



Full wwPDB NMR Structure Validation Report ⓘ

Mar 8, 2026 – 02:52 PM UTC

PDB ID : 2KFK / pdb_00002kfk
Title : Solution structure of Bem1p PB1 domain complexed with Cdc24p PB1 domain
Authors : Kobashigawa, Y.; Yoshinaga, S.; Tandai, T.; Ogura, K.; Inagaki, F.
Deposited on : 2009-02-23

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

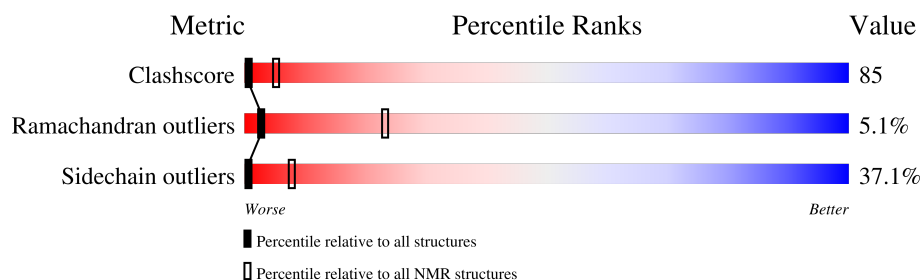
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

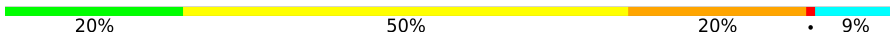
The overall completeness of chemical shifts assignment was not calculated.

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	78	
2	B	86	

2 Ensemble composition and analysis

This entry contains 20 models. Model 18 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:5-A:78, B:104-B:139, B:145-B:186 (152)	0.30	18

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 4 clusters and 1 single-model cluster was found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2700 atoms, of which 1360 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Bud emergence protein 1.

Mol	Chain	Residues	Atoms						Trace
1	A	78	Total	C	H	N	O	S	0
			1297	407	662	105	121	2	

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	PRO	-	expression tag	UNP P29366
A	2	HIS	-	expression tag	UNP P29366
A	3	MET	-	expression tag	UNP P29366

- Molecule 2 is a protein called Cell division control protein 24.

Mol	Chain	Residues	Atoms						Trace
2	B	86	Total	C	H	N	O	S	0
			1403	455	698	111	137	2	

There are 14 discrepancies between the modelled and reference sequences:

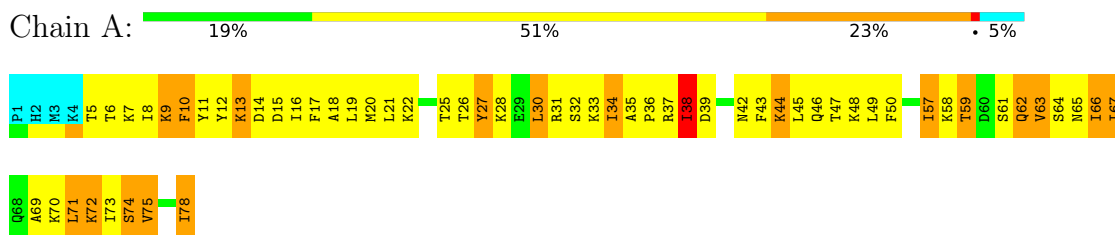
Chain	Residue	Modelled	Actual	Comment	Reference
B	101	PRO	-	expression tag	UNP P11433
B	102	LEU	-	expression tag	UNP P11433
B	103	GLY	-	expression tag	UNP P11433
B	?	-	ASN	deletion	UNP P11433
B	?	-	ASN	deletion	UNP P11433
B	?	-	SER	deletion	UNP P11433
B	?	-	ASN	deletion	UNP P11433
B	?	-	ASN	deletion	UNP P11433
B	?	-	THR	deletion	UNP P11433
B	?	-	SER	deletion	UNP P11433
B	?	-	SER	deletion	UNP P11433
B	?	-	ASN	deletion	UNP P11433
B	?	-	ASN	deletion	UNP P11433
B	?	-	ASN	deletion	UNP P11433

4 Residue-property plots

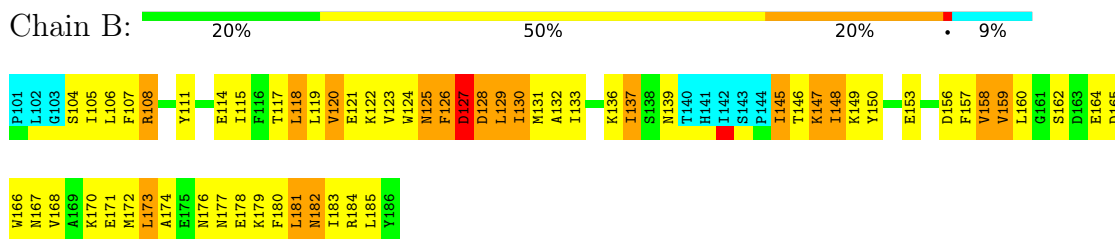
4.1 Average score per residue in the NMR ensemble

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: Bud emergence protein 1



- Molecule 2: Cell division control protein 24

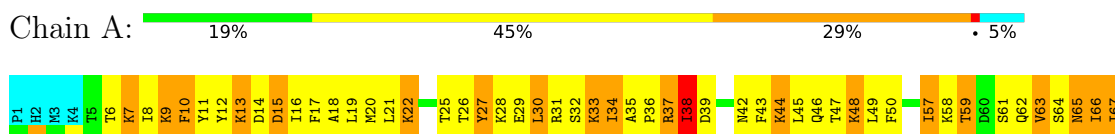


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

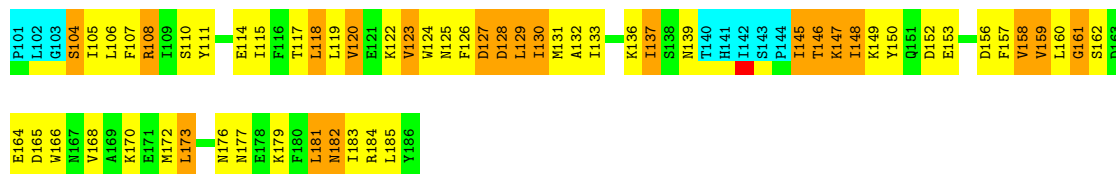
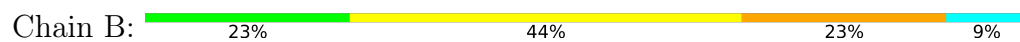
4.2.1 Score per residue for model 1

- Molecule 1: Bud emergence protein 1



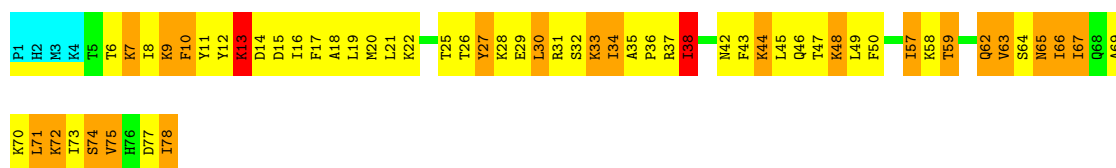
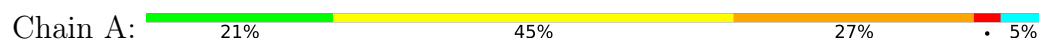


- Molecule 2: Cell division control protein 24

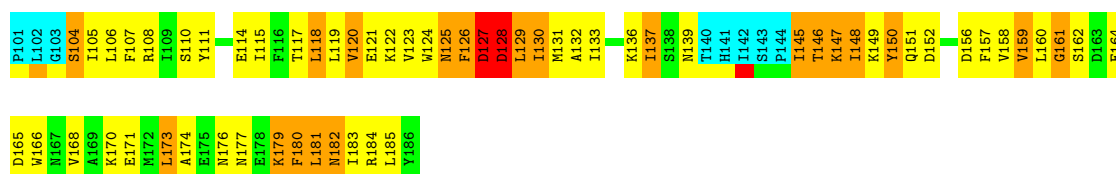
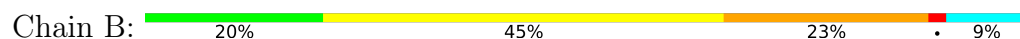


4.2.2 Score per residue for model 2

- Molecule 1: Bud emergence protein 1

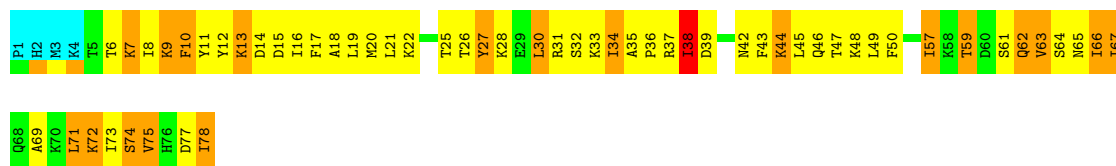


- Molecule 2: Cell division control protein 24

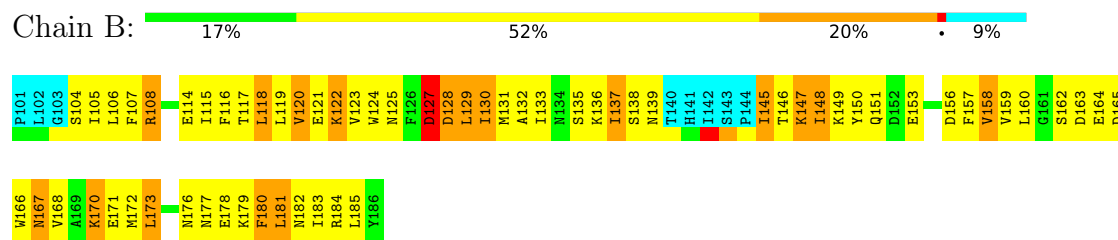


4.2.3 Score per residue for model 3

- Molecule 1: Bud emergence protein 1

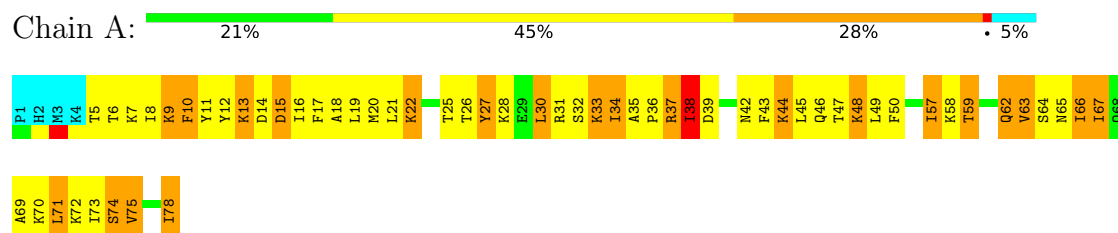


- Molecule 2: Cell division control protein 24

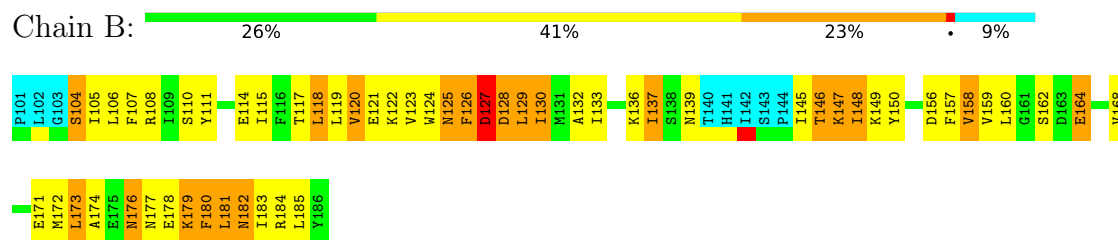


4.2.4 Score per residue for model 4

- Molecule 1: Bud emergence protein 1

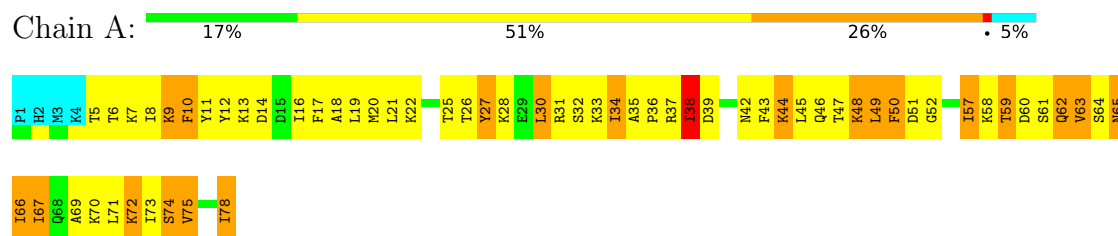


- Molecule 2: Cell division control protein 24

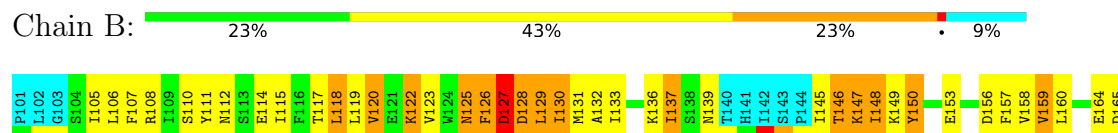


4.2.5 Score per residue for model 5

- Molecule 1: Bud emergence protein 1



- Molecule 2: Cell division control protein 24

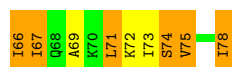
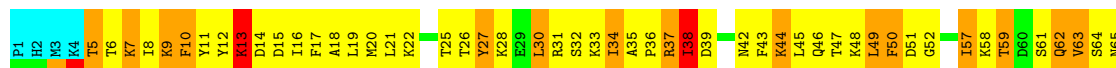




4.2.6 Score per residue for model 6

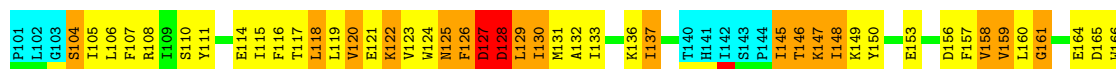
- Molecule 1: Bud emergence protein 1

Chain A: 18% 47% 27% 5%



- Molecule 2: Cell division control protein 24

Chain B: 21% 41% 27% 9%



4.2.7 Score per residue for model 7

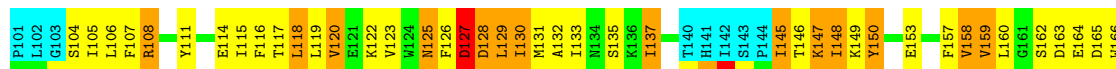
- Molecule 1: Bud emergence protein 1

Chain A: 18% 47% 28% 5%



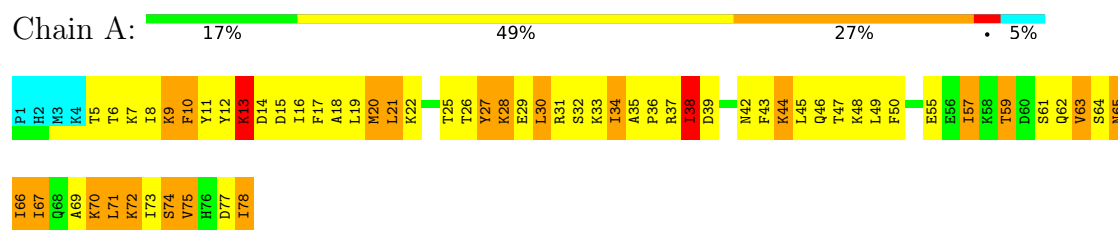
- Molecule 2: Cell division control protein 24

Chain B: 23% 42% 24% 9%

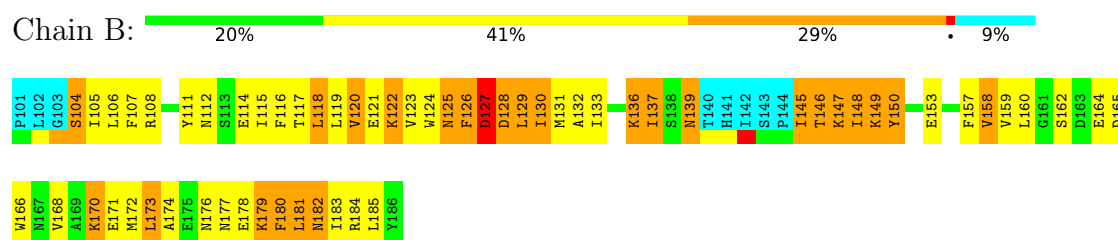


4.2.8 Score per residue for model 8

- Molecule 1: Bud emergence protein 1

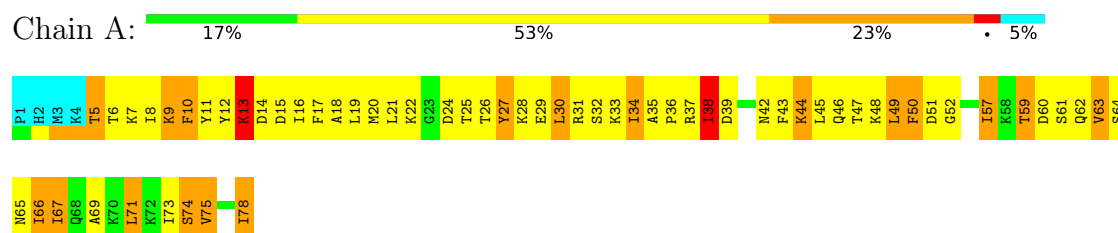


- Molecule 2: Cell division control protein 24



4.2.9 Score per residue for model 9

- Molecule 1: Bud emergence protein 1



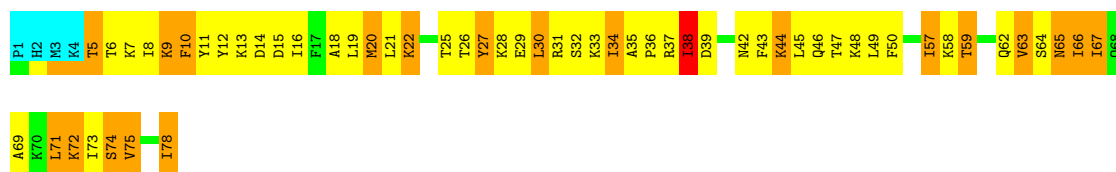
- Molecule 2: Cell division control protein 24



4.2.10 Score per residue for model 10

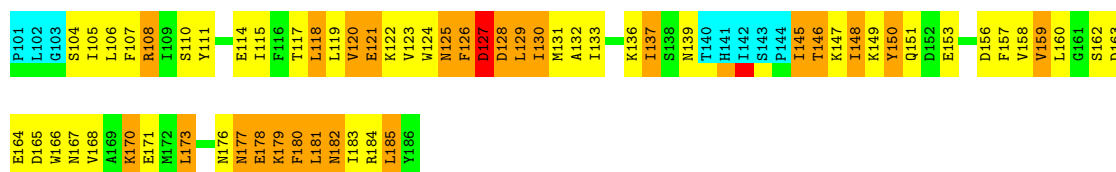
- Molecule 1: Bud emergence protein 1

Chain A: 22% 46% 26% 5%



• Molecule 2: Cell division control protein 24

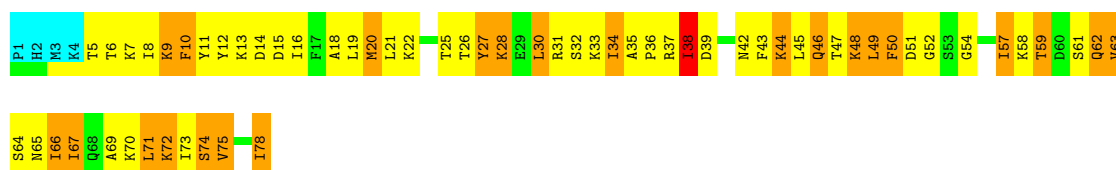
Chain B: 19% 43% 28% 9%



4.2.11 Score per residue for model 11

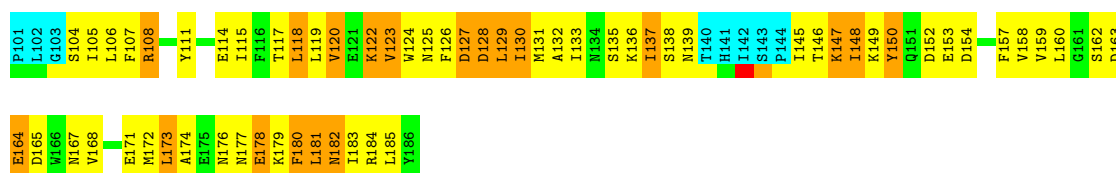
• Molecule 1: Bud emergence protein 1

Chain A: 17% 47% 29% 5%



• Molecule 2: Cell division control protein 24

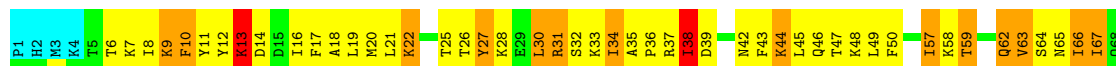
Chain B: 19% 50% 22% 9%



4.2.12 Score per residue for model 12

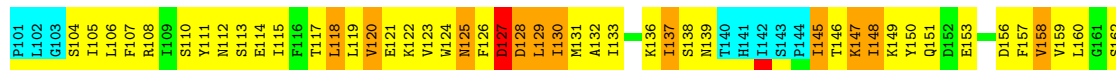
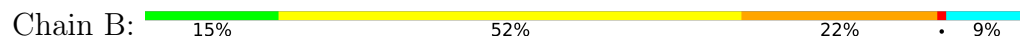
• Molecule 1: Bud emergence protein 1

Chain A: 23% 45% 24% 5%



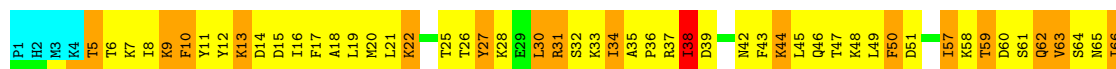
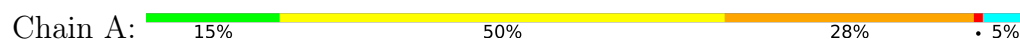


• Molecule 2: Cell division control protein 24



4.2.13 Score per residue for model 13

• Molecule 1: Bud emergence protein 1

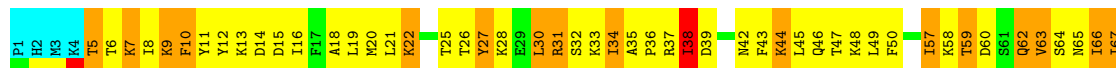
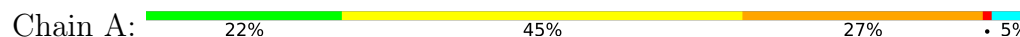


• Molecule 2: Cell division control protein 24



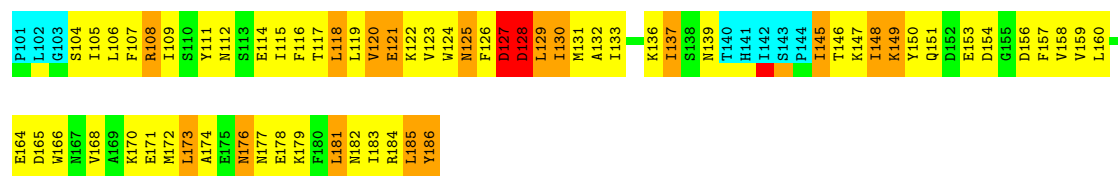
4.2.14 Score per residue for model 14

• Molecule 1: Bud emergence protein 1



• Molecule 2: Cell division control protein 24

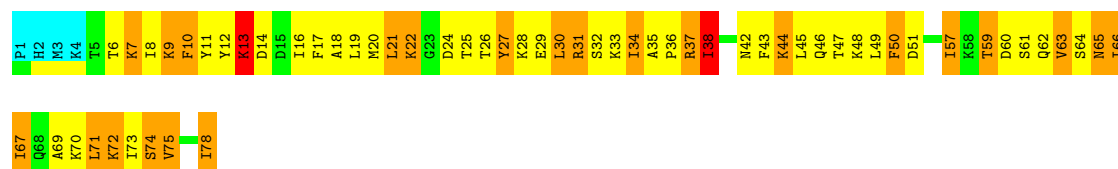
Chain B: 



4.2.15 Score per residue for model 15

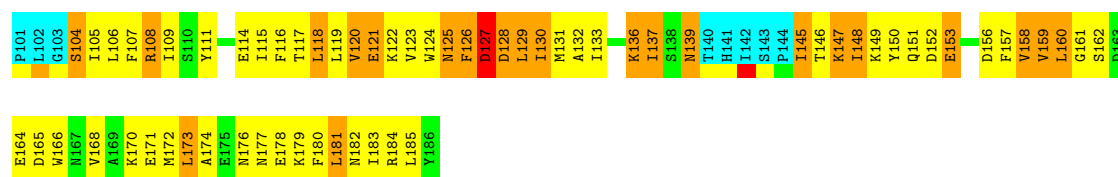
- Molecule 1: Bud emergence protein 1

Chain A: 



- Molecule 2: Cell division control protein 24

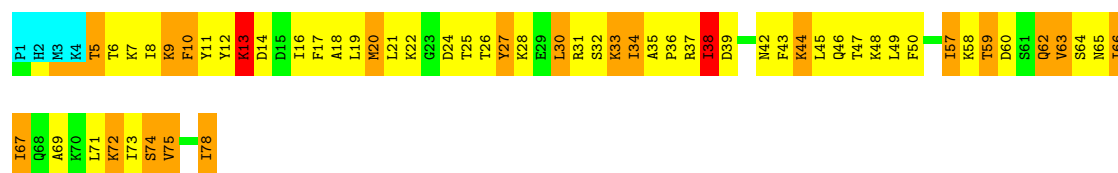
Chain B: 



4.2.16 Score per residue for model 16

- Molecule 1: Bud emergence protein 1

Chain A: 



- Molecule 2: Cell division control protein 24

Chain B: 

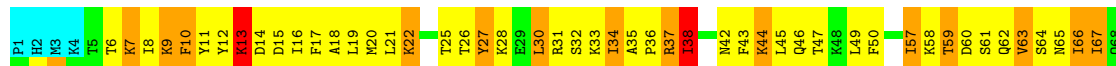




4.2.17 Score per residue for model 17

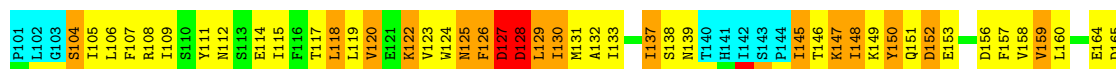
- Molecule 1: Bud emergence protein 1

Chain A: 22% 47% 23% 5%



- Molecule 2: Cell division control protein 24

Chain B: 19% 45% 24% 9%



4.2.18 Score per residue for model 18 (medoid)

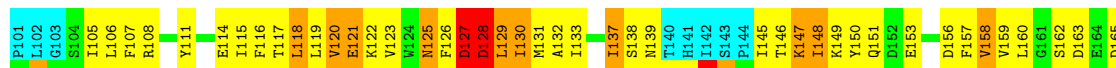
- Molecule 1: Bud emergence protein 1

Chain A: 21% 49% 23% 5%



- Molecule 2: Cell division control protein 24

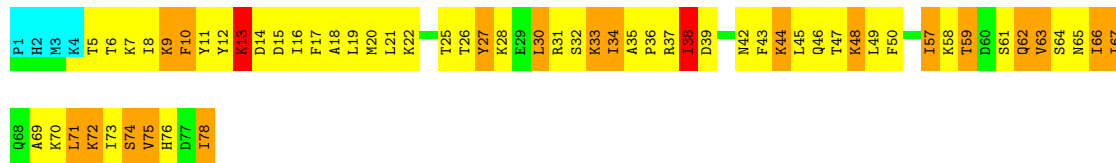
Chain B: 21% 50% 17% 9%



4.2.19 Score per residue for model 19

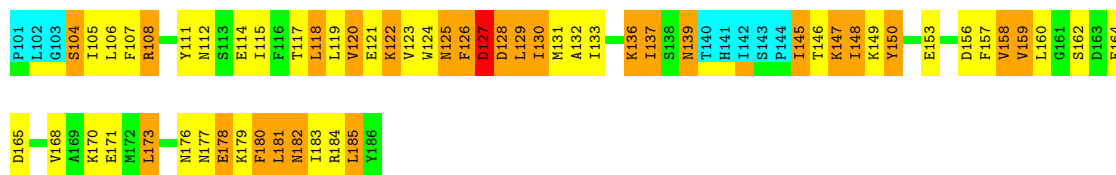
- Molecule 1: Bud emergence protein 1

Chain A: 18% 50% 24% • 5%



- Molecule 2: Cell division control protein 24

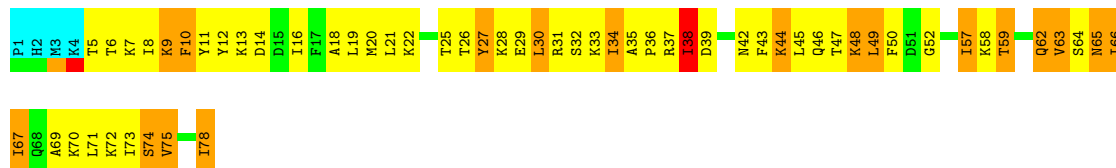
Chain B: 23% 37% 29% • 9%



4.2.20 Score per residue for model 20

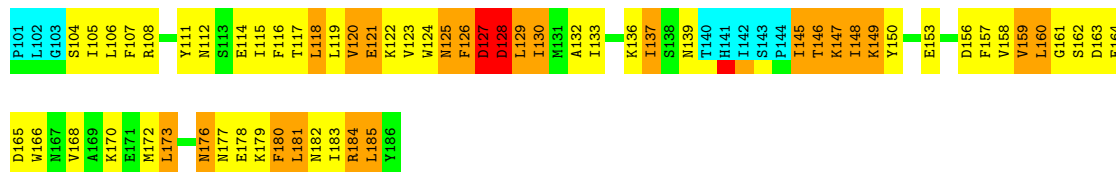
- Molecule 1: Bud emergence protein 1

Chain A: 21% 50% 23% • 5%



- Molecule 2: Cell division control protein 24

Chain B: 20% 44% 24% • 9%



5 Refinement protocol and experimental data overview ⓘ

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

Of the 50 calculated structures, 20 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
CNS	refinement	1.1
CYANA	structure solution	2.0

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	0.43±0.01	0±0/609 (0.0± 0.0%)	0.77±0.01	0±0/814 (0.0± 0.0%)
2	B	0.34±0.01	0±0/659 (0.0± 0.0%)	0.71±0.01	0±0/888 (0.0± 0.0%)
All	All	0.39	0/25360 (0.0%)	0.74	1/34040 (0.0%)

There are no bond-length outliers.

All unique angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
2	B	128	ASP	N-CA-C	-6.61	103.08	112.13	13	1

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	601	623	622	132±6
2	B	648	637	637	93±6
All	All	24980	25200	25180	4287

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 85.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:21:LEU:HD12	1:A:25:THR:HG21	1.03	1.31	15	20
2:B:106:LEU:HD21	2:B:117:THR:HB	0.94	1.38	19	20
1:A:34:ILE:HG22	1:A:38:ILE:HD11	0.91	1.40	15	20
2:B:129:LEU:HD12	2:B:130:ILE:N	0.90	1.82	1	20
1:A:27:TYR:HA	1:A:63:VAL:HG11	0.89	1.44	3	20
1:A:19:LEU:HD13	1:A:33:LYS:HB3	0.87	1.46	6	20
1:A:45:LEU:CD1	1:A:75:VAL:HB	0.86	2.00	18	20
1:A:19:LEU:HD22	1:A:33:LYS:HG2	0.85	1.47	1	6
1:A:10:PHE:CE1	1:A:19:LEU:HD12	0.85	2.06	8	20
2:B:164:GLU:O	2:B:168:VAL:HG23	0.85	1.71	3	18
2:B:114:GLU:O	2:B:115:ILE:HD13	0.85	1.71	11	20
2:B:118:LEU:HD21	2:B:132:ALA:HB1	0.82	1.51	4	20
1:A:69:ALA:HB1	1:A:71:LEU:HD21	0.81	1.52	1	4
1:A:19:LEU:HD21	1:A:37:ARG:HD2	0.81	1.53	1	3
2:B:146:THR:HG23	2:B:147:LYS:CE	0.81	2.06	4	5
1:A:22:LYS:O	1:A:67:ILE:HD12	0.80	1.77	19	20
2:B:173:LEU:O	2:B:173:LEU:HD22	0.79	1.78	17	14
2:B:104:SER:CB	2:B:119:LEU:HD11	0.79	2.07	6	9
1:A:66:ILE:HG21	1:A:73:ILE:HD11	0.77	1.55	1	20
2:B:133:ILE:HG22	2:B:137:ILE:HG12	0.77	1.55	17	20
1:A:18:ALA:HB2	2:B:158:VAL:HG21	0.77	1.54	4	12
1:A:19:LEU:HD21	1:A:37:ARG:HD3	0.76	1.58	2	18
1:A:47:THR:CG2	1:A:66:ILE:HD11	0.76	2.10	11	20
1:A:35:ALA:HB2	1:A:43:PHE:CE1	0.74	2.17	15	20
1:A:18:ALA:HB2	2:B:158:VAL:HG11	0.74	1.58	16	8
1:A:6:THR:O	1:A:20:MET:HA	0.74	1.81	17	20
2:B:105:ILE:HD11	2:B:122:LYS:HD3	0.74	1.57	9	7
2:B:160:LEU:HD23	2:B:165:ASP:HB3	0.74	1.59	10	10
2:B:153:GLU:HG2	2:B:172:MET:HE1	0.73	1.58	16	3
2:B:108:ARG:HD2	2:B:117:THR:HG22	0.73	1.60	20	1
1:A:19:LEU:HD21	1:A:37:ARG:CD	0.73	2.13	6	2
1:A:16:ILE:HD12	2:B:157:PHE:O	0.73	1.83	9	20
2:B:160:LEU:HD23	2:B:165:ASP:CG	0.72	2.08	1	7
1:A:19:LEU:HD13	1:A:33:LYS:CB	0.72	2.13	17	11
1:A:34:ILE:HD13	1:A:45:LEU:HG	0.72	1.60	6	20
2:B:172:MET:HE3	2:B:172:MET:O	0.72	1.84	20	5
2:B:104:SER:HB2	2:B:119:LEU:HD11	0.72	1.62	10	11
1:A:7:LYS:HE2	1:A:18:ALA:HB1	0.71	1.62	1	1
1:A:10:PHE:CD2	1:A:34:ILE:HD11	0.71	2.20	6	20
2:B:172:MET:HE2	2:B:172:MET:O	0.71	1.85	6	1
1:A:34:ILE:CG2	1:A:38:ILE:HD11	0.71	2.15	15	20
2:B:127:ASP:O	2:B:130:ILE:N	0.71	2.23	9	19

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:10:PHE:CE2	1:A:73:ILE:HD12	0.70	2.21	15	20
1:A:10:PHE:CZ	1:A:19:LEU:HD12	0.70	2.21	11	20
2:B:111:TYR:OH	2:B:145:ILE:HD11	0.70	1.85	19	1
1:A:27:TYR:CA	1:A:63:VAL:HG11	0.69	2.17	12	20
2:B:133:ILE:HD13	2:B:148:ILE:HG21	0.69	1.64	14	20
2:B:105:ILE:HD11	2:B:122:LYS:HG2	0.69	1.65	14	10
1:A:5:THR:OG1	1:A:20:MET:HB2	0.69	1.87	6	3
2:B:173:LEU:HD22	2:B:173:LEU:C	0.69	2.13	20	13
2:B:129:LEU:HD12	2:B:129:LEU:C	0.68	2.13	13	20
2:B:104:SER:HB3	2:B:119:LEU:HD11	0.68	1.65	6	3
2:B:106:LEU:HD21	2:B:117:THR:CB	0.68	2.17	19	1
1:A:42:ASN:C	1:A:78:ILE:HG13	0.68	2.14	19	1
1:A:35:ALA:HB2	1:A:43:PHE:CZ	0.67	2.24	15	20
2:B:118:LEU:HD13	2:B:120:VAL:HG12	0.67	1.66	13	20
2:B:146:THR:HG23	2:B:147:LYS:HD2	0.67	1.67	5	10
1:A:16:ILE:O	2:B:159:VAL:HG13	0.67	1.89	10	10
2:B:173:LEU:HD21	2:B:179:LYS:H	0.66	1.50	15	16
2:B:118:LEU:HD23	2:B:136:LYS:HD3	0.66	1.66	15	1
1:A:25:THR:HG23	1:A:29:GLU:CG	0.66	2.21	15	5
2:B:146:THR:HG23	2:B:147:LYS:HE2	0.65	1.67	4	5
2:B:106:LEU:CD2	2:B:117:THR:HB	0.65	2.20	19	13
1:A:5:THR:OG1	1:A:20:MET:HE2	0.65	1.91	11	1
2:B:145:ILE:HG21	2:B:185:LEU:HD11	0.64	1.69	2	6
1:A:21:LEU:CD1	1:A:25:THR:HG21	0.64	2.19	8	6
2:B:173:LEU:C	2:B:173:LEU:CD2	0.64	2.70	5	15
1:A:16:ILE:O	2:B:159:VAL:HG23	0.64	1.92	14	8
2:B:149:LYS:HE2	2:B:159:VAL:HG12	0.64	1.70	6	4
2:B:105:ILE:HD12	2:B:105:ILE:N	0.64	2.08	4	19
2:B:106:LEU:HD11	2:B:117:THR:CG2	0.64	2.23	13	8
1:A:6:THR:HG21	1:A:67:ILE:O	0.63	1.93	8	20
2:B:160:LEU:CD2	2:B:165:ASP:HB3	0.63	2.24	17	12
1:A:5:THR:HG22	1:A:22:LYS:HD2	0.63	1.70	18	3
2:B:118:LEU:HD11	2:B:132:ALA:CB	0.63	2.23	14	20
1:A:16:ILE:O	2:B:159:VAL:N	0.63	2.31	13	20
1:A:21:LEU:HD13	1:A:33:LYS:CE	0.63	2.24	8	4
1:A:7:LYS:HD3	2:B:172:MET:HE3	0.63	1.69	18	1
1:A:66:ILE:HG13	1:A:71:LEU:HD12	0.63	1.71	20	4
2:B:149:LYS:CE	2:B:159:VAL:HG12	0.63	2.23	15	2
1:A:9:LYS:HD3	1:A:16:ILE:HG21	0.62	1.69	2	20
1:A:25:THR:HG23	1:A:29:GLU:HG2	0.62	1.71	20	6
2:B:108:ARG:HD3	2:B:117:THR:HG22	0.62	1.71	19	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:158:VAL:HG23	2:B:159:VAL:N	0.62	2.07	7	12
2:B:145:ILE:HG22	2:B:146:THR:H	0.62	1.54	4	3
1:A:16:ILE:CG2	2:B:158:VAL:HG22	0.62	2.25	16	4
2:B:105:ILE:HD11	2:B:122:LYS:CG	0.62	2.24	20	12
1:A:21:LEU:N	1:A:21:LEU:HD22	0.62	2.09	15	19
1:A:34:ILE:HG21	1:A:45:LEU:HD11	0.62	1.72	10	20
2:B:160:LEU:HD23	2:B:165:ASP:CB	0.62	2.23	10	12
1:A:21:LEU:HD12	1:A:25:THR:CG2	0.62	2.19	15	5
2:B:106:LEU:C	2:B:106:LEU:HD13	0.62	2.20	4	20
1:A:57:ILE:HG22	1:A:62:GLN:HB2	0.61	1.71	2	20
2:B:148:ILE:HD11	2:B:160:LEU:HD12	0.61	1.70	4	11
1:A:6:THR:CG2	1:A:8:ILE:HD11	0.61	2.26	8	20
1:A:45:LEU:HD12	1:A:75:VAL:HB	0.61	1.73	12	20
2:B:181:LEU:HD12	2:B:183:ILE:HG12	0.61	1.70	11	20
1:A:32:SER:O	1:A:36:PRO:CD	0.61	2.48	15	20
2:B:104:SER:OG	2:B:119:LEU:HD11	0.61	1.96	7	1
2:B:181:LEU:HD13	2:B:182:ASN:N	0.61	2.11	20	20
2:B:107:PHE:HE2	2:B:129:LEU:HD13	0.61	1.56	13	10
1:A:69:ALA:HB3	1:A:71:LEU:HG	0.60	1.73	16	4
1:A:8:ILE:HG21	1:A:66:ILE:CG2	0.60	2.26	12	20
2:B:160:LEU:C	2:B:160:LEU:HD13	0.60	2.21	3	15
1:A:65:ASN:O	1:A:69:ALA:HB3	0.60	1.95	8	20
2:B:176:ASN:OD1	2:B:178:GLU:HG2	0.60	1.96	9	8
1:A:16:ILE:HD13	2:B:156:ASP:OD2	0.60	1.96	6	1
2:B:118:LEU:HD11	2:B:132:ALA:HB3	0.60	1.74	14	20
2:B:173:LEU:HD12	2:B:173:LEU:O	0.60	1.97	7	3
1:A:10:PHE:HA	1:A:73:ILE:O	0.59	1.97	5	20
1:A:44:LYS:O	1:A:75:VAL:HA	0.59	1.97	15	20
2:B:145:ILE:HG22	2:B:146:THR:N	0.59	2.12	4	20
1:A:34:ILE:HD13	1:A:45:LEU:CG	0.59	2.27	20	20
1:A:69:ALA:CB	1:A:71:LEU:HD21	0.59	2.27	1	4
1:A:57:ILE:HG21	1:A:66:ILE:HD13	0.59	1.74	3	20
2:B:173:LEU:O	2:B:177:ASN:N	0.59	2.36	13	20
1:A:12:TYR:O	1:A:14:ASP:N	0.59	2.35	17	20
1:A:16:ILE:O	2:B:159:VAL:HG12	0.59	1.98	2	1
1:A:21:LEU:HD13	1:A:33:LYS:HE3	0.59	1.74	20	4
1:A:18:ALA:CB	2:B:158:VAL:HG21	0.58	2.27	4	7
1:A:63:VAL:O	1:A:67:ILE:HG23	0.58	1.98	11	20
2:B:171:GLU:O	2:B:174:ALA:HB3	0.58	1.97	17	13
2:B:145:ILE:O	2:B:146:THR:CB	0.58	2.50	10	1
1:A:17:PHE:CD1	2:B:159:VAL:HG13	0.58	2.34	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:149:LYS:HA	2:B:158:VAL:O	0.58	1.98	16	20
1:A:64:SER:O	1:A:67:ILE:HG12	0.58	1.99	16	20
2:B:118:LEU:HD22	2:B:119:LEU:N	0.58	2.14	2	19
2:B:150:TYR:N	2:B:158:VAL:O	0.58	2.36	16	16
2:B:173:LEU:HD11	2:B:179:LYS:H	0.58	1.59	7	4
2:B:107:PHE:CZ	2:B:129:LEU:HD22	0.57	2.35	13	1
1:A:59:THR:O	1:A:63:VAL:HG13	0.57	1.99	11	20
1:A:65:ASN:O	1:A:69:ALA:CB	0.57	2.52	8	20
2:B:108:ARG:CD	2:B:117:THR:HG22	0.57	2.29	19	8
1:A:9:LYS:CD	1:A:16:ILE:HG21	0.57	2.29	3	9
2:B:105:ILE:N	2:B:105:ILE:HD12	0.57	2.14	19	1
1:A:12:TYR:HA	1:A:75:VAL:O	0.57	1.99	1	20
1:A:10:PHE:CE2	1:A:34:ILE:HD11	0.57	2.34	16	20
1:A:16:ILE:HB	2:B:157:PHE:O	0.57	1.98	14	20
1:A:11:TYR:O	1:A:75:VAL:N	0.57	2.38	17	20
2:B:145:ILE:O	2:B:146:THR:HB	0.57	1.99	10	1
2:B:173:LEU:HD21	2:B:179:LYS:N	0.57	2.15	8	7
1:A:57:ILE:CG2	1:A:66:ILE:HD13	0.57	2.29	11	20
2:B:173:LEU:HD23	2:B:178:GLU:CG	0.56	2.30	6	6
1:A:20:MET:HB3	2:B:168:VAL:HG22	0.56	1.77	14	9
2:B:149:LYS:HD3	2:B:159:VAL:HG22	0.56	1.77	3	3
2:B:107:PHE:CE2	2:B:129:LEU:HD13	0.56	2.36	13	7
2:B:163:ASP:O	2:B:167:ASN:ND2	0.56	2.39	18	4
1:A:6:THR:HG21	1:A:8:ILE:HD11	0.56	1.76	8	20
1:A:18:ALA:CB	2:B:158:VAL:HG11	0.56	2.31	16	5
1:A:62:GLN:O	1:A:66:ILE:HD12	0.56	2.01	11	20
1:A:69:ALA:CB	1:A:71:LEU:HD13	0.56	2.31	10	16
2:B:108:ARG:O	2:B:183:ILE:N	0.56	2.39	14	20
1:A:16:ILE:HG22	2:B:158:VAL:HG22	0.56	1.76	17	3
1:A:12:TYR:CE1	1:A:38:ILE:HG21	0.55	2.36	15	20
2:B:104:SER:HB2	2:B:119:LEU:HD21	0.55	1.78	2	4
2:B:173:LEU:HD13	2:B:178:GLU:HG3	0.55	1.76	19	2
1:A:20:MET:HE3	2:B:167:ASN:HB2	0.55	1.78	16	1
2:B:164:GLU:O	2:B:167:ASN:ND2	0.55	2.39	3	1
2:B:136:LYS:O	2:B:139:ASN:OD1	0.55	2.25	19	1
1:A:21:LEU:HD13	1:A:33:LYS:HD2	0.55	1.76	15	1
1:A:10:PHE:CE1	1:A:19:LEU:CD1	0.55	2.89	20	20
2:B:107:PHE:HZ	2:B:129:LEU:HD22	0.55	1.60	13	1
2:B:128:ASP:O	2:B:132:ALA:N	0.54	2.39	11	12
2:B:150:TYR:HB2	2:B:181:LEU:HD11	0.54	1.79	6	10
1:A:30:LEU:O	1:A:33:LYS:HB2	0.54	2.02	19	10

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:173:LEU:HD23	2:B:178:GLU:HG2	0.54	1.77	20	6
1:A:17:PHE:CG	1:A:37:ARG:NH1	0.54	2.75	15	1
2:B:127:ASP:O	2:B:128:ASP:C	0.54	2.50	18	19
1:A:38:ILE:CD1	1:A:43:PHE:CG	0.54	2.90	11	20
2:B:180:PHE:CG	2:B:181:LEU:N	0.54	2.76	19	12
1:A:34:ILE:HD13	1:A:45:LEU:CD1	0.54	2.32	11	20
2:B:160:LEU:HD22	2:B:161:GLY:N	0.54	2.17	2	5
2:B:105:ILE:HD11	2:B:122:LYS:HE2	0.53	1.80	13	1
1:A:19:LEU:HD22	1:A:33:LYS:HD3	0.53	1.80	16	3
1:A:21:LEU:HD22	1:A:21:LEU:N	0.53	2.18	19	1
2:B:114:GLU:O	2:B:115:ILE:CD1	0.53	2.54	8	4
2:B:173:LEU:HD12	2:B:173:LEU:C	0.53	2.29	7	2
1:A:9:LYS:HD3	1:A:16:ILE:CG2	0.53	2.34	2	17
1:A:19:LEU:HB3	1:A:33:LYS:HD3	0.53	1.79	19	3
1:A:42:ASN:O	1:A:42:ASN:ND2	0.53	2.42	2	19
2:B:107:PHE:CE1	2:B:120:VAL:CG1	0.53	2.92	15	20
2:B:173:LEU:HD23	2:B:173:LEU:O	0.53	2.04	10	2
1:A:34:ILE:O	1:A:38:ILE:CG1	0.53	2.56	15	20
1:A:45:LEU:HB2	1:A:57:ILE:HG12	0.52	1.80	11	20
1:A:19:LEU:CD1	1:A:33:LYS:HB3	0.52	2.28	17	7
1:A:27:TYR:HB2	1:A:63:VAL:HG11	0.52	1.81	11	20
1:A:63:VAL:CG2	1:A:64:SER:N	0.52	2.72	20	20
2:B:119:LEU:HD13	2:B:120:VAL:N	0.52	2.19	2	4
2:B:145:ILE:O	2:B:146:THR:HG22	0.52	2.04	10	1
2:B:111:TYR:CZ	2:B:145:ILE:HD11	0.52	2.39	19	1
1:A:66:ILE:CG1	1:A:71:LEU:HB2	0.52	2.35	16	20
1:A:9:LYS:CD	1:A:16:ILE:CG2	0.52	2.88	18	9
1:A:34:ILE:HG22	1:A:38:ILE:CD1	0.52	2.27	15	20
2:B:159:VAL:CG2	2:B:160:LEU:N	0.52	2.73	2	1
1:A:7:LYS:HE3	1:A:18:ALA:HB1	0.52	1.81	15	2
1:A:64:SER:O	1:A:67:ILE:N	0.52	2.43	2	20
1:A:16:ILE:O	2:B:158:VAL:HA	0.52	2.05	13	14
2:B:121:GLU:HB2	2:B:124:TRP:CD1	0.52	2.40	12	13
1:A:16:ILE:O	2:B:158:VAL:CA	0.51	2.58	13	12
2:B:129:LEU:O	2:B:132:ALA:HB3	0.51	2.05	1	20
2:B:137:ILE:HG21	2:B:185:LEU:HD11	0.51	1.83	8	10
2:B:158:VAL:CG2	2:B:159:VAL:N	0.51	2.74	7	12
1:A:9:LYS:HG3	1:A:10:PHE:N	0.51	2.20	9	16
1:A:7:LYS:NZ	1:A:18:ALA:HB1	0.51	2.20	3	2
1:A:9:LYS:C	1:A:10:PHE:CD1	0.51	2.88	4	20
2:B:137:ILE:HG21	2:B:185:LEU:CD1	0.51	2.35	1	10

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:105:ILE:O	2:B:119:LEU:HD12	0.51	2.05	3	3
1:A:10:PHE:CE2	1:A:30:LEU:HD21	0.51	2.40	3	20
1:A:47:THR:HG22	1:A:66:ILE:CG1	0.51	2.35	11	20
1:A:47:THR:HG22	1:A:66:ILE:HG12	0.51	1.82	11	17
2:B:158:VAL:HG12	2:B:159:VAL:N	0.51	2.20	5	8
2:B:149:LYS:CE	2:B:149:LYS:CA	0.51	2.88	8	1
2:B:111:TYR:CD2	2:B:185:LEU:HB3	0.51	2.40	20	6
2:B:150:TYR:CE2	2:B:158:VAL:HG22	0.51	2.41	13	1
1:A:12:TYR:CD2	1:A:13:LYS:HG2	0.51	2.40	12	8
2:B:137:ILE:HD13	2:B:145:ILE:HG13	0.51	1.83	4	9
2:B:159:VAL:HG23	2:B:160:LEU:N	0.51	2.20	7	6
2:B:173:LEU:C	2:B:173:LEU:HD22	0.51	2.31	13	1
2:B:173:LEU:C	2:B:173:LEU:HD12	0.51	2.31	3	2
1:A:16:ILE:HG22	2:B:158:VAL:HB	0.51	1.83	8	10
1:A:5:THR:HG21	1:A:20:MET:SD	0.51	2.46	10	1
1:A:17:PHE:CD2	1:A:37:ARG:HB3	0.50	2.40	9	5
2:B:159:VAL:HG12	2:B:160:LEU:N	0.50	2.21	14	8
1:A:47:THR:O	1:A:54:GLY:HA3	0.50	2.07	11	2
1:A:43:PHE:N	1:A:78:ILE:HD11	0.50	2.21	20	19
1:A:21:LEU:N	1:A:21:LEU:CD2	0.50	2.75	15	5
2:B:137:ILE:HD13	2:B:145:ILE:CG1	0.50	2.36	18	7
1:A:38:ILE:HD12	1:A:43:PHE:CG	0.50	2.42	11	20
1:A:66:ILE:HG23	1:A:71:LEU:O	0.50	2.07	5	20
2:B:111:TYR:CD1	2:B:111:TYR:C	0.50	2.90	6	14
1:A:15:ASP:OD1	1:A:17:PHE:CE2	0.50	2.64	7	1
1:A:25:THR:HG23	1:A:29:GLU:CD	0.50	2.31	15	1
1:A:35:ALA:N	1:A:36:PRO:HD2	0.50	2.22	15	20
1:A:6:THR:CG2	1:A:8:ILE:CD1	0.50	2.90	8	20
1:A:12:TYR:CE2	1:A:38:ILE:HG22	0.50	2.42	10	19
1:A:44:LYS:HD2	1:A:78:ILE:HG23	0.50	1.83	20	13
1:A:16:ILE:HB	2:B:158:VAL:HA	0.50	1.84	19	12
1:A:35:ALA:CB	1:A:43:PHE:CE1	0.50	2.92	15	20
2:B:107:PHE:CE1	2:B:120:VAL:HG11	0.50	2.42	15	17
1:A:47:THR:CG2	1:A:66:ILE:CD1	0.50	2.89	11	1
2:B:148:ILE:HG22	2:B:185:LEU:HD21	0.50	1.82	11	1
2:B:125:ASN:HA	2:B:166:TRP:CE2	0.50	2.41	2	12
1:A:12:TYR:CD2	1:A:13:LYS:CG	0.50	2.95	8	1
2:B:149:LYS:HE3	2:B:159:VAL:HA	0.50	1.83	16	1
1:A:27:TYR:CB	1:A:63:VAL:HG11	0.49	2.37	11	20
2:B:173:LEU:O	2:B:176:ASN:N	0.49	2.40	7	1
1:A:8:ILE:CD1	1:A:70:LYS:HA	0.49	2.37	8	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:46:GLN:O	1:A:74:SER:N	0.49	2.43	1	20
1:A:47:THR:HG22	1:A:66:ILE:HD11	0.49	1.84	19	19
2:B:164:GLU:HA	2:B:167:ASN:OD1	0.49	2.07	3	1
1:A:34:ILE:O	1:A:37:ARG:HG3	0.49	2.06	15	1
1:A:35:ALA:HB3	1:A:36:PRO:CD	0.49	2.37	3	20
1:A:66:ILE:HA	1:A:71:LEU:HG	0.49	1.84	1	4
2:B:129:LEU:C	2:B:129:LEU:CD1	0.49	2.85	14	20
2:B:150:TYR:CD1	2:B:181:LEU:CD1	0.49	2.96	3	6
2:B:107:PHE:CD1	2:B:120:VAL:HG13	0.49	2.42	13	20
2:B:106:LEU:HD11	2:B:117:THR:HG22	0.49	1.85	9	7
2:B:145:ILE:O	2:B:146:THR:CG2	0.49	2.61	10	1
1:A:17:PHE:CD2	2:B:159:VAL:CG2	0.49	2.96	15	5
2:B:107:PHE:CD1	2:B:120:VAL:CG1	0.49	2.96	13	20
2:B:149:LYS:HG2	2:B:157:PHE:HB3	0.49	1.83	16	1
2:B:150:TYR:CE1	2:B:151:GLN:O	0.49	2.66	18	7
1:A:5:THR:CG2	1:A:22:LYS:HD2	0.49	2.37	18	2
1:A:26:THR:HG21	1:A:28:LYS:CE	0.49	2.38	11	2
2:B:145:ILE:CG2	2:B:185:LEU:HD11	0.49	2.37	20	1
2:B:153:GLU:CG	2:B:172:MET:HE1	0.49	2.33	16	2
1:A:66:ILE:HG21	1:A:73:ILE:CD1	0.49	2.34	1	20
2:B:107:PHE:CD2	2:B:133:ILE:HD11	0.48	2.43	14	5
2:B:111:TYR:CE2	2:B:186:TYR:CD1	0.48	3.01	5	1
1:A:6:THR:CG2	1:A:67:ILE:O	0.48	2.61	8	20
1:A:9:LYS:C	1:A:10:PHE:CG	0.48	2.90	6	20
1:A:27:TYR:CD2	1:A:28:LYS:N	0.48	2.82	14	19
1:A:19:LEU:HB3	1:A:33:LYS:CD	0.48	2.38	19	3
2:B:149:LYS:HE3	2:B:159:VAL:HG12	0.48	1.85	7	3
1:A:49:LEU:O	1:A:52:GLY:N	0.48	2.46	11	1
1:A:57:ILE:CG2	1:A:66:ILE:CD1	0.48	2.92	11	20
2:B:173:LEU:C	2:B:173:LEU:HD23	0.48	2.32	10	2
1:A:43:PHE:HA	1:A:76:HIS:O	0.48	2.09	19	1
1:A:73:ILE:HG22	1:A:74:SER:N	0.48	2.23	10	20
2:B:124:TRP:HB3	2:B:128:ASP:HB3	0.48	1.85	11	2
1:A:20:MET:C	1:A:21:LEU:HD22	0.48	2.34	15	5
2:B:157:PHE:CD1	2:B:184:ARG:CD	0.48	2.97	17	5
2:B:149:LYS:HG3	2:B:158:VAL:O	0.48	2.08	8	2
2:B:146:THR:HG23	2:B:147:LYS:NZ	0.48	2.24	4	4
2:B:181:LEU:CD1	2:B:182:ASN:N	0.48	2.76	20	20
1:A:27:TYR:CD2	1:A:28:LYS:HG3	0.48	2.43	18	11
2:B:118:LEU:HG	2:B:132:ALA:C	0.48	2.34	5	19
1:A:7:LYS:NZ	2:B:168:VAL:HG11	0.48	2.23	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:69:ALA:HB3	1:A:71:LEU:CG	0.48	2.38	16	4
2:B:124:TRP:CE3	2:B:128:ASP:CG	0.48	2.92	1	2
2:B:108:ARG:NH1	2:B:115:ILE:HG21	0.48	2.24	13	5
2:B:173:LEU:HD21	2:B:178:GLU:N	0.48	2.23	4	8
1:A:7:LYS:HE3	2:B:168:VAL:HG11	0.48	1.85	15	1
1:A:34:ILE:HG12	1:A:37:ARG:NH2	0.48	2.23	15	1
1:A:30:LEU:HD22	1:A:30:LEU:O	0.48	2.09	13	14
1:A:15:ASP:OD2	2:B:159:VAL:HG11	0.48	2.09	2	1
1:A:9:LYS:HD2	1:A:11:TYR:CE1	0.48	2.44	14	12
2:B:163:ASP:O	2:B:167:ASN:OD1	0.48	2.32	3	1
2:B:120:VAL:O	2:B:120:VAL:CG2	0.47	2.62	13	11
2:B:114:GLU:C	2:B:115:ILE:HD13	0.47	2.33	2	5
1:A:21:LEU:CD1	1:A:33:LYS:HD2	0.47	2.39	6	11
2:B:157:PHE:CD2	2:B:184:ARG:CD	0.47	2.96	6	8
2:B:118:LEU:HD13	2:B:118:LEU:O	0.47	2.10	13	10
2:B:145:ILE:HG22	2:B:147:LYS:H	0.47	1.69	18	1
1:A:30:LEU:O	1:A:30:LEU:HD22	0.47	2.09	4	6
2:B:145:ILE:CG2	2:B:146:THR:N	0.47	2.78	11	6
1:A:26:THR:HG21	1:A:28:LYS:NZ	0.47	2.24	5	1
2:B:137:ILE:HG21	2:B:185:LEU:CD2	0.47	2.39	17	4
2:B:129:LEU:O	2:B:133:ILE:N	0.47	2.44	13	11
1:A:66:ILE:HA	1:A:71:LEU:HB2	0.47	1.86	3	16
2:B:160:LEU:HD22	2:B:165:ASP:HB3	0.47	1.87	3	1
2:B:166:TRP:CZ2	2:B:170:LYS:CE	0.47	2.97	10	3
1:A:21:LEU:CD1	1:A:33:LYS:CE	0.47	2.93	7	6
2:B:181:LEU:CD1	2:B:181:LEU:C	0.47	2.88	14	20
2:B:120:VAL:HG22	2:B:120:VAL:O	0.47	2.10	1	13
1:A:69:ALA:CB	1:A:71:LEU:CD2	0.47	2.92	1	4
2:B:119:LEU:HD13	2:B:119:LEU:C	0.47	2.35	4	3
1:A:69:ALA:HB1	1:A:71:LEU:CD1	0.47	2.40	11	16
1:A:5:THR:HG22	1:A:22:LYS:CD	0.47	2.39	6	1
2:B:125:ASN:O	2:B:126:PHE:C	0.46	2.58	9	16
2:B:160:LEU:HD13	2:B:160:LEU:O	0.46	2.09	3	2
2:B:130:ILE:HG22	2:B:131:MET:N	0.46	2.25	13	1
1:A:26:THR:C	1:A:63:VAL:HG21	0.46	2.36	16	20
2:B:120:VAL:O	2:B:120:VAL:HG22	0.46	2.10	13	6
2:B:150:TYR:CD1	2:B:181:LEU:HD11	0.46	2.46	3	3
1:A:27:TYR:CG	1:A:28:LYS:N	0.46	2.83	6	13
2:B:106:LEU:HD11	2:B:117:THR:HB	0.46	1.86	3	4
1:A:47:THR:HG22	1:A:66:ILE:CD1	0.46	2.41	4	19
2:B:160:LEU:HD23	2:B:165:ASP:OD1	0.46	2.10	1	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:50:PHE:CD1	1:A:51:ASP:N	0.46	2.84	5	5
1:A:9:LYS:CD	2:B:156:ASP:OD2	0.46	2.64	6	1
1:A:62:GLN:C	1:A:66:ILE:HD12	0.46	2.36	11	17
1:A:73:ILE:CG2	1:A:74:SER:N	0.46	2.79	2	20
2:B:105:ILE:N	2:B:105:ILE:CD1	0.46	2.79	1	10
2:B:111:TYR:O	2:B:112:ASN:HB2	0.46	2.11	14	3
2:B:149:LYS:HB3	2:B:184:ARG:HG2	0.46	1.87	16	1
1:A:6:THR:HG22	1:A:8:ILE:HG12	0.46	1.88	5	19
1:A:12:TYR:CD2	1:A:38:ILE:CG2	0.46	2.99	10	9
1:A:26:THR:CG2	1:A:28:LYS:CE	0.46	2.94	11	2
2:B:110:SER:OG	2:B:115:ILE:HD12	0.46	2.11	16	1
1:A:48:LYS:N	1:A:72:LYS:O	0.45	2.49	19	17
2:B:173:LEU:HA	2:B:176:ASN:HD21	0.45	1.70	7	1
2:B:136:LYS:O	2:B:139:ASN:ND2	0.45	2.48	10	4
1:A:12:TYR:CZ	1:A:13:LYS:CE	0.45	2.99	9	5
1:A:34:ILE:CG1	1:A:37:ARG:NH2	0.45	2.80	15	1
2:B:118:LEU:HD22	2:B:118:LEU:C	0.45	2.36	2	17
1:A:21:LEU:CD1	1:A:33:LYS:CD	0.45	2.94	3	1
1:A:34:ILE:O	1:A:38:ILE:HG12	0.45	2.11	15	8
2:B:183:ILE:HG22	2:B:184:ARG:N	0.45	2.26	13	14
2:B:158:VAL:CG1	2:B:159:VAL:N	0.45	2.79	16	8
2:B:173:LEU:O	2:B:176:ASN:ND2	0.45	2.50	7	1
1:A:27:TYR:CZ	1:A:31:ARG:HD3	0.45	2.47	15	3
2:B:123:VAL:HG22	2:B:124:TRP:CD1	0.45	2.47	11	2
1:A:17:PHE:CD1	2:B:159:VAL:HB	0.45	2.47	8	3
2:B:178:GLU:O	2:B:179:LYS:CB	0.45	2.64	10	2
2:B:150:TYR:CD1	2:B:158:VAL:HG22	0.45	2.47	7	1
2:B:126:PHE:CE1	2:B:163:ASP:N	0.45	2.84	13	1
2:B:106:LEU:HD11	2:B:117:THR:CB	0.45	2.42	3	3
2:B:126:PHE:CE1	2:B:130:ILE:CG1	0.45	2.99	15	1
1:A:27:TYR:CE1	1:A:31:ARG:HD2	0.45	2.47	12	5
1:A:17:PHE:CE1	2:B:159:VAL:CG2	0.45	3.00	2	1
2:B:180:PHE:CD2	2:B:181:LEU:N	0.45	2.85	2	4
1:A:57:ILE:HG22	1:A:63:VAL:N	0.45	2.26	11	2
2:B:150:TYR:CE2	2:B:158:VAL:CG2	0.45	3.00	13	1
1:A:9:LYS:HE2	2:B:158:VAL:HG11	0.45	1.88	19	1
2:B:149:LYS:CD	2:B:159:VAL:HG22	0.45	2.42	18	4
1:A:10:PHE:CD1	1:A:10:PHE:N	0.45	2.84	4	15
1:A:27:TYR:CG	1:A:58:LYS:O	0.45	2.70	2	15
1:A:17:PHE:CD1	1:A:37:ARG:HB3	0.45	2.46	17	3
2:B:104:SER:OG	2:B:119:LEU:HD21	0.45	2.11	7	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:160:LEU:CD2	2:B:165:ASP:HB2	0.45	2.42	12	2
2:B:173:LEU:HD23	2:B:178:GLU:HG3	0.45	1.89	16	2
1:A:12:TYR:O	1:A:15:ASP:N	0.45	2.50	19	3
2:B:176:ASN:N	2:B:176:ASN:ND2	0.45	2.65	20	5
1:A:10:PHE:O	1:A:16:ILE:HA	0.44	2.13	4	11
1:A:31:ARG:HG3	1:A:43:PHE:CE2	0.44	2.46	2	4
2:B:130:ILE:HA	2:B:133:ILE:HD12	0.44	1.88	1	2
2:B:118:LEU:CD1	2:B:120:VAL:HG12	0.44	2.41	17	16
1:A:31:ARG:HG2	1:A:43:PHE:CZ	0.44	2.48	19	6
2:B:180:PHE:CD1	2:B:181:LEU:N	0.44	2.85	4	2
1:A:17:PHE:CD2	2:B:159:VAL:HG22	0.44	2.47	6	3
1:A:35:ALA:HB3	1:A:36:PRO:HD3	0.44	1.89	17	19
2:B:105:ILE:C	2:B:119:LEU:HD12	0.44	2.37	3	2
1:A:17:PHE:HB2	1:A:37:ARG:CZ	0.44	2.42	15	1
1:A:7:LYS:CE	1:A:18:ALA:HB1	0.44	2.41	3	4
1:A:38:ILE:HD12	1:A:43:PHE:CB	0.44	2.42	19	19
2:B:181:LEU:CD1	2:B:183:ILE:CD1	0.44	2.96	13	16
1:A:11:TYR:HD1	1:A:73:ILE:O	0.44	1.95	5	2
1:A:45:LEU:CB	1:A:57:ILE:HG12	0.44	2.42	11	20
1:A:13:LYS:N	1:A:13:LYS:CD	0.44	2.80	9	4
1:A:12:TYR:HB2	1:A:75:VAL:HG13	0.44	1.89	17	3
2:B:186:TYR:CD1	2:B:186:TYR:C	0.44	2.95	14	1
1:A:45:LEU:CD1	1:A:75:VAL:CB	0.44	2.88	9	18
1:A:45:LEU:HB2	1:A:57:ILE:CG1	0.44	2.43	11	20
2:B:173:LEU:O	2:B:173:LEU:HD12	0.44	2.13	1	1
2:B:126:PHE:O	2:B:130:ILE:HB	0.44	2.12	2	11
2:B:137:ILE:O	2:B:139:ASN:N	0.44	2.50	3	5
1:A:6:THR:HG22	1:A:8:ILE:CD1	0.44	2.43	8	3
1:A:10:PHE:N	1:A:10:PHE:CD1	0.44	2.85	9	5
1:A:17:PHE:CD2	2:B:159:VAL:HB	0.44	2.47	18	2
1:A:57:ILE:HA	1:A:62:GLN:NE2	0.44	2.28	13	19
2:B:127:ASP:O	2:B:129:LEU:N	0.44	2.51	9	8
1:A:21:LEU:HD11	1:A:33:LYS:HD2	0.44	1.89	18	7
2:B:176:ASN:HB2	2:B:178:GLU:HG2	0.44	1.90	7	1
1:A:26:THR:HG22	1:A:27:TYR:N	0.44	2.28	16	4
2:B:157:PHE:CD1	2:B:184:ARG:HD2	0.44	2.48	11	1
1:A:12:TYR:CE2	1:A:38:ILE:CG2	0.44	3.01	10	11
1:A:12:TYR:CD2	1:A:13:LYS:HG3	0.44	2.48	8	7
2:B:110:SER:HB2	2:B:115:ILE:HD12	0.44	1.88	2	2
2:B:137:ILE:C	2:B:139:ASN:N	0.44	2.76	17	11
2:B:106:LEU:HB3	2:B:180:PHE:CD2	0.44	2.48	16	5

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:149:LYS:CA	2:B:158:VAL:O	0.43	2.66	2	9
1:A:12:TYR:N	1:A:15:ASP:O	0.43	2.46	4	5
2:B:145:ILE:HG21	2:B:185:LEU:HD22	0.43	1.89	16	1
1:A:13:LYS:CD	1:A:13:LYS:N	0.43	2.81	17	1
2:B:179:LYS:C	2:B:180:PHE:CD1	0.43	2.96	5	2
2:B:173:LEU:CD2	2:B:178:GLU:HG2	0.43	2.43	9	1
1:A:10:PHE:HB2	1:A:37:ARG:NH2	0.43	2.29	15	1
2:B:109:ILE:CG2	2:B:185:LEU:HD22	0.43	2.43	17	1
1:A:34:ILE:CG2	1:A:38:ILE:CD1	0.43	2.96	4	5
1:A:27:TYR:O	1:A:30:LEU:N	0.43	2.51	11	6
2:B:146:THR:HG22	2:B:147:LYS:HE2	0.43	1.91	7	1
2:B:109:ILE:N	2:B:116:PHE:O	0.43	2.42	14	4
2:B:150:TYR:CD1	2:B:150:TYR:C	0.43	2.97	16	7
1:A:27:TYR:O	1:A:28:LYS:C	0.43	2.62	16	13
1:A:5:THR:HG22	1:A:22:LYS:CG	0.43	2.44	20	1
1:A:65:ASN:O	1:A:69:ALA:N	0.43	2.46	11	20
2:B:157:PHE:CD2	2:B:184:ARG:HD2	0.43	2.49	19	6
1:A:17:PHE:CD1	1:A:17:PHE:N	0.43	2.86	15	1
1:A:6:THR:HG21	1:A:67:ILE:HA	0.43	1.91	20	15
2:B:176:ASN:N	2:B:176:ASN:HD22	0.43	2.12	20	2
2:B:107:PHE:CE2	2:B:133:ILE:HD11	0.43	2.49	14	1
1:A:17:PHE:CE2	2:B:159:VAL:HG21	0.43	2.49	15	1
1:A:64:SER:HA	1:A:67:ILE:HD13	0.42	1.90	12	14
2:B:176:ASN:ND2	2:B:176:ASN:N	0.42	2.67	17	2
2:B:105:ILE:HD11	2:B:122:LYS:CD	0.42	2.44	18	2
2:B:111:TYR:CG	2:B:185:LEU:HB3	0.42	2.49	19	1
1:A:16:ILE:O	2:B:159:VAL:CG1	0.42	2.66	2	1
2:B:106:LEU:C	2:B:106:LEU:CD1	0.42	2.91	4	10
2:B:160:LEU:CD2	2:B:165:ASP:CB	0.42	2.97	3	2
2:B:150:TYR:O	2:B:150:TYR:CG	0.42	2.71	20	2
2:B:147:LYS:C	2:B:147:LYS:CE	0.42	2.93	17	1
1:A:31:ARG:O	1:A:43:PHE:CZ	0.42	2.73	1	3
2:B:137:ILE:CD1	2:B:145:ILE:CG1	0.42	2.98	4	5
1:A:27:TYR:CE2	1:A:28:LYS:CD	0.42	3.02	8	6
1:A:10:PHE:O	1:A:16:ILE:HG23	0.42	2.14	19	2
2:B:137:ILE:CD1	2:B:145:ILE:HB	0.42	2.44	20	1
1:A:9:LYS:NZ	2:B:154:ASP:CB	0.42	2.82	9	1
2:B:124:TRP:O	2:B:166:TRP:NE1	0.42	2.52	16	3
1:A:27:TYR:HB2	1:A:63:VAL:CG1	0.42	2.43	2	5
1:A:57:ILE:HD11	1:A:73:ILE:HG23	0.42	1.91	15	4
2:B:149:LYS:O	2:B:183:ILE:HA	0.42	2.14	7	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:160:LEU:C	2:B:160:LEU:CD1	0.42	2.91	3	1
2:B:164:GLU:CG	2:B:165:ASP:N	0.42	2.83	7	2
2:B:152:ASP:OD1	2:B:172:MET:HE3	0.42	2.15	11	1
2:B:151:GLN:HB2	2:B:157:PHE:CE2	0.42	2.50	3	1
2:B:118:LEU:HD13	2:B:118:LEU:C	0.42	2.40	14	5
2:B:145:ILE:HG13	2:B:185:LEU:HD21	0.42	1.91	2	1
1:A:12:TYR:CD1	1:A:13:LYS:HD3	0.42	2.50	9	2
2:B:180:PHE:CD1	2:B:180:PHE:C	0.42	2.98	3	1
2:B:145:ILE:CG2	2:B:146:THR:H	0.42	2.24	4	1
1:A:49:LEU:HD12	1:A:52:GLY:H	0.42	1.75	5	4
1:A:15:ASP:OD2	1:A:17:PHE:CZ	0.42	2.73	7	1
1:A:27:TYR:CE2	1:A:28:LYS:HD2	0.42	2.50	17	6
1:A:9:LYS:CG	1:A:10:PHE:N	0.42	2.82	19	1
1:A:12:TYR:CZ	1:A:38:ILE:CG2	0.42	3.03	15	2
1:A:17:PHE:HB3	1:A:37:ARG:NH1	0.42	2.29	7	1
2:B:124:TRP:O	2:B:166:TRP:CZ2	0.41	2.73	1	1
2:B:181:LEU:HD12	2:B:183:ILE:CG1	0.41	2.44	13	5
2:B:129:LEU:O	2:B:133:ILE:HG13	0.41	2.14	11	1
2:B:149:LYS:HE3	2:B:159:VAL:CA	0.41	2.45	16	1
1:A:34:ILE:HG21	1:A:75:VAL:HG21	0.41	1.92	15	5
2:B:166:TRP:CZ2	2:B:170:LYS:HE2	0.41	2.51	8	2
1:A:12:TYR:CZ	1:A:13:LYS:HD3	0.41	2.50	19	1
2:B:124:TRP:HB3	2:B:128:ASP:CB	0.41	2.45	11	2
1:A:45:LEU:HD13	1:A:75:VAL:HB	0.41	1.82	7	3
1:A:12:TYR:CZ	1:A:13:LYS:HE2	0.41	2.50	17	2
2:B:130:ILE:HD13	2:B:133:ILE:HD12	0.41	1.92	18	1
1:A:38:ILE:CD1	1:A:75:VAL:HG21	0.41	2.45	16	16
2:B:111:TYR:HD2	2:B:185:LEU:HB3	0.41	1.75	6	2
1:A:12:TYR:CE1	1:A:38:ILE:CG2	0.41	3.02	15	1
2:B:145:ILE:HD12	2:B:185:LEU:CD1	0.41	2.46	20	1
2:B:173:LEU:CD1	2:B:178:GLU:N	0.41	2.83	3	2
2:B:110:SER:CB	2:B:115:ILE:CD1	0.41	2.99	10	5
2:B:116:PHE:CZ	2:B:139:ASN:OD1	0.41	2.74	3	2
1:A:15:ASP:OD1	1:A:15:ASP:C	0.41	2.64	7	1
1:A:50:PHE:CE2	1:A:51:ASP:OD1	0.41	2.74	7	2
2:B:166:TRP:O	2:B:170:LYS:HD3	0.41	2.15	10	2
1:A:12:TYR:CE1	1:A:13:LYS:HD3	0.41	2.50	19	1
2:B:111:TYR:CZ	2:B:112:ASN:OD1	0.41	2.74	12	1
1:A:16:ILE:O	2:B:158:VAL:HB	0.41	2.16	19	1
1:A:69:ALA:HB1	1:A:71:LEU:CD2	0.41	2.34	1	2
2:B:107:PHE:HB2	2:B:118:LEU:O	0.41	2.15	7	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:105:ILE:CD1	2:B:122:LYS:HG2	0.41	2.45	8	1
1:A:45:LEU:CB	1:A:57:ILE:CG1	0.41	2.99	11	1
2:B:111:TYR:CE1	2:B:112:ASN:OD1	0.41	2.74	14	1
1:A:17:PHE:CD2	1:A:37:ARG:NH1	0.41	2.89	15	1
1:A:34:ILE:HA	1:A:37:ARG:HG2	0.41	1.93	11	2
1:A:25:THR:HG23	1:A:29:GLU:HB3	0.41	1.93	9	1
1:A:17:PHE:CB	1:A:37:ARG:CZ	0.41	2.99	15	1
2:B:121:GLU:OE2	2:B:124:TRP:CZ2	0.41	2.74	15	1
1:A:17:PHE:C	2:B:158:VAL:HG13	0.41	2.41	16	1
2:B:152:ASP:OD1	2:B:172:MET:HE2	0.41	2.15	16	1
1:A:38:ILE:CD1	1:A:43:PHE:CB	0.41	2.99	19	1
1:A:69:ALA:CB	1:A:71:LEU:CD1	0.41	2.98	11	8
2:B:118:LEU:HG	2:B:133:ILE:HA	0.41	1.93	4	4
2:B:151:GLN:O	2:B:182:ASN:ND2	0.41	2.54	10	1
2:B:104:SER:HA	2:B:121:GLU:HA	0.41	1.93	15	1
2:B:150:TYR:CE2	2:B:152:ASP:OD2	0.41	2.74	17	1
1:A:7:LYS:NZ	2:B:168:VAL:HG12	0.40	2.32	2	1
2:B:126:PHE:CE2	2:B:127:ASP:OD1	0.40	2.75	2	1
2:B:149:LYS:HB2	2:B:184:ARG:HG2	0.40	1.92	2	1
1:A:21:LEU:HD11	1:A:33:LYS:CD	0.40	2.46	3	1
2:B:111:TYR:CE2	2:B:112:ASN:OD1	0.40	2.74	8	1
2:B:173:LEU:CD2	2:B:178:GLU:CG	0.40	2.99	9	2
2:B:111:TYR:CE1	2:B:112:ASN:CG	0.40	2.99	17	1
1:A:49:LEU:O	1:A:50:PHE:C	0.40	2.64	5	2
2:B:150:TYR:CD1	2:B:182:ASN:OD1	0.40	2.74	12	1
1:A:11:TYR:CD2	1:A:16:ILE:HG12	0.40	2.51	13	1
2:B:176:ASN:OD1	2:B:178:GLU:CG	0.40	2.67	20	1
1:A:17:PHE:CD1	2:B:159:VAL:CG1	0.40	3.03	2	1
2:B:166:TRP:CE3	2:B:167:ASN:OD1	0.40	2.74	6	1
2:B:173:LEU:O	2:B:174:ALA:C	0.40	2.65	7	1
2:B:111:TYR:CE2	2:B:112:ASN:CG	0.40	2.99	12	2
2:B:116:PHE:CE1	2:B:139:ASN:OD1	0.40	2.73	18	2
2:B:121:GLU:CD	2:B:124:TRP:CZ2	0.40	2.99	12	1
2:B:146:THR:HG23	2:B:147:LYS:CD	0.40	2.46	19	1
2:B:150:TYR:HA	2:B:182:ASN:O	0.40	2.16	20	1
1:A:7:LYS:NZ	1:A:19:LEU:O	0.40	2.55	1	1
1:A:48:LYS:O	1:A:72:LYS:HG3	0.40	2.16	2	1
1:A:33:LYS:O	1:A:37:ARG:HG2	0.40	2.17	3	1
1:A:31:ARG:CG	1:A:43:PHE:CE2	0.40	3.05	2	1
2:B:150:TYR:CE1	2:B:182:ASN:OD1	0.40	2.74	12	1
2:B:109:ILE:HG13	2:B:183:ILE:HB	0.40	1.92	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:126:PHE:CE1	2:B:130:ILE:HG12	0.40	2.51	15	1
2:B:186:TYR:OXT	2:B:186:TYR:CG	0.40	2.75	17	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	73/78 (94%)	60±1 (82±2%)	10±1 (13±2%)	4±0 (5±1%)	3	25
2	B	77/86 (90%)	64±2 (84±2%)	8±2 (11±3%)	4±1 (5±2%)	2	22
All	All	3000/3280 (91%)	2484 (83%)	363 (12%)	153 (5%)	3	23

All 13 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	13	LYS	20
1	A	38	ILE	20
1	A	50	PHE	20
2	B	127	ASP	19
2	B	128	ASP	19
2	B	145	ILE	15
2	B	126	PHE	13
1	A	5	THR	10
2	B	138	SER	7
2	B	112	ASN	4
2	B	161	GLY	3
2	B	179	LYS	2
2	B	146	THR	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	68/72 (94%)	42±2 (62±3%)	26±2 (38±3%)	0	7
2	B	73/80 (91%)	46±2 (63±3%)	27±2 (37±3%)	1	8
All	All	2820/3040 (93%)	1773 (63%)	1047 (37%)	1	8

All 89 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	7	LYS	20
1	A	9	LYS	20
1	A	10	PHE	20
1	A	27	TYR	20
1	A	30	LEU	20
1	A	34	ILE	20
1	A	38	ILE	20
1	A	44	LYS	20
1	A	49	LEU	20
1	A	57	ILE	20
1	A	59	THR	20
1	A	63	VAL	20
1	A	66	ILE	20
1	A	67	ILE	20
1	A	74	SER	20
1	A	75	VAL	20
1	A	78	ILE	20
2	B	118	LEU	20
2	B	120	VAL	20
2	B	123	VAL	20
2	B	125	ASN	20
2	B	130	ILE	20
2	B	137	ILE	20
2	B	147	LYS	20
2	B	148	ILE	20
2	B	173	LEU	20
2	B	176	ASN	20
2	B	181	LEU	20
2	B	129	LEU	19
2	B	131	MET	18
2	B	170	LYS	18
1	A	39	ASP	17

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Mol	Chain	Res	Type	Models (Total)
2	B	153	GLU	17
2	B	127	ASP	17
2	B	136	LYS	16
2	B	156	ASP	16
1	A	71	LEU	16
1	A	72	LYS	15
2	B	162	SER	14
2	B	182	ASN	13
1	A	62	GLN	13
1	A	61	SER	12
2	B	158	VAL	12
2	B	159	VAL	12
2	B	180	PHE	12
2	B	108	ARG	11
2	B	122	LYS	11
2	B	104	SER	10
1	A	13	LYS	10
1	A	70	LYS	10
2	B	150	TYR	9
2	B	185	LEU	9
1	A	15	ASP	8
1	A	22	LYS	8
1	A	48	LYS	8
1	A	65	ASN	8
2	B	139	ASN	7
2	B	146	THR	7
2	B	128	ASP	7
2	B	167	ASN	7
1	A	31	ARG	7
1	A	60	ASP	7
2	B	121	GLU	7
1	A	33	LYS	6
1	A	37	ARG	5
2	B	152	ASP	5
1	A	77	ASP	5
2	B	179	LYS	5
2	B	178	GLU	4
1	A	24	ASP	4
1	A	20	MET	4
2	B	149	LYS	4
2	B	163	ASP	4
2	B	135	SER	3

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Mol	Chain	Res	Type	Models (Total)
2	B	171	GLU	3
2	B	177	ASN	3
2	B	172	MET	3
1	A	28	LYS	3
2	B	160	LEU	3
2	B	164	GLU	2
2	B	110	SER	2
1	A	21	LEU	2
2	B	154	ASP	2
2	B	184	ARG	2
1	A	55	GLU	1
1	A	46	GLN	1
2	B	113	SER	1
2	B	186	TYR	1
2	B	151	GLN	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