



Full wwPDB NMR Structure Validation Report ⓘ

Mar 8, 2026 – 08:13 AM UTC

PDB ID : 2KLR / pdb_00002klr
Title : Solid-state NMR structure of the alpha-crystallin domain in alphaB-crystallin oligomers
Authors : Jehle, S.; Rajagopal, P.; Markovic, S.; Bardiaux, B.; Kuehne, R.; Higman, V.A.; Klevit, R.E.; van Rossum, B.; Oschkinat, H.
Deposited on : 2009-07-08

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

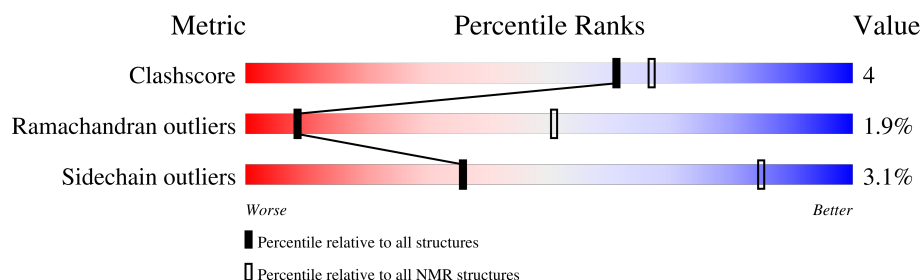
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLID-STATE 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	175	
1	B	175	

2 Ensemble composition and analysis

This entry contains 10 models. Model 10 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:73-A:150, B:73-B:150 (156)	1.38	10

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 and 2 single-model clusters were found.

Cluster number	Models
1	2, 4, 8, 10
2	1, 3, 6, 9
Single-model clusters	5; 7

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 2670 atoms, of which 1338 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Alpha-crystallin B chain.

Mol	Chain	Residues	Atoms					Trace
1	A	82	Total	C	H	N	O	0
			1335	417	669	123	126	
1	B	82	Total	C	H	N	O	0
			1335	417	669	123	126	

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: Alpha-crystallin B chain



- Molecule 1: Alpha-crystallin B chain

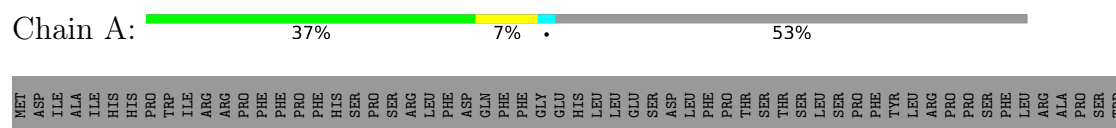


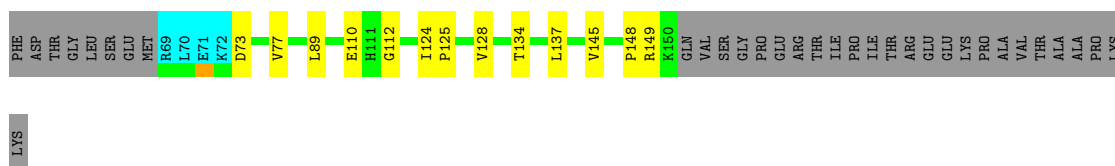
4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

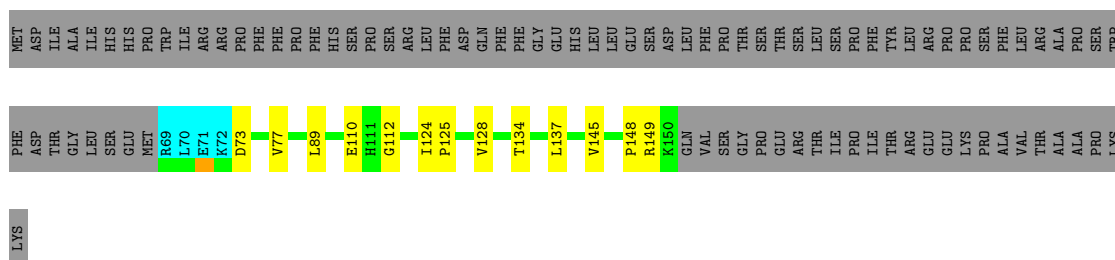
- Molecule 1: Alpha-crystallin B chain





- Molecule 1: Alpha-crystallin B chain

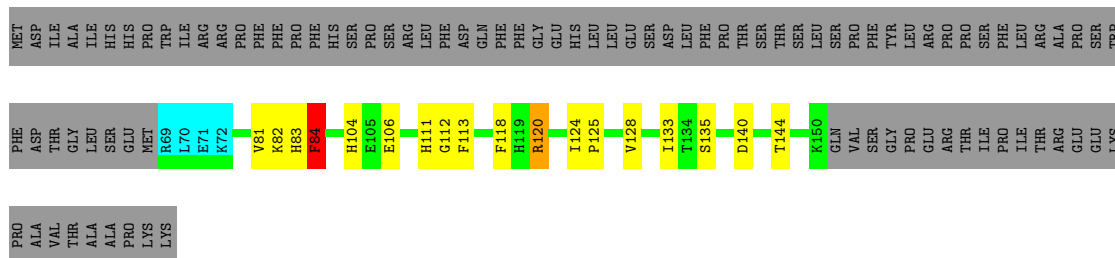
Chain B: 37% 7% 53%



4.2.2 Score per residue for model 2

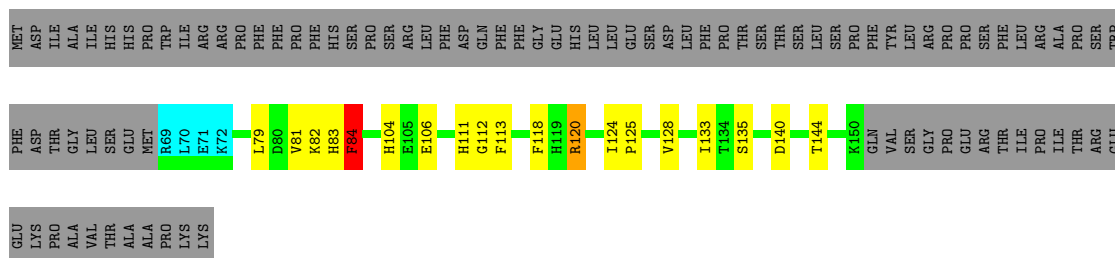
- Molecule 1: Alpha-crystallin B chain

Chain A: 34% 9% 53%



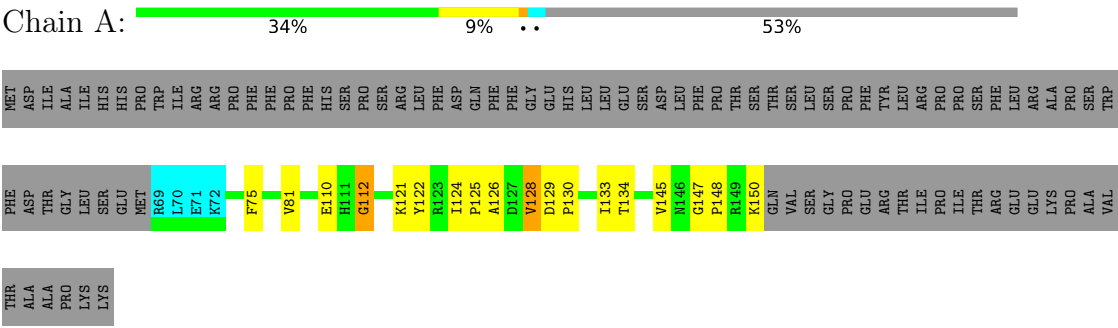
- Molecule 1: Alpha-crystallin B chain

Chain B: 34% 10% 53%

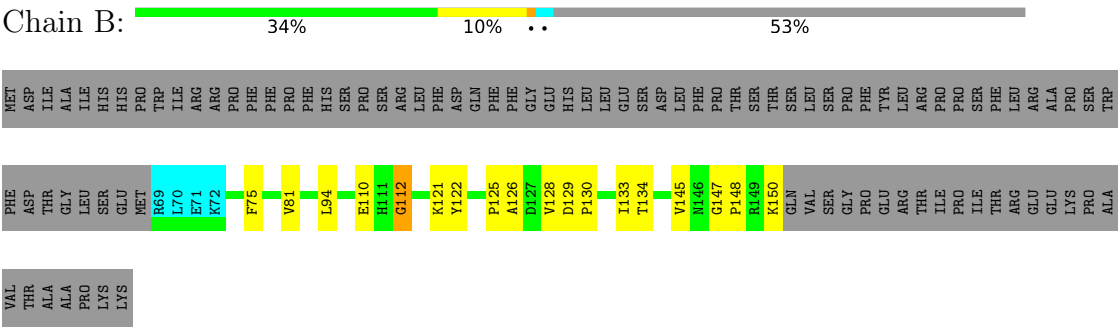


4.2.3 Score per residue for model 3

- Molecule 1: Alpha-crystallin B chain

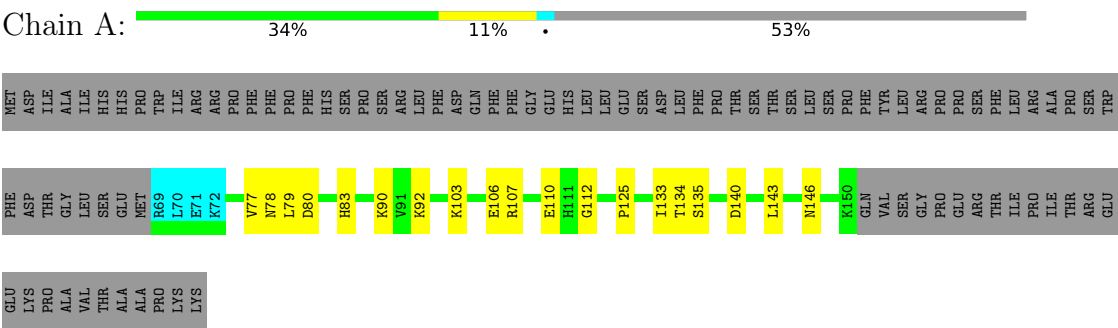


• Molecule 1: Alpha-crystallin B chain

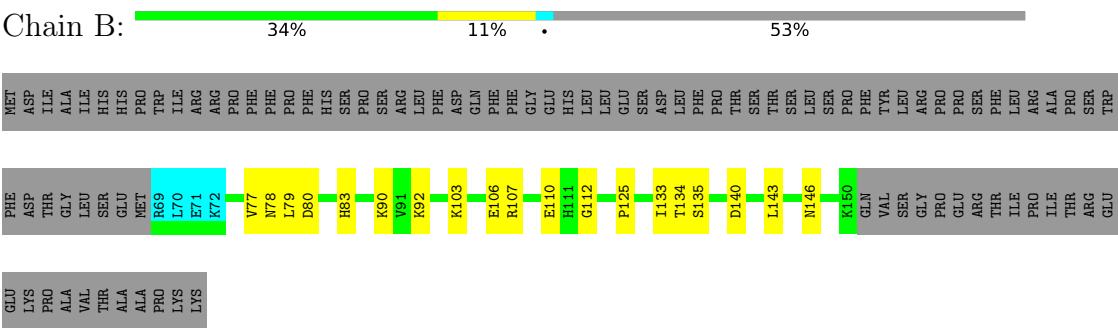


4.2.4 Score per residue for model 4

• Molecule 1: Alpha-crystallin B chain

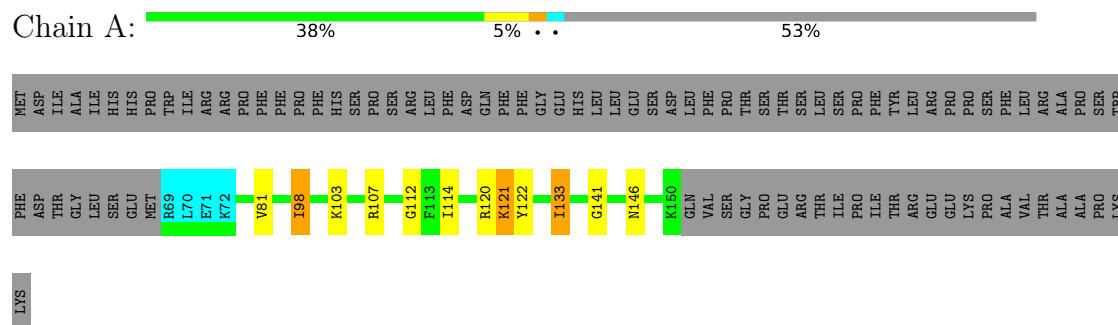


• Molecule 1: Alpha-crystallin B chain

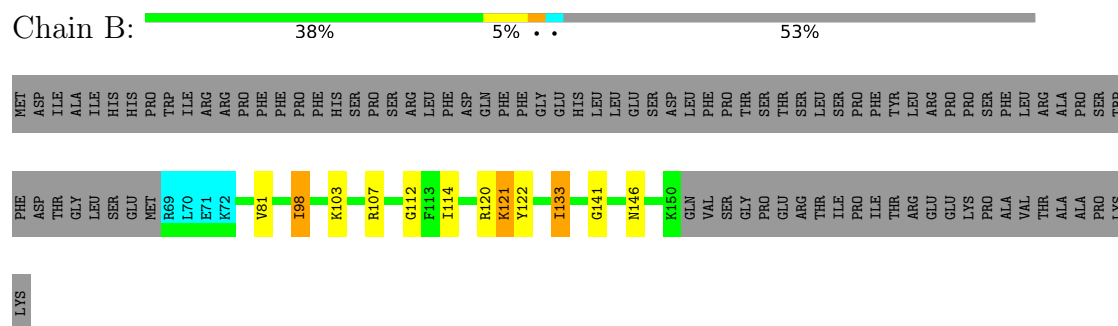


4.2.5 Score per residue for model 5

- Molecule 1: Alpha-crystallin B chain

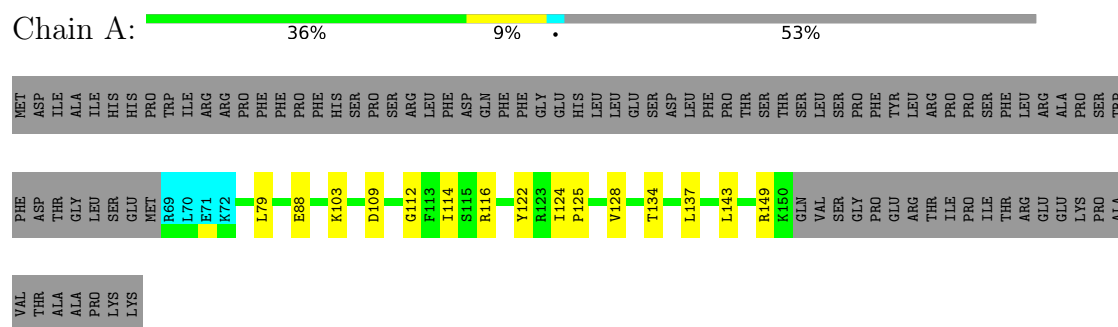


- Molecule 1: Alpha-crystallin B chain

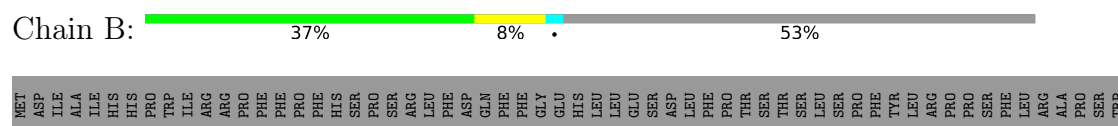


4.2.6 Score per residue for model 6

- Molecule 1: Alpha-crystallin B chain



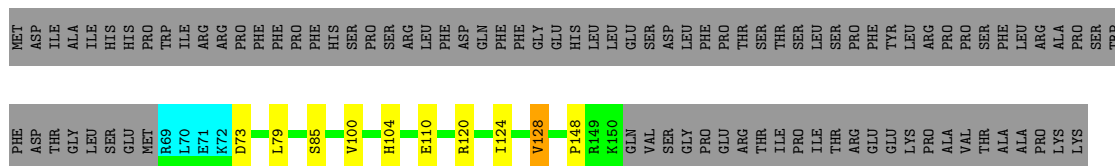
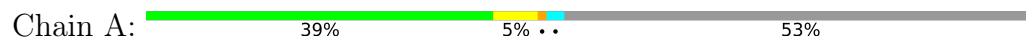
- Molecule 1: Alpha-crystallin B chain



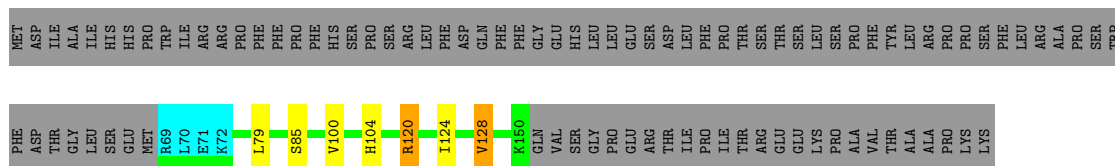


4.2.7 Score per residue for model 7

- Molecule 1: Alpha-crystallin B chain

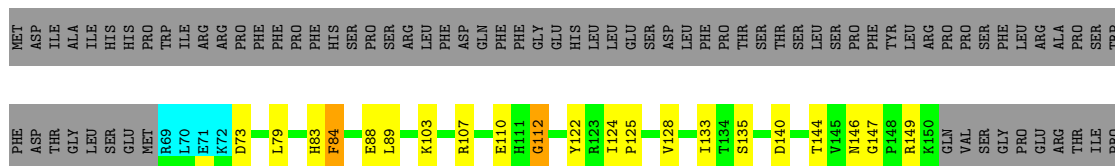
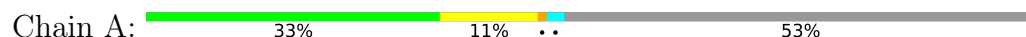


- Molecule 1: Alpha-crystallin B chain

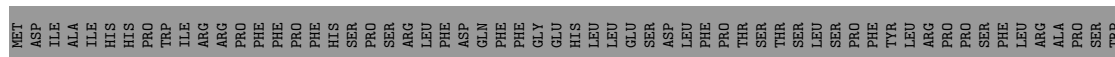
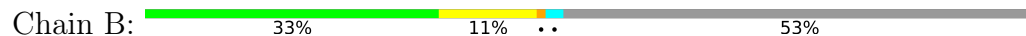


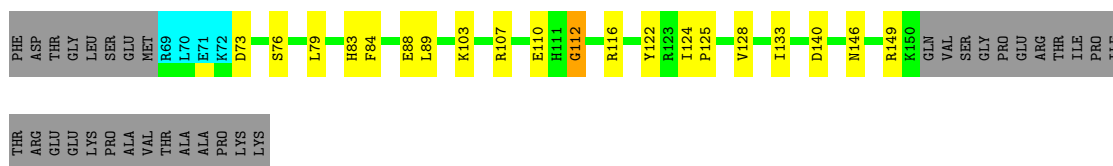
4.2.8 Score per residue for model 8

- Molecule 1: Alpha-crystallin B chain



- Molecule 1: Alpha-crystallin B chain

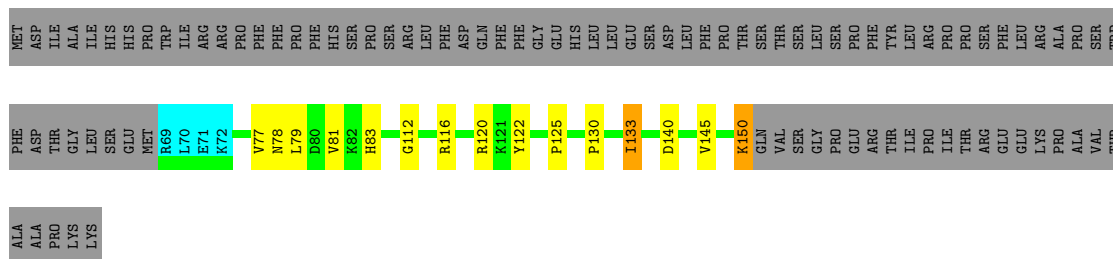




4.2.9 Score per residue for model 9

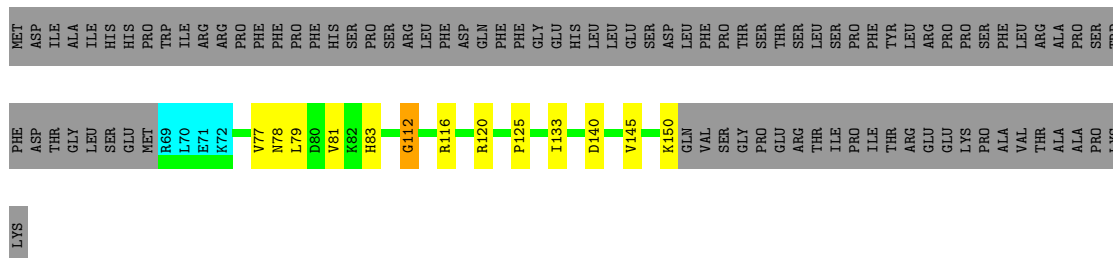
- Molecule 1: Alpha-crystallin B chain

Chain A: 36% 7% 53%



- Molecule 1: Alpha-crystallin B chain

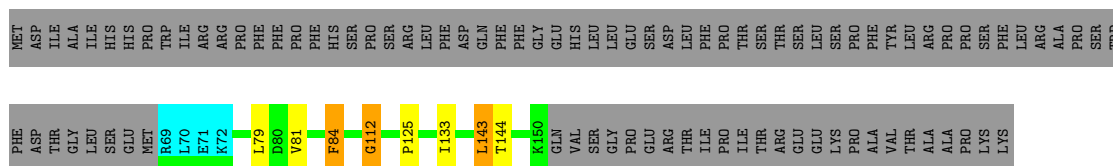
Chain B: 37% 7% 53%



4.2.10 Score per residue for model 10 (medoid)

- Molecule 1: Alpha-crystallin B chain

Chain A: 40% 5% 53%



- Molecule 1: Alpha-crystallin B chain

Chain B: 39% 5% 53%

MET	ASP	ILE	ALA	ILE	HIS	HIS	PRO	TRP	ILE	ARG	ARG	PRO	PHE	PHE	PHE	HIS	SER	PRO	SER	ARG	LEU	PHE	ASP	GLN	PHE	PHE	GLY	GLU	HIS	LEU	LEU	GLU	SER	SER	ASP	LEU	PHE	PRO	THR	SER	THR	SER	LEU	SER	PRO	PHE	TYR	LEU	ARG	PRO	PRO	SER	PHE	LEU	ARG	ALA	PRO	SER	TRP
PHE	ASP	THR	GLY	LEU	SER	GLU	MET	R69	L70	E71	K72	V77	N78	L79	D80	V81	F84	G112	K121	P125	I133	L143	T144	K150	GLN	VAL	SER	GLY	PRO	GLU	ARG	THR	THR	ILE	PRO	ILE	THR	ARG	GLU	GLU	LYS	PRO	ALA	VAL	THR	ALA	ALA	PRO	LYS	LYS									

5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing, molecular dynamics, torsion angle dynamics*.

Of the 200 calculated structures, 10 were deposited, based on the following criterion: *structures with the lowest energy*.

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

Software name	Classification	Version
X-PLOR NIH	structure solution	
ARIA	structure solution	2.2
ARIA	refinement	2.2
CNS	structure solution	1.2
CNS	refinement	1.2
SOLARIA	structure solution	1

No chemical shift data was provided. Note: This is a solid-state NMR structure, where hydrogen atoms are typically not assigned a chemical shift value, which may lead to lower completeness of assignment measure.

6 Model quality (i)

6.1 Standard geometry (i)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	1.41±0.07	3±2/642 (0.4± 0.4%)	1.01±0.04	0±0/865 (0.1± 0.1%)
1	B	1.41±0.06	3±2/642 (0.5± 0.4%)	1.00±0.05	0±0/865 (0.0± 0.1%)
All	All	1.41	56/12840 (0.4%)	1.01	9/17300 (0.1%)

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	0.3±0.5
1	B	0.0±0.0	0.4±0.7
All	All	0	7

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	B	80	ASP	C-N	-8.70	1.25	1.33	4	1
1	A	80	ASP	C-N	-8.56	1.26	1.33	4	1
1	A	112	GLY	N-CA	-7.61	1.37	1.45	2	2
1	B	112	GLY	N-CA	-7.57	1.37	1.45	2	2
1	B	112	GLY	CA-C	-7.31	1.45	1.51	2	5
1	A	112	GLY	CA-C	-7.24	1.45	1.51	2	4
1	B	79	LEU	N-CA	-6.40	1.38	1.46	4	4
1	A	112	GLY	C-N	-6.37	1.24	1.33	1	6
1	A	79	LEU	N-CA	-6.26	1.38	1.46	4	3
1	B	112	GLY	C-N	-6.21	1.24	1.33	2	6
1	A	78	ASN	CA-C	-5.79	1.45	1.52	9	1
1	A	98	ILE	CA-CB	5.69	1.61	1.54	5	1
1	B	98	ILE	CA-CB	5.67	1.61	1.54	5	1
1	A	111	HIS	C-N	-5.62	1.26	1.33	2	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	B	78	ASN	CA-C	-5.61	1.46	1.52	9	1
1	A	78	ASN	C-N	-5.52	1.26	1.33	9	2
1	B	78	ASN	C-N	-5.49	1.26	1.33	4	2
1	B	113	PHE	N-CA	-5.47	1.39	1.45	2	1
1	A	113	PHE	N-CA	-5.40	1.39	1.45	2	1
1	B	111	HIS	C-N	-5.38	1.26	1.33	2	1
1	B	75	PHE	N-CA	-5.38	1.39	1.45	3	1
1	B	77	VAL	C-N	-5.35	1.26	1.33	4	2
1	A	77	VAL	C-N	-5.32	1.26	1.33	4	1
1	B	116	ARG	C-N	-5.22	1.26	1.33	9	1
1	A	75	PHE	N-CA	-5.17	1.39	1.45	3	1
1	B	120	ARG	CA-C	-5.11	1.46	1.52	7	1
1	A	116	ARG	C-N	-5.10	1.26	1.33	9	1
1	B	76	SER	N-CA	-5.09	1.40	1.46	8	1
1	A	116	ARG	N-CA	-5.07	1.39	1.45	9	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	128	VAL	N-CA-C	5.36	115.83	108.12	3	2
1	B	128	VAL	N-CA-C	5.34	115.82	108.12	3	2
1	B	84	PHE	CA-CB-CG	5.23	119.03	113.80	2	2
1	A	84	PHE	CA-CB-CG	5.23	119.03	113.80	2	3

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	116	ARG	Sidechain	2
1	A	120	ARG	Sidechain	1
1	B	120	ARG	Sidechain	1
1	A	116	ARG	Sidechain	1
1	A	107	ARG	Sidechain	1
1	B	107	ARG	Sidechain	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	629	626	621	6±2
1	B	629	626	621	6±2
All	All	12580	12520	12420	110

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:B:83:HIS:HA	1:B:140:ASP:O	0.66	1.91	8	4
1:A:83:HIS:HA	1:A:140:ASP:O	0.64	1.91	8	4
1:A:122:TYR:CD2	1:B:112:GLY:HA3	0.59	2.33	3	3
1:B:98:ILE:HG23	1:B:122:TYR:HB2	0.58	1.75	5	1
1:A:98:ILE:HG23	1:A:122:TYR:HB2	0.58	1.75	5	1
1:A:73:ASP:O	1:A:148:PRO:HA	0.56	2.00	1	2
1:A:129:ASP:H	1:A:148:PRO:HB3	0.56	1.61	3	1
1:B:129:ASP:H	1:B:148:PRO:HB3	0.55	1.61	3	1
1:B:73:ASP:O	1:B:148:PRO:HA	0.55	2.01	1	1
1:B:79:LEU:HD22	1:B:122:TYR:CE1	0.55	2.37	6	1
1:A:112:GLY:HA3	1:B:122:TYR:CD2	0.54	2.38	3	2
1:A:81:VAL:HG12	1:A:82:LYS:O	0.54	2.02	2	1
1:A:79:LEU:HD22	1:A:122:TYR:CZ	0.54	2.38	8	1
1:B:81:VAL:HG12	1:B:82:LYS:O	0.54	2.02	2	1
1:B:133:ILE:HA	1:B:146:ASN:O	0.54	2.03	8	3
1:B:81:VAL:HB	1:B:141:GLY:O	0.54	2.02	5	1
1:A:133:ILE:HA	1:A:146:ASN:O	0.53	2.02	8	3
1:B:77:VAL:HB	1:B:145:VAL:CG2	0.53	2.33	1	2
1:A:79:LEU:HD22	1:A:122:TYR:CE1	0.53	2.37	6	1
1:A:112:GLY:HA2	1:B:121:LYS:O	0.53	2.04	3	2
1:B:79:LEU:HD22	1:B:122:TYR:CZ	0.53	2.38	8	1
1:A:77:VAL:HB	1:A:145:VAL:CG2	0.52	2.33	1	2
1:B:124:ILE:O	1:B:128:VAL:HB	0.52	2.04	2	5
1:A:81:VAL:HB	1:A:141:GLY:O	0.52	2.03	5	1
1:A:124:ILE:O	1:A:128:VAL:HB	0.52	2.03	2	6
1:A:121:LYS:O	1:B:112:GLY:HA2	0.52	2.03	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:106:GLU:HB3	1:A:107:ARG:NH1	0.50	2.21	4	1
1:A:81:VAL:HA	1:A:120:ARG:CZ	0.50	2.37	2	1
1:B:106:GLU:HB3	1:B:107:ARG:NH1	0.50	2.22	4	1
1:B:81:VAL:HA	1:B:120:ARG:CZ	0.49	2.38	2	1
1:A:79:LEU:HD13	1:A:100:VAL:HB	0.48	1.85	7	1
1:B:107:ARG:HG2	1:B:114:ILE:HD11	0.48	1.85	5	1
1:A:110:GLU:CD	1:A:110:GLU:H	0.48	2.16	4	1
1:A:79:LEU:HA	1:A:120:ARG:HD2	0.48	1.86	7	1
1:B:110:GLU:H	1:B:110:GLU:CD	0.48	2.16	4	1
1:A:84:PHE:HA	1:A:104:HIS:HB2	0.47	1.86	2	1
1:B:79:LEU:HA	1:B:120:ARG:HD2	0.47	1.86	7	1
1:B:79:LEU:HD13	1:B:100:VAL:HB	0.47	1.86	7	1
1:A:107:ARG:HG2	1:A:114:ILE:HD11	0.47	1.85	5	1
1:B:84:PHE:HA	1:B:104:HIS:HB2	0.47	1.86	2	1
1:A:143:LEU:HD12	1:A:144:THR:N	0.47	2.24	10	1
1:B:143:LEU:HD12	1:B:144:THR:N	0.46	2.25	10	1
1:A:126:ALA:O	1:A:150:LYS:HA	0.44	2.13	3	1
1:A:147:GLY:N	1:A:148:PRO:HD2	0.44	2.28	3	1
1:B:126:ALA:O	1:B:150:LYS:HA	0.44	2.12	3	1
1:A:130:PRO:HA	1:A:133:ILE:CG2	0.44	2.43	3	1
1:A:128:VAL:HG13	1:A:133:ILE:CD1	0.44	2.43	2	1
1:B:106:GLU:HB3	1:B:107:ARG:CZ	0.44	2.42	4	1
1:B:128:VAL:HG13	1:B:133:ILE:CD1	0.43	2.43	2	1
1:A:106:GLU:HB3	1:A:107:ARG:CZ	0.43	2.43	4	1
1:A:81:VAL:HB	1:A:120:ARG:NH1	0.43	2.29	9	1
1:B:81:VAL:HB	1:B:120:ARG:NH1	0.43	2.28	9	1
1:B:135:SER:HA	1:B:144:THR:O	0.43	2.13	2	1
1:B:147:GLY:N	1:B:148:PRO:HD2	0.43	2.28	3	1
1:A:135:SER:HA	1:A:144:THR:O	0.43	2.14	2	2
1:B:134:THR:O	1:B:145:VAL:HA	0.42	2.14	3	2
1:B:88:GLU:HB3	1:B:103:LYS:O	0.42	2.15	6	1
1:B:130:PRO:HA	1:B:133:ILE:CG2	0.42	2.44	3	1
1:A:88:GLU:HB3	1:A:103:LYS:O	0.42	2.15	6	1
1:A:106:GLU:C	1:A:107:ARG:HD3	0.42	2.39	4	1
1:A:120:ARG:NH2	1:B:114:ILE:HD12	0.42	2.30	5	1
1:A:134:THR:O	1:A:145:VAL:HA	0.42	2.15	3	2
1:B:106:GLU:C	1:B:107:ARG:HD3	0.42	2.40	4	1
1:A:88:GLU:HB3	1:A:103:LYS:HG3	0.42	1.91	8	1
1:B:88:GLU:HB3	1:B:103:LYS:HG3	0.42	1.91	8	1
1:B:135:SER:HB2	1:B:143:LEU:HD11	0.41	1.92	4	1
1:A:121:LYS:HD2	1:A:121:LYS:O	0.41	2.15	5	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:121:LYS:O	1:B:121:LYS:HD2	0.41	2.15	5	1
1:B:77:VAL:HB	1:B:145:VAL:HG23	0.41	1.91	1	1
1:A:114:ILE:HD12	1:B:120:ARG:NH2	0.41	2.31	5	1
1:B:84:PHE:HB3	1:B:118:PHE:CE1	0.41	2.51	2	1
1:A:150:LYS:NZ	1:A:150:LYS:HB2	0.41	2.30	9	1
1:A:84:PHE:HB3	1:A:118:PHE:CE1	0.41	2.51	2	1
1:B:85:SER:OG	1:B:104:HIS:HB2	0.41	2.16	7	1
1:A:120:ARG:HG2	1:B:114:ILE:CB	0.40	2.47	5	1
1:A:135:SER:HB2	1:A:143:LEU:HD11	0.40	1.93	4	1
1:B:109:ASP:HB3	1:B:112:GLY:O	0.40	2.16	6	1
1:A:85:SER:OG	1:A:104:HIS:HB2	0.40	2.16	7	1
1:A:109:ASP:HB3	1:A:112:GLY:O	0.40	2.16	6	1
1:A:133:ILE:HG12	1:A:147:GLY:HA3	0.40	1.94	8	1
1:A:130:PRO:HA	1:A:133:ILE:CG1	0.40	2.47	9	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	77/175 (44%)	70±1 (91±2%)	5±2 (7±2%)	2±1 (2±1%)	8	51
1	B	77/175 (44%)	70±1 (91±2%)	5±2 (7±2%)	1±1 (2±1%)	9	52
All	All	1540/3500 (44%)	1407 (91%)	104 (7%)	29 (2%)	8	51

All 10 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	125	PRO	8
1	B	125	PRO	8
1	A	110	GLU	4
1	B	110	GLU	3
1	A	106	GLU	1
1	B	106	GLU	1
1	A	73	ASP	1

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Mol	Chain	Res	Type	Models (Total)
1	B	73	ASP	1
1	A	81	VAL	1
1	B	81	VAL	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	72/161 (45%)	70±1 (97±2%)	2±1 (3±2%)	36	85
1	B	72/161 (45%)	70±1 (97±1%)	2±1 (3±1%)	36	85
All	All	1440/3220 (45%)	1396 (97%)	44 (3%)	36	85

All 26 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	84	PHE	3
1	B	84	PHE	3
1	A	133	ILE	3
1	B	133	ILE	3
1	A	89	LEU	2
1	A	137	LEU	2
1	B	89	LEU	2
1	B	137	LEU	2
1	A	103	LYS	2
1	A	134	THR	2
1	B	103	LYS	2
1	B	134	THR	2
1	A	143	LEU	2
1	B	143	LEU	2
1	A	81	VAL	1
1	B	81	VAL	1
1	B	94	LEU	1
1	A	90	LYS	1
1	A	92	LYS	1
1	B	90	LYS	1
1	B	92	LYS	1

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Mol	Chain	Res	Type	Models (Total)
1	A	121	LYS	1
1	B	121	LYS	1
1	A	114	ILE	1
1	A	150	LYS	1
1	B	150	LYS	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