



Full wwPDB NMR Structure Validation Report ⓘ

Apr 26, 2016 – 10:42 PM BST

PDB ID : 2K4A
Title : FGF-1-C2A binary complex structure: a key component in the fibroblast growthfactor non-classical pathway
Authors : Yu, C.; Mohan, S.K.
Deposited on : 2008-06-02

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
<http://wwpdb.org/validation/2016/NMRValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

Cyrange : Kirchner and Güntert (2011)
NmrClust : Kelley et al. (1996)
MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : rb-20027457
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027457

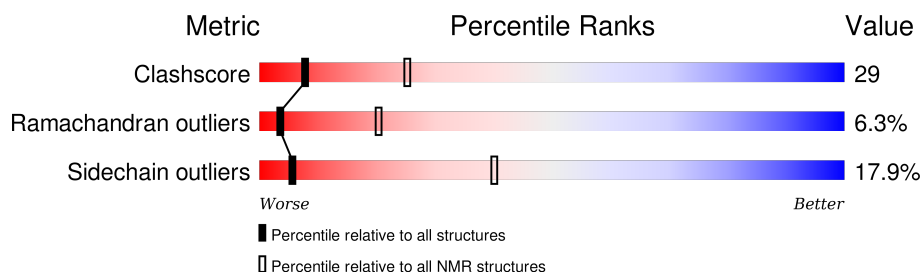
1 Overall quality at a glance

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	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

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	128	
2	B	133	

2 Ensemble composition and analysis

This entry contains 20 models. Model 1 is the overall representative, medoid model (most similar to other models).

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:1-A:128, B:1-B:133 (261)	0.35	1

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 2 single-model clusters were found.

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

3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 4201 atoms, of which 2093 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Synaptotagmin-1.

Mol	Chain	Residues	Atoms						Trace
1	A	128	Total	C	H	N	O	S	0
			2086	677	1042	169	195	3	

- Molecule 2 is a protein called Heparin-binding growth factor 1.

Mol	Chain	Residues	Atoms						Trace
2	B	133	Total	C	H	N	O	S	0
			2115	671	1051	185	204	4	

4 Residue-property plots

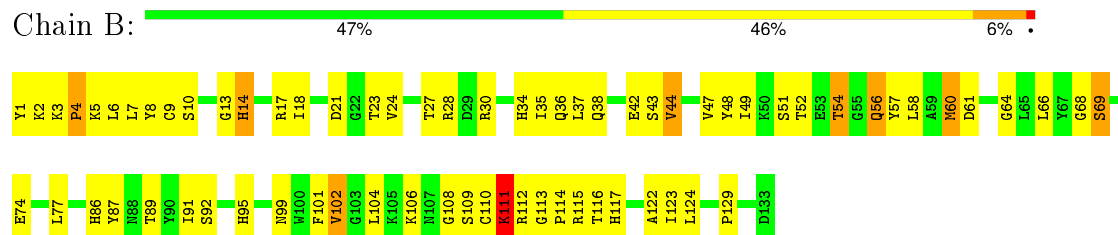
4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA and DNA 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: Synaptotagmin-1



- Molecule 2: Heparin-binding growth factor 1

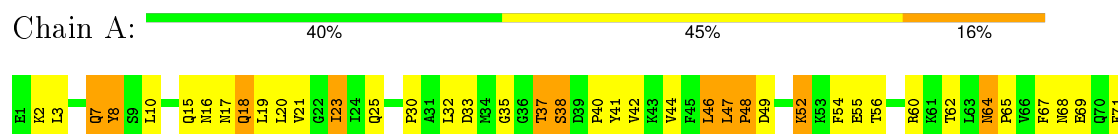


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1 (medoid)

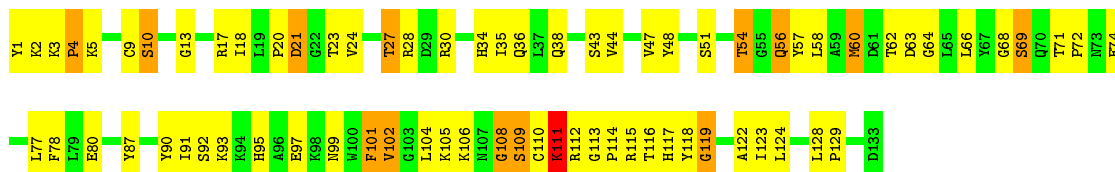
- Molecule 1: Synaptotagmin-1





• Molecule 2: Heparin-binding growth factor 1

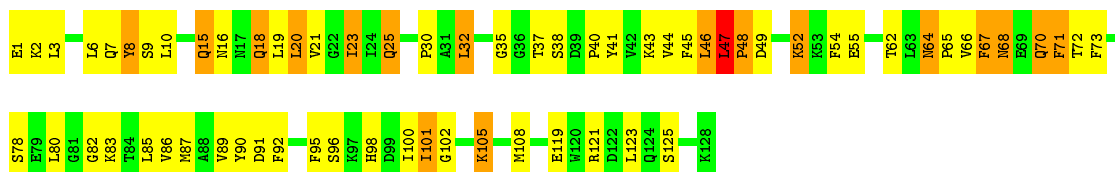
Chain B: 45% 44% 10%



4.2.2 Score per residue for model 2

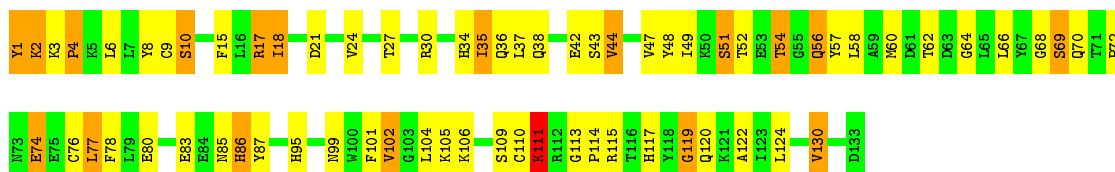
• Molecule 1: Synaptotagmin-1

Chain A: 48% 38% 13%



• Molecule 2: Heparin-binding growth factor 1

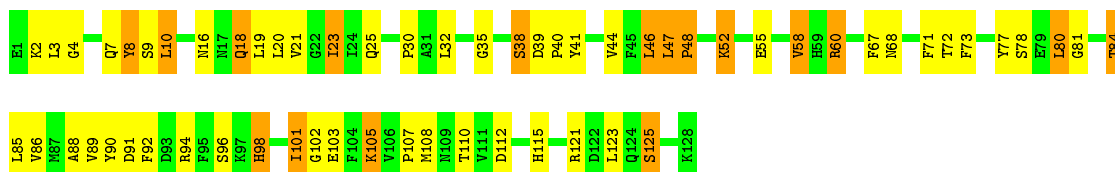
Chain B: 49% 37% 14%



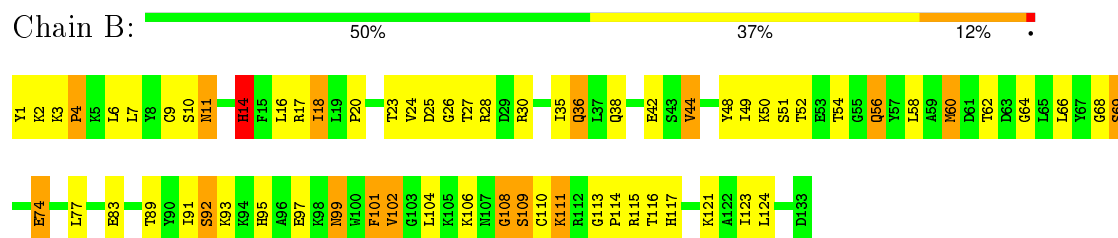
4.2.3 Score per residue for model 3

• Molecule 1: Synaptotagmin-1

Chain A: 52% 34% 13%

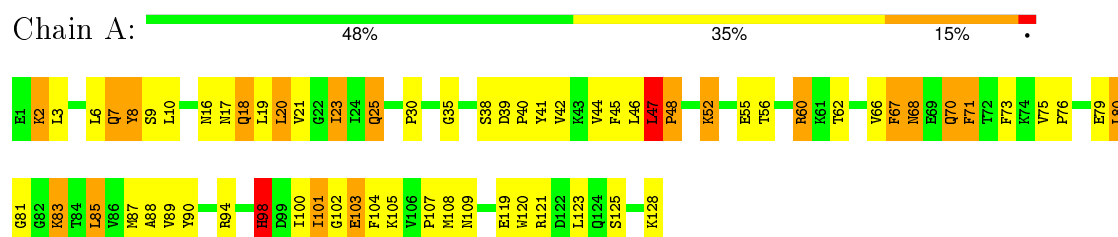


• Molecule 2: Heparin-binding growth factor 1

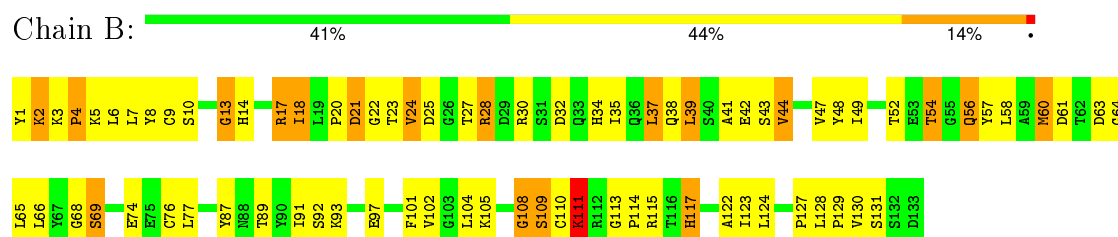


4.2.4 Score per residue for model 4

- Molecule 1: Synaptotagmin-1

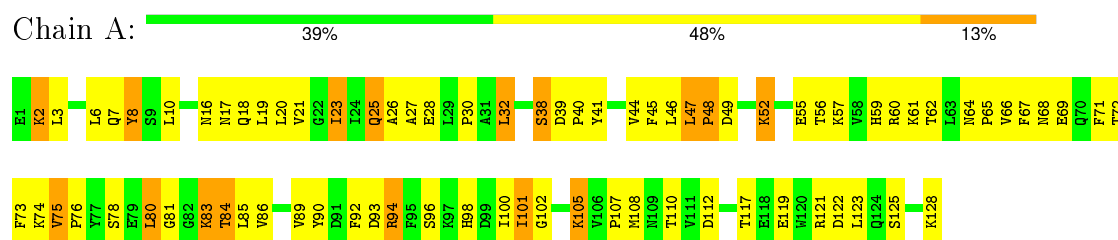


- Molecule 2: Heparin-binding growth factor 1

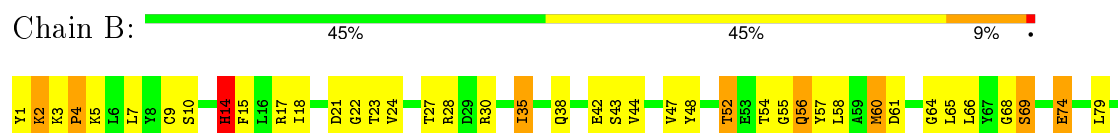


4.2.5 Score per residue for model 5

- Molecule 1: Synaptotagmin-1



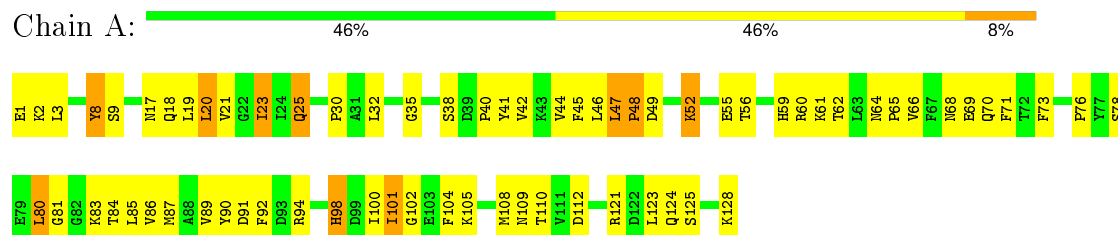
- Molecule 2: Heparin-binding growth factor 1



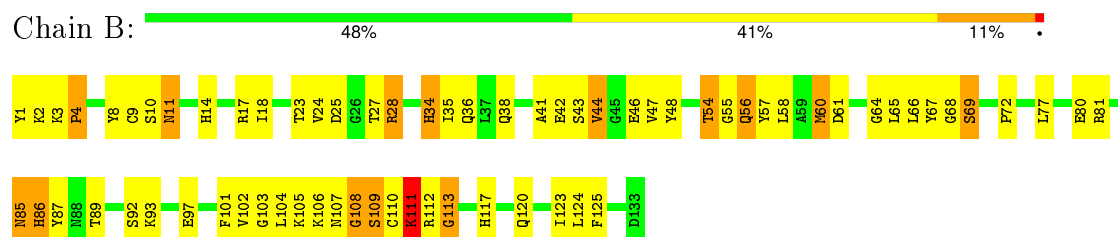


4.2.6 Score per residue for model 6

- Molecule 1: Synaptotagmin-1

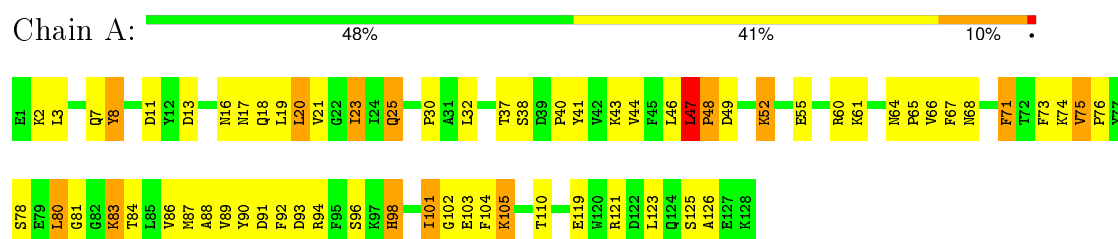


- Molecule 2: Heparin-binding growth factor 1

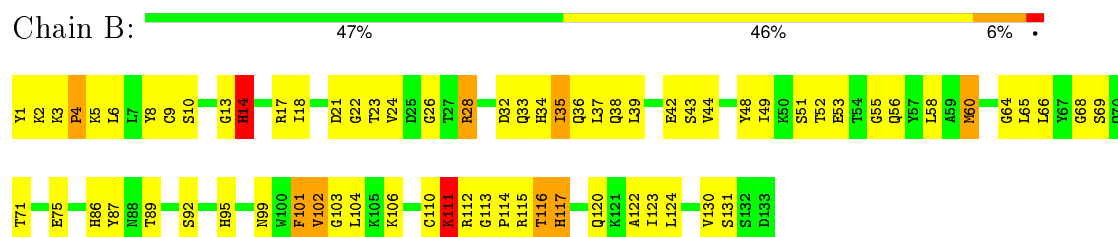


4.2.7 Score per residue for model 7

- Molecule 1: Synaptotagmin-1

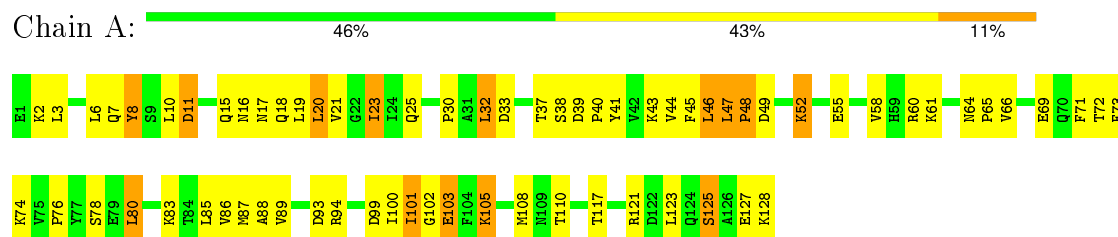


- Molecule 2: Heparin-binding growth factor 1

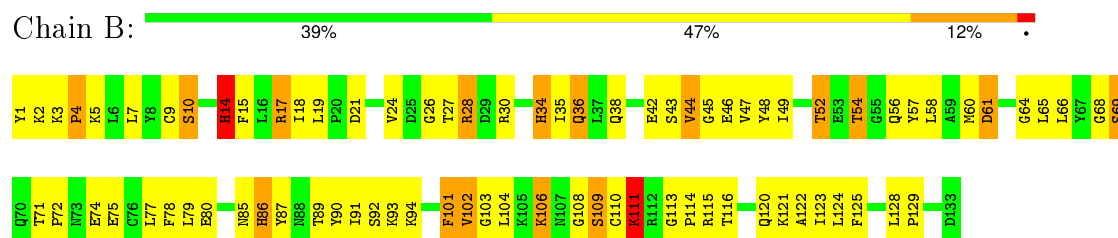


4.2.8 Score per residue for model 8

• Molecule 1: Synaptotagmin-1

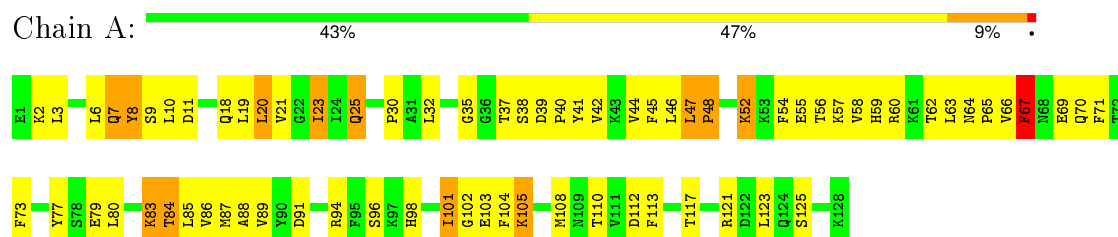


• Molecule 2: Heparin-binding growth factor 1

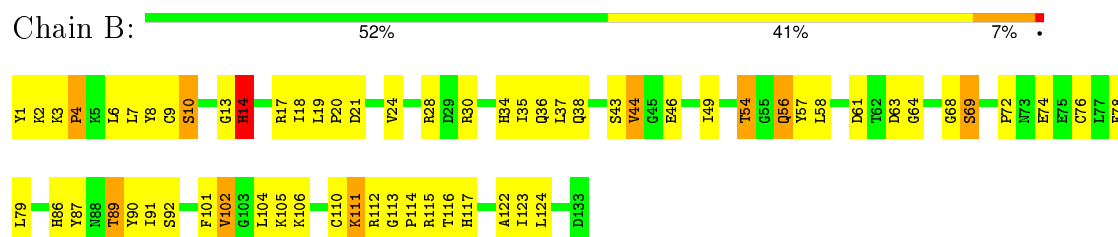


4.2.9 Score per residue for model 9

• Molecule 1: Synaptotagmin-1

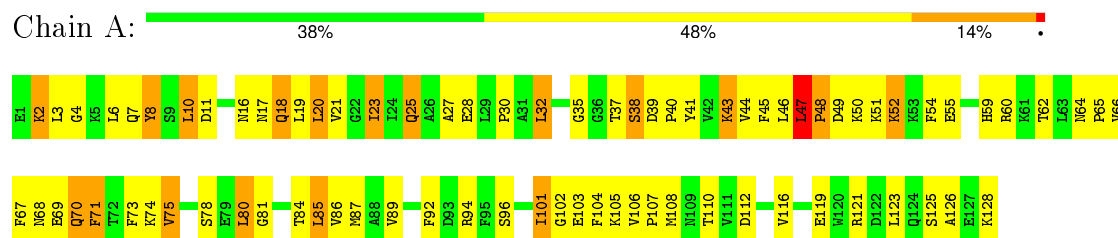


• Molecule 2: Heparin-binding growth factor 1

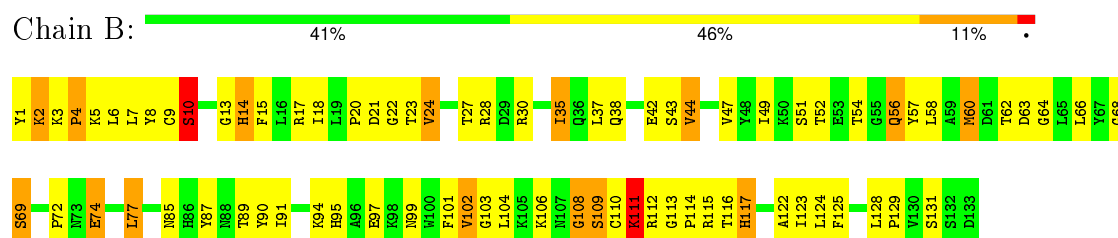


4.2.10 Score per residue for model 10

• Molecule 1: Synaptotagmin-1

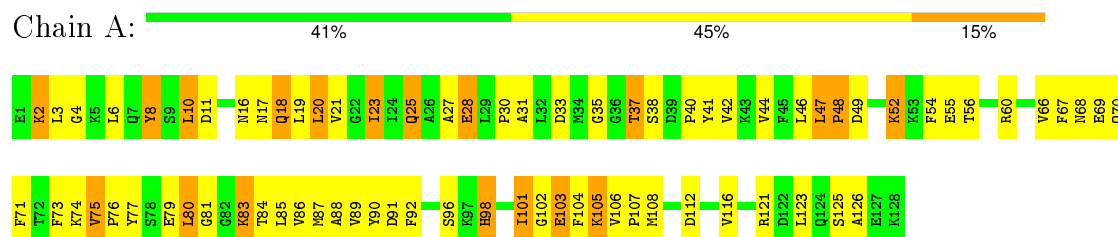


- Molecule 2: Heparin-binding growth factor 1

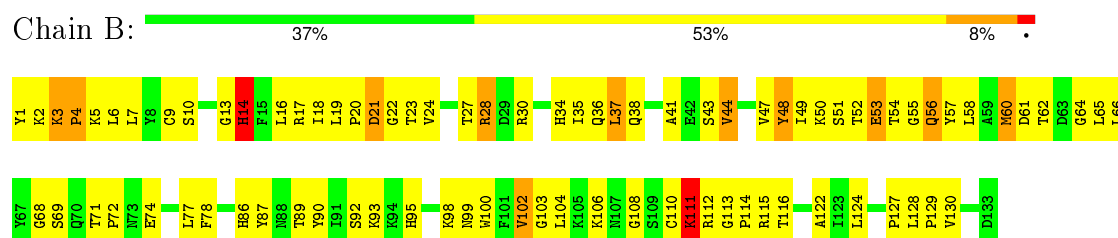


4.2.11 Score per residue for model 11

- Molecule 1: Synaptotagmin-1

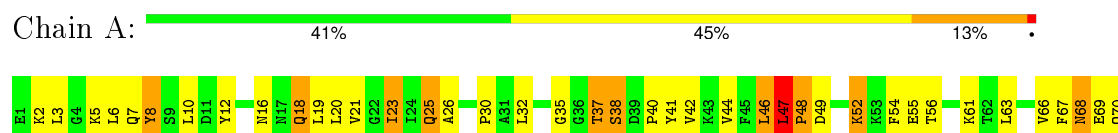


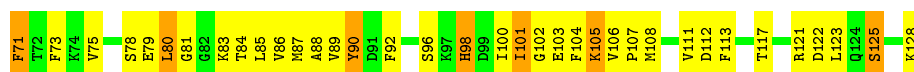
- Molecule 2: Heparin-binding growth factor 1



4.2.12 Score per residue for model 12

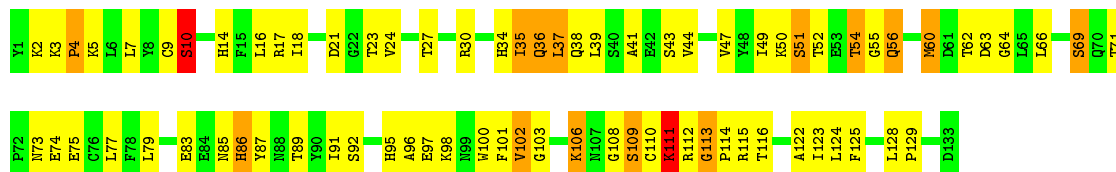
- Molecule 1: Synaptotagmin-1





- Molecule 2: Heparin-binding growth factor 1

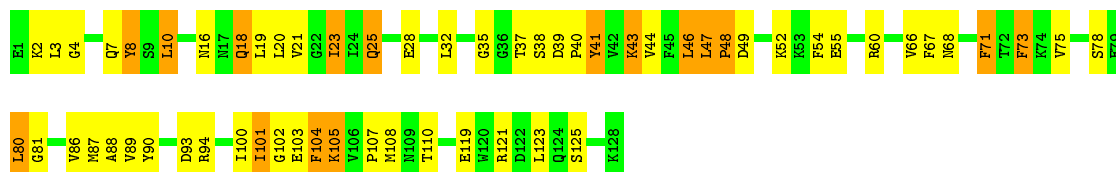
Chain B: 43% 45% 11%



4.2.13 Score per residue for model 13

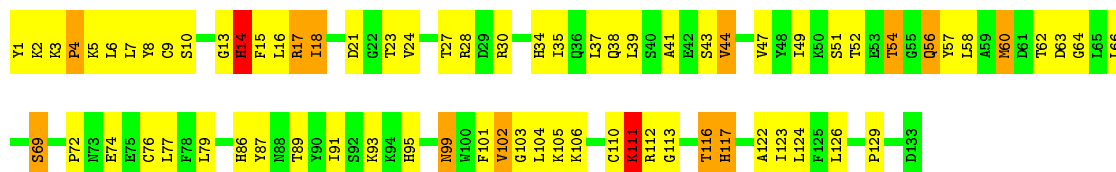
- Molecule 1: Synaptotagmin-1

Chain A: 53% 34% 13%



- Molecule 2: Heparin-binding growth factor 1

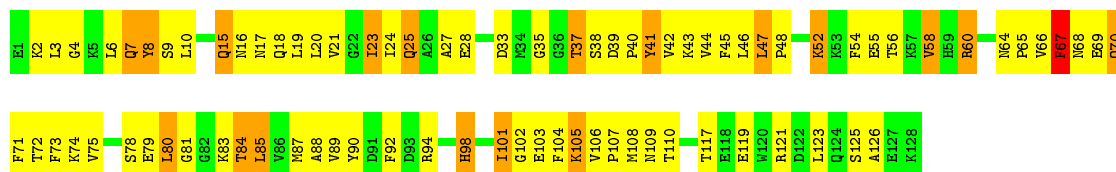
Chain B: 45% 44% 9%



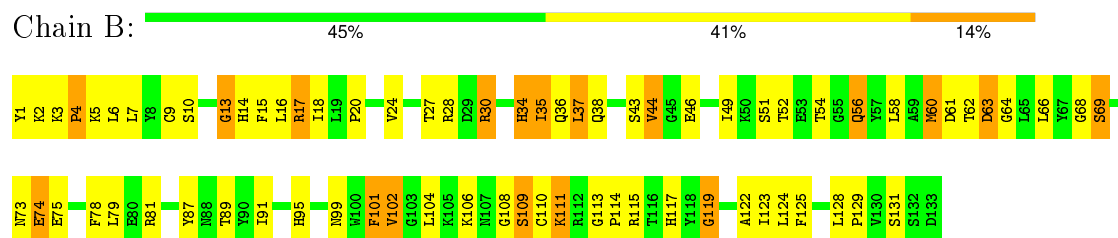
4.2.14 Score per residue for model 14

- Molecule 1: Synaptotagmin-1

Chain A: 36% 49% 14%

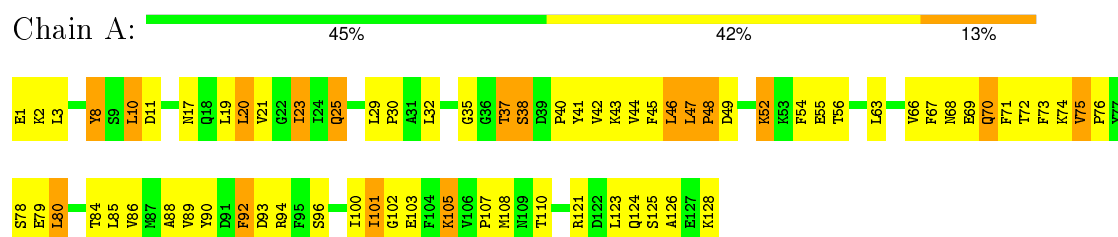


- Molecule 2: Heparin-binding growth factor 1

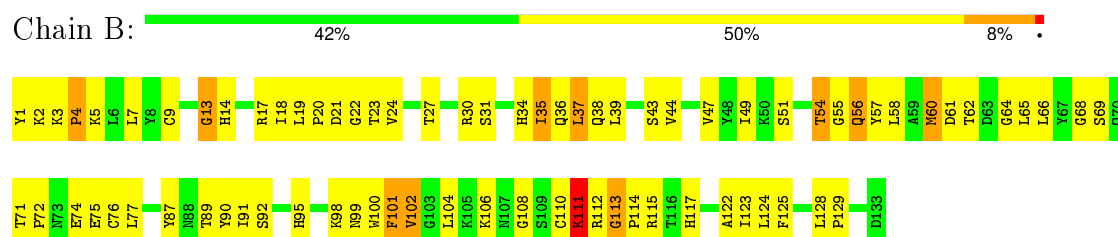


4.2.15 Score per residue for model 15

- Molecule 1: Synaptotagmin-1

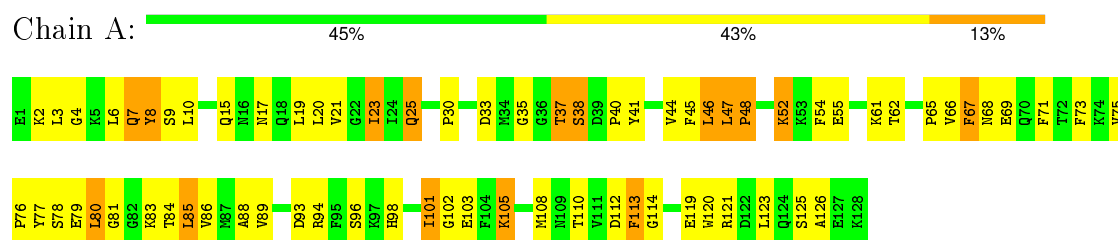


- Molecule 2: Heparin-binding growth factor 1

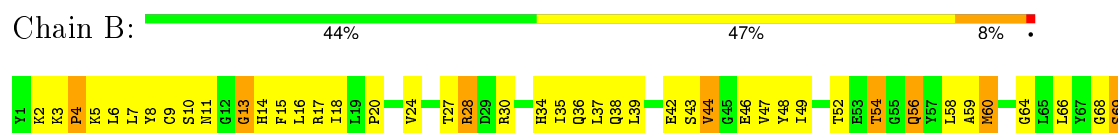


4.2.16 Score per residue for model 16

- Molecule 1: Synaptotagmin-1



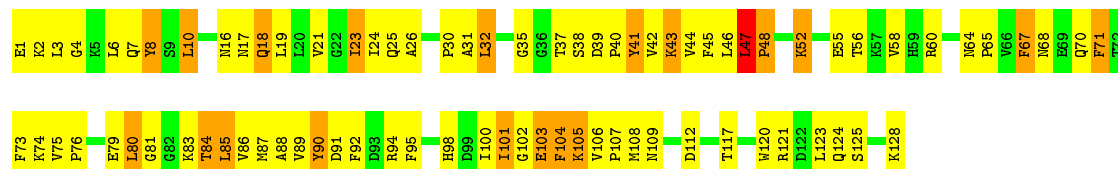
- Molecule 2: Heparin-binding growth factor 1



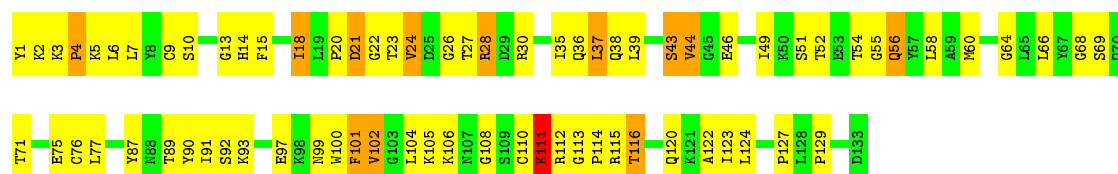


4.2.17 Score per residue for model 17

- Molecule 1: Synaptotagmin-1

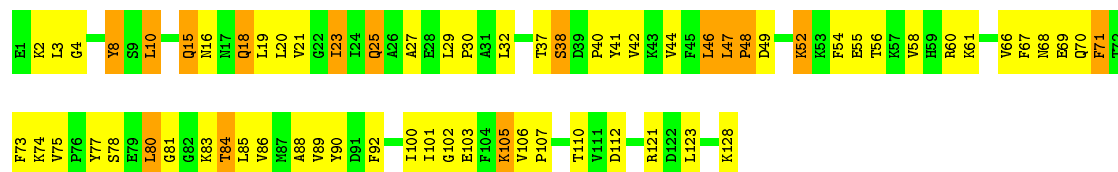


- Molecule 2: Heparin-binding growth factor 1

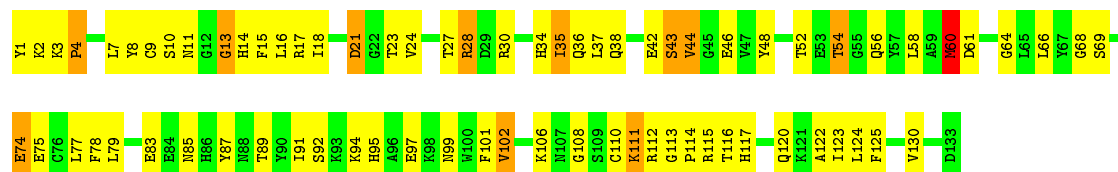


4.2.18 Score per residue for model 18

- Molecule 1: Synaptotagmin-1

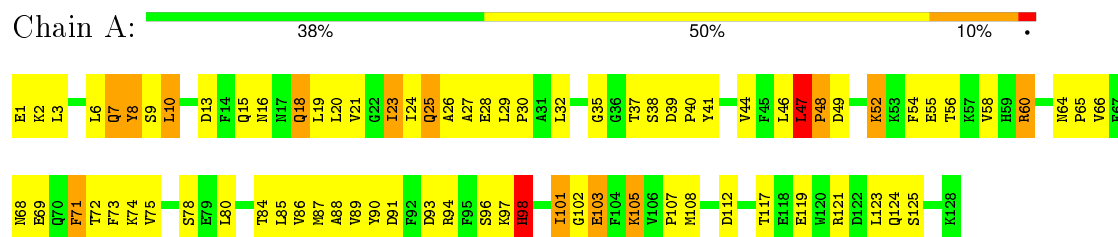


- Molecule 2: Heparin-binding growth factor 1

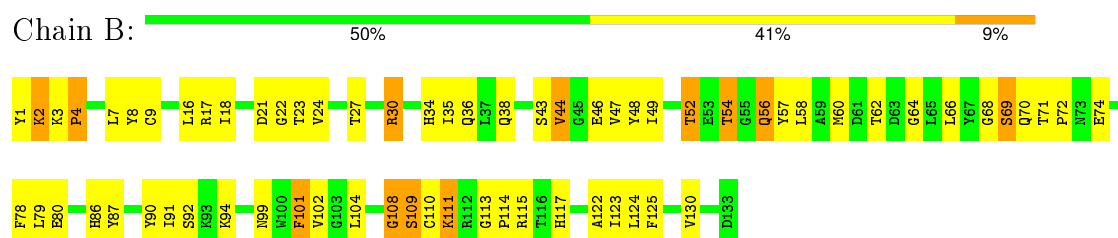


4.2.19 Score per residue for model 19

- Molecule 1: Synaptotagmin-1

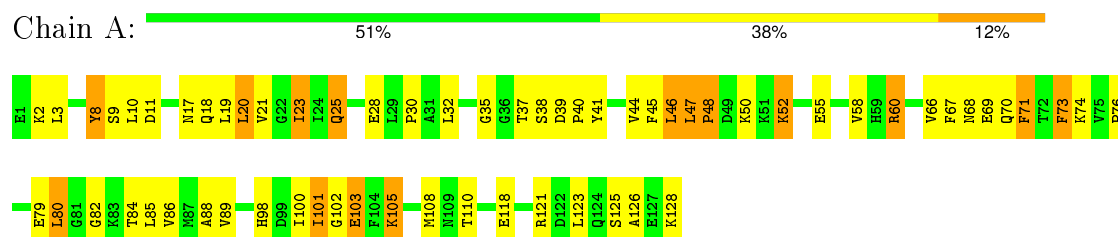


- Molecule 2: Heparin-binding growth factor 1

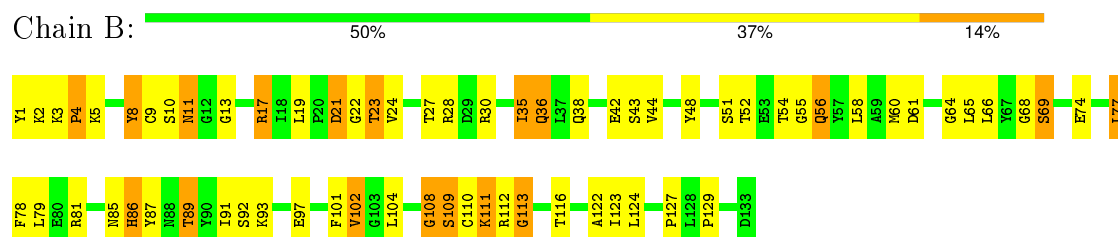


4.2.20 Score per residue for model 20

- Molecule 1: Synaptotagmin-1



- Molecule 2: Heparin-binding growth factor 1



5 Refinement protocol and experimental data overview

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

Of the 200 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
ARIA	structure solution	1.2
HADDOCK	structure solution	2.0
HADDOCK	refinement	2.0
ARIA/CNS	refinement	

No chemical shift data was provided. No validations of the models with respect to experimental NMR restraints is performed at this time.

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	1044	1042	1039	64±8
2	B	1064	1051	1046	64±8
All	All	42160	41860	41700	2459

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 29.

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:HD11	1:A:80:LEU:HD12	0.87	1.45	5	14
1:A:3:LEU:HD22	1:A:102:GLY:HA3	0.83	1.49	11	19
1:A:56:THR:HG22	1:A:69:GLU:HG2	0.81	1.53	18	8
1:A:80:LEU:HD23	1:A:81:GLY:N	0.79	1.92	3	12
2:B:101:PHE:CD2	2:B:123:ILE:HG21	0.79	2.13	9	14
1:A:44:VAL:HG21	1:A:71:PHE:CD1	0.77	2.15	11	16
1:A:19:LEU:HD11	1:A:80:LEU:CD1	0.76	2.10	17	5
1:A:44:VAL:HG11	1:A:71:PHE:CG	0.76	2.15	7	16
1:A:77:TYR:O	1:A:80:LEU:HD22	0.76	1.80	3	3
2:B:101:PHE:HD2	2:B:123:ILE:HG21	0.73	1.42	16	8
2:B:77:LEU:HD11	2:B:93:LYS:HB3	0.71	1.62	4	7
2:B:60:MET:HG3	2:B:66:LEU:HD22	0.71	1.62	13	8
1:A:3:LEU:HD12	1:A:30:PRO:CD	0.71	2.16	15	13

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:3:LEU:HD21	1:A:89:VAL:HG11	0.70	1.63	20	9
1:A:10:LEU:HD22	1:A:21:VAL:HG13	0.69	1.63	14	13
1:A:44:VAL:HG12	1:A:87:MET:SD	0.69	2.28	4	10
1:A:3:LEU:HD23	1:A:123:LEU:HB3	0.69	1.62	12	19
1:A:3:LEU:HD21	1:A:89:VAL:CG1	0.69	2.18	20	13
1:A:47:LEU:HB2	1:A:48:PRO:HD2	0.68	1.63	11	20
1:A:89:VAL:HG13	1:A:101:ILE:HG12	0.68	1.64	16	18
2:B:3:LYS:CB	2:B:4:PRO:HD2	0.68	2.19	12	20
2:B:3:LYS:HB3	2:B:4:PRO:HD2	0.67	1.66	11	20
2:B:66:LEU:HD12	2:B:110:CYS:HB2	0.67	1.65	1	9
1:A:47:LEU:CB	1:A:48:PRO:HD2	0.66	2.21	11	20
2:B:64:GLY:HA2	2:B:113:GLY:HA3	0.66	1.68	17	20
1:A:107:PRO:CB	2:B:27:THR:HG23	0.66	2.21	15	8
1:A:110:THR:HG21	2:B:28:ARG:CG	0.66	2.21	18	2
1:A:3:LEU:HD12	1:A:30:PRO:HD3	0.65	1.69	17	8
2:B:58:LEU:HA	2:B:68:GLY:HA3	0.65	1.69	10	16
2:B:42:GLU:HG3	2:B:48:TYR:CD1	0.64	2.27	8	7
1:A:44:VAL:HG21	1:A:71:PHE:CE1	0.64	2.26	6	8
2:B:24:VAL:HG13	2:B:110:CYS:SG	0.64	2.32	15	14
2:B:34:HIS:HA	2:B:54:THR:HG23	0.64	1.69	9	11
1:A:8:TYR:CZ	1:A:121:ARG:HB2	0.64	2.28	2	20
2:B:1:TYR:CE2	2:B:3:LYS:HD2	0.63	2.28	20	16
1:A:48:PRO:HD3	2:B:23:THR:HB	0.63	1.70	15	14
1:A:44:VAL:HG11	1:A:71:PHE:CD1	0.63	2.28	7	1
1:A:86:VAL:HG12	1:A:105:LYS:CB	0.63	2.24	18	11
2:B:24:VAL:CG1	2:B:110:CYS:SG	0.63	2.87	6	16
1:A:3:LEU:HD22	1:A:102:GLY:CA	0.63	2.23	2	12
1:A:23:ILE:HD11	1:A:69:GLU:HB3	0.62	1.70	19	3
1:A:110:THR:HG21	2:B:28:ARG:CD	0.62	2.25	10	7
1:A:41:TYR:CE1	1:A:55:GLU:HG3	0.61	2.30	6	19
1:A:86:VAL:HG12	1:A:105:LYS:HE3	0.61	1.72	2	3
1:A:48:PRO:HB3	2:B:21:ASP:CB	0.61	2.25	19	9
1:A:107:PRO:HB2	2:B:27:THR:HG23	0.61	1.72	11	5
2:B:104:LEU:CB	2:B:110:CYS:HA	0.61	2.26	17	6
1:A:6:LEU:HB2	1:A:123:LEU:HD21	0.61	1.71	17	3
1:A:85:LEU:HD22	1:A:108:MET:SD	0.61	2.36	10	8
2:B:9:CYS:O	2:B:10:SER:HB3	0.60	1.96	12	9
2:B:9:CYS:SG	2:B:122:ALA:HA	0.60	2.36	11	14
1:A:6:LEU:HB2	1:A:123:LEU:HD11	0.60	1.74	12	4
2:B:102:VAL:O	2:B:123:ILE:HG23	0.60	1.97	16	1
2:B:64:GLY:HA2	2:B:113:GLY:CA	0.60	2.26	17	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:101:PHE:HD1	2:B:123:ILE:HG21	0.60	1.57	19	1
2:B:104:LEU:HB2	2:B:110:CYS:HA	0.60	1.73	17	3
2:B:101:PHE:CE2	2:B:123:ILE:HG21	0.59	2.32	5	3
1:A:86:VAL:HG12	1:A:105:LYS:HB3	0.59	1.74	17	12
2:B:14:HIS:HB2	2:B:106:LYS:HE2	0.59	1.73	10	3
2:B:3:LYS:CG	2:B:4:PRO:HD2	0.59	2.28	19	17
1:A:8:TYR:CE1	1:A:121:ARG:HB2	0.59	2.33	8	8
2:B:101:PHE:CD1	2:B:123:ILE:HG21	0.59	2.32	19	1
2:B:112:ARG:HA	2:B:112:ARG:NE	0.59	2.12	15	3
1:A:19:LEU:HB2	1:A:73:PHE:HB3	0.59	1.73	12	19
1:A:8:TYR:CE2	1:A:121:ARG:HB2	0.59	2.32	12	4
2:B:1:TYR:O	2:B:130:VAL:HG23	0.58	1.98	2	5
1:A:80:LEU:HG	1:A:108:MET:HB3	0.58	1.76	2	12
2:B:14:HIS:HB2	2:B:106:LYS:CE	0.58	2.28	8	2
1:A:8:TYR:CB	1:A:23:ILE:HA	0.58	2.28	16	10
1:A:110:THR:HG21	2:B:28:ARG:HG3	0.58	1.74	18	3
2:B:9:CYS:O	2:B:10:SER:CB	0.58	2.52	8	2
2:B:17:ARG:O	2:B:24:VAL:HG23	0.58	1.99	13	11
1:A:110:THR:HG21	2:B:28:ARG:HD3	0.58	1.76	5	2
1:A:85:LEU:HB2	1:A:108:MET:SD	0.58	2.39	2	3
1:A:89:VAL:HG13	1:A:101:ILE:HG13	0.58	1.74	18	2
1:A:25:GLN:CB	1:A:66:VAL:HA	0.58	2.29	5	6
1:A:89:VAL:HG13	1:A:101:ILE:CG1	0.58	2.29	4	8
1:A:103:GLU:HG2	1:A:126:ALA:HB2	0.58	1.74	16	3
1:A:90:TYR:CZ	1:A:98:HIS:HB2	0.57	2.34	11	8
1:A:48:PRO:HD3	2:B:23:THR:CB	0.57	2.29	15	6
2:B:56:GLN:HB2	2:B:69:SER:HA	0.57	1.76	6	9
2:B:60:MET:HG3	2:B:66:LEU:CD2	0.57	2.30	3	14
1:A:3:LEU:CD2	1:A:123:LEU:HB3	0.57	2.29	19	8
2:B:4:PRO:HB3	2:B:36:GLN:HB2	0.57	1.76	3	13
2:B:90:TYR:O	2:B:102:VAL:HG12	0.57	1.98	17	2
2:B:14:HIS:HB2	2:B:106:LYS:HD2	0.57	1.76	17	3
1:A:1:GLU:HG2	1:A:124:GLN:HB3	0.57	1.77	17	1
1:A:46:LEU:HD21	1:A:73:PHE:CZ	0.57	2.35	8	2
1:A:75:VAL:HG21	1:A:83:LYS:HG2	0.57	1.77	11	2
1:A:6:LEU:HD21	1:A:23:ILE:HB	0.57	1.76	10	5
1:A:41:TYR:CE1	1:A:55:GLU:CG	0.56	2.88	15	5
2:B:58:LEU:HA	2:B:68:GLY:CA	0.56	2.30	3	9
1:A:46:LEU:HD21	1:A:73:PHE:CE2	0.56	2.35	3	8
2:B:4:PRO:HB3	2:B:36:GLN:CB	0.56	2.29	20	8
1:A:3:LEU:HD22	1:A:102:GLY:C	0.56	2.21	6	6

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:47:LEU:CB	1:A:48:PRO:CD	0.56	2.84	14	17
1:A:15:GLN:HG2	1:A:16:ASN:N	0.56	2.16	2	4
2:B:37:LEU:HD22	2:B:49:ILE:CG2	0.56	2.30	15	4
2:B:15:PHE:O	2:B:27:THR:N	0.56	2.39	10	7
2:B:18:ILE:HD12	2:B:51:SER:HB2	0.56	1.78	10	1
1:A:23:ILE:O	1:A:68:ASN:HA	0.56	2.01	2	18
1:A:46:LEU:HG	1:A:52:LYS:HD2	0.56	1.78	7	17
2:B:5:LYS:HB3	2:B:129:PRO:HB3	0.56	1.78	20	8
2:B:128:LEU:N	2:B:129:PRO:HD3	0.55	2.16	11	9
1:A:20:LEU:HD21	1:A:70:GLN:CG	0.55	2.31	9	2
2:B:14:HIS:HB3	2:B:106:LYS:HG2	0.55	1.78	5	3
2:B:87:TYR:CB	2:B:124:LEU:HB3	0.55	2.32	1	19
2:B:91:ILE:HA	2:B:101:PHE:HB3	0.55	1.78	1	11
2:B:83:GLU:HG2	2:B:85:ASN:HB3	0.55	1.78	2	3
2:B:108:GLY:O	2:B:109:SER:CB	0.55	2.54	6	5
2:B:17:ARG:HD3	2:B:27:THR:CB	0.55	2.30	6	2
1:A:25:GLN:HB3	1:A:66:VAL:HA	0.55	1.77	5	10
1:A:8:TYR:CE2	1:A:121:ARG:CB	0.55	2.90	12	5
1:A:107:PRO:HB3	2:B:27:THR:HG23	0.55	1.78	1	5
1:A:18:GLN:OE1	1:A:72:THR:HG23	0.55	2.02	8	5
1:A:86:VAL:CG1	1:A:105:LYS:HE3	0.55	2.31	10	3
1:A:105:LYS:HD3	2:B:108:GLY:HA2	0.55	1.77	8	8
2:B:66:LEU:HB2	2:B:110:CYS:HB2	0.55	1.79	17	9
2:B:114:PRO:C	2:B:115:ARG:HD2	0.55	2.23	1	15
2:B:92:SER:HB2	2:B:101:PHE:HA	0.55	1.79	9	3
1:A:3:LEU:HB3	1:A:123:LEU:HB3	0.54	1.79	2	13
1:A:3:LEU:HD12	1:A:30:PRO:HD2	0.54	1.79	15	3
2:B:64:GLY:HA2	2:B:114:PRO:HD3	0.54	1.79	1	3
2:B:101:PHE:O	2:B:116:THR:HG22	0.54	2.02	7	1
1:A:46:LEU:HD13	1:A:84:THR:O	0.54	2.01	9	6
2:B:61:ASP:HB3	2:B:65:LEU:HB2	0.54	1.79	11	6
2:B:14:HIS:HB3	2:B:106:LYS:CG	0.54	2.31	5	2
1:A:41:TYR:HB3	1:A:58:VAL:HA	0.54	1.77	9	4
2:B:9:CYS:HB2	2:B:125:PHE:HA	0.54	1.78	10	5
2:B:13:GLY:O	2:B:14:HIS:CD2	0.54	2.61	10	9
2:B:62:THR:HA	2:B:95:HIS:CD2	0.54	2.37	11	1
1:A:88:ALA:HA	1:A:103:GLU:HB3	0.54	1.79	20	7
1:A:3:LEU:CG	1:A:123:LEU:HB3	0.54	2.33	3	10
2:B:43:SER:HB2	2:B:46:GLU:HB2	0.54	1.79	9	5
1:A:103:GLU:HG2	1:A:126:ALA:CB	0.54	2.33	14	2
1:A:75:VAL:HG21	1:A:83:LYS:HE3	0.54	1.80	7	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:45:PHE:O	1:A:46:LEU:HD22	0.54	2.02	9	10
2:B:104:LEU:HG	2:B:125:PHE:CE1	0.54	2.38	15	1
1:A:38:SER:C	1:A:40:PRO:HD3	0.54	2.23	8	20
2:B:10:SER:HB3	2:B:106:LYS:HD3	0.53	1.79	12	10
1:A:39:ASP:CG	1:A:60:ARG:HA	0.53	2.24	4	9
2:B:57:TYR:CD2	2:B:72:PRO:HB3	0.53	2.37	13	1
2:B:62:THR:HA	2:B:95:HIS:CE1	0.53	2.38	10	4
2:B:42:GLU:HG3	2:B:48:TYR:CG	0.53	2.39	6	5
1:A:3:LEU:O	1:A:3:LEU:HG	0.53	2.03	1	6
2:B:14:HIS:CG	2:B:26:GLY:O	0.53	2.62	7	3
2:B:6:LEU:HD23	2:B:7:LEU:N	0.53	2.19	9	8
1:A:20:LEU:HD11	1:A:70:GLN:HG3	0.53	1.78	9	2
1:A:21:VAL:HB	1:A:71:PHE:HB3	0.53	1.80	9	16
2:B:78:PHE:CB	2:B:90:TYR:HB3	0.53	2.34	9	1
1:A:32:LEU:HD13	1:A:32:LEU:O	0.53	2.04	2	3
1:A:23:ILE:HD13	1:A:23:ILE:H	0.53	1.64	8	4
2:B:47:VAL:HG21	2:B:90:TYR:CE1	0.53	2.39	16	2
2:B:104:LEU:HA	2:B:110:CYS:HA	0.53	1.81	9	4
1:A:45:PHE:HB2	1:A:86:VAL:HG23	0.53	1.80	16	2
2:B:37:LEU:HD22	2:B:49:ILE:HG23	0.52	1.80	17	3
2:B:20:PRO:HA	2:B:56:GLN:HG3	0.52	1.81	1	6
1:A:47:LEU:CG	1:A:48:PRO:HD2	0.52	2.35	14	5
2:B:108:GLY:O	2:B:109:SER:HB2	0.52	2.04	12	10
2:B:75:GLU:HG2	2:B:94:LYS:HG2	0.52	1.80	18	1
2:B:9:CYS:SG	2:B:124:LEU:O	0.52	2.65	17	14
2:B:16:LEU:HG	2:B:24:VAL:CG2	0.52	2.35	18	5
2:B:56:GLN:HG2	2:B:69:SER:HA	0.52	1.79	3	2
2:B:14:HIS:HB2	2:B:106:LYS:HG3	0.52	1.81	6	1
1:A:49:ASP:CB	1:A:83:LYS:HD2	0.52	2.35	7	1
2:B:60:MET:CG	2:B:66:LEU:HD22	0.52	2.34	10	4
1:A:88:ALA:HA	1:A:103:GLU:CB	0.52	2.35	15	6
1:A:67:PHE:CD1	1:A:69:GLU:HG3	0.52	2.40	14	1
2:B:41:ALA:HA	2:B:47:VAL:HA	0.52	1.81	12	5
1:A:26:ALA:HB1	1:A:123:LEU:HD11	0.52	1.82	17	2
1:A:100:ILE:HD12	1:A:128:LYS:HG3	0.52	1.82	8	5
2:B:101:PHE:C	2:B:116:THR:HG21	0.52	2.25	9	3
2:B:78:PHE:CE1	2:B:102:VAL:HB	0.52	2.40	14	3
2:B:5:LYS:HD3	2:B:129:PRO:HB2	0.52	1.82	15	6
2:B:5:LYS:CD	2:B:129:PRO:HB2	0.52	2.33	5	7
2:B:49:ILE:N	2:B:49:ILE:HD12	0.52	2.20	12	5
2:B:60:MET:HG3	2:B:66:LEU:HA	0.51	1.82	4	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:44:VAL:CG2	1:A:54:PHE:HB3	0.51	2.35	2	10
1:A:16:ASN:OD1	1:A:18:GLN:HG3	0.51	2.04	11	13
2:B:17:ARG:NH1	2:B:27:THR:OG1	0.51	2.40	18	9
2:B:66:LEU:HD13	2:B:102:VAL:HG23	0.51	1.81	13	5
2:B:17:ARG:HD2	2:B:27:THR:HB	0.51	1.82	11	1
2:B:77:LEU:CD1	2:B:93:LYS:HB3	0.51	2.35	3	9
2:B:127:PRO:C	2:B:129:PRO:HD3	0.51	2.24	5	6
2:B:104:LEU:C	2:B:104:LEU:HD12	0.51	2.25	9	7
1:A:32:LEU:O	1:A:32:LEU:HD13	0.51	2.04	6	1
2:B:10:SER:HB3	2:B:106:LYS:HB2	0.51	1.82	6	4
2:B:56:GLN:CB	2:B:69:SER:HA	0.51	2.36	13	6
2:B:3:LYS:CB	2:B:4:PRO:CD	0.51	2.87	11	18
2:B:47:VAL:HG21	2:B:90:TYR:CD1	0.51	2.41	10	3
1:A:7:GLN:HB3	1:A:120:TRP:CE3	0.51	2.41	1	3
2:B:79:LEU:HB3	2:B:91:ILE:HB	0.51	1.83	5	9
2:B:34:HIS:HA	2:B:54:THR:CG2	0.51	2.36	18	3
2:B:113:GLY:HA2	2:B:116:THR:HG23	0.51	1.81	1	1
2:B:104:LEU:HB3	2:B:110:CYS:CB	0.51	2.36	11	1
1:A:92:PHE:HA	1:A:98:HIS:CE1	0.51	2.41	17	2
2:B:49:ILE:HD13	2:B:78:PHE:CD1	0.51	2.40	19	1
1:A:80:LEU:HD12	1:A:108:MET:HE2	0.51	1.82	15	2
2:B:102:VAL:HG13	2:B:102:VAL:O	0.51	2.06	17	1
1:A:102:GLY:HA2	1:A:125:SER:HA	0.51	1.83	3	4
1:A:41:TYR:CE2	1:A:90:TYR:HB3	0.51	2.41	17	3
2:B:74:GLU:HB2	2:B:94:LYS:HD3	0.51	1.83	8	2
1:A:47:LEU:HB2	1:A:48:PRO:CD	0.51	2.36	14	2
2:B:117:HIS:O	2:B:119:GLY:N	0.50	2.44	1	3
1:A:77:TYR:HB2	1:A:113:PHE:CZ	0.50	2.41	16	2
1:A:3:LEU:HB3	1:A:123:LEU:C	0.50	2.26	14	12
2:B:57:TYR:HB3	2:B:72:PRO:HB3	0.50	1.84	8	9
2:B:65:LEU:HD12	2:B:65:LEU:N	0.50	2.21	20	1
2:B:62:THR:HA	2:B:95:HIS:NE2	0.50	2.21	10	4
2:B:76:CYS:SG	2:B:76:CYS:O	0.50	2.69	13	3
2:B:10:SER:HB2	2:B:106:LYS:HB2	0.50	1.83	2	1
2:B:13:GLY:C	2:B:28:ARG:HB3	0.50	2.26	11	1
1:A:3:LEU:HD13	1:A:102:GLY:HA3	0.50	1.84	7	3
2:B:5:LYS:HA	2:B:129:PRO:HB3	0.50	1.83	8	2
2:B:78:PHE:CD2	2:B:92:SER:HA	0.50	2.41	9	1
2:B:22:GLY:HA2	2:B:68:GLY:O	0.50	2.06	4	9
2:B:100:TRP:CZ2	2:B:114:PRO:HA	0.50	2.42	12	3
2:B:68:GLY:O	2:B:69:SER:HB2	0.50	2.06	14	8

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:77:LEU:HD11	2:B:94:LYS:HB2	0.50	1.82	10	1
1:A:11:ASP:HB3	1:A:116:VAL:HG12	0.50	1.84	10	1
2:B:14:HIS:HB2	2:B:106:LYS:HE3	0.50	1.84	8	1
2:B:113:GLY:HA2	2:B:116:THR:CG2	0.50	2.37	1	1
2:B:60:MET:SD	2:B:66:LEU:HA	0.50	2.46	17	1
1:A:64:ASN:N	1:A:65:PRO:HD3	0.50	2.21	14	10
1:A:107:PRO:HG2	2:B:27:THR:HA	0.50	1.83	5	2
1:A:71:PHE:C	1:A:71:PHE:CD1	0.50	2.84	2	4
2:B:60:MET:HG3	2:B:66:LEU:HD23	0.50	1.83	11	3
1:A:46:LEU:CD2	1:A:52:LYS:HD2	0.49	2.36	1	12
2:B:91:ILE:HA	2:B:101:PHE:CB	0.49	2.36	5	8
2:B:13:GLY:O	2:B:28:ARG:HB3	0.49	2.06	4	7
2:B:101:PHE:CD2	2:B:123:ILE:HD13	0.49	2.43	15	2
1:A:44:VAL:HG11	1:A:71:PHE:CD2	0.49	2.42	6	3
1:A:103:GLU:HG2	1:A:126:ALA:HB3	0.49	1.83	1	2
2:B:81:ARG:HB3	2:B:89:THR:HB	0.49	1.84	14	3
2:B:17:ARG:C	2:B:24:VAL:HG23	0.49	2.28	19	3
2:B:77:LEU:HD12	2:B:77:LEU:O	0.49	2.07	10	2
1:A:11:ASP:CB	1:A:116:VAL:HG12	0.49	2.36	11	1
1:A:20:LEU:HD11	1:A:70:GLN:HG2	0.49	1.83	6	2
2:B:18:ILE:HD12	2:B:51:SER:CB	0.49	2.37	10	1
1:A:19:LEU:HB2	1:A:73:PHE:CB	0.49	2.37	5	13
1:A:88:ALA:CB	1:A:103:GLU:HB3	0.49	2.37	8	7
1:A:48:PRO:HG2	2:B:19:LEU:HD22	0.49	1.83	20	3
1:A:46:LEU:HD11	1:A:73:PHE:CE1	0.49	2.42	11	1
2:B:103:GLY:N	2:B:116:THR:CG2	0.49	2.76	11	1
2:B:44:VAL:HG13	2:B:44:VAL:O	0.49	2.08	17	7
2:B:105:LYS:HG2	2:B:112:ARG:HD2	0.49	1.83	17	1
2:B:7:LEU:HD11	2:B:125:PHE:HB2	0.49	1.85	8	5
1:A:30:PRO:HD2	1:A:101:ILE:HB	0.49	1.84	19	6
1:A:4:GLY:HA3	1:A:28:GLU:HB2	0.49	1.85	13	3
2:B:35:ILE:H	2:B:35:ILE:HD13	0.49	1.68	15	1
2:B:51:SER:HB2	2:B:55:GLY:HA3	0.49	1.83	7	2
1:A:10:LEU:O	1:A:117:THR:N	0.49	2.41	17	5
2:B:44:VAL:O	2:B:44:VAL:HG13	0.49	2.08	20	8
1:A:10:LEU:HD22	1:A:21:VAL:HA	0.49	1.84	1	3
1:A:100:ILE:HD12	1:A:128:LYS:CB	0.49	2.38	15	3
1:A:42:VAL:CG2	1:A:56:THR:HG21	0.49	2.37	15	6
2:B:85:ASN:O	2:B:86:HIS:HB3	0.49	2.08	6	5
2:B:114:PRO:HB2	2:B:115:ARG:HD2	0.49	1.85	1	8
2:B:49:ILE:HD12	2:B:49:ILE:N	0.49	2.23	11	9

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:104:LEU:HD12	2:B:104:LEU:C	0.49	2.28	2	5
1:A:85:LEU:C	1:A:85:LEU:HD23	0.49	2.28	9	3
1:A:79:GLU:O	1:A:83:LYS:HD3	0.49	2.08	14	2
1:A:4:GLY:C	1:A:123:LEU:HB2	0.48	2.28	16	3
1:A:43:LYS:HE2	1:A:90:TYR:HB2	0.48	1.85	15	3
1:A:5:LYS:HG2	1:A:122:ASP:HA	0.48	1.84	12	1
1:A:3:LEU:HD11	1:A:89:VAL:HG11	0.48	1.84	15	2
2:B:103:GLY:HA3	2:B:123:ILE:HG12	0.48	1.84	7	2
2:B:74:GLU:O	2:B:77:LEU:HG	0.48	2.07	2	4
1:A:8:TYR:HB3	1:A:23:ILE:HA	0.48	1.84	5	3
1:A:86:VAL:O	1:A:86:VAL:HG23	0.48	2.08	2	1
1:A:37:THR:HA	1:A:63:LEU:HD21	0.48	1.84	12	2
1:A:48:PRO:CD	2:B:23:THR:HB	0.48	2.38	10	6
1:A:10:LEU:CD2	1:A:21:VAL:HG13	0.48	2.39	9	5
2:B:16:LEU:HG	2:B:24:VAL:HG21	0.48	1.83	13	4
2:B:20:PRO:HA	2:B:56:GLN:CG	0.48	2.37	10	8
2:B:48:TYR:CE1	2:B:77:LEU:HB3	0.48	2.43	18	2
1:A:47:LEU:HG	1:A:48:PRO:HD2	0.48	1.83	14	1
1:A:52:LYS:HE2	1:A:73:PHE:CD1	0.48	2.44	20	5
1:A:85:LEU:HD21	1:A:106:VAL:HG12	0.48	1.85	17	3
2:B:87:TYR:HB3	2:B:124:LEU:HB3	0.48	1.86	10	8
1:A:90:TYR:CE2	1:A:98:HIS:HB2	0.48	2.43	1	1
1:A:25:GLN:CB	1:A:66:VAL:HG12	0.48	2.38	10	9
1:A:85:LEU:HD23	1:A:86:VAL:N	0.48	2.23	20	2
2:B:14:HIS:CB	2:B:106:LYS:HE2	0.48	2.39	10	1
2:B:99:ASN:OD1	2:B:117:HIS:NE2	0.48	2.47	7	3
2:B:124:LEU:HD22	2:B:124:LEU:N	0.48	2.24	8	6
2:B:60:MET:N	2:B:60:MET:SD	0.48	2.86	2	2
2:B:51:SER:CB	2:B:55:GLY:HA3	0.48	2.39	7	1
2:B:9:CYS:HB2	2:B:125:PHE:CD1	0.48	2.44	8	1
1:A:44:VAL:HG22	1:A:54:PHE:HB3	0.48	1.85	12	2
2:B:66:LEU:HD11	2:B:103:GLY:O	0.48	2.09	12	2
2:B:5:LYS:NZ	2:B:39:LEU:O	0.48	2.45	17	3
1:A:44:VAL:HG12	1:A:87:MET:CE	0.48	2.39	6	4
1:A:3:LEU:HD23	1:A:123:LEU:HD22	0.48	1.86	17	2
2:B:9:CYS:SG	2:B:122:ALA:C	0.48	2.92	10	3
2:B:49:ILE:CG2	2:B:58:LEU:HD22	0.48	2.39	10	1
2:B:49:ILE:HB	2:B:58:LEU:HD12	0.48	1.85	16	1
1:A:49:ASP:CG	1:A:83:LYS:HD2	0.48	2.30	6	1
2:B:128:LEU:N	2:B:129:PRO:CD	0.47	2.76	1	4
2:B:37:LEU:HD13	2:B:49:ILE:HG23	0.47	1.84	16	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:56:GLN:N	2:B:56:GLN:OE1	0.47	2.47	17	3
2:B:10:SER:HB2	2:B:106:LYS:CB	0.47	2.38	3	1
1:A:48:PRO:HG3	2:B:21:ASP:HB3	0.47	1.86	5	3
1:A:105:LYS:HD3	2:B:108:GLY:CA	0.47	2.39	5	2
2:B:105:LYS:HG3	2:B:112:ARG:HD2	0.47	1.84	5	1
1:A:3:LEU:HG	1:A:3:LEU:O	0.47	2.08	8	3
2:B:63:ASP:HB2	2:B:95:HIS:CD2	0.47	2.45	1	1
1:A:4:GLY:CA	1:A:28:GLU:HB2	0.47	2.39	13	1
1:A:29:LEU:HD13	1:A:40:PRO:HB2	0.47	1.86	15	1
1:A:17:ASN:OD1	1:A:76:PRO:HA	0.47	2.09	11	11
2:B:60:MET:HB3	2:B:66:LEU:HD23	0.47	1.86	17	2
2:B:1:TYR:CZ	2:B:3:LYS:HD2	0.47	2.45	18	3
2:B:3:LYS:HB3	2:B:4:PRO:CD	0.47	2.38	10	14
1:A:47:LEU:HD11	1:A:82:GLY:O	0.47	2.08	20	1
1:A:87:MET:HG2	1:A:104:PHE:CE2	0.47	2.45	10	3
1:A:86:VAL:HG12	1:A:105:LYS:HB2	0.47	1.86	6	2
2:B:35:ILE:HD13	2:B:35:ILE:H	0.47	1.70	2	4
2:B:78:PHE:CE2	2:B:102:VAL:HB	0.47	2.45	9	5
1:A:19:LEU:HD12	1:A:73:PHE:CD2	0.47	2.44	20	1
1:A:62:THR:CG2	1:A:65:PRO:HB3	0.47	2.40	2	1
2:B:14:HIS:HA	2:B:28:ARG:HB3	0.47	1.87	17	1
2:B:51:SER:HB3	2:B:54:THR:HB	0.47	1.86	17	1
1:A:90:TYR:CD1	1:A:100:ILE:HG12	0.47	2.44	13	6
1:A:6:LEU:HD23	1:A:104:PHE:CE1	0.47	2.45	12	1
2:B:47:VAL:HG13	2:B:80:GLU:HB2	0.47	1.86	6	2
1:A:41:TYR:CE1	1:A:43:LYS:HE2	0.47	2.45	10	1
1:A:105:LYS:HB3	1:A:105:LYS:NZ	0.47	2.25	15	1
2:B:102:VAL:N	2:B:116:THR:HG21	0.47	2.25	11	6
1:A:92:PHE:O	1:A:94:ARG:CD	0.47	2.63	10	7
1:A:57:LYS:HB3	1:A:67:PHE:CZ	0.47	2.44	9	2
1:A:43:LYS:HE3	1:A:90:TYR:HB2	0.47	1.86	2	1
2:B:86:HIS:ND1	2:B:86:HIS:O	0.47	2.48	19	2
2:B:10:SER:O	2:B:11:ASN:C	0.47	2.52	6	3
1:A:58:VAL:HG13	1:A:58:VAL:O	0.47	2.10	3	3
2:B:7:LEU:HD11	2:B:125:PHE:CB	0.47	2.40	8	2
1:A:105:LYS:CD	2:B:108:GLY:HA2	0.47	2.39	8	2
2:B:24:VAL:HG11	2:B:110:CYS:SG	0.46	2.51	11	2
2:B:75:GLU:HG2	2:B:94:LYS:HE2	0.46	1.88	18	1
1:A:100:ILE:HB	1:A:128:LYS:HB2	0.46	1.86	15	2
1:A:104:PHE:CE1	1:A:121:ARG:HG2	0.46	2.44	4	1
2:B:110:CYS:C	2:B:111:LYS:HD2	0.46	2.31	16	12

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:47:LEU:HG	1:A:48:PRO:CD	0.46	2.40	10	10
1:A:84:THR:HG23	1:A:107:PRO:HA	0.46	1.87	14	3
1:A:23:ILE:H	1:A:23:ILE:HD13	0.46	1.69	19	1
2:B:8:TYR:HB2	2:B:14:HIS:O	0.46	2.11	4	1
2:B:98:LYS:HD3	2:B:100:TRP:CH2	0.46	2.45	11	2
2:B:77:LEU:HD12	2:B:93:LYS:HB3	0.46	1.86	3	1
2:B:37:LEU:CD1	2:B:39:LEU:HB2	0.46	2.40	12	3
2:B:101:PHE:N	2:B:116:THR:HG22	0.46	2.26	17	1
1:A:47:LEU:HD23	1:A:49:ASP:HB3	0.46	1.88	2	2
2:B:47:VAL:CG1	2:B:80:GLU:HB2	0.46	2.41	1	3
2:B:17:ARG:HG3	2:B:35:ILE:HG23	0.46	1.87	20	1
2:B:5:LYS:CA	2:B:129:PRO:HB3	0.46	2.41	8	1
2:B:92:SER:O	2:B:96:ALA:N	0.46	2.48	12	1
2:B:17:ARG:HG3	2:B:34:HIS:HB3	0.46	1.87	14	1
2:B:89:THR:HG23	2:B:124:LEU:HD13	0.46	1.86	9	1
2:B:18:ILE:HD12	2:B:55:GLY:HA3	0.46	1.86	6	1
2:B:86:HIS:O	2:B:86:HIS:ND1	0.46	2.48	11	4
2:B:18:ILE:HG13	2:B:51:SER:CB	0.46	2.41	3	3
2:B:75:GLU:OE1	2:B:95:HIS:NE2	0.46	2.48	12	2
2:B:73:ASN:O	2:B:75:GLU:N	0.46	2.44	14	1
1:A:100:ILE:HD12	1:A:128:LYS:HB2	0.46	1.88	5	2
2:B:17:ARG:HG3	2:B:35:ILE:CG2	0.46	2.41	2	1
1:A:46:LEU:HD12	1:A:49:ASP:HB3	0.46	1.88	12	1
1:A:7:GLN:HB2	1:A:120:TRP:CE3	0.46	2.46	4	1
1:A:49:ASP:HB2	1:A:83:LYS:HE3	0.46	1.87	8	1
2:B:66:LEU:CD1	2:B:110:CYS:HB2	0.46	2.40	1	1
2:B:10:SER:HB3	2:B:106:LYS:CB	0.46	2.41	6	2
1:A:32:LEU:N	1:A:38:SER:OG	0.45	2.48	20	6
1:A:62:THR:HG21	1:A:65:PRO:HB3	0.45	1.86	2	4
2:B:4:PRO:HG3	2:B:36:GLN:NE2	0.45	2.25	8	1
2:B:9:CYS:O	2:B:106:LYS:HD3	0.45	2.11	17	5
2:B:112:ARG:NE	2:B:112:ARG:HA	0.45	2.27	12	4
2:B:100:TRP:CB	2:B:116:THR:HB	0.45	2.41	17	1
1:A:59:HIS:HB3	1:A:62:THR:HG21	0.45	1.88	6	2
1:A:47:LEU:O	1:A:49:ASP:N	0.45	2.49	13	10
2:B:42:GLU:N	2:B:46:GLU:O	0.45	2.48	16	3
1:A:3:LEU:HD11	1:A:29:LEU:HD22	0.45	1.88	19	1
2:B:28:ARG:O	2:B:28:ARG:HG3	0.45	2.12	1	1
1:A:81:GLY:O	1:A:109:ASN:HB2	0.45	2.11	17	4
1:A:67:PHE:CD1	1:A:69:GLU:CG	0.45	3.00	16	3
1:A:87:MET:HB3	1:A:104:PHE:CD2	0.45	2.46	13	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:18:ILE:CD1	2:B:51:SER:CB	0.45	2.94	10	1
2:B:86:HIS:O	2:B:86:HIS:CG	0.45	2.69	16	1
2:B:101:PHE:CZ	2:B:117:HIS:HA	0.45	2.47	4	1
1:A:58:VAL:HG21	1:A:92:PHE:CD2	0.45	2.46	17	1
2:B:6:LEU:HD21	2:B:15:PHE:CD1	0.45	2.46	2	3
2:B:14:HIS:CB	2:B:106:LYS:HG2	0.45	2.42	3	1
2:B:8:TYR:HB2	2:B:106:LYS:HD2	0.45	1.88	10	1
1:A:89:VAL:HG22	1:A:101:ILE:HD11	0.45	1.87	19	5
2:B:5:LYS:HA	2:B:129:PRO:CB	0.45	2.41	8	1
2:B:30:ARG:HD3	2:B:30:ARG:O	0.45	2.12	19	1
1:A:44:VAL:HG22	1:A:54:PHE:O	0.45	2.12	10	1
1:A:101:ILE:O	1:A:125:SER:HA	0.45	2.11	16	1
2:B:3:LYS:O	2:B:4:PRO:C	0.45	2.55	10	14
1:A:42:VAL:HA	1:A:89:VAL:HA	0.45	1.89	6	3
1:A:33:ASP:N	1:A:37:THR:O	0.45	2.46	8	5
1:A:7:GLN:HB2	1:A:24:ILE:HB	0.45	1.87	17	1
1:A:85:LEU:HD23	1:A:85:LEU:C	0.45	2.31	12	2
2:B:79:LEU:O	2:B:90:TYR:HA	0.45	2.12	5	2
1:A:71:PHE:CD1	1:A:71:PHE:C	0.45	2.90	10	3
2:B:14:HIS:O	2:B:106:LYS:HD2	0.45	2.12	15	1
2:B:101:PHE:N	2:B:101:PHE:CD1	0.45	2.84	17	4
1:A:79:GLU:HB3	1:A:83:LYS:HE3	0.45	1.88	11	1
1:A:47:LEU:HG	1:A:48:PRO:N	0.45	2.26	10	6
2:B:58:LEU:HD23	2:B:60:MET:HE2	0.45	1.88	13	1
1:A:10:LEU:HB2	1:A:117:THR:HB	0.45	1.89	9	2
1:A:56:THR:HG22	1:A:69:GLU:CG	0.45	2.42	1	1
2:B:101:PHE:CD1	2:B:123:ILE:HD13	0.45	2.46	14	1
1:A:107:PRO:CG	2:B:27:THR:HA	0.45	2.42	13	2
2:B:102:VAL:O	2:B:102:VAL:HG13	0.45	2.12	12	1
1:A:41:TYR:CB	1:A:58:VAL:HA	0.45	2.41	19	1
1:A:93:ASP:O	1:A:94:ARG:HB2	0.45	2.12	8	8
2:B:56:GLN:CG	2:B:69:SER:HA	0.45	2.42	3	2
1:A:80:LEU:HD23	1:A:80:LEU:C	0.45	2.32	2	1
2:B:56:GLN:OE1	2:B:56:GLN:N	0.45	2.49	4	3
2:B:30:ARG:HA	2:B:35:ILE:HD11	0.45	1.87	14	1
1:A:64:ASN:N	1:A:65:PRO:CD	0.44	2.80	9	4
1:A:103:GLU:CG	1:A:126:ALA:HB2	0.44	2.42	11	1
1:A:45:PHE:CB	1:A:86:VAL:HG23	0.44	2.43	16	3
2:B:18:ILE:HB	2:B:56:GLN:OE1	0.44	2.12	17	2
2:B:85:ASN:O	2:B:86:HIS:CB	0.44	2.64	8	4
1:A:25:GLN:CD	1:A:66:VAL:HG12	0.44	2.33	14	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:4:PRO:HB3	2:B:36:GLN:CD	0.44	2.32	19	1
2:B:7:LEU:HG	2:B:125:PHE:HB3	0.44	1.87	10	2
1:A:47:LEU:HD13	2:B:17:ARG:NH2	0.44	2.27	10	1
2:B:57:TYR:CG	2:B:72:PRO:HB3	0.44	2.47	9	1
1:A:41:TYR:C	1:A:41:TYR:CD1	0.44	2.90	15	1
2:B:74:GLU:CB	2:B:94:LYS:HD3	0.44	2.42	19	1
2:B:10:SER:CB	2:B:106:LYS:HD3	0.44	2.42	10	1
1:A:47:LEU:HD21	1:A:83:LYS:HG3	0.44	1.89	14	1
1:A:124:GLN:OE1	2:B:105:LYS:NZ	0.44	2.46	6	1
1:A:91:ASP:CG	1:A:92:PHE:N	0.44	2.69	2	5
2:B:9:CYS:SG	2:B:10:SER:N	0.44	2.90	9	7
1:A:97:LYS:O	1:A:98:HIS:C	0.44	2.55	19	1
1:A:6:LEU:HD22	1:A:104:PHE:CE2	0.44	2.47	10	1
1:A:19:LEU:HD12	1:A:73:PHE:CD1	0.44	2.48	14	1
2:B:60:MET:SD	2:B:60:MET:N	0.44	2.91	20	3
2:B:109:SER:HB3	2:B:111:LYS:HD3	0.44	1.88	1	2
1:A:52:LYS:NZ	1:A:74:LYS:O	0.44	2.50	15	5
1:A:46:LEU:HD23	1:A:52:LYS:HD2	0.44	1.90	2	1
1:A:42:VAL:HG23	1:A:56:THR:CG2	0.44	2.43	15	1
2:B:58:LEU:HB2	2:B:76:CYS:SG	0.44	2.53	4	1
1:A:45:PHE:HB2	1:A:86:VAL:CG2	0.44	2.43	6	1
2:B:53:GLU:N	2:B:53:GLU:CD	0.44	2.71	7	1
2:B:48:TYR:HB3	2:B:50:LYS:HE3	0.44	1.89	11	1
2:B:109:SER:C	2:B:111:LYS:HD2	0.44	2.33	2	1
2:B:35:ILE:HD13	2:B:35:ILE:N	0.44	2.28	12	1
2:B:3:LYS:HG2	2:B:4:PRO:HD2	0.44	1.89	9	3
2:B:58:LEU:HD23	2:B:60:MET:CE	0.44	2.43	11	3
1:A:8:TYR:HA	1:A:23:ILE:HA	0.44	1.89	17	1
2:B:4:PRO:HB3	2:B:36:GLN:HB3	0.44	1.89	20	1
1:A:59:HIS:HB3	1:A:62:THR:CG2	0.44	2.42	5	1
1:A:42:VAL:HG23	1:A:56:THR:HG21	0.44	1.90	9	2
2:B:32:ASP:O	2:B:34:HIS:N	0.44	2.49	7	2
1:A:104:PHE:C	1:A:104:PHE:CD1	0.43	2.92	17	1
1:A:6:LEU:HD13	1:A:26:ALA:HB2	0.43	1.90	12	2
1:A:91:ASP:O	1:A:98:HIS:HB3	0.43	2.13	2	1
1:A:46:LEU:HD13	1:A:47:LEU:H	0.43	1.72	12	1
2:B:13:GLY:O	2:B:28:ARG:CB	0.43	2.66	18	1
1:A:43:LYS:HE3	1:A:90:TYR:HB3	0.43	1.90	7	1
1:A:4:GLY:HA3	1:A:27:ALA:O	0.43	2.14	11	4
1:A:47:LEU:O	1:A:48:PRO:C	0.43	2.57	4	6
2:B:112:ARG:O	2:B:113:GLY:C	0.43	2.56	20	8

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:3:LEU:HD13	1:A:101:ILE:C	0.43	2.33	3	1
1:A:44:VAL:HG21	1:A:71:PHE:CG	0.43	2.48	13	1
1:A:8:TYR:CE1	1:A:119:GLU:HG2	0.43	2.48	13	1
1:A:3:LEU:CD2	1:A:102:GLY:HA3	0.43	2.41	8	1
2:B:48:TYR:CE1	2:B:77:LEU:CB	0.43	3.00	8	1
1:A:3:LEU:HD11	1:A:101:ILE:HG13	0.43	1.89	4	1
1:A:89:VAL:HG12	1:A:102:GLY:O	0.43	2.14	15	3
2:B:99:ASN:O	2:B:117:HIS:CD2	0.43	2.71	3	1
1:A:16:ASN:OD1	1:A:18:GLN:CG	0.43	2.67	13	1
1:A:107:PRO:HB3	2:B:27:THR:OG1	0.43	2.14	12	1
1:A:79:GLU:HB3	1:A:83:LYS:HE2	0.43	1.91	9	1
1:A:32:LEU:HD12	1:A:93:ASP:HB3	0.43	1.88	15	1
2:B:101:PHE:O	2:B:102:VAL:C	0.43	2.56	17	4
1:A:8:TYR:CE1	1:A:119:GLU:HG3	0.43	2.49	5	3
1:A:44:VAL:HG11	1:A:71:PHE:HB2	0.43	1.90	19	1
2:B:103:GLY:CA	2:B:123:ILE:HG12	0.43	2.44	6	1
2:B:83:GLU:HG2	2:B:85:ASN:CB	0.43	2.43	2	1
2:B:63:ASP:O	2:B:64:GLY:C	0.43	2.57	10	2
2:B:6:LEU:HD23	2:B:6:LEU:C	0.43	2.33	14	1
1:A:100:ILE:HD12	1:A:128:LYS:HB3	0.43	1.91	4	1
1:A:87:MET:HB3	1:A:104:PHE:CE2	0.43	2.48	4	5
1:A:104:PHE:CD1	1:A:121:ARG:HG2	0.43	2.48	17	1
1:A:19:LEU:CB	1:A:73:PHE:HB3	0.43	2.44	20	3
2:B:10:SER:HB2	2:B:106:LYS:HB3	0.43	1.89	3	1
2:B:17:ARG:HA	2:B:35:ILE:HG22	0.43	1.91	10	3
1:A:25:GLN:HB2	1:A:65:PRO:O	0.43	2.14	14	2
2:B:10:SER:O	2:B:106:LYS:HD3	0.43	2.14	16	1
2:B:113:GLY:HA2	2:B:116:THR:OG1	0.43	2.14	7	1
2:B:17:ARG:CG	2:B:34:HIS:HB3	0.43	2.44	7	1
1:A:3:LEU:CB	1:A:123:LEU:HB3	0.43	2.44	5	3
2:B:121:LYS:HD2	2:B:124:LEU:HD23	0.43	1.91	8	2
1:A:1:GLU:HG2	1:A:124:GLN:HB2	0.43	1.90	19	1
2:B:9:CYS:C	2:B:106:LYS:HD3	0.43	2.34	10	1
1:A:7:GLN:NE2	1:A:25:GLN:HG3	0.43	2.29	9	1
2:B:68:GLY:O	2:B:69:SER:CB	0.43	2.67	9	1
2:B:112:ARG:O	2:B:115:ARG:N	0.43	2.43	9	1
1:A:87:MET:HB3	1:A:104:PHE:CZ	0.43	2.48	9	1
2:B:66:LEU:HB2	2:B:110:CYS:CB	0.43	2.43	17	1
1:A:39:ASP:CB	1:A:92:PHE:HB3	0.43	2.44	17	1
2:B:37:LEU:HA	2:B:51:SER:HA	0.43	1.90	14	1
2:B:57:TYR:O	2:B:69:SER:N	0.43	2.52	6	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:17:ARG:O	2:B:24:VAL:HA	0.43	2.14	8	1
1:A:79:GLU:HB3	1:A:83:LYS:CE	0.43	2.44	9	1
2:B:54:THR:HG22	2:B:55:GLY:N	0.43	2.29	15	1
2:B:4:PRO:HD3	2:B:53:GLU:HG2	0.42	1.91	11	2
2:B:14:HIS:HB2	2:B:106:LYS:CD	0.42	2.42	17	3
1:A:79:GLU:HG2	1:A:83:LYS:HE2	0.42	1.91	4	3
2:B:28:ARG:O	2:B:28:ARG:NE	0.42	2.47	20	2
1:A:82:GLY:C	1:A:83:LYS:HD2	0.42	2.34	2	1
1:A:7:GLN:O	1:A:24:ILE:N	0.42	2.51	19	2
1:A:41:TYR:CD1	1:A:41:TYR:C	0.42	2.93	4	2
1:A:18:GLN:HB2	1:A:73:PHE:O	0.42	2.14	5	1
2:B:60:MET:HE2	2:B:102:VAL:CG2	0.42	2.45	12	1
1:A:39:ASP:HA	1:A:59:HIS:O	0.42	2.14	10	1
2:B:59:ALA:HB3	2:B:68:GLY:HA2	0.42	1.91	16	1
2:B:17:ARG:HD2	2:B:27:THR:CB	0.42	2.44	11	1
2:B:14:HIS:CB	2:B:106:LYS:CG	0.42	2.98	18	2
2:B:73:ASN:OD1	2:B:74:GLU:N	0.42	2.49	12	1
1:A:13:ASP:OD1	1:A:15:GLN:HG2	0.42	2.14	19	1
1:A:103:GLU:OE2	1:A:105:LYS:NZ	0.42	2.51	11	2
2:B:110:CYS:O	2:B:111:LYS:NZ	0.42	2.51	2	1
1:A:87:MET:HB3	1:A:104:PHE:CE1	0.42	2.49	12	1
1:A:126:ALA:HB1	1:A:128:LYS:HE3	0.42	1.91	10	1
2:B:6:LEU:HB3	2:B:128:LEU:HB2	0.42	1.91	4	1
1:A:85:LEU:HD21	1:A:106:VAL:CG1	0.42	2.45	18	3
2:B:1:TYR:N	2:B:132:SER:OG	0.42	2.50	5	1
1:A:39:ASP:HB2	1:A:92:PHE:C	0.42	2.35	5	1
2:B:55:GLY:C	2:B:56:GLN:HG3	0.42	2.35	5	1
2:B:78:PHE:HB3	2:B:90:TYR:HB3	0.42	1.90	8	1
1:A:49:ASP:HB2	1:A:83:LYS:HG3	0.42	1.92	1	1
1:A:19:LEU:CD1	1:A:80:LEU:HD12	0.42	2.44	16	1
1:A:64:ASN:O	1:A:64:ASN:ND2	0.42	2.53	17	1
2:B:66:LEU:HB2	2:B:110:CYS:O	0.42	2.14	2	1
1:A:110:THR:HG23	2:B:28:ARG:CZ	0.42	2.44	8	1
1:A:46:LEU:HD22	1:A:85:LEU:HA	0.42	1.92	1	1
1:A:29:LEU:HD13	1:A:40:PRO:CB	0.42	2.45	18	1
2:B:61:ASP:HB2	2:B:67:TYR:CE2	0.42	2.50	6	1
2:B:20:PRO:HA	2:B:56:GLN:HG2	0.42	1.92	11	2
2:B:78:PHE:CD1	2:B:92:SER:HA	0.42	2.49	20	1
2:B:2:LYS:CG	2:B:2:LYS:O	0.42	2.67	2	2
1:A:23:ILE:HG12	1:A:23:ILE:O	0.42	2.14	10	2
1:A:74:LYS:O	1:A:75:VAL:HG23	0.42	2.15	15	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:25:GLN:HG2	1:A:66:VAL:HG12	0.42	1.91	18	1
1:A:42:VAL:N	1:A:56:THR:OG1	0.42	2.52	4	1
1:A:19:LEU:HD13	1:A:108:MET:HE2	0.42	1.92	6	1
2:B:51:SER:HB3	2:B:55:GLY:HA3	0.42	1.90	20	1
2:B:103:GLY:N	2:B:116:THR:HG23	0.42	2.29	13	1
2:B:10:SER:O	2:B:106:LYS:NZ	0.42	2.49	12	2
1:A:110:THR:CG2	2:B:28:ARG:HB3	0.42	2.45	14	1
2:B:6:LEU:C	2:B:6:LEU:HD23	0.42	2.35	7	2
2:B:17:ARG:CD	2:B:27:THR:HB	0.42	2.45	11	1
1:A:31:ALA:HA	1:A:38:SER:OG	0.42	2.14	17	2
1:A:88:ALA:HA	1:A:103:GLU:HA	0.42	1.92	17	2
2:B:17:ARG:CA	2:B:35:ILE:HG22	0.42	2.45	14	2
2:B:42:GLU:HB2	2:B:48:TYR:CE1	0.42	2.50	5	1
1:A:1:GLU:OE2	2:B:105:LYS:NZ	0.42	2.53	2	1
2:B:105:LYS:HB2	2:B:109:SER:O	0.42	2.15	2	1
2:B:75:GLU:HA	2:B:94:LYS:HB2	0.42	1.92	8	1
1:A:1:GLU:HB2	1:A:124:GLN:CG	0.42	2.44	15	1
2:B:60:MET:CG	2:B:66:LEU:CD2	0.41	2.97	15	3
2:B:75:GLU:OE2	2:B:95:HIS:NE2	0.41	2.50	7	1
1:A:11:ASP:HB3	1:A:20:LEU:HB3	0.41	1.91	7	1
1:A:106:VAL:HG23	1:A:107:PRO:HD2	0.41	1.91	17	1
1:A:44:VAL:HG21	1:A:71:PHE:CZ	0.41	2.50	5	1
2:B:17:ARG:HG2	2:B:34:HIS:HB3	0.41	1.91	2	1
2:B:45:GLY:O	2:B:79:LEU:HA	0.41	2.15	8	1
1:A:110:THR:HG21	2:B:28:ARG:HB3	0.41	1.92	14	1
1:A:64:ASN:ND2	1:A:64:ASN:O	0.41	2.53	6	1
1:A:47:LEU:HD22	1:A:84:THR:N	0.41	2.30	11	1
1:A:11:ASP:O	1:A:20:LEU:N	0.41	2.52	20	2
2:B:35:ILE:N	2:B:35:ILE:HD13	0.41	2.30	18	1
2:B:39:LEU:HD13	2:B:49:ILE:HG13	0.41	1.91	15	1
1:A:20:LEU:HD22	1:A:70:GLN:NE2	0.41	2.30	15	1
1:A:102:GLY:HA2	1:A:126:ALA:N	0.41	2.29	7	2
1:A:2:LYS:HD3	1:A:2:LYS:O	0.41	2.14	10	4
1:A:47:LEU:HD22	1:A:84:THR:H	0.41	1.75	17	1
2:B:14:HIS:CB	2:B:106:LYS:HD2	0.41	2.43	17	2
2:B:58:LEU:O	2:B:76:CYS:SG	0.41	2.77	17	2
2:B:5:LYS:HG2	2:B:129:PRO:CB	0.41	2.45	13	1
1:A:47:LEU:HD12	2:B:19:LEU:CD2	0.41	2.45	9	2
1:A:3:LEU:HD12	1:A:30:PRO:CG	0.41	2.45	9	1
2:B:18:ILE:N	2:B:34:HIS:O	0.41	2.48	11	1
1:A:27:ALA:O	1:A:28:GLU:HB2	0.41	2.16	19	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:86:HIS:CE1	2:B:126:LEU:HD11	0.41	2.50	13	1
1:A:121:ARG:HD3	1:A:122:ASP:N	0.41	2.30	12	1
1:A:23:ILE:CD1	1:A:69:GLU:HB3	0.41	2.43	19	1
1:A:47:LEU:CD2	1:A:83:LYS:HG3	0.41	2.45	9	1
1:A:105:LYS:HD2	1:A:105:LYS:N	0.41	2.31	15	1
2:B:105:LYS:NZ	2:B:107:ASN:OD1	0.41	2.51	6	1
2:B:99:ASN:HB3	2:B:117:HIS:CE1	0.41	2.51	10	1
1:A:62:THR:O	1:A:62:THR:HG23	0.41	2.15	4	1
1:A:39:ASP:HB3	1:A:92:PHE:HB3	0.41	1.93	17	1
1:A:85:LEU:O	1:A:105:LYS:HA	0.41	2.16	20	1
2:B:74:GLU:HA	2:B:77:LEU:HD23	0.41	1.93	3	1
2:B:83:GLU:HB2	2:B:124:LEU:HD12	0.41	1.92	3	1
2:B:30:ARG:O	2:B:30:ARG:HD2	0.41	2.15	5	1
1:A:41:TYR:CE1	1:A:43:LYS:HE3	0.41	2.51	8	1
2:B:7:LEU:HD11	2:B:125:PHE:HB3	0.41	1.93	19	1
1:A:20:LEU:HD22	1:A:70:GLN:CD	0.41	2.36	10	1
2:B:14:HIS:HB2	2:B:106:LYS:CG	0.41	2.46	18	1
2:B:76:CYS:O	2:B:76:CYS:SG	0.41	2.78	15	1
1:A:88:ALA:CA	1:A:103:GLU:HB3	0.41	2.46	14	1
1:A:20:LEU:HD21	1:A:70:GLN:HG3	0.41	1.92	14	1
1:A:103:GLU:OE1	1:A:105:LYS:NZ	0.41	2.50	7	1
2:B:104:LEU:HB3	2:B:110:CYS:HB3	0.41	1.92	3	2
1:A:84:THR:HG21	2:B:25:ASP:CB	0.41	2.46	3	1
1:A:16:ASN:OD1	1:A:17:ASN:N	0.41	2.53	10	1
2:B:105:LYS:CG	2:B:112:ARG:HD2	0.41	2.45	1	1
1:A:38:SER:HB3	1:A:63:LEU:CD2	0.41	2.46	9	1
1:A:85:LEU:HD23	1:A:106:VAL:O	0.41	2.16	18	1
2:B:104:LEU:CB	2:B:110:CYS:HB3	0.41	2.45	14	1
1:A:16:ASN:O	1:A:17:ASN:HB2	0.41	2.15	14	1
1:A:104:PHE:O	1:A:105:LYS:NZ	0.41	2.54	7	1
2:B:61:ASP:HB3	2:B:65:LEU:CB	0.41	2.45	5	1
2:B:115:ARG:HD2	2:B:115:ARG:N	0.41	2.31	19	2
1:A:58:VAL:HG21	1:A:92:PHE:HB3	0.41	1.92	18	1
1:A:13:ASP:HB3	1:A:16:ASN:OD1	0.41	2.16	7	1
1:A:44:VAL:HG11	1:A:71:PHE:CB	0.41	2.45	20	1
1:A:60:ARG:O	1:A:60:ARG:CD	0.41	2.69	20	1
1:A:9:SER:HB3	1:A:118:GLU:HG2	0.41	1.93	20	1
1:A:30:PRO:HB2	1:A:101:ILE:HG22	0.41	1.91	5	1
2:B:98:LYS:HD3	2:B:100:TRP:CZ3	0.41	2.51	12	1
2:B:63:ASP:HB2	2:B:95:HIS:CG	0.41	2.51	14	1
2:B:101:PHE:CE2	2:B:123:ILE:HD13	0.41	2.51	4	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:14:HIS:CD2	2:B:106:LYS:HG2	0.41	2.51	6	1
1:A:110:THR:HG21	2:B:28:ARG:NE	0.41	2.31	6	1
1:A:105:LYS:N	1:A:105:LYS:HD2	0.41	2.30	11	1
2:B:14:HIS:ND1	2:B:26:GLY:O	0.41	2.41	17	1
2:B:17:ARG:HD3	2:B:27:THR:OG1	0.41	2.16	2	1
2:B:103:GLY:N	2:B:116:THR:HG22	0.41	2.31	8	1
1:A:80:LEU:HD12	1:A:108:MET:CE	0.41	2.46	10	1
2:B:74:GLU:HB2	2:B:77:LEU:HD23	0.41	1.93	15	1
2:B:99:ASN:HB3	2:B:117:HIS:CD2	0.41	2.50	14	1
2:B:65:LEU:N	2:B:65:LEU:HD12	0.40	2.30	7	1
2:B:130:VAL:HG23	2:B:130:VAL:O	0.40	2.16	11	1
2:B:8:TYR:CD1	2:B:9:CYS:N	0.40	2.89	20	1
2:B:64:GLY:CA	2:B:113:GLY:HA3	0.40	2.43	16	2
1:A:112:ASP:O	1:A:114:GLY:N	0.40	2.50	16	1
2:B:101:PHE:CD1	2:B:101:PHE:N	0.40	2.89	6	1
1:A:106:VAL:CG2	1:A:107:PRO:HD2	0.40	2.46	17	1
1:A:6:LEU:C	1:A:6:LEU:HD23	0.40	2.37	19	1
1:A:89:VAL:CG1	1:A:101:ILE:HG12	0.40	2.46	15	1
2:B:9:CYS:SG	2:B:122:ALA:O	0.40	2.79	7	2
1:A:25:GLN:CG	1:A:66:VAL:HG12	0.40	2.45	11	1
1:A:76:PRO:O	1:A:80:LEU:N	0.40	2.48	20	1
2:B:16:LEU:HD13	2:B:104:LEU:HD22	0.40	1.94	3	1
2:B:51:SER:HB2	2:B:55:GLY:CA	0.40	2.46	12	1
1:A:10:LEU:HD22	1:A:21:VAL:CG1	0.40	2.45	17	1
2:B:101:PHE:C	2:B:116:THR:HG22	0.40	2.37	17	1
2:B:91:ILE:O	2:B:93:LYS:N	0.40	2.54	17	1
2:B:49:ILE:CD1	2:B:49:ILE:N	0.40	2.84	19	1
1:A:8:TYR:CD1	1:A:8:TYR:N	0.40	2.89	17	1
2:B:8:TYR:HB3	2:B:15:PHE:HA	0.40	1.94	2	1
1:A:3:LEU:HD21	1:A:89:VAL:HG12	0.40	1.93	12	1
1:A:12:TYR:CD1	1:A:113:PHE:HB2	0.40	2.51	12	1
1:A:106:VAL:HG22	1:A:111:VAL:HG21	0.40	1.92	12	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	126/128 (98%)	96±3 (76±3%)	24±4 (19±3%)	6±1 (5±1%)	5	28
2	B	131/133 (98%)	91±4 (70±3%)	29±4 (22±3%)	10±2 (8±1%)	2	15
All	All	5140/5220 (98%)	3755 (73%)	1062 (21%)	323 (6%)	3	20

All 46 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
2	B	4	PRO	20
2	B	102	VAL	20
2	B	69	SER	20
2	B	111	LYS	20
1	A	47	LEU	20
1	A	48	PRO	19
2	B	44	VAL	19
1	A	52	LYS	18
2	B	54	THR	16
1	A	35	GLY	16
1	A	67	PHE	13
2	B	108	GLY	13
2	B	109	SER	10
2	B	52	THR	8
2	B	14	HIS	8
1	A	74	LYS	8
1	A	61	LYS	7
2	B	86	HIS	6
2	B	10	SER	6
2	B	13	GLY	6
2	B	11	ASN	5
2	B	61	ASP	4
2	B	113	GLY	4
1	A	28	GLU	3
1	A	98	HIS	3
2	B	119	GLY	3
2	B	2	LYS	3
2	B	70	GLN	2
1	A	95	PHE	2
2	B	74	GLU	2
1	A	58	VAL	2
2	B	60	MET	2
1	A	75	VAL	2
2	B	118	TYR	1
1	A	50	LYS	1

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Mol	Chain	Res	Type	Models (Total)
1	A	112	ASP	1
1	A	127	GLU	1
1	A	115	HIS	1
2	B	131	SER	1
2	B	55	GLY	1
1	A	113	PHE	1
2	B	3	LYS	1
2	B	33	GLN	1
1	A	94	ARG	1
2	B	92	SER	1
2	B	39	LEU	1

6.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the sidechain conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	115/115 (100%)	93±3 (80±2%)	22±3 (20±2%)	5	37
2	B	117/117 (100%)	98±2 (84±2%)	19±2 (16±2%)	6	43
All	All	4640/4640 (100%)	3808 (82%)	832 (18%)	5	40

All 107 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)
2	B	2	LYS	20
1	A	8	TYR	20
2	B	35	ILE	20
2	B	111	LYS	20
2	B	56	GLN	20
2	B	38	GLN	20
1	A	23	ILE	20
1	A	2	LYS	20
1	A	25	GLN	19
1	A	101	ILE	19
1	A	105	LYS	18
1	A	20	LEU	18
1	A	80	LEU	18

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Mol	Chain	Res	Type	Models (Total)
1	A	125	SER	18
2	B	30	ARG	17
2	B	18	ILE	16
2	B	43	SER	16
1	A	78	SER	15
1	A	37	THR	15
1	A	18	GLN	15
2	B	89	THR	15
1	A	7	GLN	14
1	A	75	VAL	13
2	B	60	MET	13
1	A	96	SER	12
1	A	84	THR	12
2	B	74	GLU	12
2	B	21	ASP	12
1	A	60	ARG	12
1	A	32	LEU	12
2	B	92	SER	11
2	B	52	THR	11
2	B	97	GLU	10
1	A	71	PHE	10
2	B	117	HIS	10
1	A	46	LEU	10
1	A	98	HIS	10
1	A	70	GLN	9
2	B	99	ASN	9
1	A	112	ASP	9
2	B	101	PHE	9
2	B	71	THR	9
1	A	10	LEU	8
2	B	28	ARG	8
2	B	14	HIS	8
1	A	38	SER	8
2	B	37	LEU	8
2	B	8	TYR	8
1	A	9	SER	8
2	B	17	ARG	8
1	A	119	GLU	7
1	A	67	PHE	7
1	A	83	LYS	7
2	B	120	GLN	7
1	A	47	LEU	7

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Mol	Chain	Res	Type	Models (Total)
1	A	103	GLU	6
1	A	85	LEU	6
1	A	15	GLN	5
2	B	116	THR	5
2	B	36	GLN	5
1	A	43	LYS	4
2	B	63	ASP	4
2	B	51	SER	4
2	B	77	LEU	4
2	B	131	SER	4
2	B	105	LYS	3
2	B	62	THR	3
2	B	48	TYR	3
1	A	68	ASN	3
2	B	24	VAL	3
1	A	91	ASP	3
2	B	34	HIS	3
1	A	11	ASP	3
1	A	72	THR	3
1	A	41	TYR	3
2	B	112	ARG	2
1	A	104	PHE	2
2	B	54	THR	2
1	A	90	TYR	2
2	B	10	SER	2
1	A	69	GLU	2
1	A	94	ARG	2
2	B	85	ASN	2
2	B	115	ARG	2
2	B	50	LYS	2
2	B	106	LYS	2
1	A	73	PHE	2
1	A	110	THR	2
1	A	64	ASN	2
2	B	25	ASP	2
1	A	79	GLU	2
2	B	53	GLU	1
2	B	23	THR	1
1	A	51	LYS	1
1	A	122	ASP	1
2	B	1	TYR	1
2	B	42	GLU	1

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Mol	Chain	Res	Type	Models (Total)
2	B	130	VAL	1
1	A	52	LYS	1
1	A	92	PHE	1
2	B	75	GLU	1
2	B	27	THR	1
1	A	1	GLU	1
2	B	95	HIS	1
1	A	50	LYS	1
1	A	99	ASP	1
2	B	31	SER	1

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided