



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 06:09 AM UTC

PDB ID : 1TKV / pdb_00001tkv
Title : Solution Structure of T4 AsiA Dimer
Authors : Lambert, L.J.; Schirf, V.; Demeler, B.; Cadene, M.; Werner, M.H.
Deposited on : 2004-06-09

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

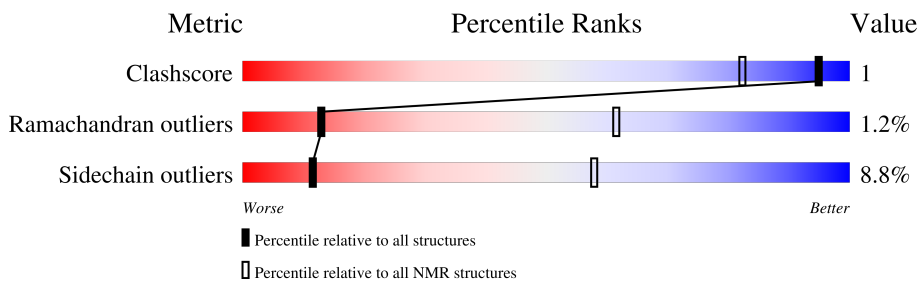
1 Overall quality at a glance i

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	90	 72% 21% . . .
1	B	90	 76% 18% . . .

2 Ensemble composition and analysis

This entry contains 23 models. Model 13 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:4-A:88, B:4-B:88 (170)	0.52	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 3 single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2948 atoms, of which 1476 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called 10 kDa anti-sigma factor.

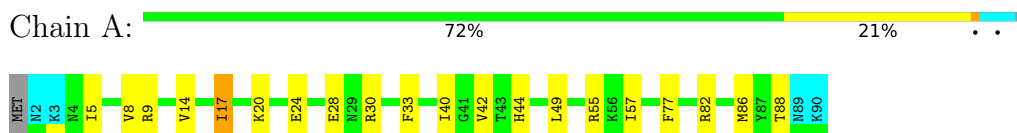
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	89	1474	466	738	127	142	1	0
1	B	89	1474	466	738	127	142	1	0

4 Residue-property plots

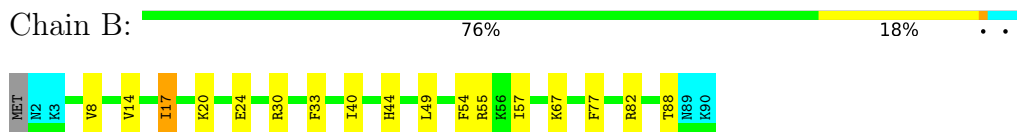
4.1 Average score per residue in the NMR ensemble

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: 10 kDa anti-sigma factor



- Molecule 1: 10 kDa anti-sigma factor

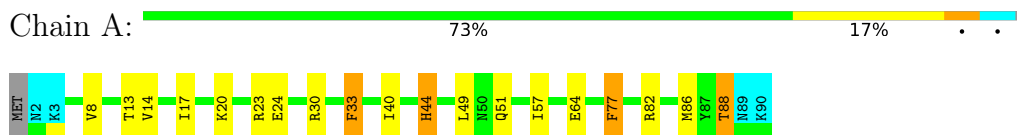


4.2 Scores per residue for each member of the ensemble

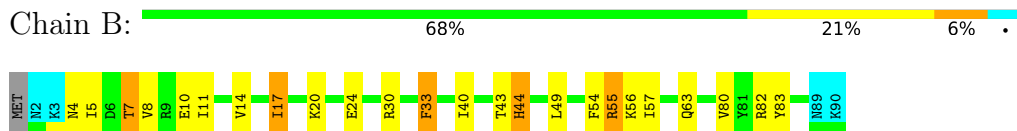
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: 10 kDa anti-sigma factor

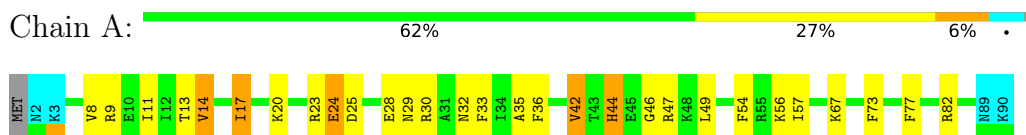


- Molecule 1: 10 kDa anti-sigma factor

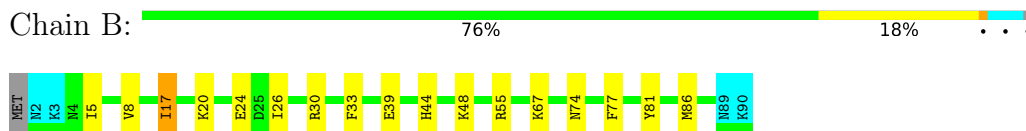


4.2.2 Score per residue for model 2

- Molecule 1: 10 kDa anti-sigma factor

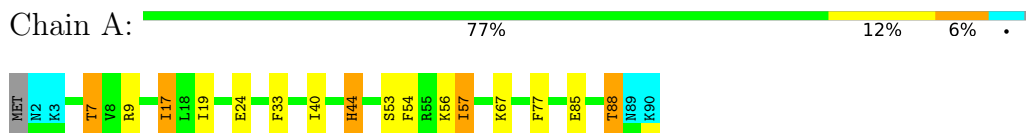


- Molecule 1: 10 kDa anti-sigma factor

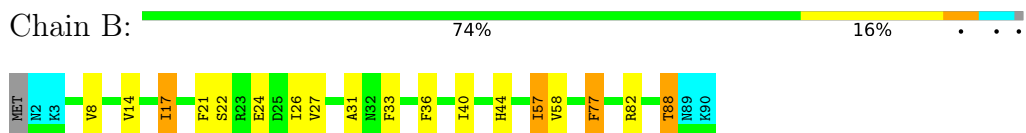


4.2.3 Score per residue for model 3

- Molecule 1: 10 kDa anti-sigma factor

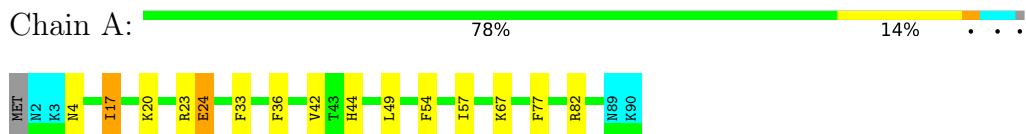


- Molecule 1: 10 kDa anti-sigma factor

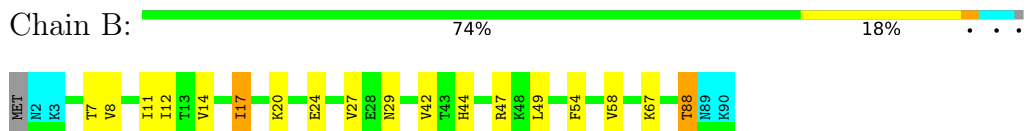


4.2.4 Score per residue for model 4

- Molecule 1: 10 kDa anti-sigma factor

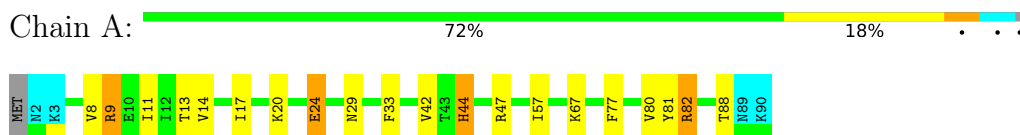


- Molecule 1: 10 kDa anti-sigma factor

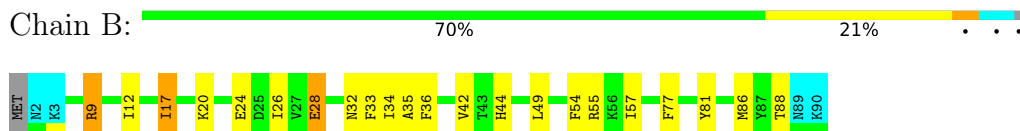


4.2.5 Score per residue for model 5

- Molecule 1: 10 kDa anti-sigma factor

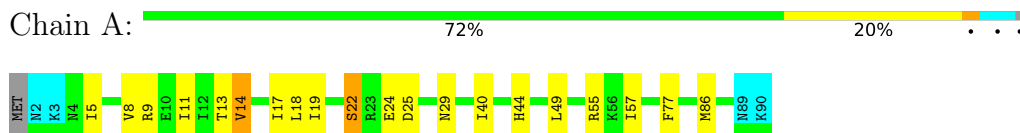


- Molecule 1: 10 kDa anti-sigma factor

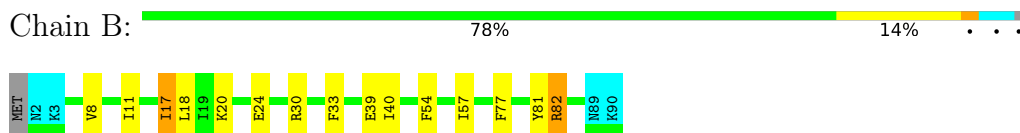


4.2.6 Score per residue for model 6

- Molecule 1: 10 kDa anti-sigma factor

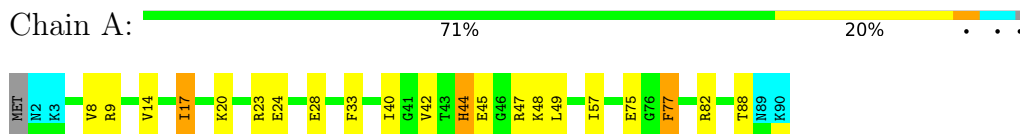


- Molecule 1: 10 kDa anti-sigma factor

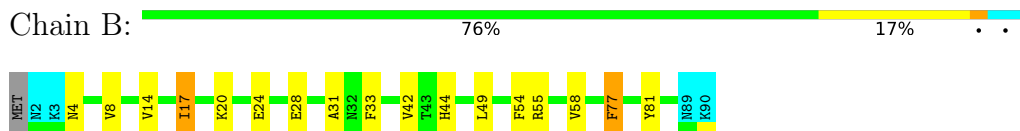


4.2.7 Score per residue for model 7

- Molecule 1: 10 kDa anti-sigma factor

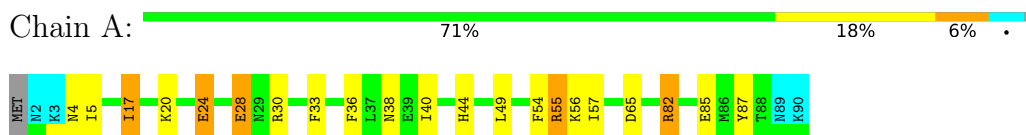


- Molecule 1: 10 kDa anti-sigma factor

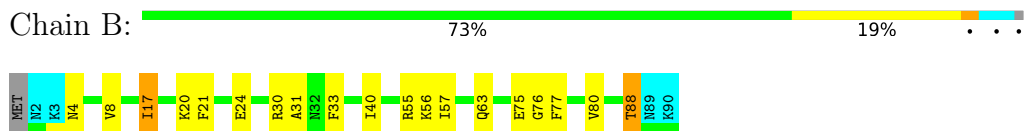


4.2.8 Score per residue for model 8

- Molecule 1: 10 kDa anti-sigma factor

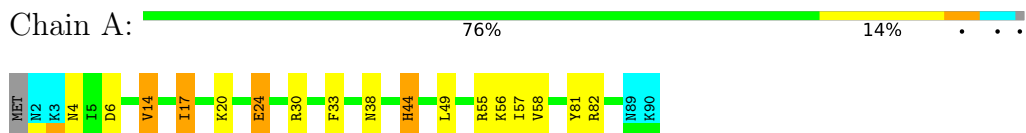


- Molecule 1: 10 kDa anti-sigma factor

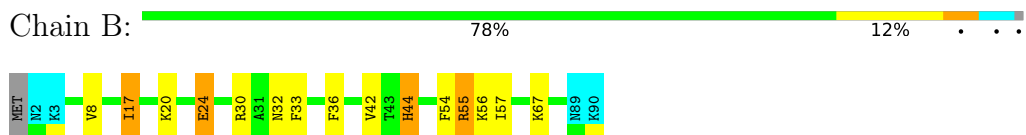


4.2.9 Score per residue for model 9

- Molecule 1: 10 kDa anti-sigma factor

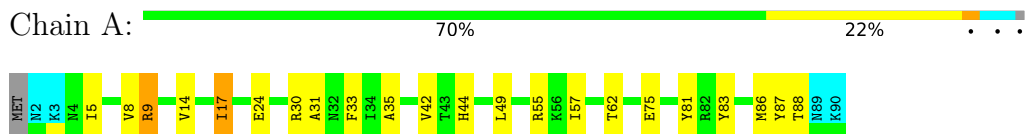


- Molecule 1: 10 kDa anti-sigma factor

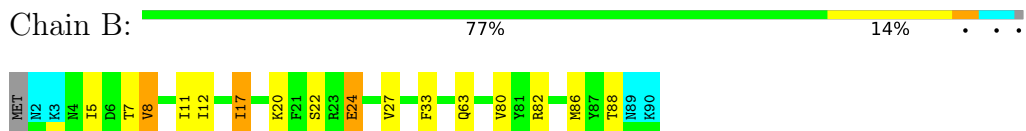


4.2.10 Score per residue for model 10

- Molecule 1: 10 kDa anti-sigma factor

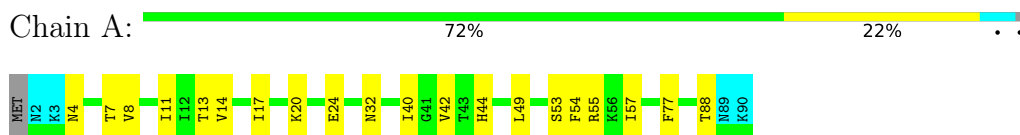


- Molecule 1: 10 kDa anti-sigma factor

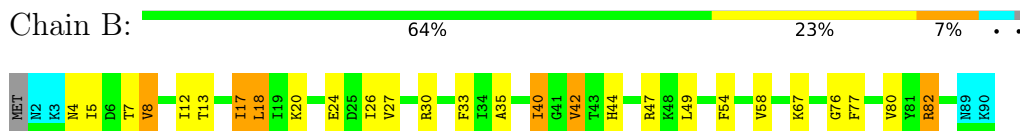


4.2.11 Score per residue for model 11

- Molecule 1: 10 kDa anti-sigma factor

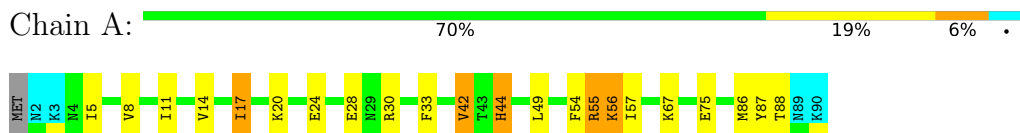


- Molecule 1: 10 kDa anti-sigma factor

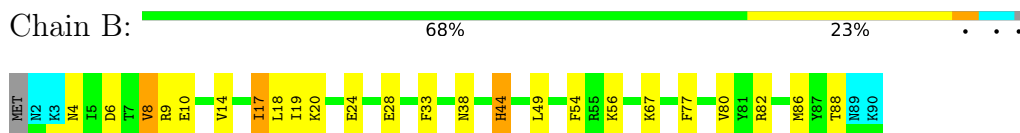


4.2.12 Score per residue for model 12

- Molecule 1: 10 kDa anti-sigma factor

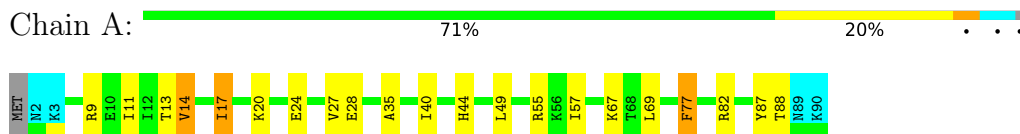


- Molecule 1: 10 kDa anti-sigma factor

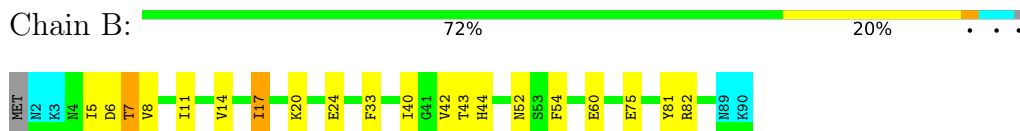


4.2.13 Score per residue for model 13 (medoid)

- Molecule 1: 10 kDa anti-sigma factor

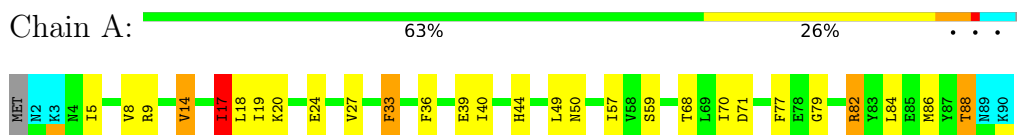


- Molecule 1: 10 kDa anti-sigma factor

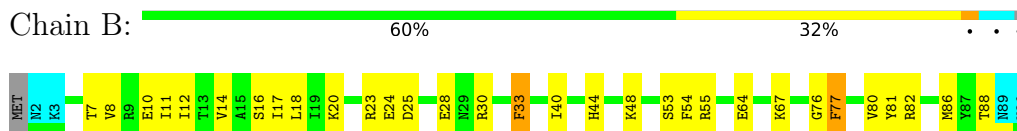


4.2.14 Score per residue for model 14

- Molecule 1: 10 kDa anti-sigma factor

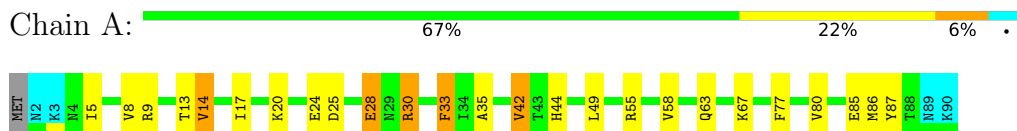


- Molecule 1: 10 kDa anti-sigma factor

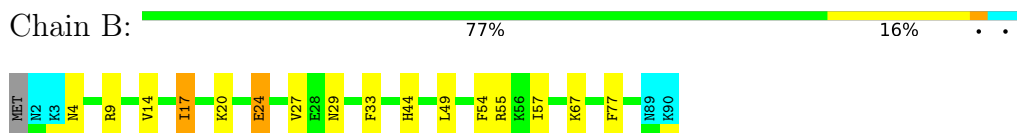


4.2.15 Score per residue for model 15

- Molecule 1: 10 kDa anti-sigma factor

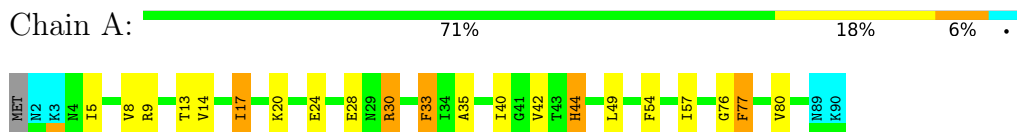


- Molecule 1: 10 kDa anti-sigma factor

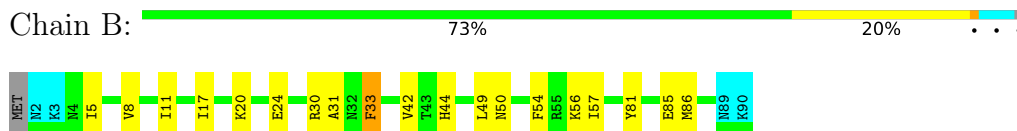


4.2.16 Score per residue for model 16

- Molecule 1: 10 kDa anti-sigma factor

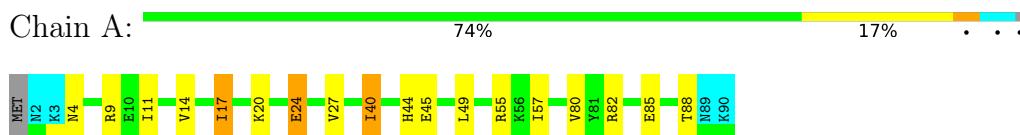


- Molecule 1: 10 kDa anti-sigma factor

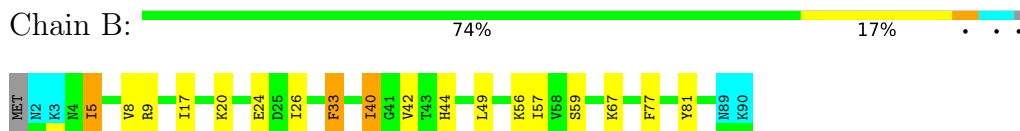


4.2.17 Score per residue for model 17

- Molecule 1: 10 kDa anti-sigma factor

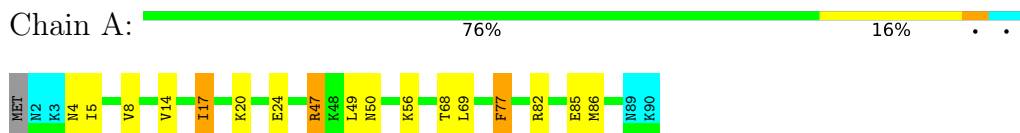


- Molecule 1: 10 kDa anti-sigma factor

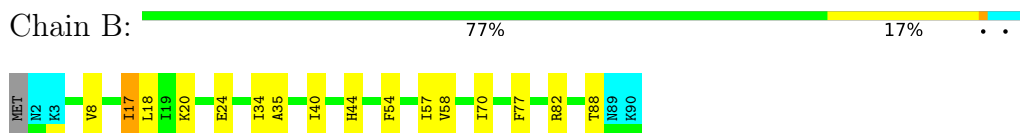


4.2.18 Score per residue for model 18

- Molecule 1: 10 kDa anti-sigma factor

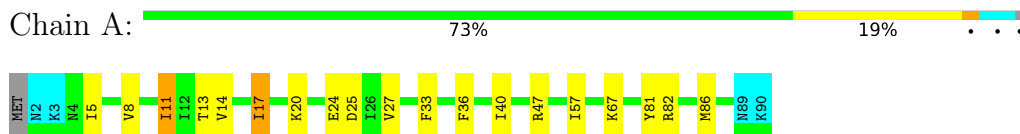


- Molecule 1: 10 kDa anti-sigma factor

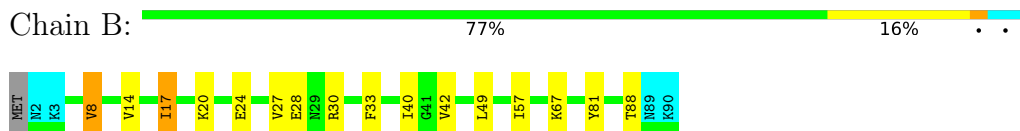


4.2.19 Score per residue for model 19

- Molecule 1: 10 kDa anti-sigma factor

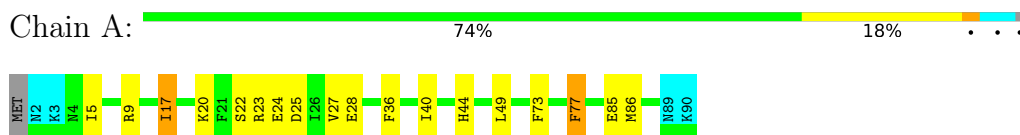


- Molecule 1: 10 kDa anti-sigma factor

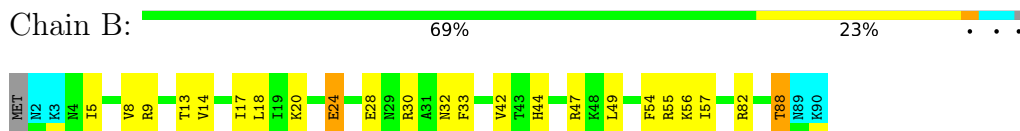


4.2.20 Score per residue for model 20

- Molecule 1: 10 kDa anti-sigma factor

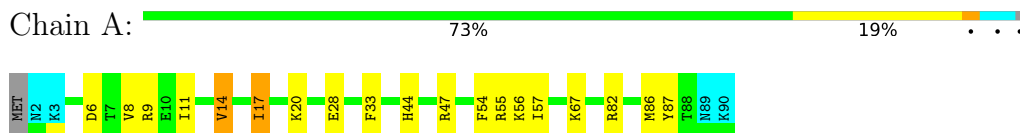


- Molecule 1: 10 kDa anti-sigma factor

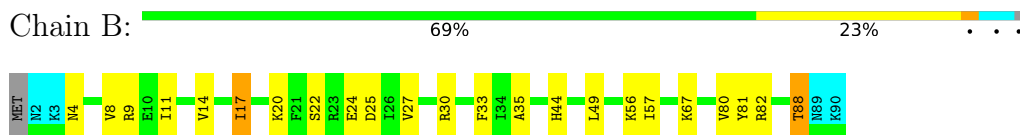


4.2.21 Score per residue for model 21

- Molecule 1: 10 kDa anti-sigma factor

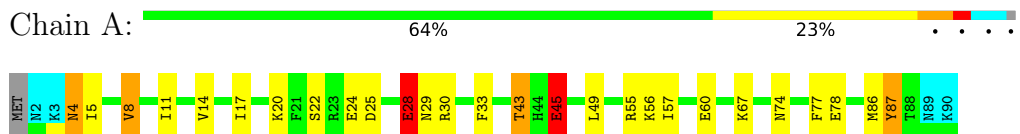


- Molecule 1: 10 kDa anti-sigma factor

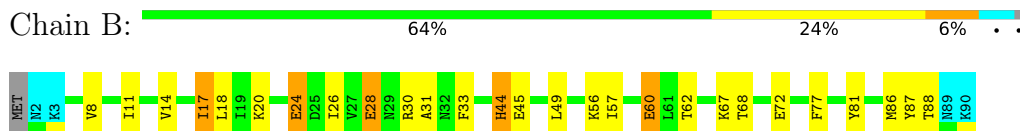


4.2.22 Score per residue for model 22

- Molecule 1: 10 kDa anti-sigma factor



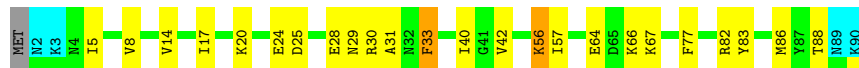
- Molecule 1: 10 kDa anti-sigma factor



4.2.23 Score per residue for model 23

- Molecule 1: 10 kDa anti-sigma factor

Chain A:  68% 24%



- Molecule 1: 10 kDa anti-sigma factor

Chain B:  70% 21%



5 Refinement protocol and experimental data overview

The models were refined using the following method: *dynamical simulated annealing/restrained energy minimization*.

Of the 100 calculated structures, 23 were deposited, based on the following criterion: *structures with acceptable covalent geometry, structures with the least restraint violations*.

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

Software name	Classification	Version
Xplor-NIH	refinement	2.9.3

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	1.47±0.01	0±0/713 (0.0± 0.0%)	1.84±0.06	16±5/960 (1.6± 0.5%)
1	B	1.47±0.01	0±0/713 (0.0± 0.0%)	1.84±0.06	16±5/960 (1.7± 0.5%)
All	All	1.47	0/32798 (0.0%)	1.84	733/44160 (1.7%)

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	2.5±1.3
1	B	0.0±0.0	2.3±1.4
All	All	0	109

There are no bond-length outliers.

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	B	58	VAL	N-CA-C	11.20	122.03	110.72	18	5
1	B	14	VAL	O-C-N	-9.88	112.29	121.87	20	1
1	B	72	GLU	N-CA-C	9.86	125.11	112.89	22	1
1	B	86	MET	N-CA-C	9.15	124.24	112.34	22	1
1	A	42	VAL	N-CA-CB	8.77	120.84	111.46	4	8
1	A	9	ARG	N-CA-C	8.68	120.75	111.28	13	4
1	A	14	VAL	O-C-N	-8.65	113.48	121.87	14	1
1	B	14	VAL	N-CA-C	8.65	119.44	110.62	20	4
1	B	77	PHE	CA-CB-CG	8.49	122.29	113.80	14	2
1	A	5	ILE	N-CA-C	7.92	118.70	110.62	14	5
1	B	33	PHE	CA-C-N	7.84	131.59	120.53	11	20
1	B	33	PHE	C-N-CA	7.84	131.59	120.53	11	20
1	B	11	ILE	CA-C-N	7.82	130.68	120.60	13	10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	B	11	ILE	C-N-CA	7.82	130.68	120.60	13	10
1	B	42	VAL	N-CA-CB	7.77	119.45	110.82	19	8
1	A	4	ASN	CA-C-N	7.74	130.59	120.60	17	7
1	A	4	ASN	C-N-CA	7.74	130.59	120.60	17	7
1	B	8	VAL	CA-C-N	7.71	130.82	120.65	18	20
1	B	8	VAL	C-N-CA	7.71	130.82	120.65	18	20
1	B	81	TYR	CA-C-N	-7.69	108.62	120.31	22	1
1	B	81	TYR	C-N-CA	-7.69	108.62	120.31	22	1
1	A	8	VAL	CA-C-N	7.63	130.84	120.54	6	15
1	A	8	VAL	C-N-CA	7.63	130.84	120.54	6	15
1	B	5	ILE	N-CA-C	7.55	118.29	110.36	16	4
1	B	17	ILE	CB-CA-C	-7.50	102.28	111.88	15	13
1	A	24	GLU	CA-C-N	7.45	131.14	120.79	22	6
1	A	24	GLU	C-N-CA	7.45	131.14	120.79	22	6
1	B	39	GLU	N-CA-CB	7.44	120.85	110.07	2	1
1	B	55	ARG	N-CA-C	7.35	121.89	112.34	14	1
1	A	17	ILE	CB-CA-C	-7.30	102.53	111.88	4	14
1	B	17	ILE	N-CA-CB	7.19	118.49	110.51	15	4
1	B	53	SER	CA-C-N	-7.19	111.10	120.44	14	1
1	B	53	SER	C-N-CA	-7.19	111.10	120.44	14	1
1	B	18	LEU	CA-C-O	7.17	128.22	119.97	14	1
1	A	28	GLU	N-CA-C	7.10	121.04	112.38	2	6
1	B	57	ILE	N-CA-C	7.01	117.12	110.53	3	5
1	A	59	SER	N-CA-C	6.89	118.68	111.03	14	1
1	B	62	THR	N-CA-CB	6.84	119.82	110.38	22	1
1	A	78	GLU	N-CA-C	6.82	119.30	111.11	22	1
1	A	69	LEU	N-CA-C	6.76	119.50	111.33	18	1
1	B	57	ILE	N-CA-CB	6.74	117.99	110.51	23	14
1	B	54	PHE	CA-C-N	-6.68	109.87	120.63	14	1
1	B	54	PHE	C-N-CA	-6.68	109.87	120.63	14	1
1	B	54	PHE	CA-CB-CG	-6.67	107.13	113.80	13	13
1	A	19	ILE	CA-C-N	6.67	129.55	120.54	6	1
1	A	19	ILE	C-N-CA	6.67	129.55	120.54	6	1
1	A	13	THR	CA-C-N	6.66	129.19	120.60	2	7
1	A	13	THR	C-N-CA	6.66	129.19	120.60	2	7
1	A	24	GLU	N-CA-C	6.63	120.47	112.38	5	2
1	A	57	ILE	N-CA-CB	6.62	118.29	110.55	4	20
1	A	35	ALA	CA-C-N	6.58	128.99	120.44	2	5
1	A	35	ALA	C-N-CA	6.58	128.99	120.44	2	5
1	A	8	VAL	O-C-N	-6.56	115.48	121.91	21	1
1	A	86	MET	N-CA-C	6.49	120.06	111.75	1	8

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	B	4	ASN	CA-C-N	6.46	128.93	120.60	7	7
1	B	4	ASN	C-N-CA	6.46	128.93	120.60	7	7
1	B	27	VAL	CA-C-N	6.44	129.79	120.38	21	7
1	B	27	VAL	C-N-CA	6.44	129.79	120.38	21	7
1	A	5	ILE	CA-C-N	6.41	129.73	120.38	10	5
1	A	5	ILE	C-N-CA	6.41	129.73	120.38	10	5
1	B	57	ILE	CA-CB-CG1	6.38	121.25	110.40	3	1
1	B	54	PHE	N-CA-C	6.34	117.86	111.07	14	1
1	A	57	ILE	N-CA-C	6.33	117.01	110.36	13	3
1	B	28	GLU	N-CA-C	6.32	120.55	112.34	12	6
1	A	54	PHE	CA-CB-CG	-6.30	107.50	113.80	21	7
1	A	46	GLY	N-CA-C	-6.27	106.66	115.32	2	1
1	B	88	THR	CA-CB-OG1	6.26	118.99	109.60	20	4
1	B	56	LYS	N-CA-C	6.24	117.75	111.07	1	2
1	B	7	THR	CB-CA-C	-6.23	100.45	110.79	23	1
1	A	77	PHE	CB-CA-C	6.21	120.72	110.90	16	5
1	A	6	ASP	CA-CB-CG	-6.18	106.42	112.60	9	1
1	A	82	ARG	CB-CA-C	-6.18	100.54	110.79	7	11
1	A	11	ILE	CA-C-N	6.17	129.24	120.53	5	9
1	A	11	ILE	C-N-CA	6.17	129.24	120.53	5	9
1	A	27	VAL	CA-C-N	6.17	129.39	120.38	14	5
1	A	27	VAL	C-N-CA	6.17	129.39	120.38	14	5
1	A	33	PHE	CA-C-N	6.11	128.83	120.46	23	16
1	A	33	PHE	C-N-CA	6.11	128.83	120.46	23	16
1	B	76	GLY	CA-C-N	-6.09	110.17	120.58	14	1
1	B	76	GLY	C-N-CA	-6.09	110.17	120.58	14	1
1	A	36	PHE	CA-CB-CG	6.07	119.87	113.80	14	4
1	A	9	ARG	N-CA-CB	-6.06	101.21	110.12	13	2
1	B	60	GLU	N-CA-C	6.04	123.66	110.80	22	1
1	A	64	GLU	N-CA-C	6.02	118.61	111.33	23	2
1	B	82	ARG	CA-C-N	6.02	128.66	120.54	18	5
1	B	82	ARG	C-N-CA	6.02	128.66	120.54	18	5
1	A	56	LYS	N-CA-C	6.00	117.90	111.36	18	6
1	A	79	GLY	O-C-N	-6.00	116.43	122.19	14	1
1	B	55	ARG	N-CA-CB	-5.99	101.00	110.22	9	2
1	A	74	ASN	N-CA-C	5.98	123.54	110.80	22	1
1	B	5	ILE	CA-C-N	5.96	128.18	120.44	2	3
1	B	5	ILE	C-N-CA	5.96	128.18	120.44	2	3
1	A	55	ARG	N-CA-CB	-5.96	101.05	110.28	11	3
1	A	19	ILE	CB-CA-C	5.94	119.82	112.04	3	1
1	A	25	ASP	N-CA-C	5.93	119.34	111.75	20	5

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	29	ASN	CA-C-N	5.90	128.50	120.54	22	5
1	A	29	ASN	C-N-CA	5.90	128.50	120.54	22	5
1	B	86	MET	CG-SD-CE	-5.88	87.98	100.90	5	5
1	A	14	VAL	N-CA-C	5.87	116.52	110.36	6	5
1	A	22	SER	CA-C-N	5.84	131.88	123.17	6	1
1	A	22	SER	C-N-CA	5.84	131.88	123.17	6	1
1	A	43	THR	N-CA-C	5.84	118.08	110.43	22	1
1	A	58	VAL	N-CA-C	5.83	117.36	111.00	15	2
1	B	4	ASN	CA-CB-CG	-5.83	106.77	112.60	21	1
1	B	26	ILE	CA-C-N	5.82	128.43	120.46	2	5
1	B	26	ILE	C-N-CA	5.82	128.43	120.46	2	5
1	B	31	ALA	N-CA-C	5.82	117.62	111.28	3	5
1	B	18	LEU	N-CA-C	5.82	118.57	111.82	14	1
1	A	47	ARG	N-CA-C	-5.82	100.22	109.76	2	1
1	A	25	ASP	CA-C-N	5.81	128.10	120.60	22	1
1	A	25	ASP	C-N-CA	5.81	128.10	120.60	22	1
1	B	12	ILE	CA-C-N	5.80	128.63	120.28	10	1
1	B	12	ILE	C-N-CA	5.80	128.63	120.28	10	1
1	B	70	ILE	N-CA-C	5.75	115.94	110.42	18	1
1	B	35	ALA	CA-C-N	5.74	127.97	120.28	21	4
1	B	35	ALA	C-N-CA	5.74	127.97	120.28	21	4
1	A	82	ARG	N-CA-CB	5.73	118.47	109.94	4	2
1	A	38	ASN	CA-CB-CG	-5.72	106.88	112.60	8	2
1	A	6	ASP	N-CA-C	5.70	117.95	111.11	21	1
1	A	68	THR	N-CA-C	-5.66	103.93	111.24	18	1
1	A	31	ALA	N-CA-C	5.64	117.23	111.14	10	2
1	B	32	ASN	CA-C-N	5.63	128.10	120.44	9	3
1	B	32	ASN	C-N-CA	5.63	128.10	120.44	9	3
1	A	86	MET	CG-SD-CE	-5.63	88.51	100.90	19	5
1	B	36	PHE	CA-CB-CG	5.61	119.41	113.80	5	3
1	A	87	TYR	N-CA-C	5.57	119.70	113.01	22	1
1	B	86	MET	N-CA-CB	5.57	118.91	110.28	16	1
1	B	38	ASN	CA-C-N	5.56	127.73	120.28	12	1
1	B	38	ASN	C-N-CA	5.56	127.73	120.28	12	1
1	A	82	ARG	CA-C-N	5.54	128.17	120.63	18	1
1	A	82	ARG	C-N-CA	5.54	128.17	120.63	18	1
1	B	50	ASN	CA-CB-CG	-5.54	107.06	112.60	16	1
1	B	88	THR	CB-CA-C	5.52	119.71	110.44	8	1
1	B	5	ILE	N-CA-CB	5.50	116.61	110.51	20	1
1	B	63	GLN	CA-C-N	5.48	127.62	120.28	1	3
1	B	63	GLN	C-N-CA	5.48	127.62	120.28	1	3

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	88	THR	CB-CA-C	5.48	119.98	110.72	3	1
1	B	59	SER	N-CA-C	5.48	117.11	111.03	17	1
1	A	77	PHE	CA-CB-CG	5.47	119.27	113.80	20	3
1	A	22	SER	N-CA-C	5.47	118.94	111.39	20	1
1	B	60	GLU	CA-C-N	5.47	129.46	121.31	13	1
1	B	60	GLU	C-N-CA	5.47	129.46	121.31	13	1
1	A	32	ASN	CA-C-N	5.47	127.55	120.44	11	2
1	A	32	ASN	C-N-CA	5.47	127.55	120.44	11	2
1	B	85	GLU	N-CA-C	5.46	117.23	111.28	16	1
1	B	82	ARG	CB-CA-C	-5.46	102.08	110.81	10	6
1	B	44	HIS	N-CA-CB	5.45	119.70	110.49	13	1
1	B	77	PHE	CB-CA-C	5.44	119.42	110.88	22	3
1	B	56	LYS	N-CA-CB	5.44	117.89	110.01	17	1
1	A	71	ASP	N-CA-C	5.43	119.08	112.23	14	1
1	B	83	TYR	CA-C-N	5.42	127.80	120.38	1	1
1	B	83	TYR	C-N-CA	5.42	127.80	120.38	1	1
1	B	16	SER	N-CA-CB	5.41	119.23	110.40	14	1
1	A	50	ASN	CA-CB-CG	-5.41	107.19	112.60	18	2
1	A	84	LEU	N-CA-C	5.41	117.60	111.11	14	1
1	B	29	ASN	CA-C-N	5.40	128.92	120.82	15	3
1	B	29	ASN	C-N-CA	5.40	128.92	120.82	15	3
1	A	53	SER	N-CA-C	5.39	118.96	112.38	11	2
1	B	26	ILE	N-CA-CB	5.37	117.45	110.57	11	1
1	A	70	ILE	CB-CA-C	-5.37	104.81	112.22	14	1
1	B	12	ILE	N-CA-C	5.36	116.73	110.62	4	1
1	B	75	GLU	CA-C-N	5.36	125.78	119.94	13	2
1	B	75	GLU	C-N-CA	5.36	125.78	119.94	13	2
1	A	84	LEU	CB-CA-C	-5.34	102.27	110.81	14	1
1	B	24	GLU	CA-C-N	5.34	127.69	120.38	20	6
1	B	24	GLU	C-N-CA	5.34	127.69	120.38	20	6
1	B	44	HIS	CA-CB-CG	5.33	119.13	113.80	22	1
1	B	52	ASN	CA-C-O	5.32	125.90	119.31	13	1
1	B	7	THR	CA-C-N	5.31	127.74	120.46	13	1
1	B	7	THR	C-N-CA	5.31	127.74	120.46	13	1
1	A	63	GLN	N-CA-C	5.29	117.05	111.28	15	1
1	B	6	ASP	CA-C-N	5.29	127.63	120.65	13	1
1	B	6	ASP	C-N-CA	5.29	127.63	120.65	13	1
1	A	17	ILE	N-CA-CB	5.29	116.73	110.55	11	2
1	B	48	LYS	N-CA-C	5.28	117.89	110.23	14	1
1	A	69	LEU	CA-C-N	5.27	127.68	120.46	13	1
1	A	69	LEU	C-N-CA	5.27	127.68	120.46	13	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	B	43	THR	N-CA-CB	5.27	119.22	110.63	13	1
1	B	10	GLU	N-CA-C	5.27	117.93	111.82	1	2
1	B	68	THR	N-CA-C	5.26	117.01	111.28	22	1
1	A	77	PHE	N-CA-C	-5.26	105.61	112.23	22	1
1	A	36	PHE	CB-CA-C	5.25	119.77	110.85	20	1
1	B	25	ASP	N-CA-C	5.23	118.44	111.75	14	1
1	A	88	THR	CA-CB-OG1	5.21	117.42	109.60	1	2
1	A	73	PHE	N-CA-CB	5.19	119.17	111.23	2	1
1	B	18	LEU	N-CA-CB	5.18	117.74	110.12	11	1
1	B	18	LEU	O-C-N	-5.18	115.90	122.27	14	1
1	A	47	ARG	N-CA-CB	-5.18	102.81	111.20	18	1
1	B	10	GLU	CA-C-O	5.17	126.44	120.90	14	1
1	A	36	PHE	N-CA-C	5.17	117.59	111.33	8	1
1	B	82	ARG	N-CA-CB	5.17	117.65	109.94	10	1
1	A	62	THR	N-CA-CB	5.17	117.83	110.44	10	1
1	B	34	ILE	CA-C-N	-5.15	112.98	120.29	5	1
1	B	34	ILE	C-N-CA	-5.15	112.98	120.29	5	1
1	A	40	ILE	CA-CB-CG1	5.14	119.14	110.40	17	1
1	A	39	GLU	N-CA-C	5.14	116.69	111.14	14	1
1	B	64	GLU	N-CA-C	5.14	116.88	111.28	14	1
1	B	12	ILE	CA-CB-CG2	5.14	119.23	110.50	10	1
1	B	30	ARG	N-CA-CB	5.14	117.52	110.07	14	1
1	B	40	ILE	CA-CB-CG1	5.13	119.12	110.40	17	1
1	A	27	VAL	N-CA-C	5.12	116.45	110.62	14	1
1	B	26	ILE	CB-CA-C	-5.11	103.82	112.16	22	1
1	A	30	ARG	N-CA-CB	5.11	118.11	109.78	2	1
1	B	13	THR	O-C-N	5.11	127.33	122.07	20	1
1	A	85	GLU	N-CA-C	5.10	117.23	111.11	8	2
1	A	51	GLN	N-CA-C	5.10	116.53	111.07	1	1
1	A	45	GLU	CA-C-N	5.08	131.38	121.41	22	1
1	A	45	GLU	C-N-CA	5.08	131.38	121.41	22	1
1	A	75	GLU	CA-C-N	5.08	125.48	119.94	10	1
1	A	75	GLU	C-N-CA	5.08	125.48	119.94	10	1
1	B	14	VAL	CB-CA-C	-5.07	105.30	112.14	20	1
1	A	56	LYS	N-CA-CB	5.04	117.31	110.01	3	1
1	B	8	VAL	CB-CA-C	-5.04	105.44	112.04	6	1
1	A	8	VAL	N-CA-CB	-5.02	103.71	110.54	1	1
1	A	66	LYS	CA-C-N	5.02	127.01	120.28	23	1
1	A	66	LYS	C-N-CA	5.02	127.01	120.28	23	1
1	A	62	THR	CA-CB-OG1	5.01	117.11	109.60	10	1

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	B	30	ARG	Sidechain	10
1	B	81	TYR	Sidechain	10
1	A	9	ARG	Sidechain	9
1	A	55	ARG	Sidechain	8
1	A	30	ARG	Sidechain	7
1	B	55	ARG	Sidechain	7
1	B	9	ARG	Sidechain	7
1	A	33	PHE	Sidechain	6
1	A	82	ARG	Sidechain	6
1	B	33	PHE	Sidechain	5
1	A	47	ARG	Sidechain	5
1	A	23	ARG	Sidechain	4
1	A	81	TYR	Sidechain	4
1	B	82	ARG	Sidechain	4
1	B	47	ARG	Sidechain	3
1	A	88	THR	Peptide	3
1	B	88	THR	Peptide	2
1	B	48	LYS	Peptide	1
1	A	48	LYS	Peptide	1
1	A	54	PHE	Sidechain	1
1	B	23	ARG	Sidechain	1
1	A	73	PHE	Sidechain	1
1	A	43	THR	Peptide	1
1	A	83	TYR	Sidechain	1
1	B	36	PHE	Sidechain	1
1	B	83	TYR	Sidechain	1

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	703	700	699	3±1
1	B	703	700	699	3±1
All	All	32338	32200	32154	74

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 1.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:17:ILE:HG21	1:B:17:ILE:HG21	0.62	1.70	23	22
1:A:14:VAL:HG13	1:B:17:ILE:HD11	0.58	1.74	13	17
1:A:17:ILE:HD11	1:B:14:VAL:HG13	0.55	1.79	1	5
1:B:17:ILE:HG23	1:B:21:PHE:CE2	0.51	2.40	3	3
1:A:76:GLY:O	1:A:80:VAL:HG22	0.48	2.08	16	1
1:B:12:ILE:HG23	1:B:80:VAL:HG21	0.46	1.88	11	2
1:B:76:GLY:O	1:B:80:VAL:HG22	0.46	2.09	8	2
1:A:44:HIS:CD2	1:A:45:GLU:CD	0.45	2.95	7	1
1:A:28:GLU:HA	1:A:87:TYR:CD2	0.45	2.47	21	6
1:B:44:HIS:CD2	1:B:45:GLU:CD	0.45	2.94	22	1
1:A:17:ILE:CG2	1:B:17:ILE:HG21	0.43	2.44	10	3
1:B:28:GLU:HA	1:B:87:TYR:CD2	0.43	2.48	22	1
1:A:44:HIS:CD2	1:A:45:GLU:HG3	0.43	2.48	17	1
1:A:7:THR:OG1	1:A:44:HIS:CE1	0.43	2.72	3	1
1:A:17:ILE:HG21	1:B:17:ILE:CG2	0.41	2.41	23	2
1:A:17:ILE:CD1	1:B:14:VAL:HG13	0.41	2.46	13	1
1:B:40:ILE:HD11	1:B:42:VAL:HG23	0.41	1.93	11	1
1:A:28:GLU:HA	1:A:87:TYR:CG	0.41	2.51	21	1
1:A:83:TYR:O	1:A:87:TYR:CD2	0.41	2.74	10	1
1:B:34:ILE:HD11	1:B:54:PHE:CD2	0.40	2.51	18	1
1:B:7:THR:OG1	1:B:44:HIS:CE1	0.40	2.74	1	1

6.3 Torsion angles [i](#)

6.3.1 Protein backbone [i](#)

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	85/90 (94%)	76±2 (90±2%)	8±2 (9±2%)	1±1 (1±1%)	14	63
1	B	85/90 (94%)	76±2 (90±3%)	8±2 (9±3%)	1±1 (1±1%)	13	61
All	All	3910/4140 (94%)	3512 (90%)	353 (9%)	45 (1%)	13	61

All 12 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	B	44	HIS	16
1	A	44	HIS	15
1	B	22	SER	3
1	A	22	SER	2
1	A	28	GLU	2
1	B	74	ASN	1
1	B	28	GLU	1
1	A	25	ASP	1
1	B	25	ASP	1
1	A	45	GLU	1
1	A	60	GLU	1
1	B	60	GLU	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	78/83 (94%)	71±2 (91±2%)	7±2 (9±2%)	11	58
1	B	78/83 (94%)	71±2 (91±2%)	7±2 (9±2%)	11	58
All	All	3588/3818 (94%)	3273 (91%)	315 (9%)	11	58

All 57 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	B	20	LYS	22
1	B	24	GLU	22
1	A	20	LYS	20
1	A	24	GLU	20
1	A	49	LEU	18
1	A	77	PHE	14
1	B	49	LEU	14
1	A	40	ILE	13
1	B	77	PHE	13
1	B	67	LYS	12
1	B	40	ILE	11
1	A	67	LYS	11
1	B	88	THR	11

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Mol	Chain	Res	Type	Models (Total)
1	A	88	THR	8
1	B	7	THR	7
1	B	17	ILE	6
1	B	18	LEU	6
1	B	56	LYS	6
1	B	8	VAL	5
1	A	42	VAL	5
1	B	80	VAL	4
1	A	17	ILE	4
1	A	85	GLU	4
1	A	30	ARG	4
1	A	56	LYS	4
1	A	14	VAL	3
1	B	14	VAL	3
1	A	80	VAL	3
1	B	55	ARG	2
1	A	7	THR	2
1	A	9	ARG	2
1	B	9	ARG	2
1	B	42	VAL	2
1	A	13	THR	2
1	A	18	LEU	2
1	A	75	GLU	2
1	A	5	ILE	2
1	A	55	ARG	2
1	B	30	ARG	2
1	A	11	ILE	2
1	A	28	GLU	2
1	B	43	THR	1
1	A	57	ILE	1
1	B	57	ILE	1
1	A	23	ARG	1
1	B	12	ILE	1
1	B	28	GLU	1
1	B	39	GLU	1
1	A	65	ASP	1
1	B	13	THR	1
1	B	6	ASP	1
1	B	19	ILE	1
1	A	19	ILE	1
1	A	68	THR	1
1	A	4	ASN	1

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Mol	Chain	Res	Type	Models (Total)
1	A	8	VAL	1
1	A	45	GLU	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