



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 06:48 AM UTC

PDB ID : 1JGN / pdb_00001jgn
Title : Solution structure of the C-terminal PABC domain of human poly(A)-binding protein in complex with the peptide from Paip2
Authors : Kozlov, G.; Siddiqui, N.; Coillet-Matillon, S.; Ekiel, I.; Gehring, K.
Deposited on : 2001-06-26

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

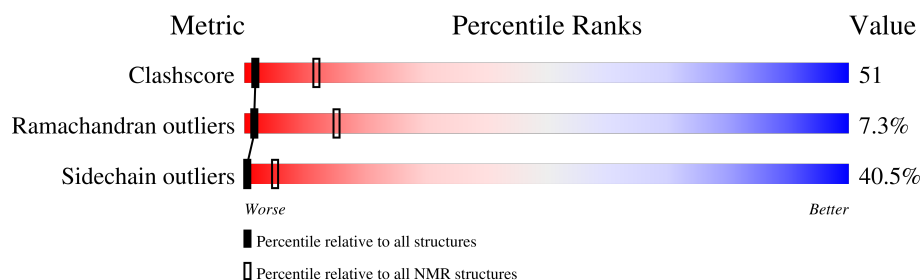
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and a grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	98	
2	B	22	

2 Ensemble composition and analysis

This entry contains 30 models. Model 16 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:8-A:15, A:19-A:79, B:5-B:13, B:18-B:19 (80)	0.36	16

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 4 single-model clusters were found.

Cluster number	Models
1	2, 3, 5, 7, 9, 11, 12, 14, 15, 16, 17, 19, 20, 24, 25, 27, 30
2	1, 10, 18
3	6, 8
4	28, 29
5	13, 21
Single-model clusters	4; 22; 23; 26

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1815 atoms, of which 924 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called polyadenylate-binding protein 1.

Mol	Chain	Residues	Atoms						Trace
1	A	98	Total	C	H	N	O	S	0
			1469	452	747	124	141	5	

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	cloning artifact	UNP P11940
A	2	PRO	-	cloning artifact	UNP P11940
A	3	LEU	-	cloning artifact	UNP P11940
A	4	GLY	-	cloning artifact	UNP P11940
A	5	SER	-	cloning artifact	UNP P11940

- Molecule 2 is a protein called polyadenylate-binding protein-interacting protein 2.

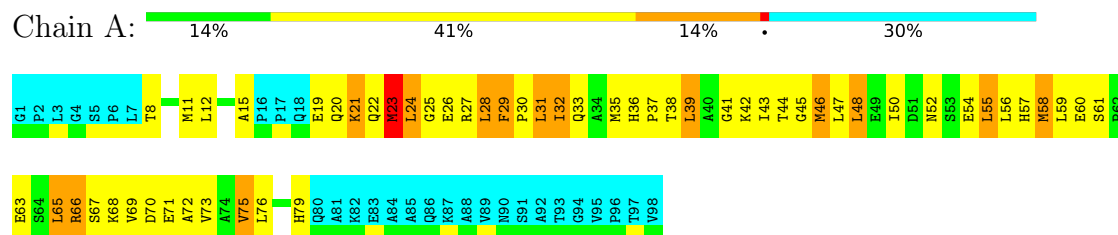
Mol	Chain	Residues	Atoms					Trace
2	B	22	Total	C	H	N	O	0
			346	109	177	29	31	

4 Residue-property plots

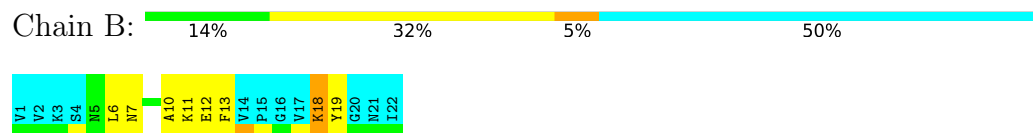
4.1 Average score per residue in the NMR ensemble

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: polyadenylate-binding protein 1



- Molecule 2: polyadenylate-binding protein-interacting protein 2

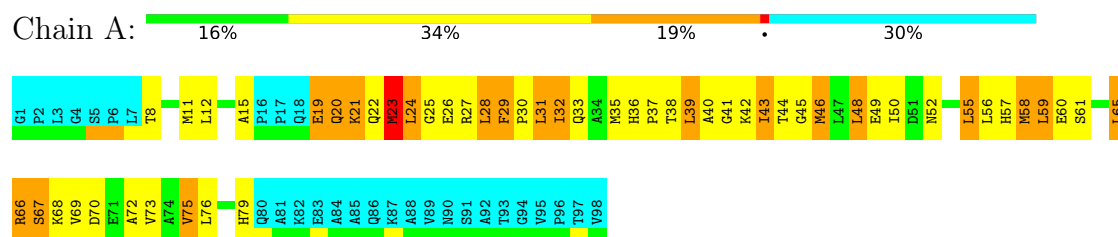


4.2 Scores per residue for each member of the ensemble

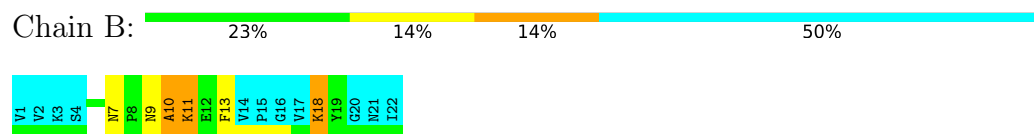
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: polyadenylate-binding protein 1

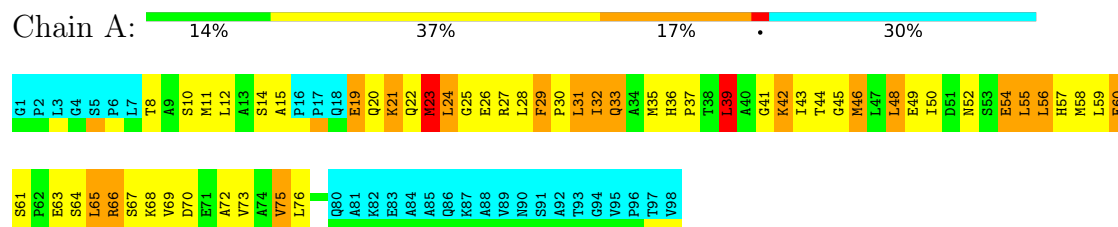


- Molecule 2: polyadenylate-binding protein-interacting protein 2

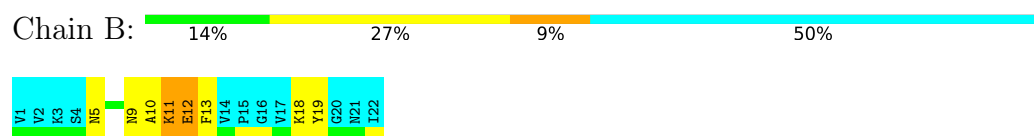


4.2.2 Score per residue for model 2

- Molecule 1: polyadenylate-binding protein 1

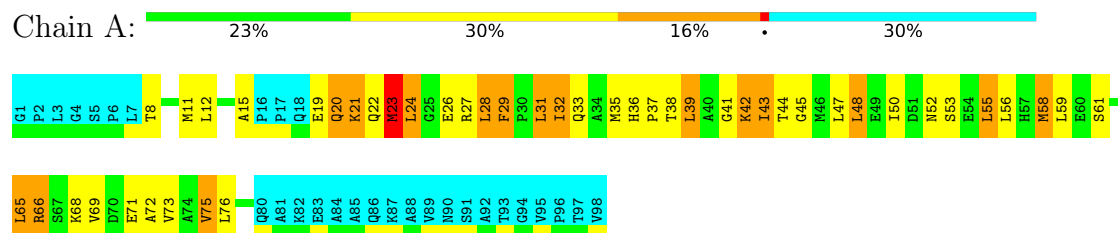


- Molecule 2: polyadenylate-binding protein-interacting protein 2

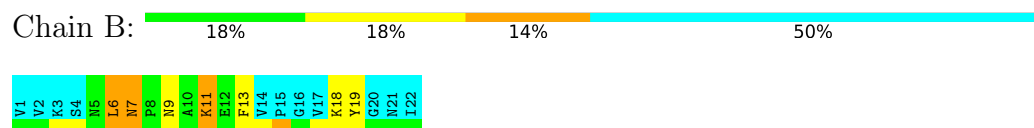


4.2.3 Score per residue for model 3

- Molecule 1: polyadenylate-binding protein 1

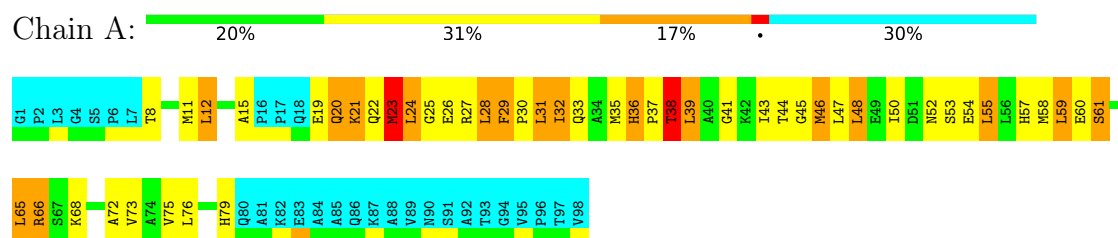


- Molecule 2: polyadenylate-binding protein-interacting protein 2

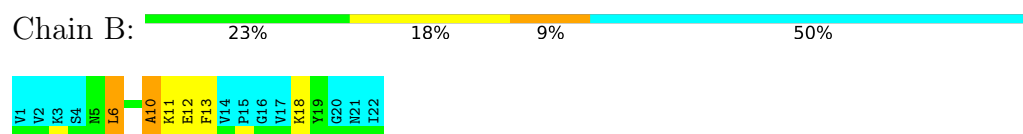


4.2.4 Score per residue for model 4

- Molecule 1: polyadenylate-binding protein 1

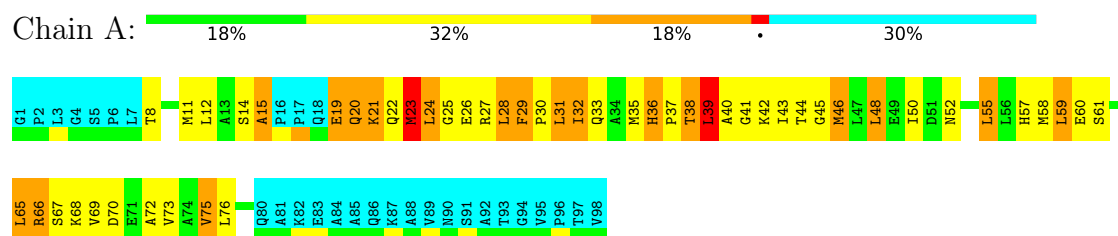


- Molecule 2: polyadenylate-binding protein-interacting protein 2

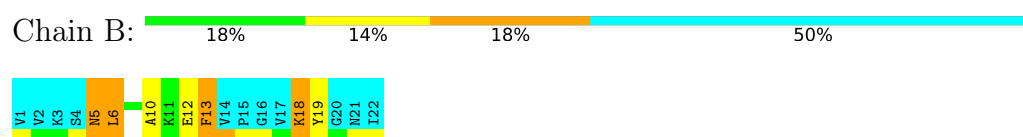


4.2.5 Score per residue for model 5

- Molecule 1: polyadenylate-binding protein 1

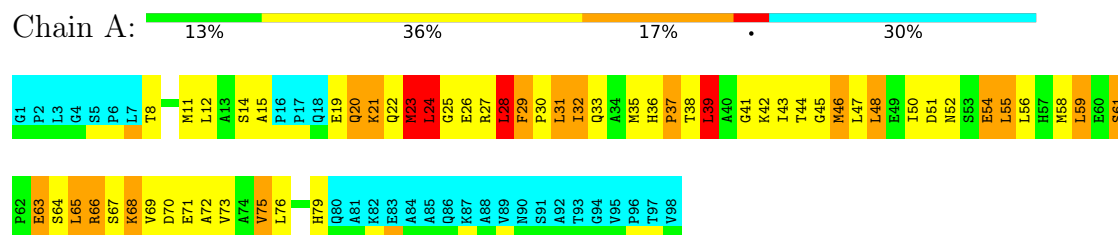


- Molecule 2: polyadenylate-binding protein-interacting protein 2

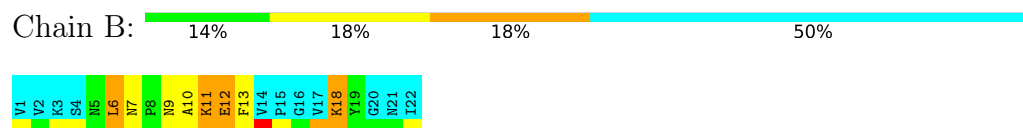


4.2.6 Score per residue for model 6

- Molecule 1: polyadenylate-binding protein 1

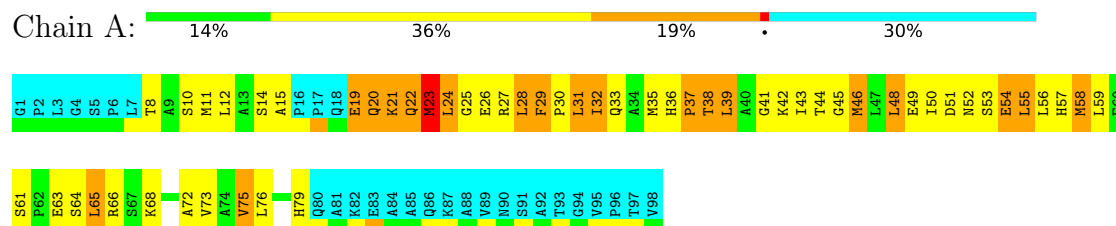


- Molecule 2: polyadenylate-binding protein-interacting protein 2

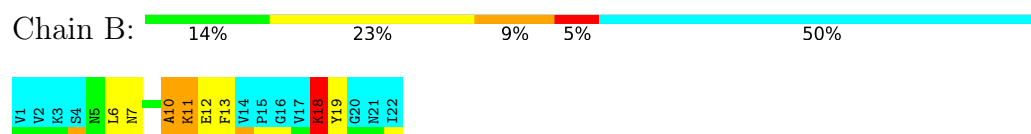


4.2.7 Score per residue for model 7

- Molecule 1: polyadenylate-binding protein 1

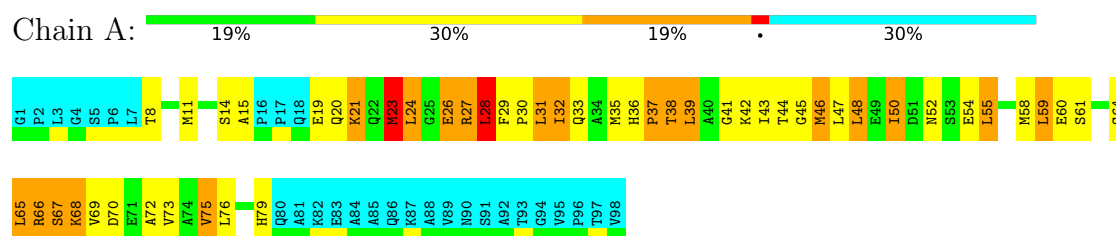


- Molecule 2: polyadenylate-binding protein-interacting protein 2

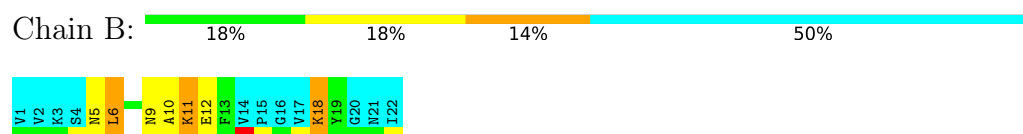


4.2.8 Score per residue for model 8

- Molecule 1: polyadenylate-binding protein 1

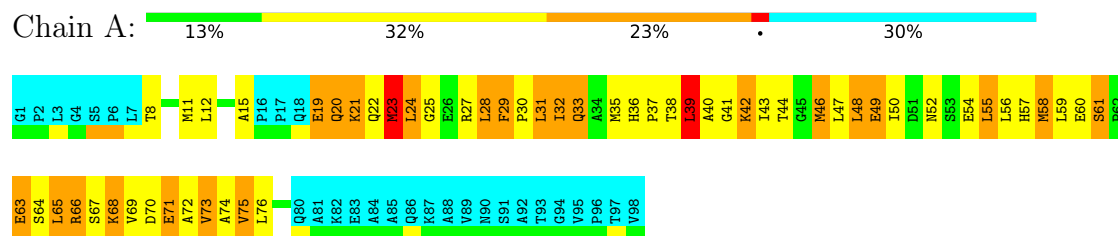


- Molecule 2: polyadenylate-binding protein-interacting protein 2

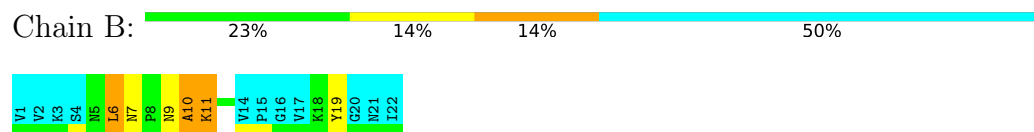


4.2.9 Score per residue for model 9

- Molecule 1: polyadenylate-binding protein 1

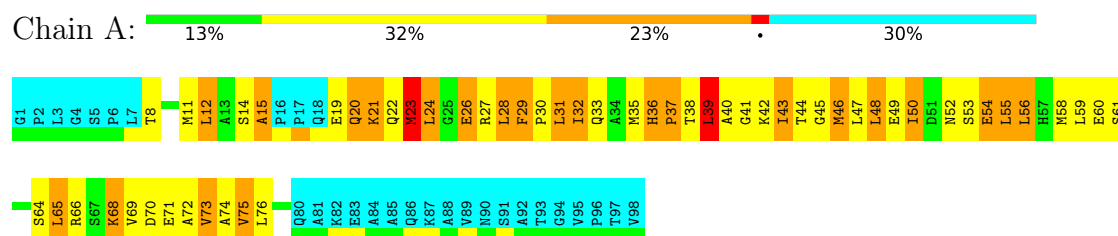


- Molecule 2: polyadenylate-binding protein-interacting protein 2

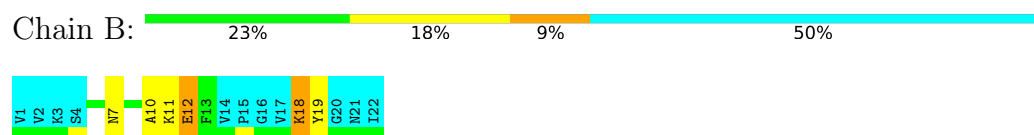


4.2.10 Score per residue for model 10

- Molecule 1: polyadenylate-binding protein 1

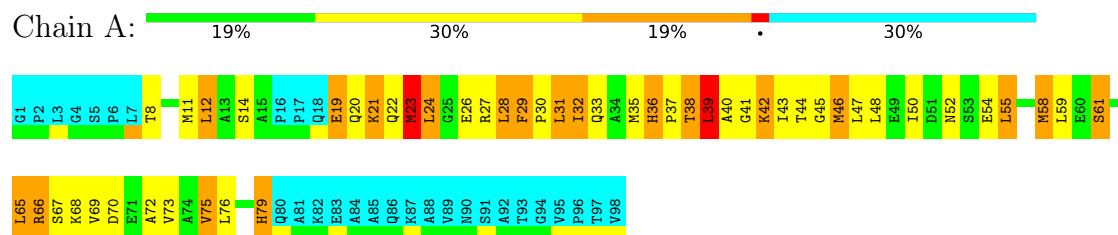


- Molecule 2: polyadenylate-binding protein-interacting protein 2

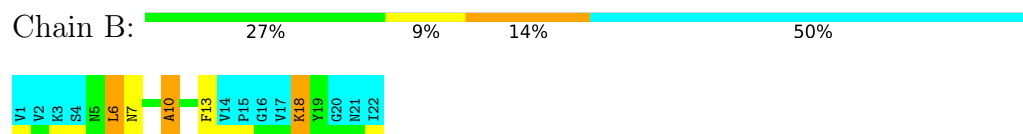


4.2.11 Score per residue for model 11

- Molecule 1: polyadenylate-binding protein 1

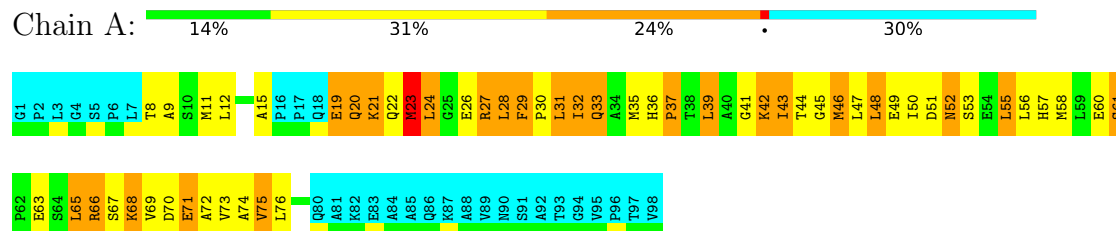


- Molecule 2: polyadenylate-binding protein-interacting protein 2

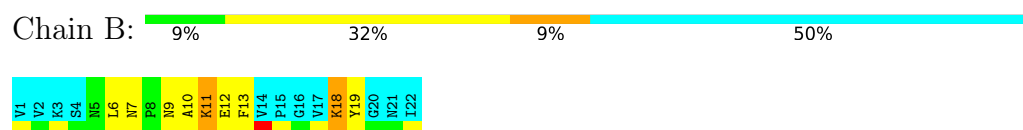


4.2.12 Score per residue for model 12

- Molecule 1: polyadenylate-binding protein 1

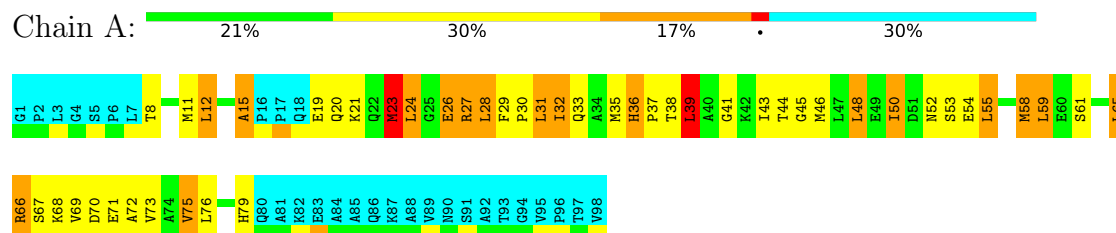


- Molecule 2: polyadenylate-binding protein-interacting protein 2

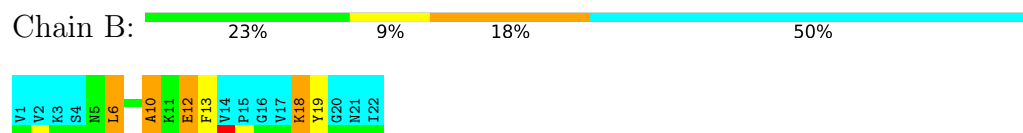


4.2.13 Score per residue for model 13

- Molecule 1: polyadenylate-binding protein 1

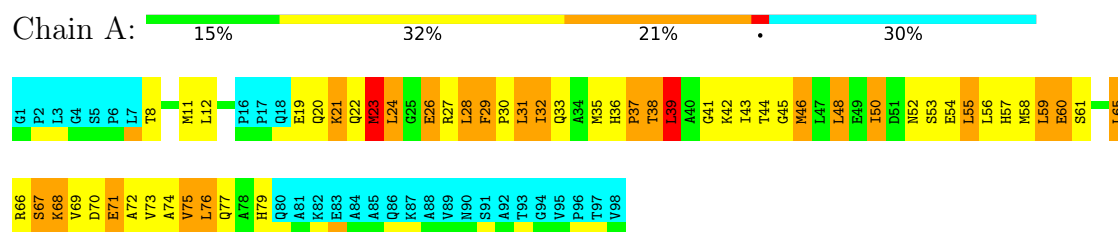


- Molecule 2: polyadenylate-binding protein-interacting protein 2

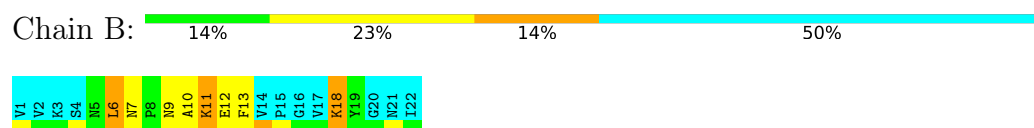


4.2.14 Score per residue for model 14

- Molecule 1: polyadenylate-binding protein 1

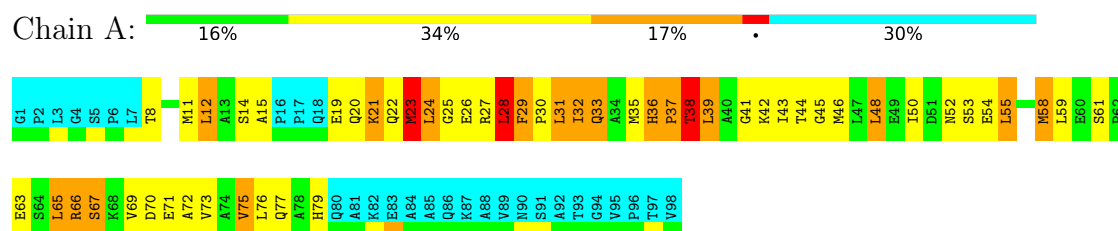


- Molecule 2: polyadenylate-binding protein-interacting protein 2

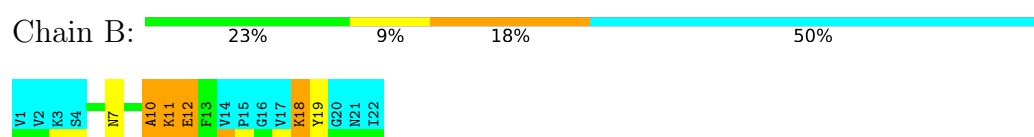


4.2.15 Score per residue for model 15

- Molecule 1: polyadenylate-binding protein 1

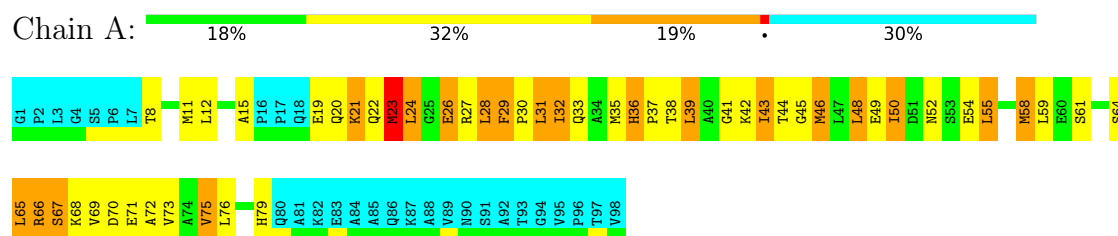


- Molecule 2: polyadenylate-binding protein-interacting protein 2



4.2.16 Score per residue for model 16 (medoid)

- Molecule 1: polyadenylate-binding protein 1



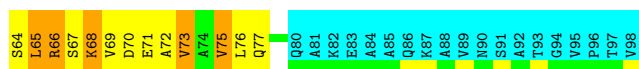
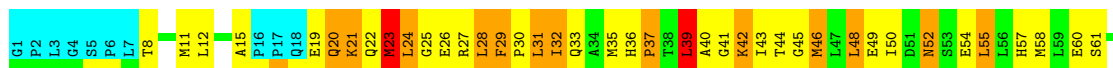
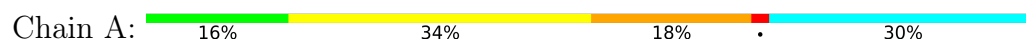
- Molecule 2: polyadenylate-binding protein-interacting protein 2





4.2.17 Score per residue for model 17

- Molecule 1: polyadenylate-binding protein 1

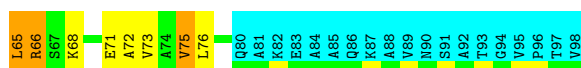
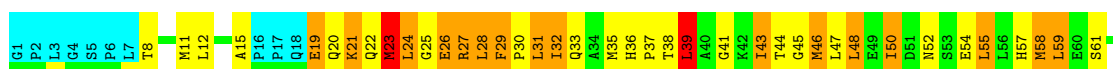


- Molecule 2: polyadenylate-binding protein-interacting protein 2

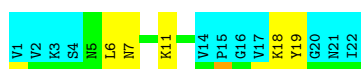
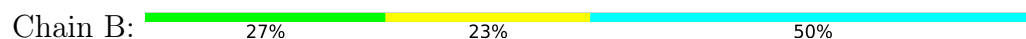


4.2.18 Score per residue for model 18

- Molecule 1: polyadenylate-binding protein 1



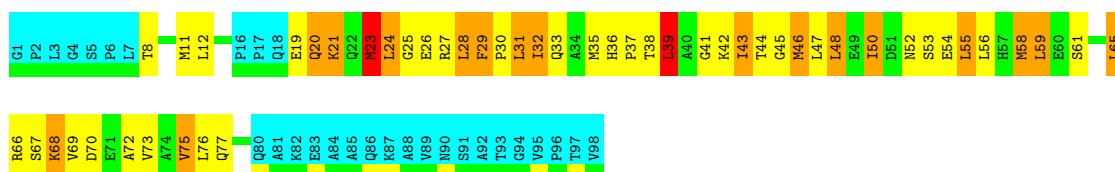
- Molecule 2: polyadenylate-binding protein-interacting protein 2



4.2.19 Score per residue for model 19

- Molecule 1: polyadenylate-binding protein 1





- Molecule 2: polyadenylate-binding protein-interacting protein 2

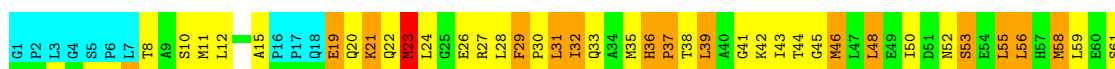
Chain B: 32% 14% 5% 50%



4.2.20 Score per residue for model 20

- Molecule 1: polyadenylate-binding protein 1

Chain A: 16% 34% 19% . 30%



- Molecule 2: polyadenylate-binding protein-interacting protein 2

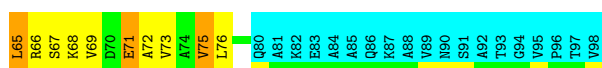
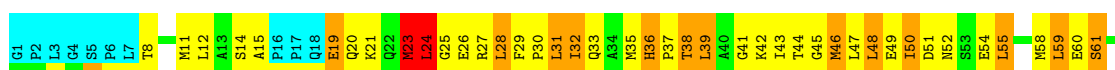
Chain B: 9% 41% 50%



4.2.21 Score per residue for model 21

- Molecule 1: polyadenylate-binding protein 1

Chain A: 17% 35% 16% . 30%



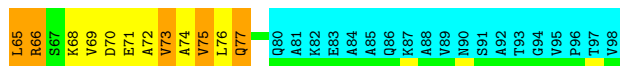
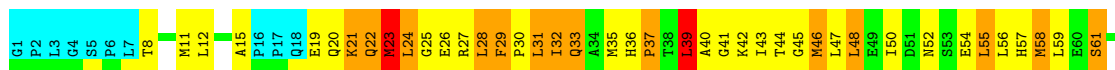
- Molecule 2: polyadenylate-binding protein-interacting protein 2

Chain B: 5% 23% 18% 5% 50%

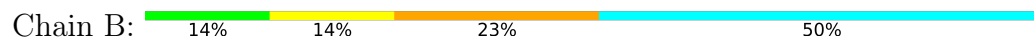


4.2.22 Score per residue for model 22

- Molecule 1: polyadenylate-binding protein 1

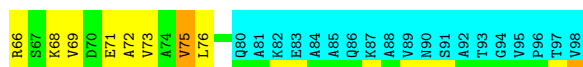


- Molecule 2: polyadenylate-binding protein-interacting protein 2

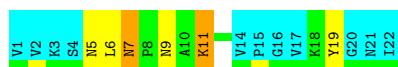


4.2.23 Score per residue for model 23

- Molecule 1: polyadenylate-binding protein 1



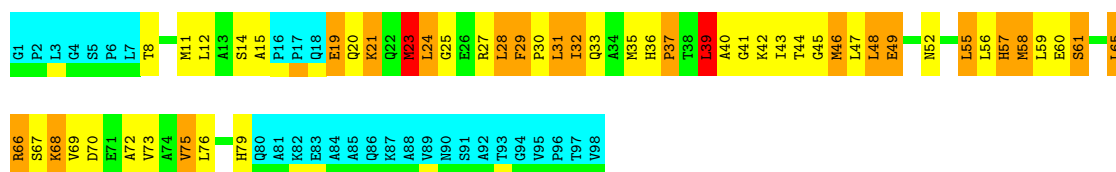
- Molecule 2: polyadenylate-binding protein-interacting protein 2



4.2.24 Score per residue for model 24

- Molecule 1: polyadenylate-binding protein 1





- Molecule 2: polyadenylate-binding protein-interacting protein 2

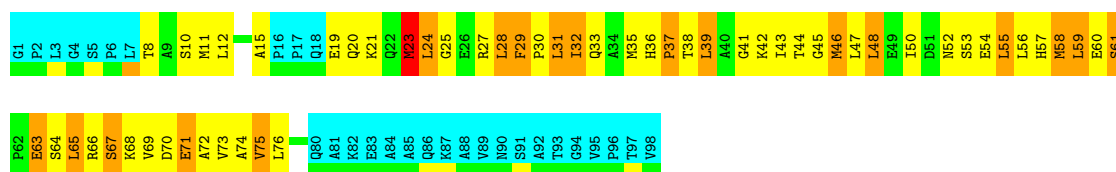
Chain B: 14% 23% 14% 50%



4.2.25 Score per residue for model 25

- Molecule 1: polyadenylate-binding protein 1

Chain A: 13% 38% 18% 30%



- Molecule 2: polyadenylate-binding protein-interacting protein 2

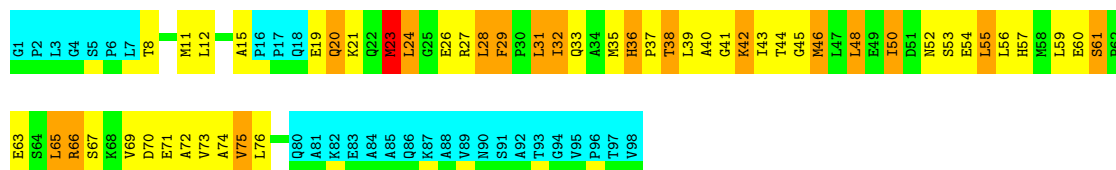
Chain B: 9% 23% 14% 5% 50%



4.2.26 Score per residue for model 26

- Molecule 1: polyadenylate-binding protein 1

Chain A: 18% 34% 17% 30%



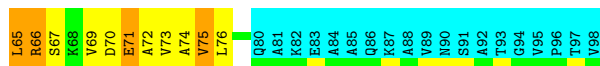
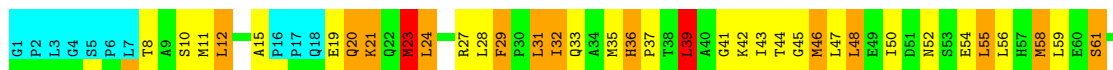
- Molecule 2: polyadenylate-binding protein-interacting protein 2

Chain B: 5% 27% 18% 50%

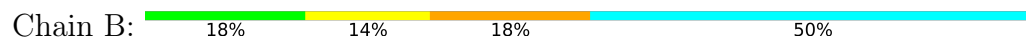


4.2.27 Score per residue for model 27

- Molecule 1: polyadenylate-binding protein 1

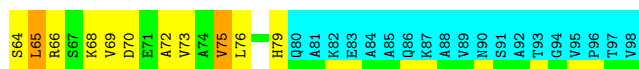


- Molecule 2: polyadenylate-binding protein-interacting protein 2

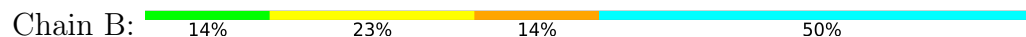


4.2.28 Score per residue for model 28

- Molecule 1: polyadenylate-binding protein 1



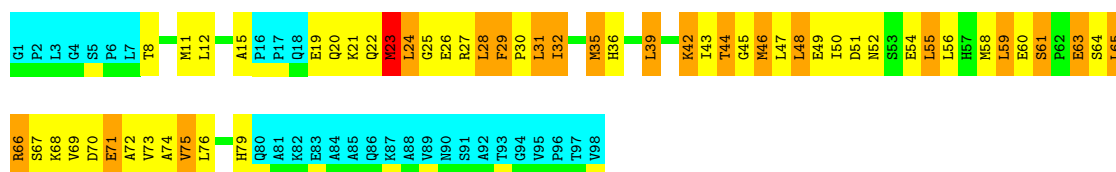
- Molecule 2: polyadenylate-binding protein-interacting protein 2



4.2.29 Score per residue for model 29

- Molecule 1: polyadenylate-binding protein 1





- Molecule 2: polyadenylate-binding protein-interacting protein 2

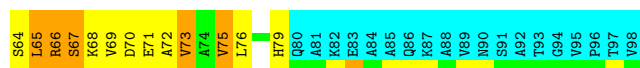
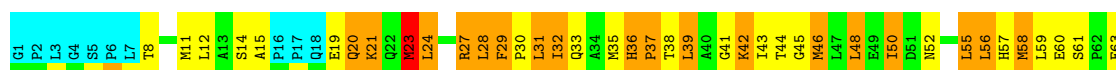
Chain B: 9% 23% 18% 50%



4.2.30 Score per residue for model 30

- Molecule 1: polyadenylate-binding protein 1

Chain A: 17% 29% 23% 30%



- Molecule 2: polyadenylate-binding protein-interacting protein 2

Chain B: 23% 23% 5% 50%



5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing*.

Of the 200 calculated structures, 30 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
ARIA	structure solution	0.9
CNS	refinement	0.9

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

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 [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	524	541	541	64±7
2	B	94	91	91	5±3
All	All	18540	18960	18960	1931

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 51.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:39:LEU:HD23	1:A:76:LEU:HD12	0.89	1.43	28	1
1:A:46:MET:HE3	1:A:47:LEU:HD23	0.86	1.47	21	1
1:A:38:THR:HG21	1:A:73:VAL:HG13	0.85	1.48	13	11
1:A:31:LEU:HD13	1:A:65:LEU:HD21	0.84	1.49	15	19
1:A:28:LEU:HD22	1:A:44:THR:HA	0.82	1.52	1	16
1:A:46:MET:HE1	1:A:75:VAL:HG21	0.80	1.52	21	1
1:A:58:MET:HE2	1:A:68:LYS:HD3	0.79	1.54	16	4
1:A:15:ALA:HB1	1:A:19:GLU:HB3	0.78	1.55	7	19
1:A:43:ILE:HD11	1:A:76:LEU:HD11	0.78	1.55	11	3
1:A:43:ILE:HG21	1:A:72:ALA:HB1	0.78	1.55	22	4
1:A:27:ARG:HG3	1:A:31:LEU:HD21	0.75	1.58	26	22
1:A:52:ASN:HA	1:A:55:LEU:HD23	0.75	1.56	18	30
1:A:50:ILE:HD11	1:A:68:LYS:NZ	0.75	1.97	7	5

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:36:HIS:N	1:A:37:PRO:HD2	0.74	1.98	25	28
1:A:42:LYS:HB3	1:A:76:LEU:HD11	0.74	1.57	12	8
1:A:27:ARG:HG2	1:A:59:LEU:HD22	0.74	1.58	27	3
1:A:12:LEU:HD21	1:A:24:LEU:HB3	0.73	1.59	18	8
1:A:27:ARG:HG2	1:A:59:LEU:HD12	0.73	1.61	14	4
1:A:45:GLY:HA2	2:B:13:PHE:HB2	0.72	1.60	22	19
1:A:46:MET:HB3	2:B:10:ALA:HB1	0.72	1.61	4	10
1:A:24:LEU:HD12	1:A:27:ARG:HD3	0.72	1.61	1	12
1:A:27:ARG:HB2	1:A:55:LEU:HD13	0.72	1.62	23	29
1:A:28:LEU:HD12	1:A:44:THR:HA	0.71	1.58	29	2
1:A:46:MET:HE3	1:A:75:VAL:HG23	0.71	1.61	14	3
1:A:11:MET:HE3	1:A:23:MET:HB3	0.71	1.63	14	27
1:A:28:LEU:HB3	1:A:44:THR:OG1	0.70	1.86	29	1
1:A:58:MET:HE3	1:A:65:LEU:HA	0.70	1.63	20	6
1:A:43:ILE:HA	1:A:46:MET:HE2	0.70	1.62	21	1
1:A:32:ILE:HG21	1:A:43:ILE:HD12	0.70	1.64	11	14
1:A:41:GLY:O	1:A:44:THR:HG22	0.69	1.87	23	7
1:A:46:MET:CE	1:A:75:VAL:HG21	0.69	2.18	21	1
1:A:43:ILE:CG2	1:A:72:ALA:HB1	0.69	2.17	11	30
1:A:47:LEU:HD23	2:B:7:ASN:HB2	0.68	1.63	24	5
1:A:59:LEU:H	1:A:59:LEU:HD13	0.68	1.48	23	4
1:A:27:ARG:HE	1:A:59:LEU:HD13	0.68	1.48	18	2
1:A:27:ARG:NE	1:A:59:LEU:HD12	0.67	2.04	16	6
1:A:42:LYS:HE3	2:B:12:GLU:HG3	0.67	1.66	14	4
1:A:31:LEU:HD13	1:A:69:VAL:HG21	0.66	1.68	8	1
1:A:31:LEU:HD22	1:A:65:LEU:HG	0.66	1.66	2	19
1:A:31:LEU:HB3	1:A:65:LEU:HD11	0.66	1.66	10	17
1:A:50:ILE:HD11	1:A:68:LYS:HZ1	0.65	1.50	16	3
1:A:72:ALA:O	1:A:75:VAL:HG22	0.65	1.90	22	29
1:A:27:ARG:NE	1:A:59:LEU:HD13	0.65	2.06	5	4
1:A:65:LEU:C	1:A:65:LEU:HD22	0.65	2.17	11	19
1:A:38:THR:HG22	1:A:39:LEU:HD12	0.65	1.67	26	1
1:A:8:THR:HG21	1:A:29:PHE:HB3	0.65	1.68	22	14
1:A:39:LEU:HB3	1:A:76:LEU:HD22	0.65	1.69	22	1
1:A:8:THR:HA	1:A:11:MET:HG3	0.64	1.69	5	30
1:A:12:LEU:HD12	1:A:23:MET:HE2	0.64	1.69	13	7
1:A:12:LEU:HD12	1:A:23:MET:SD	0.64	2.32	12	6
1:A:23:MET:SD	1:A:24:LEU:HD22	0.64	2.33	5	17
1:A:43:ILE:HD13	1:A:73:VAL:HA	0.64	1.69	10	8
1:A:28:LEU:HD23	1:A:32:ILE:HD11	0.64	1.70	1	12
1:A:12:LEU:HD11	1:A:24:LEU:HB3	0.63	1.69	29	12

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:27:ARG:NH1	1:A:30:PRO:HB2	0.63	2.07	24	3
1:A:42:LYS:NZ	2:B:12:GLU:HA	0.62	2.10	2	2
1:A:65:LEU:HD13	1:A:66:ARG:N	0.61	2.10	9	18
1:A:28:LEU:HD21	1:A:47:LEU:HB2	0.61	1.71	23	4
1:A:42:LYS:HZ1	2:B:12:GLU:HA	0.61	1.56	17	1
1:A:43:ILE:HD11	1:A:76:LEU:CD1	0.61	2.26	11	4
1:A:27:ARG:NE	1:A:59:LEU:HD22	0.61	2.11	30	2
1:A:75:VAL:HG12	2:B:5:ASN:HB2	0.61	1.71	20	4
1:A:50:ILE:HD11	1:A:68:LYS:HZ2	0.61	1.56	19	1
1:A:15:ALA:O	1:A:19:GLU:HB2	0.61	1.96	10	3
1:A:58:MET:HE2	1:A:68:LYS:HE2	0.61	1.70	20	4
1:A:19:GLU:HG3	1:A:22:GLN:HB3	0.60	1.73	17	8
1:A:15:ALA:HB1	1:A:19:GLU:HB2	0.60	1.73	16	5
1:A:68:LYS:HG2	2:B:6:LEU:HD21	0.60	1.74	14	3
1:A:20:GLN:HG2	1:A:21:LYS:N	0.60	2.12	18	3
1:A:47:LEU:HD13	2:B:7:ASN:HB2	0.59	1.73	21	2
1:A:31:LEU:HD13	1:A:65:LEU:CD1	0.59	2.28	23	8
1:A:21:LYS:HD2	1:A:22:GLN:N	0.58	2.12	18	2
1:A:39:LEU:HB2	1:A:76:LEU:HD13	0.58	1.75	26	2
1:A:15:ALA:HB1	1:A:19:GLU:CB	0.58	2.28	24	10
1:A:15:ALA:HB3	1:A:23:MET:HE3	0.58	1.75	6	10
1:A:65:LEU:HG	1:A:66:ARG:N	0.58	2.14	8	11
2:B:9:ASN:O	2:B:11:LYS:HG2	0.58	1.98	20	8
1:A:12:LEU:HA	1:A:23:MET:HE2	0.58	1.74	16	9
1:A:31:LEU:HD13	1:A:65:LEU:HD11	0.58	1.76	23	7
1:A:70:ASP:O	1:A:73:VAL:HG22	0.57	1.99	14	17
1:A:58:MET:SD	1:A:68:LYS:HE2	0.57	2.39	14	6
1:A:46:MET:HA	2:B:11:LYS:C	0.57	2.24	21	5
1:A:46:MET:HB2	2:B:10:ALA:HB1	0.57	1.75	21	3
1:A:20:GLN:HE21	1:A:21:LYS:HG3	0.57	1.59	19	1
1:A:24:LEU:HG	1:A:27:ARG:HA	0.57	1.75	23	9
1:A:32:ILE:CG2	1:A:43:ILE:HD12	0.57	2.30	10	12
1:A:47:LEU:HD11	2:B:6:LEU:HB3	0.57	1.74	21	3
1:A:46:MET:HE2	1:A:75:VAL:HG23	0.57	1.77	10	8
1:A:39:LEU:HB3	1:A:76:LEU:HG	0.57	1.76	24	7
1:A:51:ASP:HB2	1:A:54:GLU:HB2	0.57	1.76	29	4
1:A:24:LEU:HD23	1:A:25:GLY:N	0.56	2.15	18	18
1:A:45:GLY:O	1:A:48:LEU:HB2	0.56	2.01	30	25
1:A:63:GLU:HG3	1:A:64:SER:N	0.56	2.15	29	6
1:A:27:ARG:CZ	1:A:59:LEU:HD12	0.56	2.31	11	5
1:A:26:GLU:HG2	1:A:50:ILE:HG22	0.56	1.75	18	5

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:24:LEU:HB2	1:A:30:PRO:CG	0.56	2.31	29	10
1:A:20:GLN:HE22	2:B:18:LYS:HB3	0.56	1.61	19	2
1:A:39:LEU:O	1:A:43:ILE:HG13	0.56	2.01	25	11
1:A:29:PHE:CZ	1:A:41:GLY:HA2	0.56	2.36	10	28
1:A:31:LEU:CD1	1:A:69:VAL:HG21	0.56	2.31	8	2
1:A:33:GLN:O	1:A:36:HIS:HB3	0.56	2.01	14	28
1:A:36:HIS:N	1:A:37:PRO:CD	0.56	2.69	26	28
1:A:38:THR:HG22	1:A:76:LEU:HD12	0.56	1.75	7	5
1:A:58:MET:SD	1:A:65:LEU:HD23	0.55	2.42	14	3
1:A:28:LEU:HA	1:A:31:LEU:HD11	0.55	1.79	8	2
1:A:27:ARG:NH2	1:A:59:LEU:HD13	0.55	2.16	24	1
1:A:46:MET:HE3	1:A:75:VAL:HG12	0.55	1.77	4	1
1:A:61:SER:O	1:A:65:LEU:HB3	0.55	2.02	1	29
1:A:47:LEU:HD11	2:B:6:LEU:HB2	0.55	1.77	3	1
1:A:20:GLN:HA	1:A:23:MET:HE3	0.54	1.78	12	2
1:A:58:MET:SD	1:A:65:LEU:HA	0.54	2.43	4	4
1:A:25:GLY:HA3	1:A:48:LEU:HD21	0.54	1.79	24	2
1:A:65:LEU:HD22	1:A:65:LEU:O	0.54	2.02	2	19
1:A:20:GLN:NE2	1:A:21:LYS:HG3	0.54	2.18	19	6
1:A:24:LEU:HD23	1:A:24:LEU:C	0.54	2.27	18	11
1:A:12:LEU:HD21	1:A:27:ARG:NH1	0.54	2.18	7	5
1:A:42:LYS:HE2	1:A:46:MET:SD	0.54	2.43	25	8
1:A:69:VAL:O	1:A:72:ALA:HB3	0.54	2.02	26	23
1:A:68:LYS:HG2	2:B:6:LEU:HD22	0.53	1.80	17	1
1:A:27:ARG:O	1:A:31:LEU:HG	0.53	2.03	23	30
1:A:21:LYS:HD3	1:A:22:GLN:N	0.53	2.18	3	12
1:A:52:ASN:O	1:A:55:LEU:HG	0.53	2.03	18	23
1:A:23:MET:SD	1:A:24:LEU:HG	0.53	2.44	11	12
1:A:24:LEU:HD22	1:A:27:ARG:HD3	0.53	1.81	8	3
1:A:56:LEU:HA	1:A:59:LEU:HD21	0.53	1.80	20	3
1:A:35:MET:HE3	1:A:69:VAL:CG1	0.53	2.33	29	1
1:A:43:ILE:HG12	1:A:76:LEU:CD2	0.53	2.34	28	8
1:A:24:LEU:C	1:A:24:LEU:HD23	0.53	2.29	21	7
1:A:27:ARG:O	1:A:30:PRO:HG2	0.53	2.04	23	12
1:A:21:LYS:HB3	2:B:18:LYS:HA	0.53	1.80	19	7
1:A:38:THR:HG21	1:A:73:VAL:CG1	0.53	2.34	21	1
1:A:68:LYS:O	2:B:6:LEU:HD11	0.52	2.03	5	4
1:A:31:LEU:HD13	1:A:65:LEU:CD2	0.52	2.31	7	8
1:A:25:GLY:CA	1:A:48:LEU:HD21	0.52	2.34	23	2
1:A:35:MET:HE1	1:A:66:ARG:HG2	0.52	1.79	29	1
2:B:7:ASN:HD22	2:B:9:ASN:H	0.52	1.47	21	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:42:LYS:HE3	1:A:46:MET:SD	0.52	2.44	17	2
1:A:28:LEU:CD2	1:A:44:THR:HA	0.52	2.34	23	2
1:A:20:GLN:HE22	1:A:21:LYS:HG3	0.52	1.64	3	2
1:A:42:LYS:HE3	2:B:12:GLU:HG2	0.52	1.80	26	4
1:A:38:THR:O	1:A:39:LEU:HG	0.52	2.04	26	1
1:A:24:LEU:HB2	1:A:30:PRO:HG2	0.52	1.81	28	3
1:A:39:LEU:HD23	1:A:76:LEU:HD11	0.51	1.80	21	2
1:A:32:ILE:HD13	1:A:43:ILE:HB	0.51	1.81	22	1
1:A:27:ARG:CG	1:A:59:LEU:HD23	0.51	2.35	1	5
1:A:35:MET:C	1:A:37:PRO:HD2	0.51	2.30	30	25
1:A:52:ASN:CA	1:A:55:LEU:HD23	0.51	2.35	26	21
1:A:28:LEU:HA	1:A:31:LEU:CD1	0.51	2.35	30	3
1:A:44:THR:O	1:A:48:LEU:HB2	0.51	2.05	24	1
1:A:24:LEU:HB2	1:A:30:PRO:CD	0.51	2.35	19	14
1:A:67:SER:O	1:A:70:ASP:HB3	0.51	2.05	19	16
1:A:42:LYS:HZ3	2:B:12:GLU:HA	0.51	1.64	2	1
1:A:72:ALA:HA	1:A:75:VAL:HG13	0.51	1.82	9	5
1:A:42:LYS:HB3	1:A:76:LEU:HD12	0.51	1.81	20	2
1:A:55:LEU:C	1:A:55:LEU:HD12	0.51	2.31	3	30
1:A:27:ARG:NH2	1:A:59:LEU:HD22	0.51	2.21	29	2
1:A:58:MET:HB3	1:A:65:LEU:HG	0.51	1.82	19	4
1:A:35:MET:HE2	1:A:66:ARG:HE	0.51	1.66	6	1
1:A:68:LYS:HG2	2:B:6:LEU:HD11	0.51	1.80	21	1
1:A:42:LYS:HG3	1:A:46:MET:HE2	0.51	1.82	6	1
1:A:27:ARG:HB2	1:A:55:LEU:CD1	0.50	2.36	29	12
1:A:42:LYS:HZ2	1:A:45:GLY:HA3	0.50	1.66	17	2
1:A:35:MET:HE2	1:A:66:ARG:HG2	0.50	1.82	14	1
1:A:43:ILE:HA	1:A:76:LEU:HD21	0.50	1.83	10	1
1:A:27:ARG:HD2	1:A:30:PRO:HG2	0.50	1.83	9	6
1:A:26:GLU:CD	1:A:50:ILE:HG22	0.50	2.31	29	1
1:A:72:ALA:O	1:A:75:VAL:HB	0.50	2.07	4	1
1:A:29:PHE:O	1:A:33:GLN:N	0.50	2.45	26	20
1:A:42:LYS:O	1:A:46:MET:HG2	0.50	2.06	9	2
1:A:46:MET:HA	2:B:11:LYS:O	0.50	2.07	21	2
1:A:42:LYS:HG3	1:A:46:MET:HE3	0.50	1.82	22	1
1:A:59:LEU:HD22	1:A:60:GLU:N	0.49	2.22	25	3
1:A:44:THR:C	1:A:48:LEU:HD12	0.49	2.33	23	1
1:A:75:VAL:O	1:A:79:HIS:HB2	0.49	2.08	11	1
1:A:46:MET:CE	1:A:47:LEU:HD23	0.49	2.30	21	1
1:A:49:GLU:N	1:A:49:GLU:OE1	0.49	2.45	23	1
1:A:46:MET:SD	1:A:75:VAL:HG21	0.49	2.48	22	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:29:PHE:CD2	1:A:44:THR:HG21	0.49	2.41	14	22
1:A:31:LEU:HD22	1:A:65:LEU:CD1	0.49	2.37	8	1
1:A:42:LYS:NZ	2:B:12:GLU:HG2	0.49	2.23	10	3
1:A:49:GLU:OE2	2:B:11:LYS:HG3	0.49	2.07	24	6
1:A:50:ILE:HD11	1:A:68:LYS:HE3	0.49	1.85	9	6
1:A:28:LEU:HA	1:A:31:LEU:HD12	0.49	1.84	7	6
1:A:24:LEU:HD12	1:A:27:ARG:CD	0.49	2.37	5	8
1:A:35:MET:HE3	1:A:69:VAL:HB	0.49	1.84	29	1
1:A:43:ILE:HG12	1:A:76:LEU:HG	0.49	1.85	10	2
1:A:55:LEU:O	1:A:58:MET:HB2	0.49	2.08	23	3
1:A:58:MET:SD	1:A:58:MET:N	0.48	2.86	7	9
1:A:59:LEU:H	1:A:59:LEU:HD23	0.48	1.68	27	3
1:A:54:GLU:O	1:A:58:MET:HG2	0.48	2.08	2	8
1:A:47:LEU:HD21	2:B:6:LEU:HB3	0.48	1.86	9	1
1:A:28:LEU:CD1	1:A:44:THR:HA	0.48	2.37	29	1
1:A:48:LEU:C	1:A:50:ILE:H	0.48	2.15	23	8
1:A:31:LEU:HD22	1:A:65:LEU:CG	0.48	2.38	22	10
1:A:53:SER:O	1:A:56:LEU:HG	0.48	2.08	23	3
1:A:42:LYS:HE2	1:A:46:MET:HG2	0.48	1.85	21	1
2:B:11:LYS:HG2	2:B:19:TYR:CE2	0.48	2.43	28	3
1:A:71:GLU:HG2	2:B:5:ASN:HB3	0.48	1.85	21	1
1:A:32:ILE:HG23	1:A:69:VAL:CG1	0.48	2.39	28	5
1:A:11:MET:HE3	1:A:23:MET:CB	0.48	2.38	12	5
1:A:28:LEU:C	1:A:30:PRO:HD2	0.48	2.34	8	8
1:A:35:MET:HE1	1:A:69:VAL:HB	0.48	1.86	23	6
1:A:28:LEU:HD11	1:A:47:LEU:HB2	0.47	1.86	4	2
1:A:20:GLN:NE2	2:B:18:LYS:HG2	0.47	2.23	18	1
1:A:56:LEU:O	1:A:59:LEU:HG	0.47	2.10	2	3
1:A:56:LEU:HA	1:A:59:LEU:HD11	0.47	1.85	23	3
1:A:61:SER:HB3	1:A:64:SER:HB2	0.47	1.87	17	8
1:A:46:MET:HG3	1:A:47:LEU:HD12	0.47	1.85	24	1
1:A:30:PRO:O	1:A:33:GLN:HB3	0.47	2.10	15	13
1:A:12:LEU:HD13	1:A:27:ARG:HH12	0.47	1.70	3	1
1:A:28:LEU:HD12	1:A:32:ILE:HD11	0.47	1.85	3	5
1:A:24:LEU:HB2	1:A:30:PRO:HG3	0.47	1.87	6	5
1:A:27:ARG:CZ	1:A:59:LEU:HD13	0.47	2.39	20	4
1:A:12:LEU:HD21	1:A:27:ARG:HH12	0.47	1.70	7	1
1:A:50:ILE:HD11	1:A:68:LYS:HZ3	0.47	1.67	7	1
1:A:43:ILE:HG12	1:A:76:LEU:HD21	0.47	1.87	18	1
1:A:48:LEU:HD13	2:B:13:PHE:CG	0.47	2.45	24	1
1:A:27:ARG:NH2	1:A:59:LEU:HG	0.47	2.25	28	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:28:LEU:O	1:A:32:ILE:HG13	0.47	2.10	26	15
1:A:29:PHE:N	1:A:30:PRO:HD2	0.47	2.25	5	12
1:A:32:ILE:O	1:A:40:ALA:HB2	0.47	2.10	22	6
1:A:48:LEU:HG	2:B:13:PHE:CD1	0.47	2.44	12	4
1:A:47:LEU:N	1:A:47:LEU:HD22	0.46	2.25	4	4
1:A:21:LYS:HD3	1:A:22:GLN:H	0.46	1.69	3	1
1:A:12:LEU:HD11	1:A:24:LEU:HD23	0.46	1.86	26	1
1:A:45:GLY:CA	2:B:13:PHE:HB2	0.46	2.40	2	5
1:A:31:LEU:HD22	1:A:65:LEU:CD2	0.46	2.41	21	4
1:A:45:GLY:O	2:B:11:LYS:O	0.46	2.33	29	5
1:A:59:LEU:HD12	1:A:59:LEU:O	0.46	2.11	24	1
1:A:20:GLN:HE21	1:A:21:LYS:HB2	0.46	1.69	26	1
1:A:20:GLN:NE2	2:B:18:LYS:HG3	0.46	2.26	26	1
1:A:51:ASP:HB2	1:A:54:GLU:HB3	0.46	1.88	23	1
1:A:76:LEU:C	1:A:76:LEU:HD13	0.46	2.36	23	8
1:A:12:LEU:CD2	1:A:24:LEU:HB3	0.46	2.40	5	1
1:A:36:HIS:ND1	1:A:37:PRO:HD3	0.46	2.25	7	10
1:A:24:LEU:HB2	1:A:30:PRO:HD3	0.46	1.88	19	1
1:A:48:LEU:HD13	2:B:13:PHE:CD1	0.46	2.46	24	1
1:A:32:ILE:HA	1:A:35:MET:HG3	0.46	1.88	29	1
1:A:47:LEU:O	1:A:50:ILE:HB	0.45	2.11	23	2
1:A:46:MET:CB	2:B:10:ALA:HB1	0.45	2.41	21	1
1:A:32:ILE:O	1:A:36:HIS:N	0.45	2.50	21	28
1:A:43:ILE:HG12	1:A:76:LEU:HB2	0.45	1.88	23	3
1:A:71:GLU:O	1:A:74:ALA:HB3	0.45	2.12	22	9
1:A:63:GLU:O	1:A:67:SER:HB2	0.45	2.11	15	1
1:A:48:LEU:HG	2:B:13:PHE:CE1	0.45	2.46	5	2
1:A:19:GLU:HA	1:A:19:GLU:OE1	0.45	2.10	11	2
1:A:21:LYS:HE2	1:A:22:GLN:N	0.45	2.26	4	2
1:A:20:GLN:HE22	2:B:18:LYS:HG2	0.45	1.71	3	4
1:A:38:THR:CG2	1:A:73:VAL:HG13	0.45	2.40	3	4
1:A:19:GLU:O	1:A:22:GLN:HB3	0.45	2.12	29	5
1:A:42:LYS:HZ3	2:B:12:GLU:HG2	0.45	1.71	17	1
1:A:47:LEU:H	1:A:47:LEU:HD22	0.45	1.72	19	1
1:A:53:SER:HA	1:A:56:LEU:HD21	0.45	1.88	20	1
1:A:56:LEU:HA	1:A:59:LEU:CD2	0.45	2.42	20	1
1:A:36:HIS:O	1:A:37:PRO:C	0.44	2.60	15	21
1:A:42:LYS:HB3	1:A:76:LEU:CD1	0.44	2.42	14	1
1:A:19:GLU:OE1	1:A:19:GLU:HA	0.44	2.13	20	2
1:A:58:MET:HE3	1:A:65:LEU:CA	0.44	2.42	25	1
1:A:47:LEU:HD11	2:B:7:ASN:H	0.44	1.73	21	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:27:ARG:HE	1:A:59:LEU:HD22	0.44	1.71	30	2
1:A:49:GLU:HB2	2:B:7:ASN:ND2	0.44	2.27	1	2
1:A:60:GLU:O	1:A:60:GLU:HG3	0.44	2.11	14	3
1:A:19:GLU:C	1:A:23:MET:HE3	0.44	2.37	28	2
1:A:57:HIS:HA	1:A:60:GLU:HG2	0.44	1.88	2	2
1:A:65:LEU:C	1:A:65:LEU:CD2	0.44	2.90	18	10
1:A:20:GLN:O	1:A:24:LEU:HD12	0.44	2.13	26	5
1:A:42:LYS:HG3	1:A:46:MET:SD	0.44	2.52	11	2
1:A:24:LEU:HD13	1:A:26:GLU:O	0.44	2.13	10	5
1:A:15:ALA:HB3	1:A:23:MET:HE2	0.44	1.89	21	1
1:A:20:GLN:HE22	2:B:18:LYS:HG3	0.43	1.73	27	1
1:A:35:MET:HE2	1:A:66:ARG:HG3	0.43	1.90	15	2
1:A:74:ALA:O	1:A:77:GLN:HG3	0.43	2.14	22	2
1:A:24:LEU:CD1	1:A:27:ARG:HD3	0.43	2.42	6	1
1:A:55:LEU:HG	1:A:56:LEU:N	0.43	2.28	20	4
1:A:43:ILE:CD1	1:A:76:LEU:HD11	0.43	2.42	22	1
1:A:35:MET:HE1	1:A:66:ARG:HA	0.43	1.91	29	1
1:A:35:MET:HE2	1:A:66:ARG:NE	0.43	2.27	6	1
1:A:24:LEU:HA	1:A:30:PRO:HD3	0.43	1.90	8	1
1:A:58:MET:N	1:A:58:MET:SD	0.43	2.92	18	2
1:A:15:ALA:CB	1:A:23:MET:HE2	0.43	2.44	21	2
1:A:49:GLU:OE2	2:B:9:ASN:HB3	0.43	2.14	9	1
1:A:22:GLN:OE1	1:A:23:MET:HG3	0.43	2.14	11	1
1:A:36:HIS:C	1:A:38:THR:N	0.43	2.77	26	4
1:A:46:MET:HE1	1:A:75:VAL:CG2	0.43	2.35	21	1
1:A:36:HIS:O	1:A:38:THR:N	0.43	2.52	15	9
1:A:12:LEU:HD23	1:A:23:MET:HE2	0.43	1.90	3	1
1:A:68:LYS:O	2:B:6:LEU:HD22	0.43	2.14	24	2
1:A:42:LYS:HZ1	2:B:12:GLU:HG2	0.43	1.73	10	1
1:A:58:MET:CA	1:A:65:LEU:HB2	0.43	2.44	27	3
1:A:58:MET:HA	1:A:65:LEU:HB2	0.42	1.91	24	5
1:A:54:GLU:O	1:A:58:MET:SD	0.42	2.77	25	1
1:A:48:LEU:HB3	2:B:19:TYR:HD2	0.42	1.73	26	1
1:A:39:LEU:O	1:A:40:ALA:C	0.42	2.61	1	4
1:A:46:MET:O	2:B:10:ALA:HA	0.42	2.13	4	1
1:A:26:GLU:O	1:A:55:LEU:HD13	0.42	2.14	8	1
1:A:39:LEU:HB3	1:A:76:LEU:CD1	0.42	2.44	14	1
1:A:28:LEU:O	1:A:29:PHE:C	0.42	2.62	23	6
1:A:43:ILE:HG22	1:A:72:ALA:HB1	0.42	1.90	1	1
1:A:49:GLU:HG3	2:B:9:ASN:O	0.42	2.15	12	1
1:A:49:GLU:HG2	2:B:11:LYS:HZ2	0.42	1.74	23	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:26:GLU:HB3	1:A:55:LEU:HB2	0.42	1.91	26	1
1:A:11:MET:O	1:A:15:ALA:N	0.42	2.53	22	4
1:A:26:GLU:HB3	1:A:55:LEU:CB	0.42	2.45	26	1
1:A:46:MET:HB2	1:A:47:LEU:HD22	0.42	1.91	4	1
1:A:27:ARG:CZ	1:A:59:LEU:HG	0.42	2.45	8	1
1:A:20:GLN:CG	1:A:21:LYS:N	0.42	2.82	17	3
2:B:7:ASN:ND2	2:B:8:PRO:HD2	0.42	2.29	21	1
1:A:47:LEU:CD1	2:B:7:ASN:HB2	0.42	2.44	21	2
1:A:28:LEU:HB3	1:A:44:THR:CB	0.42	2.43	29	1
1:A:49:GLU:CD	2:B:11:LYS:HG3	0.42	2.40	29	1
1:A:27:ARG:O	1:A:30:PRO:HD2	0.42	2.15	16	1
1:A:59:LEU:H	1:A:59:LEU:CD1	0.42	2.25	23	1
1:A:57:HIS:O	1:A:60:GLU:HG2	0.42	2.15	1	3
1:A:8:THR:O	1:A:12:LEU:HG	0.41	2.15	3	3
1:A:27:ARG:NE	1:A:59:LEU:HD23	0.41	2.30	4	1
1:A:28:LEU:HD13	1:A:48:LEU:HA	0.41	1.90	19	1
1:A:75:VAL:CG2	1:A:76:LEU:N	0.41	2.83	21	1
1:A:9:ALA:O	1:A:12:LEU:HB2	0.41	2.15	12	2
1:A:46:MET:HE2	1:A:75:VAL:CG2	0.41	2.45	16	2
1:A:24:LEU:HD12	1:A:27:ARG:HH11	0.41	1.75	17	1
1:A:42:LYS:CE	2:B:12:GLU:HG2	0.41	2.45	25	1
1:A:49:GLU:OE1	2:B:9:ASN:HB3	0.41	2.15	29	1
1:A:58:MET:HE1	1:A:68:LYS:HD3	0.41	1.91	21	1
1:A:32:ILE:HD13	1:A:43:ILE:CG2	0.41	2.46	9	2
1:A:21:LYS:H	1:A:21:LYS:HG3	0.41	1.52	1	1
2:B:9:ASN:HB3	2:B:11:LYS:NZ	0.41	2.31	25	1
1:A:66:ARG:O	1:A:70:ASP:HB2	0.41	2.16	6	1
1:A:59:LEU:CD2	1:A:59:LEU:C	0.41	2.94	14	2
1:A:42:LYS:NZ	1:A:45:GLY:HA3	0.41	2.30	17	1
1:A:44:THR:HG23	1:A:45:GLY:N	0.41	2.31	22	1
1:A:27:ARG:HB2	1:A:55:LEU:HB2	0.41	1.92	8	1
1:A:23:MET:H	1:A:23:MET:HG3	0.41	1.52	12	1
1:A:56:LEU:O	1:A:60:GLU:HG3	0.41	2.16	26	1
1:A:15:ALA:CB	1:A:23:MET:HE3	0.41	2.45	3	1
1:A:71:GLU:O	1:A:75:VAL:HG13	0.41	2.16	6	1
1:A:46:MET:HG3	1:A:47:LEU:HD22	0.41	1.93	8	1
1:A:20:GLN:CD	1:A:21:LYS:HG3	0.41	2.41	9	1
1:A:26:GLU:OE1	1:A:51:ASP:C	0.41	2.64	29	1
1:A:40:ALA:O	1:A:44:THR:HG22	0.41	2.16	5	2
1:A:11:MET:HE3	1:A:23:MET:HA	0.40	1.93	21	2
1:A:71:GLU:HB3	2:B:6:LEU:HD13	0.40	1.93	25	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:21:LYS:HB3	1:A:21:LYS:HE3	0.40	1.67	26	1
1:A:27:ARG:HD2	1:A:30:PRO:CG	0.40	2.46	11	1
1:A:57:HIS:O	1:A:61:SER:HB2	0.40	2.16	22	1
1:A:26:GLU:HG2	1:A:51:ASP:O	0.40	2.17	12	1
1:A:32:ILE:HD12	1:A:43:ILE:HB	0.40	1.92	28	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	69/98 (70%)	55±1 (80±2%)	10±2 (15±2%)	3±1 (5±1%)	3	24
2	B	11/22 (50%)	6±1 (58±9%)	2±2 (21±14%)	2±1 (22±13%)	0	2
All	All	2400/3600 (67%)	1846 (77%)	379 (16%)	175 (7%)	1	15

All 11 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	23	MET	30
1	A	39	LEU	29
2	B	10	ALA	22
2	B	18	LYS	18
1	A	28	LEU	17
2	B	12	GLU	16
2	B	6	LEU	15
1	A	37	PRO	14
1	A	38	THR	8
1	A	15	ALA	3
1	A	24	LEU	3

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	58/79 (73%)	34±3 (58±4%)	24±3 (42±4%)	0	4
2	B	10/19 (53%)	7±1 (67±11%)	3±1 (33±11%)	1	12
All	All	2040/2940 (69%)	1214 (60%)	826 (40%)	0	5

All 56 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	23	MET	30
1	A	31	LEU	30
1	A	32	ILE	30
1	A	48	LEU	30
1	A	55	LEU	30
1	A	65	LEU	30
1	A	21	LYS	29
1	A	46	MET	29
1	A	75	VAL	29
1	A	20	GLN	28
1	A	24	LEU	28
1	A	29	PHE	26
1	A	50	ILE	23
1	A	26	GLU	22
1	A	39	LEU	22
1	A	66	ARG	21
2	B	11	LYS	21
2	B	19	TYR	19
1	A	58	MET	18
2	B	18	LYS	18
1	A	54	GLU	17
1	A	28	LEU	16
1	A	71	GLU	15
2	B	7	ASN	15
1	A	36	HIS	15
1	A	19	GLU	14
1	A	59	LEU	14
1	A	42	LYS	13

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Mol	Chain	Res	Type	Models (Total)
1	A	61	SER	13
1	A	68	LYS	13
1	A	67	SER	12
1	A	53	SER	12
1	A	57	HIS	12
2	B	6	LEU	12
1	A	56	LEU	11
1	A	14	SER	11
1	A	63	GLU	11
1	A	60	GLU	9
1	A	43	ILE	8
2	B	12	GLU	7
1	A	33	GLN	6
1	A	49	GLU	6
1	A	12	LEU	6
1	A	38	THR	6
2	B	5	ASN	5
1	A	27	ARG	5
1	A	73	VAL	5
1	A	77	GLN	5
1	A	76	LEU	4
1	A	10	SER	3
2	B	13	PHE	3
1	A	52	ASN	3
1	A	22	GLN	2
1	A	79	HIS	2
1	A	35	MET	1
1	A	44	THR	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