



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 03:34 AM UTC

PDB ID : 1C7V / pdb_00001c7v
Title : NMR SOLUTION STRUCTURE OF THE CALCIUM-BOUND C-TERMINAL DOMAIN (W81-S161) OF CALCIUM VECTOR PROTEIN FROM AMPHIOXUS
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Deposited on : 2000-03-27

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

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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

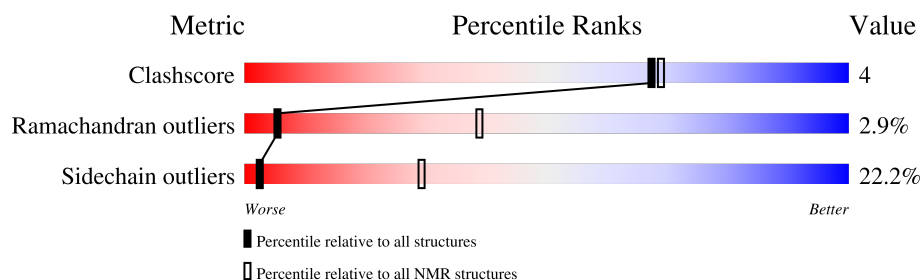
1 Overall quality at a glance

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	81	

2 Ensemble composition and analysis

This entry contains 20 models. Model 7 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:89-A:114, A:123-A:152 (56)	1.07	7

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 2 clusters. No single-model clusters were found.

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

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 1049 atoms, of which 510 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called CALCIUM VECTOR PROTEIN.

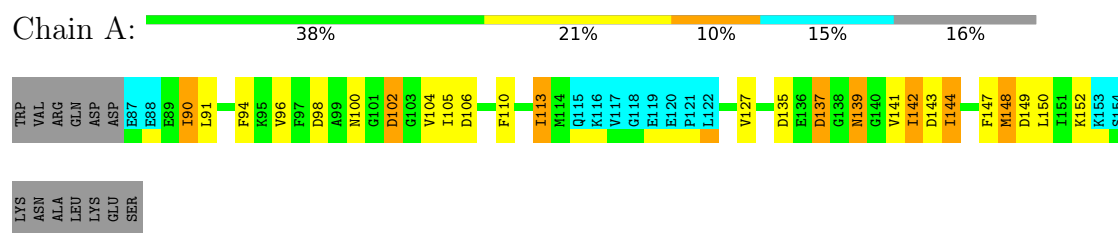
Mol	Chain	Residues	Atoms						Trace
1	A	68	Total	C	H	N	O	S	0
			1049	341	510	80	115	3	

4 Residue-property plots [i](#)

4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

• Molecule 1: CALCIUM VECTOR PROTEIN

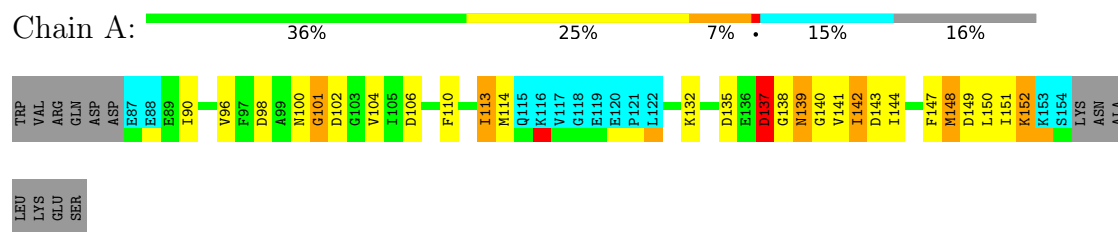


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

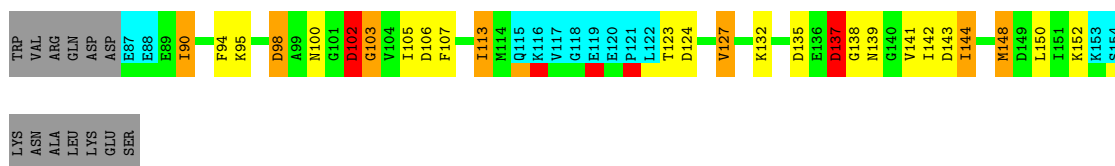
• Molecule 1: CALCIUM VECTOR PROTEIN



4.2.2 Score per residue for model 2

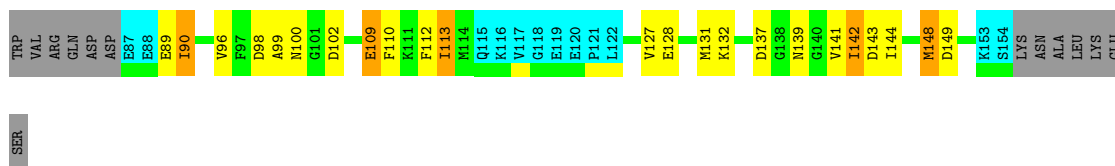
• Molecule 1: CALCIUM VECTOR PROTEIN





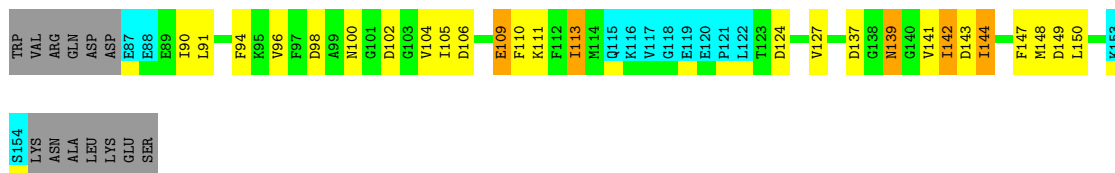
4.2.3 Score per residue for model 3

- Molecule 1: CALCIUM VECTOR PROTEIN



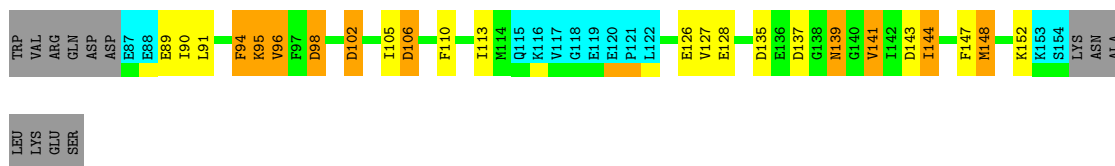
4.2.4 Score per residue for model 4

- Molecule 1: CALCIUM VECTOR PROTEIN



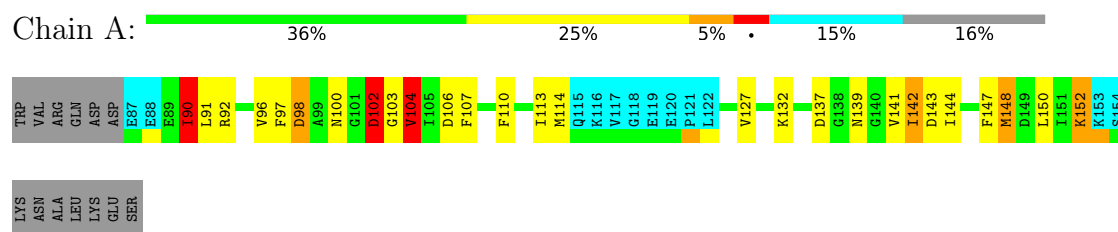
4.2.5 Score per residue for model 5

- Molecule 1: CALCIUM VECTOR PROTEIN



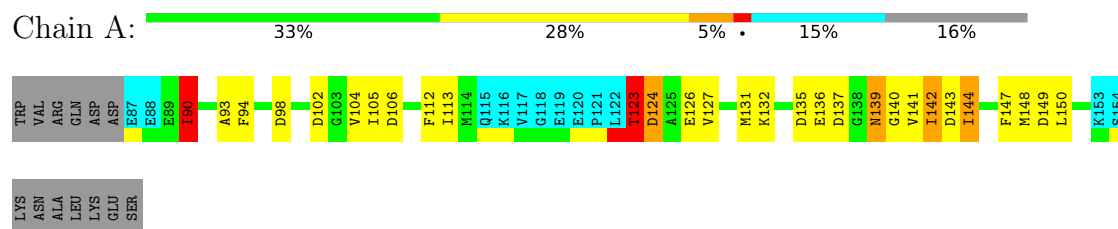
4.2.6 Score per residue for model 6

- Molecule 1: CALCIUM VECTOR PROTEIN



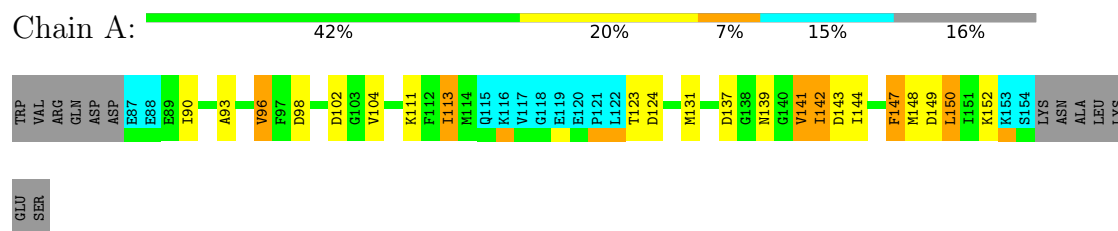
4.2.7 Score per residue for model 7 (medoid)

- Molecule 1: CALCIUM VECTOR PROTEIN



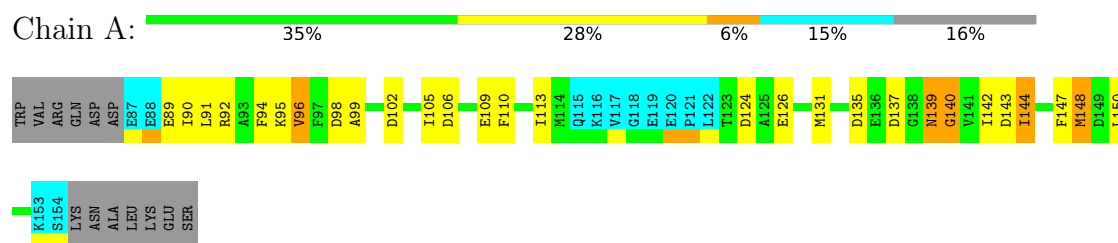
4.2.8 Score per residue for model 8

- Molecule 1: CALCIUM VECTOR PROTEIN



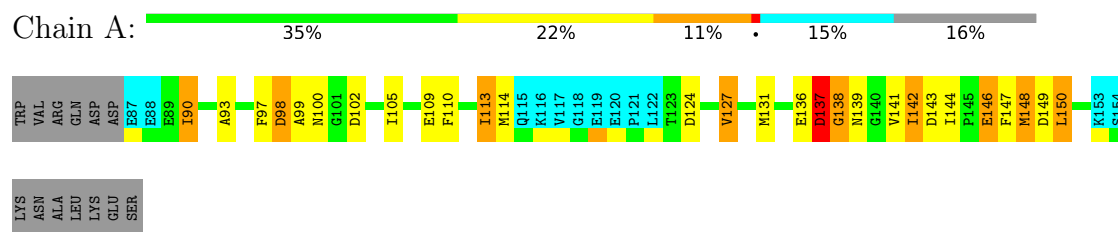
4.2.9 Score per residue for model 9

- Molecule 1: CALCIUM VECTOR PROTEIN



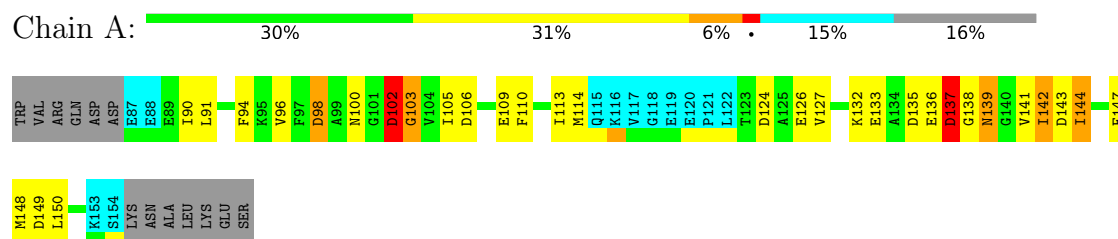
4.2.10 Score per residue for model 10

• Molecule 1: CALCIUM VECTOR PROTEIN



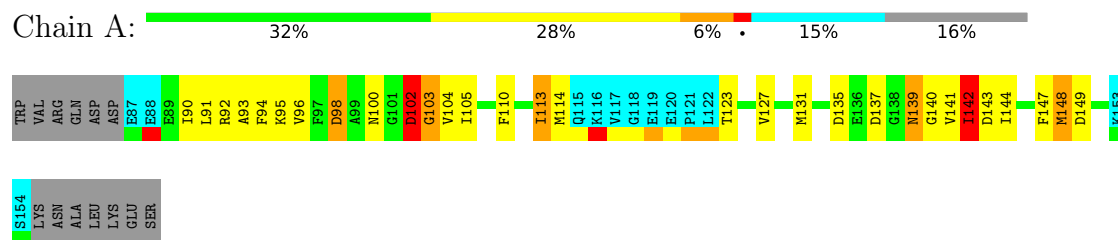
4.2.11 Score per residue for model 11

• Molecule 1: CALCIUM VECTOR PROTEIN



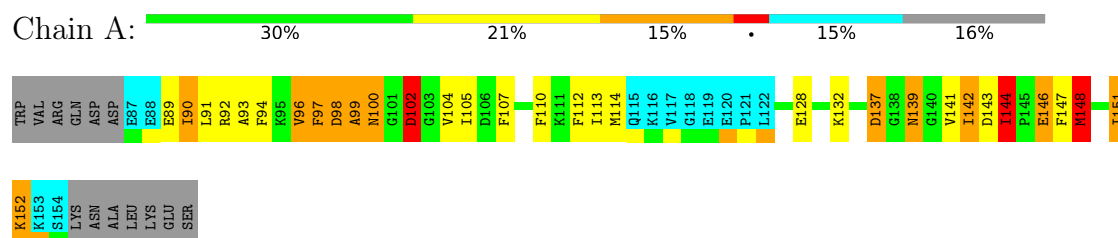
4.2.12 Score per residue for model 12

• Molecule 1: CALCIUM VECTOR PROTEIN



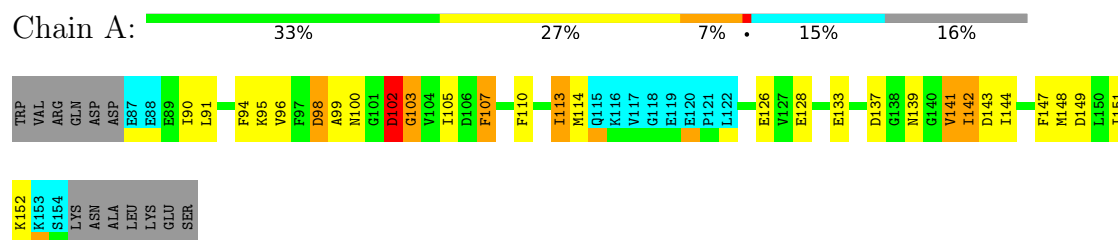
4.2.13 Score per residue for model 13

• Molecule 1: CALCIUM VECTOR PROTEIN



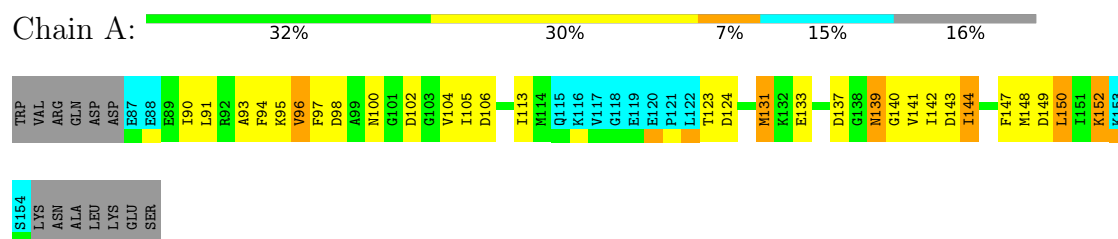
4.2.14 Score per residue for model 14

• Molecule 1: CALCIUM VECTOR PROTEIN



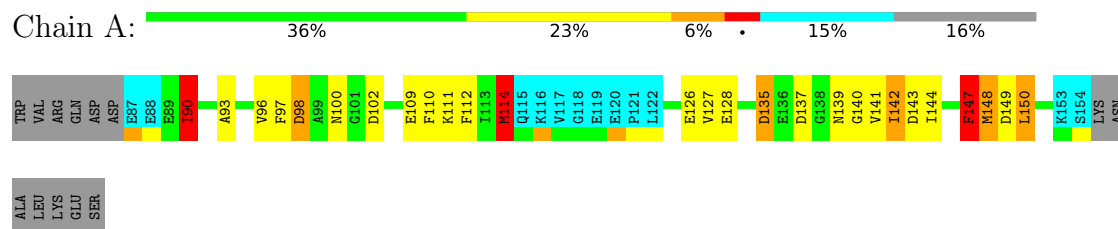
4.2.15 Score per residue for model 15

• Molecule 1: CALCIUM VECTOR PROTEIN



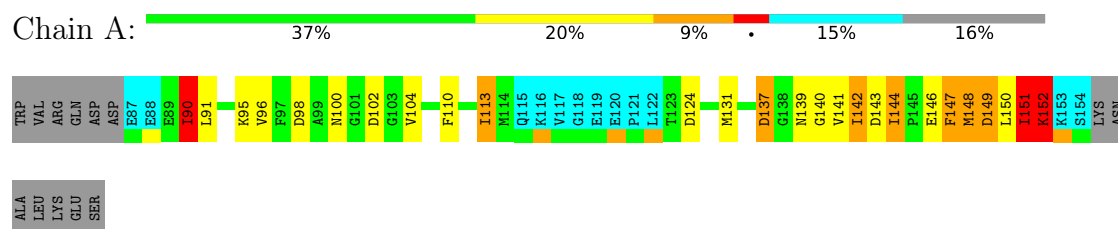
4.2.16 Score per residue for model 16

• Molecule 1: CALCIUM VECTOR PROTEIN



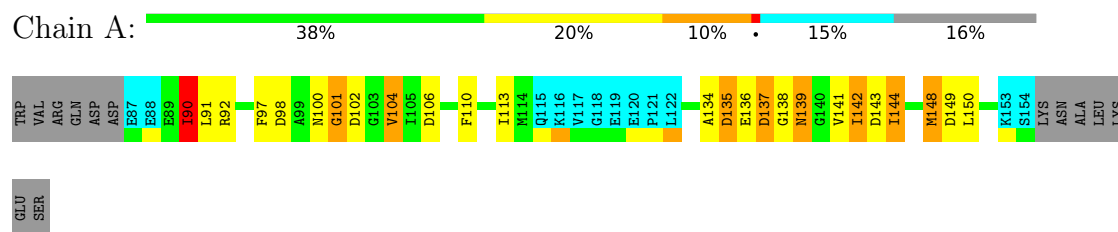
4.2.17 Score per residue for model 17

• Molecule 1: CALCIUM VECTOR PROTEIN



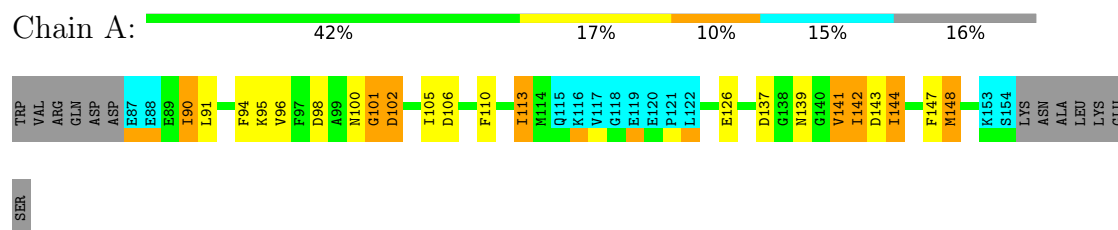
4.2.18 Score per residue for model 18

• Molecule 1: CALCIUM VECTOR PROTEIN



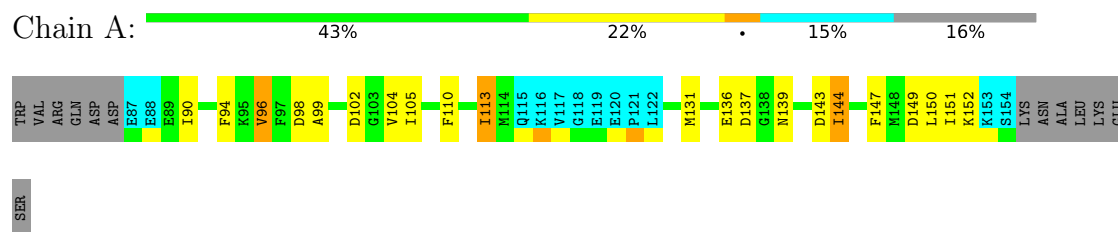
4.2.19 Score per residue for model 19

• Molecule 1: CALCIUM VECTOR PROTEIN



4.2.20 Score per residue for model 20

• Molecule 1: CALCIUM VECTOR PROTEIN



5 Refinement protocol and experimental data overview

The models were refined using the following method: *Distance Geometry (DGII) and Simulated Annealing (Discover)*.

Of the 100 calculated structures, 20 were deposited, based on the following criterion: *agreement with NOE restraints, geometry quality*.

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

Software name	Classification	Version
Felix	structure solution	97.0
Felix	refinement	97.0

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	1.88±0.05	3±1/451 (0.7± 0.3%)	2.25±0.22	22±5/606 (3.6± 0.8%)
All	All	1.88	66/9020 (0.7%)	2.26	441/12120 (3.6%)

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	0.8±0.8
All	All	0	16

All unique bond 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(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	144	ILE	CA-CB	11.09	1.59	1.54	20	12
1	A	144	ILE	CA-C	7.65	1.59	1.52	2	16
1	A	89	GLU	CA-C	7.23	1.62	1.52	5	1
1	A	151	ILE	CA-C	7.07	1.61	1.52	17	5
1	A	137	ASP	CA-C	6.38	1.61	1.52	13	1
1	A	100	ASN	CA-C	6.02	1.60	1.52	13	1
1	A	96	VAL	CA-C	5.91	1.60	1.52	15	6
1	A	109	GLU	CA-C	5.84	1.60	1.52	16	1
1	A	106	ASP	CA-C	5.80	1.59	1.52	6	1
1	A	103	GLY	C-N	5.72	1.36	1.33	6	1
1	A	139	ASN	CA-C	5.64	1.59	1.52	10	1
1	A	104	VAL	CA-C	5.59	1.58	1.52	6	1
1	A	136	GLU	CA-C	5.58	1.60	1.52	11	3
1	A	149	ASP	CA-C	5.53	1.59	1.52	16	1
1	A	104	VAL	CA-CB	5.51	1.59	1.55	6	1
1	A	99	ALA	N-CA	5.43	1.53	1.46	13	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	102	ASP	CA-C	5.35	1.59	1.52	19	2
1	A	148	MET	CA-C	5.32	1.59	1.52	9	2
1	A	146	GLU	CA-C	5.32	1.59	1.52	17	1
1	A	152	LYS	N-CA	5.24	1.52	1.46	17	1
1	A	95	LYS	CA-C	5.24	1.60	1.52	14	2
1	A	147	PHE	CA-C	5.23	1.59	1.52	13	1
1	A	112	PHE	CA-C	5.23	1.59	1.53	3	1
1	A	135	ASP	C-N	-5.18	1.26	1.33	18	1
1	A	90	ILE	CA-CB	5.12	1.60	1.54	16	1
1	A	126	GLU	CA-C	5.08	1.59	1.52	5	1

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	135	ASP	O-C-N	-46.14	62.84	122.82	18	1
1	A	135	ASP	CA-C-O	-21.49	96.31	120.60	18	1
1	A	137	ASP	O-C-N	-17.99	98.67	122.59	11	2
1	A	98	ASP	CA-CB-CG	14.59	127.19	112.60	19	20
1	A	97	PHE	CA-CB-CG	14.39	128.19	113.80	15	3
1	A	90	ILE	N-CA-CB	14.23	126.73	110.65	16	2
1	A	143	ASP	CA-CB-CG	13.43	126.03	112.60	9	20
1	A	147	PHE	CA-CB-CG	12.28	126.08	113.80	16	3
1	A	139	ASN	CA-CB-CG	11.21	123.81	112.60	18	20
1	A	112	PHE	CA-CB-CG	11.15	124.95	113.80	13	3
1	A	110	PHE	CA-CB-CG	11.11	124.91	113.80	18	15
1	A	149	ASP	CA-C-O	-10.85	109.30	120.90	17	1
1	A	90	ILE	CA-CB-CG1	10.23	127.79	110.40	16	8
1	A	139	ASN	O-C-N	-10.20	105.21	121.53	9	1
1	A	148	MET	CA-CB-CG	9.95	134.00	114.10	1	7
1	A	106	ASP	CA-CB-CG	9.44	122.04	112.60	1	10
1	A	141	VAL	CA-CB-CG2	9.40	126.39	110.40	5	8
1	A	102	ASP	N-CA-CB	9.30	126.21	110.49	13	2
1	A	113	ILE	N-CA-C	9.19	121.32	111.58	4	11
1	A	100	ASN	N-CA-CB	9.16	125.97	110.49	13	3
1	A	109	GLU	N-CA-CB	-9.13	96.77	110.01	16	2
1	A	102	ASP	CA-CB-CG	9.05	121.65	112.60	2	18
1	A	127	VAL	CA-CB-CG1	8.83	125.41	110.40	2	4
1	A	102	ASP	CA-C-N	8.58	134.47	122.29	6	1
1	A	102	ASP	C-N-CA	8.58	134.47	122.29	6	1
1	A	144	ILE	N-CA-CB	8.27	115.71	110.50	2	8

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	137	ASP	CA-C-O	-8.21	108.77	120.51	11	2
1	A	90	ILE	CA-CB-CG2	8.07	124.22	110.50	18	1
1	A	90	ILE	CB-CA-C	-7.89	101.69	111.87	16	2
1	A	144	ILE	CA-C-O	-7.88	113.41	118.69	3	6
1	A	134	ALA	N-CA-C	7.86	122.87	113.28	18	1
1	A	94	PHE	CA-CB-CG	7.84	121.64	113.80	5	1
1	A	152	LYS	N-CA-C	7.69	122.24	113.02	20	5
1	A	137	ASP	N-CA-C	7.67	122.64	113.28	4	7
1	A	90	ILE	N-CA-C	-7.60	103.28	110.42	16	2
1	A	104	VAL	N-CA-C	7.59	116.96	107.70	6	2
1	A	137	ASP	CA-CB-CG	7.57	120.17	112.60	4	15
1	A	103	GLY	N-CA-C	7.43	130.80	113.18	2	4
1	A	141	VAL	CA-CB-CG1	7.41	123.00	110.40	19	1
1	A	150	LEU	N-CA-CB	-7.36	98.87	110.28	2	4
1	A	124	ASP	CA-CB-CG	7.25	119.86	112.60	2	7
1	A	100	ASN	CA-C-N	7.21	135.55	121.41	13	1
1	A	100	ASN	C-N-CA	7.21	135.55	121.41	13	1
1	A	139	ASN	CA-C-O	-7.19	107.48	120.42	9	1
1	A	142	ILE	N-CA-C	7.18	118.37	108.89	18	8
1	A	99	ALA	CA-C-N	7.16	135.22	121.54	13	1
1	A	99	ALA	C-N-CA	7.16	135.22	121.54	13	1
1	A	102	ASP	CA-C-O	-7.09	109.07	121.62	11	4
1	A	139	ASN	OD1-CG-ND2	-7.06	115.54	122.60	2	9
1	A	100	ASN	CA-CB-CG	7.04	119.64	112.60	2	12
1	A	144	ILE	CB-CA-C	-6.92	107.02	114.35	2	5
1	A	142	ILE	CA-CB-CG1	6.85	122.05	110.40	1	15
1	A	148	MET	CA-C-O	-6.84	113.66	120.70	3	5
1	A	124	ASP	N-CA-C	6.83	125.34	110.80	7	4
1	A	98	ASP	CB-CA-C	6.78	119.93	110.16	14	3
1	A	114	MET	N-CA-C	6.78	119.25	111.11	11	2
1	A	135	ASP	N-CA-CB	6.74	120.40	110.35	7	3
1	A	123	THR	N-CA-CB	6.74	121.49	110.37	8	4
1	A	140	GLY	N-CA-C	6.72	129.10	113.18	9	1
1	A	152	LYS	CA-C-N	6.71	134.35	121.54	14	3
1	A	152	LYS	C-N-CA	6.71	134.35	121.54	14	3
1	A	146	GLU	CB-CG-CD	6.66	123.92	112.60	10	3
1	A	135	ASP	CA-CB-CG	6.63	119.23	112.60	2	4
1	A	93	ALA	CA-C-O	-6.63	114.07	120.90	15	6
1	A	141	VAL	N-CA-CB	6.57	121.36	110.86	8	3
1	A	137	ASP	N-CA-CB	6.53	121.52	110.49	13	5
1	A	141	VAL	CB-CA-C	6.49	119.99	110.77	16	8

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	98	ASP	N-CA-CB	-6.35	99.97	109.69	17	2
1	A	107	PHE	N-CA-C	6.25	117.79	110.97	6	2
1	A	95	LYS	CB-CA-C	6.23	123.33	110.31	5	8
1	A	144	ILE	CA-CB-CG1	6.19	120.92	110.40	2	3
1	A	99	ALA	N-CA-C	6.18	117.68	111.07	14	1
1	A	149	ASP	CA-C-N	6.15	128.44	120.44	15	2
1	A	149	ASP	C-N-CA	6.15	128.44	120.44	15	2
1	A	151	ILE	CA-CB-CG1	6.10	120.77	110.40	13	1
1	A	136	GLU	N-CA-C	5.99	119.85	112.54	20	1
1	A	150	LEU	CB-CA-C	5.97	120.70	110.79	1	4
1	A	124	ASP	N-CA-CB	-5.97	100.40	110.49	7	1
1	A	151	ILE	CA-C-N	5.94	132.89	121.54	17	1
1	A	151	ILE	C-N-CA	5.94	132.89	121.54	17	1
1	A	131	MET	CA-C-O	-5.83	114.24	120.42	10	3
1	A	107	PHE	CA-CB-CG	5.81	119.61	113.80	14	1
1	A	137	ASP	CB-CA-C	5.80	121.96	110.42	2	3
1	A	100	ASN	OD1-CG-ND2	-5.75	116.85	122.60	1	7
1	A	123	THR	CA-CB-OG1	5.75	118.23	109.60	7	1
1	A	151	ILE	CB-CA-C	5.75	119.55	112.02	14	1
1	A	123	THR	N-CA-C	-5.62	99.36	108.41	12	1
1	A	149	ASP	CA-CB-CG	5.55	118.15	112.60	11	2
1	A	100	ASN	CA-C-O	5.55	128.44	120.51	13	1
1	A	127	VAL	CA-CB-CG2	5.50	119.75	110.40	10	1
1	A	97	PHE	N-CA-C	5.49	119.08	112.38	15	2
1	A	104	VAL	CA-CB-CG1	5.44	119.64	110.40	6	2
1	A	136	GLU	CB-CG-CD	5.41	121.80	112.60	7	1
1	A	96	VAL	CA-CB-CG1	5.41	119.59	110.40	5	10
1	A	137	ASP	CA-C-N	5.41	132.01	121.41	13	1
1	A	137	ASP	C-N-CA	5.41	132.01	121.41	13	1
1	A	123	THR	CA-C-N	5.38	131.82	121.54	7	1
1	A	123	THR	C-N-CA	5.38	131.82	121.54	7	1
1	A	93	ALA	N-CA-CB	5.36	117.84	110.07	16	1
1	A	149	ASP	N-CA-CB	5.31	117.85	109.94	17	1
1	A	102	ASP	O-C-N	-5.30	115.85	123.06	2	1
1	A	142	ILE	CB-CA-C	-5.30	104.26	111.15	16	1
1	A	96	VAL	N-CA-C	-5.25	105.60	110.53	5	1
1	A	136	GLU	CA-C-N	5.24	131.55	121.54	11	1
1	A	136	GLU	C-N-CA	5.24	131.55	121.54	11	1
1	A	96	VAL	N-CA-CB	5.22	118.42	110.58	1	1
1	A	101	GLY	N-CA-C	5.22	125.55	113.18	18	3
1	A	126	GLU	CA-CB-CG	5.20	124.50	114.10	11	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	89	GLU	CA-C-O	5.19	127.94	120.51	5	1
1	A	139	ASN	CB-CG-ND2	5.19	124.19	116.40	17	3
1	A	109	GLU	CB-CG-CD	5.09	121.26	112.60	3	1
1	A	90	ILE	CA-C-O	-5.09	114.42	120.78	13	1
1	A	113	ILE	CA-CB-CG2	5.08	119.13	110.50	20	1
1	A	105	ILE	CB-CA-C	-5.07	104.27	111.21	10	1
1	A	105	ILE	N-CA-C	-5.05	100.48	107.80	20	1
1	A	138	GLY	N-CA-C	5.04	125.13	113.18	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	A	102	ASP	Mainchain	4
1	A	147	PHE	Sidechain	2
1	A	137	ASP	Mainchain	2
1	A	92	ARG	Sidechain	2
1	A	123	THR	Mainchain	1
1	A	139	ASN	Mainchain	1
1	A	97	PHE	Sidechain	1
1	A	107	PHE	Sidechain	1
1	A	110	PHE	Sidechain	1
1	A	135	ASP	Mainchain	1

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	444	417	417	3±2
All	All	8880	8340	8339	65

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 4.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:98:ASP:OD2	1:A:102:ASP:HA	1.14	1.41	6	2
1:A:90:ILE:HG23	1:A:148:MET:HE3	0.90	1.39	18	1
1:A:98:ASP:OD1	1:A:104:VAL:O	0.74	2.04	6	1
1:A:90:ILE:HD11	1:A:152:LYS:HE2	0.62	1.70	6	1
1:A:90:ILE:HG23	1:A:147:PHE:HE2	0.62	1.52	19	2
1:A:135:ASP:OD2	1:A:140:GLY:N	0.57	2.37	12	2
1:A:90:ILE:CG2	1:A:148:MET:HE3	0.56	2.22	18	2
1:A:135:ASP:OD2	1:A:140:GLY:HA2	0.56	2.01	1	1
1:A:90:ILE:HG21	1:A:148:MET:HG2	0.55	1.76	10	2
1:A:90:ILE:HG23	1:A:147:PHE:CE2	0.55	2.36	11	12
1:A:90:ILE:HD12	1:A:90:ILE:H	0.54	1.60	6	1
1:A:99:ALA:HB2	1:A:109:GLU:HG2	0.54	1.78	10	2
1:A:90:ILE:HG23	1:A:147:PHE:CE1	0.54	2.38	6	1
1:A:135:ASP:OD2	1:A:140:GLY:CA	0.53	2.56	1	1
1:A:114:MET:HE2	1:A:127:VAL:HG21	0.51	1.83	12	1
1:A:94:PHE:CE2	1:A:105:ILE:HD12	0.50	2.42	13	11
1:A:131:MET:SD	1:A:142:ILE:HD11	0.49	2.47	12	2
1:A:94:PHE:HB2	1:A:147:PHE:CD2	0.49	2.42	5	2
1:A:90:ILE:HG21	1:A:148:MET:HB3	0.49	1.84	6	2
1:A:102:ASP:CG	1:A:144:ILE:HD11	0.48	2.34	13	1
1:A:90:ILE:HD11	1:A:152:LYS:CE	0.48	2.37	6	1
1:A:98:ASP:OD2	1:A:102:ASP:OD1	0.47	2.31	13	1
1:A:149:ASP:O	1:A:152:LYS:HA	0.46	2.10	17	1
1:A:89:GLU:O	1:A:93:ALA:HB3	0.44	2.11	13	1
1:A:131:MET:HE1	1:A:140:GLY:O	0.44	2.12	7	2
1:A:148:MET:N	1:A:148:MET:SD	0.44	2.91	18	1
1:A:111:LYS:O	1:A:114:MET:SD	0.44	2.76	16	1
1:A:90:ILE:HG23	1:A:147:PHE:HE1	0.43	1.70	6	1
1:A:131:MET:HE1	1:A:140:GLY:C	0.43	2.38	17	1
1:A:94:PHE:HB2	1:A:147:PHE:CE2	0.43	2.48	5	2
1:A:90:ILE:HG21	1:A:148:MET:HE3	0.42	1.91	15	1
1:A:98:ASP:C	1:A:100:ASN:H	0.41	2.23	13	1
1:A:99:ALA:CB	1:A:109:GLU:HG2	0.40	2.46	3	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	56/81 (69%)	51±2 (91±4%)	4±2 (6±3%)	2±1 (3±2%)	5	39
All	All	1120/1620 (69%)	1018 (91%)	70 (6%)	32 (3%)	5	39

All 14 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	137	ASP	6
1	A	138	GLY	5
1	A	103	GLY	4
1	A	101	GLY	3
1	A	98	ASP	2
1	A	114	MET	2
1	A	99	ALA	2
1	A	139	ASN	2
1	A	89	GLU	1
1	A	102	ASP	1
1	A	124	ASP	1
1	A	140	GLY	1
1	A	151	ILE	1
1	A	152	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	47/70 (67%)	37±2 (78±5%)	10±2 (22±5%)	2	29
All	All	940/1400 (67%)	731 (78%)	209 (22%)	2	29

All 36 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	113	ILE	18
1	A	142	ILE	16
1	A	148	MET	14
1	A	144	ILE	12

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Mol	Chain	Res	Type	Models (Total)
1	A	91	LEU	12
1	A	150	LEU	12
1	A	104	VAL	11
1	A	141	VAL	11
1	A	102	ASP	8
1	A	96	VAL	8
1	A	132	LYS	7
1	A	139	ASN	7
1	A	90	ILE	7
1	A	149	ASP	6
1	A	127	VAL	6
1	A	152	LYS	5
1	A	128	GLU	5
1	A	98	ASP	5
1	A	126	GLU	5
1	A	114	MET	4
1	A	137	ASP	3
1	A	92	ARG	3
1	A	133	GLU	3
1	A	109	GLU	2
1	A	111	LYS	2
1	A	106	ASP	2
1	A	131	MET	2
1	A	97	PHE	2
1	A	146	GLU	2
1	A	124	ASP	2
1	A	151	ILE	2
1	A	95	LYS	1
1	A	123	THR	1
1	A	89	GLU	1
1	A	107	PHE	1
1	A	147	PHE	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