	2.48	8	7
1:A:44:HIS:CD2	1:A:44:HIS:N	0.46	2.84	22	1
1:A:92:LEU:HD12	1:A:92:LEU:O	0.46	2.11	6	8
1:A:39:PHE:CE2	1:A:129:ARG:O	0.46	2.68	21	2
1:A:97:VAL:CG1	1:A:98:THR:N	0.46	2.78	5	3
1:A:144:LYS:C	1:A:146:ILE:N	0.46	2.73	13	10
1:A:123:TRP:C	1:A:124:TYR:CD1	0.46	2.93	1	1
1:A:106:ARG:HH21	1:A:108:GLU:CD	0.46	2.17	9	1
1:A:106:ARG:NH2	1:A:116:ARG:HH21	0.46	2.09	12	1
1:A:31:ARG:NH2	1:A:55:LYS:O	0.46	2.49	14	1
1:A:105:GLU:OE2	1:A:115:TYR:CE1	0.46	2.69	28	1
1:A:116:ARG:CZ	1:A:121:THR:O	0.46	2.64	3	1
1:A:86:LYS:HZ3	1:A:90:ARG:HH21	0.46	1.53	4	1
1:A:144:LYS:C	1:A:146:ILE:H	0.46	2.19	12	6
1:A:49:VAL:HG21	1:A:83:LEU:HD11	0.46	1.86	7	1
1:A:81:ARG:HB3	1:A:93:ALA:HB1	0.46	1.87	13	1
1:A:135:LEU:C	1:A:137:SER:H	0.46	2.18	26	10
1:A:98:THR:OG1	1:A:101:CYS:SG	0.45	2.71	6	2
1:A:81:ARG:NH1	1:A:95:LYS:CG	0.45	2.79	8	1
1:A:127:LEU:O	1:A:134:LYS:NZ	0.45	2.49	23	1
1:A:49:VAL:HG21	1:A:83:LEU:CD1	0.45	2.42	1	3
1:A:69:ARG:O	1:A:69:ARG:CG	0.45	2.63	11	2
1:A:46:ASP:OD1	1:A:48:ARG:N	0.45	2.49	22	3
1:A:110:ASN:C	1:A:112:TYR:N	0.45	2.74	7	23
1:A:85:MET:SD	1:A:91:LEU:CD1	0.45	3.05	7	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:92:LEU:CD1	1:A:92:LEU:C	0.45	2.90	16	5
1:A:87:GLU:CD	1:A:88:ASP:N	0.45	2.75	29	1
1:A:98:THR:C	1:A:100:GLU:N	0.44	2.73	25	3
1:A:99:ASP:C	1:A:101:CYS:H	0.44	2.21	14	2
1:A:57:ASP:C	1:A:59:HIS:N	0.44	2.75	7	21
1:A:72:VAL:O	1:A:72:VAL:CG2	0.44	2.65	9	1
1:A:144:LYS:CB	1:A:144:LYS:NZ	0.44	2.80	27	1
1:A:61:LYS:O	1:A:77:VAL:HG23	0.44	2.13	19	3
1:A:31:ARG:CB	1:A:31:ARG:CZ	0.44	2.96	11	3
1:A:48:ARG:CZ	1:A:48:ARG:CB	0.44	2.95	15	3
1:A:51:GLY:N	1:A:131:GLY:O	0.44	2.50	16	1
1:A:110:ASN:OD1	1:A:110:ASN:N	0.44	2.51	19	2
1:A:100:GLU:O	1:A:101:CYS:SG	0.44	2.72	28	2
1:A:44:HIS:NE2	1:A:50:ASP:CG	0.44	2.76	3	2
1:A:81:ARG:HE	1:A:95:LYS:NZ	0.44	2.10	3	1
1:A:59:HIS:CD2	1:A:78:SER:OG	0.44	2.71	6	1
1:A:67:GLU:OE1	1:A:102:PHE:CD1	0.44	2.71	17	1
1:A:37:GLY:O	1:A:53:ARG:CZ	0.44	2.65	20	1
1:A:54:GLU:O	1:A:57:ASP:OD1	0.44	2.36	12	1
1:A:59:HIS:C	1:A:61:LYS:H	0.44	2.21	26	2
1:A:108:GLU:C	1:A:110:ASN:H	0.44	2.20	30	1
1:A:113:ASN:O	1:A:146:ILE:O	0.43	2.36	1	1
1:A:31:ARG:CZ	1:A:31:ARG:CB	0.43	2.96	21	7
1:A:100:GLU:C	1:A:101:CYS:SG	0.43	3.01	29	3
1:A:42:ARG:NH2	1:A:59:HIS:ND1	0.43	2.66	18	1
1:A:53:ARG:O	1:A:54:GLU:C	0.43	2.60	16	4
1:A:37:GLY:O	1:A:38:GLY:C	0.43	2.61	22	21
1:A:100:GLU:N	1:A:100:GLU:OE1	0.43	2.52	24	1
1:A:55:LYS:O	1:A:60:ILE:CD1	0.43	2.66	29	1
1:A:116:ARG:NE	1:A:124:TYR:CZ	0.43	2.86	29	1
1:A:47:GLY:O	1:A:48:ARG:C	0.43	2.62	29	12
1:A:110:ASN:O	1:A:112:TYR:CD2	0.43	2.72	2	2
1:A:36:ASN:OD1	1:A:144:LYS:O	0.43	2.37	5	8
1:A:53:ARG:CZ	1:A:53:ARG:CB	0.43	2.95	13	6
1:A:33:TYR:CD1	1:A:40:PHE:CZ	0.43	3.07	24	1
1:A:36:ASN:OD1	1:A:145:ALA:N	0.43	2.52	29	1
1:A:109:SER:C	1:A:111:ASN:H	0.43	2.21	4	3
1:A:67:GLU:O	1:A:71:VAL:O	0.43	2.36	10	1
1:A:55:LYS:CD	1:A:55:LYS:O	0.43	2.67	17	1
1:A:134:LYS:O	1:A:135:LEU:C	0.43	2.62	16	16
1:A:67:GLU:O	1:A:68:GLU:O	0.43	2.37	3	2

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:137:SER:C	1:A:139:THR:H	0.43	2.22	4	2
1:A:46:ASP:OD1	1:A:46:ASP:C	0.43	2.62	18	5
1:A:56:SER:O	1:A:57:ASP:C	0.43	2.61	1	2
1:A:65:GLN:O	1:A:73:SER:N	0.43	2.52	17	1
1:A:99:ASP:OD1	1:A:99:ASP:N	0.43	2.52	18	1
1:A:63:GLN:OE1	1:A:64:LEU:O	0.43	2.37	22	1
1:A:143:GLN:C	1:A:145:ALA:N	0.43	2.76	26	2
1:A:50:ASP:OD1	1:A:50:ASP:C	0.43	2.62	12	2
1:A:134:LYS:O	1:A:135:LEU:O	0.43	2.37	12	1
1:A:77:VAL:O	1:A:80:ASN:OD1	0.43	2.37	14	1
1:A:88:ASP:CG	1:A:89:GLY:H	0.43	2.22	20	1
1:A:108:GLU:C	1:A:110:ASN:N	0.43	2.76	30	1
1:A:77:VAL:O	1:A:78:SER:C	0.42	2.62	10	6
1:A:36:ASN:OD1	1:A:144:LYS:C	0.42	2.62	10	4
1:A:110:ASN:C	1:A:110:ASN:HD22	0.42	2.21	21	1
1:A:98:THR:O	1:A:101:CYS:SG	0.42	2.77	1	1
1:A:130:THR:O	1:A:132:GLN:OE1	0.42	2.37	10	5
1:A:141:PRO:C	1:A:143:GLN:H	0.42	2.21	13	3
1:A:138:LYS:N	1:A:138:LYS:CD	0.42	2.82	20	1
1:A:105:GLU:OE2	1:A:115:TYR:CZ	0.42	2.72	4	1
1:A:108:GLU:OE2	1:A:112:TYR:O	0.42	2.38	23	1
1:A:54:GLU:OE2	1:A:56:SER:OG	0.42	2.37	1	1
1:A:40:PHE:O	1:A:41:LEU:C	0.42	2.61	14	7
1:A:50:ASP:O	1:A:50:ASP:OD1	0.42	2.37	17	2
1:A:116:ARG:NH2	1:A:124:TYR:OH	0.42	2.53	25	1
1:A:130:THR:C	1:A:132:GLN:OE1	0.42	2.62	25	2
1:A:63:GLN:OE1	1:A:63:GLN:O	0.42	2.37	11	1
1:A:54:GLU:OE2	1:A:57:ASP:OD1	0.42	2.37	22	1
1:A:63:GLN:OE1	1:A:63:GLN:C	0.42	2.62	22	1
1:A:143:GLN:C	1:A:145:ALA:H	0.42	2.23	22	3
1:A:106:ARG:O	1:A:114:THR:OG1	0.42	2.37	3	2
1:A:90:ARG:CZ	1:A:90:ARG:CB	0.42	2.96	13	1
1:A:130:THR:CG2	1:A:132:GLN:OE1	0.42	2.64	25	1
1:A:85:MET:O	1:A:100:GLU:OE1	0.42	2.38	2	2
1:A:98:THR:O	1:A:99:ASP:C	0.42	2.62	6	3
1:A:54:GLU:OE2	1:A:57:ASP:OD2	0.42	2.37	11	1
1:A:122:SER:O	1:A:123:TRP:CG	0.42	2.73	25	1
1:A:103:PHE:CD1	1:A:103:PHE:N	0.42	2.87	28	1
1:A:63:GLN:NE2	1:A:75:LYS:NZ	0.42	2.68	8	1
1:A:118:ARG:O	1:A:119:LYS:C	0.42	2.63	13	1
1:A:78:SER:O	1:A:80:ASN:ND2	0.42	2.52	25	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:63:GLN:CD	1:A:63:GLN:C	0.42	2.87	12	2
1:A:137:SER:OG	1:A:138:LYS:NZ	0.42	2.43	12	1
1:A:85:MET:O	1:A:100:GLU:CD	0.42	2.63	9	1
1:A:109:SER:C	1:A:110:ASN:OD1	0.42	2.63	11	1
1:A:102:PHE:O	1:A:103:PHE:CD1	0.42	2.73	17	1
1:A:36:ASN:CG	1:A:144:LYS:O	0.41	2.63	12	7
1:A:104:PHE:N	1:A:116:ARG:O	0.41	2.53	18	1
1:A:108:GLU:OE1	1:A:114:THR:HG23	0.41	2.15	20	1
1:A:114:THR:HG22	1:A:146:ILE:O	0.41	2.15	22	1
1:A:30:LYS:HE2	1:A:64:LEU:HD22	0.41	1.91	2	1
1:A:112:TYR:C	1:A:113:ASN:OD1	0.41	2.63	9	2
1:A:108:GLU:CD	1:A:112:TYR:O	0.41	2.62	23	1
1:A:50:ASP:OD1	1:A:50:ASP:N	0.41	2.53	2	1
1:A:78:SER:O	1:A:80:ASN:OD1	0.41	2.38	22	3
1:A:42:ARG:NE	1:A:59:HIS:CB	0.41	2.83	19	1
1:A:110:ASN:ND2	1:A:110:ASN:H	0.41	2.13	28	1
1:A:54:GLU:C	1:A:56:SER:N	0.41	2.78	7	1
1:A:97:VAL:O	1:A:97:VAL:HG12	0.41	2.16	10	2
1:A:136:GLY:O	1:A:139:THR:OG1	0.41	2.37	11	1
1:A:110:ASN:C	1:A:111:ASN:OD1	0.41	2.64	14	2
1:A:67:GLU:O	1:A:67:GLU:OE1	0.41	2.38	14	1
1:A:87:GLU:OE1	1:A:87:GLU:C	0.41	2.63	2	1
1:A:113:ASN:HD22	1:A:113:ASN:H	0.41	1.54	7	1
1:A:65:GLN:OE1	1:A:66:ALA:O	0.41	2.38	26	1
1:A:67:GLU:O	1:A:67:GLU:CD	0.41	2.63	11	1
1:A:86:LYS:HE3	1:A:92:LEU:HD21	0.41	1.92	13	1
1:A:50:ASP:OD2	1:A:52:VAL:CG1	0.41	2.69	16	1
1:A:59:HIS:C	1:A:61:LYS:N	0.41	2.78	23	1
1:A:42:ARG:CZ	1:A:57:ASP:OD2	0.41	2.69	24	1
1:A:79:ALA:O	1:A:80:ASN:C	0.41	2.63	29	1
1:A:84:ALA:HB3	1:A:92:LEU:HD12	0.41	1.93	26	2
1:A:46:ASP:CG	1:A:48:ARG:NH1	0.40	2.79	15	1
1:A:72:VAL:CG2	1:A:103:PHE:O	0.40	2.69	17	1
1:A:32:LEU:HD12	1:A:41:LEU:HD23	0.40	1.92	18	1
1:A:42:ARG:CZ	1:A:57:ASP:OD1	0.40	2.69	2	1
1:A:67:GLU:O	1:A:68:GLU:C	0.40	2.62	10	1
1:A:34:CYS:O	1:A:35:LYS:C	0.40	2.64	16	2
1:A:112:TYR:C	1:A:113:ASN:HD22	0.40	2.25	30	1
1:A:71:VAL:HG12	1:A:72:VAL:N	0.40	2.30	1	1
1:A:111:ASN:N	1:A:111:ASN:HD22	0.40	2.13	23	1
1:A:110:ASN:ND2	1:A:110:ASN:C	0.40	2.79	21	1

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:63:GLN:N	1:A:75:LYS:O	0.40	2.55	9	1
1:A:63:GLN:C	1:A:63:GLN:OE1	0.40	2.64	12	1
1:A:56:SER:O	1:A:56:SER:OG	0.40	2.38	17	1
1:A:102:PHE:C	1:A:103:PHE:CD2	0.40	3.00	18	1
1:A:82:TYR:CG	1:A:97:VAL:HG22	0.40	2.52	27	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	126/155 (81%)	106±3 (84±2%)	17±3 (14±2%)	2±1 (2±1%)	8	51
All	All	3780/4650 (81%)	3189 (84%)	518 (14%)	73 (2%)	8	51

All 18 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	38	GLY	23
1	A	77	VAL	11
1	A	54	GLU	7
1	A	68	GLU	5
1	A	128	LYS	3
1	A	69	ARG	3
1	A	121	THR	3
1	A	48	ARG	3
1	A	119	LYS	3
1	A	135	LEU	2
1	A	110	ASN	2
1	A	58	PRO	2
1	A	86	LYS	1
1	A	141	PRO	1
1	A	88	ASP	1
1	A	89	GLY	1
1	A	129	ARG	1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Models (Total)
1	A	70	GLY	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	108/129 (84%)	97±3 (90±2%)	11±3 (10±2%)	10	54
All	All	3240/3870 (84%)	2920 (90%)	320 (10%)	10	54

All 54 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	60	ILE	16
1	A	130	THR	16
1	A	118	ARG	16
1	A	57	ASP	15
1	A	55	LYS	14
1	A	96	SER	14
1	A	78	SER	12
1	A	87	GLU	11
1	A	99	ASP	11
1	A	48	ARG	9
1	A	56	SER	9
1	A	116	ARG	9
1	A	109	SER	9
1	A	31	ARG	9
1	A	144	LYS	9
1	A	35	LYS	8
1	A	106	ARG	8
1	A	94	SER	8
1	A	111	ASN	8
1	A	129	ARG	7
1	A	67	GLU	6
1	A	113	ASN	6
1	A	69	ARG	5
1	A	100	GLU	5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Models (Total)
1	A	146	ILE	5
1	A	110	ASN	5
1	A	85	MET	4
1	A	105	GLU	4
1	A	54	GLU	4
1	A	90	ARG	4
1	A	95	LYS	4
1	A	30	LYS	3
1	A	65	GLN	3
1	A	88	ASP	3
1	A	143	GLN	3
1	A	63	GLN	3
1	A	128	LYS	3
1	A	151	MET	3
1	A	53	ARG	3
1	A	81	ARG	3
1	A	107	LEU	3
1	A	61	LYS	3
1	A	108	GLU	2
1	A	138	LYS	2
1	A	75	LYS	2
1	A	52	VAL	2
1	A	132	GLN	2
1	A	72	VAL	1
1	A	68	GLU	1
1	A	28	ASP	1
1	A	71	VAL	1
1	A	134	LYS	1
1	A	119	LYS	1
1	A	86	LYS	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 [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