



# Full wwPDB NMR Structure Validation Report ⓘ

Apr 20, 2024 – 09:03 AM EDT

PDB ID : 2N64  
BMRB ID : 25750  
Title : NMR Structure of the C-terminal Coiled-Coil Domain of CIN85  
Authors : Habeck, M.; Becker, S.; Griesinger, C.; Wong, L.E.  
Deposited on : 2015-08-11

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.

The types of validation reports are described at

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

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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)  
wwPDB-RCI : v\_1n\_11\_5\_13\_A (Berjanski et al., 2005)  
PANAV : Wang et al. (2010)  
wwPDB-ShiftChecker : v1.2  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36.2

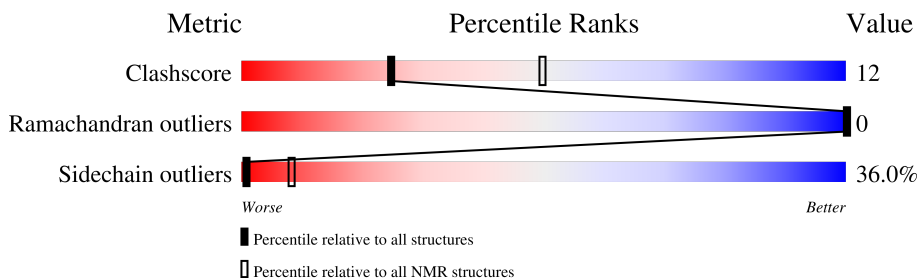
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*SOLUTION NMR*

The overall completeness of chemical shifts assignment is 20%.

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	75	
1	B	75	
1	C	75	

## 2 Ensemble composition and analysis i

This entry contains 30 models. Model 20 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *closest to the average*.

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:605-A:664, B:605-B:664, C:605-C:665 (181)	0.73	20

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 6 clusters and 1 single-model cluster was found.

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

### 3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3771 atoms, of which 1944 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called SH3 domain-containing kinase-binding protein 1.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
1	A	75	1257	373	648	113	119	4	0
1	B	75	1257	373	648	113	119	4	0
1	C	75	1257	373	648	113	119	4	0

There are 9 discrepancies between the modelled and reference sequences:

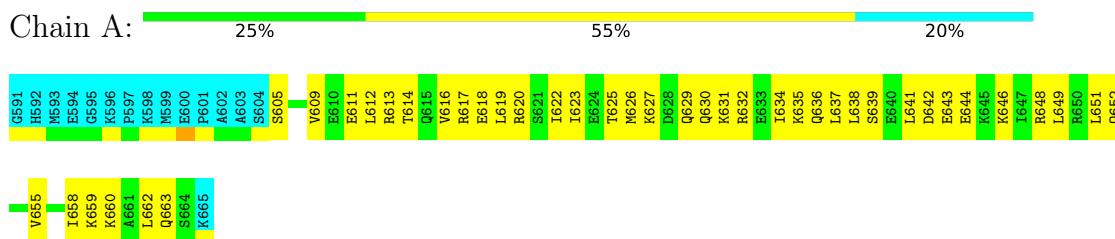
Chain	Residue	Modelled	Actual	Comment	Reference
A	591	GLY	-	expression tag	UNP Q96B97
A	592	HIS	-	expression tag	UNP Q96B97
A	593	MET	-	expression tag	UNP Q96B97
B	591	GLY	-	expression tag	UNP Q96B97
B	592	HIS	-	expression tag	UNP Q96B97
B	593	MET	-	expression tag	UNP Q96B97
C	591	GLY	-	expression tag	UNP Q96B97
C	592	HIS	-	expression tag	UNP Q96B97
C	593	MET	-	expression tag	UNP Q96B97

## 4 Residue-property plots

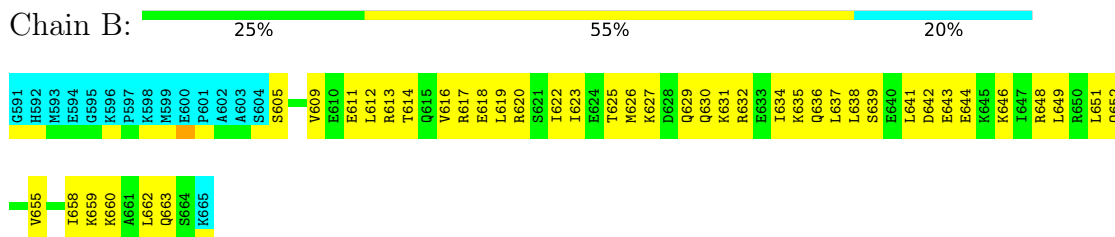
### 4.1 Average score per residue in the NMR ensemble

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

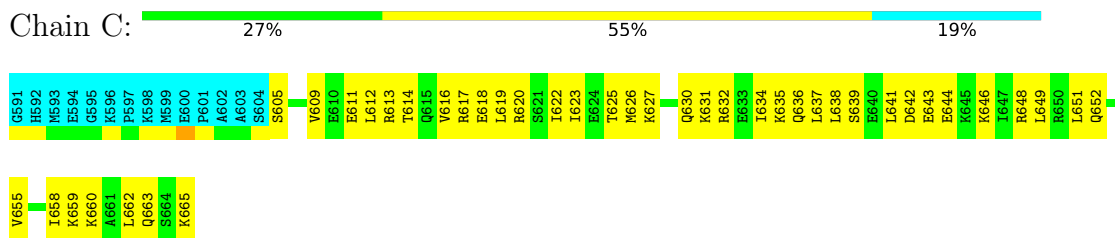
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

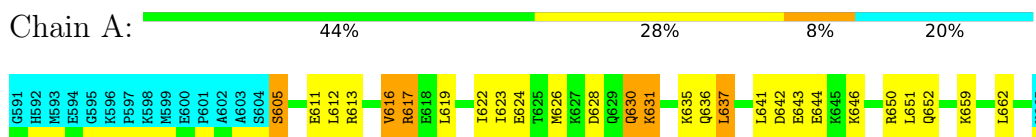


### 4.2 Scores per residue for each member of the ensemble

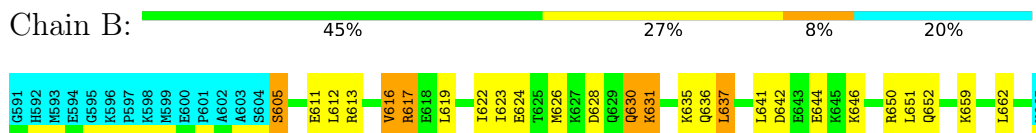
Colouring as in section 4.1 above.

### 4.2.1 Score per residue for model 1

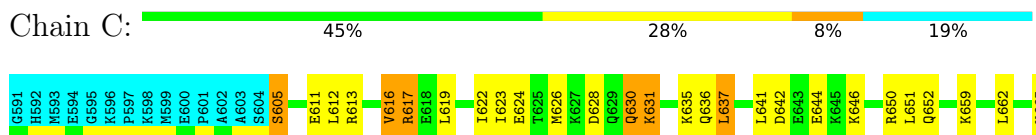
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

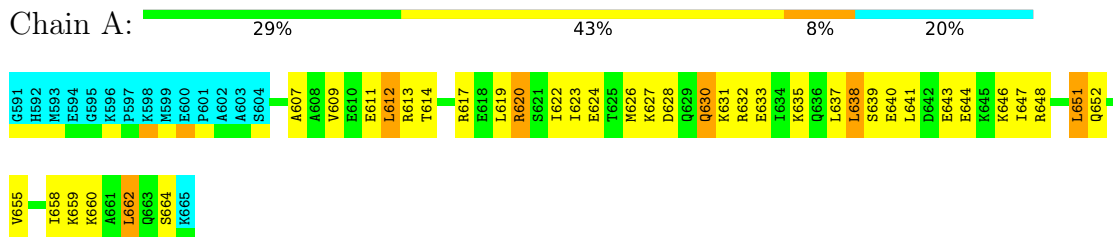


- Molecule 1: SH3 domain-containing kinase-binding protein 1

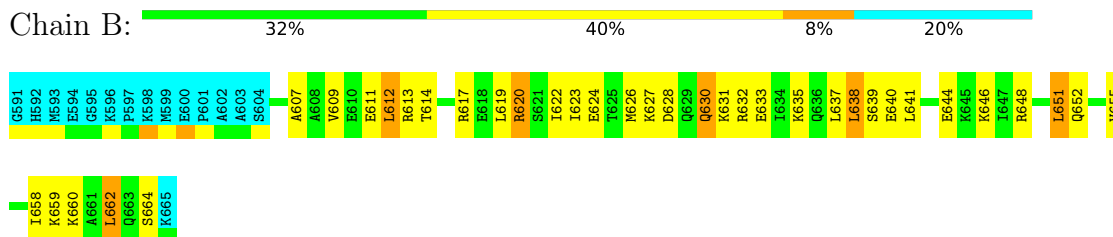


### 4.2.2 Score per residue for model 2

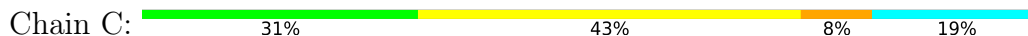
- Molecule 1: SH3 domain-containing kinase-binding protein 1

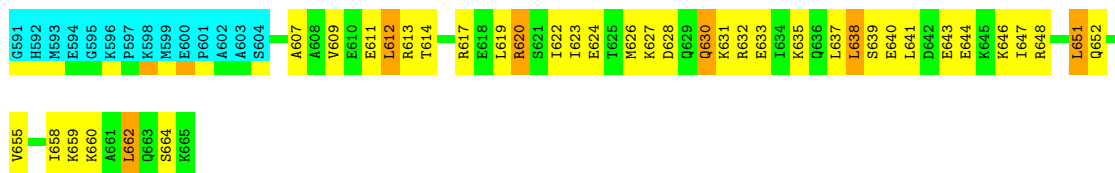


- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

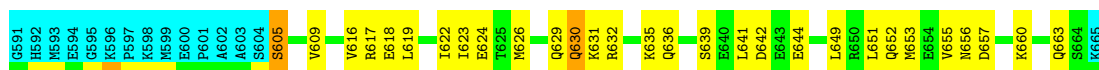




### 4.2.3 Score per residue for model 3

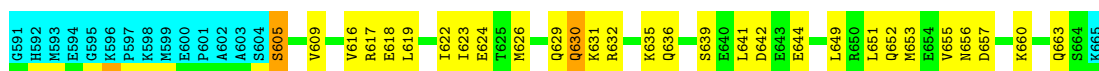
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain A: 41% 36% 0% 20%



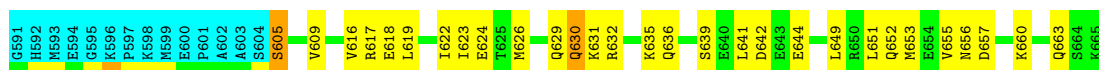
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain B: 41% 36% 0% 20%



- Molecule 1: SH3 domain-containing kinase-binding protein 1

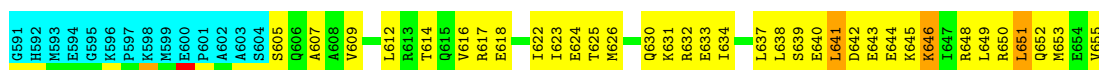
Chain C: 43% 36% 0% 19%



### 4.2.4 Score per residue for model 4

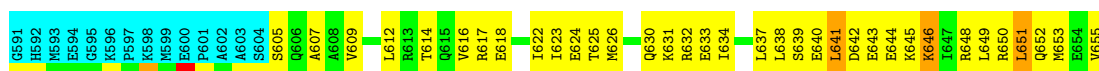
- Molecule 1: SH3 domain-containing kinase-binding protein 1

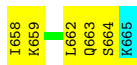
Chain A: 27% 49% 0% 20%



- Molecule 1: SH3 domain-containing kinase-binding protein 1

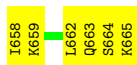
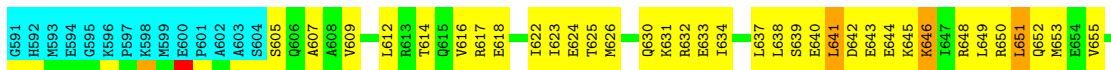
Chain B: 27% 49% 0% 20%





- Molecule 1: SH3 domain-containing kinase-binding protein 1

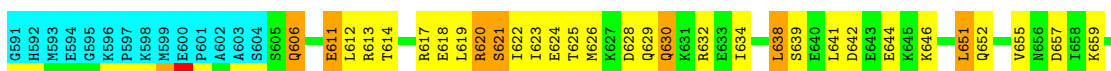
Chain C: 27% 51% 19%



#### 4.2.5 Score per residue for model 5

- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain A: 39% 32% 9% 20%



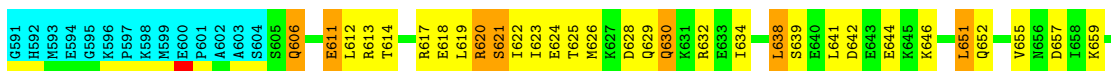
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain B: 39% 32% 9% 20%



- Molecule 1: SH3 domain-containing kinase-binding protein 1

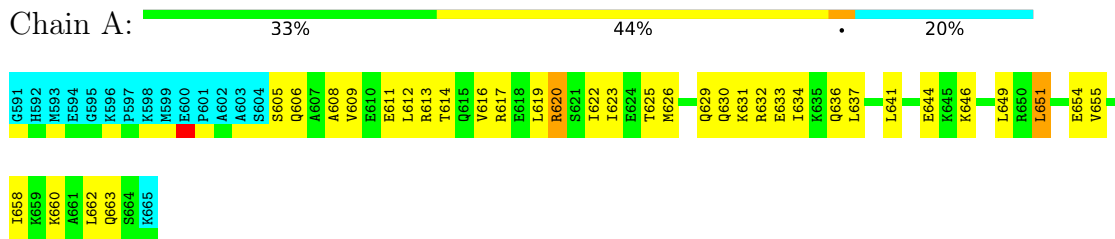
Chain C: 39% 33% 9% 19%



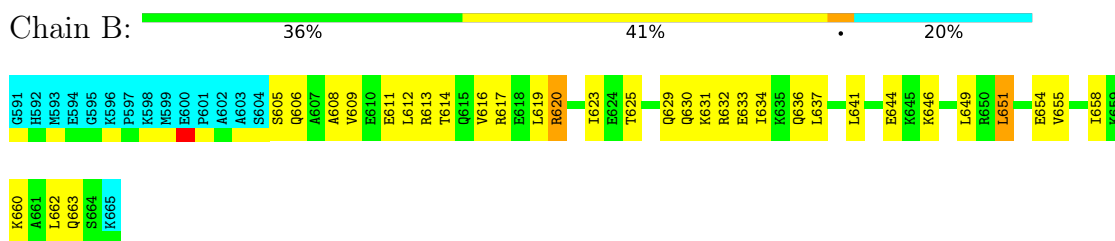
#### 4.2.6 Score per residue for model 6

- Molecule 1: SH3 domain-containing kinase-binding protein 1

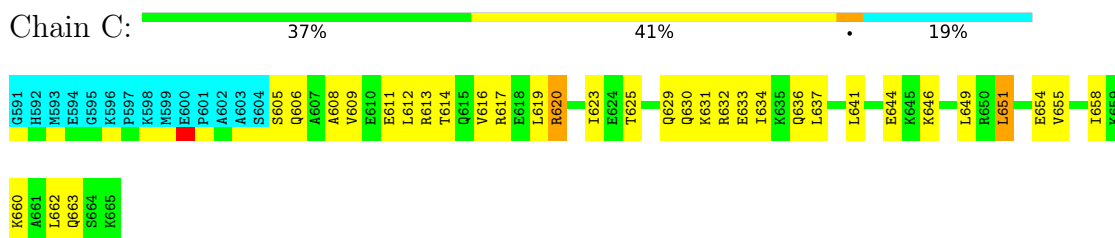




- Molecule 1: SH3 domain-containing kinase-binding protein 1

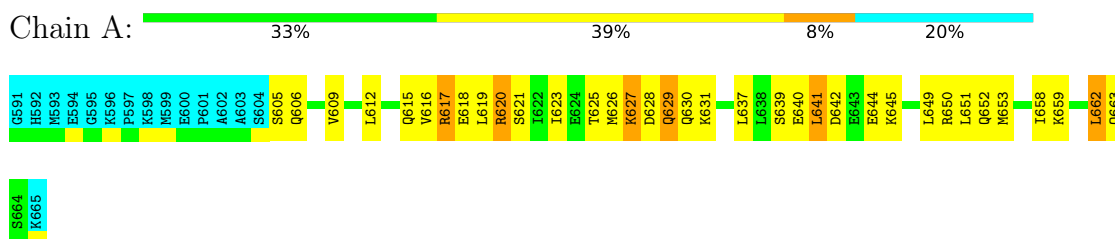


- Molecule 1: SH3 domain-containing kinase-binding protein 1

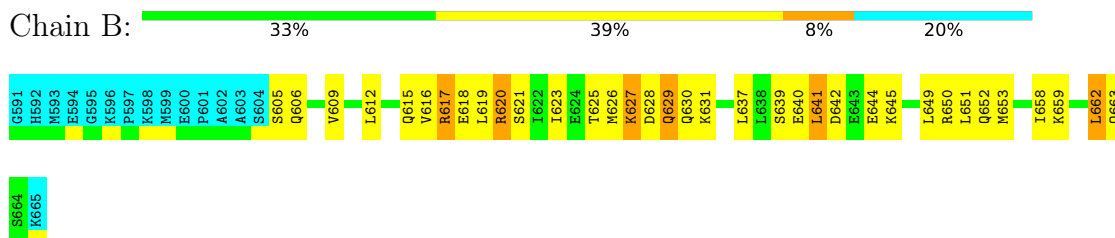


#### 4.2.7 Score per residue for model 7

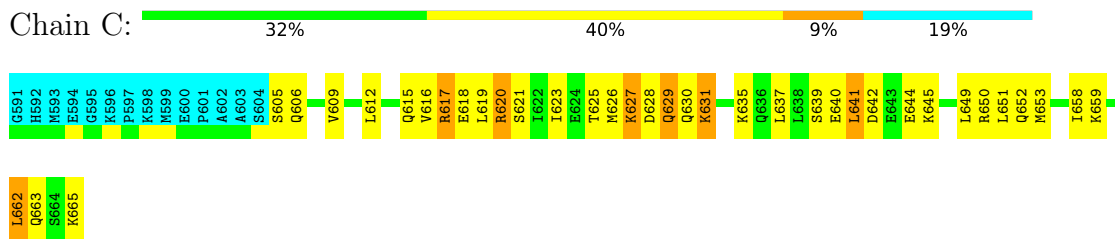
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

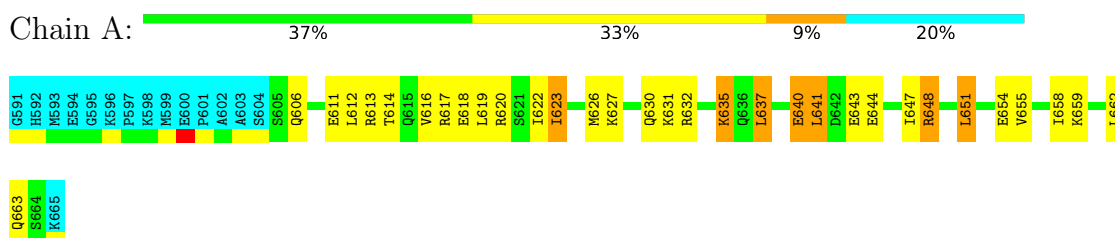


- Molecule 1: SH3 domain-containing kinase-binding protein 1

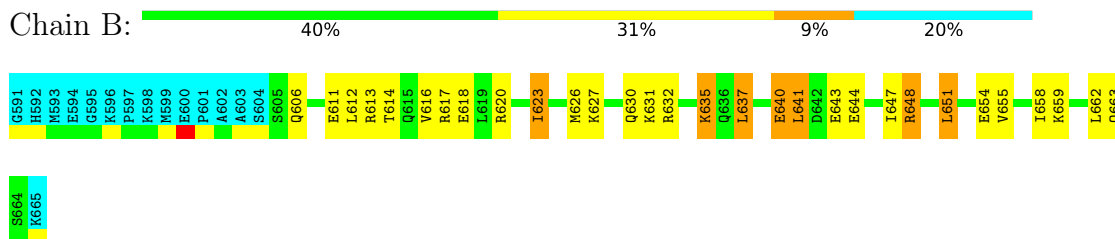


#### 4.2.8 Score per residue for model 8

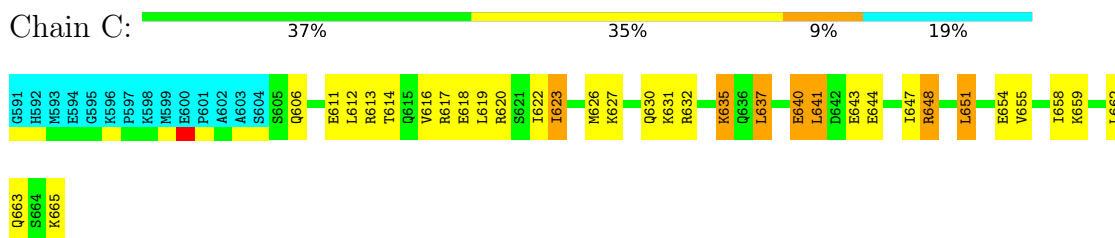
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1



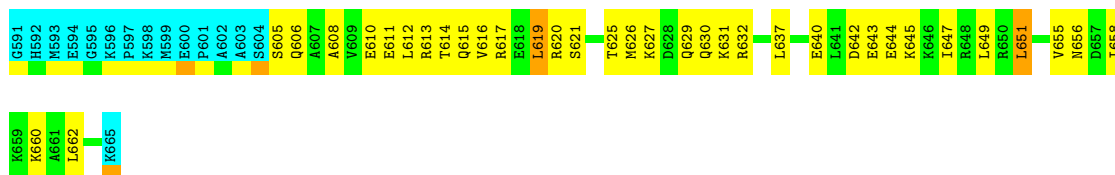
- Molecule 1: SH3 domain-containing kinase-binding protein 1



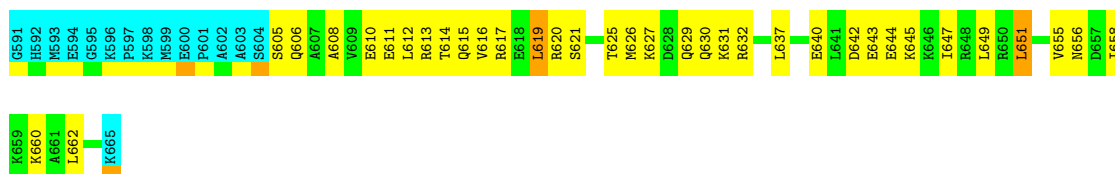
#### 4.2.9 Score per residue for model 9

- Molecule 1: SH3 domain-containing kinase-binding protein 1

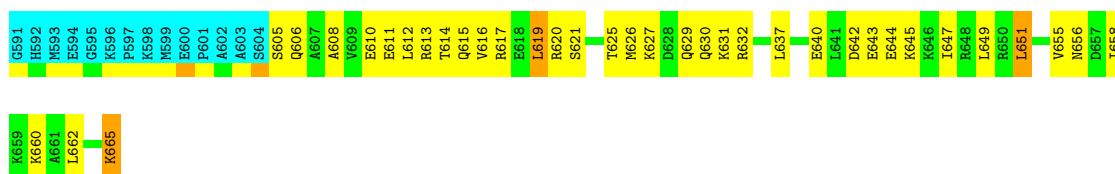




- Molecule 1: SH3 domain-containing kinase-binding protein 1

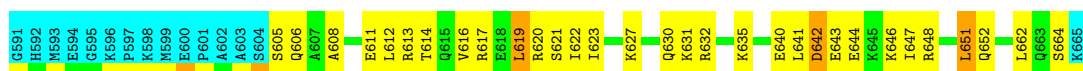


- Molecule 1: SH3 domain-containing kinase-binding protein 1

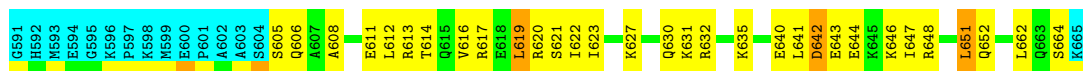
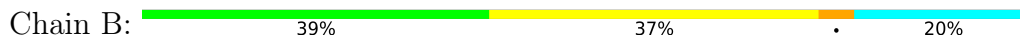


#### 4.2.10 Score per residue for model 10

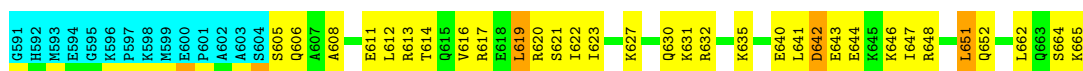
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

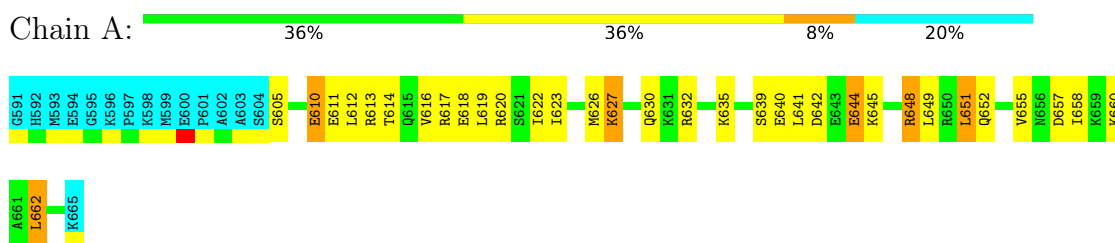


- Molecule 1: SH3 domain-containing kinase-binding protein 1

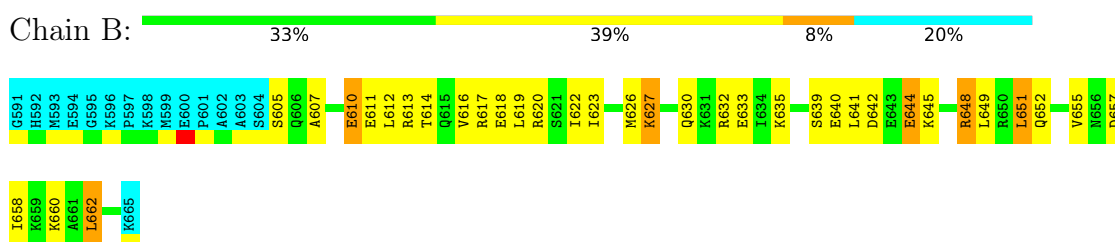


### 4.2.11 Score per residue for model 11

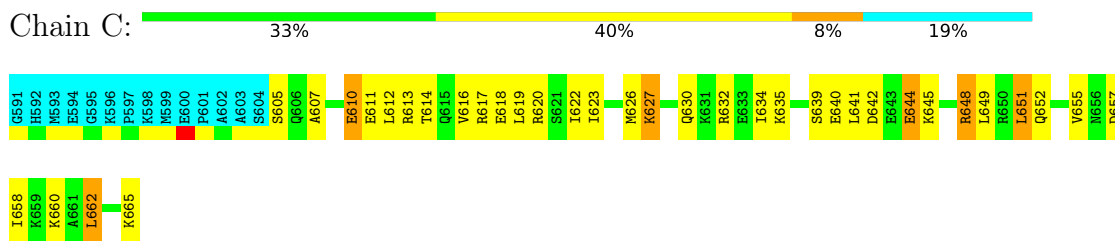
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

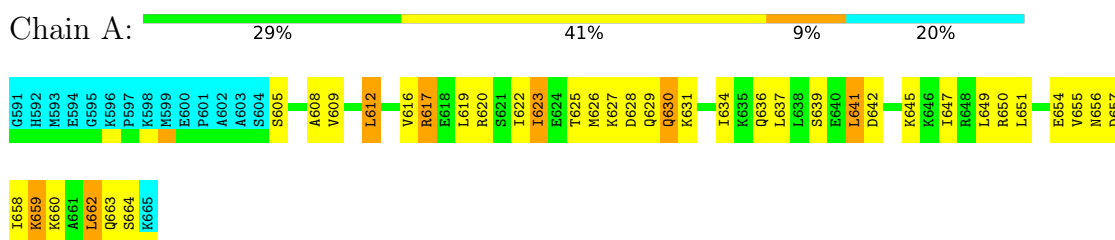


- Molecule 1: SH3 domain-containing kinase-binding protein 1



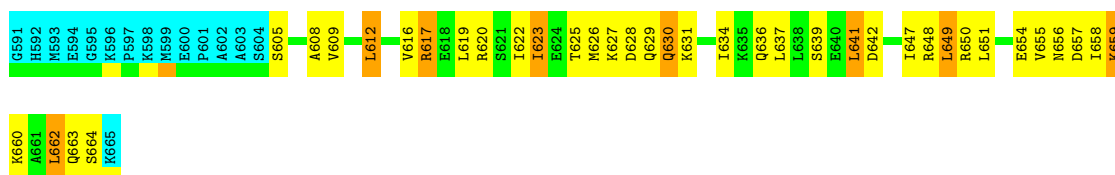
### 4.2.12 Score per residue for model 12

- Molecule 1: SH3 domain-containing kinase-binding protein 1



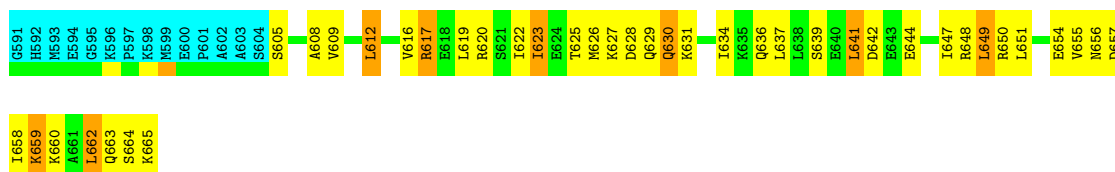
- Molecule 1: SH3 domain-containing kinase-binding protein 1





- Molecule 1: SH3 domain-containing kinase-binding protein 1

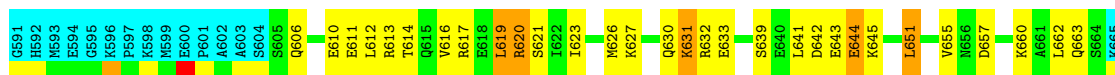
Chain C: 28% 43% 11% 19%



#### 4.2.13 Score per residue for model 13

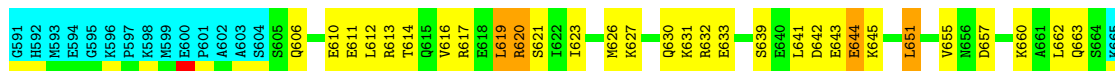
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain A: 40% 33% 7% 20%



- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain B: 40% 35% 5% 20%



- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain C: 43% 33% 5% 19%

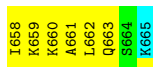


#### 4.2.14 Score per residue for model 14

- Molecule 1: SH3 domain-containing kinase-binding protein 1

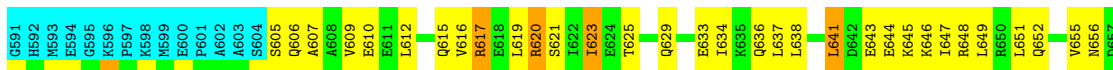
Chain A: 29% 45% 5% 20%





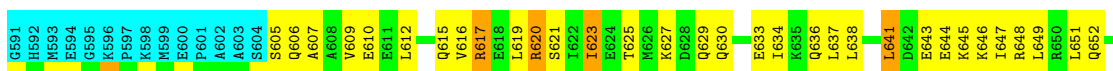
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain B: 29% 45% 5% 20%



- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain C: 27% 49% 5% 19%



#### 4.2.15 Score per residue for model 15

- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain A: 43% 32% 5% 20%



- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain B: 43% 32% 5% 20%



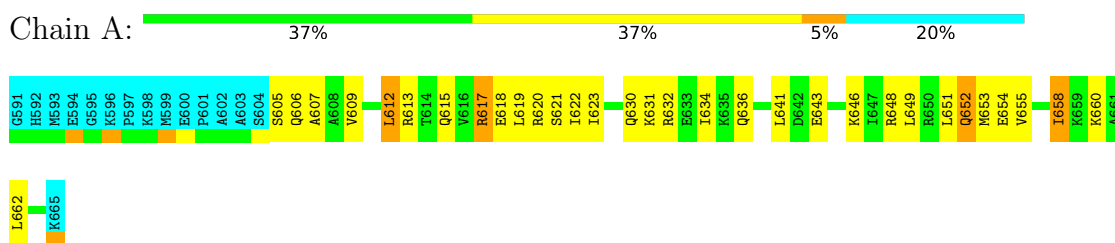
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain C: 44% 32% 5% 19%

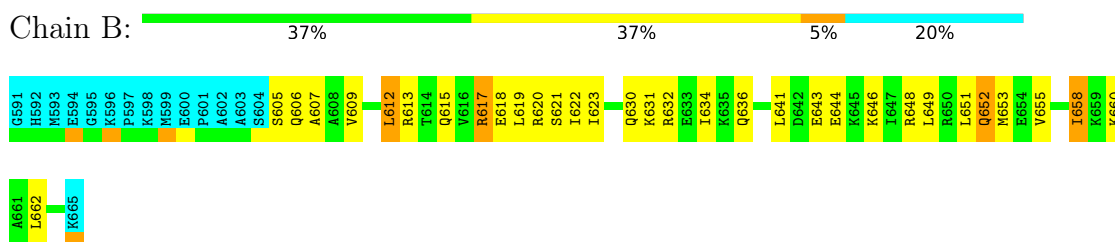


### 4.2.16 Score per residue for model 16

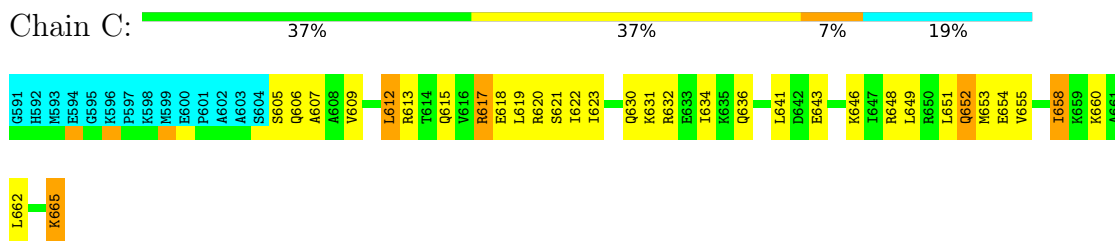
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

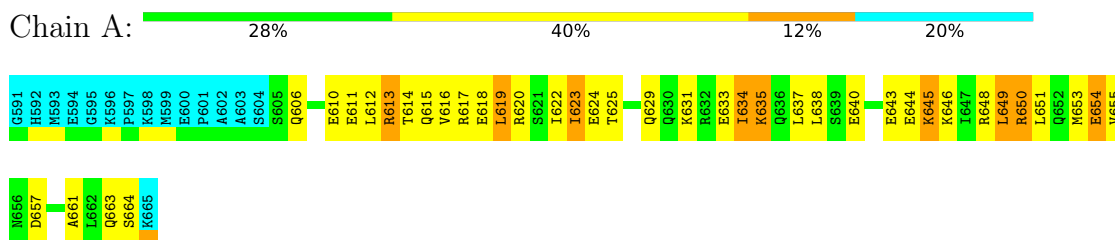


- Molecule 1: SH3 domain-containing kinase-binding protein 1



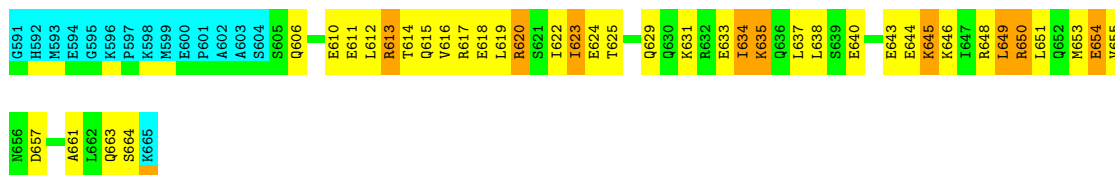
### 4.2.17 Score per residue for model 17

- Molecule 1: SH3 domain-containing kinase-binding protein 1



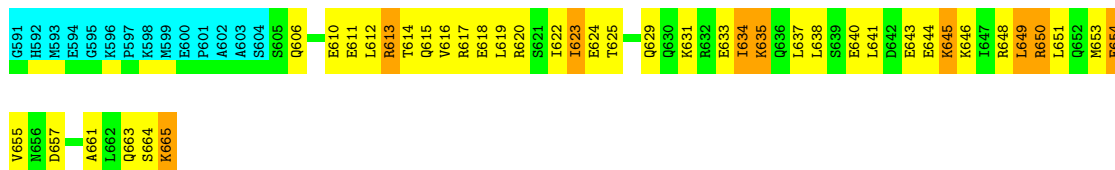
- Molecule 1: SH3 domain-containing kinase-binding protein 1





- Molecule 1: SH3 domain-containing kinase-binding protein 1

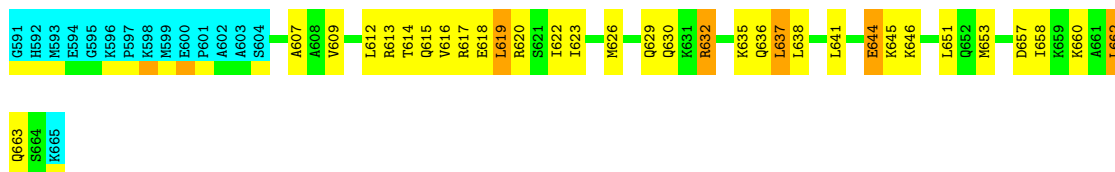
Chain C: 27% 43% 12% 19%



#### 4.2.18 Score per residue for model 18

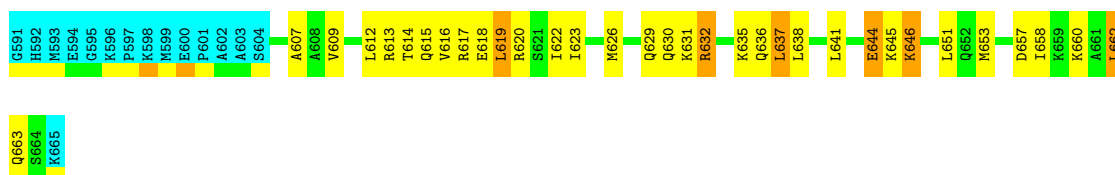
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain A: 37% 36% 7% 20%



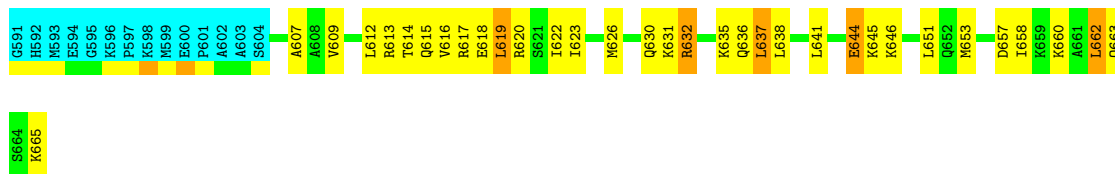
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain B: 36% 36% 8% 20%



- Molecule 1: SH3 domain-containing kinase-binding protein 1

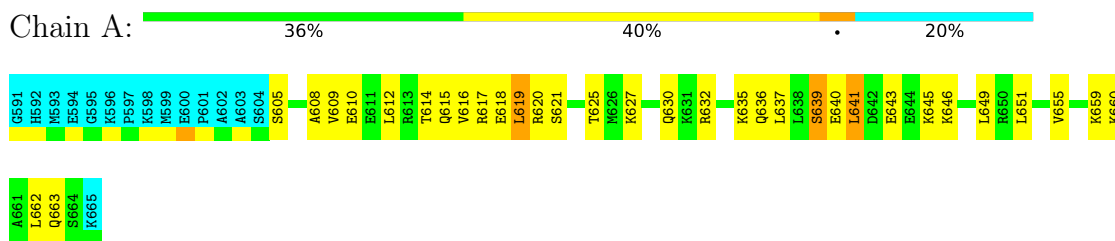
Chain C: 37% 37% 7% 19%



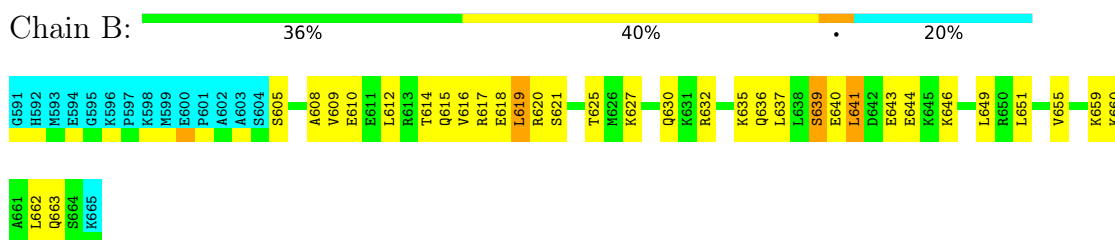


### 4.2.19 Score per residue for model 19

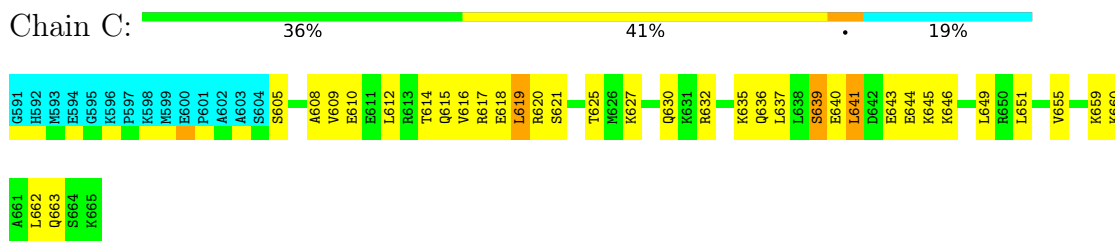
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

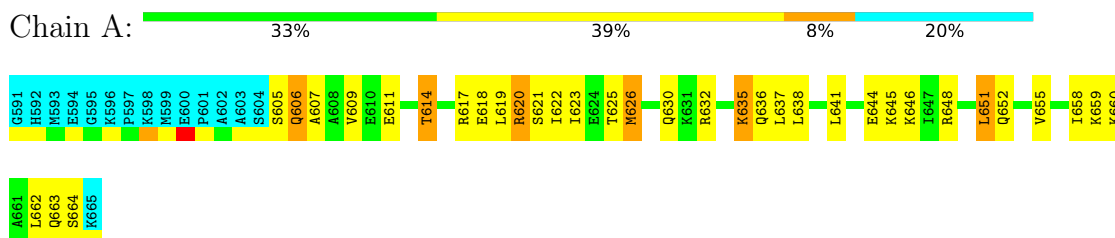


- Molecule 1: SH3 domain-containing kinase-binding protein 1



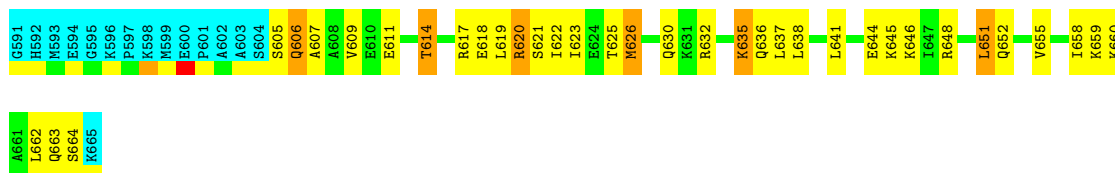
### 4.2.20 Score per residue for model 20 (medoid)

- Molecule 1: SH3 domain-containing kinase-binding protein 1



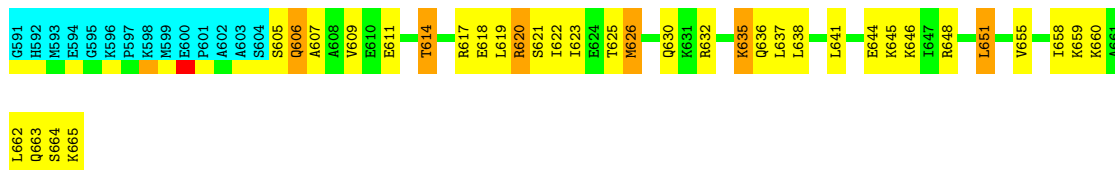
- Molecule 1: SH3 domain-containing kinase-binding protein 1





- Molecule 1: SH3 domain-containing kinase-binding protein 1

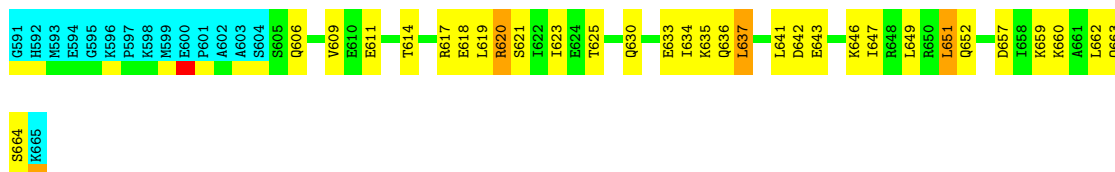
Chain C: 35% 39% 8% 19%



#### 4.2.21 Score per residue for model 21

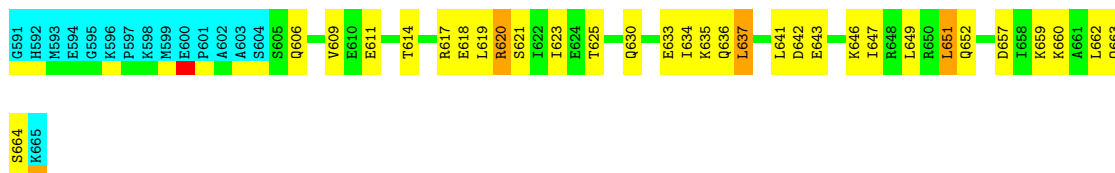
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain A: 39% 37% 20%



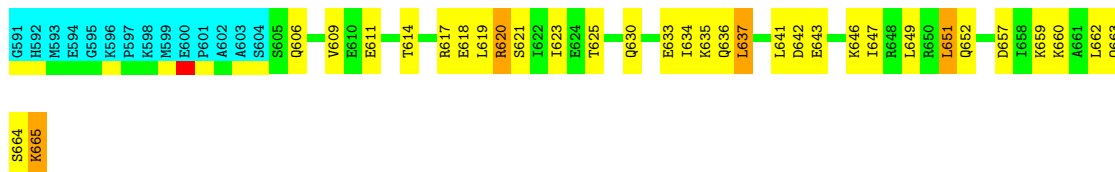
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain B: 39% 37% 20%



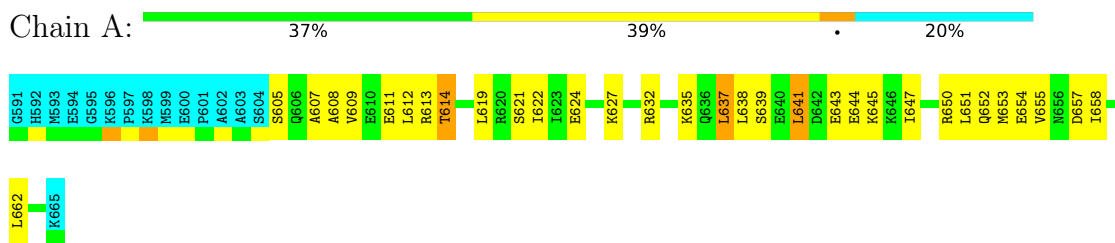
- Molecule 1: SH3 domain-containing kinase-binding protein 1

Chain C: 39% 37% 5% 19%

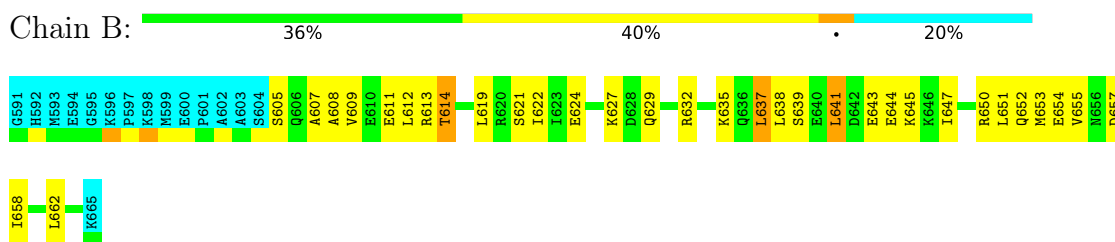


#### 4.2.22 Score per residue for model 22

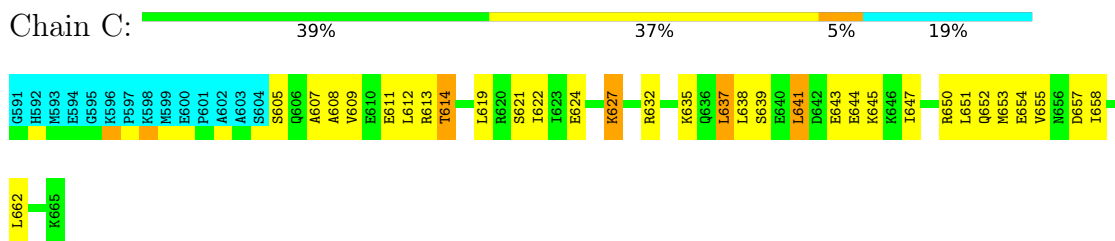
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

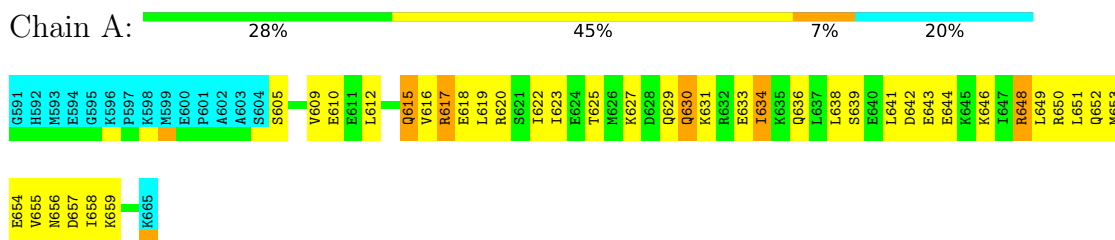


- Molecule 1: SH3 domain-containing kinase-binding protein 1



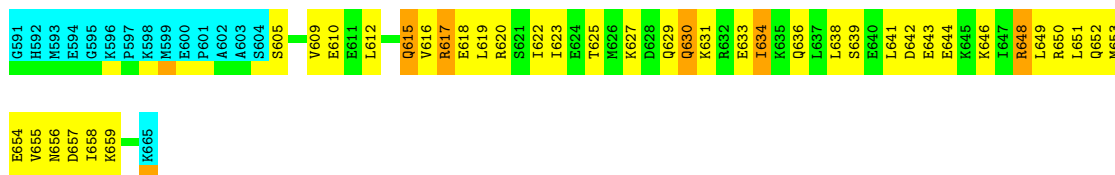
#### 4.2.23 Score per residue for model 23

- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1



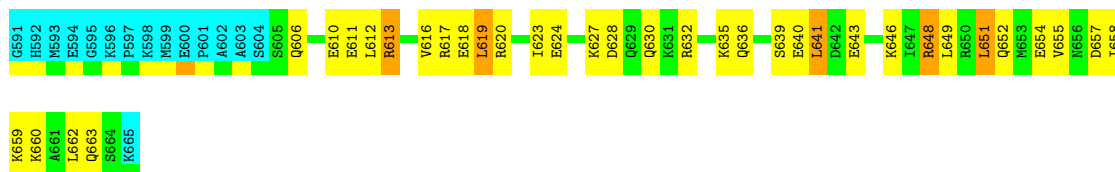


- Molecule 1: SH3 domain-containing kinase-binding protein 1

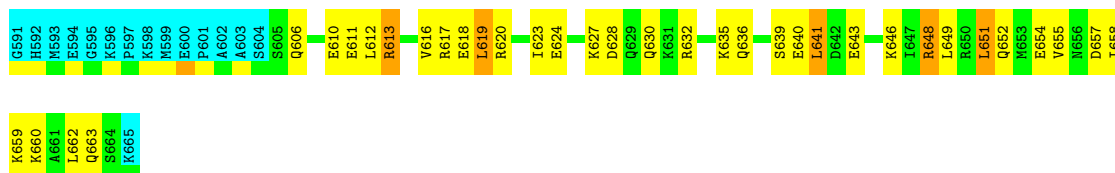
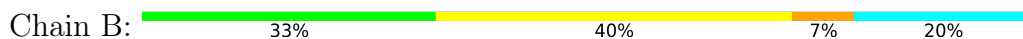


#### 4.2.24 Score per residue for model 24

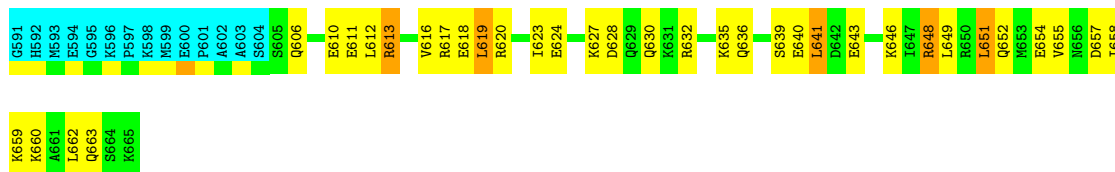
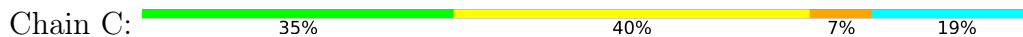
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

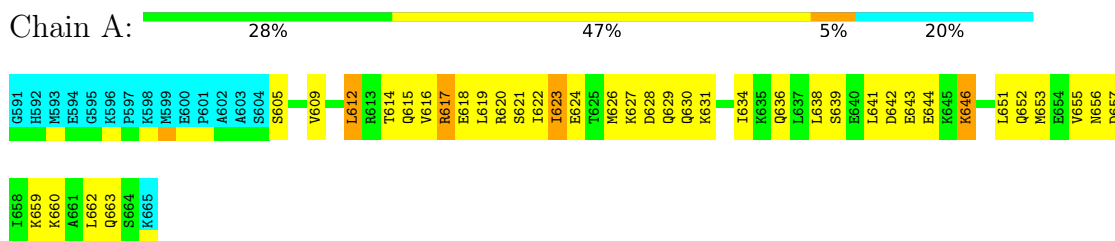


- Molecule 1: SH3 domain-containing kinase-binding protein 1

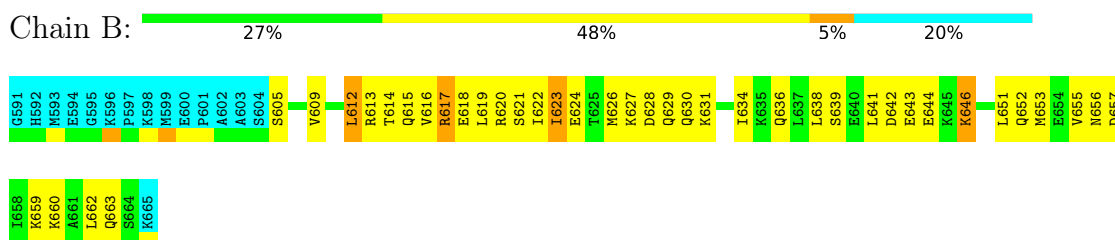


### 4.2.25 Score per residue for model 25

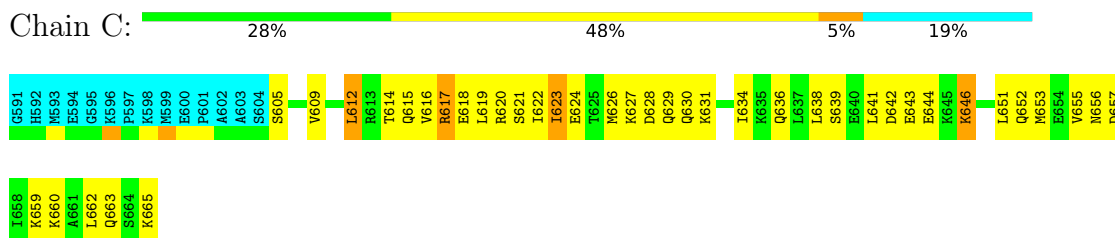
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

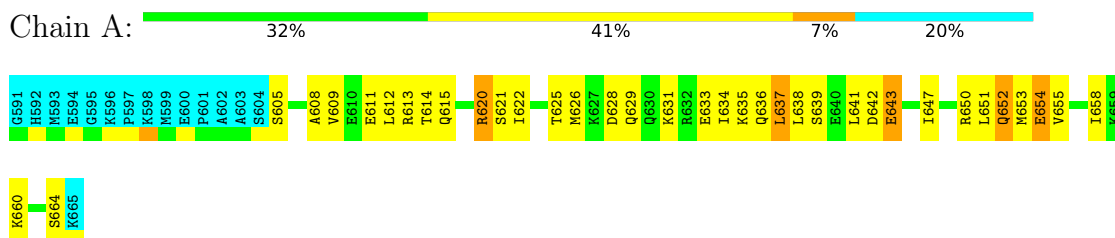


- Molecule 1: SH3 domain-containing kinase-binding protein 1



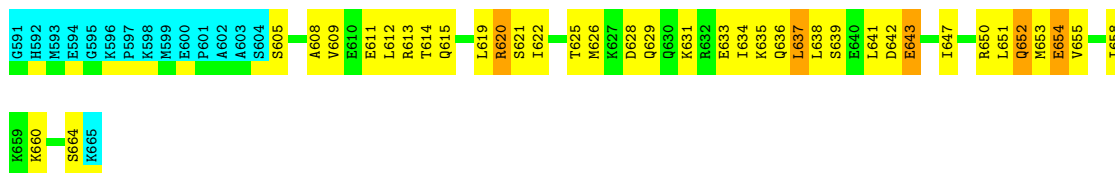
### 4.2.26 Score per residue for model 26

- Molecule 1: SH3 domain-containing kinase-binding protein 1

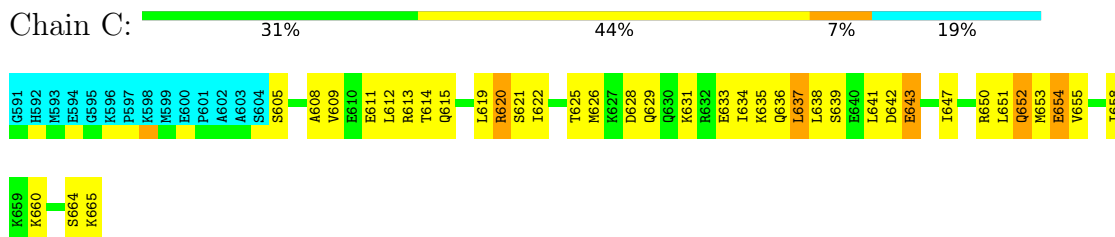


- Molecule 1: SH3 domain-containing kinase-binding protein 1



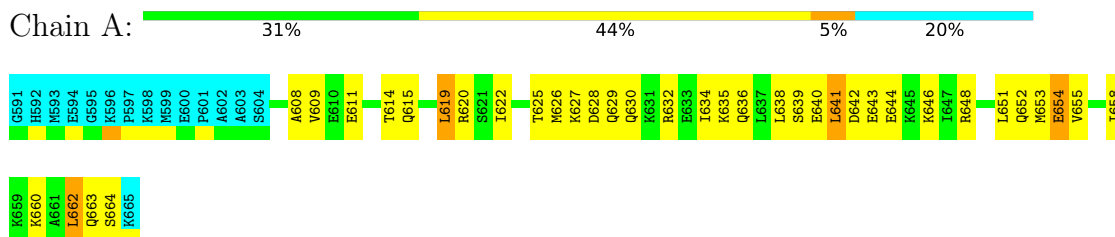


- Molecule 1: SH3 domain-containing kinase-binding protein 1

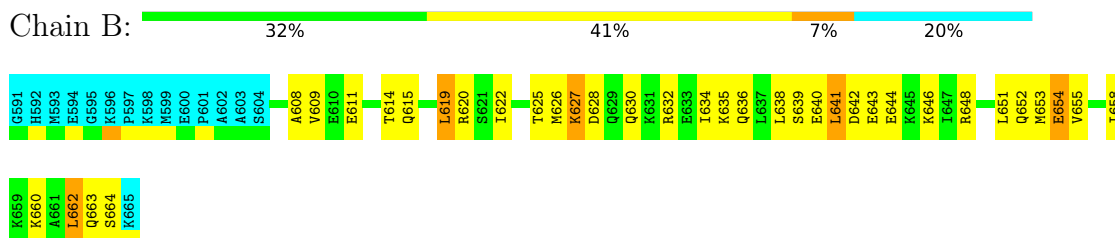


#### 4.2.27 Score per residue for model 27

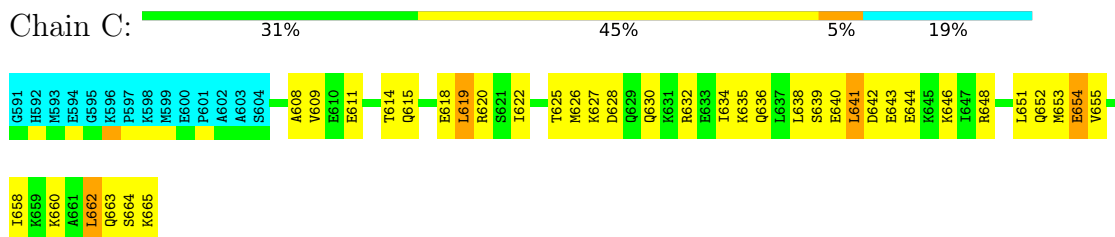
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

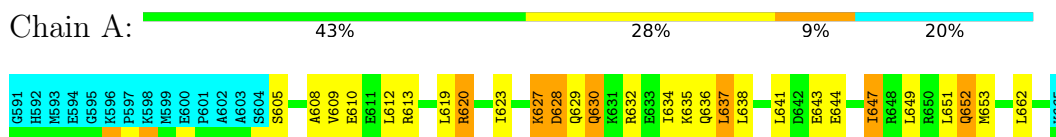


- Molecule 1: SH3 domain-containing kinase-binding protein 1

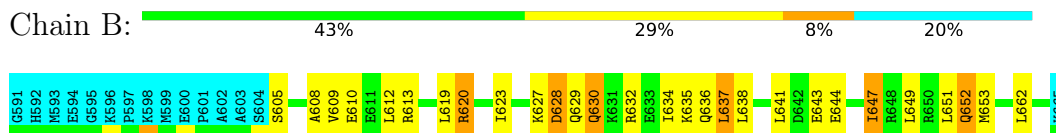


#### 4.2.28 Score per residue for model 28

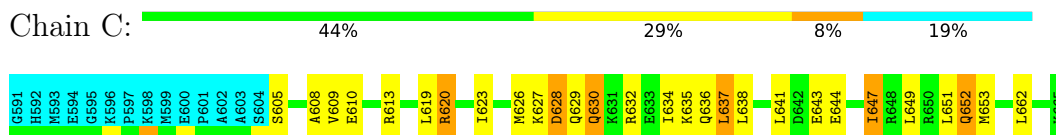
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

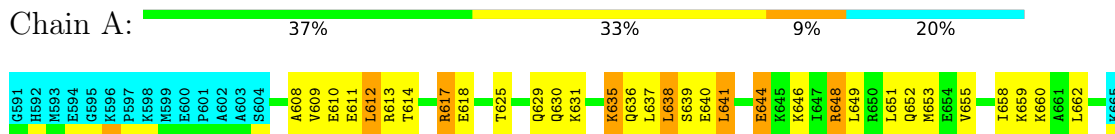


- Molecule 1: SH3 domain-containing kinase-binding protein 1

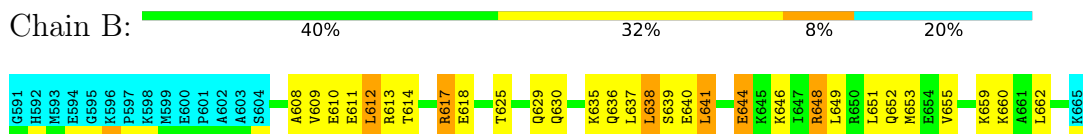


#### 4.2.29 Score per residue for model 29

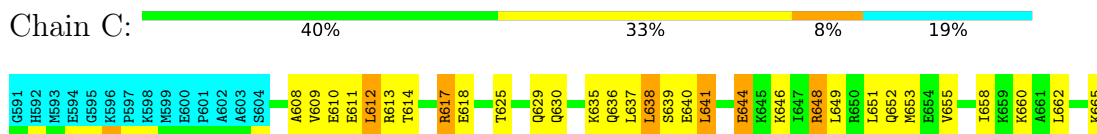
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1

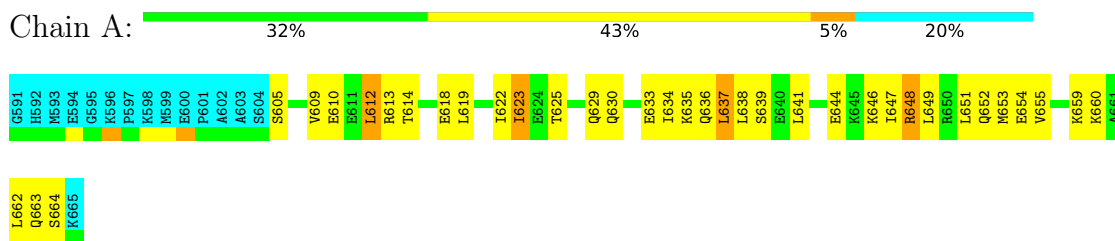


- Molecule 1: SH3 domain-containing kinase-binding protein 1

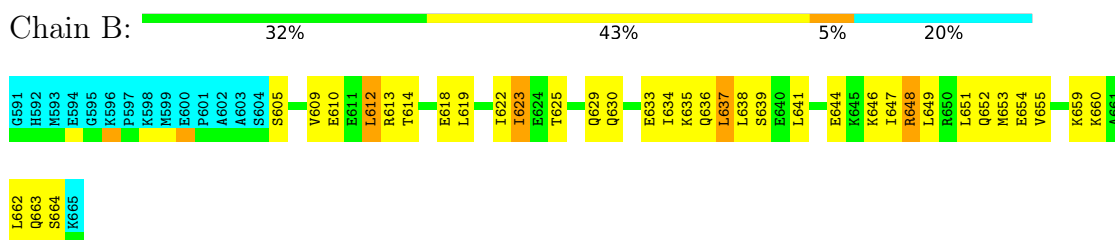


### 4.2.30 Score per residue for model 30

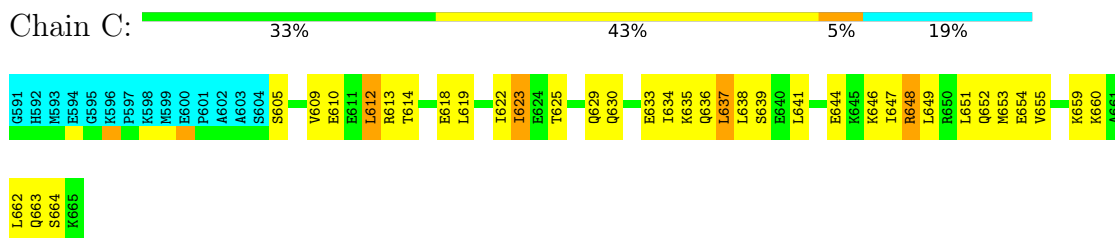
- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1



- Molecule 1: SH3 domain-containing kinase-binding protein 1





## 5 Refinement protocol and experimental data overview

The models were refined using the following method: *torsion angle dynamics*.

Of the 2440 calculated structures, 30 were deposited, based on the following criterion: *every 84th structure after convergence*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
ISD (Inferential structure determination)	refinement	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	657
Number of shifts mapped to atoms	657
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	20%

## 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	499	534	534	18±5
1	B	499	534	534	18±5
1	C	509	547	547	18±5
All	All	45210	48450	48450	1093

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 12.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:609:VAL:HG11	1:C:608:ALA:HB1	0.86	1.47	29	6
1:A:662:LEU:HD22	1:C:658:ILE:HG13	0.85	1.48	20	3
1:B:608:ALA:HB1	1:C:609:VAL:HG11	0.84	1.47	29	6
1:A:608:ALA:HB1	1:B:609:VAL:HG11	0.84	1.47	29	6
1:A:658:ILE:HG13	1:B:662:LEU:HD22	0.83	1.49	20	3
1:B:658:ILE:HG13	1:C:662:LEU:HD22	0.82	1.49	20	3
1:B:612:LEU:HD21	1:C:613:ARG:HA	0.81	1.53	8	10
1:B:651:LEU:O	1:B:655:VAL:HG23	0.81	1.76	27	21
1:A:651:LEU:O	1:A:655:VAL:HG23	0.80	1.76	27	21
1:C:651:LEU:O	1:C:655:VAL:HG23	0.80	1.76	27	21
1:B:619:LEU:HA	1:B:622:ILE:HD12	0.80	1.52	23	10
1:A:612:LEU:HD21	1:B:613:ARG:HA	0.80	1.53	8	10

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:619:LEU:HA	1:A:622:ILE:HD12	0.80	1.52	23	10
1:C:619:LEU:HA	1:C:622:ILE:HD12	0.80	1.52	23	10
1:C:665:LYS:OXT	1:C:665:LYS:HG3	0.80	1.75	8	6
1:A:613:ARG:HA	1:C:612:LEU:HD21	0.78	1.53	8	10
1:A:612:LEU:O	1:A:616:VAL:HG23	0.74	1.82	19	15
1:C:612:LEU:O	1:C:616:VAL:HG23	0.74	1.82	19	15
1:B:612:LEU:O	1:B:616:VAL:HG23	0.73	1.82	19	15
1:A:620:ARG:HA	1:A:623:ILE:HD12	0.73	1.61	20	7
1:B:620:ARG:HA	1:B:623:ILE:HD12	0.71	1.61	20	7
1:C:620:ARG:HA	1:C:623:ILE:HD12	0.71	1.61	20	7
1:A:645:LYS:O	1:A:649:LEU:HD12	0.71	1.85	17	1
1:B:645:LYS:O	1:B:649:LEU:HD12	0.71	1.85	17	1
1:C:645:LYS:O	1:C:649:LEU:HD12	0.70	1.85	17	1
1:A:655:VAL:HG21	1:C:651:LEU:HD11	0.69	1.64	30	6
1:C:605:SER:O	1:C:609:VAL:HG23	0.68	1.88	14	13
1:B:619:LEU:O	1:B:623:ILE:HD12	0.68	1.88	7	6
1:A:605:SER:O	1:A:609:VAL:HG23	0.68	1.88	14	13
1:A:619:LEU:O	1:A:623:ILE:HD12	0.68	1.88	7	6
1:B:605:SER:O	1:B:609:VAL:HG23	0.68	1.88	4	13
1:A:615:GLN:O	1:A:619:LEU:HD12	0.68	1.89	9	1
1:A:635:LYS:HA	1:A:638:LEU:HD12	0.68	1.66	20	1
1:C:615:GLN:O	1:C:619:LEU:HD12	0.67	1.89	9	1
1:A:651:LEU:HD11	1:B:655:VAL:HG21	0.67	1.64	30	6
1:C:635:LYS:HA	1:C:638:LEU:HD12	0.67	1.66	20	1
1:B:651:LEU:HD11	1:C:655:VAL:HG21	0.67	1.64	30	6
1:C:634:ILE:HA	1:C:637:LEU:HD12	0.67	1.67	17	1
1:B:615:GLN:O	1:B:619:LEU:HD12	0.67	1.89	9	2
1:B:634:ILE:HA	1:B:637:LEU:HD12	0.67	1.67	17	1
1:B:635:LYS:HA	1:B:638:LEU:HD12	0.67	1.66	20	1
1:A:662:LEU:HD11	1:C:662:LEU:HG	0.66	1.67	22	2
1:B:619:LEU:HD21	1:C:620:ARG:HG3	0.66	1.68	2	3
1:B:662:LEU:HG	1:C:662:LEU:HD11	0.66	1.67	22	2
1:C:619:LEU:O	1:C:623:ILE:HD12	0.66	1.88	7	5
1:B:606:GLN:HA	1:B:606:GLN:OE1	0.65	1.91	20	1
1:A:606:GLN:OE1	1:A:606:GLN:HA	0.65	1.91	20	1
1:A:634:ILE:HA	1:A:637:LEU:HD12	0.65	1.67	17	1
1:A:620:ARG:HG3	1:C:619:LEU:HD21	0.64	1.68	2	3
1:A:662:LEU:HG	1:B:662:LEU:HD11	0.64	1.67	22	2
1:C:606:GLN:HA	1:C:606:GLN:OE1	0.64	1.91	20	1
1:B:612:LEU:HD22	1:C:616:VAL:HG21	0.64	1.69	24	4
1:A:619:LEU:HD21	1:B:620:ARG:HG3	0.64	1.68	2	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:605:SER:O	1:A:609:VAL:HG22	0.64	1.92	3	1
1:A:612:LEU:HD22	1:B:616:VAL:HG21	0.63	1.69	24	4
1:C:665:LYS:OXT	1:C:665:LYS:CG	0.63	2.47	8	7
1:C:605:SER:O	1:C:609:VAL:HG22	0.63	1.92	3	1
1:A:661:ALA:HB3	1:B:662:LEU:HD21	0.63	1.71	14	1
1:B:605:SER:O	1:B:609:VAL:HG22	0.63	1.92	3	1
1:A:651:LEU:HD22	1:B:652:GLN:HG2	0.63	1.71	1	6
1:B:661:ALA:HB3	1:C:662:LEU:HD21	0.63	1.71	14	1
1:A:623:ILE:HD13	1:C:619:LEU:HD21	0.63	1.71	25	1
1:A:616:VAL:HG21	1:C:612:LEU:HD22	0.63	1.69	24	4
1:A:619:LEU:HD23	1:B:623:ILE:HD11	0.63	1.71	13	3
1:C:634:ILE:O	1:C:638:LEU:HD12	0.62	1.94	26	4
1:A:641:LEU:HD21	1:C:641:LEU:HA	0.62	1.71	27	5
1:B:641:LEU:HA	1:C:641:LEU:HD21	0.62	1.71	27	5
1:A:651:LEU:HD13	1:B:652:GLN:HG3	0.62	1.71	28	2
1:B:651:LEU:HD22	1:C:652:GLN:HG2	0.62	1.71	1	6
1:A:641:LEU:HA	1:B:641:LEU:HD21	0.62	1.71	27	5
1:A:634:ILE:O	1:A:638:LEU:HD12	0.62	1.94	26	4
1:A:662:LEU:HD21	1:C:661:ALA:HB3	0.62	1.71	14	1
1:B:651:LEU:HD13	1:C:652:GLN:HG3	0.62	1.71	28	2
1:A:618:GLU:O	1:A:622:ILE:HG13	0.62	1.95	17	3
1:B:619:LEU:HD23	1:C:623:ILE:HD11	0.61	1.71	13	3
1:A:613:ARG:HG2	1:C:612:LEU:HD21	0.61	1.72	26	3
1:A:652:GLN:HG2	1:C:651:LEU:HD22	0.61	1.71	1	6
1:B:634:ILE:O	1:B:638:LEU:HD12	0.61	1.94	26	4
1:A:652:GLN:HG3	1:C:651:LEU:HD13	0.61	1.71	28	2
1:B:618:GLU:O	1:B:622:ILE:HG13	0.61	1.95	17	3
1:C:621:SER:O	1:C:625:THR:HG23	0.61	1.96	26	3
1:A:621:SER:O	1:A:625:THR:HG23	0.61	1.96	26	3
1:A:619:LEU:HD21	1:B:623:ILE:HD13	0.61	1.71	25	1
1:B:619:LEU:HD21	1:C:623:ILE:HD13	0.61	1.71	25	1
1:A:623:ILE:HD11	1:C:619:LEU:HD23	0.61	1.71	13	3
1:A:612:LEU:HD21	1:B:613:ARG:HG2	0.61	1.72	26	3
1:B:621:SER:O	1:B:625:THR:HG23	0.60	1.96	26	3
1:B:662:LEU:HD12	1:C:662:LEU:CD1	0.60	2.26	11	1
1:C:618:GLU:O	1:C:622:ILE:HG13	0.60	1.95	17	3
1:B:612:LEU:HD21	1:C:613:ARG:HG2	0.60	1.72	26	3
1:A:662:LEU:HD12	1:B:662:LEU:CD1	0.60	2.26	11	1
1:A:662:LEU:CD1	1:C:662:LEU:HD12	0.60	2.26	11	1
1:A:626:MET:SD	1:B:623:ILE:HG22	0.58	2.39	13	4
1:A:637:LEU:HB2	1:C:637:LEU:HD22	0.58	1.75	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:616:VAL:HG11	1:C:615:GLN:HB3	0.58	1.75	14	2
1:B:615:GLN:HB3	1:C:616:VAL:HG11	0.58	1.75	23	2
1:A:609:VAL:HG11	1:C:608:ALA:CB	0.58	2.27	29	2
1:B:626:MET:SD	1:C:623:ILE:HG22	0.58	2.39	13	4
1:B:613:ARG:O	1:B:617:ARG:HD2	0.58	1.99	1	1
1:B:635:LYS:HE3	1:B:635:LYS:HA	0.57	1.77	8	1
1:A:615:GLN:HB3	1:B:616:VAL:HG11	0.57	1.75	23	2
1:B:608:ALA:CB	1:C:609:VAL:HG11	0.57	2.27	29	3
1:A:613:ARG:O	1:A:617:ARG:HD2	0.57	1.99	1	1
1:B:622:ILE:HG22	1:C:623:ILE:HG21	0.57	1.76	1	2
1:C:613:ARG:O	1:C:617:ARG:HD2	0.57	1.99	1	1
1:A:634:ILE:HG21	1:C:633:GLU:HB3	0.57	1.77	30	6
1:A:623:ILE:HD11	1:C:623:ILE:HD11	0.57	1.77	12	3
1:A:623:ILE:HG22	1:C:626:MET:SD	0.57	2.38	13	4
1:A:635:LYS:HE3	1:A:635:LYS:HA	0.57	1.76	8	1
1:B:637:LEU:HD22	1:C:637:LEU:HB2	0.57	1.75	17	1
1:A:623:ILE:HD11	1:B:623:ILE:HD11	0.57	1.77	12	3
1:B:623:ILE:HD11	1:C:623:ILE:HD11	0.57	1.77	12	3
1:A:633:GLU:HB3	1:B:634:ILE:HG21	0.57	1.77	30	6
1:A:623:ILE:HG21	1:C:622:ILE:HG22	0.56	1.76	1	2
1:A:662:LEU:HD22	1:C:658:ILE:HG23	0.56	1.77	12	1
1:A:658:ILE:HG23	1:B:662:LEU:HD22	0.56	1.77	12	1
1:A:623:ILE:HG12	1:B:623:ILE:HD13	0.56	1.77	24	2
1:A:622:ILE:HG22	1:B:623:ILE:HG21	0.56	1.76	1	2
1:A:637:LEU:HD22	1:B:637:LEU:HB2	0.56	1.75	17	1
1:B:623:ILE:HG12	1:C:623:ILE:HD13	0.56	1.77	24	2
1:A:608:ALA:CB	1:B:609:VAL:HG11	0.56	2.27	29	3
1:A:631:LYS:O	1:A:635:LYS:HD2	0.56	2.01	1	2
1:C:631:LYS:O	1:C:635:LYS:HD2	0.56	2.01	1	1
1:B:637:LEU:HD12	1:C:634:ILE:HG23	0.56	1.78	21	4
1:A:658:ILE:HD12	1:B:659:LYS:HA	0.56	1.78	20	2
1:C:654:GLU:O	1:C:658:ILE:HG13	0.56	2.01	23	5
1:C:635:LYS:HE3	1:C:635:LYS:HA	0.55	1.77	8	1
1:C:643:GLU:O	1:C:647:ILE:HG13	0.55	2.02	21	3
1:A:643:GLU:O	1:A:647:ILE:HG13	0.55	2.02	21	3
1:B:654:GLU:O	1:B:658:ILE:HG13	0.55	2.01	23	4
1:B:633:GLU:HB3	1:C:634:ILE:HG21	0.55	1.77	30	7
1:A:654:GLU:O	1:A:658:ILE:HG13	0.55	2.01	23	5
1:A:645:LYS:HE3	1:C:644:GLU:HG2	0.55	1.79	17	1
1:B:622:ILE:HG22	1:B:626:MET:SD	0.55	2.42	27	5
1:B:625:THR:O	1:B:629:GLN:HB2	0.55	2.02	14	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:637:LEU:HD12	1:B:634:ILE:HG23	0.55	1.78	21	4
1:B:658:ILE:HD12	1:C:659:LYS:HA	0.55	1.78	20	2
1:A:625:THR:O	1:A:629:GLN:HB2	0.55	2.02	14	3
1:C:625:THR:O	1:C:629:GLN:HB2	0.55	2.02	14	3
1:B:644:GLU:HG2	1:C:645:LYS:HE3	0.55	1.79	17	1
1:A:655:VAL:HG21	1:C:651:LEU:HD21	0.55	1.79	19	1
1:C:611:GLU:O	1:C:614:THR:HG22	0.55	2.02	26	1
1:B:631:LYS:O	1:B:635:LYS:HD2	0.54	2.01	1	1
1:B:658:ILE:HG23	1:C:662:LEU:HD22	0.54	1.77	12	1
1:A:634:ILE:HG23	1:C:637:LEU:HD12	0.54	1.78	21	4
1:A:623:ILE:HD13	1:C:623:ILE:HG12	0.54	1.77	24	2
1:C:659:LYS:HG2	1:C:663:GLN:NE2	0.54	2.17	30	1
1:B:643:GLU:O	1:B:647:ILE:HG13	0.54	2.02	21	3
1:B:659:LYS:HG2	1:B:663:GLN:NE2	0.54	2.17	30	1
1:A:651:LEU:HD21	1:B:655:VAL:HG21	0.54	1.79	19	1
1:A:622:ILE:HG22	1:A:626:MET:SD	0.54	2.43	20	5
1:B:644:GLU:HA	1:B:647:ILE:HD12	0.53	1.80	30	1
1:C:622:ILE:HG22	1:C:626:MET:SD	0.53	2.42	27	5
1:B:611:GLU:O	1:B:614:THR:HG22	0.53	2.02	26	1
1:A:644:GLU:HA	1:A:647:ILE:HD12	0.53	1.80	30	1
1:C:648:ARG:CZ	1:C:648:ARG:HB3	0.53	2.32	23	4
1:A:659:LYS:HA	1:C:658:ILE:HD12	0.53	1.78	20	2
1:B:651:LEU:HD21	1:C:655:VAL:HG21	0.53	1.79	19	1
1:A:659:LYS:HG2	1:A:663:GLN:NE2	0.53	2.17	30	1
1:B:648:ARG:CZ	1:B:648:ARG:HB3	0.53	2.32	23	4
1:B:658:ILE:CG1	1:C:662:LEU:HD13	0.53	2.34	11	1
1:A:644:GLU:HG2	1:B:645:LYS:HE3	0.53	1.79	17	1
1:A:640:GLU:HB3	1:B:641:LEU:HD13	0.53	1.81	27	3
1:A:637:LEU:HD11	1:B:638:LEU:HD12	0.53	1.81	2	1
1:C:627:LYS:O	1:C:630:GLN:HG3	0.53	2.04	25	2
1:A:648:ARG:CZ	1:A:648:ARG:HB3	0.53	2.32	23	4
1:B:640:GLU:HB3	1:C:641:LEU:HD13	0.53	1.81	27	3
1:C:644:GLU:HA	1:C:647:ILE:HD12	0.52	1.80	30	1
1:B:637:LEU:HD11	1:C:638:LEU:HD12	0.52	1.81	2	1
1:A:658:ILE:CG1	1:B:662:LEU:HD13	0.52	2.34	11	1
1:A:645:LYS:HG2	1:C:644:GLU:HG3	0.52	1.81	18	3
1:A:611:GLU:O	1:A:614:THR:HG22	0.52	2.02	26	1
1:B:627:LYS:O	1:B:630:GLN:HG3	0.52	2.04	25	2
1:A:641:LEU:HD13	1:C:640:GLU:HB3	0.52	1.81	27	3
1:B:644:GLU:HG3	1:C:645:LYS:HG2	0.52	1.81	18	3
1:A:638:LEU:HD12	1:C:637:LEU:HD11	0.52	1.81	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:662:LEU:HD13	1:C:658:ILE:CG1	0.52	2.34	11	1
1:C:622:ILE:CG2	1:C:626:MET:SD	0.52	2.98	27	1
1:A:640:GLU:O	1:A:644:GLU:HG3	0.52	2.05	17	1
1:A:644:GLU:HG3	1:B:645:LYS:HG2	0.51	1.81	18	2
1:A:627:LYS:O	1:A:630:GLN:HG3	0.51	2.05	23	2
1:A:622:ILE:CG2	1:A:626:MET:SD	0.51	2.98	27	1
1:B:622:ILE:CG2	1:B:626:MET:SD	0.51	2.98	27	1
1:C:640:GLU:O	1:C:644:GLU:HG3	0.51	2.05	17	1
1:B:640:GLU:O	1:B:644:GLU:HG3	0.51	2.05	17	1
1:B:629:GLN:O	1:B:633:GLU:HB2	0.51	2.06	17	1
1:B:640:GLU:HG2	1:C:641:LEU:HD22	0.51	1.82	27	2
1:A:655:VAL:HA	1:A:658:ILE:HD12	0.51	1.83	12	1
1:B:631:LYS:O	1:B:634:ILE:HG13	0.51	2.05	23	4
1:C:631:LYS:O	1:C:634:ILE:HG13	0.51	2.05	23	4
1:A:641:LEU:HD22	1:C:640:GLU:HG2	0.51	1.82	27	1
1:B:644:GLU:HG2	1:C:645:LYS:HD3	0.50	1.82	11	1
1:C:629:GLN:O	1:C:633:GLU:HB2	0.50	2.06	17	1
1:C:613:ARG:HB2	1:C:613:ARG:NH1	0.50	2.21	6	1
1:A:640:GLU:HG2	1:B:641:LEU:HD22	0.50	1.82	27	1
1:A:644:GLU:HG2	1:B:645:LYS:HD3	0.50	1.82	11	1
1:B:655:VAL:HA	1:B:658:ILE:HD12	0.50	1.83	12	1
1:B:613:ARG:HB2	1:B:613:ARG:NH1	0.50	2.21	6	1
1:C:643:GLU:O	1:C:647:ILE:HD12	0.50	2.06	26	4
1:A:631:LYS:O	1:A:634:ILE:HG13	0.50	2.05	23	4
1:A:654:GLU:OE2	1:A:658:ILE:HD11	0.50	2.07	26	1
1:B:619:LEU:HD11	1:C:616:VAL:HG13	0.50	1.84	3	1
1:B:654:GLU:OE2	1:B:658:ILE:HD11	0.50	2.07	26	1
1:A:613:ARG:HB2	1:A:613:ARG:NH1	0.50	2.21	6	1
1:A:643:GLU:O	1:A:647:ILE:HD12	0.50	2.06	26	4
1:B:643:GLU:O	1:B:647:ILE:HD12	0.50	2.06	26	4
1:A:629:GLN:O	1:A:633:GLU:HB2	0.50	2.06	17	1
1:A:613:ARG:HB2	1:A:613:ARG:HH11	0.49	1.67	6	1
1:A:662:LEU:HD12	1:B:662:LEU:HD11	0.49	1.84	11	1
1:B:613:ARG:HB2	1:B:613:ARG:HH11	0.49	1.67	6	1
1:A:626:MET:HG2	1:B:627:LYS:HE2	0.49	1.84	9	1
1:B:630:GLN:HG2	1:B:631:LYS:N	0.49	2.21	12	1
1:A:619:LEU:HD11	1:B:616:VAL:HG13	0.49	1.84	3	1
1:A:645:LYS:HD3	1:C:644:GLU:HG2	0.49	1.82	11	1
1:C:630:GLN:HG2	1:C:631:LYS:N	0.49	2.21	12	1
1:A:623:ILE:CG2	1:A:624:GLU:N	0.49	2.75	17	1
1:C:661:ALA:O	1:C:665:LYS:HG2	0.49	2.07	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:658:ILE:HG13	1:C:662:LEU:HD13	0.49	1.85	9	2
1:C:654:GLU:OE2	1:C:658:ILE:HD11	0.49	2.07	26	1
1:C:655:VAL:HA	1:C:658:ILE:HD12	0.49	1.83	12	1
1:B:623:ILE:CG2	1:B:624:GLU:N	0.49	2.75	17	1
1:B:640:GLU:OE1	1:C:641:LEU:HD13	0.49	2.08	24	1
1:A:616:VAL:HG13	1:C:619:LEU:HD11	0.49	1.84	3	1
1:B:662:LEU:HD12	1:C:662:LEU:HD11	0.49	1.84	11	1
1:A:609:VAL:HG13	1:C:612:LEU:HD23	0.49	1.84	29	1
1:A:651:LEU:HD21	1:B:652:GLN:HA	0.49	1.85	21	5
1:B:626:MET:HG2	1:C:627:LYS:HE2	0.49	1.84	9	1
1:A:644:GLU:OE2	1:B:645:LYS:HA	0.49	2.08	13	1
1:A:645:LYS:HA	1:C:644:GLU:OE2	0.49	2.08	13	1
1:C:623:ILE:CG2	1:C:624:GLU:N	0.49	2.75	17	1
1:C:613:ARG:HB2	1:C:613:ARG:HH11	0.48	1.67	6	1
1:A:630:GLN:HG2	1:A:631:LYS:N	0.48	2.21	12	1
1:A:644:GLU:HB2	1:B:645:LYS:HE2	0.48	1.85	14	2
1:A:612:LEU:HG	1:B:612:LEU:HB3	0.48	1.85	30	4
1:B:612:LEU:HD23	1:C:609:VAL:HG13	0.48	1.84	29	1
1:A:641:LEU:HD13	1:C:640:GLU:OE1	0.48	2.08	24	1
1:B:611:GLU:O	1:B:615:GLN:HG3	0.48	2.08	27	1
1:A:640:GLU:OE1	1:B:641:LEU:HD13	0.48	2.08	24	1
1:C:642:ASP:O	1:C:646:LYS:HD2	0.48	2.09	1	1
1:A:662:LEU:HD13	1:C:658:ILE:HG13	0.48	1.85	9	2
1:B:644:GLU:HB2	1:C:645:LYS:HE2	0.48	1.85	14	2
1:A:611:GLU:O	1:A:615:GLN:HG3	0.48	2.08	27	1
1:A:662:LEU:HD11	1:C:662:LEU:HD12	0.48	1.84	11	1
1:B:644:GLU:OE2	1:C:645:LYS:HA	0.48	2.08	13	1
1:B:651:LEU:HD21	1:C:652:GLN:HA	0.48	1.85	21	4
1:B:612:LEU:HG	1:C:612:LEU:HB3	0.48	1.86	30	3
1:C:611:GLU:O	1:C:615:GLN:HG3	0.48	2.08	27	1
1:C:665:LYS:OXT	1:C:665:LYS:HG2	0.47	2.08	21	2
1:A:658:ILE:HG13	1:B:662:LEU:HD13	0.47	1.85	9	2
1:A:652:GLN:HA	1:C:651:LEU:HD21	0.47	1.85	21	5
1:A:627:LYS:HE2	1:C:626:MET:HG2	0.47	1.84	9	2
1:A:654:GLU:HB3	1:B:655:VAL:HG11	0.47	1.86	17	1
1:B:626:MET:SD	1:C:627:LYS:HG3	0.47	2.50	11	2
1:A:612:LEU:HB3	1:C:612:LEU:HG	0.47	1.85	30	3
1:A:651:LEU:HD11	1:B:655:VAL:CG2	0.47	2.40	27	5
1:A:655:VAL:CG2	1:C:651:LEU:HD11	0.47	2.38	30	5
1:B:651:LEU:HD11	1:C:655:VAL:CG2	0.47	2.40	27	5
1:A:612:LEU:HD23	1:B:609:VAL:HG13	0.47	1.84	29	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:642:ASP:O	1:A:646:LYS:HD2	0.47	2.09	1	1
1:A:637:LEU:HD23	1:B:641:LEU:HD12	0.47	1.86	12	4
1:A:627:LYS:HG3	1:C:626:MET:SD	0.47	2.50	11	2
1:A:651:LEU:HD11	1:B:652:GLN:HB2	0.47	1.86	21	1
1:B:630:GLN:HA	1:B:630:GLN:OE1	0.47	2.10	28	1
1:A:646:LYS:O	1:A:650:ARG:HG3	0.47	2.10	4	2
1:A:619:LEU:HD13	1:B:620:ARG:HD2	0.47	1.86	7	2
1:B:622:ILE:HG23	1:B:626:MET:SD	0.47	2.50	26	1
1:A:630:GLN:OE1	1:A:630:GLN:HA	0.47	2.10	28	1
1:B:642:ASP:O	1:B:646:LYS:HD2	0.47	2.09	1	1
1:A:641:LEU:HD12	1:C:637:LEU:HD23	0.47	1.87	22	4
1:A:626:MET:SD	1:B:627:LYS:HG3	0.47	2.50	11	2
1:B:643:GLU:O	1:B:647:ILE:HG22	0.47	2.10	14	1
1:C:630:GLN:HA	1:C:630:GLN:OE1	0.47	2.10	28	1
1:B:646:LYS:O	1:B:650:ARG:HG3	0.47	2.10	4	2
1:B:612:LEU:HD21	1:C:613:ARG:HB2	0.47	1.87	16	1
1:C:639:SER:O	1:C:643:GLU:HG2	0.47	2.10	19	1
1:A:623:ILE:HG12	1:B:623:ILE:HG12	0.47	1.87	18	3
1:A:652:GLN:HB2	1:C:651:LEU:HD11	0.47	1.86	21	1
1:A:623:ILE:HG12	1:C:623:ILE:HG12	0.47	1.87	8	3
1:B:619:LEU:HD13	1:C:620:ARG:HD2	0.47	1.86	7	2
1:A:643:GLU:O	1:A:647:ILE:HG22	0.47	2.10	14	1
1:B:639:SER:O	1:B:643:GLU:HG2	0.47	2.10	19	1
1:A:620:ARG:HB2	1:A:620:ARG:HH11	0.47	1.70	26	1
1:C:646:LYS:O	1:C:650:ARG:HG3	0.46	2.10	4	2
1:B:622:ILE:O	1:B:626:MET:HG3	0.46	2.10	4	1
1:A:613:ARG:CA	1:C:612:LEU:HD21	0.46	2.39	5	2
1:A:655:VAL:HG11	1:C:654:GLU:HB3	0.46	1.86	17	1
1:C:620:ARG:HH11	1:C:620:ARG:HB2	0.46	1.70	26	1
1:A:620:ARG:HD2	1:C:619:LEU:HD13	0.46	1.86	7	2
1:A:622:ILE:HG23	1:A:626:MET:SD	0.46	2.50	26	1
1:C:622:ILE:HG23	1:C:626:MET:SD	0.46	2.50	26	1
1:B:623:ILE:HG12	1:C:623:ILE:HG12	0.46	1.87	8	3
1:A:637:LEU:HG	1:B:638:LEU:HG	0.46	1.87	26	3
1:B:630:GLN:HG2	1:C:630:GLN:HB3	0.46	1.88	1	1
1:A:622:ILE:O	1:A:626:MET:HG3	0.46	2.10	4	2
1:C:622:ILE:O	1:C:626:MET:HG3	0.46	2.10	4	1
1:B:654:GLU:HB3	1:C:655:VAL:HG11	0.46	1.86	17	1
1:A:645:LYS:HE2	1:C:644:GLU:HB2	0.46	1.85	14	2
1:B:611:GLU:O	1:B:614:THR:HG23	0.46	2.11	22	1
1:B:637:LEU:HD23	1:C:641:LEU:HD12	0.46	1.86	12	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:660:LYS:HZ2	1:A:660:LYS:HA	0.46	1.71	15	1
1:A:613:ARG:HB2	1:C:612:LEU:HD21	0.46	1.87	16	1
1:A:611:GLU:O	1:A:614:THR:HG23	0.46	2.11	22	1
1:A:612:LEU:HD21	1:B:613:ARG:CA	0.46	2.39	5	1
1:A:638:LEU:HG	1:C:637:LEU:HG	0.46	1.88	18	3
1:B:637:LEU:HG	1:C:638:LEU:HG	0.46	1.87	26	3
1:B:620:ARG:HB2	1:B:620:ARG:HH11	0.46	1.70	26	1
1:B:662:LEU:HD23	1:B:663:GLN:HE21	0.46	1.70	4	1
1:A:630:GLN:HG2	1:B:630:GLN:HB3	0.46	1.88	1	1
1:C:640:GLU:HA	1:C:643:GLU:OE1	0.46	2.11	17	1
1:B:629:GLN:O	1:B:633:GLU:HG3	0.46	2.11	23	1
1:A:651:LEU:HD22	1:B:651:LEU:HD22	0.45	1.88	10	1
1:A:651:LEU:HD22	1:C:651:LEU:HD22	0.45	1.88	10	1
1:C:643:GLU:O	1:C:647:ILE:HG22	0.45	2.10	14	1
1:A:630:GLN:HB3	1:C:630:GLN:HG2	0.45	1.88	1	1
1:B:651:LEU:HD22	1:C:651:LEU:HD22	0.45	1.88	10	1
1:A:639:SER:O	1:A:643:GLU:HG2	0.45	2.10	19	1
1:A:623:ILE:HD13	1:C:619:LEU:HD22	0.45	1.88	5	1
1:B:616:VAL:HA	1:B:619:LEU:HD12	0.45	1.87	1	1
1:C:665:LYS:CG	1:C:665:LYS:OXT	0.45	2.63	9	1
1:C:611:GLU:O	1:C:614:THR:HG23	0.45	2.11	22	1
1:C:629:GLN:O	1:C:633:GLU:HG3	0.45	2.11	23	1
1:A:662:LEU:HD23	1:A:663:GLN:HE21	0.45	1.70	4	1
1:A:623:ILE:CG1	1:B:623:ILE:HD13	0.45	2.41	24	1
1:A:651:LEU:HD21	1:B:652:GLN:CA	0.45	2.41	5	1
1:B:627:LYS:O	1:B:630:GLN:HG2	0.45	2.12	7	1
1:B:649:LEU:HA	1:B:652:GLN:HG3	0.45	1.89	7	1
1:A:612:LEU:HD21	1:B:613:ARG:HB2	0.45	1.87	16	1
1:A:631:LYS:O	1:A:635:LYS:HB2	0.45	2.12	17	1
1:B:651:LEU:HD11	1:C:652:GLN:HB2	0.45	1.86	21	1
1:C:642:ASP:O	1:C:646:LYS:HG3	0.45	2.12	10	1
1:A:649:LEU:HA	1:A:652:GLN:HG3	0.45	1.89	7	1
1:B:623:ILE:CG1	1:C:623:ILE:HD13	0.45	2.41	24	1
1:B:611:GLU:O	1:B:615:GLN:HG2	0.45	2.12	26	1
1:A:627:LYS:O	1:A:630:GLN:HG2	0.45	2.12	7	1
1:C:645:LYS:O	1:C:649:LEU:HG	0.45	2.12	9	1
1:B:646:LYS:HB3	1:B:650:ARG:NH1	0.45	2.27	17	1
1:A:652:GLN:CA	1:C:651:LEU:HD21	0.45	2.41	5	1
1:A:623:ILE:HD13	1:C:623:ILE:CG1	0.45	2.41	24	1
1:A:654:GLU:O	1:A:658:ILE:HG22	0.45	2.12	24	1
1:B:654:GLU:O	1:B:658:ILE:HG22	0.45	2.12	24	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:645:LYS:O	1:B:649:LEU:HG	0.44	2.12	9	1
1:A:642:ASP:O	1:A:646:LYS:HG3	0.44	2.12	10	1
1:B:640:GLU:HA	1:B:643:GLU:OE1	0.44	2.11	17	1
1:C:662:LEU:HA	1:C:665:LYS:HD3	0.44	1.88	21	1
1:C:611:GLU:O	1:C:615:GLN:HG2	0.44	2.12	26	1
1:C:616:VAL:HA	1:C:619:LEU:HD12	0.44	1.87	1	1
1:C:662:LEU:HD23	1:C:663:GLN:HE21	0.44	1.70	4	1
1:B:630:GLN:NE2	1:C:630:GLN:HB3	0.44	2.27	5	1
1:C:631:LYS:O	1:C:635:LYS:HB2	0.44	2.12	17	2
1:A:626:MET:HG2	1:B:630:GLN:NE2	0.44	2.27	25	2
1:C:627:LYS:O	1:C:630:GLN:HG2	0.44	2.12	7	2
1:A:640:GLU:HA	1:A:643:GLU:OE1	0.44	2.11	17	1
1:A:611:GLU:O	1:A:615:GLN:HG2	0.44	2.12	26	1
1:A:616:VAL:HA	1:A:619:LEU:HD12	0.44	1.87	1	1
1:A:619:LEU:HD22	1:B:623:ILE:HD13	0.44	1.88	5	1
1:B:642:ASP:O	1:B:646:LYS:HG3	0.44	2.12	10	1
1:A:648:ARG:HH11	1:A:652:GLN:HG3	0.44	1.72	14	1
1:B:655:VAL:O	1:B:659:LYS:HB2	0.44	2.13	15	1
1:B:637:LEU:HD21	1:C:638:LEU:HA	0.44	1.89	15	2
1:B:651:LEU:HD21	1:C:652:GLN:CA	0.44	2.41	5	1
1:A:645:LYS:O	1:A:649:LEU:HG	0.44	2.12	9	1
1:C:654:GLU:O	1:C:658:ILE:HG22	0.44	2.12	24	1
1:A:630:GLN:NE2	1:B:630:GLN:HB3	0.44	2.27	5	1
1:A:630:GLN:HB3	1:C:630:GLN:NE2	0.44	2.27	5	1
1:A:630:GLN:NE2	1:C:626:MET:HG2	0.44	2.27	25	2
1:A:659:LYS:HD2	1:A:663:GLN:HE21	0.44	1.73	12	1
1:B:631:LYS:O	1:B:635:LYS:HB2	0.44	2.12	17	1
1:C:646:LYS:HB3	1:C:650:ARG:NH1	0.44	2.27	17	1
1:B:626:MET:HG2	1:C:630:GLN:NE2	0.44	2.28	7	2
1:A:629:GLN:O	1:A:633:GLU:HG3	0.44	2.11	23	1
1:C:628:ASP:HB3	1:C:632:ARG:NH2	0.44	2.28	28	1
1:C:619:LEU:O	1:C:623:ILE:HG13	0.44	2.13	2	2
1:C:659:LYS:HD2	1:C:663:GLN:HE21	0.44	1.73	12	1
1:B:619:LEU:O	1:B:623:ILE:HG13	0.43	2.13	2	2
1:C:648:ARG:HH11	1:C:652:GLN:HG3	0.43	1.72	14	1
1:A:655:VAL:O	1:A:659:LYS:HB2	0.43	2.13	15	1
1:C:656:ASN:O	1:C:660:LYS:HG3	0.43	2.13	3	1
1:B:637:LEU:HD22	1:C:637:LEU:CB	0.43	2.43	17	1
1:B:615:GLN:HG2	1:B:619:LEU:CD1	0.43	2.44	19	1
1:A:619:LEU:O	1:A:623:ILE:HG13	0.43	2.13	2	3
1:B:612:LEU:HD21	1:C:613:ARG:CA	0.43	2.39	5	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:659:LYS:HD2	1:B:663:GLN:HE21	0.43	1.73	12	1
1:A:646:LYS:HB3	1:A:650:ARG:NH1	0.43	2.27	17	1
1:B:651:LEU:CD2	1:C:655:VAL:HG21	0.43	2.43	19	1
1:C:615:GLN:HG2	1:C:619:LEU:CD1	0.43	2.44	19	1
1:A:628:ASP:HB3	1:A:632:ARG:NH2	0.43	2.28	28	1
1:A:658:ILE:HD11	1:B:659:LYS:HG2	0.43	1.90	12	1
1:B:651:LEU:HG	1:C:652:GLN:HG3	0.43	1.91	24	1
1:C:649:LEU:HA	1:C:652:GLN:HG3	0.43	1.89	7	1
1:A:627:LYS:HG3	1:C:626:MET:HG3	0.43	1.91	12	1
1:B:648:ARG:HA	1:C:648:ARG:HD3	0.43	1.91	11	1
1:A:626:MET:HG3	1:B:627:LYS:HG3	0.43	1.91	12	1
1:A:659:LYS:HG2	1:C:658:ILE:HD11	0.43	1.90	12	1
1:B:628:ASP:HB3	1:B:632:ARG:NH2	0.43	2.28	28	1
1:B:656:ASN:O	1:B:660:LYS:HG3	0.43	2.13	3	1
1:A:655:VAL:HG21	1:C:651:LEU:CD2	0.43	2.43	19	1
1:B:619:LEU:HD22	1:C:623:ILE:HD13	0.43	1.88	5	1
1:B:648:ARG:HH11	1:B:652:GLN:HG3	0.43	1.72	14	1
1:A:632:ARG:NE	1:A:632:ARG:HA	0.43	2.29	18	1
1:A:651:LEU:CD2	1:B:655:VAL:HG21	0.43	2.43	19	1
1:A:619:LEU:HD22	1:B:620:ARG:HD2	0.43	1.91	14	1
1:A:637:LEU:HD21	1:B:638:LEU:HA	0.43	1.90	15	2
1:A:649:LEU:O	1:A:653:MET:HG2	0.43	2.14	23	1
1:A:630:GLN:HG2	1:B:630:GLN:HG2	0.43	1.91	2	2
1:A:656:ASN:O	1:A:660:LYS:HG3	0.43	2.13	3	1
1:A:648:ARG:HD3	1:C:648:ARG:HA	0.43	1.90	11	1
1:C:655:VAL:O	1:C:659:LYS:HB2	0.43	2.13	15	1
1:A:615:GLN:HG2	1:A:619:LEU:CD1	0.43	2.44	19	1
1:C:632:ARG:NE	1:C:632:ARG:HA	0.42	2.29	18	1
1:B:630:GLN:HG2	1:C:630:GLN:HG2	0.42	1.91	2	2
1:A:623:ILE:CG2	1:C:626:MET:SD	0.42	3.08	4	1
1:A:638:LEU:HA	1:C:637:LEU:HD21	0.42	1.90	29	2
1:B:632:ARG:NE	1:B:632:ARG:HA	0.42	2.29	18	1
1:B:649:LEU:O	1:B:653:MET:HG2	0.42	2.14	23	1
1:A:652:GLN:HG3	1:C:651:LEU:HG	0.42	1.91	24	1
1:B:658:ILE:HD11	1:C:659:LYS:HG2	0.42	1.90	12	1
1:A:631:LYS:HA	1:A:631:LYS:HD3	0.42	1.56	13	1
1:A:637:LEU:CB	1:C:637:LEU:HD22	0.42	2.43	17	1
1:C:649:LEU:O	1:C:653:MET:HG2	0.42	2.14	23	1
1:A:651:LEU:HG	1:B:652:GLN:HG3	0.42	1.91	24	1
1:A:630:GLN:HG2	1:C:630:GLN:HG2	0.42	1.91	2	2
1:A:645:LYS:HE3	1:C:640:GLU:HG2	0.42	1.92	7	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:626:MET:HG3	1:C:627:LYS:HG3	0.42	1.91	12	1
1:C:660:LYS:HA	1:C:660:LYS:HZ2	0.42	1.73	15	1
1:A:648:ARG:HD3	1:C:644:GLU:HG2	0.42	1.92	27	1
1:A:626:MET:SD	1:B:623:ILE:CG2	0.42	3.07	4	1
1:A:655:VAL:HG21	1:C:651:LEU:CD1	0.42	2.44	27	1
1:A:641:LEU:O	1:A:645:LYS:HG3	0.42	2.15	7	2
1:A:620:ARG:HD2	1:C:619:LEU:HD22	0.42	1.91	14	1
1:B:660:LYS:HZ2	1:B:660:LYS:HA	0.42	1.74	15	1
1:A:658:ILE:HD11	1:B:655:VAL:HG13	0.42	1.90	16	1
1:B:611:GLU:O	1:B:614:THR:CG2	0.42	2.68	29	1
1:C:643:GLU:HG2	1:C:647:ILE:HD11	0.42	1.92	2	1
1:A:662:LEU:HD13	1:C:658:ILE:CG2	0.42	2.45	7	1
1:B:643:GLU:OE1	1:B:643:GLU:HA	0.42	2.14	15	1
1:B:658:ILE:HD11	1:C:655:VAL:HG13	0.42	1.90	16	1
1:A:659:LYS:HE3	1:C:658:ILE:HB	0.42	1.92	20	1
1:B:619:LEU:HD11	1:C:620:ARG:HB2	0.42	1.92	24	1
1:C:611:GLU:O	1:C:614:THR:CG2	0.42	2.68	29	1
1:C:641:LEU:O	1:C:645:LYS:HG3	0.42	2.14	4	2
1:A:648:ARG:HA	1:B:648:ARG:HD3	0.42	1.91	11	1
1:C:644:GLU:HA	1:C:647:ILE:CG2	0.42	2.45	14	1
1:C:644:GLU:O	1:C:648:ARG:HB2	0.42	2.15	29	1
1:B:626:MET:SD	1:C:623:ILE:CG2	0.42	3.07	4	1
1:A:658:ILE:HB	1:B:659:LYS:HE3	0.42	1.92	20	1
1:B:641:LEU:O	1:B:645:LYS:HG3	0.41	2.15	7	2
1:A:619:LEU:O	1:A:622:ILE:HB	0.41	2.15	8	1
1:A:648:ARG:NH2	1:C:647:ILE:HD13	0.41	2.30	8	1
1:A:658:ILE:CG2	1:B:662:LEU:HD13	0.41	2.45	7	1
1:B:619:LEU:HD11	1:C:623:ILE:HD13	0.41	1.92	12	2
1:B:647:ILE:HD13	1:C:648:ARG:NH2	0.41	2.31	8	1
1:B:658:ILE:HB	1:C:659:LYS:HE3	0.41	1.92	20	1
1:A:620:ARG:HB2	1:C:619:LEU:HD11	0.41	1.92	24	1
1:B:658:ILE:CG2	1:C:662:LEU:HD13	0.41	2.45	7	1
1:B:629:GLN:HB3	1:C:627:LYS:HZ1	0.41	1.74	22	1
1:B:651:LEU:CD1	1:C:655:VAL:HG21	0.41	2.44	27	1
1:A:623:ILE:HD13	1:C:619:LEU:HD11	0.41	1.92	12	2
1:B:644:GLU:HA	1:B:647:ILE:CG2	0.41	2.45	14	1
1:A:611:GLU:O	1:A:614:THR:CG2	0.41	2.68	29	1
1:A:644:GLU:O	1:A:648:ARG:HB2	0.41	2.15	29	1
1:A:644:GLU:HA	1:A:647:ILE:CG2	0.41	2.45	14	1
1:B:619:LEU:HD22	1:C:620:ARG:HD2	0.41	1.91	14	1
1:A:647:ILE:HG22	1:B:652:GLN:HE22	0.41	1.75	26	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:644:GLU:HG2	1:B:648:ARG:HD3	0.41	1.92	27	1
1:B:640:GLU:HG2	1:C:645:LYS:HE3	0.41	1.92	7	1
1:B:654:GLU:OE2	1:C:655:VAL:HG12	0.41	2.16	23	1
1:B:644:GLU:O	1:B:648:ARG:HB2	0.41	2.15	29	1
1:A:640:GLU:HG2	1:B:645:LYS:HE3	0.41	1.92	7	1
1:A:637:LEU:CD2	1:B:637:LEU:HB3	0.41	2.46	1	1
1:B:648:ARG:HH12	1:B:649:LEU:HD12	0.41	1.76	12	1
1:B:606:GLN:O	1:B:610:GLU:HG2	0.41	2.16	13	1
1:B:646:LYS:HA	1:B:646:LYS:HE2	0.41	1.93	18	1
1:A:638:LEU:HG	1:C:637:LEU:CG	0.41	2.46	26	1
1:B:637:LEU:CD2	1:C:637:LEU:HB3	0.41	2.46	1	1
1:A:643:GLU:HG2	1:A:647:ILE:HD11	0.41	1.92	2	1
1:C:612:LEU:HD12	1:C:612:LEU:HA	0.41	1.72	12	1
1:A:606:GLN:O	1:A:610:GLU:HG2	0.41	2.16	13	1
1:A:643:GLU:HA	1:A:643:GLU:OE1	0.41	2.14	15	1
1:C:643:GLU:OE1	1:C:643:GLU:HA	0.41	2.14	15	1
1:A:629:GLN:HG3	1:B:631:LYS:HZ1	0.41	1.75	18	1
1:A:654:GLU:OE2	1:B:655:VAL:HG12	0.41	2.16	23	1
1:A:655:VAL:HG12	1:C:654:GLU:OE2	0.41	2.16	23	1
1:A:642:ASP:O	1:A:646:LYS:HD3	0.41	2.15	25	1
1:C:619:LEU:O	1:C:622:ILE:HB	0.41	2.15	8	1
1:A:645:LYS:HD3	1:C:644:GLU:HG3	0.41	1.93	12	1
1:A:637:LEU:HD22	1:B:637:LEU:CB	0.41	2.43	17	1
1:A:644:GLU:OE2	1:B:644:GLU:HB2	0.41	2.16	17	1
1:C:642:ASP:O	1:C:646:LYS:HD3	0.41	2.15	25	1
1:A:652:GLN:HE22	1:C:647:ILE:HG22	0.41	1.75	26	1
1:B:629:GLN:HG3	1:C:631:LYS:HZ1	0.40	1.76	18	1
1:C:615:GLN:O	1:C:619:LEU:HG	0.40	2.17	26	1
1:B:644:GLU:HG2	1:C:648:ARG:HD3	0.40	1.92	27	1
1:B:648:ARG:NH1	1:B:648:ARG:HB3	0.40	2.31	27	1
1:A:647:ILE:HD13	1:B:648:ARG:NH2	0.40	2.31	8	1
1:A:655:VAL:HG13	1:C:658:ILE:HD11	0.40	1.90	16	1
1:A:644:GLU:HB2	1:C:644:GLU:OE2	0.40	2.16	17	1
1:C:665:LYS:HG3	1:C:665:LYS:OXT	0.40	2.16	18	1
1:A:637:LEU:HB3	1:C:637:LEU:CD2	0.40	2.46	1	1
1:A:637:LEU:CG	1:B:638:LEU:HG	0.40	2.46	26	1
1:B:637:LEU:CG	1:C:638:LEU:HG	0.40	2.46	26	1
1:C:648:ARG:NH1	1:C:648:ARG:HB3	0.40	2.32	27	1
1:B:633:GLU:HB2	1:C:634:ILE:HD13	0.40	1.93	30	1
1:A:643:GLU:HA	1:A:646:LYS:HE2	0.40	1.94	10	1
1:A:662:LEU:HD21	1:C:661:ALA:CB	0.40	2.45	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:619:LEU:HG	1:B:620:ARG:HD2	0.40	1.94	17	1
1:A:619:LEU:HD11	1:B:620:ARG:HB2	0.40	1.92	24	1
1:B:615:GLN:O	1:B:619:LEU:HG	0.40	2.16	26	1
1:A:629:GLN:HB3	1:B:627:LYS:NZ	0.40	2.31	27	1
1:A:659:LYS:HD2	1:C:658:ILE:HD11	0.40	1.94	29	1
1:A:643:GLU:HA	1:A:646:LYS:HD2	0.40	1.94	1	1
1:C:648:ARG:HH12	1:C:649:LEU:HD12	0.40	1.76	12	1
1:A:612:LEU:HD22	1:B:613:ARG:HE	0.40	1.77	25	1
1:B:642:ASP:O	1:B:646:LYS:HD3	0.40	2.15	25	1
1:A:648:ARG:HB3	1:A:648:ARG:NH1	0.40	2.32	27	1
1:A:658:ILE:HD11	1:B:659:LYS:HD2	0.40	1.94	29	1

## 6.3 Torsion angles [i](#)

### 6.3.1 Protein backbone [i](#)

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	60/75 (80%)	60±0 (100±0%)	0±0 (0±0%)	0±0 (0±0%)	100	100
1	B	60/75 (80%)	60±0 (100±0%)	0±0 (0±0%)	0±0 (0±0%)	100	100
1	C	60/75 (80%)	60±0 (100±0%)	0±0 (0±0%)	0±0 (0±0%)	100	100
All	All	5400/6750 (80%)	5394 (100%)	6 (0%)	0 (0%)	100	100

There are no Ramachandran outliers.

### 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	57/68 (84%)	37±3 (64±5%)	20±3 (36±5%)	1	8

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	57/68 (84%)	36±3 (64±5%)	20±3 (36±5%)	1 8
1	C	58/68 (85%)	37±3 (64±5%)	21±3 (36±5%)	1 8
All	All	5160/6120 (84%)	3302 (64%)	1858 (36%)	1 8

All 166 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	641	LEU	27
1	B	641	LEU	27
1	C	641	LEU	27
1	A	617	ARG	25
1	B	617	ARG	25
1	C	617	ARG	25
1	A	630	GLN	24
1	B	630	GLN	24
1	C	630	GLN	24
1	A	636	GLN	19
1	A	662	LEU	19
1	B	636	GLN	19
1	B	644	GLU	19
1	B	662	LEU	19
1	C	636	GLN	19
1	C	662	LEU	19
1	A	620	ARG	19
1	A	660	LYS	19
1	B	620	ARG	19
1	B	660	LYS	19
1	C	620	ARG	19
1	C	660	LYS	19
1	A	644	GLU	18
1	C	644	GLU	18
1	A	632	ARG	18
1	A	639	SER	18
1	B	632	ARG	18
1	B	639	SER	18
1	C	632	ARG	18
1	C	639	SER	18
1	A	635	LYS	17
1	A	646	LYS	17
1	B	635	LYS	17

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Mol	Chain	Res	Type	Models (Total)
1	B	646	LYS	17
1	C	635	LYS	17
1	C	646	LYS	17
1	A	614	THR	16
1	B	614	THR	16
1	C	614	THR	16
1	A	651	LEU	15
1	B	651	LEU	15
1	C	651	LEU	15
1	C	665	LYS	15
1	A	649	LEU	14
1	A	653	MET	14
1	A	663	GLN	14
1	B	649	LEU	14
1	B	653	MET	14
1	B	663	GLN	14
1	C	649	LEU	14
1	C	653	MET	14
1	C	663	GLN	14
1	A	637	LEU	13
1	A	659	LYS	13
1	B	637	LEU	13
1	B	659	LYS	13
1	C	637	LEU	13
1	C	659	LYS	13
1	A	611	GLU	13
1	B	611	GLU	13
1	C	611	GLU	13
1	A	642	ASP	13
1	B	642	ASP	13
1	C	642	ASP	13
1	A	619	LEU	13
1	B	619	LEU	13
1	C	619	LEU	13
1	A	657	ASP	12
1	B	657	ASP	12
1	C	657	ASP	12
1	A	618	GLU	12
1	B	618	GLU	12
1	C	618	GLU	12
1	A	606	GLN	12
1	B	606	GLN	12

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Mol	Chain	Res	Type	Models (Total)
1	C	606	GLN	12
1	A	631	LYS	11
1	B	631	LYS	11
1	C	631	LYS	11
1	A	627	LYS	11
1	A	648	ARG	11
1	B	627	LYS	11
1	B	648	ARG	11
1	C	627	LYS	11
1	C	648	ARG	11
1	A	621	SER	11
1	B	621	SER	11
1	C	621	SER	11
1	A	628	ASP	10
1	B	628	ASP	10
1	C	628	ASP	10
1	A	652	GLN	10
1	A	664	SER	10
1	B	652	GLN	10
1	B	664	SER	10
1	C	652	GLN	10
1	C	664	SER	10
1	A	629	GLN	10
1	B	629	GLN	10
1	C	629	GLN	10
1	A	625	THR	10
1	B	625	THR	10
1	C	625	THR	10
1	A	612	LEU	9
1	A	624	GLU	9
1	B	612	LEU	9
1	B	624	GLU	9
1	C	612	LEU	9
1	C	624	GLU	9
1	A	638	LEU	9
1	B	638	LEU	9
1	C	638	LEU	9
1	A	640	GLU	8
1	B	640	GLU	8
1	C	640	GLU	8
1	A	643	GLU	8
1	B	643	GLU	8

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Mol	Chain	Res	Type	Models (Total)
1	C	643	GLU	8
1	A	605	SER	7
1	B	605	SER	7
1	C	605	SER	7
1	A	654	GLU	7
1	B	654	GLU	7
1	C	654	GLU	7
1	A	615	GLN	6
1	A	650	ARG	6
1	B	615	GLN	6
1	B	650	ARG	6
1	C	615	GLN	6
1	C	650	ARG	6
1	A	623	ILE	6
1	B	623	ILE	6
1	C	623	ILE	6
1	A	610	GLU	6
1	B	610	GLU	6
1	C	610	GLU	6
1	A	656	ASN	5
1	B	656	ASN	5
1	C	656	ASN	5
1	A	658	ILE	4
1	B	658	ILE	4
1	C	658	ILE	4
1	A	626	MET	3
1	B	626	MET	3
1	C	626	MET	3
1	A	609	VAL	3
1	B	609	VAL	3
1	C	609	VAL	3
1	A	633	GLU	2
1	B	633	GLU	2
1	C	633	GLU	2
1	A	647	ILE	2
1	B	647	ILE	2
1	C	647	ILE	2
1	A	613	ARG	2
1	A	634	ILE	2
1	B	613	ARG	2
1	B	634	ILE	2
1	C	613	ARG	2

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Mol	Chain	Res	Type	Models (Total)
1	C	634	ILE	2
1	A	616	VAL	1
1	B	616	VAL	1
1	C	616	VAL	1
1	A	645	LYS	1
1	B	645	LYS	1
1	C	645	LYS	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 i

The completeness of assignment taking into account all chemical shift lists is 20% for the well-defined parts and 20% for the entire structure.

### 7.1 Chemical shift list 1

File name: working\_cs.cif

Chemical shift list name: *assigned\_chem\_shift\_list*

#### 7.1.1 Bookkeeping i

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	657
Number of shifts mapped to atoms	657
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

#### 7.1.2 Chemical shift referencing i

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	73	$-0.45 \pm 0.09$	None needed ( $< 0.5$ ppm)
$^{13}\text{C}_\beta$	71	$0.88 \pm 0.14$	Should be checked
$^{13}\text{C}'$	71	$2.56 \pm 0.06$	Should be applied
$^{15}\text{N}$	71	$-0.30 \pm 0.12$	None needed ( $< 0.5$ ppm)

#### 7.1.3 Completeness of resonance assignments i

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 20%, i.e. 553 atoms were assigned a chemical shift out of a possible 2736. 0 out of 33 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	$^1\text{H}$	$^{13}\text{C}$	$^{15}\text{N}$
Backbone	238/905 (26%)	60/362 (17%)	118/362 (33%)	60/181 (33%)
Sidechain	315/1831 (17%)	191/1169 (16%)	116/562 (21%)	8/100 (8%)

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	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Overall	553/2736 (20%)	251/1531 (16%)	234/924 (25%)	68/281 (24%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 20%, i.e. 657 atoms were assigned a chemical shift out of a possible 3276. 0 out of 33 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	<sup>1</sup> H	<sup>13</sup> C	<sup>15</sup> N
Backbone	286/1119 (26%)	71/450 (16%)	144/450 (32%)	71/219 (32%)
Sidechain	371/2133 (17%)	230/1365 (17%)	133/660 (20%)	8/108 (7%)
Aromatic	0/24 (0%)	0/12 (0%)	0/6 (0%)	0/6 (0%)
Overall	657/3276 (20%)	301/1827 (16%)	277/1116 (25%)	79/333 (24%)

#### 7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

#### 7.1.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:

