



# Full wwPDB NMR Structure Validation Report ⓘ

Feb 16, 2022 – 07:25 AM EST

PDB ID : 1MGS  
Title : THE SOLUTION STRUCTURE OF MELANOMA GROWTH STIMULATING ACTIVITY  
Authors : Fairbrother, W.J.  
Deposited on : 1994-07-19

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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

MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
RCI : v\_1n\_11\_5\_13\_A (Berjanski et al., 2005)  
PANAV : Wang et al. (2010)  
ShiftChecker : 2.26  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.26

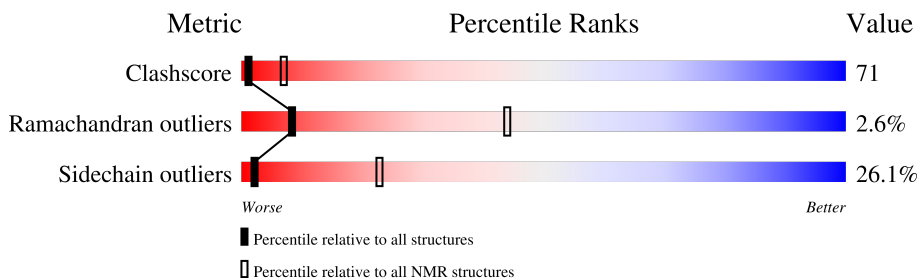
# 1 Overall quality at a glance

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	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

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  $\geq 3$ , 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	73	
1	B	73	

## 2 Ensemble composition and analysis i

This entry contains 25 models. Model 3 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:9-A:69, B:9-B:69 (122)	0.30	3

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 3 clusters and 1 single-model cluster was found.

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

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2250 atoms, of which 1156 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY.

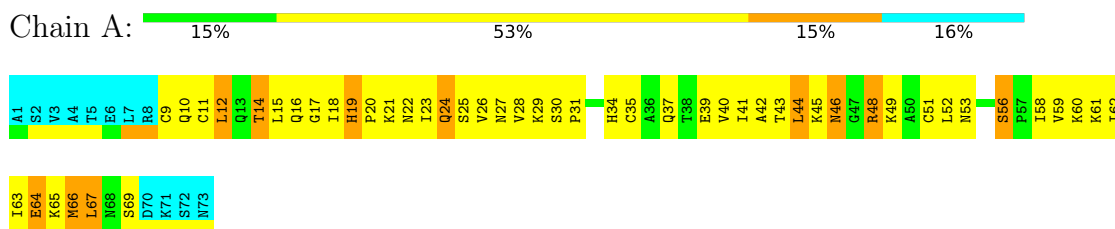
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	73	1125	337	578	102	103	5	0
1	B	73	1125	337	578	102	103	5	0

## 4 Residue-property plots [i](#)

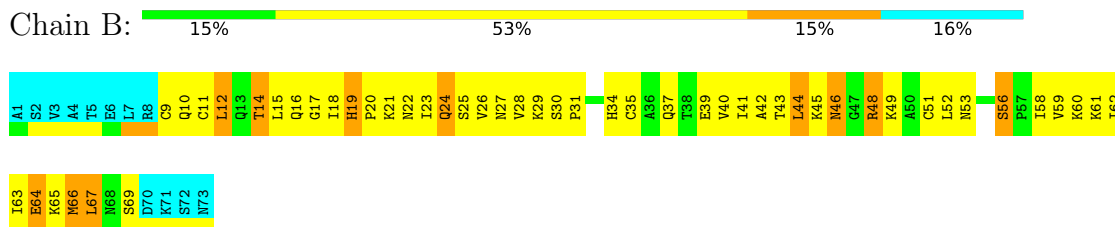
### 4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY



- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

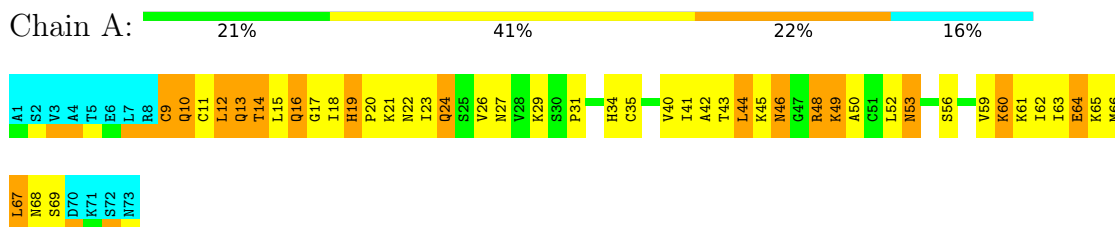


### 4.2 Scores per residue for each member of the ensemble

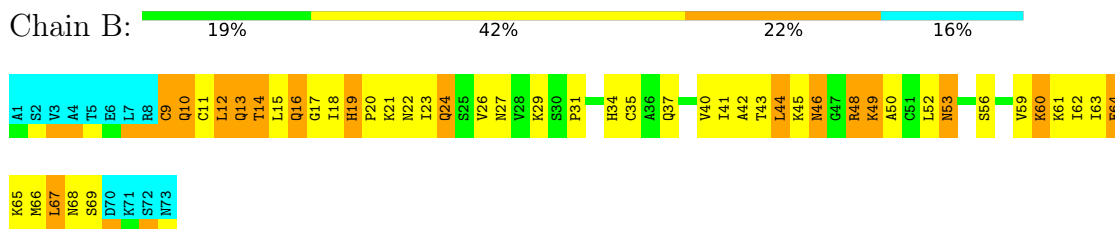
Colouring as in section [4.1](#) above.

#### 4.2.1 Score per residue for model 1

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

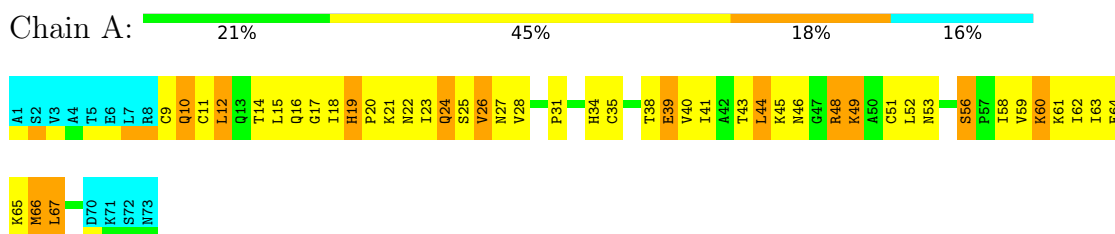


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

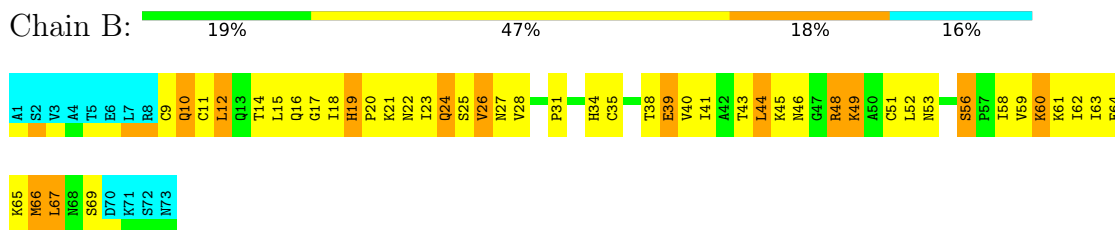


#### 4.2.2 Score per residue for model 2

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

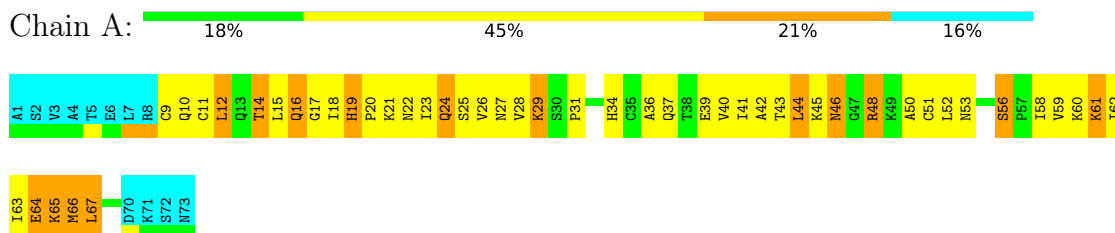


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

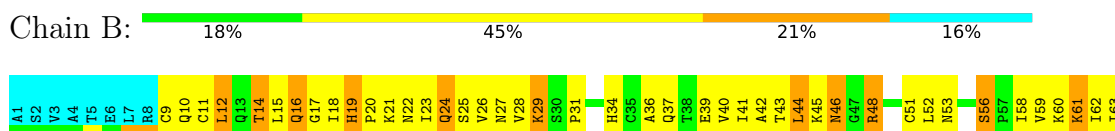


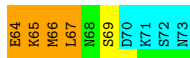
#### 4.2.3 Score per residue for model 3 (medoid)

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY



- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

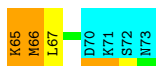




#### 4.2.4 Score per residue for model 4

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

Chain A: 22% 42% 18% 16%



- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

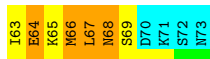
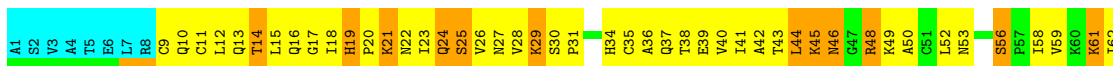
Chain B: 22% 42% 18% 16%



#### 4.2.5 Score per residue for model 5

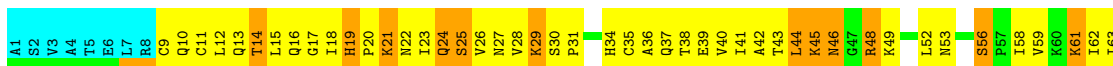
- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

Chain A: 11% 51% 22% 16%



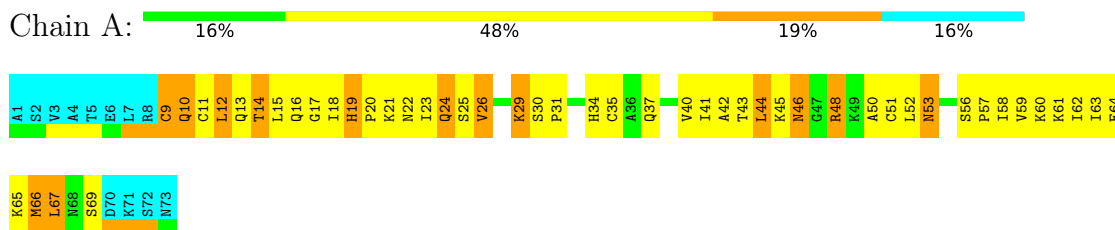
- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

Chain B: 12% 49% 22% 16%

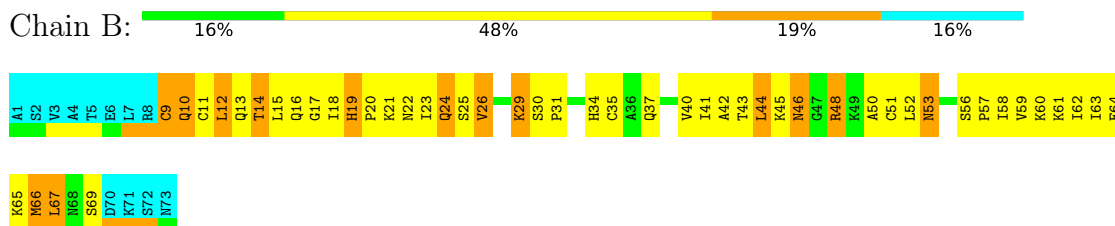


### 4.2.6 Score per residue for model 6

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

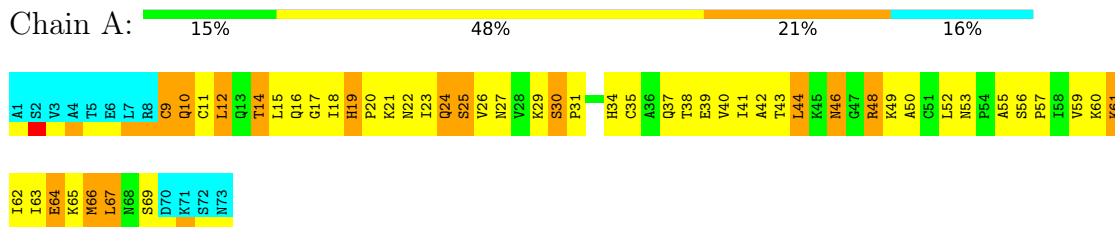


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

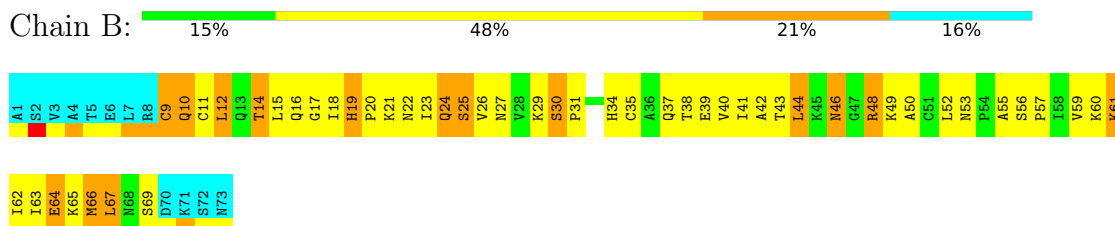


### 4.2.7 Score per residue for model 7

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

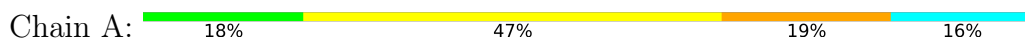


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

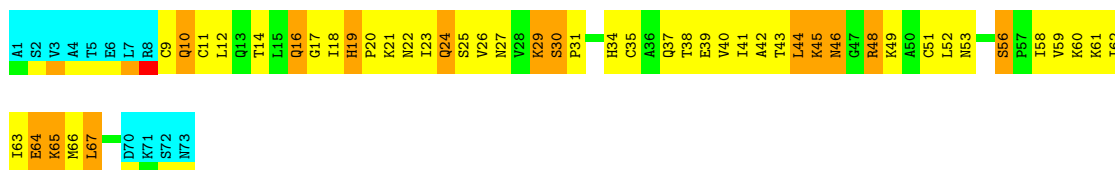


### 4.2.8 Score per residue for model 8

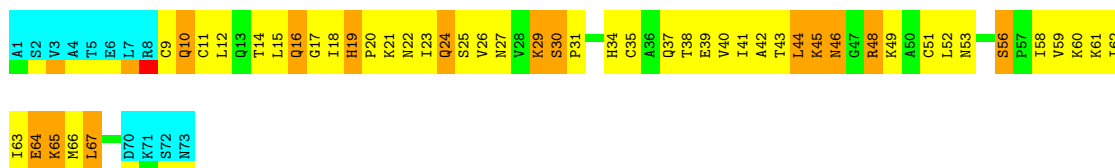
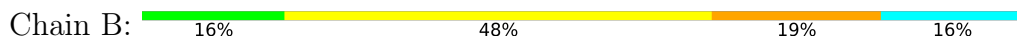
- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY





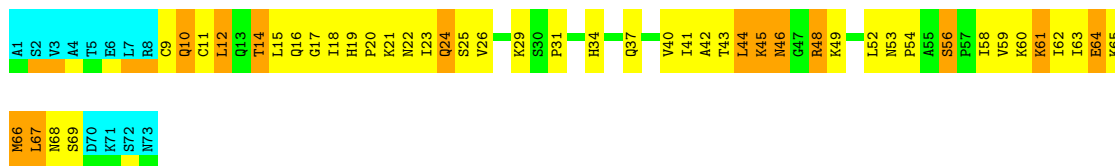
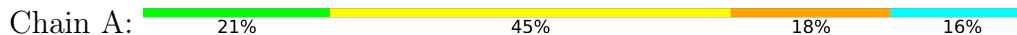


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

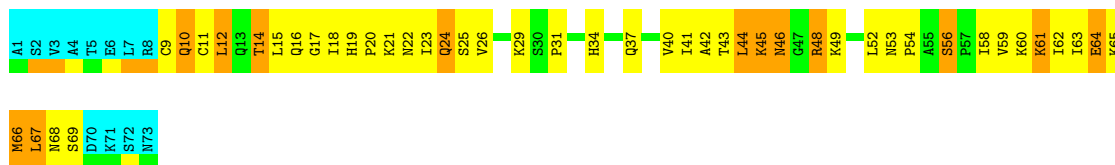
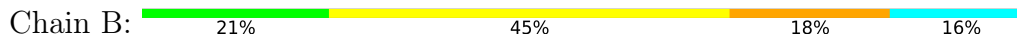


#### 4.2.9 Score per residue for model 9

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

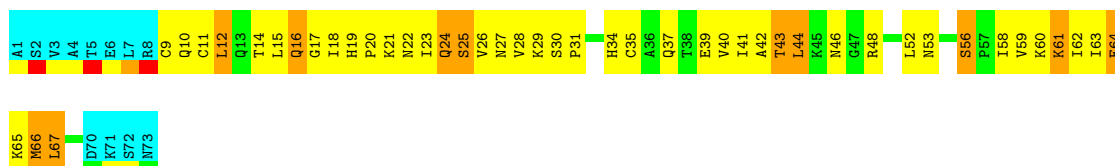


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

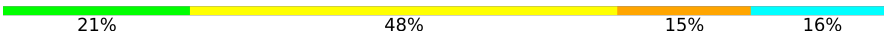


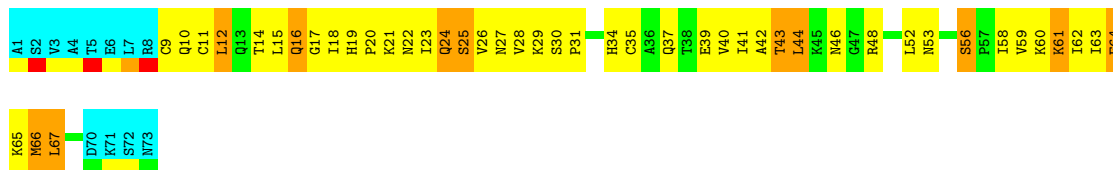
#### 4.2.10 Score per residue for model 10

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY




- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

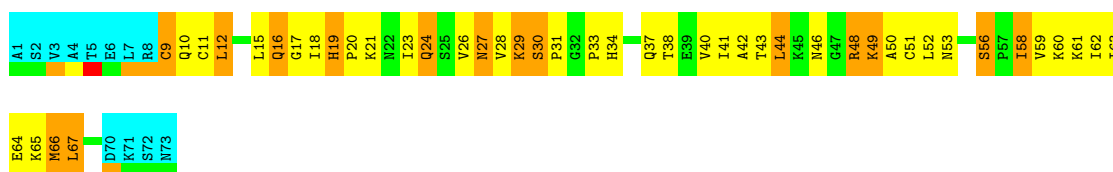
Chain B: 



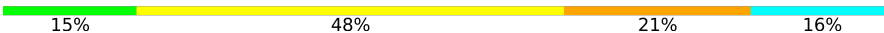
#### 4.2.11 Score per residue for model 11

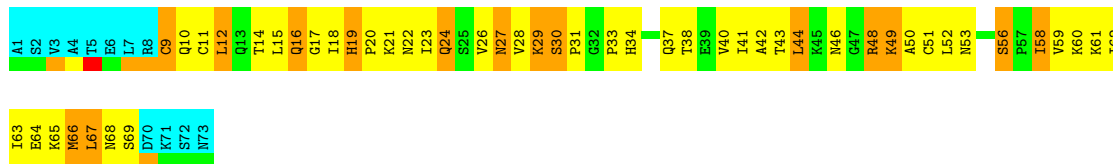
- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

Chain A: 



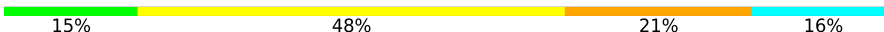
- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

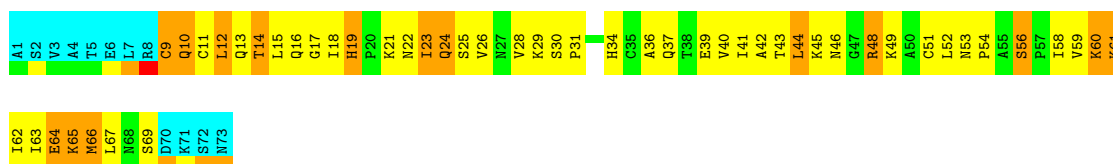
Chain B: 



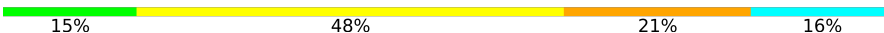
#### 4.2.12 Score per residue for model 12

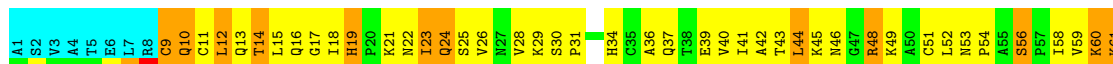
- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

Chain A: 



- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

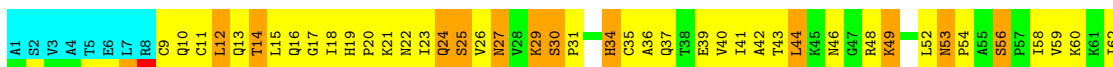
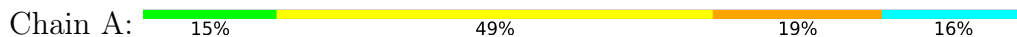
Chain B: 



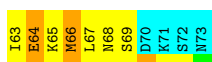
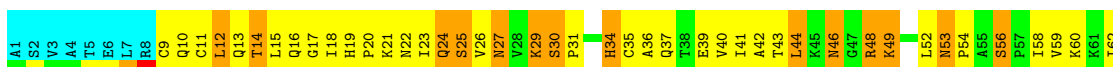
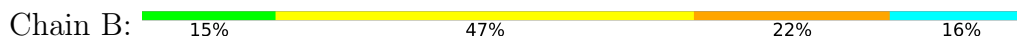


#### 4.2.13 Score per residue for model 13

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

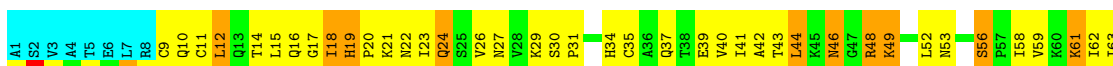


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

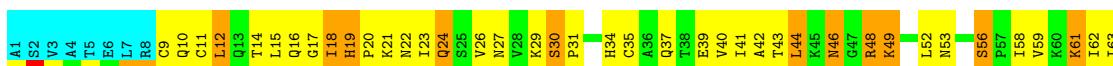
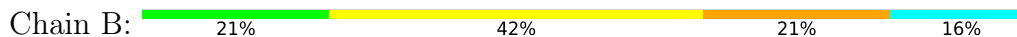


#### 4.2.14 Score per residue for model 14

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

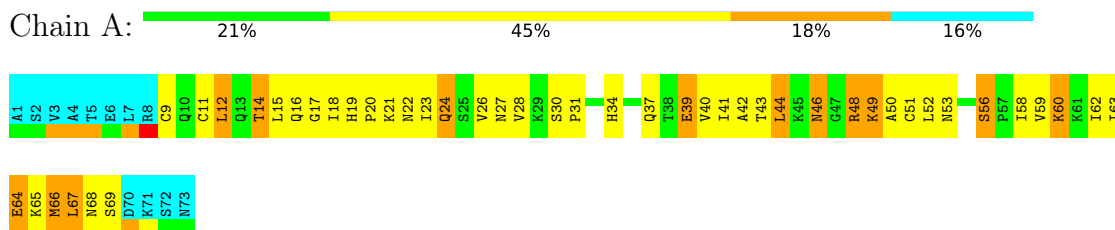


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

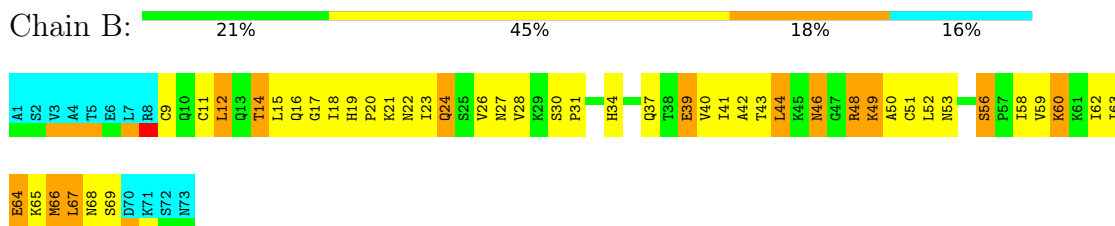


### 4.2.15 Score per residue for model 15

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

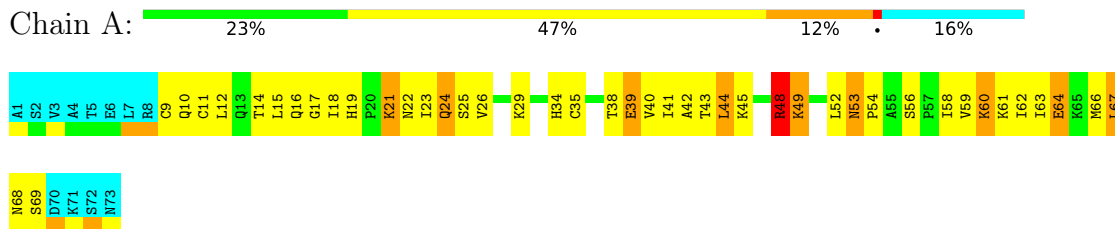


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

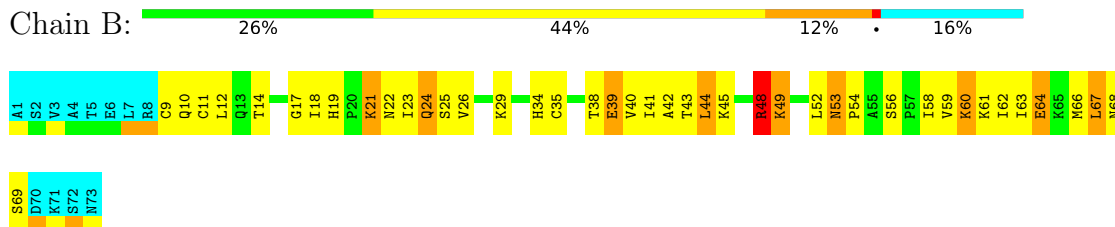


### 4.2.16 Score per residue for model 16

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

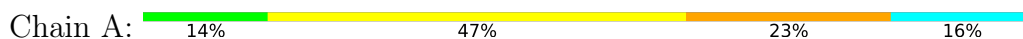


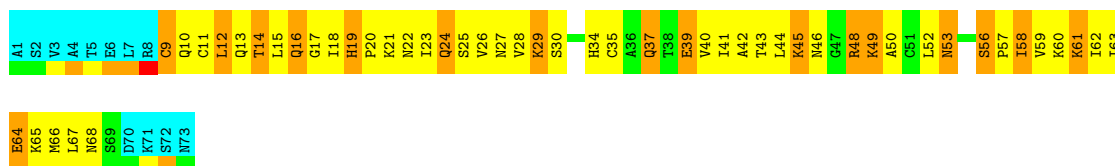
- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY



### 4.2.17 Score per residue for model 17

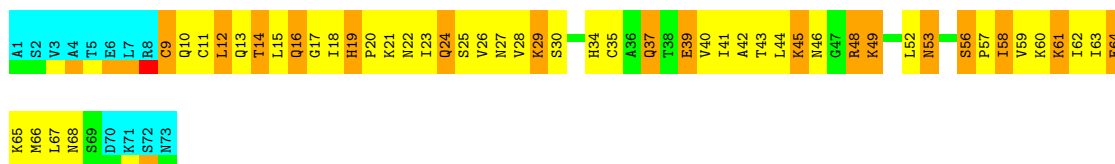
- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY





- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

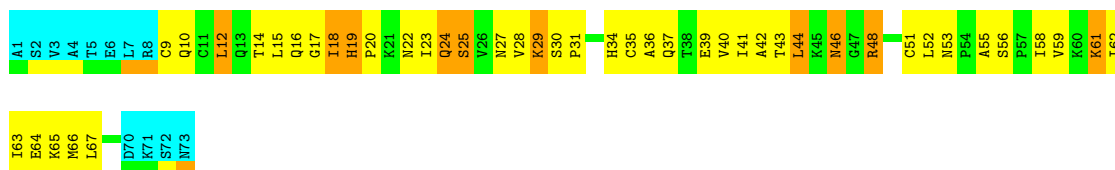
Chain B: 15% 45% 23% 16%



#### 4.2.18 Score per residue for model 18

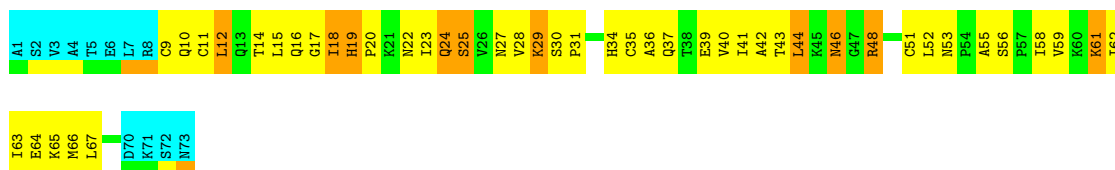
- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

Chain A: 22% 48% 14% 16%



- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

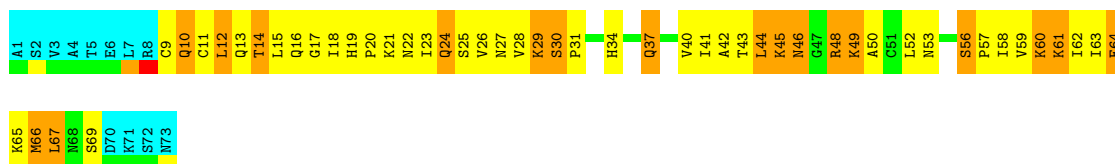
Chain B: 21% 49% 14% 16%



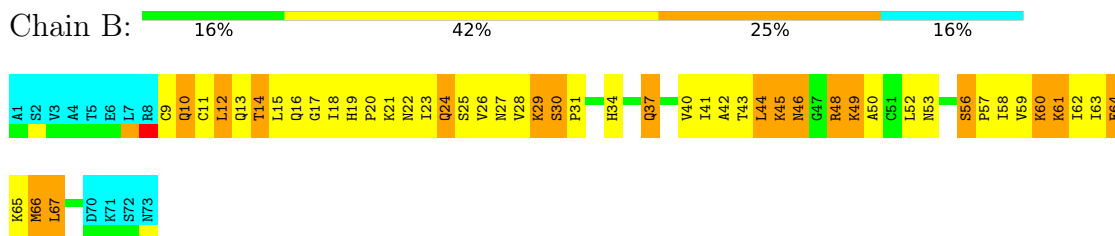
#### 4.2.19 Score per residue for model 19

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

Chain A: 15% 44% 25% 16%

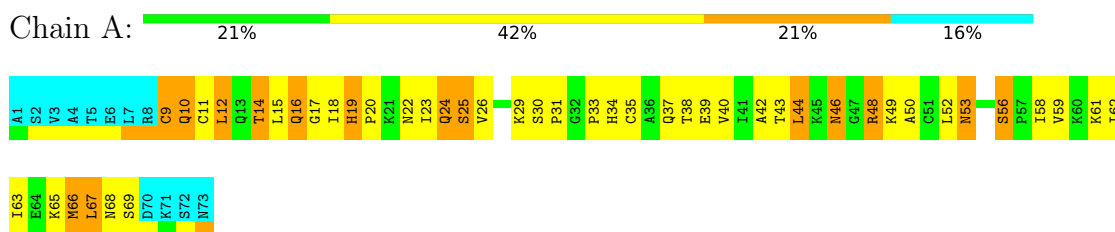


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

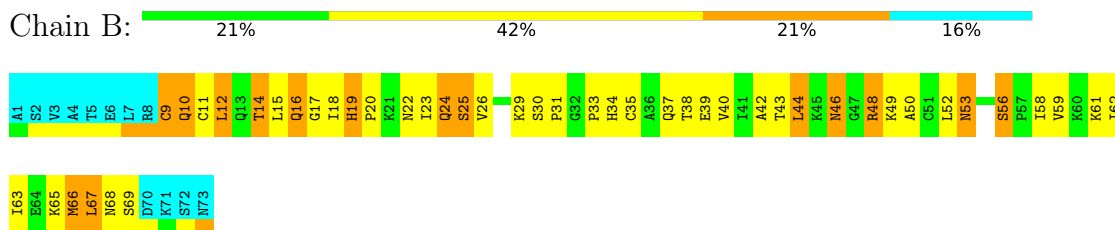


#### 4.2.20 Score per residue for model 20

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

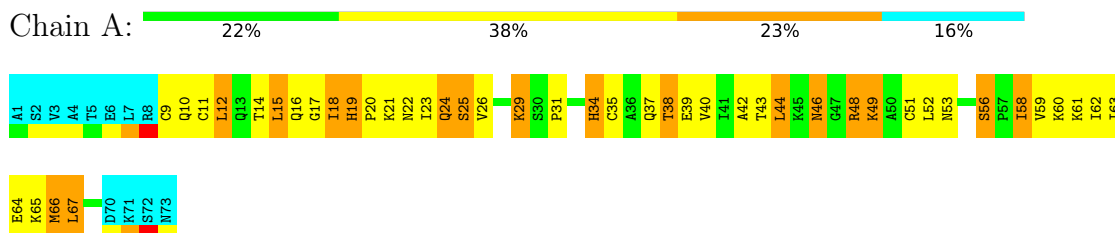


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

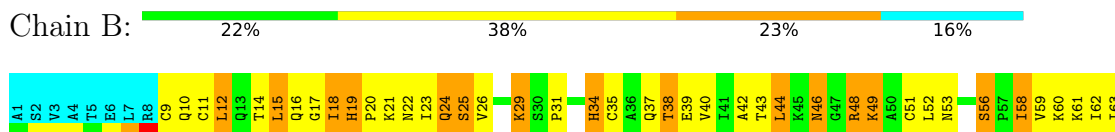


#### 4.2.21 Score per residue for model 21

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY



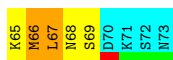
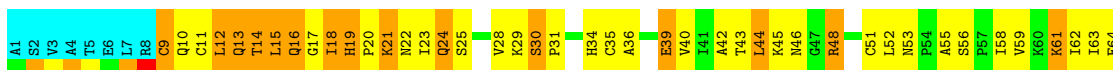
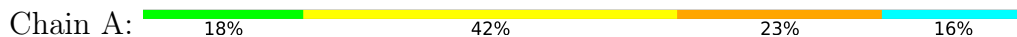
- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY



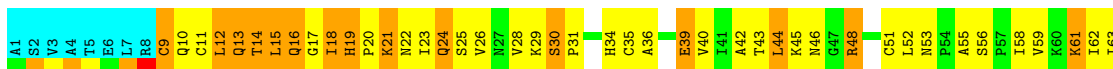
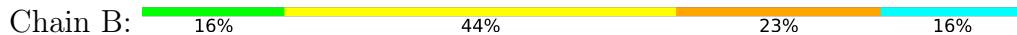


#### 4.2.22 Score per residue for model 22

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

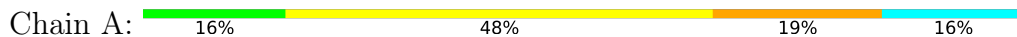


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

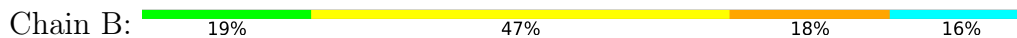


#### 4.2.23 Score per residue for model 23

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

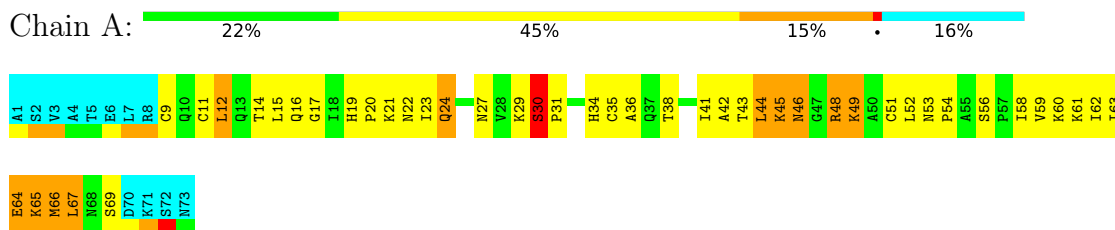


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

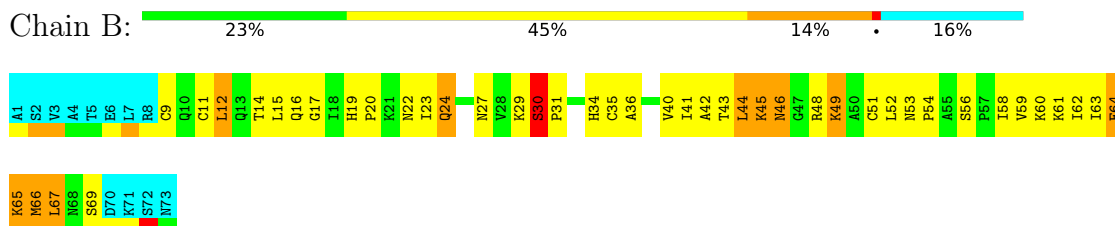


#### 4.2.24 Score per residue for model 24

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

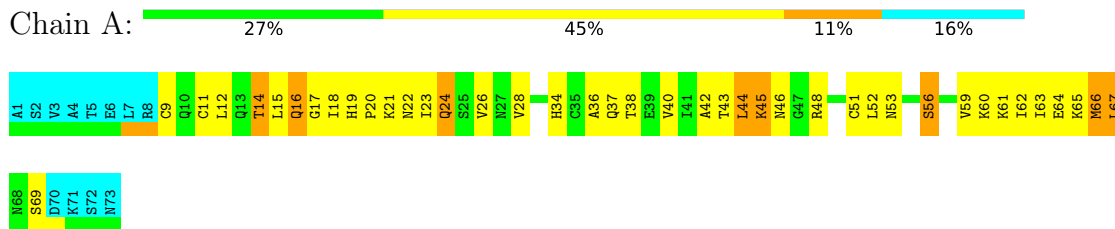


- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY

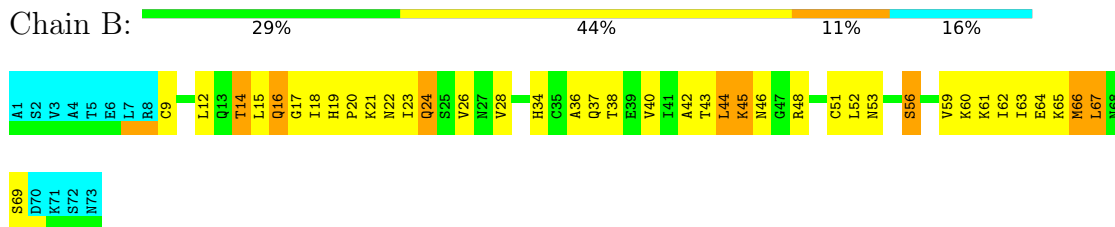


#### 4.2.25 Score per residue for model 25

- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY



- Molecule 1: HUMAN MELANOMA GROWTH STIMULATORY ACTIVITY





## 5 Refinement protocol and experimental data overview

Of the ? calculated structures, 25 were deposited, based on the following criterion: ?.

The authors did not provide any information on software used for structure solution, optimization or refinement.

No chemical shift data was provided.

## 6 Model quality [i](#)

### 6.1 Standard geometry [i](#)

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

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	1.0±0.2
1	B	0.0±0.0	1.0±0.2
All	All	0	48

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	A	48	ARG	Sidechain	24
1	B	48	ARG	Sidechain	24

### 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	457	487	487	72±8
1	B	457	487	487	72±9
All	All	22850	24350	24350	3362

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 71.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:67:LEU:HD11	1:B:63:ILE:HG21	1.13	1.16	23	10
1:A:63:ILE:HG21	1:B:67:LEU:HD11	1.11	1.21	4	11
1:A:67:LEU:HD11	1:B:63:ILE:CG2	1.01	1.86	17	25
1:A:63:ILE:CG2	1:B:67:LEU:HD11	1.01	1.84	4	25
1:A:40:VAL:HG11	1:A:63:ILE:HD11	0.98	1.00	18	21
1:B:40:VAL:CG1	1:B:63:ILE:HD11	0.98	1.87	18	4
1:A:40:VAL:CG1	1:A:63:ILE:HD11	0.98	1.87	18	4
1:B:40:VAL:HG11	1:B:63:ILE:HD11	0.96	1.00	18	22
1:B:52:LEU:HD13	1:B:62:ILE:HD13	0.94	1.39	7	12
1:A:18:ILE:CD1	1:A:23:ILE:HD11	0.91	1.96	4	12
1:B:22:ASN:O	1:B:23:ILE:HD13	0.91	1.65	12	7
1:A:52:LEU:HD13	1:A:62:ILE:HD13	0.91	1.40	20	12
1:A:22:ASN:O	1:A:23:ILE:HD13	0.91	1.65	12	7
1:B:18:ILE:CD1	1:B:23:ILE:HD11	0.91	1.96	4	12
1:A:67:LEU:HD11	1:B:63:ILE:HG22	0.90	1.42	16	18
1:B:40:VAL:HG11	1:B:63:ILE:CD1	0.89	1.95	18	5
1:A:22:ASN:O	1:A:44:LEU:HD23	0.89	1.68	16	8
1:B:22:ASN:O	1:B:44:LEU:HD23	0.89	1.68	16	8
1:B:15:LEU:HD12	1:B:51:CYS:O	0.88	1.68	21	3
1:A:63:ILE:HG22	1:B:67:LEU:HD11	0.88	1.42	16	18
1:A:15:LEU:HD12	1:A:51:CYS:O	0.88	1.68	21	3
1:A:23:ILE:HG23	1:A:42:ALA:HB1	0.86	1.47	11	24
1:B:17:GLY:HA2	1:B:52:LEU:HD22	0.86	1.47	15	18
1:B:23:ILE:HG23	1:B:42:ALA:HB1	0.86	1.47	11	24
1:A:17:GLY:HA2	1:A:52:LEU:HD22	0.86	1.48	19	18
1:B:12:LEU:O	1:B:12:LEU:HD13	0.85	1.69	16	2
1:A:40:VAL:HG11	1:A:63:ILE:CD1	0.85	1.96	18	4
1:A:23:ILE:CG2	1:A:42:ALA:HB1	0.85	2.02	11	19
1:A:12:LEU:O	1:A:12:LEU:HD13	0.85	1.69	16	2
1:B:59:VAL:HA	1:B:62:ILE:HD12	0.85	1.47	1	12
1:B:23:ILE:CG2	1:B:42:ALA:HB1	0.84	2.02	11	18
1:A:59:VAL:HA	1:A:62:ILE:HD12	0.84	1.47	1	12
1:A:18:ILE:HD11	1:A:23:ILE:CG1	0.82	2.04	7	1
1:B:18:ILE:HD11	1:B:23:ILE:CG1	0.82	2.04	7	1
1:B:9:CYS:HB2	1:B:12:LEU:HD22	0.79	1.55	23	13
1:A:9:CYS:HB2	1:A:12:LEU:HD22	0.79	1.55	23	13
1:A:18:ILE:HD13	1:A:18:ILE:N	0.78	1.94	23	5
1:A:17:GLY:O	1:A:18:ILE:HG23	0.78	1.78	11	13
1:B:17:GLY:O	1:B:18:ILE:HG23	0.77	1.78	11	13
1:B:18:ILE:N	1:B:18:ILE:HD13	0.77	1.94	21	5
1:A:18:ILE:HG22	1:A:52:LEU:HD11	0.74	1.59	16	2
1:A:16:GLN:O	1:A:59:VAL:HG23	0.74	1.83	17	14

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:52:LEU:HB3	1:B:59:VAL:HG22	0.74	1.59	22	25
1:B:16:GLN:O	1:B:59:VAL:HG23	0.74	1.83	17	14
1:A:52:LEU:HB3	1:A:59:VAL:HG22	0.73	1.59	22	25
1:B:18:ILE:HG22	1:B:52:LEU:HD11	0.72	1.59	16	2
1:A:18:ILE:HD13	1:A:23:ILE:HD11	0.71	1.63	6	4
1:A:44:LEU:HD12	1:A:48:ARG:HB2	0.70	1.61	16	7
1:B:44:LEU:HD12	1:B:48:ARG:HB2	0.70	1.62	16	7
1:A:28:VAL:HG22	1:A:40:VAL:HG13	0.70	1.63	17	1
1:B:18:ILE:HD13	1:B:23:ILE:HD11	0.70	1.63	6	5
1:A:63:ILE:HG21	1:B:67:LEU:CD1	0.70	2.13	17	15
1:B:52:LEU:CD1	1:B:62:ILE:HD13	0.70	2.17	11	4
1:B:28:VAL:HG22	1:B:40:VAL:HG13	0.69	1.63	17	1
1:A:18:ILE:CD1	1:A:52:LEU:HD21	0.68	2.19	23	4
1:B:28:VAL:HG13	1:B:40:VAL:HG22	0.68	1.66	23	8
1:B:18:ILE:CD1	1:B:52:LEU:HD21	0.68	2.19	23	4
1:A:52:LEU:CD1	1:A:62:ILE:HD13	0.67	2.17	11	3
1:A:67:LEU:CD1	1:B:63:ILE:HG21	0.67	2.13	17	15
1:A:28:VAL:HG13	1:A:40:VAL:HG22	0.66	1.68	19	8
1:A:23:ILE:HG21	1:A:42:ALA:HB1	0.66	1.67	10	8
1:B:23:ILE:HG21	1:B:42:ALA:HB1	0.66	1.67	10	6
1:A:40:VAL:HG21	1:A:54:PRO:HB3	0.66	1.67	12	2
1:B:27:ASN:O	1:B:41:ILE:HD12	0.65	1.90	3	11
1:B:18:ILE:HD11	1:B:23:ILE:HG12	0.65	1.66	7	1
1:B:23:ILE:HD13	1:B:62:ILE:HG23	0.65	1.68	25	6
1:A:18:ILE:HD13	1:A:18:ILE:H	0.65	1.52	22	4
1:A:27:ASN:O	1:A:41:ILE:HD12	0.65	1.91	3	11
1:B:63:ILE:HD12	1:B:63:ILE:N	0.64	2.08	17	24
1:A:64:GLU:OE2	1:B:67:LEU:HD13	0.64	1.92	18	1
1:A:26:VAL:O	1:B:28:VAL:HG23	0.64	1.93	25	2
1:A:63:ILE:HD12	1:A:63:ILE:N	0.64	2.08	17	24
1:A:67:LEU:HD13	1:B:64:GLU:OE2	0.64	1.92	18	1
1:A:18:ILE:HD11	1:A:23:ILE:HG12	0.64	1.66	7	1
1:B:40:VAL:HG21	1:B:54:PRO:HB3	0.64	1.67	12	2
1:B:18:ILE:HD12	1:B:19:HIS:O	0.64	1.92	7	1
1:B:23:ILE:HG12	1:B:42:ALA:HB1	0.63	1.69	25	3
1:B:18:ILE:HD13	1:B:18:ILE:H	0.63	1.52	22	4
1:A:23:ILE:HD13	1:A:62:ILE:CG2	0.63	2.24	25	5
1:A:28:VAL:HG23	1:B:26:VAL:O	0.63	1.93	25	2
1:B:23:ILE:CD1	1:B:62:ILE:HG23	0.63	2.24	24	6
1:A:14:THR:HG22	1:A:52:LEU:O	0.63	1.94	19	10
1:A:18:ILE:HD12	1:A:19:HIS:O	0.63	1.92	7	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:23:ILE:CD1	1:A:62:ILE:HG23	0.63	2.24	13	6
1:B:14:THR:HG22	1:B:52:LEU:O	0.63	1.94	19	10
1:A:23:ILE:HD13	1:A:62:ILE:HG23	0.63	1.69	25	6
1:B:23:ILE:HD13	1:B:62:ILE:CG2	0.62	2.24	25	5
1:B:53:ASN:O	1:B:59:VAL:HG21	0.62	1.94	7	25
1:A:44:LEU:HD12	1:A:48:ARG:CB	0.62	2.25	16	3
1:A:23:ILE:HG12	1:A:42:ALA:HB1	0.62	1.69	25	3
1:A:53:ASN:O	1:A:59:VAL:HG21	0.62	1.95	18	25
1:A:24:GLN:HG2	1:A:43:THR:HG22	0.61	1.73	6	4
1:B:18:ILE:HG22	1:B:52:LEU:CD1	0.61	2.24	16	3
1:B:44:LEU:HD12	1:B:48:ARG:CB	0.61	2.25	16	3
1:A:18:ILE:HG22	1:A:52:LEU:CD1	0.61	2.24	16	3
1:B:52:LEU:HD13	1:B:62:ILE:CD1	0.61	2.24	11	2
1:A:25:SER:O	1:A:26:VAL:HG23	0.60	1.97	12	9
1:B:24:GLN:HG2	1:B:43:THR:HG22	0.60	1.73	6	4
1:B:25:SER:O	1:B:26:VAL:HG23	0.59	1.97	12	9
1:B:58:ILE:CD1	1:B:62:ILE:CD1	0.59	2.81	21	2
1:A:67:LEU:CD1	1:B:63:ILE:CG2	0.59	2.81	20	19
1:B:58:ILE:O	1:B:62:ILE:HD13	0.59	1.98	18	13
1:A:29:LYS:HZ3	1:A:29:LYS:HB3	0.59	1.57	6	1
1:A:63:ILE:CG2	1:B:67:LEU:CD1	0.59	2.81	25	19
1:A:18:ILE:CD1	1:A:23:ILE:CD1	0.59	2.81	17	12
1:A:63:ILE:HG22	1:B:67:LEU:CD1	0.59	2.28	6	10
1:A:53:ASN:HD22	1:A:55:ALA:HB3	0.59	1.58	22	2
1:B:18:ILE:HG22	1:B:52:LEU:CD2	0.59	2.28	12	1
1:B:18:ILE:CD1	1:B:23:ILE:CD1	0.58	2.81	11	12
1:A:58:ILE:CD1	1:A:62:ILE:CD1	0.58	2.81	17	2
1:B:63:ILE:N	1:B:63:ILE:CD1	0.58	2.67	7	18
1:B:18:ILE:HD13	1:B:23:ILE:CD1	0.58	2.28	8	4
1:B:19:HIS:CB	1:B:22:ASN:ND2	0.58	2.67	24	4
1:A:19:HIS:CB	1:A:22:ASN:ND2	0.58	2.67	25	4
1:A:12:LEU:HD13	1:A:12:LEU:C	0.58	2.19	16	1
1:A:63:ILE:N	1:A:63:ILE:CD1	0.58	2.67	7	17
1:B:12:LEU:HD13	1:B:12:LEU:C	0.58	2.19	16	1
1:A:52:LEU:HD13	1:A:62:ILE:CD1	0.58	2.24	11	2
1:A:58:ILE:O	1:A:62:ILE:HD13	0.58	1.98	18	13
1:A:18:ILE:HG22	1:A:52:LEU:CD2	0.58	2.28	12	1
1:B:22:ASN:O	1:B:44:LEU:CD2	0.57	2.53	14	22
1:B:58:ILE:CG2	1:B:59:VAL:N	0.57	2.67	21	17
1:B:29:LYS:CE	1:B:34:HIS:NE2	0.57	2.67	23	2
1:A:58:ILE:O	1:A:58:ILE:HD13	0.57	1.99	11	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:22:ASN:O	1:A:44:LEU:CD2	0.57	2.53	14	22
1:A:67:LEU:CD1	1:B:63:ILE:HG22	0.57	2.28	6	11
1:B:18:ILE:CD1	1:B:23:ILE:CG1	0.57	2.81	7	4
1:B:18:ILE:N	1:B:18:ILE:CD1	0.57	2.67	21	5
1:B:18:ILE:HD11	1:B:52:LEU:HD21	0.57	1.76	23	1
1:A:58:ILE:CG2	1:A:59:VAL:N	0.57	2.67	17	17
1:B:58:ILE:HD13	1:B:58:ILE:O	0.57	1.99	11	3
1:A:18:ILE:CD1	1:A:52:LEU:CD2	0.57	2.82	23	1
1:A:18:ILE:HD13	1:A:23:ILE:CD1	0.57	2.28	8	4
1:A:29:LYS:CE	1:A:34:HIS:NE2	0.57	2.67	23	2
1:B:53:ASN:HD22	1:B:55:ALA:HB3	0.57	1.58	22	2
1:B:26:VAL:HG23	1:B:66:MET:HG3	0.56	1.77	13	5
1:B:53:ASN:O	1:B:59:VAL:CB	0.56	2.53	11	25
1:A:43:THR:HG22	1:A:43:THR:O	0.56	2.00	10	5
1:A:53:ASN:O	1:A:59:VAL:CB	0.56	2.53	11	25
1:A:58:ILE:HG23	1:A:59:VAL:N	0.56	2.15	19	16
1:B:18:ILE:CD1	1:B:52:LEU:CD2	0.56	2.83	23	1
1:A:18:ILE:HD11	1:A:52:LEU:HD21	0.56	1.76	23	1
1:B:58:ILE:HG23	1:B:59:VAL:N	0.56	2.16	5	16
1:B:43:THR:O	1:B:43:THR:HG22	0.56	2.00	10	8
1:A:26:VAL:HG23	1:A:66:MET:HG3	0.56	1.77	13	5
1:A:53:ASN:O	1:A:59:VAL:CG2	0.55	2.54	12	25
1:B:53:ASN:O	1:B:59:VAL:CG2	0.55	2.54	12	25
1:B:53:ASN:N	1:B:59:VAL:HG21	0.55	2.17	4	23
1:A:19:HIS:CG	1:A:20:PRO:HD2	0.55	2.37	19	22
1:B:49:LYS:CD	1:B:49:LYS:N	0.55	2.70	16	6
1:B:19:HIS:CG	1:B:20:PRO:HD2	0.55	2.37	9	22
1:A:49:LYS:N	1:A:49:LYS:CD	0.54	2.70	16	14
1:A:53:ASN:N	1:A:59:VAL:HG21	0.54	2.17	4	23
1:A:18:ILE:CD1	1:A:23:ILE:CG1	0.54	2.81	7	3
1:A:34:HIS:N	1:A:34:HIS:ND1	0.54	2.56	11	16
1:A:18:ILE:H	1:A:62:ILE:HD11	0.54	1.63	25	4
1:A:24:GLN:CG	1:A:43:THR:O	0.54	2.56	5	14
1:B:21:LYS:O	1:B:45:LYS:CD	0.54	2.56	9	6
1:B:29:LYS:HZ3	1:B:29:LYS:HB3	0.54	1.63	6	1
1:B:58:ILE:O	1:B:62:ILE:CD1	0.54	2.56	18	11
1:A:18:ILE:HD13	1:A:23:ILE:CG1	0.54	2.33	11	3
1:A:58:ILE:O	1:A:62:ILE:CD1	0.53	2.56	8	11
1:B:21:LYS:O	1:B:45:LYS:CE	0.53	2.56	12	10
1:A:24:GLN:CB	1:A:43:THR:O	0.53	2.57	11	11
1:A:21:LYS:O	1:A:45:LYS:CD	0.53	2.57	8	6

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:21:LYS:O	1:A:45:LYS:CE	0.53	2.57	12	10
1:B:24:GLN:CG	1:B:43:THR:O	0.53	2.56	21	14
1:B:29:LYS:CG	1:B:29:LYS:O	0.53	2.57	17	4
1:B:24:GLN:CB	1:B:43:THR:O	0.53	2.57	15	11
1:B:19:HIS:ND1	1:B:20:PRO:HD2	0.53	2.18	9	12
1:B:18:ILE:CG2	1:B:52:LEU:HD11	0.53	2.33	16	1
1:B:34:HIS:N	1:B:34:HIS:ND1	0.53	2.56	11	17
1:A:16:GLN:O	1:A:58:ILE:CG2	0.53	2.57	19	14
1:B:16:GLN:O	1:B:58:ILE:CG2	0.53	2.57	14	13
1:B:58:ILE:CD1	1:B:62:ILE:HD11	0.53	2.34	17	2
1:B:18:ILE:HD13	1:B:23:ILE:CG1	0.53	2.33	11	3
1:A:14:THR:CG2	1:A:52:LEU:O	0.53	2.57	5	7
1:B:14:THR:CG2	1:B:52:LEU:O	0.53	2.57	5	7
1:A:40:VAL:HG11	1:A:63:ILE:CG1	0.53	2.33	14	1
1:A:54:PRO:O	1:A:60:LYS:CE	0.53	2.57	24	5
1:B:40:VAL:HG11	1:B:63:ILE:CG1	0.53	2.33	14	1
1:B:23:ILE:HG22	1:B:24:GLN:N	0.53	2.19	19	6
1:A:25:SER:O	1:A:26:VAL:CG2	0.53	2.57	12	9
1:B:54:PRO:O	1:B:60:LYS:CE	0.53	2.57	24	5
1:A:58:ILE:HD13	1:A:62:ILE:CD1	0.53	2.34	21	2
1:B:58:ILE:HD13	1:B:62:ILE:CD1	0.53	2.34	21	2
1:A:23:ILE:HG22	1:A:24:GLN:N	0.53	2.19	19	6
1:A:19:HIS:ND1	1:A:20:PRO:HD2	0.53	2.19	9	13
1:A:31:PRO:CB	1:A:35:CYS:O	0.53	2.57	6	4
1:B:31:PRO:CB	1:B:36:ALA:O	0.53	2.57	22	5
1:B:29:LYS:CD	1:B:29:LYS:O	0.53	2.57	5	1
1:B:44:LEU:CB	1:B:46:ASN:OD1	0.53	2.57	12	2
1:A:18:ILE:N	1:A:18:ILE:CD1	0.53	2.67	21	5
1:A:61:LYS:CE	1:A:64:GLU:OE2	0.53	2.57	23	1
1:B:60:LYS:O	1:B:64:GLU:CG	0.52	2.57	7	3
1:A:31:PRO:CB	1:A:36:ALA:O	0.52	2.57	22	5
1:B:61:LYS:O	1:B:65:LYS:CG	0.52	2.58	6	6
1:A:18:ILE:O	1:A:62:ILE:CG1	0.52	2.57	13	2
1:A:29:LYS:CG	1:A:29:LYS:O	0.52	2.57	17	5
1:B:25:SER:O	1:B:26:VAL:CG2	0.52	2.57	12	9
1:B:31:PRO:CB	1:B:35:CYS:O	0.52	2.57	20	4
1:B:53:ASN:ND2	1:B:55:ALA:HB3	0.52	2.19	22	3
1:A:11:CYS:SG	1:A:41:ILE:CG1	0.52	2.98	12	3
1:A:67:LEU:CD1	1:B:64:GLU:OE2	0.52	2.57	18	1
1:A:53:ASN:ND2	1:A:55:ALA:HB3	0.52	2.19	22	2
1:B:18:ILE:O	1:B:62:ILE:CG1	0.52	2.57	13	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:18:ILE:H	1:B:62:ILE:HD11	0.52	1.63	25	4
1:A:64:GLU:OE2	1:B:67:LEU:CD1	0.52	2.57	18	1
1:A:61:LYS:O	1:A:65:LYS:CG	0.52	2.57	21	6
1:A:43:THR:OG1	1:A:49:LYS:CD	0.52	2.57	12	1
1:B:43:THR:OG1	1:B:49:LYS:CD	0.52	2.57	12	1
1:A:18:ILE:CG2	1:A:52:LEU:HD11	0.52	2.33	16	1
1:B:10:GLN:OE1	1:B:29:LYS:NZ	0.52	2.43	6	1
1:B:11:CYS:SG	1:B:41:ILE:CG1	0.52	2.98	12	3
1:A:44:LEU:CB	1:A:46:ASN:OD1	0.52	2.57	12	2
1:A:58:ILE:CD1	1:A:62:ILE:HD11	0.52	2.34	21	2
1:A:21:LYS:O	1:A:45:LYS:NZ	0.52	2.43	25	6
1:B:41:ILE:CG1	1:B:51:CYS:SG	0.52	2.98	8	3
1:A:29:LYS:O	1:A:29:LYS:CD	0.52	2.57	5	1
1:B:29:LYS:O	1:B:29:LYS:CG	0.52	2.57	5	1
1:B:22:ASN:HB3	1:B:44:LEU:HD22	0.51	1.81	16	6
1:A:41:ILE:CG1	1:A:51:CYS:SG	0.51	2.98	8	3
1:B:61:LYS:CE	1:B:64:GLU:OE2	0.51	2.57	23	1
1:A:60:LYS:O	1:A:64:GLU:CG	0.51	2.57	7	2
1:B:21:LYS:O	1:B:45:LYS:NZ	0.51	2.43	16	6
1:B:19:HIS:HB2	1:B:22:ASN:ND2	0.51	2.21	13	4
1:A:18:ILE:CD1	1:A:23:ILE:HG13	0.51	2.36	7	1
1:B:10:GLN:N	1:B:39:GLU:OE1	0.51	2.43	12	1
1:A:22:ASN:HB3	1:A:44:LEU:HD22	0.51	1.81	12	6
1:B:18:ILE:CD1	1:B:23:ILE:HG13	0.51	2.36	7	1
1:B:24:GLN:N	1:B:43:THR:O	0.51	2.44	12	7
1:A:18:ILE:HD12	1:A:23:ILE:HD11	0.51	1.81	5	3
1:A:10:GLN:OE1	1:A:29:LYS:NZ	0.51	2.43	6	1
1:A:23:ILE:CD1	1:A:62:ILE:CG2	0.51	2.89	10	4
1:A:10:GLN:N	1:A:39:GLU:OE1	0.51	2.43	12	1
1:B:24:GLN:HG3	1:B:25:SER:N	0.50	2.21	3	15
1:A:58:ILE:HD13	1:A:58:ILE:O	0.50	2.06	21	1
1:B:10:GLN:HG3	1:B:29:LYS:CE	0.50	2.37	4	3
1:A:24:GLN:HG3	1:A:25:SER:N	0.50	2.22	7	15
1:A:10:GLN:HG3	1:A:29:LYS:CE	0.50	2.36	4	3
1:A:24:GLN:CG	1:A:43:THR:HG22	0.50	2.37	6	3
1:A:24:GLN:N	1:A:43:THR:O	0.50	2.44	12	7
1:A:19:HIS:HB2	1:A:22:ASN:ND2	0.50	2.21	13	4
1:A:29:LYS:HE2	1:A:34:HIS:NE2	0.50	2.21	23	1
1:B:19:HIS:HB3	1:B:22:ASN:ND2	0.50	2.22	25	4
1:B:24:GLN:CG	1:B:43:THR:HG22	0.50	2.37	6	3
1:A:61:LYS:CE	1:A:64:GLU:CD	0.49	2.81	6	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:29:LYS:HE2	1:B:41:ILE:CD1	0.49	2.37	6	1
1:B:29:LYS:HE3	1:B:41:ILE:CD1	0.49	2.38	14	1
1:B:61:LYS:CE	1:B:64:GLU:CD	0.49	2.81	6	2
1:B:29:LYS:HE2	1:B:34:HIS:NE2	0.49	2.21	23	1
1:A:52:LEU:HB3	1:A:59:VAL:CG2	0.49	2.38	16	9
1:B:34:HIS:ND1	1:B:34:HIS:N	0.49	2.60	18	4
1:A:16:GLN:O	1:A:59:VAL:CG2	0.49	2.60	21	1
1:A:29:LYS:HE2	1:A:41:ILE:CD1	0.49	2.37	6	1
1:A:61:LYS:NZ	1:A:64:GLU:OE2	0.49	2.43	6	1
1:A:29:LYS:HE3	1:A:34:HIS:NE2	0.49	2.23	9	3
1:B:23:ILE:CD1	1:B:62:ILE:CG2	0.49	2.89	10	4
1:A:29:LYS:HE3	1:A:41:ILE:CD1	0.49	2.38	14	1
1:A:15:LEU:HB2	1:A:52:LEU:HD23	0.48	1.85	13	2
1:B:18:ILE:HD12	1:B:23:ILE:HD11	0.48	1.84	20	6
1:A:10:GLN:NE2	1:A:29:LYS:HD3	0.48	2.23	8	1
1:A:19:HIS:HB3	1:A:22:ASN:ND2	0.48	2.22	25	4
1:B:65:LYS:O	1:B:69:SER:N	0.48	2.46	13	3
1:B:29:LYS:HE3	1:B:34:HIS:NE2	0.48	2.23	9	3
1:A:10:GLN:CG	1:A:29:LYS:CD	0.48	2.92	17	1
1:B:10:GLN:CG	1:B:29:LYS:CE	0.48	2.92	4	1
1:B:45:LYS:HZ2	1:B:45:LYS:HB3	0.48	1.68	5	2
1:B:10:GLN:OE1	1:B:29:LYS:CE	0.48	2.61	6	1
1:B:10:GLN:NE2	1:B:29:LYS:HD3	0.48	2.23	8	1
1:A:65:LYS:O	1:A:69:SER:N	0.48	2.46	13	3
1:A:18:ILE:HD12	1:A:18:ILE:O	0.48	2.09	8	9
1:B:9:CYS:CA	1:B:39:GLU:OE2	0.48	2.62	2	1
1:A:24:GLN:HG3	1:A:43:THR:O	0.48	2.09	3	14
1:A:10:GLN:CG	1:A:29:LYS:CE	0.48	2.92	4	1
1:A:18:ILE:HD11	1:A:23:ILE:HD11	0.48	1.84	20	2
1:A:10:GLN:OE1	1:A:29:LYS:CE	0.48	2.61	6	1
1:B:61:LYS:NZ	1:B:64:GLU:OE2	0.48	2.43	6	1
1:A:31:PRO:CA	1:A:35:CYS:O	0.48	2.62	6	8
1:B:17:GLY:CA	1:B:52:LEU:HD22	0.48	2.39	13	2
1:B:49:LYS:HD2	1:B:49:LYS:N	0.48	2.24	1	9
1:A:52:LEU:CB	1:A:59:VAL:HG22	0.48	2.38	11	2
1:A:49:LYS:N	1:A:49:LYS:HD2	0.48	2.24	1	11
1:A:61:LYS:CE	1:A:61:LYS:HA	0.47	2.39	21	10
1:B:18:ILE:HD12	1:B:18:ILE:O	0.47	2.09	8	8
1:B:61:LYS:HA	1:B:61:LYS:CE	0.47	2.39	2	7
1:B:52:LEU:HB3	1:B:59:VAL:CG2	0.47	2.38	16	8
1:A:10:GLN:HB2	1:A:29:LYS:CD	0.47	2.39	20	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:10:GLN:HB2	1:B:29:LYS:CD	0.47	2.39	20	3
1:A:40:VAL:HG23	1:A:54:PRO:HG3	0.47	1.86	13	1
1:B:58:ILE:HD13	1:B:62:ILE:HD13	0.47	1.86	17	2
1:B:24:GLN:HG3	1:B:43:THR:O	0.47	2.09	3	14
1:B:39:GLU:OE1	1:B:39:GLU:N	0.47	2.48	8	1
1:B:16:GLN:O	1:B:59:VAL:CG2	0.47	2.60	21	1
1:A:10:GLN:HG3	1:A:29:LYS:NZ	0.47	2.24	22	1
1:B:31:PRO:CA	1:B:35:CYS:O	0.47	2.62	6	8
1:B:43:THR:O	1:B:44:LEU:O	0.47	2.32	12	21
1:A:64:GLU:HG3	1:A:65:LYS:N	0.47	2.24	6	6
1:A:17:GLY:CA	1:A:52:LEU:HD22	0.47	2.39	11	3
1:B:15:LEU:HB2	1:B:52:LEU:HD23	0.47	1.86	13	1
1:B:10:GLN:O	1:B:11:CYS:SG	0.47	2.72	16	1
1:A:10:GLN:NE2	1:A:29:LYS:HD2	0.47	2.25	19	3
1:B:64:GLU:HG3	1:B:65:LYS:N	0.47	2.24	6	6
1:B:61:LYS:CE	1:B:61:LYS:HA	0.47	2.39	6	3
1:A:39:GLU:OE1	1:A:39:GLU:N	0.47	2.48	8	1
1:A:18:ILE:HG13	1:A:52:LEU:CD2	0.47	2.40	9	1
1:B:49:LYS:N	1:B:49:LYS:HD2	0.47	2.25	19	2
1:B:49:LYS:N	1:B:49:LYS:CD	0.47	2.78	1	7
1:B:61:LYS:O	1:B:65:LYS:CB	0.47	2.63	24	9
1:B:18:ILE:HG13	1:B:52:LEU:CD2	0.47	2.40	9	1
1:B:58:ILE:C	1:B:58:ILE:CD1	0.47	2.82	11	1
1:A:10:GLN:O	1:A:11:CYS:SG	0.47	2.73	16	1
1:B:10:GLN:CG	1:B:29:LYS:CD	0.47	2.92	17	1
1:A:20:PRO:HG3	1:A:65:LYS:CE	0.47	2.40	18	1
1:B:10:GLN:HG3	1:B:29:LYS:NZ	0.47	2.24	22	1
1:A:43:THR:O	1:A:44:LEU:O	0.47	2.33	1	21
1:B:20:PRO:HG3	1:B:65:LYS:CD	0.47	2.40	3	2
1:B:46:ASN:ND2	1:B:48:ARG:HD2	0.47	2.24	17	2
1:A:41:ILE:N	1:A:41:ILE:HD12	0.47	2.25	10	1
1:A:58:ILE:C	1:A:58:ILE:CD1	0.47	2.82	11	1
1:B:53:ASN:O	1:B:56:SER:OG	0.47	2.33	13	11
1:A:28:VAL:O	1:B:66:MET:CE	0.47	2.63	18	1
1:A:9:CYS:CA	1:A:39:GLU:OE2	0.47	2.62	2	1
1:A:60:LYS:O	1:A:64:GLU:OE2	0.47	2.33	11	3
1:A:46:ASN:ND2	1:A:48:ARG:HD2	0.47	2.24	17	2
1:A:58:ILE:HD13	1:A:62:ILE:HD13	0.47	1.86	17	2
1:A:22:ASN:O	1:A:44:LEU:HD22	0.46	2.10	5	12
1:B:10:GLN:NE2	1:B:29:LYS:HD2	0.46	2.25	19	3
1:A:61:LYS:O	1:A:65:LYS:CB	0.46	2.63	24	9

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:20:PRO:HG3	1:A:65:LYS:CD	0.46	2.40	3	2
1:A:37:GLN:O	1:A:39:GLU:OE2	0.46	2.34	15	6
1:A:61:LYS:N	1:A:61:LYS:HD2	0.46	2.25	19	3
1:B:61:LYS:N	1:B:61:LYS:HD2	0.46	2.25	19	3
1:A:37:GLN:O	1:A:39:GLU:OE1	0.46	2.34	14	5
1:A:53:ASN:C	1:A:59:VAL:HG11	0.46	2.31	11	5
1:B:29:LYS:O	1:B:30:SER:O	0.46	2.33	13	4
1:B:53:ASN:OD1	1:B:55:ALA:HB3	0.46	2.11	18	1
1:A:17:GLY:O	1:A:18:ILE:CG2	0.46	2.61	1	4
1:B:60:LYS:O	1:B:64:GLU:OE2	0.46	2.33	2	3
1:A:53:ASN:O	1:A:56:SER:OG	0.46	2.34	5	11
1:B:41:ILE:HD12	1:B:41:ILE:N	0.46	2.25	10	1
1:A:61:LYS:HB3	1:A:65:LYS:NZ	0.46	2.25	1	1
1:A:27:ASN:OD1	1:B:27:ASN:OD1	0.46	2.34	11	6
1:A:43:THR:O	1:A:43:THR:HG22	0.46	2.09	5	3
1:A:29:LYS:O	1:A:30:SER:O	0.46	2.34	13	3
1:B:40:VAL:HG23	1:B:54:PRO:HG3	0.46	1.86	13	1
1:B:37:GLN:O	1:B:39:GLU:OE1	0.46	2.33	21	5
1:B:53:ASN:C	1:B:59:VAL:HG11	0.46	2.31	11	6
1:A:14:THR:OG1	1:A:53:ASN:OD1	0.46	2.34	25	1
1:B:37:GLN:O	1:B:39:GLU:OE2	0.46	2.34	21	6
1:A:38:THR:O	1:A:39:GLU:OE1	0.46	2.34	7	1
1:A:14:THR:HG21	1:A:53:ASN:HD22	0.46	1.71	16	2
1:A:38:THR:O	1:A:39:GLU:OE2	0.46	2.34	5	2
1:B:38:THR:O	1:B:39:GLU:OE2	0.46	2.34	21	2
1:A:44:LEU:HB2	1:A:46:ASN:OD1	0.46	2.11	12	4
1:A:10:GLN:HB3	1:A:41:ILE:HD11	0.46	1.87	16	1
1:A:37:GLN:OE1	1:A:37:GLN:CA	0.46	2.63	7	2
1:A:63:ILE:HB	1:B:67:LEU:HD11	0.46	1.87	11	1
1:A:28:VAL:CG2	1:A:40:VAL:HG13	0.46	2.39	17	1
1:B:17:GLY:O	1:B:18:ILE:CG2	0.46	2.61	1	4
1:A:22:ASN:O	1:A:23:ILE:CD1	0.46	2.57	6	1
1:B:44:LEU:HB2	1:B:46:ASN:OD1	0.46	2.11	22	4
1:A:67:LEU:HD11	1:B:63:ILE:HB	0.46	1.88	11	1
1:B:24:GLN:HB3	1:B:43:THR:O	0.46	2.11	25	4
1:A:12:LEU:C	1:A:12:LEU:CD1	0.46	2.84	16	1
1:B:12:LEU:C	1:B:12:LEU:CD1	0.46	2.84	16	1
1:A:66:MET:CE	1:B:28:VAL:O	0.46	2.63	18	1
1:B:38:THR:O	1:B:39:GLU:OE1	0.45	2.34	7	1
1:A:17:GLY:HA3	1:A:58:ILE:CG2	0.45	2.41	12	4
1:B:17:GLY:HA3	1:B:58:ILE:CG2	0.45	2.41	12	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:20:PRO:HG3	1:B:65:LYS:CE	0.45	2.40	18	1
1:A:15:LEU:CD1	1:A:50:ALA:HB1	0.45	2.42	7	11
1:B:60:LYS:O	1:B:64:GLU:HG2	0.45	2.11	9	6
1:A:24:GLN:HB3	1:A:43:THR:O	0.45	2.11	25	4
1:B:61:LYS:HB3	1:B:65:LYS:NZ	0.45	2.25	1	1
1:A:35:CYS:SG	1:A:39:GLU:OE2	0.45	2.74	2	8
1:B:35:CYS:SG	1:B:39:GLU:OE2	0.45	2.75	2	8
1:A:35:CYS:SG	1:A:39:GLU:OE1	0.45	2.75	16	2
1:A:43:THR:HG23	1:A:49:LYS:HG3	0.45	1.88	16	1
1:A:13:GLN:O	1:A:14:THR:HG23	0.45	2.11	1	2
1:A:24:GLN:HB2	1:A:43:THR:O	0.45	2.12	1	1
1:A:29:LYS:O	1:A:29:LYS:HG3	0.45	2.12	17	3
1:B:57:PRO:O	1:B:61:LYS:HD2	0.45	2.12	19	3
1:A:35:CYS:SG	1:A:37:GLN:OE1	0.45	2.74	8	2
1:B:52:LEU:CB	1:B:59:VAL:HG22	0.45	2.38	22	3
1:B:24:GLN:HB2	1:B:43:THR:O	0.45	2.12	1	1
1:B:25:SER:OG	1:B:43:THR:CB	0.45	2.65	12	5
1:B:22:ASN:O	1:B:44:LEU:HD22	0.45	2.11	4	10
1:A:54:PRO:O	1:A:60:LYS:HE3	0.45	2.12	4	4
1:B:54:PRO:O	1:B:60:LYS:HE3	0.45	2.12	12	4
1:A:60:LYS:O	1:A:64:GLU:HG2	0.45	2.11	9	6
1:A:33:PRO:HD2	1:A:34:HIS:CE1	0.45	2.47	11	2
1:A:53:ASN:OD1	1:A:55:ALA:HB3	0.45	2.10	18	1
1:A:20:PRO:HA	1:A:23:ILE:HD12	0.45	1.88	11	2
1:B:13:GLN:O	1:B:14:THR:HG23	0.45	2.11	1	2
1:A:66:MET:SD	1:A:69:SER:OG	0.45	2.74	25	4
1:B:58:ILE:O	1:B:58:ILE:CD1	0.45	2.64	11	1
1:A:14:THR:OG1	1:A:53:ASN:ND2	0.45	2.50	17	4
1:B:9:CYS:O	1:B:10:GLN:C	0.45	2.55	16	17
1:B:37:GLN:OE1	1:B:37:GLN:CA	0.45	2.64	6	2
1:B:20:PRO:HA	1:B:23:ILE:HD12	0.45	1.88	11	2
1:B:31:PRO:HB3	1:B:37:GLN:N	0.45	2.27	4	6
1:B:33:PRO:HD2	1:B:34:HIS:CE1	0.45	2.47	11	2
1:A:22:ASN:HB3	1:A:44:LEU:CD2	0.45	2.42	16	3
1:B:14:THR:OG1	1:B:53:ASN:ND2	0.45	2.50	17	4
1:B:10:GLN:HB3	1:B:41:ILE:HD11	0.45	1.87	16	1
1:B:35:CYS:SG	1:B:39:GLU:OE1	0.45	2.74	16	2
1:A:9:CYS:O	1:A:10:GLN:C	0.45	2.55	8	16
1:B:53:ASN:O	1:B:59:VAL:HB	0.45	2.11	25	24
1:A:25:SER:OG	1:A:43:THR:OG1	0.45	2.34	2	2
1:B:66:MET:SD	1:B:69:SER:OG	0.45	2.74	25	5

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:23:ILE:O	1:A:45:LYS:NZ	0.45	2.49	24	2
1:A:66:MET:O	1:A:69:SER:OG	0.45	2.34	6	3
1:A:41:ILE:HG22	1:A:42:ALA:N	0.45	2.26	9	2
1:B:9:CYS:SG	1:B:37:GLN:OE1	0.45	2.74	10	1
1:B:14:THR:HG21	1:B:53:ASN:HD22	0.45	1.71	16	2
1:B:35:CYS:SG	1:B:37:GLN:OE1	0.45	2.74	8	3
1:A:31:PRO:HB3	1:A:37:GLN:N	0.45	2.27	14	6
1:B:41:ILE:HG12	1:B:51:CYS:SG	0.45	2.52	8	2
1:B:29:LYS:O	1:B:29:LYS:HD3	0.45	2.12	5	1
1:A:19:HIS:O	1:A:23:ILE:HD11	0.45	2.12	9	1
1:A:9:CYS:SG	1:A:37:GLN:OE1	0.45	2.74	10	1
1:A:30:SER:HB3	1:B:66:MET:CE	0.45	2.41	24	1
1:B:11:CYS:SG	1:B:41:ILE:HG13	0.44	2.52	12	8
1:B:15:LEU:CD1	1:B:50:ALA:HB1	0.44	2.42	7	8
1:B:25:SER:OG	1:B:43:THR:OG1	0.44	2.34	2	3
1:B:56:SER:O	1:B:60:LYS:HG3	0.44	2.12	13	12
1:A:25:SER:OG	1:A:43:THR:CB	0.44	2.64	12	4
1:A:10:GLN:NE2	1:A:29:LYS:CD	0.44	2.80	8	1
1:A:58:ILE:O	1:A:58:ILE:CD1	0.44	2.64	11	1
1:B:22:ASN:HB3	1:B:44:LEU:CD2	0.44	2.42	16	3
1:A:46:ASN:ND2	1:A:48:ARG:HB2	0.44	2.28	15	13
1:B:56:SER:O	1:B:60:LYS:CG	0.44	2.66	21	3
1:B:27:ASN:O	1:B:41:ILE:CD1	0.44	2.63	3	2
1:B:57:PRO:O	1:B:61:LYS:HG2	0.44	2.12	6	1
1:B:29:LYS:O	1:B:29:LYS:HG3	0.44	2.13	14	2
1:B:57:PRO:O	1:B:61:LYS:CG	0.44	2.66	6	1
1:A:61:LYS:O	1:A:65:LYS:HB3	0.44	2.13	7	1
1:B:41:ILE:HG22	1:B:42:ALA:N	0.44	2.26	9	2
1:A:19:HIS:CD2	1:A:21:LYS:CG	0.44	3.00	15	1
1:A:15:LEU:O	1:A:16:GLN:C	0.44	2.56	25	22
1:A:22:ASN:O	1:A:44:LEU:HA	0.44	2.12	14	15
1:A:53:ASN:O	1:A:59:VAL:HB	0.44	2.13	21	24
1:A:61:LYS:O	1:A:65:LYS:HB2	0.44	2.13	2	10
1:A:9:CYS:HB2	1:A:12:LEU:CD2	0.44	2.43	24	5
1:A:57:PRO:O	1:A:61:LYS:HD2	0.44	2.12	19	3
1:A:28:VAL:CG1	1:A:40:VAL:HG22	0.44	2.41	11	2
1:B:22:ASN:O	1:B:44:LEU:HA	0.44	2.12	14	12
1:A:56:SER:O	1:A:60:LYS:HG3	0.44	2.12	21	13
1:B:61:LYS:O	1:B:65:LYS:HB2	0.44	2.13	2	6
1:B:26:VAL:HG13	1:B:41:ILE:O	0.44	2.13	4	1
1:A:43:THR:OG1	1:A:49:LYS:HD2	0.44	2.13	12	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:18:ILE:CG2	1:B:52:LEU:HD21	0.44	2.42	12	1
1:B:43:THR:HG23	1:B:49:LYS:HG3	0.44	1.88	16	1
1:A:43:THR:O	1:A:44:LEU:C	0.44	2.56	16	19
1:A:61:LYS:CA	1:A:61:LYS:HE2	0.44	2.43	21	2
1:B:61:LYS:O	1:B:65:LYS:HB3	0.44	2.13	7	1
1:A:53:ASN:O	1:A:53:ASN:OD1	0.44	2.36	11	1
1:A:61:LYS:O	1:A:65:LYS:HG2	0.44	2.13	14	1
1:B:58:ILE:HD13	1:B:58:ILE:C	0.44	2.33	17	2
1:A:66:MET:CE	1:B:30:SER:HB3	0.44	2.41	24	1
1:B:14:THR:OG1	1:B:53:ASN:OD1	0.44	2.34	25	1
1:A:26:VAL:CG2	1:A:66:MET:HG3	0.44	2.43	7	15
1:B:46:ASN:ND2	1:B:48:ARG:HB2	0.44	2.28	21	13
1:A:34:HIS:ND1	1:A:34:HIS:N	0.44	2.64	13	7
1:B:25:SER:CB	1:B:43:THR:HB	0.44	2.42	22	3
1:A:18:ILE:O	1:A:19:HIS:C	0.44	2.56	14	9
1:A:58:ILE:HD13	1:A:58:ILE:C	0.44	2.33	17	2
1:B:61:LYS:HA	1:B:64:GLU:CG	0.44	2.43	19	1
1:A:9:CYS:O	1:A:11:CYS:N	0.44	2.51	2	14
1:B:9:CYS:O	1:B:11:CYS:N	0.44	2.51	10	14
1:B:63:ILE:O	1:B:64:GLU:C	0.44	2.56	18	8
1:B:43:THR:O	1:B:44:LEU:C	0.44	2.56	25	17
1:A:57:PRO:O	1:A:61:LYS:CG	0.44	2.66	6	1
1:B:19:HIS:CD2	1:B:21:LYS:CG	0.44	3.00	15	1
1:A:16:GLN:O	1:A:58:ILE:HG21	0.44	2.13	19	1
1:A:29:LYS:NZ	1:A:39:GLU:HG3	0.44	2.28	20	1
1:A:11:CYS:SG	1:A:41:ILE:HG13	0.44	2.52	12	8
1:A:31:PRO:HA	1:A:35:CYS:O	0.44	2.13	20	11
1:B:15:LEU:O	1:B:16:GLN:C	0.44	2.57	9	21
1:A:31:PRO:HB3	1:A:36:ALA:O	0.44	2.13	24	4
1:A:26:VAL:HG13	1:A:41:ILE:O	0.44	2.13	4	1
1:A:41:ILE:HG12	1:A:51:CYS:SG	0.44	2.53	8	2
1:A:63:ILE:O	1:A:65:LYS:N	0.44	2.51	12	1
1:B:11:CYS:O	1:B:12:LEU:C	0.43	2.57	17	21
1:A:25:SER:CB	1:A:43:THR:HB	0.43	2.42	2	3
1:B:10:GLN:NE2	1:B:29:LYS:CD	0.43	2.80	8	1
1:A:41:ILE:CG2	1:A:42:ALA:N	0.43	2.81	9	1
1:A:18:ILE:CG2	1:A:52:LEU:HD21	0.43	2.42	12	1
1:B:63:ILE:O	1:B:65:LYS:N	0.43	2.51	12	1
1:A:11:CYS:O	1:A:12:LEU:C	0.43	2.57	1	19
1:A:63:ILE:O	1:A:64:GLU:C	0.43	2.56	18	8
1:B:26:VAL:CG2	1:B:66:MET:HG3	0.43	2.43	5	16

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:31:PRO:HA	1:B:35:CYS:O	0.43	2.13	20	11
1:B:31:PRO:HB3	1:B:36:ALA:O	0.43	2.13	13	3
1:B:53:ASN:O	1:B:53:ASN:OD1	0.43	2.36	11	1
1:B:29:LYS:NZ	1:B:39:GLU:HG3	0.43	2.28	20	1
1:A:41:ILE:HG13	1:A:51:CYS:SG	0.43	2.54	6	6
1:A:22:ASN:C	1:A:23:ILE:HD13	0.43	2.29	12	1
1:A:23:ILE:HG22	1:A:24:GLN:H	0.43	1.74	14	1
1:B:20:PRO:HA	1:B:23:ILE:CG1	0.43	2.43	14	1
1:A:57:PRO:O	1:A:61:LYS:HG2	0.43	2.12	6	1
1:B:18:ILE:O	1:B:19:HIS:C	0.43	2.57	10	9
1:B:43:THR:OG1	1:B:49:LYS:HD2	0.43	2.13	12	1
1:B:23:ILE:HG22	1:B:24:GLN:H	0.43	1.73	14	1
1:B:61:LYS:O	1:B:65:LYS:HG2	0.43	2.13	14	1
1:A:27:ASN:ND2	1:B:27:ASN:ND2	0.43	2.66	19	1
1:A:61:LYS:HA	1:A:64:GLU:CG	0.43	2.43	19	1
1:A:37:GLN:CD	1:A:37:GLN:N	0.43	2.72	23	1
1:A:56:SER:O	1:A:60:LYS:CG	0.43	2.66	21	3
1:B:9:CYS:HB2	1:B:12:LEU:CD2	0.43	2.43	24	4
1:B:65:LYS:O	1:B:66:MET:C	0.43	2.57	1	9
1:A:15:LEU:CD1	1:A:51:CYS:O	0.43	2.61	22	1
1:B:68:ASN:O	1:B:69:SER:C	0.43	2.57	22	8
1:B:41:ILE:HG13	1:B:51:CYS:SG	0.43	2.54	6	6
1:B:61:LYS:HE2	1:B:61:LYS:CA	0.43	2.43	21	2
1:B:41:ILE:CG2	1:B:42:ALA:N	0.43	2.81	9	1
1:B:17:GLY:HA3	1:B:58:ILE:HG21	0.43	1.91	10	1
1:A:20:PRO:HA	1:A:23:ILE:CG1	0.43	2.43	14	1
1:A:27:ASN:OD1	1:B:27:ASN:CG	0.43	2.57	15	7
1:A:29:LYS:O	1:A:29:LYS:HD3	0.43	2.12	5	1
1:B:16:GLN:O	1:B:58:ILE:HG21	0.43	2.13	19	1
1:A:9:CYS:HB2	1:A:12:LEU:HD23	0.43	1.91	25	1
1:A:20:PRO:CB	1:A:65:LYS:HG2	0.43	2.44	4	2
1:A:27:ASN:CG	1:B:27:ASN:OD1	0.43	2.57	14	7
1:B:19:HIS:O	1:B:23:ILE:HD11	0.43	2.12	9	1
1:A:10:GLN:CB	1:A:29:LYS:HE3	0.43	2.44	11	1
1:B:10:GLN:CB	1:B:29:LYS:HE3	0.43	2.44	11	1
1:B:29:LYS:O	1:B:29:LYS:HG2	0.43	2.14	11	1
1:A:61:LYS:HA	1:A:61:LYS:HE2	0.43	1.91	21	3
1:A:65:LYS:O	1:A:66:MET:C	0.43	2.57	8	9
1:B:62:ILE:HD12	1:B:62:ILE:N	0.43	2.29	12	3
1:A:14:THR:OG1	1:A:53:ASN:CG	0.43	2.57	12	1
1:B:14:THR:OG1	1:B:53:ASN:CG	0.43	2.57	12	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:13:GLN:OE1	1:A:14:THR:N	0.43	2.52	22	1
1:A:17:GLY:C	1:A:18:ILE:HD13	0.43	2.35	23	1
1:B:37:GLN:O	1:B:38:THR:C	0.43	2.57	25	1
1:A:25:SER:OG	1:A:43:THR:HB	0.42	2.14	8	8
1:B:66:MET:O	1:B:69:SER:OG	0.42	2.34	6	2
1:A:17:GLY:HA3	1:A:58:ILE:HG21	0.42	1.91	10	1
1:B:20:PRO:CB	1:B:65:LYS:HG2	0.42	2.44	4	2
1:A:45:LYS:NZ	1:A:45:LYS:HB3	0.42	2.29	24	2
1:A:37:GLN:OE1	1:A:37:GLN:HA	0.42	2.14	7	2
1:B:22:ASN:C	1:B:23:ILE:HD13	0.42	2.29	12	1
1:B:38:THR:O	1:B:39:GLU:CD	0.42	2.57	20	1
1:B:66:MET:O	1:B:67:LEU:C	0.42	2.57	1	9
1:A:66:MET:O	1:A:67:LEU:C	0.42	2.57	2	8
1:B:18:ILE:HG23	1:B:23:ILE:HD11	0.42	1.91	13	1
1:A:68:ASN:O	1:A:69:SER:C	0.42	2.57	16	7
1:B:10:GLN:NE2	1:B:10:GLN:HA	0.42	2.29	4	1
1:A:29:LYS:HB3	1:A:29:LYS:NZ	0.42	2.29	6	1
1:B:61:LYS:HD3	1:B:64:GLU:OE2	0.42	2.15	23	1
1:A:21:LYS:O	1:A:45:LYS:HE3	0.42	2.15	22	7
1:B:18:ILE:O	1:B:62:ILE:HG13	0.42	2.15	4	2
1:B:25:SER:OG	1:B:43:THR:HB	0.42	2.14	8	7
1:B:37:GLN:OE1	1:B:37:GLN:HA	0.42	2.14	7	2
1:A:59:VAL:O	1:A:60:LYS:C	0.42	2.58	12	1
1:B:40:VAL:CG2	1:B:54:PRO:HB3	0.42	2.42	12	1
1:A:10:GLN:HB2	1:A:29:LYS:CE	0.42	2.44	20	1
1:B:45:LYS:HB3	1:B:45:LYS:NZ	0.42	2.29	5	2
1:A:29:LYS:O	1:A:30:SER:C	0.42	2.58	12	1
1:A:36:ALA:O	1:A:37:GLN:CD	0.42	2.58	25	2
1:B:36:ALA:O	1:B:37:GLN:CD	0.42	2.58	25	2
1:B:37:GLN:CD	1:B:37:GLN:N	0.42	2.72	23	1
1:B:21:LYS:O	1:B:45:LYS:HE3	0.42	2.14	2	6
1:B:61:LYS:HA	1:B:61:LYS:HE2	0.42	1.91	2	2
1:A:10:GLN:CG	1:A:29:LYS:HE3	0.42	2.45	5	2
1:B:35:CYS:HB3	1:B:39:GLU:OE2	0.42	2.15	10	1
1:A:29:LYS:O	1:A:29:LYS:HG2	0.42	2.14	11	1
1:A:61:LYS:HE3	1:A:64:GLU:CD	0.42	2.35	11	1
1:B:61:LYS:HE3	1:B:64:GLU:CD	0.42	2.35	11	1
1:A:46:ASN:OD1	1:A:46:ASN:N	0.42	2.53	3	4
1:B:29:LYS:CE	1:B:41:ILE:CD1	0.42	2.98	14	2
1:B:30:SER:CB	1:B:31:PRO:CD	0.42	2.98	19	6
1:A:10:GLN:HG3	1:A:29:LYS:CD	0.42	2.45	21	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:9:CYS:HB2	1:B:12:LEU:HD23	0.42	1.91	25	1
1:A:35:CYS:SG	1:A:39:GLU:CD	0.42	2.98	2	2
1:B:18:ILE:CG1	1:B:52:LEU:HD21	0.42	2.45	9	1
1:A:26:VAL:CG1	1:A:63:ILE:HG13	0.42	2.45	19	1
1:B:13:GLN:OE1	1:B:14:THR:N	0.42	2.52	22	1
1:A:27:ASN:O	1:A:41:ILE:CD1	0.42	2.65	13	2
1:A:30:SER:CB	1:A:31:PRO:CD	0.42	2.98	19	5
1:A:35:CYS:HB3	1:A:39:GLU:OE2	0.42	2.15	10	1
1:A:51:CYS:O	1:A:52:LEU:HD23	0.42	2.15	11	1
1:A:15:LEU:HG	1:A:51:CYS:O	0.42	2.15	25	4
1:A:20:PRO:HG2	1:A:21:LYS:CE	0.42	2.44	15	1
1:B:20:PRO:HG2	1:B:21:LYS:CE	0.42	2.44	15	1
1:B:35:CYS:CB	1:B:39:GLU:OE2	0.42	2.68	18	1
1:B:10:GLN:HB2	1:B:29:LYS:CE	0.42	2.44	20	1
1:B:31:PRO:HB3	1:B:36:ALA:C	0.41	2.35	5	2
1:A:61:LYS:HE3	1:A:64:GLU:OE2	0.41	2.15	11	1
1:A:15:LEU:O	1:A:17:GLY:N	0.41	2.54	9	7
1:B:35:CYS:SG	1:B:39:GLU:CD	0.41	2.98	2	2
1:A:29:LYS:CE	1:A:41:ILE:CD1	0.41	2.98	14	2
1:A:61:LYS:O	1:A:65:LYS:HG3	0.41	2.15	11	2
1:B:28:VAL:CG1	1:B:40:VAL:HG22	0.41	2.41	11	1
1:B:51:CYS:O	1:B:52:LEU:HD23	0.41	2.15	11	1
1:A:40:VAL:HG23	1:A:54:PRO:CG	0.41	2.45	12	1
1:A:18:ILE:HG23	1:A:23:ILE:HD11	0.41	1.91	13	1
1:B:26:VAL:CG1	1:B:63:ILE:HG13	0.41	2.45	19	1
1:A:37:GLN:O	1:A:39:GLU:CD	0.41	2.58	21	1
1:A:61:LYS:HD3	1:A:64:GLU:OE2	0.41	2.15	23	1
1:B:10:GLN:HG3	1:B:29:LYS:CB	0.41	2.45	3	1
1:A:31:PRO:HB3	1:A:36:ALA:C	0.41	2.35	5	2
1:A:38:THR:O	1:A:39:GLU:CD	0.41	2.57	20	1
1:B:10:GLN:HG3	1:B:29:LYS:CD	0.41	2.45	21	1
1:B:15:LEU:O	1:B:17:GLY:N	0.41	2.54	9	7
1:A:62:ILE:HD12	1:A:62:ILE:N	0.41	2.29	12	1
1:B:15:LEU:HG	1:B:51:CYS:O	0.41	2.15	25	4
1:B:37:GLN:O	1:B:39:GLU:CD	0.41	2.58	21	1
1:A:10:GLN:NE2	1:A:10:GLN:HA	0.41	2.29	4	1
1:A:29:LYS:HE3	1:B:24:GLN:OE1	0.41	2.15	8	1
1:B:59:VAL:O	1:B:60:LYS:C	0.41	2.58	12	1
1:A:37:GLN:O	1:A:38:THR:C	0.41	2.57	25	1
1:A:24:GLN:HB2	1:A:44:LEU:O	0.41	2.16	2	2
1:A:10:GLN:HG3	1:A:29:LYS:CB	0.41	2.45	3	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:21:LYS:HD2	1:B:22:ASN:OD1	0.41	2.15	16	1
1:A:35:CYS:CB	1:A:39:GLU:OE2	0.41	2.68	18	1
1:B:10:GLN:NE2	1:B:29:LYS:HG2	0.41	2.30	23	1
1:B:46:ASN:OD1	1:B:46:ASN:N	0.41	2.53	3	2
1:B:20:PRO:HG3	1:B:65:LYS:HD3	0.41	1.93	3	1
1:A:44:LEU:HB3	1:A:46:ASN:OD1	0.41	2.16	14	5
1:A:25:SER:HA	1:B:28:VAL:O	0.41	2.16	12	1
1:B:28:VAL:CG2	1:B:40:VAL:HG13	0.41	2.39	17	1
1:B:15:LEU:CD1	1:B:51:CYS:O	0.41	2.61	22	1
1:A:10:GLN:NE2	1:A:29:LYS:HG2	0.41	2.30	23	1
1:B:20:PRO:CB	1:B:65:LYS:HB3	0.41	2.46	23	1
1:B:54:PRO:O	1:B:60:LYS:HE2	0.41	2.16	9	1
1:A:9:CYS:HA	1:A:39:GLU:OE1	0.41	2.16	12	1
1:A:61:LYS:CD	1:A:64:GLU:OE2	0.41	2.69	23	1
1:B:61:LYS:CD	1:B:64:GLU:OE2	0.41	2.69	23	1
1:A:18:ILE:O	1:A:62:ILE:HG13	0.41	2.15	4	1
1:B:10:GLN:CG	1:B:29:LYS:HE3	0.41	2.46	4	2
1:B:18:ILE:O	1:B:18:ILE:HD12	0.41	2.16	4	1
1:B:10:GLN:HA	1:B:10:GLN:OE1	0.41	2.16	5	1
1:B:23:ILE:O	1:B:45:LYS:NZ	0.41	2.52	5	1
1:A:24:GLN:CG	1:A:43:THR:CG2	0.41	2.99	12	2
1:B:44:LEU:HB3	1:B:46:ASN:OD1	0.41	2.16	8	5
1:A:18:ILE:CG1	1:A:52:LEU:HD21	0.41	2.45	9	1
1:B:64:GLU:O	1:B:65:LYS:C	0.41	2.59	11	1
1:B:40:VAL:HG23	1:B:54:PRO:CG	0.41	2.45	12	1
1:B:19:HIS:CD2	1:B:21:LYS:HG2	0.41	2.50	15	1
1:A:21:LYS:HD2	1:A:22:ASN:OD1	0.41	2.15	16	1
1:B:29:LYS:HG3	1:B:29:LYS:O	0.41	2.16	20	1
1:A:10:GLN:CG	1:A:29:LYS:NZ	0.41	2.84	22	1
1:A:21:LYS:O	1:A:45:LYS:HE2	0.41	2.16	5	2
1:B:44:LEU:N	1:B:48:ARG:O	0.41	2.54	11	1
1:A:28:VAL:O	1:B:25:SER:HA	0.41	2.16	12	1
1:B:66:MET:O	1:B:69:SER:HB3	0.41	2.16	12	1
1:B:24:GLN:HB2	1:B:44:LEU:O	0.41	2.16	16	1
1:A:39:GLU:HA	1:A:39:GLU:OE1	0.41	2.16	20	1
1:A:24:GLN:OE1	1:B:29:LYS:HE3	0.40	2.15	8	1
1:B:61:LYS:HE3	1:B:64:GLU:OE2	0.40	2.15	11	1
1:B:29:LYS:O	1:B:30:SER:C	0.40	2.58	12	2
1:B:18:ILE:HD11	1:B:23:ILE:HD11	0.40	1.84	20	3
1:B:29:LYS:HB3	1:B:29:LYS:NZ	0.40	2.29	6	1
1:B:61:LYS:O	1:B:65:LYS:HG3	0.40	2.15	6	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:9:CYS:HA	1:B:39:GLU:OE1	0.40	2.16	12	1
1:A:54:PRO:HB2	1:A:60:LYS:NZ	0.40	2.32	16	1
1:B:10:GLN:CG	1:B:29:LYS:NZ	0.40	2.84	22	1
1:B:24:GLN:CG	1:B:43:THR:CG2	0.40	2.99	6	2
1:A:10:GLN:HB2	1:A:29:LYS:HD2	0.40	1.93	7	1
1:A:30:SER:CB	1:A:31:PRO:HD2	0.40	2.46	7	1
1:B:30:SER:CB	1:B:31:PRO:HD2	0.40	2.46	7	1
1:A:66:MET:O	1:A:69:SER:HB3	0.40	2.16	12	1
1:A:63:ILE:CD1	1:A:63:ILE:N	0.40	2.84	14	1
1:A:19:HIS:CD2	1:A:21:LYS:HG2	0.40	2.50	15	1
1:B:10:GLN:HB2	1:B:39:GLU:HB2	0.40	1.93	23	1
1:A:18:ILE:O	1:A:18:ILE:HD12	0.40	2.16	4	1
1:B:62:ILE:HG22	1:B:63:ILE:HD12	0.40	1.93	8	1
1:A:44:LEU:N	1:A:48:ARG:O	0.40	2.54	11	1
1:A:10:GLN:OE1	1:A:10:GLN:HA	0.40	2.16	14	1
1:A:20:PRO:HG3	1:A:65:LYS:HD3	0.40	1.93	3	1
1:A:10:GLN:HA	1:A:10:GLN:OE1	0.40	2.16	5	1
1:A:15:LEU:HD12	1:A:52:LEU:HD23	0.40	1.92	12	1
1:B:54:PRO:HB2	1:B:60:LYS:NZ	0.40	2.32	16	1
1:B:39:GLU:HA	1:B:39:GLU:OE1	0.40	2.16	20	1
1:A:10:GLN:CB	1:A:39:GLU:HB2	0.40	2.46	23	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	61/73 (84%)	49±2 (80±3%)	10±2 (17±3%)	2±1 (3±1%)	8	44
1	B	61/73 (84%)	49±2 (80±3%)	10±2 (17±3%)	2±1 (3±1%)	8	44
All	All	3050/3650 (84%)	2454 (80%)	518 (17%)	78 (3%)	8	44

All 10 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	44	LEU	24
1	B	44	LEU	24
1	A	9	CYS	8
1	B	9	CYS	8
1	A	30	SER	3
1	B	30	SER	3
1	A	26	VAL	2
1	B	26	VAL	2
1	A	34	HIS	2
1	B	34	HIS	2

### 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	54/64 (84%)	40±2 (74±3%)	14±2 (26±3%)	2	23
1	B	54/64 (84%)	40±2 (74±3%)	14±2 (26±3%)	2	23
All	All	2700/3200 (84%)	1995 (74%)	705 (26%)	2	23

All 66 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	24	GLN	25
1	B	24	GLN	25
1	B	14	THR	24
1	A	14	THR	23
1	A	56	SER	23
1	B	56	SER	23
1	A	12	LEU	21
1	B	12	LEU	21
1	A	66	MET	21
1	B	66	MET	21
1	A	46	ASN	20
1	A	67	LEU	20
1	B	46	ASN	20
1	B	67	LEU	20
1	A	64	GLU	19

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Mol	Chain	Res	Type	Models (Total)
1	B	64	GLU	19
1	A	19	HIS	18
1	B	19	HIS	18
1	A	61	LYS	15
1	B	61	LYS	15
1	A	49	LYS	14
1	B	49	LYS	14
1	A	30	SER	14
1	B	30	SER	14
1	A	29	LYS	13
1	B	29	LYS	13
1	A	21	LYS	12
1	B	21	LYS	12
1	A	10	GLN	11
1	B	10	GLN	11
1	A	16	GLN	9
1	A	45	LYS	9
1	B	16	GLN	9
1	B	45	LYS	9
1	A	25	SER	9
1	B	25	SER	9
1	A	60	LYS	8
1	B	60	LYS	8
1	A	65	LYS	8
1	B	65	LYS	8
1	A	13	GLN	7
1	A	53	ASN	7
1	B	13	GLN	7
1	B	53	ASN	7
1	A	39	GLU	5
1	B	39	GLU	5
1	A	18	ILE	5
1	B	18	ILE	5
1	A	58	ILE	3
1	B	58	ILE	3
1	A	48	ARG	2
1	B	48	ARG	2
1	A	68	ASN	2
1	B	68	ASN	2
1	A	27	ASN	2
1	B	27	ASN	2
1	A	37	GLN	2

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Mol	Chain	Res	Type	Models (Total)
1	B	37	GLN	2
1	A	15	LEU	2
1	B	15	LEU	2
1	A	43	THR	1
1	B	43	THR	1
1	A	23	ILE	1
1	B	23	ILE	1
1	A	38	THR	1
1	B	38	THR	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 monosaccharides 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