Full wwPDB NMR Structure Validation Report

Mar 5, 2026 – 04:30 PM UTC

PDB ID	:	1ZV6 / pdb_00001zv6
Title	:	NMR structure of the human dematin headpiece S74E mutant
Authors	:	Jiang, Z.G.; McKnight, C.J.
Deposited on	:	2005-06-01

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:

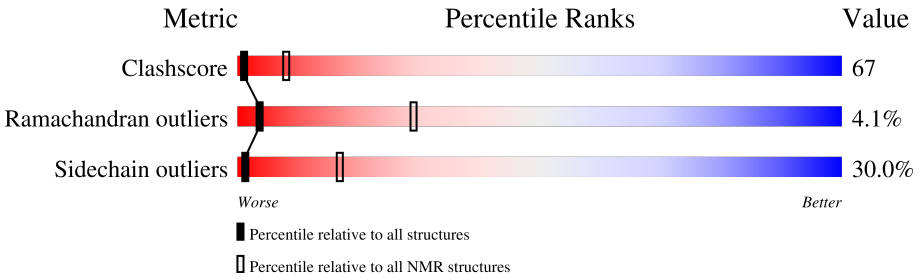
MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 <=5%

Mol	Chain	Length	Quality of chain
1	A	68	13% 66% 10% 10%

2 Ensemble composition and analysis (i)

This entry contains 21 models. Model 8 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *minimized average structure*.

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:11-A:23, A:29-A:76 (61)	0.40	8

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 1 single-model cluster was found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1143 atoms, of which 582 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called EPB49 protein.

Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	68	1143	362	582	99	97	3	0

There is a discrepancy between the modelled and reference sequences:

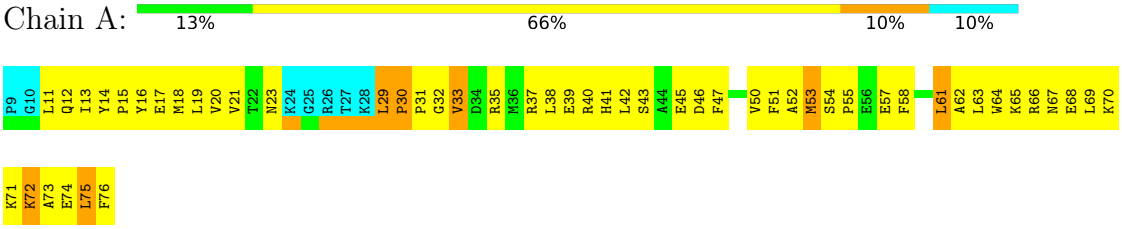
Chain	Residue	Modelled	Actual	Comment	Reference
A	74	GLU	SER	engineered mutation	GB 31127246

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: EPB49 protein

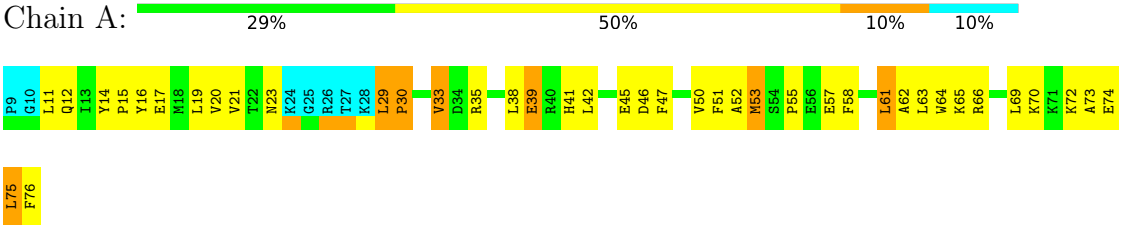


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

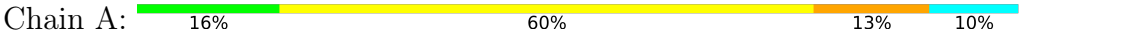
4.2.1 Score per residue for model 1

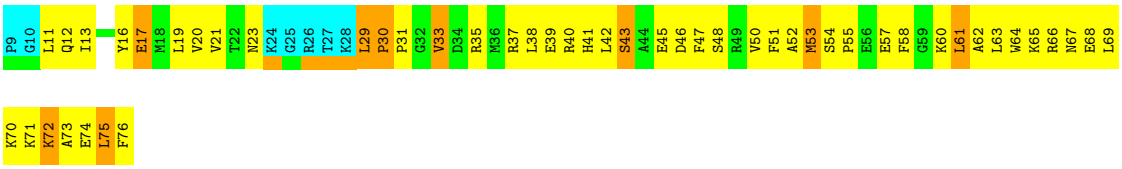
• Molecule 1: EPB49 protein



4.2.2 Score per residue for model 2

• Molecule 1: EPB49 protein

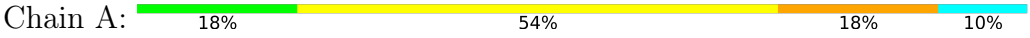


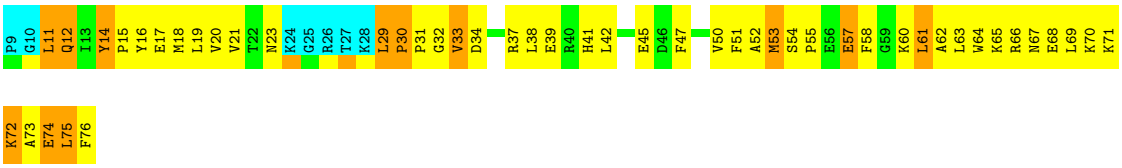


4.2.3 Score per residue for model 3

• Molecule 1: EPB49 protein

Chain A:

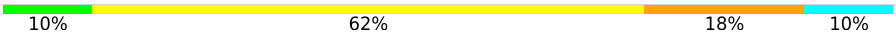


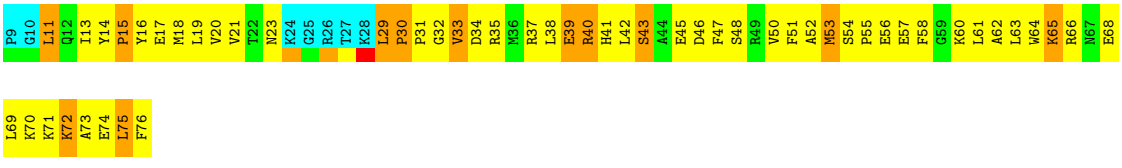


4.2.4 Score per residue for model 4

• Molecule 1: EPB49 protein

Chain A:

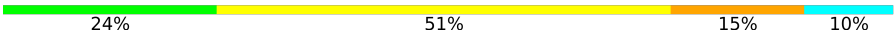


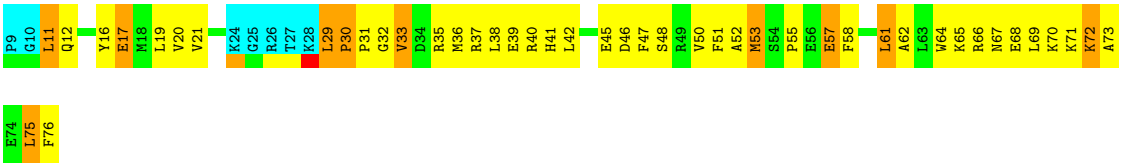


4.2.5 Score per residue for model 5

• Molecule 1: EPB49 protein

Chain A:

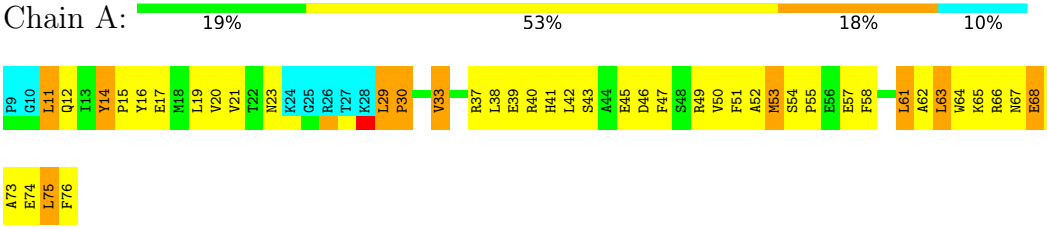




4.2.6 Score per residue for model 6

• Molecule 1: EPB49 protein

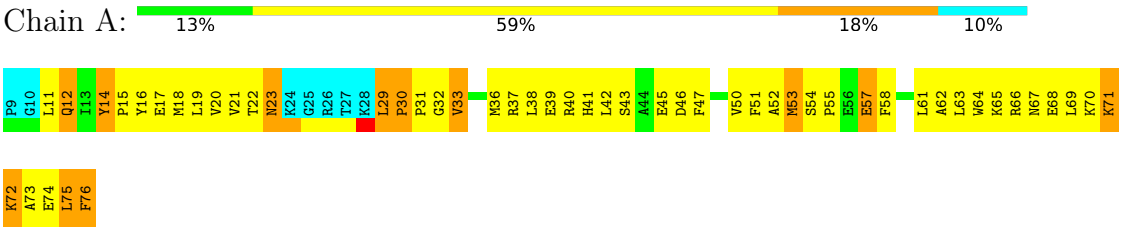
Chain A:



4.2.7 Score per residue for model 7

• Molecule 1: EPB49 protein

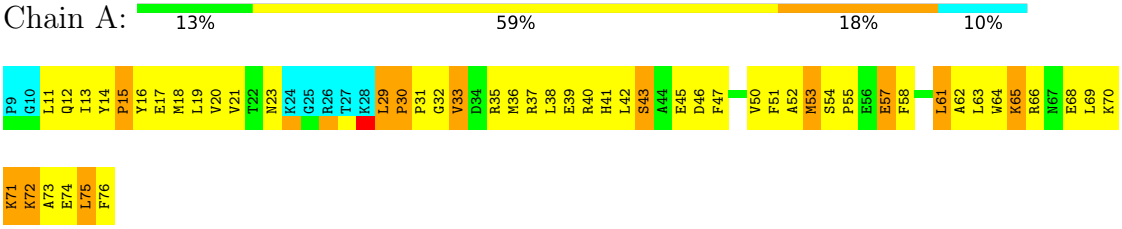
Chain A:



4.2.8 Score per residue for model 8 (medoid)

• Molecule 1: EPB49 protein

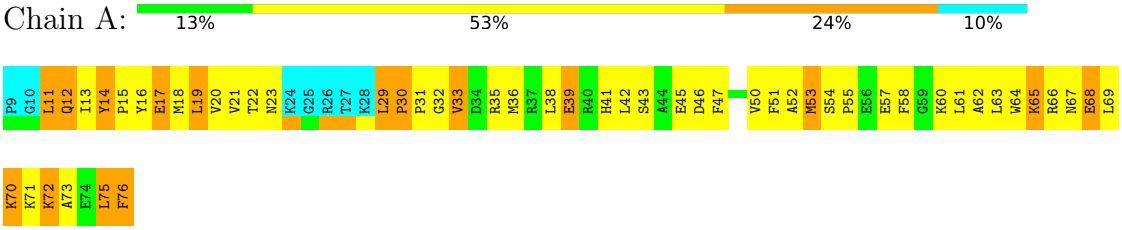
Chain A:



4.2.9 Score per residue for model 9

• Molecule 1: EPB49 protein

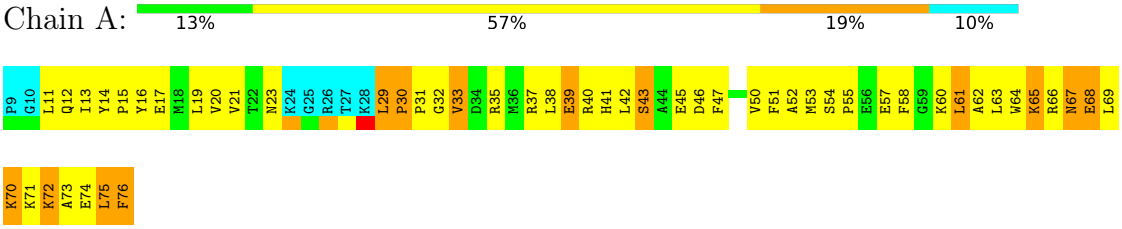
Chain A:



4.2.10 Score per residue for model 10

• Molecule 1: EPB49 protein

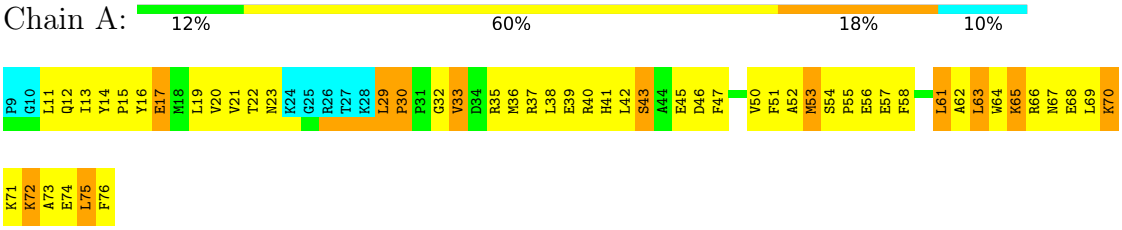
Chain A:



4.2.11 Score per residue for model 11

• Molecule 1: EPB49 protein

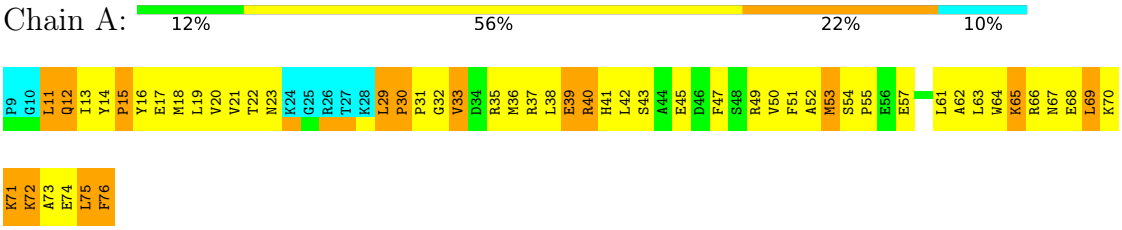
Chain A:



4.2.12 Score per residue for model 12

• Molecule 1: EPB49 protein

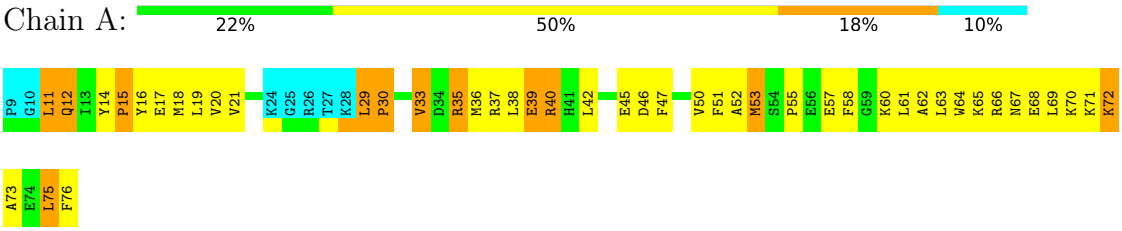
Chain A:



4.2.13 Score per residue for model 13

• Molecule 1: EPB49 protein

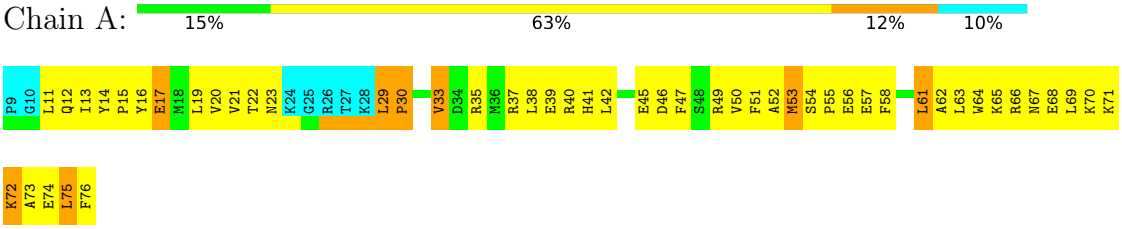
Chain A:



4.2.14 Score per residue for model 14

• Molecule 1: EPB49 protein

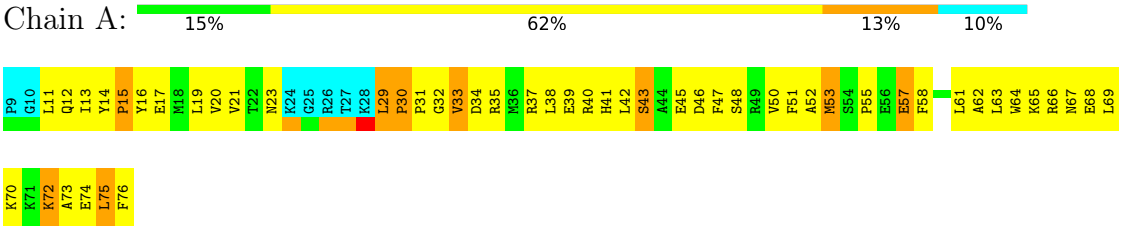
Chain A:



4.2.15 Score per residue for model 15

• Molecule 1: EPB49 protein

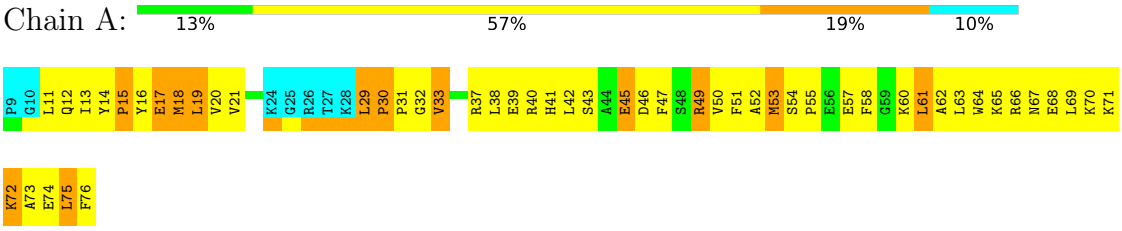
Chain A:



4.2.16 Score per residue for model 16

• Molecule 1: EPB49 protein

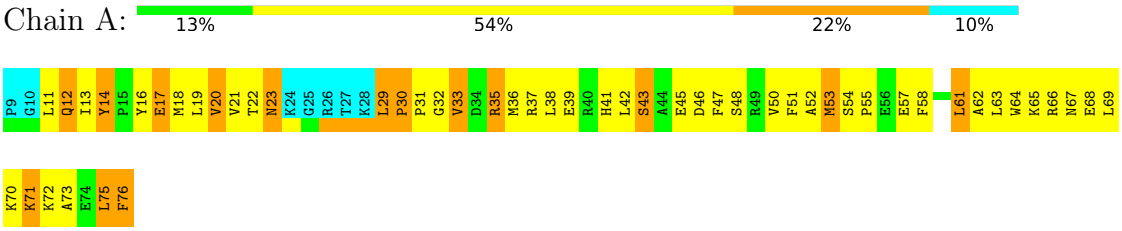
Chain A:



4.2.17 Score per residue for model 17

• Molecule 1: EPB49 protein

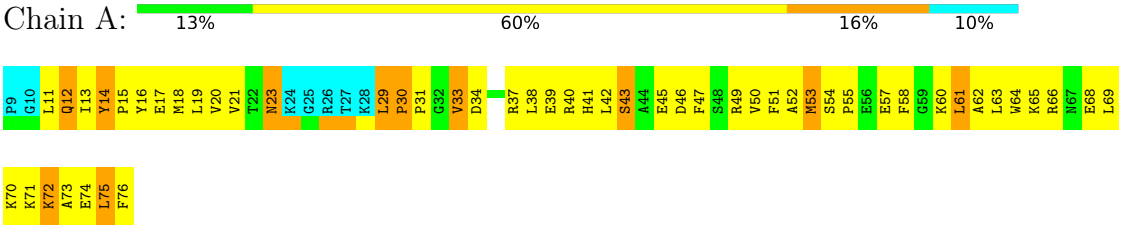
Chain A:



4.2.18 Score per residue for model 18

• Molecule 1: EPB49 protein

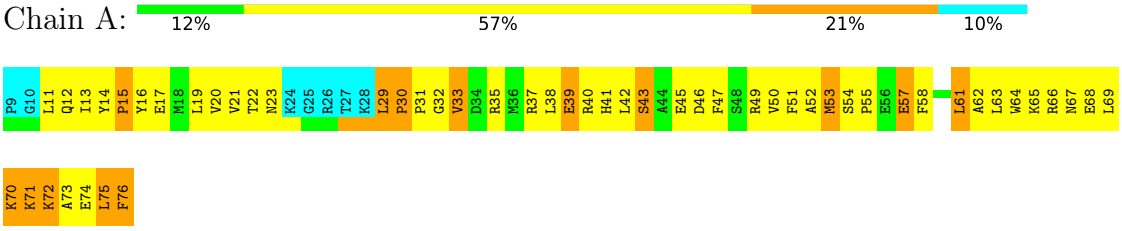
Chain A:



4.2.19 Score per residue for model 19

• Molecule 1: EPB49 protein

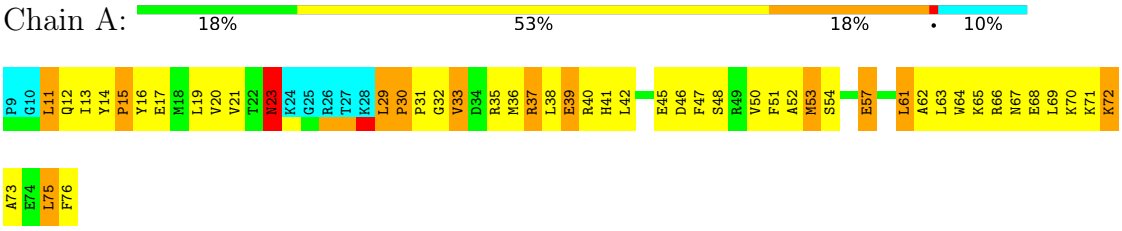
Chain A:



4.2.20 Score per residue for model 20

• Molecule 1: EPB49 protein

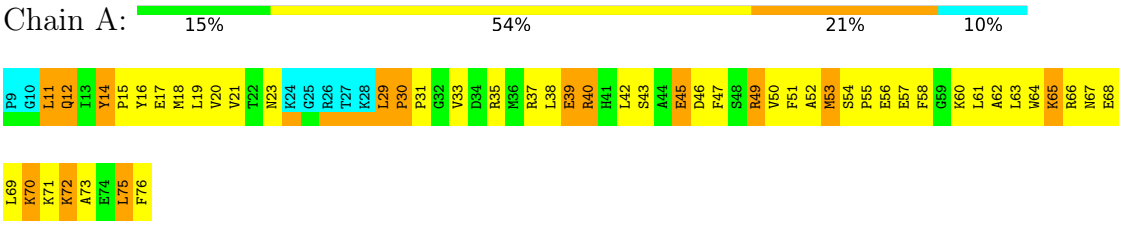
Chain A:



4.2.21 Score per residue for model 21

• Molecule 1: EPB49 protein

Chain A:



5 Refinement protocol and experimental data overview

The models were refined using the following method: *DISTANCE GEOMETRY/SIMULATED ANNEALING*.

Of the 100 calculated structures, 21 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
MOLMOL	structure solution	2K.2
CNS	refinement	1.0

No chemical shift data was provided.

6 Model quality

6.1 Standard geometry

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

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	510	521	520	69±8
All	All	10710	10941	10920	1443

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 67.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:19:LEU:HD22	1:A:33:VAL:HG11	0.95	1.31	12	16
1:A:53:MET:HE1	1:A:69:LEU:HD13	0.90	1.42	15	2
1:A:29:LEU:HD13	1:A:38:LEU:HD21	0.88	1.45	13	4
1:A:29:LEU:HD23	1:A:33:VAL:O	0.87	1.70	16	5
1:A:61:LEU:HD13	1:A:65:LYS:CD	0.85	2.01	12	3
1:A:30:PRO:O	1:A:33:VAL:HG23	0.85	1.71	21	14
1:A:21:VAL:HG22	1:A:38:LEU:HD12	0.84	1.48	12	9
1:A:19:LEU:CD2	1:A:33:VAL:HG11	0.83	2.03	15	16
1:A:29:LEU:HD21	1:A:38:LEU:HD11	0.83	1.51	2	2
1:A:29:LEU:HD23	1:A:29:LEU:N	0.82	1.88	21	5
1:A:16:TYR:CE1	1:A:75:LEU:HD21	0.82	2.09	11	20
1:A:16:TYR:CD1	1:A:50:VAL:HG21	0.81	2.10	21	18
1:A:29:LEU:HD12	1:A:33:VAL:HB	0.81	1.50	13	4
1:A:29:LEU:HD12	1:A:33:VAL:CG1	0.81	2.05	14	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:29:LEU:HD21	1:A:38:LEU:HD21	0.80	1.54	18	2
1:A:29:LEU:H	1:A:29:LEU:HD23	0.79	1.37	19	3
1:A:29:LEU:CD1	1:A:38:LEU:HD21	0.79	2.07	11	10
1:A:33:VAL:HG12	1:A:38:LEU:HD23	0.78	1.56	9	10
1:A:36:MET:HE2	1:A:37:ARG:CG	0.78	2.07	20	1
1:A:21:VAL:HG13	1:A:35:ARG:HG2	0.78	1.55	19	4
1:A:66:ARG:HA	1:A:69:LEU:HD12	0.77	1.55	11	21
1:A:14:TYR:CE1	1:A:33:VAL:HG22	0.77	2.14	9	4
1:A:30:PRO:HD2	1:A:33:VAL:HG21	0.77	1.57	10	12
1:A:19:LEU:O	1:A:38:LEU:HD13	0.77	1.80	12	8
1:A:29:LEU:HD23	1:A:33:VAL:C	0.77	2.05	15	3
1:A:46:ASP:O	1:A:50:VAL:HG23	0.76	1.79	17	19
1:A:29:LEU:HD12	1:A:33:VAL:CB	0.76	2.11	13	4
1:A:17:GLU:O	1:A:20:VAL:HG23	0.74	1.83	8	13
1:A:29:LEU:CD2	1:A:38:LEU:HD21	0.74	2.12	18	2
1:A:21:VAL:HG13	1:A:35:ARG:HD2	0.73	1.60	2	1
1:A:29:LEU:HG	1:A:38:LEU:HD21	0.72	1.59	12	5
1:A:29:LEU:H	1:A:29:LEU:HD22	0.72	1.45	9	3
1:A:33:VAL:HG12	1:A:38:LEU:CD2	0.72	2.15	9	4
1:A:29:LEU:HB2	1:A:33:VAL:HB	0.72	1.62	13	7
1:A:29:LEU:HD12	1:A:33:VAL:HG11	0.72	1.62	21	2
1:A:19:LEU:HD23	1:A:19:LEU:N	0.71	2.01	16	2
1:A:29:LEU:HD23	1:A:29:LEU:H	0.71	1.44	21	2
1:A:21:VAL:HG22	1:A:38:LEU:HD13	0.70	1.62	21	3
1:A:33:VAL:HG13	1:A:41:HIS:CG	0.70	2.21	15	10
1:A:29:LEU:HB2	1:A:33:VAL:HG21	0.69	1.62	21	1
1:A:53:MET:SD	1:A:61:LEU:HD11	0.69	2.28	10	15
1:A:29:LEU:CD1	1:A:33:VAL:HB	0.69	2.17	13	4
1:A:33:VAL:HG13	1:A:41:HIS:CD2	0.68	2.23	3	7
1:A:29:LEU:HD11	1:A:38:LEU:HD21	0.67	1.65	10	7
1:A:39:GLU:HA	1:A:42:LEU:HD12	0.67	1.65	2	21
1:A:11:LEU:HD13	1:A:11:LEU:N	0.67	2.05	3	3
1:A:16:TYR:CE2	1:A:75:LEU:HD21	0.67	2.25	20	1
1:A:16:TYR:CD2	1:A:50:VAL:HG21	0.67	2.25	20	1
1:A:61:LEU:HD13	1:A:65:LYS:HD3	0.67	1.66	12	1
1:A:29:LEU:CB	1:A:33:VAL:HB	0.66	2.21	13	11
1:A:61:LEU:HD13	1:A:65:LYS:HD2	0.66	1.67	12	3
1:A:64:TRP:CE3	1:A:65:LYS:N	0.65	2.65	6	21
1:A:29:LEU:HG	1:A:33:VAL:HB	0.65	1.69	7	12
1:A:21:VAL:HG13	1:A:35:ARG:CD	0.65	2.21	2	1
1:A:36:MET:HE2	1:A:37:ARG:HG3	0.64	1.67	20	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:53:MET:CE	1:A:61:LEU:HD11	0.64	2.22	20	1
1:A:17:GLU:HA	1:A:20:VAL:CG2	0.64	2.22	9	20
1:A:29:LEU:CG	1:A:38:LEU:HD21	0.63	2.23	18	3
1:A:19:LEU:O	1:A:38:LEU:HD22	0.62	1.94	18	8
1:A:29:LEU:N	1:A:29:LEU:CD2	0.62	2.62	21	5
1:A:16:TYR:CE1	1:A:50:VAL:HG11	0.62	2.29	5	10
1:A:53:MET:CE	1:A:69:LEU:HD13	0.62	2.23	15	2
1:A:14:TYR:HB3	1:A:15:PRO:HD2	0.62	1.72	15	11
1:A:21:VAL:HG23	1:A:75:LEU:HB3	0.62	1.71	6	14
1:A:65:LYS:HD2	1:A:69:LEU:HD11	0.62	1.72	9	4
1:A:36:MET:HE2	1:A:37:ARG:HG2	0.62	1.70	20	1
1:A:29:LEU:HD21	1:A:38:LEU:CD2	0.62	2.24	18	2
1:A:18:MET:HE3	1:A:30:PRO:HG3	0.62	1.69	9	1
1:A:19:LEU:HD21	1:A:33:VAL:HG11	0.62	1.69	2	1
1:A:19:LEU:HD23	1:A:29:LEU:HD12	0.62	1.70	7	1
1:A:29:LEU:HB3	1:A:33:VAL:HB	0.61	1.72	15	9
1:A:50:VAL:HG12	1:A:73:ALA:HB1	0.61	1.70	20	2
1:A:50:VAL:CG1	1:A:73:ALA:HB1	0.61	2.25	20	13
1:A:39:GLU:O	1:A:47:PHE:CE1	0.61	2.53	20	21
1:A:14:TYR:CD1	1:A:19:LEU:HD21	0.61	2.30	7	4
1:A:29:LEU:CG	1:A:33:VAL:HB	0.60	2.26	11	15
1:A:13:ILE:HG21	1:A:43:SER:HB3	0.60	1.73	19	6
1:A:11:LEU:HD13	1:A:41:HIS:ND1	0.60	2.11	20	4
1:A:16:TYR:HE1	1:A:75:LEU:HD21	0.60	1.55	10	18
1:A:51:PHE:O	1:A:52:ALA:HB3	0.60	1.97	5	21
1:A:19:LEU:HD22	1:A:33:VAL:CG1	0.59	2.18	12	2
1:A:29:LEU:CD1	1:A:33:VAL:HG11	0.59	2.28	21	1
1:A:29:LEU:HB2	1:A:33:VAL:CG2	0.59	2.28	19	5
1:A:21:VAL:HG23	1:A:75:LEU:CB	0.58	2.28	12	13
1:A:19:LEU:HD23	1:A:29:LEU:CD1	0.58	2.28	7	5
1:A:19:LEU:HD22	1:A:38:LEU:CD2	0.58	2.28	2	1
1:A:19:LEU:HA	1:A:29:LEU:HD12	0.58	1.73	9	1
1:A:57:GLU:O	1:A:61:LEU:HG	0.58	1.99	4	21
1:A:47:PHE:CE1	1:A:51:PHE:CD2	0.57	2.92	10	20
1:A:47:PHE:CE2	1:A:55:PRO:HA	0.57	2.33	4	20
1:A:29:LEU:HD23	1:A:34:ASP:N	0.57	2.14	15	3
1:A:16:TYR:O	1:A:20:VAL:HG23	0.57	1.99	7	3
1:A:30:PRO:CD	1:A:33:VAL:HG21	0.57	2.30	2	10
1:A:61:LEU:HD22	1:A:65:LYS:HE2	0.56	1.76	20	1
1:A:33:VAL:HG12	1:A:41:HIS:CD2	0.56	2.35	2	1
1:A:47:PHE:CE1	1:A:51:PHE:CE2	0.56	2.93	9	19

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:17:GLU:C	1:A:20:VAL:HG23	0.56	2.25	3	15
1:A:21:VAL:CG2	1:A:75:LEU:HA	0.56	2.30	3	18
1:A:16:TYR:OH	1:A:73:ALA:O	0.56	2.23	5	5
1:A:11:LEU:HD11	1:A:40:ARG:O	0.55	2.00	4	2
1:A:18:MET:C	1:A:19:LEU:HD23	0.55	2.26	9	2
1:A:29:LEU:HB2	1:A:33:VAL:CB	0.55	2.32	13	6
1:A:20:VAL:HG11	1:A:23:ASN:HB2	0.55	1.78	21	2
1:A:13:ILE:HG21	1:A:43:SER:HB2	0.55	1.79	16	2
1:A:11:LEU:HD13	1:A:41:HIS:CE1	0.55	2.37	20	1
1:A:61:LEU:HD13	1:A:65:LYS:HG2	0.55	1.78	14	1
1:A:21:VAL:HG22	1:A:38:LEU:CD1	0.55	2.32	18	12
1:A:21:VAL:HG23	1:A:75:LEU:HA	0.54	1.77	9	13
1:A:15:PRO:O	1:A:19:LEU:HD12	0.53	2.03	6	5
1:A:21:VAL:HG13	1:A:35:ARG:NE	0.53	2.19	2	2
1:A:70:LYS:C	1:A:76:PHE:HB2	0.53	2.29	6	15
1:A:11:LEU:H	1:A:11:LEU:HD22	0.53	1.64	3	3
1:A:30:PRO:HG2	1:A:33:VAL:HG21	0.53	1.80	6	3
1:A:20:VAL:HA	1:A:75:LEU:HD23	0.52	1.82	1	5
1:A:53:MET:HE1	1:A:58:PHE:CA	0.52	2.34	7	14
1:A:13:ILE:HG23	1:A:41:HIS:O	0.52	2.04	8	6
1:A:53:MET:CE	1:A:58:PHE:CD2	0.52	2.92	4	4
1:A:21:VAL:HG23	1:A:75:LEU:CA	0.52	2.35	9	11
1:A:29:LEU:HD23	1:A:34:ASP:CA	0.52	2.34	18	3
1:A:51:PHE:CE1	1:A:73:ALA:HB2	0.52	2.39	17	9
1:A:65:LYS:O	1:A:69:LEU:HG	0.52	2.04	5	19
1:A:16:TYR:CE1	1:A:50:VAL:HG21	0.52	2.39	7	1
1:A:53:MET:HE3	1:A:58:PHE:CD2	0.52	2.39	10	2
1:A:50:VAL:HG12	1:A:73:ALA:CB	0.51	2.34	20	10
1:A:21:VAL:HG13	1:A:35:ARG:HB3	0.51	1.81	13	1
1:A:16:TYR:CE2	1:A:75:LEU:CD2	0.51	2.93	20	1
1:A:17:GLU:HA	1:A:20:VAL:HG23	0.51	1.82	20	13
1:A:61:LEU:HD13	1:A:65:LYS:HE3	0.51	1.82	6	1
1:A:17:GLU:CA	1:A:20:VAL:HG23	0.51	2.36	20	13
1:A:32:GLY:O	1:A:41:HIS:CE1	0.51	2.64	16	14
1:A:70:LYS:HD2	1:A:75:LEU:HD12	0.51	1.81	10	1
1:A:65:LYS:HG3	1:A:69:LEU:HD11	0.51	1.81	17	1
1:A:13:ILE:HG21	1:A:43:SER:CB	0.51	2.35	19	4
1:A:19:LEU:HB3	1:A:38:LEU:HD22	0.50	1.82	12	4
1:A:14:TYR:N	1:A:14:TYR:CD1	0.50	2.79	6	1
1:A:69:LEU:O	1:A:73:ALA:N	0.50	2.42	10	17
1:A:19:LEU:HA	1:A:29:LEU:CD1	0.50	2.36	1	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:29:LEU:HD21	1:A:38:LEU:CD1	0.50	2.31	2	1
1:A:51:PHE:O	1:A:52:ALA:CB	0.50	2.59	5	4
1:A:12:GLN:HG3	1:A:14:TYR:OH	0.50	2.07	18	2
1:A:62:ALA:O	1:A:66:ARG:HG3	0.50	2.07	14	20
1:A:14:TYR:O	1:A:19:LEU:HD12	0.50	2.06	13	6
1:A:29:LEU:O	1:A:30:PRO:O	0.50	2.30	19	4
1:A:53:MET:SD	1:A:69:LEU:HD22	0.50	2.46	12	1
1:A:29:LEU:CG	1:A:33:VAL:HG11	0.50	2.37	21	1
1:A:23:ASN:ND2	1:A:74:GLU:HB3	0.50	2.21	2	12
1:A:29:LEU:HD22	1:A:29:LEU:N	0.49	2.20	9	3
1:A:11:LEU:HD21	1:A:40:ARG:HD3	0.49	1.82	12	1
1:A:30:PRO:HD2	1:A:33:VAL:CG2	0.49	2.37	18	3
1:A:54:SER:O	1:A:57:GLU:N	0.49	2.43	10	17
1:A:51:PHE:CD1	1:A:69:LEU:HB3	0.49	2.42	5	15
1:A:47:PHE:CD1	1:A:51:PHE:CD2	0.49	3.01	20	3
1:A:33:VAL:CG1	1:A:41:HIS:CG	0.49	2.95	2	1
1:A:71:LYS:CG	1:A:76:PHE:CG	0.49	2.96	6	11
1:A:53:MET:HE1	1:A:58:PHE:CB	0.49	2.37	1	1
1:A:68:GLU:O	1:A:72:LYS:N	0.48	2.46	6	20
1:A:53:MET:HG2	1:A:57:GLU:HB2	0.48	1.84	5	3
1:A:14:TYR:CE1	1:A:33:VAL:CG2	0.48	2.94	3	3
1:A:71:LYS:CE	1:A:76:PHE:CD1	0.48	2.96	12	1
1:A:51:PHE:CE1	1:A:73:ALA:CB	0.48	2.96	2	15
1:A:12:GLN:HG2	1:A:14:TYR:OH	0.48	2.09	3	2
1:A:29:LEU:HD21	1:A:38:LEU:CG	0.48	2.39	15	3
1:A:62:ALA:HB3	1:A:64:TRP:CE2	0.48	2.44	21	18
1:A:29:LEU:HD11	1:A:38:LEU:CD2	0.47	2.39	6	6
1:A:53:MET:HE3	1:A:58:PHE:HD2	0.47	1.68	10	1
1:A:64:TRP:CZ3	1:A:65:LYS:CG	0.47	2.97	4	2
1:A:14:TYR:CD2	1:A:30:PRO:HB2	0.47	2.43	21	1
1:A:70:LYS:HB3	1:A:76:PHE:HA	0.47	1.86	1	4
1:A:61:LEU:HD13	1:A:65:LYS:CG	0.47	2.40	14	2
1:A:51:PHE:HB3	1:A:53:MET:HE2	0.47	1.86	15	1
1:A:64:TRP:CD2	1:A:65:LYS:N	0.47	2.83	3	20
1:A:15:PRO:O	1:A:19:LEU:HG	0.47	2.10	18	5
1:A:67:ASN:O	1:A:71:LYS:N	0.47	2.48	17	15
1:A:16:TYR:CD2	1:A:17:GLU:N	0.47	2.83	16	2
1:A:19:LEU:HD13	1:A:41:HIS:CB	0.46	2.40	16	2
1:A:70:LYS:NZ	1:A:75:LEU:O	0.46	2.47	10	1
1:A:70:LYS:HD2	1:A:75:LEU:O	0.46	2.11	10	1
1:A:42:LEU:HD13	1:A:51:PHE:HE1	0.46	1.70	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:29:LEU:HG	1:A:33:VAL:CG1	0.46	2.40	21	1
1:A:30:PRO:CB	1:A:31:PRO:CD	0.46	2.94	16	12
1:A:29:LEU:CB	1:A:30:PRO:CD	0.46	2.93	9	3
1:A:65:LYS:HE3	1:A:69:LEU:HD21	0.46	1.87	10	1
1:A:14:TYR:HB2	1:A:19:LEU:HD21	0.46	1.88	1	2
1:A:64:TRP:CZ3	1:A:65:LYS:HG2	0.46	2.46	9	5
1:A:47:PHE:CE2	1:A:53:MET:SD	0.46	3.08	4	3
1:A:53:MET:HG3	1:A:57:GLU:HB2	0.46	1.87	10	1
1:A:16:TYR:CE1	1:A:75:LEU:CD2	0.45	2.98	21	1
1:A:47:PHE:CD1	1:A:51:PHE:CD1	0.45	3.04	1	1
1:A:30:PRO:CB	1:A:31:PRO:HD2	0.45	2.41	2	2
1:A:70:LYS:O	1:A:75:LEU:C	0.45	2.59	4	16
1:A:14:TYR:CB	1:A:15:PRO:CD	0.45	2.95	3	4
1:A:63:LEU:O	1:A:67:ASN:ND2	0.45	2.50	15	4
1:A:22:THR:O	1:A:23:ASN:C	0.45	2.59	17	4
1:A:61:LEU:CD1	1:A:69:LEU:HD11	0.45	2.41	7	11
1:A:53:MET:HE1	1:A:58:PHE:CD2	0.45	2.46	4	3
1:A:65:LYS:HD3	1:A:69:LEU:HD11	0.45	1.89	4	1
1:A:29:LEU:CD1	1:A:38:LEU:CD2	0.45	2.90	11	4
1:A:29:LEU:HD12	1:A:33:VAL:HG12	0.45	1.86	19	3
1:A:14:TYR:HB3	1:A:15:PRO:CD	0.45	2.41	12	2
1:A:35:ARG:HG3	1:A:38:LEU:HD12	0.45	1.87	17	1
1:A:73:ALA:O	1:A:75:LEU:HG	0.45	2.12	11	2
1:A:16:TYR:O	1:A:19:LEU:N	0.44	2.51	4	10
1:A:66:ARG:CA	1:A:69:LEU:HD12	0.44	2.39	15	9
1:A:22:THR:CG2	1:A:23:ASN:N	0.44	2.80	9	2
1:A:29:LEU:O	1:A:30:PRO:C	0.44	2.60	21	1
1:A:13:ILE:HG21	1:A:43:SER:CA	0.44	2.43	8	3
1:A:64:TRP:CE3	1:A:65:LYS:CA	0.44	3.00	15	12
1:A:70:LYS:CD	1:A:75:LEU:O	0.44	2.65	10	1
1:A:12:GLN:CG	1:A:14:TYR:OH	0.44	2.66	17	3
1:A:53:MET:HG3	1:A:54:SER:N	0.44	2.27	10	1
1:A:71:LYS:HG2	1:A:76:PHE:CG	0.44	2.48	6	7
1:A:11:LEU:N	1:A:11:LEU:CD1	0.43	2.77	21	2
1:A:39:GLU:HA	1:A:42:LEU:CD1	0.43	2.43	7	4
1:A:13:ILE:HG22	1:A:43:SER:OG	0.43	2.13	15	1
1:A:48:SER:O	1:A:52:ALA:HA	0.43	2.13	20	3
1:A:14:TYR:CG	1:A:30:PRO:CG	0.43	3.02	19	1
1:A:33:VAL:HG12	1:A:41:HIS:CG	0.43	2.49	2	1
1:A:20:VAL:HG11	1:A:23:ASN:OD1	0.43	2.14	20	1
1:A:62:ALA:CB	1:A:64:TRP:NE1	0.43	2.82	6	10

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:29:LEU:HD23	1:A:34:ASP:HA	0.43	1.89	18	1
1:A:71:LYS:HG3	1:A:76:PHE:CG	0.43	2.49	8	9
1:A:29:LEU:HD11	1:A:38:LEU:HD11	0.43	1.90	15	1
1:A:38:LEU:C	1:A:40:ARG:H	0.43	2.22	21	1
1:A:67:ASN:OD1	1:A:76:PHE:CE2	0.43	2.72	6	1
1:A:64:TRP:CZ3	1:A:65:LYS:CD	0.43	3.02	14	1
1:A:53:MET:HE1	1:A:61:LEU:HD11	0.43	1.91	20	1
1:A:12:GLN:O	1:A:14:TYR:CE1	0.43	2.72	3	4
1:A:20:VAL:HG12	1:A:21:VAL:N	0.42	2.30	12	6
1:A:19:LEU:CD2	1:A:29:LEU:CD1	0.42	2.97	6	1
1:A:61:LEU:HD12	1:A:69:LEU:HD11	0.42	1.91	21	3
1:A:51:PHE:HB3	1:A:53:MET:CE	0.42	2.44	15	1
1:A:30:PRO:CG	1:A:33:VAL:HG21	0.42	2.44	18	1
1:A:13:ILE:HG21	1:A:43:SER:HA	0.42	1.91	12	1
1:A:14:TYR:CD2	1:A:30:PRO:CB	0.42	3.03	21	1
1:A:71:LYS:CD	1:A:76:PHE:CG	0.42	3.02	10	1
1:A:16:TYR:HD1	1:A:42:LEU:HD21	0.42	1.75	3	1
1:A:19:LEU:O	1:A:38:LEU:CD2	0.42	2.68	13	1
1:A:45:GLU:O	1:A:49:ARG:CG	0.42	2.67	16	2
1:A:51:PHE:CZ	1:A:75:LEU:CD1	0.41	3.03	1	1
1:A:71:LYS:HG2	1:A:76:PHE:CD2	0.41	2.49	18	2
1:A:73:ALA:O	1:A:74:GLU:HB2	0.41	2.15	7	1
1:A:65:LYS:NZ	1:A:69:LEU:HD21	0.41	2.31	8	1
1:A:61:LEU:HD13	1:A:69:LEU:HD11	0.41	1.92	20	1
1:A:71:LYS:HG2	1:A:76:PHE:CD1	0.41	2.50	6	1
1:A:23:ASN:ND2	1:A:74:GLU:CB	0.41	2.83	10	2
1:A:12:GLN:O	1:A:14:TYR:CZ	0.41	2.72	9	1
1:A:30:PRO:HB2	1:A:31:PRO:CD	0.41	2.46	7	3
1:A:19:LEU:O	1:A:38:LEU:CD1	0.41	2.63	12	2
1:A:42:LEU:HD13	1:A:51:PHE:CE1	0.41	2.50	1	1
1:A:12:GLN:HB2	1:A:14:TYR:CZ	0.41	2.50	12	1
1:A:19:LEU:CB	1:A:38:LEU:HD22	0.41	2.46	16	1
1:A:29:LEU:HB3	1:A:33:VAL:CG2	0.41	2.46	7	1
1:A:64:TRP:CH2	1:A:65:LYS:HG3	0.41	2.51	12	1
1:A:29:LEU:CG	1:A:33:VAL:CG1	0.41	2.99	21	1
1:A:70:LYS:HB3	1:A:76:PHE:HB2	0.41	1.91	6	1
1:A:22:THR:HG23	1:A:23:ASN:N	0.41	2.30	12	1
1:A:19:LEU:CD2	1:A:29:LEU:HD12	0.40	2.46	6	1
1:A:54:SER:O	1:A:57:GLU:HB2	0.40	2.16	9	1
1:A:12:GLN:OE1	1:A:14:TYR:CE2	0.40	2.74	13	1
1:A:14:TYR:O	1:A:19:LEU:CD1	0.40	2.69	15	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:53:MET:HE3	1:A:53:MET:HB2	0.40	1.77	1	1
1:A:76:PHE:OXT	1:A:76:PHE:CD1	0.40	2.75	10	1
1:A:50:VAL:HG11	1:A:73:ALA:HB1	0.40	1.93	6	1
1:A:14:TYR:CG	1:A:30:PRO:HG3	0.40	2.52	7	1
1:A:14:TYR:CB	1:A:15:PRO:HD2	0.40	2.46	15	1
1:A:76:PHE:CD2	1:A:76:PHE:OXT	0.40	2.74	15	1
1:A:64:TRP:CZ3	1:A:65:LYS:HD3	0.40	2.52	17	1
1:A:23:ASN:ND2	1:A:74:GLU:HA	0.40	2.32	3	1
1:A:33:VAL:HG13	1:A:41:HIS:CB	0.40	2.46	18	1
1:A:76:PHE:CD2	1:A:76:PHE:O	0.40	2.74	20	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	60/68 (88%)	46±2 (76±3%)	12±2 (20±3%)	2±1 (4±1%)	4	29
All	All	1260/1428 (88%)	958 (76%)	250 (20%)	52 (4%)	4	29

All 6 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	30	PRO	21
1	A	61	LEU	14
1	A	15	PRO	11
1	A	23	ASN	4
1	A	20	VAL	1
1	A	31	PRO	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	55/60 (92%)	38±2 (70±4%)	17±2 (30±4%)	1 17
All	All	1155/1260 (92%)	808 (70%)	347 (30%)	1 17

All 34 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	11	LEU	21
1	A	29	LEU	21
1	A	45	GLU	21
1	A	75	LEU	21
1	A	12	GLN	20
1	A	33	VAL	20
1	A	53	MET	20
1	A	72	LYS	20
1	A	37	ARG	19
1	A	63	LEU	18
1	A	40	ARG	17
1	A	43	SER	12
1	A	35	ARG	10
1	A	18	MET	10
1	A	39	GLU	9
1	A	60	LYS	9
1	A	17	GLU	8
1	A	36	MET	8
1	A	14	TYR	7
1	A	57	GLU	7
1	A	65	LYS	7
1	A	49	ARG	7
1	A	71	LYS	6
1	A	76	PHE	6
1	A	70	LYS	5
1	A	56	GLU	4
1	A	48	SER	3
1	A	68	GLU	3
1	A	74	GLU	2
1	A	19	LEU	2
1	A	34	ASP	1
1	A	67	ASN	1
1	A	69	LEU	1
1	A	23	ASN	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

There are no oligosaccharides in this entry.

6.6 Ligand geometry

There are no ligands in this entry.

6.7 Other polymers

There are no such molecules in this entry.

6.8 Polymer linkage issues

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided