



Full wwPDB NMR Structure Validation Report ⓘ

Mar 27, 2026 – 09:02 AM UTC

PDB ID : 2NEF / pdb_00002nef
Title : HIV-1 NEF (REGULATORY FACTOR), NMR, 40 STRUCTURES
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Deposited on : 1997-02-12

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

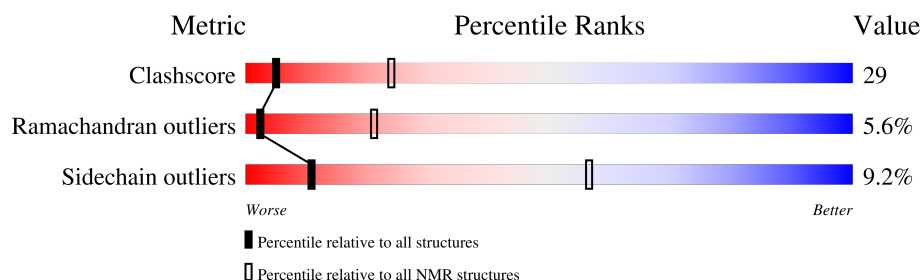
1 Overall quality at a glance

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	136	

2 Ensemble composition and analysis

This entry contains 40 models. Model 13 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:74-A:128, A:134-A:147, A:181-A:195 (84)	0.68	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 5 clusters and 9 single-model clusters were found.

Cluster number	Models
1	2, 3, 7, 9, 10, 13, 16, 17, 19, 21, 23, 25, 27, 29, 36, 40
2	1, 4, 18, 20, 22, 24, 34
3	5, 11, 32, 38
4	6, 8
5	30, 39
Single-model clusters	12; 14; 15; 26; 28; 31; 33; 35; 37

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2230 atoms, of which 1085 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called NEGATIVE FACTOR (F-PROTEIN).

Mol	Chain	Residues	Atoms						Trace
1	A	136	Total	C	H	N	O	S	0
			2230	742	1085	193	207	3	

There are 17 discrepancies between the modelled and reference sequences:

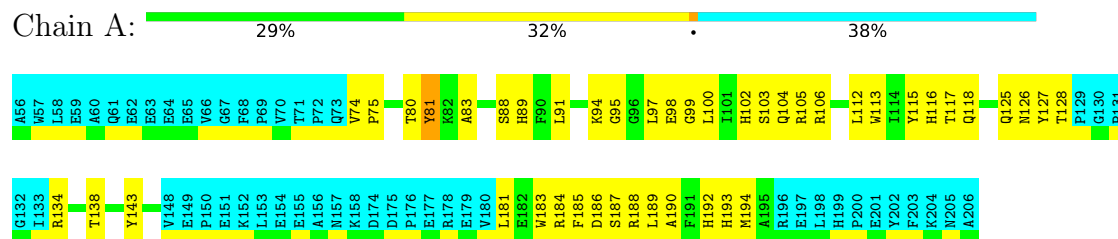
Chain	Residue	Modelled	Actual	Comment	Reference
A	65	GLU	LYS	conflict	UNP Q70627
A	?	-	GLY	deletion	UNP Q70627
A	?	-	GLU	deletion	UNP Q70627
A	?	-	ASN	deletion	UNP Q70627
A	?	-	THR	deletion	UNP Q70627
A	?	-	SER	deletion	UNP Q70627
A	?	-	LEU	deletion	UNP Q70627
A	?	-	LEU	deletion	UNP Q70627
A	?	-	HIS	deletion	UNP Q70627
A	?	-	PRO	deletion	UNP Q70627
A	?	-	VAL	deletion	UNP Q70627
A	?	-	SER	deletion	UNP Q70627
A	?	-	LEU	deletion	UNP Q70627
A	?	-	HIS	deletion	UNP Q70627
A	?	-	GLY	deletion	UNP Q70627
A	?	-	MET	deletion	UNP Q70627
A	194	MET	VAL	conflict	UNP Q70627

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: NEGATIVE FACTOR (F-PROTEIN)

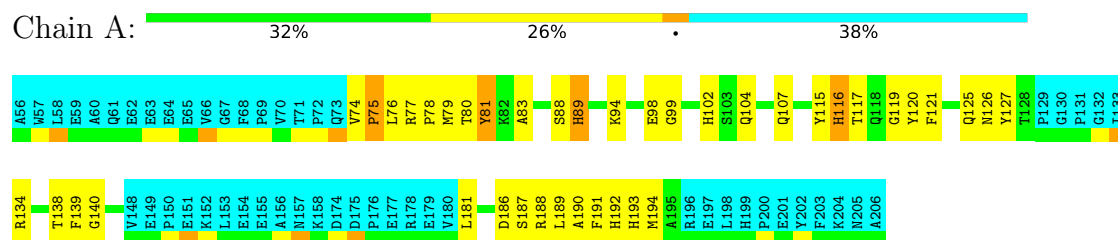


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

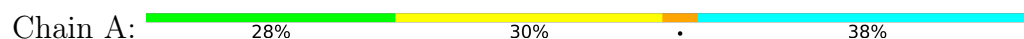
4.2.1 Score per residue for model 1

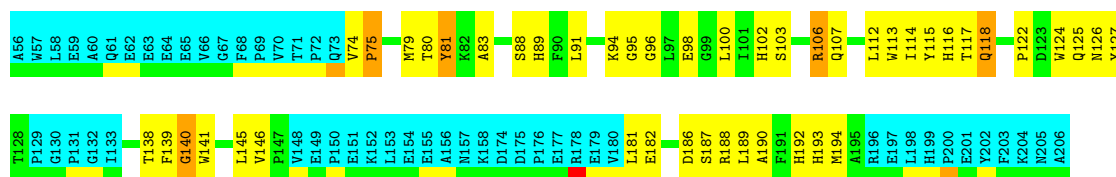
- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



4.2.2 Score per residue for model 2

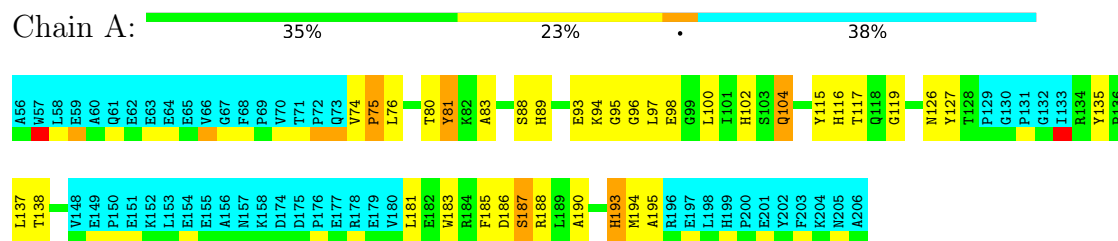
- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)





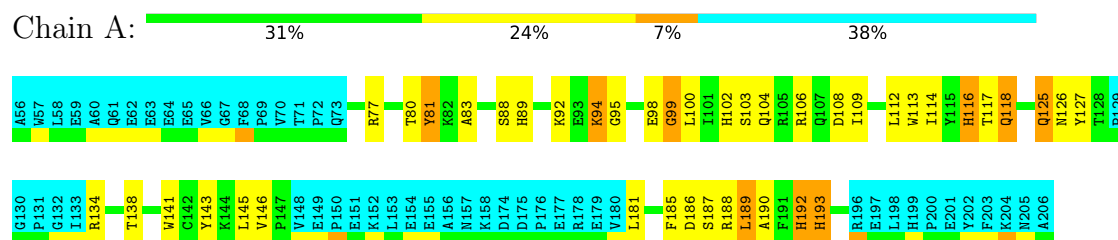
4.2.3 Score per residue for model 3

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



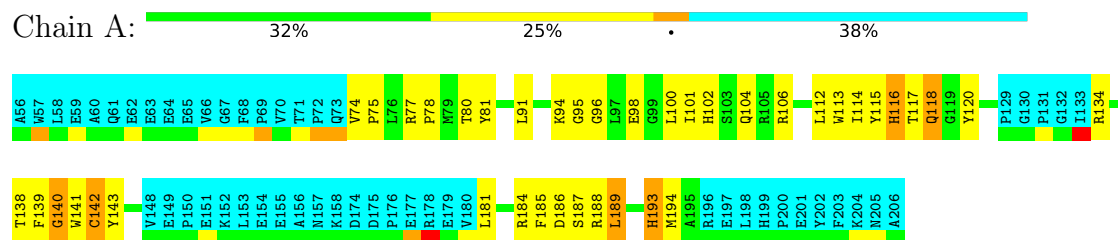
4.2.4 Score per residue for model 4

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



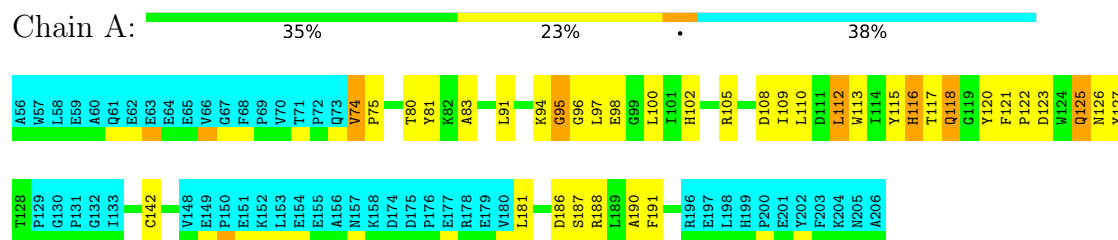
4.2.5 Score per residue for model 5

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



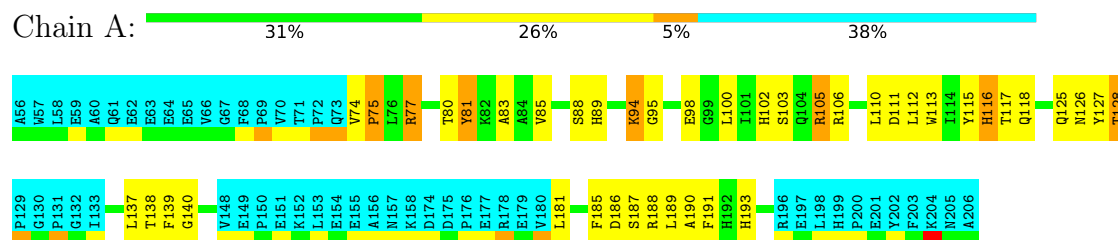
4.2.6 Score per residue for model 6

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



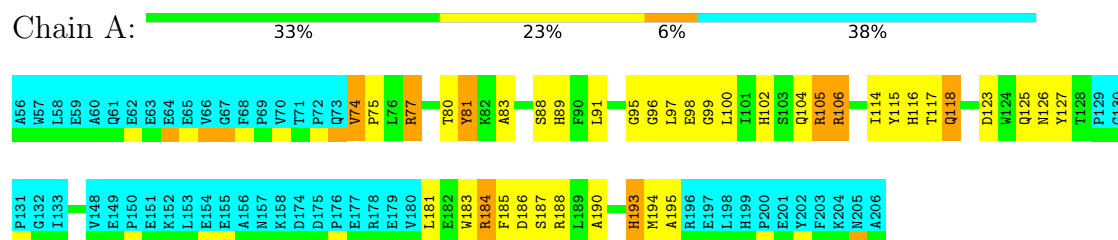
4.2.7 Score per residue for model 7

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



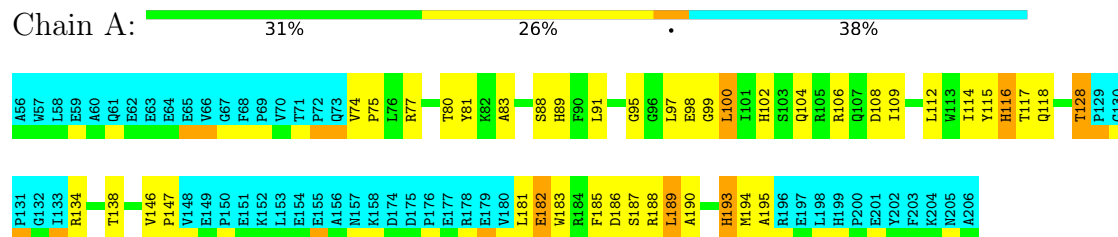
4.2.8 Score per residue for model 8

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



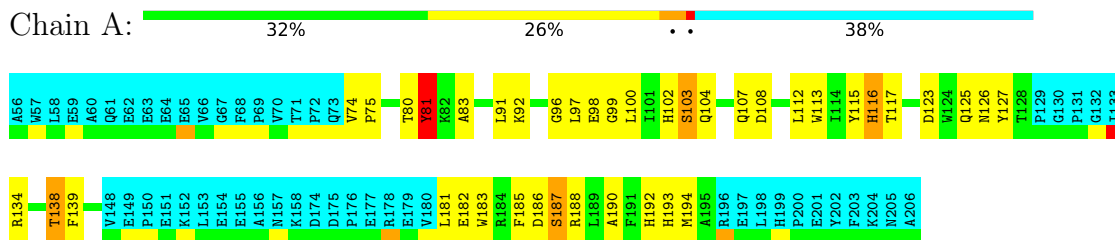
4.2.9 Score per residue for model 9

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



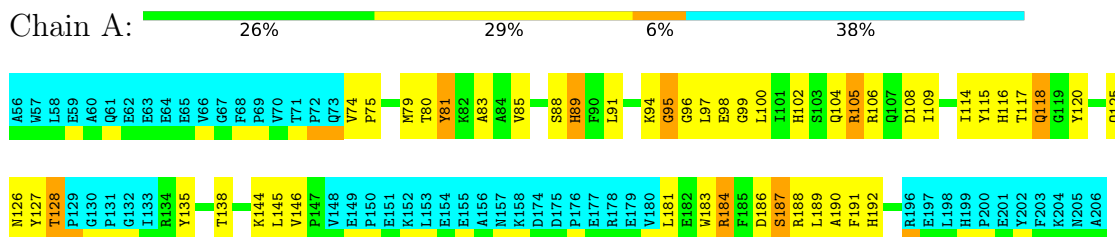
4.2.10 Score per residue for model 10

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



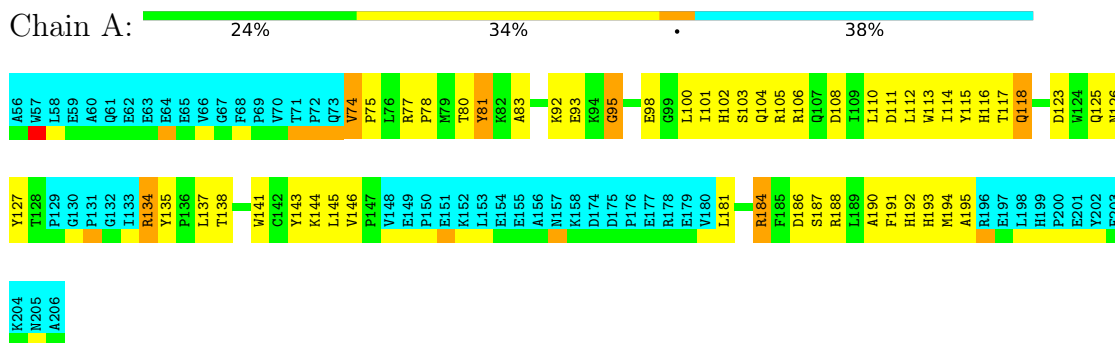
4.2.11 Score per residue for model 11

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



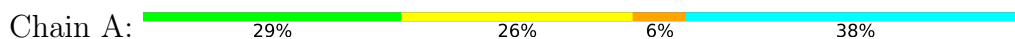
4.2.12 Score per residue for model 12

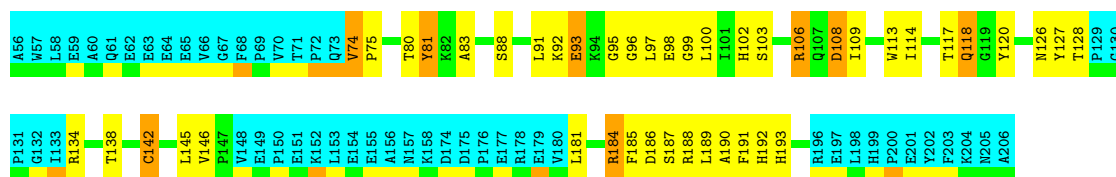
- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



4.2.13 Score per residue for model 13 (medoid)

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

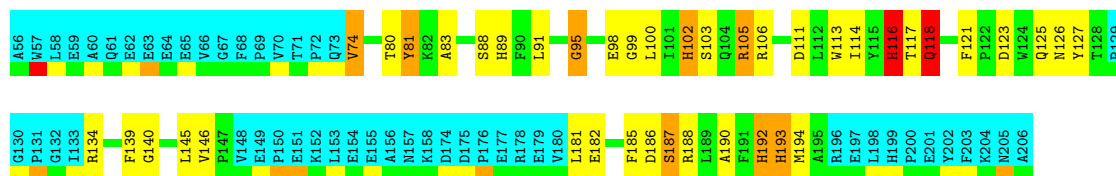




4.2.14 Score per residue for model 14

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

Chain A: 32% 23% 6% • 38%



4.2.15 Score per residue for model 15

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

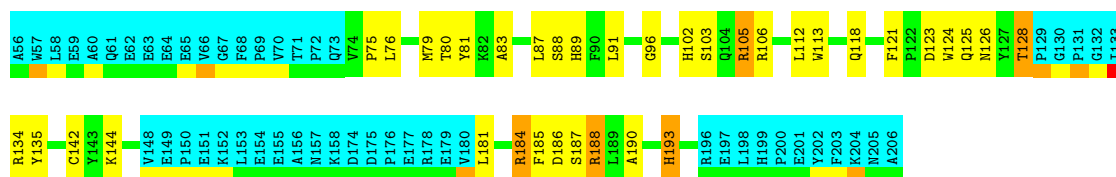
Chain A: 38% 21% • 38%



4.2.16 Score per residue for model 16

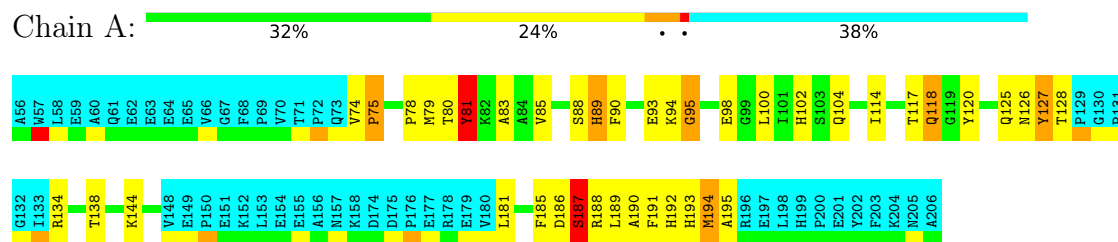
- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

Chain A: 35% 23% • 38%



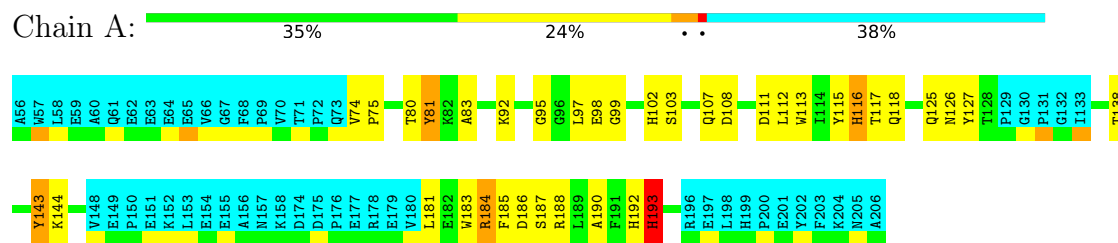
4.2.17 Score per residue for model 17

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



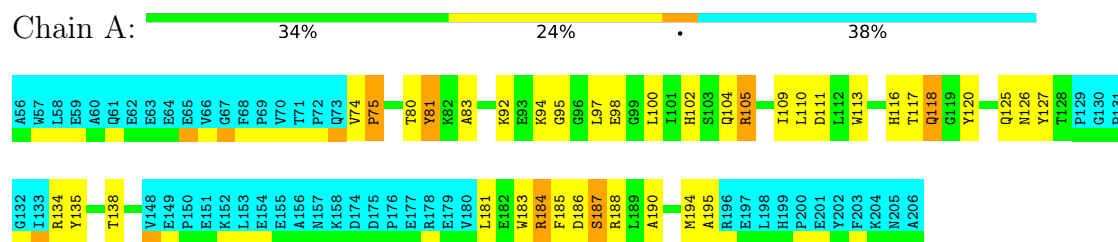
4.2.18 Score per residue for model 18

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



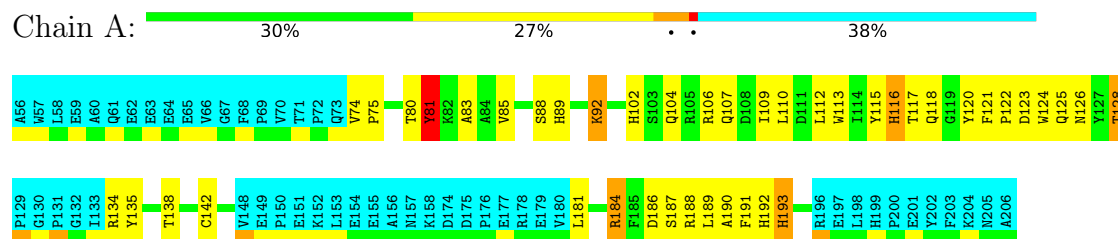
4.2.19 Score per residue for model 19

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



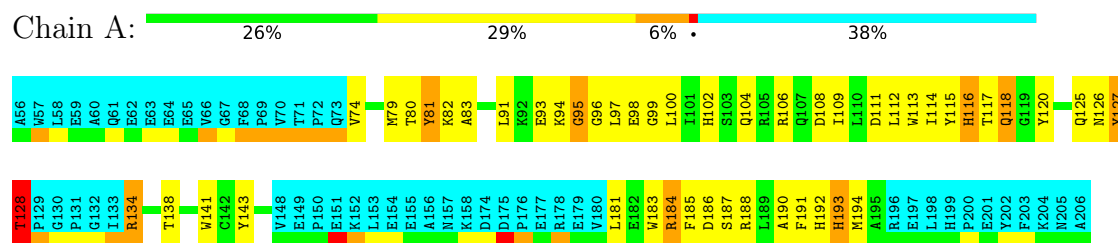
4.2.20 Score per residue for model 20

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



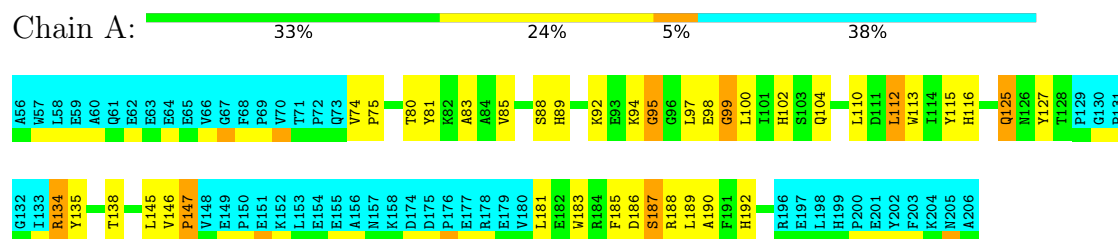
4.2.21 Score per residue for model 21

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



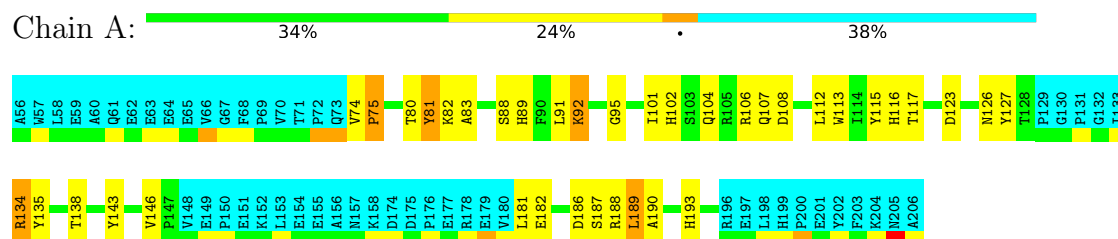
4.2.22 Score per residue for model 22

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



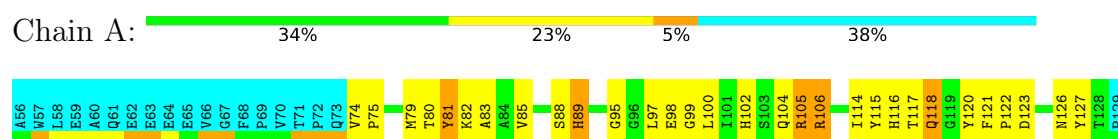
4.2.23 Score per residue for model 23

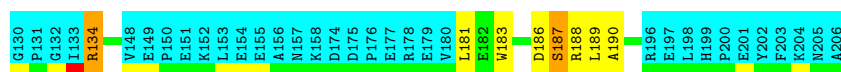
- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



4.2.24 Score per residue for model 24

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

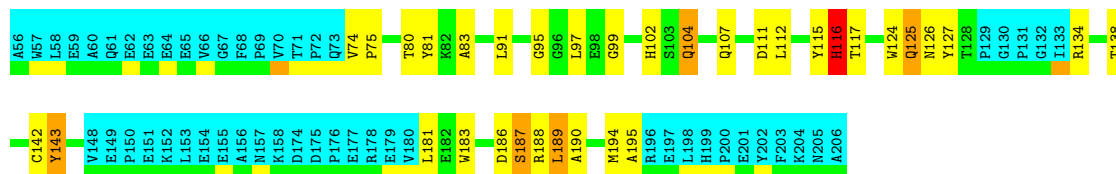




4.2.25 Score per residue for model 25

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

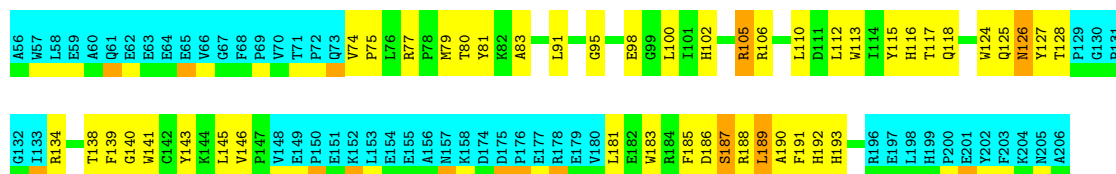
Chain A: 37% 21% . . 38%



4.2.26 Score per residue for model 26

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

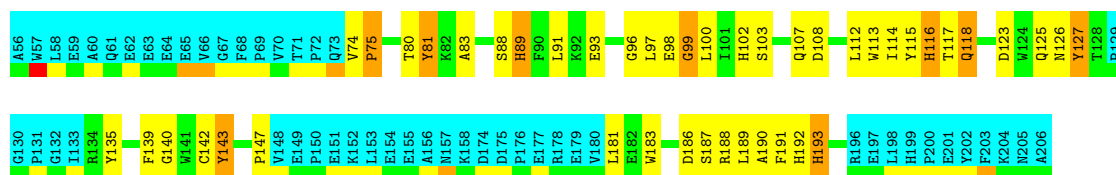
Chain A: 29% 30% . 38%



4.2.27 Score per residue for model 27

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

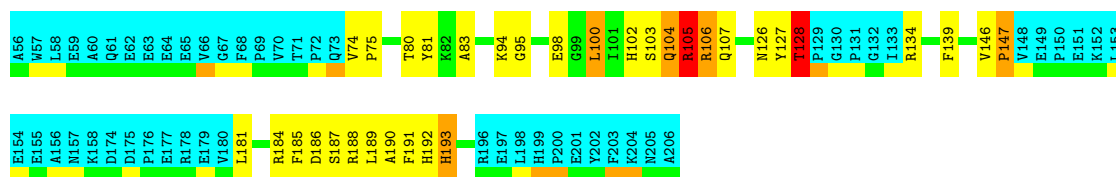
Chain A: 29% 26% 7% 38%



4.2.28 Score per residue for model 28

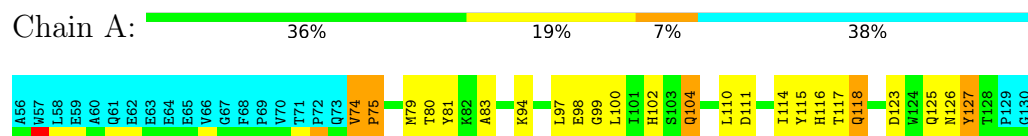
- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

Chain A: 38% 19% . . 38%



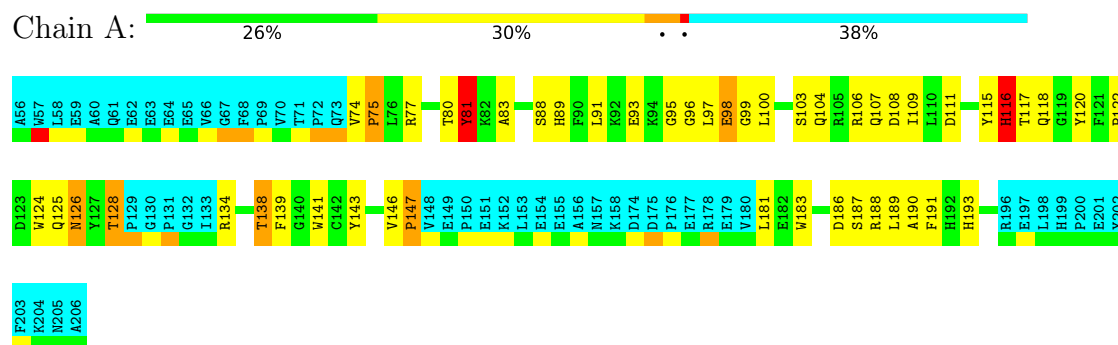
4.2.29 Score per residue for model 29

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



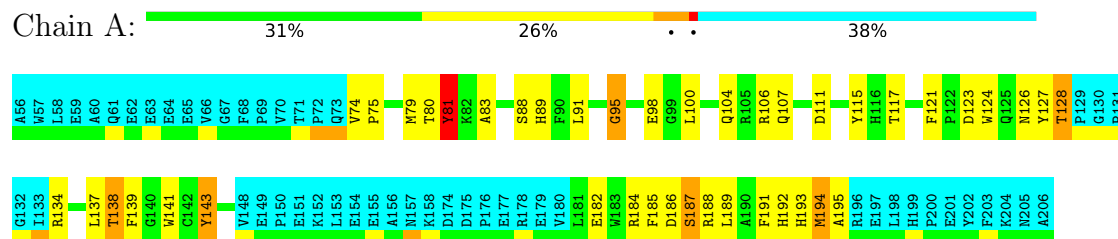
4.2.30 Score per residue for model 30

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



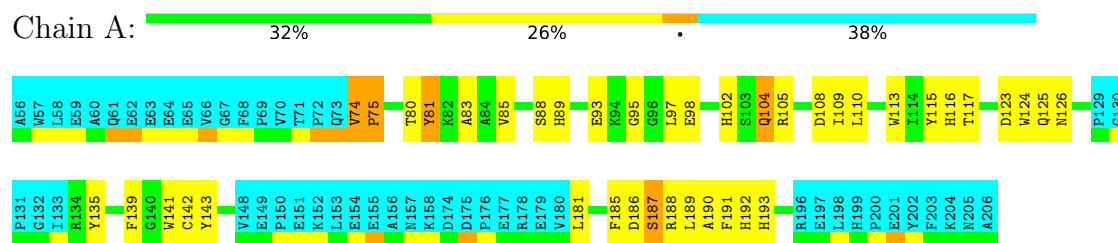
4.2.31 Score per residue for model 31

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



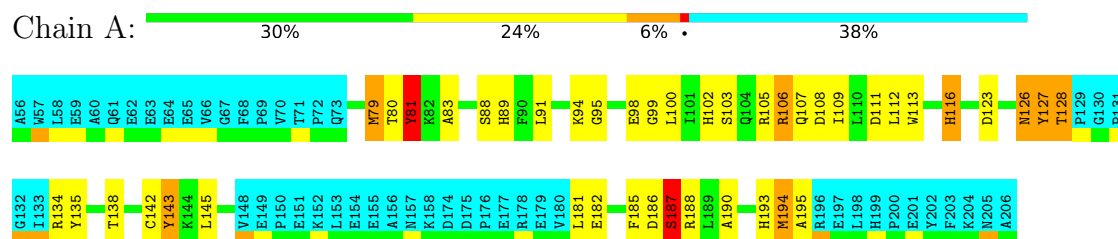
4.2.32 Score per residue for model 32

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



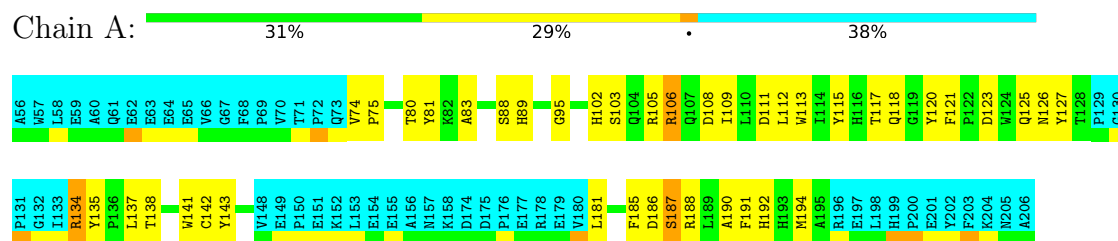
4.2.33 Score per residue for model 33

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



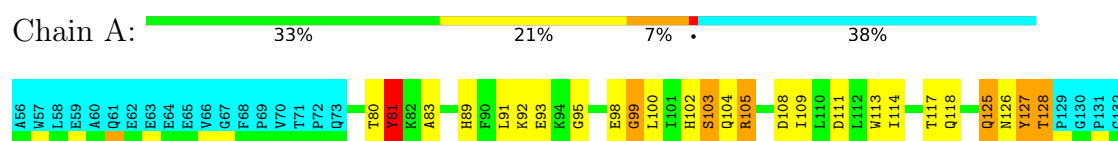
4.2.34 Score per residue for model 34

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



4.2.35 Score per residue for model 35

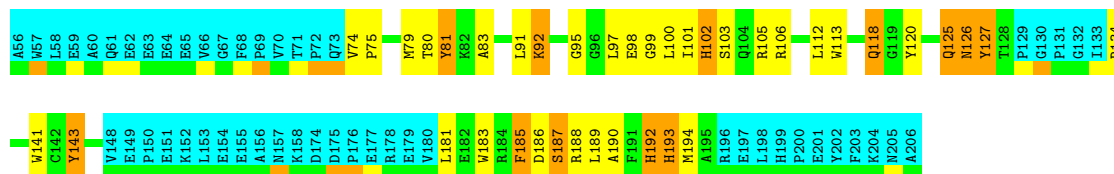
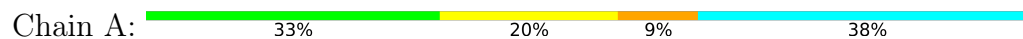
- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)





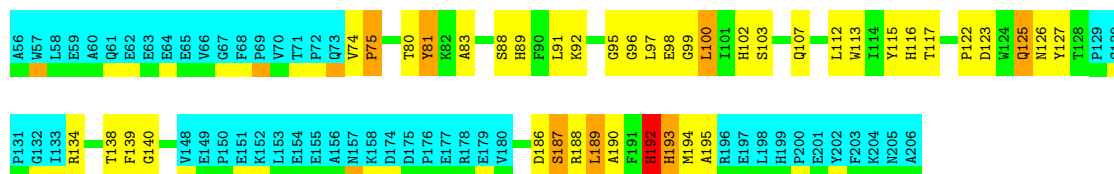
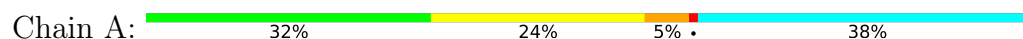
4.2.36 Score per residue for model 36

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



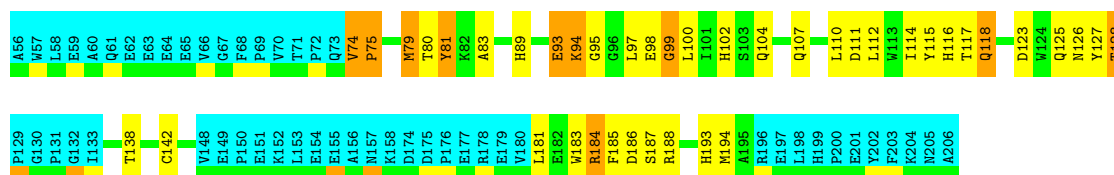
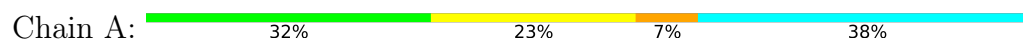
4.2.37 Score per residue for model 37

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



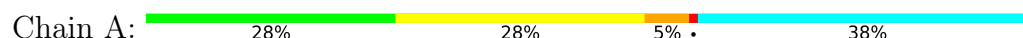
4.2.38 Score per residue for model 38

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)



4.2.39 Score per residue for model 39

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

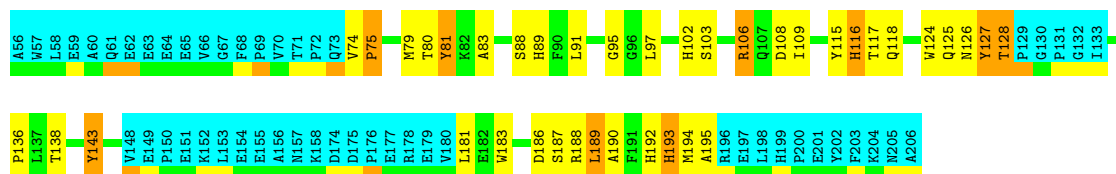




4.2.40 Score per residue for model 40

- Molecule 1: NEGATIVE FACTOR (F-PROTEIN)

Chain A: 33% 22% 7% 38%



5 Refinement protocol and experimental data overview ⓘ

The models were refined using the following method: ?.

Of the ? calculated structures, 40 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

6.1 Standard geometry

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	1.44±0.01	0±0/751 (0.0± 0.0%)	1.19±0.01	0±0/1023 (0.0± 0.0%)
All	All	1.44	5/30040 (0.0%)	1.19	9/40920 (0.0%)

All unique bond outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	95	GLY	N-CA	5.83	1.50	1.44	14	5

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	81	TYR	N-CA-C	-5.50	104.67	111.33	40	8
1	A	191	PHE	CA-CB-CG	-5.26	108.54	113.80	6	1

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	722	697	697	42±7
All	All	28880	27880	27880	1663

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 29.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:106:ARG:HE	1:A:107:GLN:N	0.74	1.79	39	2
1:A:103:SER:N	1:A:106:ARG:HH11	0.74	1.81	28	2
1:A:106:ARG:NE	1:A:106:ARG:N	0.72	2.37	28	1
1:A:192:HIS:NE2	1:A:194:MET:HE2	0.72	2.00	21	1
1:A:106:ARG:HE	1:A:106:ARG:N	0.71	1.84	39	1
1:A:194:MET:SD	1:A:195:ALA:N	0.67	2.67	17	2
1:A:106:ARG:HE	1:A:107:GLN:H	0.67	1.32	39	2
1:A:118:GLN:NE2	1:A:118:GLN:H	0.67	1.88	38	6
1:A:103:SER:N	1:A:106:ARG:NH1	0.66	2.43	39	2
1:A:106:ARG:N	1:A:106:ARG:HE	0.66	1.89	28	1
1:A:143:TYR:C	1:A:143:TYR:CD1	0.66	2.73	33	3
1:A:118:GLN:NE2	1:A:120:TYR:CG	0.65	2.65	36	8
1:A:102:HIS:NE2	1:A:106:ARG:NH2	0.65	2.45	28	1
1:A:102:HIS:CG	1:A:106:ARG:NH1	0.64	2.65	39	2
1:A:106:ARG:N	1:A:106:ARG:NE	0.64	2.45	39	1
1:A:118:GLN:NE2	1:A:120:TYR:CD2	0.64	2.65	19	10
1:A:134:ARG:NH1	1:A:145:LEU:H	0.64	1.90	12	1
1:A:105:ARG:N	1:A:106:ARG:NH2	0.64	2.46	39	1
1:A:112:LEU:HD13	1:A:112:LEU:C	0.63	2.18	12	2
1:A:192:HIS:CG	1:A:193:HIS:N	0.63	2.64	37	5
1:A:102:HIS:N	1:A:181:LEU:O	0.63	2.31	39	37
1:A:113:TRP:CE2	1:A:117:THR:HG21	0.63	2.29	19	1
1:A:135:TYR:CE2	1:A:144:LYS:NZ	0.62	2.68	11	1
1:A:143:TYR:CZ	1:A:183:TRP:CD1	0.62	2.87	18	2
1:A:141:TRP:CE2	1:A:143:TYR:CD1	0.62	2.87	4	3
1:A:97:LEU:HD11	1:A:183:TRP:CE3	0.62	2.30	10	14
1:A:143:TYR:CD1	1:A:143:TYR:O	0.62	2.52	40	5
1:A:191:PHE:CD2	1:A:192:HIS:N	0.61	2.69	28	2
1:A:142:CYS:SG	1:A:190:ALA:N	0.61	2.74	32	5
1:A:141:TRP:CZ2	1:A:143:TYR:CD2	0.61	2.88	26	5
1:A:92:LYS:NZ	1:A:143:TYR:OH	0.61	2.34	23	1
1:A:186:ASP:O	1:A:188:ARG:N	0.60	2.34	3	39
1:A:126:ASN:N	1:A:126:ASN:HD22	0.60	1.94	36	2
1:A:106:ARG:NE	1:A:107:GLN:N	0.60	2.47	39	1
1:A:125:GLN:HE21	1:A:125:GLN:N	0.60	1.95	36	1
1:A:89:HIS:N	1:A:89:HIS:HD1	0.59	1.95	1	4
1:A:127:TYR:CE1	1:A:143:TYR:CZ	0.59	2.91	40	1
1:A:100:LEU:C	1:A:100:LEU:HD13	0.59	2.22	27	5
1:A:125:GLN:N	1:A:125:GLN:NE2	0.59	2.50	36	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:126:ASN:C	1:A:127:TYR:CG	0.59	2.80	25	26
1:A:79:MET:SD	1:A:124:TRP:CG	0.59	2.96	2	2
1:A:192:HIS:NE2	1:A:194:MET:SD	0.59	2.76	1	2
1:A:194:MET:SD	1:A:194:MET:N	0.59	2.76	31	3
1:A:128:THR:OG1	1:A:134:ARG:NE	0.58	2.36	21	2
1:A:98:GLU:O	1:A:100:LEU:N	0.58	2.36	30	24
1:A:118:GLN:NE2	1:A:118:GLN:O	0.58	2.36	24	8
1:A:106:ARG:H	1:A:106:ARG:CD	0.58	2.11	28	1
1:A:143:TYR:CE2	1:A:183:TRP:NE1	0.58	2.71	18	2
1:A:79:MET:SD	1:A:124:TRP:CD1	0.58	2.96	16	5
1:A:118:GLN:H	1:A:118:GLN:CD	0.58	2.07	4	9
1:A:126:ASN:C	1:A:127:TYR:CD1	0.58	2.82	13	22
1:A:112:LEU:C	1:A:112:LEU:HD23	0.57	2.24	37	9
1:A:80:THR:OG1	1:A:83:ALA:N	0.57	2.37	8	33
1:A:79:MET:SD	1:A:120:TYR:CD1	0.57	2.98	17	3
1:A:102:HIS:CE1	1:A:106:ARG:HH12	0.57	2.16	39	2
1:A:97:LEU:HD21	1:A:183:TRP:CD1	0.57	2.34	18	3
1:A:80:THR:O	1:A:83:ALA:N	0.57	2.38	40	39
1:A:106:ARG:HE	1:A:106:ARG:CA	0.57	2.13	28	3
1:A:125:GLN:H	1:A:125:GLN:CD	0.57	2.08	6	2
1:A:81:TYR:CD1	1:A:82:LYS:N	0.56	2.73	23	1
1:A:89:HIS:N	1:A:89:HIS:ND1	0.56	2.49	17	4
1:A:126:ASN:O	1:A:127:TYR:CG	0.56	2.58	19	24
1:A:126:ASN:O	1:A:127:TYR:CD2	0.56	2.59	29	9
1:A:124:TRP:CG	1:A:125:GLN:N	0.56	2.73	26	1
1:A:106:ARG:NE	1:A:107:GLN:H	0.56	1.98	39	2
1:A:116:HIS:ND1	1:A:116:HIS:C	0.56	2.63	14	2
1:A:138:THR:OG1	1:A:139:PHE:N	0.56	2.39	10	3
1:A:116:HIS:ND1	1:A:116:HIS:N	0.56	2.52	4	3
1:A:192:HIS:CD2	1:A:192:HIS:C	0.56	2.84	37	2
1:A:102:HIS:CE1	1:A:106:ARG:HH22	0.56	2.19	28	1
1:A:102:HIS:CD2	1:A:106:ARG:NH1	0.56	2.74	39	1
1:A:91:LEU:O	1:A:96:GLY:N	0.55	2.39	30	8
1:A:118:GLN:NE2	1:A:120:TYR:CE2	0.55	2.75	20	4
1:A:109:ILE:HD12	1:A:109:ILE:N	0.55	2.16	40	4
1:A:121:PHE:N	1:A:121:PHE:CD1	0.55	2.73	24	2
1:A:103:SER:OG	1:A:104:GLN:N	0.55	2.40	30	5
1:A:135:TYR:CD2	1:A:144:LYS:NZ	0.55	2.74	11	1
1:A:105:ARG:H	1:A:106:ARG:NH2	0.55	2.00	39	1
1:A:102:HIS:CE1	1:A:107:GLN:H	0.55	2.19	2	4
1:A:106:ARG:HH21	1:A:107:GLN:CB	0.55	2.15	28	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:139:PHE:CZ	1:A:189:LEU:O	0.55	2.60	35	2
1:A:141:TRP:CZ3	1:A:143:TYR:CE1	0.55	2.94	4	3
1:A:74:VAL:HG22	1:A:118:GLN:C	0.55	2.27	14	2
1:A:139:PHE:CE2	1:A:189:LEU:O	0.55	2.60	35	1
1:A:98:GLU:CD	1:A:99:GLY:N	0.55	2.65	18	8
1:A:193:HIS:ND1	1:A:193:HIS:N	0.55	2.51	3	5
1:A:145:LEU:HD23	1:A:182:GLU:C	0.55	2.27	33	1
1:A:118:GLN:NE2	1:A:120:TYR:CZ	0.55	2.75	6	3
1:A:79:MET:HE1	1:A:124:TRP:CE2	0.55	2.36	39	1
1:A:118:GLN:H	1:A:118:GLN:NE2	0.54	2.00	17	2
1:A:118:GLN:NE2	1:A:120:TYR:CE1	0.54	2.75	6	1
1:A:182:GLU:N	1:A:182:GLU:OE1	0.54	2.41	9	2
1:A:105:ARG:H	1:A:105:ARG:CD	0.54	2.16	26	1
1:A:188:ARG:O	1:A:191:PHE:CD1	0.54	2.60	28	1
1:A:88:SER:OG	1:A:89:HIS:ND1	0.54	2.40	27	18
1:A:118:GLN:NE2	1:A:118:GLN:C	0.54	2.66	13	7
1:A:113:TRP:O	1:A:117:THR:HG23	0.54	2.03	13	8
1:A:184:ARG:NH1	1:A:184:ARG:CG	0.54	2.71	13	11
1:A:106:ARG:NH1	1:A:106:ARG:CG	0.54	2.71	33	12
1:A:125:GLN:C	1:A:126:ASN:ND2	0.54	2.66	30	11
1:A:134:ARG:NH1	1:A:134:ARG:CG	0.54	2.71	29	4
1:A:79:MET:SD	1:A:124:TRP:CD2	0.54	3.01	26	1
1:A:139:PHE:O	1:A:141:TRP:N	0.54	2.41	2	2
1:A:126:ASN:O	1:A:137:LEU:HD12	0.54	2.02	12	3
1:A:143:TYR:CZ	1:A:183:TRP:NE1	0.54	2.75	18	1
1:A:102:HIS:CD2	1:A:106:ARG:NH2	0.54	2.76	28	1
1:A:88:SER:CB	1:A:142:CYS:SG	0.54	2.96	13	1
1:A:141:TRP:CH2	1:A:143:TYR:CD2	0.54	2.95	26	2
1:A:97:LEU:HD11	1:A:183:TRP:CZ3	0.54	2.38	39	8
1:A:105:ARG:CG	1:A:105:ARG:HH11	0.54	2.16	14	11
1:A:184:ARG:CG	1:A:184:ARG:HH11	0.54	2.16	13	10
1:A:191:PHE:CG	1:A:192:HIS:N	0.54	2.75	32	6
1:A:185:PHE:CD1	1:A:185:PHE:C	0.54	2.86	3	24
1:A:102:HIS:ND1	1:A:102:HIS:C	0.54	2.66	33	9
1:A:102:HIS:O	1:A:102:HIS:ND1	0.53	2.41	35	8
1:A:101:ILE:O	1:A:106:ARG:NH1	0.53	2.41	36	3
1:A:188:ARG:O	1:A:191:PHE:CD2	0.53	2.61	7	3
1:A:186:ASP:C	1:A:188:ARG:N	0.53	2.67	28	39
1:A:193:HIS:ND1	1:A:193:HIS:O	0.53	2.41	7	13
1:A:107:GLN:O	1:A:107:GLN:NE2	0.53	2.42	2	3
1:A:77:ARG:NH1	1:A:77:ARG:CG	0.53	2.72	8	5

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:106:ARG:CG	1:A:106:ARG:HH11	0.53	2.17	34	14
1:A:105:ARG:CG	1:A:105:ARG:NH1	0.53	2.71	36	5
1:A:134:ARG:CG	1:A:134:ARG:HH11	0.53	2.16	29	4
1:A:127:TYR:HH	1:A:183:TRP:CD1	0.53	2.22	26	1
1:A:106:ARG:NE	1:A:106:ARG:H	0.53	2.02	28	1
1:A:89:HIS:CD2	1:A:187:SER:HG	0.53	2.21	33	1
1:A:135:TYR:CD1	1:A:135:TYR:N	0.53	2.75	35	10
1:A:95:GLY:O	1:A:97:LEU:N	0.53	2.42	13	3
1:A:105:ARG:NH1	1:A:105:ARG:CG	0.53	2.71	24	8
1:A:143:TYR:CD1	1:A:143:TYR:C	0.53	2.85	25	5
1:A:89:HIS:NE2	1:A:187:SER:OG	0.53	2.41	33	2
1:A:186:ASP:OD1	1:A:188:ARG:NE	0.53	2.42	1	2
1:A:141:TRP:CH2	1:A:143:TYR:CE1	0.53	2.97	4	2
1:A:108:ASP:OD1	1:A:108:ASP:N	0.53	2.42	13	5
1:A:94:LYS:O	1:A:96:GLY:N	0.53	2.42	6	3
1:A:188:ARG:O	1:A:192:HIS:CE1	0.53	2.62	4	1
1:A:125:GLN:O	1:A:126:ASN:ND2	0.52	2.42	21	11
1:A:127:TYR:CB	1:A:134:ARG:HE	0.52	2.16	12	1
1:A:105:ARG:H	1:A:106:ARG:CZ	0.52	2.18	39	2
1:A:106:ARG:HH21	1:A:107:GLN:H	0.52	1.45	28	1
1:A:134:ARG:NH1	1:A:134:ARG:O	0.52	2.42	12	1
1:A:182:GLU:CD	1:A:184:ARG:NH2	0.52	2.67	31	1
1:A:125:GLN:CD	1:A:125:GLN:N	0.52	2.67	20	1
1:A:193:HIS:CD2	1:A:193:HIS:O	0.52	2.62	2	2
1:A:125:GLN:HE21	1:A:125:GLN:H	0.52	1.47	36	1
1:A:188:ARG:HH11	1:A:188:ARG:CG	0.52	2.18	16	1
1:A:125:GLN:NE2	1:A:127:TYR:OH	0.52	2.42	27	1
1:A:192:HIS:CD2	1:A:193:HIS:N	0.52	2.78	37	1
1:A:126:ASN:O	1:A:127:TYR:CD1	0.52	2.62	24	3
1:A:118:GLN:NE2	1:A:120:TYR:CD1	0.52	2.78	6	4
1:A:106:ARG:CG	1:A:106:ARG:NH1	0.52	2.71	40	3
1:A:192:HIS:CG	1:A:193:HIS:H	0.52	2.22	13	2
1:A:118:GLN:C	1:A:118:GLN:HE21	0.52	2.13	13	4
1:A:134:ARG:HH11	1:A:134:ARG:CG	0.52	2.17	39	1
1:A:117:THR:OG1	1:A:118:GLN:NE2	0.52	2.43	4	3
1:A:118:GLN:CD	1:A:120:TYR:CZ	0.52	2.87	36	4
1:A:118:GLN:CD	1:A:120:TYR:CE2	0.52	2.87	13	6
1:A:123:ASP:N	1:A:123:ASP:OD1	0.52	2.42	8	6
1:A:126:ASN:C	1:A:127:TYR:CD2	0.52	2.88	23	1
1:A:80:THR:OG1	1:A:83:ALA:HB2	0.52	2.04	33	5
1:A:188:ARG:O	1:A:191:PHE:CE1	0.52	2.62	28	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:134:ARG:CG	1:A:134:ARG:NH1	0.52	2.71	39	1
1:A:191:PHE:CD1	1:A:192:HIS:N	0.52	2.78	13	7
1:A:115:TYR:C	1:A:117:THR:H	0.51	2.13	6	25
1:A:114:ILE:O	1:A:117:THR:OG1	0.51	2.29	38	15
1:A:97:LEU:O	1:A:98:GLU:O	0.51	2.27	30	1
1:A:191:PHE:C	1:A:191:PHE:CD1	0.51	2.88	26	1
1:A:102:HIS:C	1:A:106:ARG:NH1	0.51	2.68	28	2
1:A:103:SER:C	1:A:106:ARG:NH1	0.51	2.69	39	2
1:A:186:ASP:C	1:A:188:ARG:H	0.51	2.14	21	33
1:A:89:HIS:N	1:A:89:HIS:CD2	0.51	2.77	23	1
1:A:182:GLU:OE2	1:A:184:ARG:NH2	0.51	2.44	31	1
1:A:111:ASP:N	1:A:111:ASP:OD1	0.51	2.42	30	3
1:A:111:ASP:OD1	1:A:125:GLN:NE2	0.51	2.44	18	1
1:A:126:ASN:N	1:A:126:ASN:ND2	0.51	2.58	36	2
1:A:118:GLN:O	1:A:118:GLN:CD	0.51	2.54	9	5
1:A:188:ARG:C	1:A:190:ALA:H	0.50	2.14	28	33
1:A:188:ARG:CG	1:A:188:ARG:NH1	0.50	2.72	16	5
1:A:188:ARG:O	1:A:191:PHE:CE2	0.50	2.64	17	2
1:A:121:PHE:CD2	1:A:123:ASP:OD1	0.50	2.64	14	1
1:A:124:TRP:CD1	1:A:124:TRP:H	0.50	2.25	32	3
1:A:98:GLU:C	1:A:100:LEU:H	0.50	2.14	30	22
1:A:142:CYS:SG	1:A:142:CYS:O	0.50	2.70	33	2
1:A:88:SER:HG	1:A:89:HIS:CE1	0.50	2.24	17	2
1:A:99:GLY:N	1:A:183:TRP:O	0.50	2.41	29	3
1:A:79:MET:SD	1:A:79:MET:C	0.50	2.95	33	2
1:A:186:ASP:OD1	1:A:188:ARG:NH2	0.50	2.44	36	1
1:A:93:GLU:OE1	1:A:93:GLU:N	0.50	2.44	38	1
1:A:98:GLU:C	1:A:100:LEU:N	0.50	2.70	38	28
1:A:123:ASP:N	1:A:125:GLN:OE1	0.50	2.45	6	1
1:A:186:ASP:OD2	1:A:188:ARG:NE	0.50	2.45	12	1
1:A:88:SER:OG	1:A:89:HIS:CE1	0.50	2.65	24	2
1:A:194:MET:C	1:A:194:MET:SD	0.50	2.95	34	1
1:A:125:GLN:N	1:A:125:GLN:OE1	0.50	2.45	37	4
1:A:143:TYR:CE2	1:A:183:TRP:CD1	0.50	2.99	18	2
1:A:127:TYR:CD1	1:A:127:TYR:N	0.50	2.79	35	6
1:A:125:GLN:CD	1:A:125:GLN:H	0.50	2.13	20	1
1:A:89:HIS:ND1	1:A:89:HIS:N	0.49	2.59	27	4
1:A:118:GLN:CD	1:A:118:GLN:N	0.49	2.70	2	8
1:A:77:ARG:CG	1:A:77:ARG:HH11	0.49	2.19	8	5
1:A:91:LEU:O	1:A:95:GLY:CA	0.49	2.61	14	7
1:A:141:TRP:CD2	1:A:143:TYR:CD1	0.49	3.00	4	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:115:TYR:O	1:A:117:THR:N	0.49	2.44	6	4
1:A:88:SER:OG	1:A:89:HIS:N	0.49	2.45	31	9
1:A:107:GLN:CG	1:A:108:ASP:N	0.49	2.75	39	4
1:A:128:THR:C	1:A:134:ARG:NH1	0.49	2.71	13	1
1:A:102:HIS:ND1	1:A:102:HIS:O	0.49	2.43	36	1
1:A:143:TYR:CE1	1:A:183:TRP:CD1	0.48	3.02	18	1
1:A:127:TYR:CG	1:A:143:TYR:OH	0.48	2.61	31	1
1:A:188:ARG:CG	1:A:188:ARG:HH11	0.48	2.21	31	4
1:A:188:ARG:C	1:A:190:ALA:N	0.48	2.71	19	35
1:A:112:LEU:C	1:A:112:LEU:CD1	0.48	2.87	12	2
1:A:185:PHE:CZ	1:A:187:SER:OG	0.48	2.65	14	1
1:A:98:GLU:O	1:A:183:TRP:O	0.48	2.32	30	1
1:A:74:VAL:O	1:A:75:PRO:C	0.48	2.56	27	31
1:A:114:ILE:HG23	1:A:118:GLN:HE22	0.48	1.68	27	3
1:A:125:GLN:NE2	1:A:125:GLN:O	0.48	2.46	20	2
1:A:142:CYS:O	1:A:142:CYS:SG	0.48	2.71	20	1
1:A:112:LEU:O	1:A:115:TYR:N	0.48	2.46	22	6
1:A:184:ARG:CG	1:A:184:ARG:NH1	0.48	2.72	12	1
1:A:186:ASP:CG	1:A:188:ARG:HE	0.48	2.17	22	1
1:A:102:HIS:HE2	1:A:107:GLN:HB3	0.48	1.67	18	1
1:A:102:HIS:CB	1:A:181:LEU:O	0.48	2.62	28	2
1:A:193:HIS:CG	1:A:193:HIS:O	0.48	2.64	1	1
1:A:124:TRP:CG	1:A:125:GLN:H	0.47	2.27	26	1
1:A:127:TYR:OH	1:A:183:TRP:NE1	0.47	2.47	26	1
1:A:127:TYR:CE1	1:A:143:TYR:OH	0.47	2.62	40	1
1:A:94:LYS:CD	1:A:94:LYS:N	0.47	2.77	28	2
1:A:125:GLN:C	1:A:126:ASN:HD22	0.47	2.18	30	2
1:A:117:THR:OG1	1:A:118:GLN:N	0.47	2.48	7	2
1:A:88:SER:OG	1:A:89:HIS:CD2	0.47	2.68	23	1
1:A:74:VAL:CG2	1:A:118:GLN:O	0.47	2.62	24	3
1:A:192:HIS:H	1:A:192:HIS:CD2	0.47	2.27	22	1
1:A:80:THR:OG1	1:A:83:ALA:CB	0.47	2.63	38	5
1:A:128:THR:C	1:A:134:ARG:HH11	0.47	2.18	13	1
1:A:89:HIS:CE1	1:A:187:SER:HG	0.47	2.28	32	1
1:A:139:PHE:C	1:A:141:TRP:N	0.47	2.73	2	1
1:A:92:LYS:CD	1:A:92:LYS:C	0.47	2.88	10	2
1:A:139:PHE:CD1	1:A:140:GLY:N	0.47	2.83	1	3
1:A:145:LEU:O	1:A:146:VAL:HG23	0.47	2.10	15	8
1:A:193:HIS:CD2	1:A:193:HIS:N	0.47	2.82	39	3
1:A:191:PHE:O	1:A:193:HIS:CE1	0.47	2.67	12	1
1:A:79:MET:SD	1:A:120:TYR:CG	0.47	3.08	11	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:100:LEU:C	1:A:100:LEU:CD1	0.47	2.88	27	4
1:A:124:TRP:CE2	1:A:125:GLN:NE2	0.47	2.83	25	1
1:A:139:PHE:CG	1:A:140:GLY:N	0.46	2.83	35	3
1:A:91:LEU:O	1:A:95:GLY:O	0.46	2.34	9	11
1:A:191:PHE:O	1:A:192:HIS:CG	0.46	2.68	27	3
1:A:118:GLN:OE1	1:A:120:TYR:CE1	0.46	2.68	36	1
1:A:98:GLU:C	1:A:98:GLU:CD	0.46	2.84	14	8
1:A:189:LEU:HD23	1:A:189:LEU:O	0.46	2.10	9	7
1:A:93:GLU:N	1:A:93:GLU:CD	0.46	2.73	13	2
1:A:128:THR:OG1	1:A:134:ARG:CZ	0.46	2.63	13	1
1:A:106:ARG:HE	1:A:106:ARG:C	0.46	2.19	39	1
1:A:112:LEU:C	1:A:112:LEU:CD2	0.46	2.89	37	5
1:A:89:HIS:CE1	1:A:187:SER:OG	0.46	2.69	32	2
1:A:91:LEU:O	1:A:95:GLY:N	0.46	2.49	25	3
1:A:134:ARG:HH12	1:A:145:LEU:H	0.46	1.54	12	1
1:A:109:ILE:N	1:A:109:ILE:CD1	0.46	2.78	40	3
1:A:125:GLN:O	1:A:125:GLN:CD	0.46	2.59	37	7
1:A:140:GLY:O	1:A:142:CYS:SG	0.46	2.74	5	2
1:A:141:TRP:CH2	1:A:143:TYR:CE2	0.46	3.03	26	1
1:A:74:VAL:O	1:A:75:PRO:O	0.46	2.34	38	2
1:A:134:ARG:NH1	1:A:181:LEU:CD1	0.46	2.79	30	1
1:A:191:PHE:O	1:A:192:HIS:ND1	0.46	2.49	31	1
1:A:127:TYR:CZ	1:A:143:TYR:OH	0.45	2.62	40	1
1:A:80:THR:O	1:A:83:ALA:HB3	0.45	2.11	39	14
1:A:191:PHE:CD1	1:A:191:PHE:C	0.45	2.94	17	4
1:A:102:HIS:HE2	1:A:107:GLN:CB	0.45	2.25	18	1
1:A:111:ASP:OD2	1:A:125:GLN:NE2	0.45	2.50	35	3
1:A:79:MET:SD	1:A:120:TYR:CE1	0.45	3.09	24	1
1:A:89:HIS:NE2	1:A:187:SER:CB	0.45	2.79	31	1
1:A:92:LYS:O	1:A:95:GLY:O	0.45	2.34	35	4
1:A:124:TRP:CD2	1:A:125:GLN:N	0.45	2.85	26	1
1:A:121:PHE:O	1:A:121:PHE:CD2	0.45	2.69	24	1
1:A:186:ASP:O	1:A:187:SER:C	0.45	2.59	22	10
1:A:80:THR:O	1:A:81:TYR:C	0.45	2.60	38	38
1:A:141:TRP:CD1	1:A:143:TYR:CD1	0.45	3.04	31	2
1:A:108:ASP:O	1:A:109:ILE:C	0.45	2.60	9	9
1:A:112:LEU:HD23	1:A:112:LEU:O	0.45	2.12	37	3
1:A:146:VAL:O	1:A:147:PRO:O	0.45	2.35	30	3
1:A:102:HIS:NE2	1:A:107:GLN:OE1	0.45	2.44	23	1
1:A:193:HIS:O	1:A:193:HIS:CG	0.45	2.68	2	1
1:A:128:THR:O	1:A:134:ARG:NH1	0.44	2.40	13	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:112:LEU:HD13	1:A:112:LEU:O	0.44	2.12	12	2
1:A:98:GLU:C	1:A:98:GLU:OE1	0.44	2.61	10	1
1:A:188:ARG:N	1:A:188:ARG:CD	0.44	2.80	17	1
1:A:98:GLU:C	1:A:183:TRP:O	0.44	2.60	30	1
1:A:82:LYS:NZ	1:A:86:ASP:OD2	0.44	2.46	39	1
1:A:114:ILE:HG23	1:A:118:GLN:NE2	0.44	2.28	27	1
1:A:98:GLU:CD	1:A:98:GLU:C	0.44	2.86	4	5
1:A:104:GLN:N	1:A:106:ARG:NH1	0.44	2.66	28	1
1:A:127:TYR:CE2	1:A:143:TYR:OH	0.44	2.69	40	1
1:A:85:VAL:O	1:A:88:SER:OG	0.44	2.36	17	7
1:A:122:PRO:O	1:A:125:GLN:CD	0.44	2.60	37	1
1:A:125:GLN:OE1	1:A:125:GLN:C	0.44	2.61	2	1
1:A:102:HIS:NE2	1:A:107:GLN:CB	0.44	2.80	18	1
1:A:189:LEU:HD13	1:A:189:LEU:O	0.44	2.12	26	3
1:A:125:GLN:N	1:A:125:GLN:CD	0.44	2.76	22	2
1:A:142:CYS:SG	1:A:189:LEU:C	0.44	3.01	32	1
1:A:125:GLN:NE2	1:A:126:ASN:ND2	0.43	2.65	36	1
1:A:115:TYR:C	1:A:117:THR:N	0.43	2.76	6	12
1:A:182:GLU:CD	1:A:182:GLU:C	0.43	2.85	14	2
1:A:124:TRP:O	1:A:126:ASN:N	0.43	2.50	26	1
1:A:107:GLN:CD	1:A:107:GLN:C	0.43	2.86	2	4
1:A:111:ASP:CG	1:A:125:GLN:OE1	0.43	2.62	19	2
1:A:123:ASP:OD1	1:A:123:ASP:C	0.43	2.61	34	6
1:A:125:GLN:O	1:A:126:ASN:CG	0.43	2.61	40	12
1:A:94:LYS:O	1:A:95:GLY:C	0.43	2.61	33	9
1:A:112:LEU:O	1:A:113:TRP:C	0.43	2.60	37	17
1:A:121:PHE:O	1:A:123:ASP:CG	0.43	2.62	6	2
1:A:97:LEU:CD1	1:A:183:TRP:CE3	0.43	3.01	9	1
1:A:90:PHE:CE2	1:A:94:LYS:NZ	0.43	2.83	17	1
1:A:145:LEU:HD23	1:A:182:GLU:O	0.43	2.13	33	1
1:A:98:GLU:OE2	1:A:99:GLY:N	0.43	2.51	1	2
1:A:80:THR:OG1	1:A:120:TYR:OH	0.43	2.36	5	1
1:A:125:GLN:CD	1:A:125:GLN:C	0.43	2.86	12	2
1:A:121:PHE:O	1:A:123:ASP:OD1	0.43	2.37	31	3
1:A:97:LEU:O	1:A:98:GLU:C	0.43	2.61	15	3
1:A:146:VAL:O	1:A:182:GLU:OE1	0.43	2.36	9	1
1:A:91:LEU:O	1:A:95:GLY:C	0.43	2.62	13	3
1:A:118:GLN:OE1	1:A:120:TYR:CZ	0.43	2.72	24	2
1:A:98:GLU:OE2	1:A:184:ARG:NH2	0.43	2.51	19	1
1:A:107:GLN:OE1	1:A:125:GLN:OE1	0.43	2.36	30	1
1:A:107:GLN:NE2	1:A:111:ASP:OD2	0.43	2.42	31	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:139:PHE:CD1	1:A:139:PHE:C	0.43	2.97	35	1
1:A:125:GLN:NE2	1:A:125:GLN:H	0.43	2.10	36	1
1:A:106:ARG:CD	1:A:106:ARG:N	0.43	2.81	12	1
1:A:116:HIS:CD2	1:A:116:HIS:C	0.43	2.96	24	1
1:A:125:GLN:O	1:A:126:ASN:OD1	0.43	2.36	34	2
1:A:139:PHE:O	1:A:140:GLY:C	0.43	2.60	14	4
1:A:192:HIS:CE1	1:A:194:MET:SD	0.43	3.12	2	2
1:A:104:GLN:CD	1:A:104:GLN:C	0.43	2.87	3	3
1:A:93:GLU:N	1:A:93:GLU:OE1	0.43	2.52	13	1
1:A:81:TYR:CD2	1:A:82:LYS:N	0.43	2.87	21	1
1:A:146:VAL:O	1:A:182:GLU:CD	0.43	2.62	23	1
1:A:142:CYS:SG	1:A:190:ALA:CA	0.42	3.07	25	1
1:A:110:LEU:O	1:A:111:ASP:C	0.42	2.62	38	4
1:A:81:TYR:CE1	1:A:139:PHE:CE2	0.42	3.07	30	2
1:A:89:HIS:O	1:A:93:GLU:CD	0.42	2.62	38	1
1:A:81:TYR:C	1:A:81:TYR:CD1	0.42	2.95	4	1
1:A:184:ARG:NH1	1:A:185:PHE:O	0.42	2.53	5	1
1:A:136:PRO:O	1:A:195:ALA:CB	0.42	2.67	40	1
1:A:107:GLN:NE2	1:A:107:GLN:C	0.42	2.78	2	1
1:A:125:GLN:O	1:A:125:GLN:CG	0.42	2.67	14	1
1:A:109:ILE:O	1:A:110:LEU:C	0.42	2.62	32	3
1:A:142:CYS:HG	1:A:190:ALA:N	0.42	2.13	34	1
1:A:94:LYS:CG	1:A:95:GLY:H	0.42	2.27	38	1
1:A:145:LEU:O	1:A:146:VAL:CG2	0.42	2.67	11	6
1:A:194:MET:O	1:A:195:ALA:C	0.42	2.61	9	8
1:A:139:PHE:CD2	1:A:140:GLY:N	0.42	2.88	5	1
1:A:143:TYR:CD1	1:A:144:LYS:N	0.42	2.87	18	2
1:A:127:TYR:CE1	1:A:143:TYR:CE1	0.42	3.08	40	1
1:A:76:LEU:CD2	1:A:119:GLY:O	0.42	2.67	3	2
1:A:102:HIS:C	1:A:102:HIS:HD1	0.42	2.23	26	1
1:A:127:TYR:CD1	1:A:143:TYR:OH	0.42	2.69	40	1
1:A:194:MET:SD	1:A:194:MET:C	0.42	3.03	17	1
1:A:126:ASN:ND2	1:A:126:ASN:N	0.42	2.66	30	1
1:A:144:LYS:HZ3	1:A:146:VAL:CG2	0.42	2.28	12	1
1:A:146:VAL:HG11	1:A:184:ARG:NH2	0.42	2.30	28	1
1:A:89:HIS:O	1:A:93:GLU:OE1	0.42	2.37	38	2
1:A:89:HIS:CE1	1:A:187:SER:CB	0.42	3.03	31	1
1:A:193:HIS:CE1	1:A:195:ALA:HB3	0.42	2.50	31	1
1:A:107:GLN:HE22	1:A:111:ASP:CG	0.41	2.23	25	1
1:A:111:ASP:OD1	1:A:125:GLN:OE1	0.41	2.38	19	1
1:A:97:LEU:HD21	1:A:183:TRP:CD2	0.41	2.49	30	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:109:ILE:O	1:A:113:TRP:CB	0.41	2.68	35	1
1:A:77:ARG:CB	1:A:78:PRO:CD	0.41	2.98	1	1
1:A:81:TYR:CE1	1:A:139:PHE:CD2	0.41	3.08	15	1
1:A:88:SER:HG	1:A:89:HIS:N	0.41	2.14	24	1
1:A:126:ASN:O	1:A:126:ASN:CG	0.41	2.63	26	1
1:A:125:GLN:O	1:A:125:GLN:OE1	0.41	2.37	2	1
1:A:77:ARG:O	1:A:78:PRO:C	0.41	2.62	5	2
1:A:182:GLU:CD	1:A:182:GLU:O	0.41	2.63	14	2
1:A:189:LEU:C	1:A:191:PHE:H	0.41	2.22	30	1
1:A:111:ASP:OD1	1:A:125:GLN:CD	0.41	2.63	34	1
1:A:102:HIS:CG	1:A:106:ARG:HH12	0.41	2.33	39	1
1:A:139:PHE:CE1	1:A:189:LEU:O	0.41	2.73	7	1
1:A:92:LYS:CE	1:A:143:TYR:OH	0.41	2.68	23	1
1:A:81:TYR:CD1	1:A:81:TYR:C	0.41	2.98	3	1
1:A:124:TRP:CD1	1:A:124:TRP:N	0.41	2.89	25	2
1:A:77:ARG:HH21	1:A:118:GLN:CG	0.41	2.28	30	1
1:A:125:GLN:C	1:A:126:ASN:CG	0.41	2.89	10	1
1:A:109:ILE:HD12	1:A:109:ILE:H	0.41	1.74	19	1
1:A:191:PHE:C	1:A:192:HIS:CG	0.41	2.99	27	1
1:A:106:ARG:NH2	1:A:107:GLN:H	0.41	2.11	28	1
1:A:105:ARG:HG2	1:A:105:ARG:HH11	0.41	1.75	32	1
1:A:123:ASP:OD1	1:A:123:ASP:N	0.41	2.51	12	1
1:A:98:GLU:OE1	1:A:185:PHE:N	0.41	2.38	18	1
1:A:112:LEU:HD23	1:A:112:LEU:C	0.41	2.41	20	1
1:A:134:ARG:HA	1:A:134:ARG:HE	0.41	1.76	21	1
1:A:122:PRO:C	1:A:124:TRP:H	0.41	2.24	30	1
1:A:111:ASP:OD2	1:A:125:GLN:OE1	0.41	2.38	34	1
1:A:92:LYS:NZ	1:A:98:GLU:OE2	0.41	2.53	36	1
1:A:182:GLU:C	1:A:182:GLU:OE1	0.40	2.64	14	1
1:A:192:HIS:ND1	1:A:192:HIS:N	0.40	2.68	14	1
1:A:116:HIS:O	1:A:116:HIS:ND1	0.40	2.53	18	1
1:A:127:TYR:CD2	1:A:143:TYR:OH	0.40	2.72	36	2
1:A:106:ARG:NE	1:A:106:ARG:CA	0.40	2.84	28	1
1:A:94:LYS:CB	1:A:94:LYS:NZ	0.40	2.84	7	1
1:A:87:LEU:HD13	1:A:118:GLN:OE1	0.40	2.17	16	1
1:A:108:ASP:N	1:A:108:ASP:OD1	0.40	2.51	18	1
1:A:114:ILE:O	1:A:118:GLN:NE2	0.40	2.54	29	1
1:A:189:LEU:O	1:A:189:LEU:HD23	0.40	2.17	31	1
1:A:102:HIS:CE1	1:A:106:ARG:NH1	0.40	2.89	39	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	84/136 (62%)	61±3 (72±3%)	18±3 (22±3%)	5±1 (6±2%)	2	21
All	All	3360/5440 (62%)	2435 (72%)	737 (22%)	188 (6%)	2	21

All 24 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	187	SER	39
1	A	104	GLN	22
1	A	116	HIS	20
1	A	103	SER	18
1	A	75	PRO	16
1	A	99	GLY	16
1	A	193	HIS	8
1	A	95	GLY	7
1	A	147	PRO	6
1	A	93	GLU	5
1	A	105	ARG	5
1	A	122	PRO	4
1	A	192	HIS	4
1	A	127	TYR	3
1	A	140	GLY	2
1	A	110	LEU	2
1	A	96	GLY	2
1	A	118	GLN	2
1	A	128	THR	2
1	A	101	ILE	1
1	A	76	LEU	1
1	A	78	PRO	1
1	A	98	GLU	1
1	A	92	LYS	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	75/120 (62%)	68±2 (91±3%)	7±2 (9±3%)	11	56
All	All	3000/4800 (62%)	2724 (91%)	276 (9%)	11	56

All 38 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	81	TYR	31
1	A	138	THR	29
1	A	189	LEU	18
1	A	118	GLN	17
1	A	128	THR	15
1	A	193	HIS	13
1	A	106	ARG	12
1	A	105	ARG	12
1	A	184	ARG	12
1	A	74	VAL	9
1	A	143	TYR	9
1	A	134	ARG	8
1	A	116	HIS	7
1	A	125	GLN	7
1	A	127	TYR	7
1	A	187	SER	6
1	A	92	LYS	6
1	A	89	HIS	5
1	A	94	LYS	5
1	A	194	MET	5
1	A	126	ASN	5
1	A	121	PHE	4
1	A	77	ARG	4
1	A	79	MET	4
1	A	142	CYS	3
1	A	100	LEU	3
1	A	93	GLU	3
1	A	144	LYS	3
1	A	192	HIS	2

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Mol	Chain	Res	Type	Models (Total)
1	A	112	LEU	2
1	A	102	HIS	2
1	A	188	ARG	2
1	A	182	GLU	1
1	A	108	ASP	1
1	A	110	LEU	1
1	A	82	LYS	1
1	A	135	TYR	1
1	A	185	PHE	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