



Full wwPDB NMR Structure Validation Report ⓘ

Feb 17, 2018 – 02:36 pm GMT

PDB ID : 1JSP
Title : NMR Structure of CBP Bromodomain in complex with p53 peptide
Authors : He, Y.; Mujtaba, S.; Zeng, L.; Yan, S.; Zhou, M.-M.
Deposited on : 2001-08-17

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

Cyrange	:	Kirchner and Güntert (2011)
NmrClust	:	Kelley et al. (1996)
MolProbity	:	4.02b-467
Mogul	:	1.7.3 (157068), CSD as539be (2018)
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
ShiftChecker	:	trunk30686
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk30686

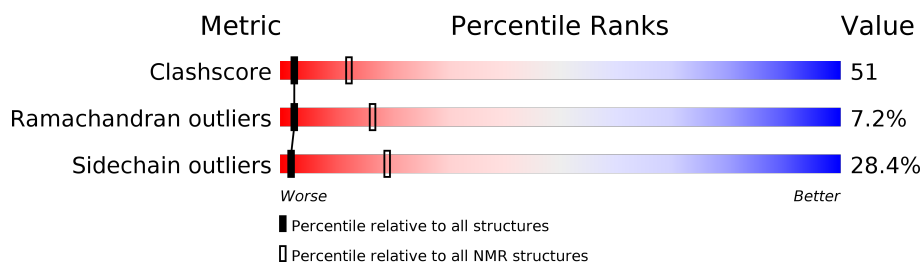
1 Overall quality at a glance

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	136279	12091
Ramachandran outliers	132675	10835
Sidechain outliers	132484	10811

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

Mol	Chain	Length	Quality of chain
1	A	20	
2	B	121	

2 Ensemble composition and analysis

This entry contains 20 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:383-A:384, B:1081- B:1197 (119)	0.65	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 5 clusters and 1 single-model cluster was found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2375 atoms, of which 1192 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called tumor protein p53.

Mol	Chain	Residues	Atoms						Trace
1	A	20	Total	C	H	N	O	S	0
			354	105	186	34	28	1	

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	382	ALY	LYS	MODIFIED RESIDUE	UNP P04637

- Molecule 2 is a protein called CREB-BINDING PROTEIN.

Mol	Chain	Residues	Atoms						Trace
2	B	121	Total	C	H	N	O	S	0
			2021	655	1006	169	185	6	

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1077	GLY	-	CLONING ARTIFACT	UNP Q92793
B	1078	SER	-	CLONING ARTIFACT	UNP Q92793
B	1079	HIS	-	CLONING ARTIFACT	UNP Q92793
B	1080	MET	-	CLONING ARTIFACT	UNP Q92793

4 Residue-property plots

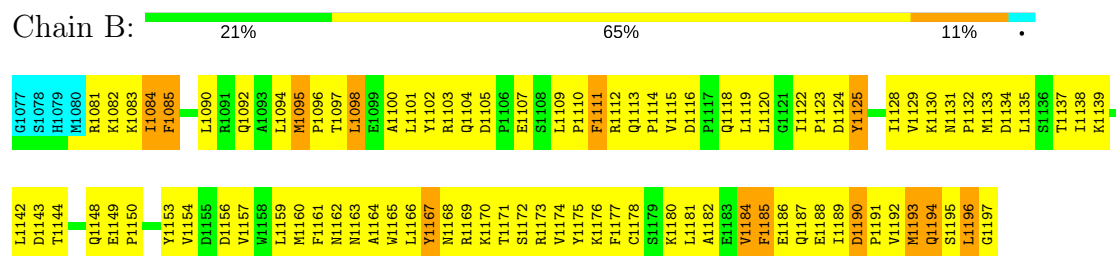
4.1 Average score per residue in the NMR ensemble

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

- Molecule 1: tumor protein p53



- Molecule 2: CREB-BINDING PROTEIN



4.2 Scores per residue for each member of the ensemble

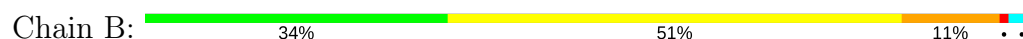
Colouring as in section 4.1 above.

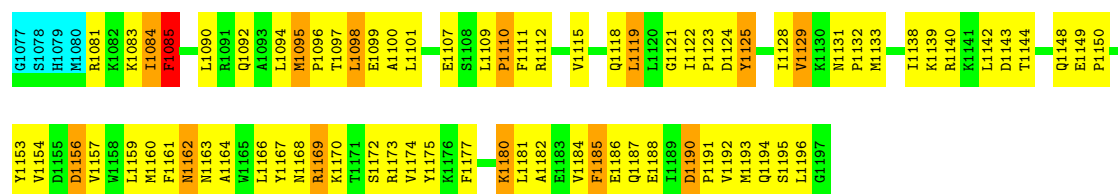
4.2.1 Score per residue for model 1

- Molecule 1: tumor protein p53



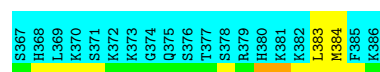
- Molecule 2: CREB-BINDING PROTEIN



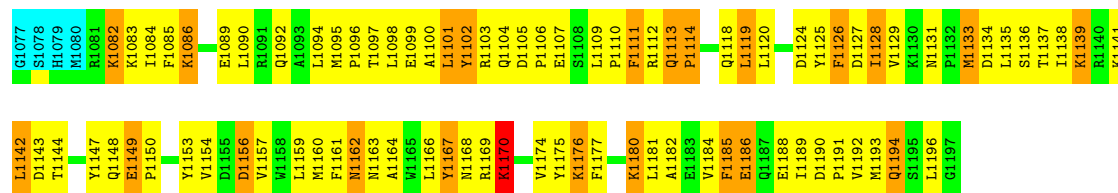


4.2.2 Score per residue for model 2

- Molecule 1: tumor protein p53

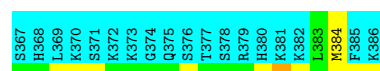


- Molecule 2: CREB-BINDING PROTEIN

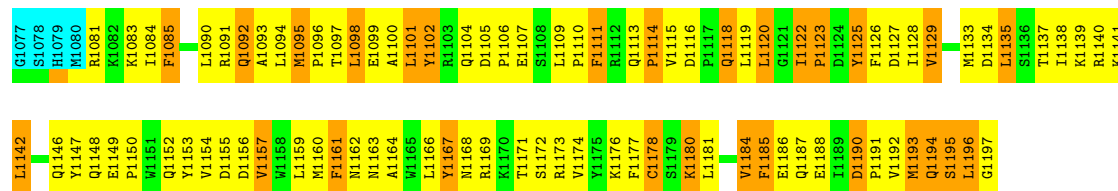


4.2.3 Score per residue for model 3 (medoid)

- Molecule 1: tumor protein p53



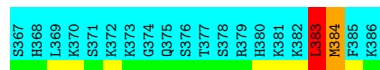
- Molecule 2: CREB-BINDING PROTEIN



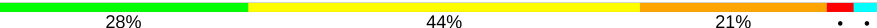
4.2.4 Score per residue for model 4

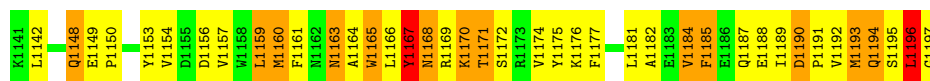
- Molecule 1: tumor protein p53

Chain A:  5% 5% 90%



- Molecule 2: CREB-BINDING PROTEIN

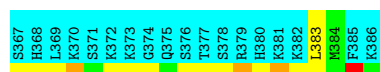
Chain B:  28% 44% 21%



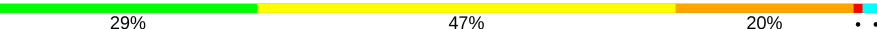
4.2.5 Score per residue for model 5

- Molecule 1: tumor protein p53

Chain A:  5% 5% 90%



- Molecule 2: CREB-BINDING PROTEIN

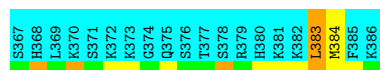
Chain B:  29% 47% 20%



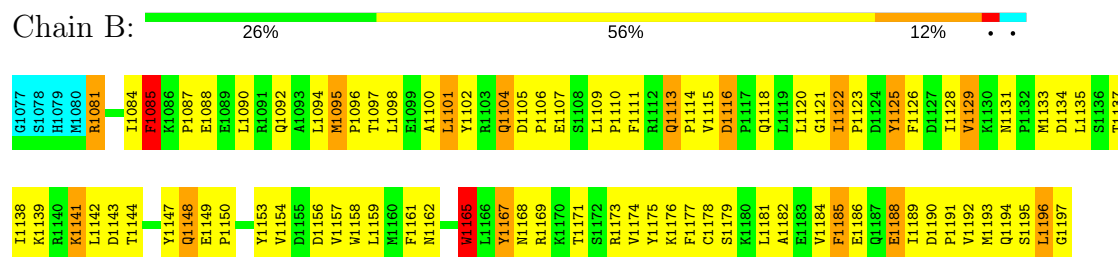
4.2.6 Score per residue for model 6

- Molecule 1: tumor protein p53

Chain A:  5% 5% 90%



- Molecule 2: CREB-BINDING PROTEIN

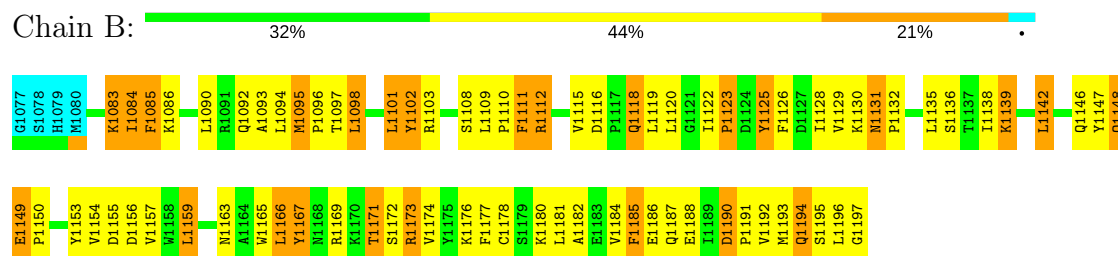


4.2.7 Score per residue for model 7

- Molecule 1: tumor protein p53



- Molecule 2: CREB-BINDING PROTEIN

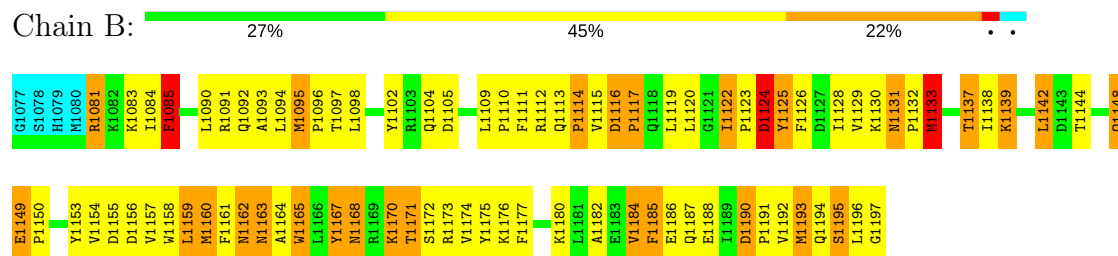


4.2.8 Score per residue for model 8

- Molecule 1: tumor protein p53



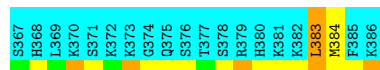
- Molecule 2: CREB-BINDING PROTEIN



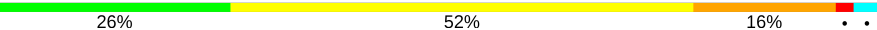
4.2.9 Score per residue for model 9

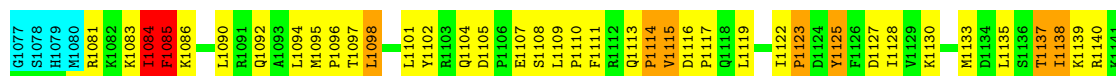
- Molecule 1: tumor protein p53

Chain A:  5% 5% 90%



- Molecule 2: CREB-BINDING PROTEIN

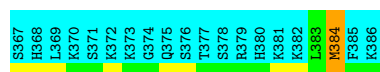
Chain B:  26% 52% 16% • •



4.2.10 Score per residue for model 10

- Molecule 1: tumor protein p53

Chain A:  5% 5% 90%



- Molecule 2: CREB-BINDING PROTEIN

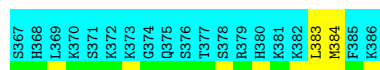
Chain B:  28% 51% 15% • •



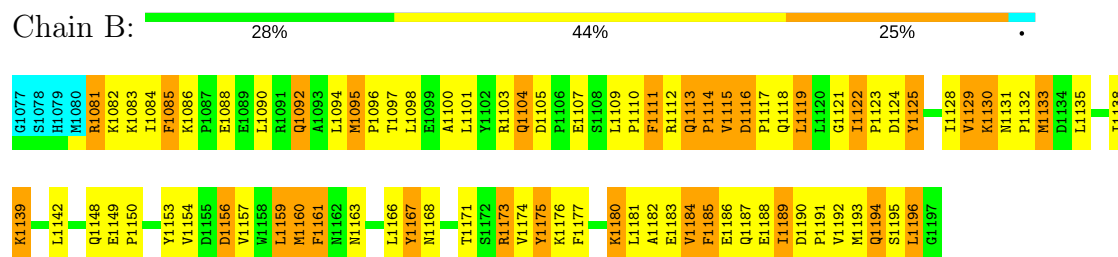
4.2.11 Score per residue for model 11

- Molecule 1: tumor protein p53

Chain A:  10% 90%



- Molecule 2: CREB-BINDING PROTEIN

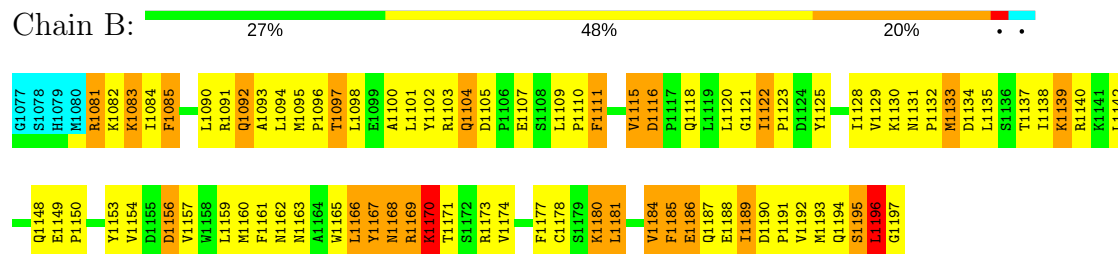


4.2.12 Score per residue for model 12

- Molecule 1: tumor protein p53



- Molecule 2: CREB-BINDING PROTEIN

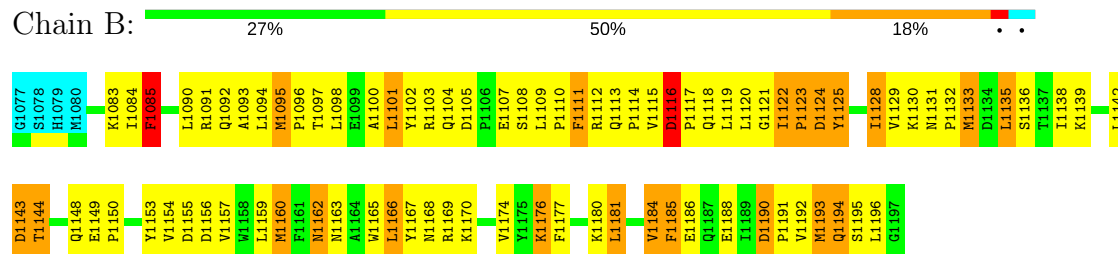


4.2.13 Score per residue for model 13

- Molecule 1: tumor protein p53



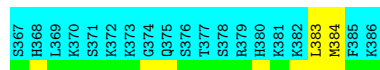
- Molecule 2: CREB-BINDING PROTEIN



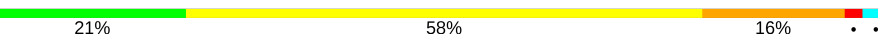
4.2.14 Score per residue for model 14

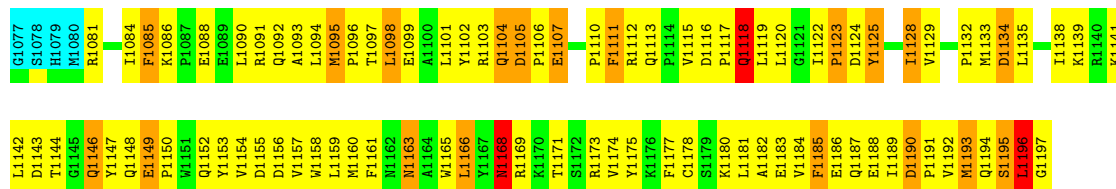
- Molecule 1: tumor protein p53

Chain A: 



- Molecule 2: CREB-BINDING PROTEIN

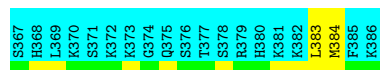
Chain B: 




4.2.15 Score per residue for model 15

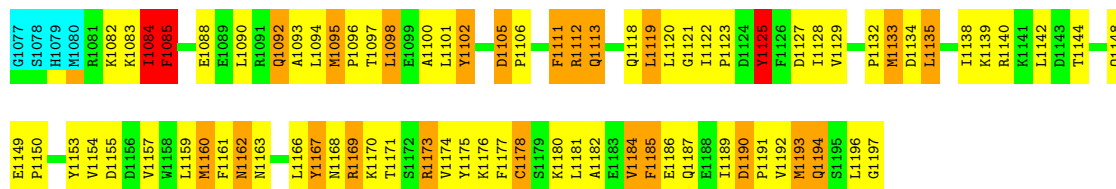
- Molecule 1: tumor protein p53

Chain A: 



- Molecule 2: CREB-BINDING PROTEIN

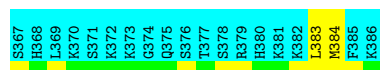
Chain B: 



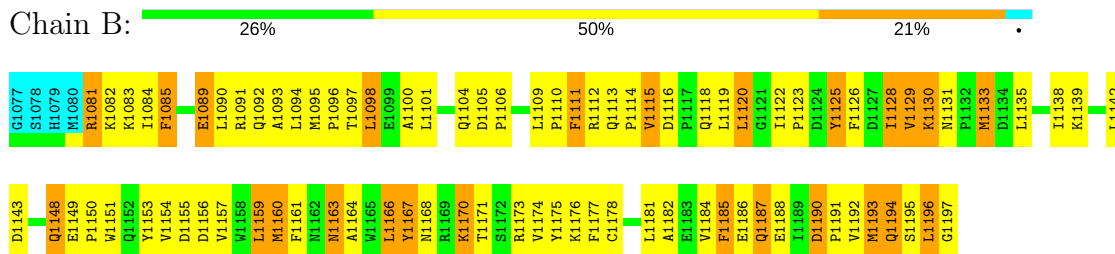
4.2.16 Score per residue for model 16

- Molecule 1: tumor protein p53

Chain A: 



- Molecule 2: CREB-BINDING PROTEIN

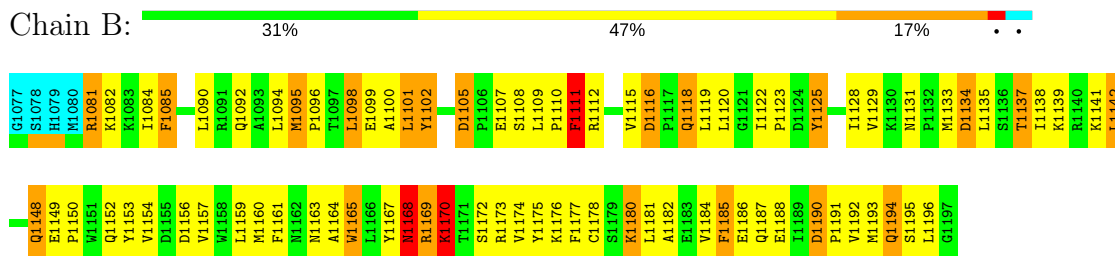


4.2.17 Score per residue for model 17

- Molecule 1: tumor protein p53



- Molecule 2: CREB-BINDING PROTEIN

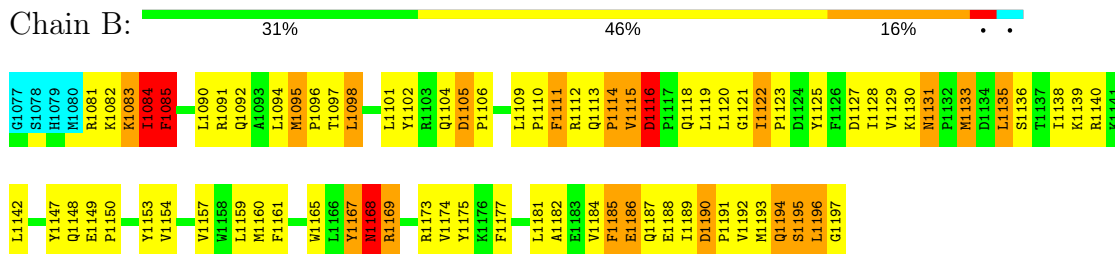


4.2.18 Score per residue for model 18

- Molecule 1: tumor protein p53



- Molecule 2: CREB-BINDING PROTEIN



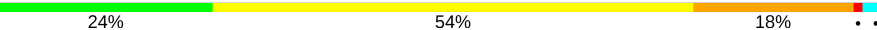
4.2.19 Score per residue for model 19

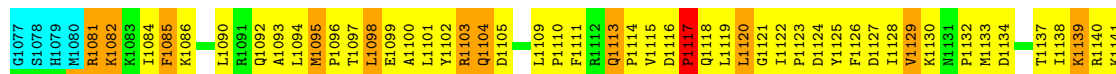
- Molecule 1: tumor protein p53

Chain A:  5% 5% 90%



- Molecule 2: CREB-BINDING PROTEIN

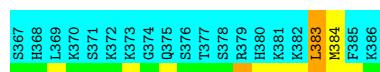
Chain B:  24% 54% 18%



4.2.20 Score per residue for model 20

- Molecule 1: tumor protein p53

Chain A:  5% 5% 90%



- Molecule 2: CREB-BINDING PROTEIN

Chain B:  26% 55% 16%



5 Refinement protocol and experimental data overview ⓘ

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

Of the 200 calculated structures, 20 were deposited, based on the following criterion: *structures with the least restraint violations, structures with the lowest energy*.

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

Software name	Classification	Version
X-PLOR	structure solution	3.1
X-PLOR	refinement	3.1

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

6 Model quality ⓘ

6.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ALY

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	16	20	20	3±3
2	B	987	980	980	102±9
All	All	20060	20000	20000	2048

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 51.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1098:LEU:HD21	2:B:1142:LEU:HD23	1.08	1.21	3	4
2:B:1098:LEU:HD11	2:B:1142:LEU:HD13	1.07	1.27	14	5
2:B:1098:LEU:HD21	2:B:1142:LEU:HD21	1.05	1.26	5	1
2:B:1098:LEU:HD21	2:B:1142:LEU:HD13	1.04	1.12	19	2
2:B:1122:ILE:HD11	2:B:1128:ILE:HD11	0.99	1.31	4	8
2:B:1098:LEU:HD11	2:B:1142:LEU:HD23	0.95	1.39	20	5
2:B:1133:MET:CE	2:B:1159:LEU:HD12	0.93	1.94	14	1
2:B:1100:ALA:HB2	2:B:1188:GLU:OE2	0.87	1.69	5	5
2:B:1098:LEU:CD2	2:B:1142:LEU:HD23	0.86	1.99	3	3
2:B:1189:ILE:HD12	2:B:1190:ASP:N	0.86	1.86	6	4
2:B:1128:ILE:HG21	2:B:1167:TYR:CD1	0.86	2.06	6	8

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1101:LEU:HA	2:B:1181:LEU:HD13	0.86	1.47	13	2
2:B:1117:PRO:CB	2:B:1119:LEU:HD23	0.86	2.01	19	1
2:B:1154:VAL:HG22	2:B:1194:GLN:NE2	0.85	1.85	12	1
2:B:1111:PHE:CG	2:B:1135:LEU:HD12	0.85	2.07	17	6
2:B:1168:ASN:ND2	2:B:1174:VAL:HG12	0.85	1.85	1	1
2:B:1117:PRO:HB2	2:B:1119:LEU:HD23	0.85	1.47	19	1
2:B:1142:LEU:HD21	2:B:1153:TYR:CE1	0.84	2.08	18	9
2:B:1098:LEU:CD1	2:B:1142:LEU:HD23	0.84	2.03	7	1
2:B:1128:ILE:HD12	2:B:1167:TYR:CD2	0.83	2.08	10	1
2:B:1084:ILE:HD13	2:B:1090:LEU:HD11	0.83	1.48	9	2
2:B:1098:LEU:HD21	2:B:1142:LEU:CD1	0.83	2.01	19	1
2:B:1111:PHE:CD2	2:B:1135:LEU:HD22	0.83	2.09	7	5
2:B:1168:ASN:OD1	2:B:1174:VAL:HG12	0.83	1.72	6	4
2:B:1115:VAL:HG13	2:B:1125:TYR:CE2	0.82	2.10	19	2
2:B:1094:LEU:HD11	2:B:1148:GLN:O	0.82	1.74	8	11
1:A:383:LEU:O	1:A:383:LEU:HD23	0.81	1.74	5	1
2:B:1184:VAL:HG22	2:B:1188:GLU:OE2	0.81	1.76	6	1
2:B:1101:LEU:HD21	2:B:1135:LEU:HD21	0.80	1.52	10	1
2:B:1128:ILE:HG22	2:B:1166:LEU:HD23	0.80	1.54	14	2
2:B:1142:LEU:HD11	2:B:1153:TYR:CE1	0.80	2.12	1	9
2:B:1108:SER:OG	2:B:1135:LEU:HD23	0.80	1.76	20	1
2:B:1098:LEU:CD1	2:B:1142:LEU:HD13	0.79	2.07	9	1
2:B:1188:GLU:O	2:B:1192:VAL:HG23	0.79	1.76	1	18
2:B:1101:LEU:HD21	2:B:1135:LEU:HD11	0.79	1.54	2	1
2:B:1142:LEU:HD11	2:B:1153:TYR:CD1	0.78	2.13	2	6
2:B:1125:TYR:CE1	2:B:1129:VAL:HG11	0.78	2.13	6	1
2:B:1139:LYS:HD2	2:B:1142:LEU:HD12	0.78	1.56	19	4
2:B:1154:VAL:HG11	2:B:1194:GLN:OE1	0.77	1.78	14	6
2:B:1181:LEU:HD23	2:B:1184:VAL:HG11	0.77	1.53	4	1
2:B:1090:LEU:HD22	2:B:1150:PRO:HD3	0.77	1.56	6	10
2:B:1098:LEU:HD11	2:B:1142:LEU:CD2	0.77	2.09	20	4
2:B:1098:LEU:HD22	2:B:1142:LEU:HD13	0.77	1.56	6	1
2:B:1083:LYS:C	2:B:1084:ILE:HD12	0.76	2.01	15	3
2:B:1084:ILE:HD12	2:B:1193:MET:HB2	0.76	1.57	3	8
1:A:383:LEU:CD2	2:B:1115:VAL:HG21	0.76	2.10	17	1
2:B:1100:ALA:HB1	2:B:1184:VAL:HG22	0.76	1.57	1	4
2:B:1138:ILE:O	2:B:1142:LEU:HD13	0.76	1.81	20	8
2:B:1138:ILE:HD12	2:B:1156:ASP:CB	0.75	2.12	16	13
2:B:1128:ILE:O	2:B:1166:LEU:HD22	0.74	1.82	16	2
2:B:1125:TYR:CE2	2:B:1129:VAL:HG11	0.74	2.18	8	2
2:B:1084:ILE:HG23	2:B:1193:MET:HB2	0.73	1.60	1	11

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1084:ILE:CG1	2:B:1090:LEU:HD21	0.73	2.12	13	20
2:B:1111:PHE:CE1	2:B:1181:LEU:HD12	0.73	2.18	9	1
2:B:1081:ARG:CG	2:B:1196:LEU:HD23	0.73	2.13	18	2
2:B:1180:LYS:O	2:B:1184:VAL:HG22	0.73	1.83	3	7
2:B:1138:ILE:HD12	2:B:1156:ASP:HB3	0.73	1.59	3	14
2:B:1115:VAL:HG13	2:B:1125:TYR:OH	0.73	1.83	3	4
2:B:1129:VAL:HG22	2:B:1163:ASN:CA	0.73	2.14	20	2
2:B:1138:ILE:HD12	2:B:1156:ASP:HB2	0.73	1.60	11	6
2:B:1100:ALA:HB2	2:B:1188:GLU:OE1	0.72	1.84	2	1
2:B:1119:LEU:O	2:B:1119:LEU:HD12	0.72	1.83	13	5
1:A:383:LEU:HD23	2:B:1115:VAL:HG23	0.72	1.59	16	1
2:B:1128:ILE:HG22	2:B:1166:LEU:HB3	0.72	1.60	2	1
2:B:1090:LEU:HD13	2:B:1149:GLU:HA	0.72	1.61	6	17
2:B:1154:VAL:HG21	2:B:1194:GLN:OE1	0.72	1.84	19	8
2:B:1122:ILE:HG12	2:B:1128:ILE:HD11	0.71	1.60	18	1
2:B:1098:LEU:HD21	2:B:1142:LEU:CD2	0.71	2.14	2	2
2:B:1131:ASN:O	2:B:1159:LEU:HD21	0.71	1.85	7	4
2:B:1115:VAL:HG21	2:B:1125:TYR:OH	0.70	1.86	11	1
2:B:1129:VAL:HG22	2:B:1163:ASN:CB	0.70	2.17	2	2
2:B:1083:LYS:C	2:B:1084:ILE:HD13	0.70	2.07	4	3
2:B:1098:LEU:HD22	2:B:1142:LEU:HD23	0.70	1.63	13	1
2:B:1128:ILE:HA	2:B:1166:LEU:HD23	0.70	1.62	15	4
2:B:1133:MET:SD	2:B:1160:MET:SD	0.70	2.89	18	1
2:B:1101:LEU:HA	2:B:1181:LEU:HD22	0.70	1.64	4	2
2:B:1138:ILE:HG23	2:B:1139:LYS:HD3	0.69	1.61	19	2
2:B:1098:LEU:HD13	2:B:1139:LYS:HE3	0.69	1.63	12	1
2:B:1139:LYS:HD3	2:B:1142:LEU:HD12	0.69	1.65	9	3
2:B:1138:ILE:HD11	2:B:1157:VAL:HG23	0.69	1.63	3	16
2:B:1119:LEU:HD12	2:B:1120:LEU:N	0.69	2.03	2	2
2:B:1098:LEU:HD13	2:B:1139:LYS:CE	0.69	2.18	12	1
2:B:1131:ASN:HB3	2:B:1159:LEU:HD11	0.69	1.64	4	2
2:B:1129:VAL:HG22	2:B:1163:ASN:HB2	0.68	1.65	2	1
2:B:1133:MET:HE1	2:B:1159:LEU:HD13	0.68	1.66	11	4
2:B:1122:ILE:HD12	2:B:1123:PRO:O	0.68	1.89	10	5
1:A:384:MET:HB2	2:B:1174:VAL:HG21	0.68	1.65	17	4
2:B:1093:ALA:HB1	2:B:1192:VAL:HG22	0.68	1.63	10	10
2:B:1081:ARG:HG3	2:B:1196:LEU:HD23	0.67	1.64	1	1
2:B:1131:ASN:O	2:B:1159:LEU:HD22	0.67	1.90	11	2
2:B:1119:LEU:HD12	2:B:1119:LEU:O	0.67	1.88	3	5
2:B:1128:ILE:HD12	2:B:1167:TYR:HB2	0.67	1.65	3	3
2:B:1115:VAL:HA	2:B:1119:LEU:HD11	0.67	1.66	13	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1111:PHE:CE1	2:B:1135:LEU:HD13	0.67	2.24	20	1
2:B:1189:ILE:HD13	2:B:1194:GLN:OE1	0.67	1.90	20	1
2:B:1129:VAL:HG22	2:B:1163:ASN:HA	0.66	1.64	20	6
2:B:1128:ILE:HG21	2:B:1167:TYR:CG	0.66	2.25	18	3
2:B:1162:ASN:O	2:B:1166:LEU:HD23	0.66	1.91	2	2
2:B:1109:LEU:O	2:B:1109:LEU:HD23	0.66	1.90	20	1
2:B:1098:LEU:HD11	2:B:1139:LYS:HD2	0.66	1.68	18	2
2:B:1125:TYR:CE1	2:B:1129:VAL:HG21	0.66	2.26	18	4
2:B:1115:VAL:HG11	2:B:1125:TYR:OH	0.66	1.91	11	2
2:B:1162:ASN:OD1	2:B:1166:LEU:HD22	0.66	1.90	15	1
2:B:1084:ILE:HD13	2:B:1084:ILE:N	0.66	2.05	10	5
2:B:1142:LEU:HD21	2:B:1153:TYR:CD1	0.66	2.25	10	4
2:B:1129:VAL:HG11	2:B:1132:PRO:HB3	0.66	1.65	7	1
2:B:1132:PRO:HA	2:B:1159:LEU:HD11	0.66	1.68	14	1
2:B:1119:LEU:HD12	2:B:1120:LEU:HB2	0.66	1.67	16	1
2:B:1196:LEU:HD23	2:B:1196:LEU:O	0.65	1.91	13	1
2:B:1115:VAL:HG21	2:B:1125:TYR:CE2	0.65	2.25	11	1
2:B:1159:LEU:O	2:B:1159:LEU:HD12	0.65	1.92	7	3
2:B:1098:LEU:HD11	2:B:1142:LEU:HD22	0.65	1.69	10	1
2:B:1133:MET:CE	2:B:1159:LEU:HD22	0.65	2.22	12	3
2:B:1084:ILE:HG12	2:B:1090:LEU:HD21	0.65	1.68	9	3
1:A:383:LEU:O	1:A:383:LEU:HD13	0.65	1.92	19	2
1:A:383:LEU:HD21	2:B:1115:VAL:HG11	0.65	1.68	17	1
2:B:1189:ILE:HD13	2:B:1195:SER:OG	0.65	1.92	6	2
2:B:1168:ASN:HB3	2:B:1174:VAL:HG12	0.64	1.66	8	2
2:B:1193:MET:CE	2:B:1196:LEU:HD22	0.64	2.22	11	2
2:B:1181:LEU:N	2:B:1181:LEU:HD23	0.64	2.08	10	1
2:B:1125:TYR:CD1	2:B:1129:VAL:HG11	0.64	2.27	6	1
1:A:383:LEU:HD12	2:B:1173:ARG:NH2	0.64	2.08	8	1
2:B:1193:MET:HE2	2:B:1196:LEU:HD13	0.64	1.70	2	6
2:B:1101:LEU:HD22	2:B:1135:LEU:HD21	0.64	1.70	3	1
2:B:1084:ILE:HG23	2:B:1192:VAL:O	0.63	1.92	10	3
2:B:1101:LEU:HD23	2:B:1139:LYS:NZ	0.63	2.09	14	2
2:B:1115:VAL:O	2:B:1115:VAL:HG22	0.63	1.93	17	3
2:B:1084:ILE:HG13	2:B:1090:LEU:HD21	0.63	1.69	14	15
2:B:1111:PHE:CD1	2:B:1135:LEU:HD12	0.63	2.29	15	1
2:B:1124:ASP:O	2:B:1128:ILE:HD11	0.63	1.94	2	2
1:A:384:MET:CB	2:B:1174:VAL:HG21	0.63	2.24	20	7
2:B:1129:VAL:HG11	2:B:1163:ASN:HB2	0.63	1.69	14	2
2:B:1101:LEU:CD2	2:B:1135:LEU:HD21	0.62	2.23	10	1
2:B:1101:LEU:HD13	2:B:1139:LYS:NZ	0.62	2.09	7	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1154:VAL:HG11	2:B:1194:GLN:NE2	0.62	2.09	20	2
1:A:383:LEU:HD21	2:B:1115:VAL:HG21	0.62	1.72	17	1
2:B:1090:LEU:HD22	2:B:1150:PRO:CD	0.62	2.23	4	2
2:B:1084:ILE:CB	2:B:1090:LEU:HD21	0.62	2.25	5	14
2:B:1084:ILE:N	2:B:1084:ILE:HD13	0.62	2.10	7	5
2:B:1090:LEU:C	2:B:1094:LEU:HD12	0.62	2.15	14	2
2:B:1139:LYS:HD3	2:B:1142:LEU:HD22	0.62	1.71	18	4
2:B:1111:PHE:CG	2:B:1135:LEU:HD22	0.62	2.29	7	2
2:B:1128:ILE:HG21	2:B:1167:TYR:CE1	0.62	2.28	1	2
2:B:1115:VAL:HG21	2:B:1125:TYR:CZ	0.62	2.28	11	1
2:B:1122:ILE:CG1	2:B:1128:ILE:HD11	0.62	2.25	15	2
2:B:1084:ILE:HB	2:B:1090:LEU:HD21	0.62	1.72	5	10
2:B:1115:VAL:HG22	2:B:1116:ASP:N	0.61	2.10	12	3
2:B:1181:LEU:N	2:B:1181:LEU:HD22	0.61	2.09	1	1
2:B:1129:VAL:HG13	2:B:1163:ASN:HB2	0.61	1.71	20	1
2:B:1122:ILE:HG13	2:B:1128:ILE:HD11	0.61	1.71	15	1
2:B:1115:VAL:O	2:B:1115:VAL:HG13	0.61	1.95	8	2
2:B:1154:VAL:HG21	2:B:1194:GLN:HB2	0.61	1.70	4	5
2:B:1193:MET:HE2	2:B:1196:LEU:HD22	0.61	1.71	11	1
2:B:1101:LEU:O	2:B:1181:LEU:HD22	0.61	1.94	19	8
2:B:1119:LEU:HD12	2:B:1120:LEU:CB	0.61	2.25	16	1
2:B:1101:LEU:HD13	2:B:1185:PHE:CB	0.61	2.25	18	3
1:A:383:LEU:CD2	2:B:1115:VAL:HG11	0.61	2.24	17	1
2:B:1101:LEU:O	2:B:1181:LEU:HD12	0.61	1.96	1	2
2:B:1128:ILE:HD12	2:B:1167:TYR:CG	0.61	2.30	10	1
2:B:1142:LEU:HD11	2:B:1153:TYR:HE1	0.61	1.52	1	3
2:B:1193:MET:CE	2:B:1196:LEU:HD13	0.60	2.26	17	7
2:B:1094:LEU:O	2:B:1098:LEU:HD23	0.60	1.97	5	4
2:B:1084:ILE:N	2:B:1084:ILE:HD12	0.60	2.10	15	2
2:B:1128:ILE:HD12	2:B:1167:TYR:CB	0.60	2.25	3	1
2:B:1168:ASN:CG	2:B:1174:VAL:HG12	0.60	2.16	2	2
2:B:1115:VAL:HG21	2:B:1122:ILE:CG2	0.60	2.27	5	2
1:A:384:MET:HB3	2:B:1174:VAL:HG21	0.60	1.72	20	2
2:B:1101:LEU:HD11	2:B:1135:LEU:HD21	0.59	1.73	13	1
2:B:1138:ILE:HG23	2:B:1139:LYS:HG2	0.59	1.74	20	4
2:B:1159:LEU:HD13	2:B:1159:LEU:C	0.59	2.17	14	1
2:B:1171:THR:O	2:B:1171:THR:HG23	0.59	1.97	8	4
2:B:1115:VAL:HG21	2:B:1122:ILE:HG23	0.59	1.72	6	3
2:B:1111:PHE:CD2	2:B:1135:LEU:HD12	0.59	2.32	17	2
2:B:1094:LEU:HD13	2:B:1148:GLN:HA	0.59	1.75	4	1
2:B:1100:ALA:HB1	2:B:1184:VAL:HG13	0.59	1.73	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:383:LEU:HD22	2:B:1115:VAL:HB	0.59	1.74	12	1
2:B:1115:VAL:HG23	2:B:1117:PRO:O	0.59	1.96	9	1
2:B:1111:PHE:CZ	2:B:1181:LEU:HD12	0.59	2.33	9	1
2:B:1161:PHE:CD2	2:B:1185:PHE:CE1	0.59	2.91	2	3
2:B:1101:LEU:HD13	2:B:1161:PHE:CZ	0.59	2.32	9	3
2:B:1139:LYS:CD	2:B:1142:LEU:HD12	0.59	2.27	9	3
2:B:1161:PHE:CE2	2:B:1185:PHE:CD1	0.59	2.91	6	10
2:B:1153:TYR:O	2:B:1157:VAL:HG23	0.59	1.98	6	1
2:B:1128:ILE:HG13	2:B:1129:VAL:HG12	0.58	1.75	3	2
2:B:1092:GLN:O	2:B:1096:PRO:CD	0.58	2.52	16	20
2:B:1092:GLN:O	2:B:1096:PRO:HD2	0.58	1.98	5	20
2:B:1171:THR:HG23	2:B:1171:THR:O	0.58	1.99	4	3
2:B:1100:ALA:HB1	2:B:1184:VAL:HB	0.58	1.75	12	4
2:B:1135:LEU:HD23	2:B:1138:ILE:CG2	0.58	2.28	9	1
2:B:1181:LEU:HD23	2:B:1184:VAL:CG1	0.58	2.29	4	1
2:B:1144:THR:HG23	2:B:1146:GLN:OE1	0.58	1.99	20	1
2:B:1122:ILE:HD11	2:B:1128:ILE:CD1	0.57	2.20	4	4
2:B:1189:ILE:HD12	2:B:1190:ASP:H	0.57	1.58	15	2
2:B:1101:LEU:CA	2:B:1181:LEU:HD13	0.57	2.25	13	1
2:B:1119:LEU:C	2:B:1119:LEU:HD12	0.57	2.20	16	1
2:B:1111:PHE:CB	2:B:1135:LEU:HD12	0.57	2.28	12	2
2:B:1090:LEU:O	2:B:1094:LEU:HD12	0.57	1.99	14	3
2:B:1189:ILE:HD12	2:B:1195:SER:OG	0.57	1.99	12	2
2:B:1122:ILE:CD1	2:B:1128:ILE:HD12	0.57	2.30	13	1
2:B:1101:LEU:HD13	2:B:1161:PHE:CE2	0.57	2.33	12	4
2:B:1101:LEU:HD13	2:B:1139:LYS:HZ1	0.57	1.59	7	1
1:A:383:LEU:HD13	2:B:1115:VAL:HB	0.57	1.77	12	1
2:B:1098:LEU:C	2:B:1098:LEU:HD12	0.57	2.19	6	1
2:B:1116:ASP:N	2:B:1117:PRO:HA	0.57	2.14	8	2
2:B:1122:ILE:HD12	2:B:1123:PRO:N	0.57	2.14	15	7
2:B:1161:PHE:CE2	2:B:1185:PHE:CD2	0.57	2.92	3	2
2:B:1128:ILE:HG21	2:B:1167:TYR:HB2	0.56	1.77	13	2
2:B:1173:ARG:HG2	2:B:1174:VAL:N	0.56	2.15	9	1
2:B:1114:PRO:O	2:B:1119:LEU:HD21	0.56	2.00	20	1
2:B:1098:LEU:CD2	2:B:1142:LEU:HD13	0.56	2.30	6	2
2:B:1190:ASP:N	2:B:1191:PRO:CD	0.56	2.69	2	19
2:B:1129:VAL:HG12	2:B:1132:PRO:HD3	0.56	1.78	10	4
2:B:1084:ILE:HD12	2:B:1084:ILE:N	0.56	2.15	9	1
2:B:1154:VAL:HG21	2:B:1194:GLN:CD	0.56	2.20	13	3
2:B:1181:LEU:N	2:B:1181:LEU:CD2	0.56	2.69	17	1
2:B:1122:ILE:HB	2:B:1123:PRO:HD2	0.56	1.78	15	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1100:ALA:HB2	2:B:1188:GLU:CD	0.55	2.22	11	2
2:B:1081:ARG:HG2	2:B:1196:LEU:HD23	0.55	1.78	18	1
2:B:1122:ILE:HD11	2:B:1128:ILE:HD12	0.55	1.77	13	1
2:B:1094:LEU:O	2:B:1098:LEU:HD12	0.55	2.01	2	2
2:B:1125:TYR:CZ	2:B:1129:VAL:HG21	0.55	2.37	8	5
2:B:1094:LEU:CD2	2:B:1153:TYR:CD2	0.55	2.90	12	18
2:B:1181:LEU:HD22	2:B:1181:LEU:N	0.55	2.17	7	2
2:B:1177:PHE:O	2:B:1181:LEU:HD23	0.55	2.01	1	2
2:B:1098:LEU:HD11	2:B:1142:LEU:CD1	0.55	2.25	9	1
2:B:1181:LEU:CD2	2:B:1181:LEU:N	0.55	2.69	7	5
2:B:1173:ARG:O	2:B:1175:TYR:N	0.55	2.40	9	1
2:B:1115:VAL:HG13	2:B:1116:ASP:N	0.54	2.17	16	1
2:B:1125:TYR:HD1	2:B:1128:ILE:HD11	0.54	1.63	10	1
2:B:1084:ILE:HD13	2:B:1090:LEU:HD21	0.54	1.79	18	1
2:B:1161:PHE:CE2	2:B:1185:PHE:CG	0.54	2.95	11	4
2:B:1189:ILE:HD12	2:B:1195:SER:CB	0.54	2.32	12	1
2:B:1164:ALA:N	2:B:1167:TYR:CE1	0.54	2.75	10	1
2:B:1159:LEU:C	2:B:1159:LEU:HD23	0.54	2.22	3	6
2:B:1129:VAL:O	2:B:1159:LEU:HD11	0.54	2.02	5	1
2:B:1170:LYS:O	2:B:1171:THR:HG23	0.54	2.02	16	1
2:B:1180:LYS:O	2:B:1184:VAL:HG12	0.54	2.03	1	3
2:B:1125:TYR:CE1	2:B:1129:VAL:CG2	0.54	2.91	18	2
2:B:1153:TYR:CE1	2:B:1157:VAL:HG21	0.53	2.37	9	14
2:B:1094:LEU:O	2:B:1098:LEU:HD13	0.53	2.03	10	3
2:B:1102:TYR:CD1	2:B:1139:LYS:CE	0.53	2.91	5	2
2:B:1111:PHE:HB3	2:B:1135:LEU:HD12	0.53	1.80	12	2
2:B:1102:TYR:CD2	2:B:1139:LYS:CE	0.53	2.92	2	1
2:B:1098:LEU:HD21	2:B:1142:LEU:HB3	0.53	1.79	10	1
2:B:1167:TYR:N	2:B:1167:TYR:CD1	0.53	2.76	10	1
2:B:1181:LEU:HA	2:B:1184:VAL:HG12	0.53	1.81	7	6
2:B:1162:ASN:C	2:B:1166:LEU:HD23	0.53	2.24	2	1
2:B:1084:ILE:CD1	2:B:1090:LEU:HD21	0.53	2.33	18	1
2:B:1129:VAL:HG22	2:B:1163:ASN:O	0.53	2.03	8	1
2:B:1122:ILE:CG1	2:B:1125:TYR:CD1	0.53	2.91	20	9
2:B:1135:LEU:HD23	2:B:1138:ILE:HG21	0.53	1.78	9	1
2:B:1185:PHE:CD1	2:B:1186:GLU:N	0.53	2.77	11	9
2:B:1161:PHE:CE1	2:B:1182:ALA:N	0.53	2.77	2	8
2:B:1185:PHE:CG	2:B:1186:GLU:N	0.53	2.76	12	8
2:B:1157:VAL:HG11	2:B:1185:PHE:CE2	0.53	2.39	1	7
2:B:1153:TYR:CE1	2:B:1157:VAL:CG2	0.53	2.92	15	14
2:B:1163:ASN:O	2:B:1164:ALA:HB2	0.53	2.01	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1122:ILE:C	2:B:1122:ILE:HD12	0.53	2.24	16	4
2:B:1133:MET:HE1	2:B:1159:LEU:HB3	0.53	1.80	9	4
2:B:1111:PHE:CE2	2:B:1160:MET:CG	0.53	2.92	2	1
2:B:1125:TYR:CD1	2:B:1129:VAL:CG1	0.53	2.92	6	1
2:B:1101:LEU:HD22	2:B:1139:LYS:CE	0.53	2.34	20	1
2:B:1157:VAL:CG1	2:B:1185:PHE:CE1	0.53	2.91	15	2
2:B:1111:PHE:CD1	2:B:1135:LEU:CD1	0.53	2.92	15	1
2:B:1116:ASP:N	2:B:1117:PRO:HD2	0.53	2.19	19	1
2:B:1100:ALA:HB2	2:B:1188:GLU:HG3	0.53	1.80	10	1
2:B:1109:LEU:N	2:B:1110:PRO:CD	0.53	2.71	16	6
2:B:1157:VAL:CG1	2:B:1185:PHE:CE2	0.53	2.92	8	9
2:B:1122:ILE:HD12	2:B:1122:ILE:C	0.53	2.24	10	5
2:B:1111:PHE:HD2	2:B:1135:LEU:HD22	0.53	1.64	2	1
2:B:1138:ILE:CD1	2:B:1157:VAL:HG23	0.52	2.34	16	8
2:B:1116:ASP:N	2:B:1117:PRO:CA	0.52	2.72	9	2
2:B:1121:GLY:O	2:B:1122:ILE:O	0.52	2.26	12	2
2:B:1122:ILE:HD11	2:B:1125:TYR:HD1	0.52	1.64	14	5
2:B:1122:ILE:N	2:B:1123:PRO:CA	0.52	2.72	7	2
2:B:1177:PHE:CD1	2:B:1178:CYS:N	0.52	2.77	20	5
2:B:1105:ASP:N	2:B:1106:PRO:HA	0.52	2.19	20	1
2:B:1167:TYR:CD1	2:B:1168:ASN:N	0.52	2.77	4	3
2:B:1100:ALA:HB3	2:B:1188:GLU:CD	0.52	2.24	6	1
2:B:1098:LEU:HD12	2:B:1098:LEU:C	0.52	2.24	13	1
2:B:1161:PHE:CE1	2:B:1182:ALA:HB2	0.52	2.39	10	3
1:A:384:MET:HE3	2:B:1167:TYR:O	0.52	2.04	12	1
2:B:1141:LYS:CB	2:B:1147:TYR:CE2	0.52	2.93	6	1
2:B:1098:LEU:HD12	2:B:1142:LEU:HD23	0.52	1.78	7	1
2:B:1113:GLN:CB	2:B:1114:PRO:CA	0.52	2.88	19	1
2:B:1129:VAL:HG21	2:B:1163:ASN:HD22	0.52	1.64	19	1
2:B:1129:VAL:HG11	2:B:1132:PRO:HG3	0.52	1.81	5	1
2:B:1113:GLN:N	2:B:1114:PRO:CD	0.52	2.73	11	5
2:B:1161:PHE:CE2	2:B:1185:PHE:CE1	0.52	2.98	9	7
1:A:384:MET:CA	2:B:1174:VAL:CG2	0.52	2.88	18	3
2:B:1170:LYS:C	2:B:1171:THR:HG23	0.52	2.25	20	1
2:B:1187:GLN:O	2:B:1191:PRO:CD	0.52	2.58	5	17
2:B:1129:VAL:CG2	2:B:1167:TYR:CE1	0.52	2.92	10	1
2:B:1100:ALA:HB3	2:B:1188:GLU:OE1	0.52	2.04	6	1
2:B:1128:ILE:CG2	2:B:1167:TYR:CD1	0.52	2.92	15	4
2:B:1102:TYR:CE1	2:B:1139:LYS:CB	0.52	2.92	20	1
2:B:1195:SER:O	2:B:1197:GLY:N	0.51	2.43	6	13
2:B:1141:LYS:O	2:B:1144:THR:HG22	0.51	2.05	20	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1105:ASP:O	2:B:1109:LEU:HD13	0.51	2.06	9	1
2:B:1128:ILE:HG22	2:B:1166:LEU:HD22	0.51	1.82	12	1
2:B:1142:LEU:HD21	2:B:1153:TYR:HE1	0.51	1.63	17	5
2:B:1101:LEU:CD1	2:B:1161:PHE:CZ	0.51	2.93	15	5
2:B:1159:LEU:HD23	2:B:1159:LEU:C	0.51	2.26	17	2
2:B:1098:LEU:HD12	2:B:1153:TYR:OH	0.51	2.06	8	1
2:B:1122:ILE:HD11	2:B:1125:TYR:CD1	0.51	2.40	14	4
2:B:1101:LEU:CD1	2:B:1161:PHE:CE2	0.51	2.93	12	3
2:B:1125:TYR:OH	2:B:1129:VAL:HG11	0.51	2.06	20	2
2:B:1168:ASN:ND2	2:B:1175:TYR:CD1	0.51	2.79	15	4
2:B:1122:ILE:N	2:B:1123:PRO:HA	0.51	2.21	17	2
2:B:1164:ALA:HA	2:B:1167:TYR:CE1	0.51	2.41	8	4
2:B:1115:VAL:HG13	2:B:1115:VAL:O	0.51	2.06	11	1
2:B:1163:ASN:O	2:B:1167:TYR:CD1	0.51	2.64	3	2
2:B:1162:ASN:OD1	2:B:1166:LEU:HD13	0.51	2.06	15	2
2:B:1098:LEU:HD21	2:B:1139:LYS:HE3	0.50	1.83	8	1
2:B:1097:THR:O	2:B:1100:ALA:HB3	0.50	2.05	13	1
2:B:1168:ASN:ND2	2:B:1175:TYR:CE2	0.50	2.80	11	1
2:B:1097:THR:HG22	2:B:1098:LEU:N	0.50	2.21	12	1
2:B:1102:TYR:CD1	2:B:1139:LYS:CG	0.50	2.94	20	1
2:B:1159:LEU:HD23	2:B:1159:LEU:O	0.50	2.05	17	2
2:B:1115:VAL:HG22	2:B:1115:VAL:O	0.50	2.06	7	1
2:B:1133:MET:HE3	2:B:1159:LEU:HD12	0.50	1.78	14	1
2:B:1153:TYR:CZ	2:B:1157:VAL:HG21	0.50	2.41	6	4
2:B:1106:PRO:HA	2:B:1109:LEU:HD12	0.50	1.82	6	2
1:A:384:MET:CB	2:B:1174:VAL:CG2	0.50	2.90	18	1
2:B:1125:TYR:CZ	2:B:1163:ASN:OD1	0.50	2.65	13	1
2:B:1090:LEU:HD13	2:B:1150:PRO:HD3	0.50	1.82	2	4
1:A:384:MET:CA	2:B:1174:VAL:HG21	0.49	2.37	7	6
2:B:1196:LEU:HD13	2:B:1196:LEU:C	0.49	2.27	14	2
1:A:384:MET:HA	2:B:1174:VAL:CG2	0.49	2.37	4	5
2:B:1102:TYR:CD1	2:B:1139:LYS:HG3	0.49	2.42	20	2
1:A:383:LEU:HD23	1:A:383:LEU:O	0.49	2.06	20	1
2:B:1160:MET:HE2	2:B:1160:MET:HA	0.49	1.84	8	2
2:B:1116:ASP:CB	2:B:1117:PRO:CD	0.49	2.90	14	2
2:B:1094:LEU:HD11	2:B:1148:GLN:HA	0.49	1.84	1	1
2:B:1167:TYR:CE1	2:B:1168:ASN:OD1	0.49	2.66	3	2
2:B:1139:LYS:HA	2:B:1142:LEU:HD22	0.49	1.84	20	1
2:B:1178:CYS:O	2:B:1182:ALA:HB2	0.49	2.06	20	2
2:B:1107:GLU:O	2:B:1181:LEU:HD12	0.49	2.08	5	3
2:B:1189:ILE:CG2	2:B:1194:GLN:NE2	0.49	2.76	12	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1181:LEU:HD23	2:B:1181:LEU:N	0.49	2.21	13	1
2:B:1108:SER:HB3	2:B:1181:LEU:HD11	0.49	1.85	7	2
2:B:1173:ARG:O	2:B:1177:PHE:CD2	0.49	2.66	17	5
2:B:1138:ILE:HD11	2:B:1157:VAL:CG2	0.49	2.37	18	4
2:B:1111:PHE:CD1	2:B:1135:LEU:HD22	0.49	2.43	20	1
2:B:1129:VAL:HG21	2:B:1132:PRO:HB3	0.49	1.85	19	1
2:B:1173:ARG:O	2:B:1177:PHE:CE2	0.49	2.66	11	5
2:B:1097:THR:HG23	2:B:1098:LEU:N	0.49	2.23	13	2
2:B:1125:TYR:CD1	2:B:1128:ILE:HD11	0.49	2.43	10	1
2:B:1115:VAL:O	2:B:1125:TYR:CD2	0.49	2.66	14	1
2:B:1125:TYR:CE2	2:B:1163:ASN:ND2	0.49	2.80	16	1
2:B:1122:ILE:HD11	2:B:1128:ILE:HD13	0.49	1.84	15	1
2:B:1119:LEU:O	2:B:1119:LEU:CD1	0.49	2.60	19	1
2:B:1175:TYR:CG	2:B:1175:TYR:O	0.49	2.66	17	3
2:B:1168:ASN:CB	2:B:1174:VAL:HG12	0.48	2.38	9	2
2:B:1102:TYR:CD1	2:B:1139:LYS:HE2	0.48	2.43	5	3
2:B:1102:TYR:CD2	2:B:1139:LYS:HE3	0.48	2.43	2	1
2:B:1184:VAL:CG1	2:B:1185:PHE:N	0.48	2.76	4	4
2:B:1111:PHE:CE2	2:B:1181:LEU:HD12	0.48	2.43	16	3
2:B:1161:PHE:CE1	2:B:1182:ALA:CA	0.48	2.96	18	11
2:B:1132:PRO:CB	2:B:1163:ASN:ND2	0.48	2.77	13	1
2:B:1133:MET:CE	2:B:1159:LEU:CD2	0.48	2.91	12	1
2:B:1111:PHE:CE2	2:B:1181:LEU:CD1	0.48	2.96	15	1
2:B:1185:PHE:CD2	2:B:1186:GLU:N	0.48	2.80	15	1
2:B:1143:ASP:O	2:B:1144:THR:HG23	0.48	2.09	13	2
2:B:1184:VAL:HG13	2:B:1185:PHE:N	0.48	2.23	14	5
2:B:1105:ASP:N	2:B:1106:PRO:CA	0.48	2.76	20	1
2:B:1157:VAL:HG11	2:B:1185:PHE:CD2	0.48	2.43	14	3
2:B:1113:GLN:N	2:B:1114:PRO:HD3	0.48	2.24	11	7
2:B:1137:THR:HG22	2:B:1138:ILE:N	0.48	2.23	10	5
2:B:1174:VAL:O	2:B:1177:PHE:CZ	0.48	2.67	2	6
2:B:1150:PRO:O	2:B:1154:VAL:HG23	0.48	2.08	9	15
2:B:1081:ARG:O	2:B:1083:LYS:N	0.48	2.46	12	1
2:B:1109:LEU:CB	2:B:1110:PRO:CD	0.48	2.92	11	9
2:B:1133:MET:N	2:B:1133:MET:CE	0.48	2.76	8	1
2:B:1167:TYR:CG	2:B:1167:TYR:O	0.48	2.67	3	2
2:B:1159:LEU:O	2:B:1159:LEU:HD23	0.48	2.09	18	3
2:B:1150:PRO:CB	2:B:1194:GLN:CB	0.48	2.91	12	1
2:B:1194:GLN:CG	2:B:1195:SER:N	0.48	2.76	6	2
2:B:1125:TYR:CZ	2:B:1129:VAL:HG11	0.48	2.44	18	5
2:B:1161:PHE:CD1	2:B:1182:ALA:HB2	0.48	2.43	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1111:PHE:CE2	2:B:1181:LEU:HD13	0.48	2.44	15	1
2:B:1175:TYR:O	2:B:1175:TYR:CG	0.48	2.67	9	5
2:B:1133:MET:O	2:B:1137:THR:HG22	0.48	2.09	2	1
2:B:1102:TYR:CG	2:B:1139:LYS:HE2	0.48	2.44	15	1
2:B:1160:MET:O	2:B:1160:MET:HE2	0.48	2.09	10	1
2:B:1170:LYS:C	2:B:1171:THR:HG22	0.48	2.29	15	1
2:B:1102:TYR:CD2	2:B:1135:LEU:O	0.48	2.67	4	1
2:B:1081:ARG:CD	2:B:1196:LEU:CB	0.48	2.91	19	1
2:B:1194:GLN:HG3	2:B:1195:SER:N	0.47	2.23	6	1
2:B:1094:LEU:HD22	2:B:1153:TYR:CD2	0.47	2.44	13	1
2:B:1115:VAL:HG22	2:B:1121:GLY:HA2	0.47	1.86	1	1
2:B:1125:TYR:CZ	2:B:1163:ASN:ND2	0.47	2.82	16	1
2:B:1128:ILE:C	2:B:1129:VAL:HG23	0.47	2.28	11	6
2:B:1098:LEU:HD11	2:B:1139:LYS:CE	0.47	2.39	1	1
2:B:1122:ILE:CD1	2:B:1128:ILE:HD11	0.47	2.39	7	2
2:B:1182:ALA:O	2:B:1185:PHE:CD1	0.47	2.67	20	1
2:B:1115:VAL:O	2:B:1125:TYR:CE2	0.47	2.67	14	2
2:B:1165:TRP:N	2:B:1168:ASN:ND2	0.47	2.61	9	2
2:B:1181:LEU:HA	2:B:1184:VAL:HG23	0.47	1.85	13	1
2:B:1137:THR:CG2	2:B:1138:ILE:N	0.47	2.77	9	4
1:A:383:LEU:CD2	2:B:1115:VAL:CG2	0.47	2.90	17	1
2:B:1095:MET:CB	2:B:1096:PRO:CD	0.47	2.92	4	2
2:B:1084:ILE:O	2:B:1085:PHE:CG	0.47	2.68	8	6
1:A:383:LEU:CG	2:B:1115:VAL:HG21	0.47	2.39	17	1
2:B:1108:SER:HA	2:B:1181:LEU:HD11	0.47	1.85	9	1
2:B:1167:TYR:CD1	2:B:1167:TYR:C	0.47	2.86	3	4
2:B:1084:ILE:O	2:B:1085:PHE:CD2	0.47	2.67	6	4
2:B:1133:MET:CG	2:B:1134:ASP:N	0.47	2.77	14	5
2:B:1113:GLN:CB	2:B:1114:PRO:HA	0.47	2.39	19	1
2:B:1117:PRO:O	2:B:1119:LEU:N	0.47	2.48	8	1
2:B:1154:VAL:CG1	2:B:1194:GLN:OE1	0.47	2.61	6	1
2:B:1174:VAL:O	2:B:1176:LYS:N	0.47	2.46	9	1
2:B:1150:PRO:O	2:B:1154:VAL:N	0.47	2.46	12	12
2:B:1084:ILE:N	2:B:1084:ILE:CD1	0.47	2.76	4	4
2:B:1122:ILE:CG1	2:B:1125:TYR:CE1	0.47	2.98	1	2
2:B:1182:ALA:HA	2:B:1185:PHE:CD1	0.47	2.45	1	4
2:B:1110:PRO:O	2:B:1177:PHE:CD1	0.47	2.68	14	2
2:B:1138:ILE:HG23	2:B:1139:LYS:N	0.47	2.24	7	3
2:B:1175:TYR:O	2:B:1175:TYR:CD2	0.47	2.68	1	1
2:B:1175:TYR:CD2	2:B:1175:TYR:O	0.47	2.67	9	3
2:B:1196:LEU:C	2:B:1196:LEU:HD13	0.47	2.30	16	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1169:ARG:CG	2:B:1170:LYS:N	0.47	2.78	17	1
2:B:1150:PRO:HB3	2:B:1194:GLN:CB	0.47	2.40	18	10
2:B:1094:LEU:HA	2:B:1097:THR:CG2	0.47	2.40	20	17
2:B:1122:ILE:HG13	2:B:1125:TYR:CD1	0.47	2.45	6	4
2:B:1122:ILE:HG12	2:B:1125:TYR:CD1	0.47	2.44	5	2
2:B:1115:VAL:HG13	2:B:1116:ASP:H	0.46	1.70	5	5
2:B:1137:THR:HG23	2:B:1138:ILE:N	0.46	2.25	2	1
2:B:1108:SER:CA	2:B:1181:LEU:HD11	0.46	2.41	9	1
2:B:1122:ILE:O	2:B:1124:ASP:N	0.46	2.48	14	2
2:B:1164:ALA:O	2:B:1168:ASN:N	0.46	2.48	10	1
2:B:1107:GLU:CB	2:B:1181:LEU:HD21	0.46	2.40	10	1
2:B:1172:SER:O	2:B:1174:VAL:N	0.46	2.48	7	1
2:B:1116:ASP:O	2:B:1118:GLN:N	0.46	2.47	19	3
2:B:1187:GLN:HG2	2:B:1188:GLU:N	0.46	2.25	20	1
2:B:1116:ASP:CB	2:B:1117:PRO:HD3	0.46	2.41	14	1
2:B:1141:LYS:HD2	2:B:1147:TYR:CZ	0.46	2.45	14	1
2:B:1157:VAL:HG11	2:B:1185:PHE:CE1	0.46	2.46	3	2
2:B:1115:VAL:HG11	2:B:1167:TYR:CE2	0.46	2.45	3	1
2:B:1102:TYR:CG	2:B:1139:LYS:CE	0.46	2.99	15	1
2:B:1129:VAL:HG13	2:B:1129:VAL:O	0.46	2.10	19	1
2:B:1094:LEU:HD11	2:B:1149:GLU:N	0.46	2.25	6	1
2:B:1165:TRP:HA	2:B:1168:ASN:ND2	0.46	2.26	12	1
2:B:1129:VAL:CG1	2:B:1163:ASN:ND2	0.46	2.78	4	1
1:A:383:LEU:HD22	2:B:1173:ARG:NH1	0.46	2.25	6	1
1:A:384:MET:HE1	2:B:1169:ARG:CD	0.46	2.41	1	1
2:B:1129:VAL:HG21	2:B:1163:ASN:ND2	0.46	2.25	19	1
2:B:1190:ASP:N	2:B:1191:PRO:HD2	0.46	2.25	16	15
2:B:1098:LEU:CD2	2:B:1139:LYS:NZ	0.46	2.79	11	1
2:B:1101:LEU:HD13	2:B:1185:PHE:HB3	0.46	1.87	18	1
2:B:1110:PRO:HG3	2:B:1177:PHE:CG	0.46	2.46	19	1
2:B:1168:ASN:O	2:B:1169:ARG:CG	0.46	2.64	18	2
2:B:1114:PRO:C	2:B:1119:LEU:HD21	0.46	2.31	13	1
2:B:1121:GLY:O	2:B:1122:ILE:HG13	0.46	2.10	13	2
2:B:1135:LEU:O	2:B:1139:LYS:CG	0.46	2.64	3	4
2:B:1122:ILE:CD1	2:B:1122:ILE:C	0.46	2.85	17	2
2:B:1094:LEU:CD1	2:B:1148:GLN:O	0.46	2.63	14	2
2:B:1133:MET:O	2:B:1134:ASP:CB	0.46	2.64	17	2
2:B:1157:VAL:HG12	2:B:1185:PHE:CE2	0.46	2.46	2	1
2:B:1133:MET:HE2	2:B:1159:LEU:HD12	0.46	1.81	14	1
2:B:1135:LEU:O	2:B:1139:LYS:HG2	0.46	2.11	18	1
2:B:1116:ASP:N	2:B:1117:PRO:CD	0.46	2.79	19	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1173:ARG:HG3	2:B:1174:VAL:HG23	0.46	1.88	8	1
2:B:1192:VAL:O	2:B:1193:MET:CB	0.46	2.63	17	13
2:B:1094:LEU:CD1	2:B:1148:GLN:HA	0.46	2.41	20	6
2:B:1150:PRO:CB	2:B:1194:GLN:HB2	0.46	2.41	12	2
1:A:384:MET:CG	1:A:384:MET:O	0.45	2.64	20	1
2:B:1111:PHE:CZ	2:B:1178:CYS:SG	0.45	3.10	16	2
2:B:1115:VAL:HG13	2:B:1125:TYR:CZ	0.45	2.45	19	1
2:B:1122:ILE:HD12	2:B:1124:ASP:N	0.45	2.26	8	1
2:B:1115:VAL:CG2	2:B:1125:TYR:CE2	0.45	2.99	11	1
2:B:1174:VAL:HA	2:B:1177:PHE:CE2	0.45	2.46	9	14
2:B:1125:TYR:O	2:B:1126:PHE:CB	0.45	2.64	3	2
2:B:1118:GLN:O	2:B:1120:LEU:N	0.45	2.49	19	2
2:B:1102:TYR:CB	2:B:1139:LYS:CE	0.45	2.94	6	2
2:B:1115:VAL:HG21	2:B:1122:ILE:HG21	0.45	1.88	12	1
2:B:1107:GLU:O	2:B:1111:PHE:CE1	0.45	2.69	5	1
2:B:1169:ARG:O	2:B:1175:TYR:CD2	0.45	2.69	15	1
2:B:1083:LYS:O	2:B:1085:PHE:N	0.45	2.49	15	6
2:B:1084:ILE:CD1	2:B:1084:ILE:N	0.45	2.78	9	2
2:B:1138:ILE:O	2:B:1142:LEU:CB	0.45	2.65	5	1
2:B:1149:GLU:O	2:B:1149:GLU:CG	0.45	2.64	9	1
2:B:1125:TYR:CE2	2:B:1129:VAL:CG1	0.45	2.96	8	1
2:B:1095:MET:CB	2:B:1096:PRO:HD3	0.45	2.42	13	13
2:B:1084:ILE:CG1	2:B:1150:PRO:HG2	0.45	2.41	5	9
2:B:1169:ARG:O	2:B:1171:THR:N	0.45	2.49	9	1
2:B:1122:ILE:HB	2:B:1123:PRO:CD	0.45	2.41	12	3
2:B:1160:MET:O	2:B:1160:MET:CE	0.45	2.65	19	2
2:B:1093:ALA:O	2:B:1097:THR:HG22	0.45	2.12	16	2
2:B:1160:MET:HE3	2:B:1160:MET:HA	0.45	1.86	5	1
2:B:1081:ARG:HD3	2:B:1196:LEU:HD23	0.45	1.87	17	1
2:B:1101:LEU:C	2:B:1101:LEU:HD23	0.45	2.32	20	1
2:B:1161:PHE:CE2	2:B:1185:PHE:CE2	0.45	3.04	19	1
2:B:1094:LEU:O	2:B:1097:THR:HG22	0.45	2.11	1	3
2:B:1100:ALA:CB	2:B:1184:VAL:HG22	0.45	2.36	1	1
2:B:1102:TYR:CD2	2:B:1139:LYS:HG3	0.45	2.47	4	4
2:B:1096:PRO:O	2:B:1099:GLU:CG	0.45	2.65	20	1
2:B:1117:PRO:O	2:B:1118:GLN:CB	0.45	2.64	14	1
2:B:1098:LEU:HD11	2:B:1142:LEU:CG	0.45	2.42	16	1
2:B:1101:LEU:HD11	2:B:1157:VAL:HG13	0.45	1.88	17	2
2:B:1109:LEU:N	2:B:1110:PRO:HD2	0.45	2.27	19	4
2:B:1189:ILE:HD12	2:B:1195:SER:HB2	0.45	1.89	12	1
2:B:1139:LYS:HD2	2:B:1142:LEU:HD23	0.45	1.88	5	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1116:ASP:HB3	2:B:1117:PRO:CD	0.45	2.42	14	1
1:A:383:LEU:CD2	2:B:1115:VAL:HG23	0.45	2.38	16	1
2:B:1082:LYS:O	2:B:1084:ILE:CD1	0.44	2.65	12	6
2:B:1115:VAL:CG2	2:B:1116:ASP:N	0.44	2.80	12	1
2:B:1119:LEU:CD1	2:B:1119:LEU:O	0.44	2.65	17	1
2:B:1111:PHE:CE1	2:B:1178:CYS:SG	0.44	3.10	14	2
2:B:1150:PRO:HA	2:B:1153:TYR:HB3	0.44	1.88	9	1
2:B:1128:ILE:CD1	2:B:1167:TYR:CD2	0.44	3.00	9	1
1:A:383:LEU:HD11	2:B:1114:PRO:C	0.44	2.31	18	1
2:B:1111:PHE:CE1	2:B:1160:MET:SD	0.44	3.10	4	1
2:B:1157:VAL:O	2:B:1161:PHE:N	0.44	2.51	11	3
2:B:1107:GLU:HB3	2:B:1181:LEU:HD21	0.44	1.87	1	1
2:B:1133:MET:HB2	2:B:1137:THR:HG21	0.44	1.88	17	1
2:B:1116:ASP:HB3	2:B:1117:PRO:HD3	0.44	1.88	14	1
2:B:1083:LYS:O	2:B:1084:ILE:C	0.44	2.55	18	3
2:B:1133:MET:HE2	2:B:1133:MET:N	0.44	2.28	19	1
2:B:1094:LEU:CD2	2:B:1153:TYR:CE2	0.44	3.00	11	2
2:B:1157:VAL:HG12	2:B:1161:PHE:HD2	0.44	1.72	11	1
2:B:1129:VAL:O	2:B:1130:LYS:CB	0.44	2.66	16	2
2:B:1109:LEU:CB	2:B:1110:PRO:HD3	0.44	2.43	9	9
2:B:1143:ASP:C	2:B:1144:THR:HG23	0.44	2.32	10	1
2:B:1125:TYR:OH	2:B:1163:ASN:ND2	0.44	2.50	16	6
2:B:1190:ASP:HB2	2:B:1191:PRO:HD3	0.44	1.88	12	1
2:B:1107:GLU:O	2:B:1181:LEU:HD11	0.44	2.11	2	2
2:B:1181:LEU:HA	2:B:1184:VAL:CG2	0.44	2.42	12	2
2:B:1110:PRO:HG2	2:B:1177:PHE:CD1	0.44	2.47	19	2
1:A:384:MET:HE2	2:B:1167:TYR:O	0.44	2.12	16	1
2:B:1170:LYS:O	2:B:1171:THR:CG2	0.44	2.65	15	1
2:B:1115:VAL:HG22	2:B:1117:PRO:HB3	0.44	1.88	11	1
2:B:1160:MET:CE	2:B:1160:MET:HA	0.44	2.43	11	4
2:B:1168:ASN:CG	2:B:1175:TYR:CD2	0.44	2.91	11	1
2:B:1090:LEU:CB	2:B:1148:GLN:O	0.44	2.66	7	6
2:B:1129:VAL:HG23	2:B:1163:ASN:HB2	0.44	1.89	3	1
2:B:1159:LEU:C	2:B:1159:LEU:HD12	0.44	2.31	10	2
2:B:1196:LEU:CD1	2:B:1196:LEU:C	0.44	2.85	4	2
2:B:1115:VAL:HA	2:B:1119:LEU:HD21	0.44	1.90	16	1
2:B:1133:MET:O	2:B:1160:MET:CE	0.44	2.66	11	3
2:B:1094:LEU:O	2:B:1098:LEU:CD2	0.44	2.66	6	3
2:B:1165:TRP:CD1	2:B:1165:TRP:N	0.44	2.84	13	1
2:B:1084:ILE:HD12	2:B:1193:MET:CB	0.44	2.42	1	1
2:B:1169:ARG:O	2:B:1170:LYS:CG	0.44	2.66	12	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1168:ASN:OD1	2:B:1174:VAL:CG1	0.44	2.66	9	2
2:B:1108:SER:OG	2:B:1109:LEU:N	0.44	2.50	17	1
2:B:1190:ASP:HB3	2:B:1191:PRO:HD3	0.44	1.90	11	1
2:B:1141:LYS:HB3	2:B:1147:TYR:CE2	0.44	2.47	6	1
2:B:1146:GLN:C	2:B:1147:TYR:CD1	0.44	2.92	10	2
2:B:1122:ILE:CG1	2:B:1128:ILE:CD1	0.44	2.96	15	1
2:B:1150:PRO:O	2:B:1154:VAL:CB	0.43	2.66	11	5
2:B:1116:ASP:HB2	2:B:1117:PRO:HD2	0.43	1.90	13	1
2:B:1111:PHE:CE1	2:B:1181:LEU:HG	0.43	2.48	7	1
2:B:1123:PRO:O	2:B:1125:TYR:N	0.43	2.51	20	2
2:B:1131:ASN:CB	2:B:1159:LEU:HD11	0.43	2.42	18	1
2:B:1116:ASP:OD1	2:B:1132:PRO:CG	0.43	2.66	19	1
2:B:1084:ILE:HG13	2:B:1090:LEU:CD2	0.43	2.43	13	6
2:B:1102:TYR:CB	2:B:1139:LYS:HE3	0.43	2.43	6	2
2:B:1081:ARG:NH2	2:B:1195:SER:O	0.43	2.51	6	2
2:B:1129:VAL:HG11	2:B:1163:ASN:CB	0.43	2.43	13	1
2:B:1095:MET:O	2:B:1098:LEU:CB	0.43	2.66	9	5
2:B:1133:MET:O	2:B:1137:THR:CG2	0.43	2.66	2	1
2:B:1129:VAL:CG2	2:B:1163:ASN:O	0.43	2.67	8	3
2:B:1188:GLU:O	2:B:1192:VAL:CG2	0.43	2.66	16	10
2:B:1098:LEU:O	2:B:1102:TYR:N	0.43	2.51	10	9
2:B:1167:TYR:C	2:B:1167:TYR:CD1	0.43	2.91	11	1
2:B:1164:ALA:O	2:B:1168:ASN:CB	0.43	2.67	10	1
2:B:1107:GLU:HB3	2:B:1181:LEU:CD2	0.43	2.43	12	2
1:A:383:LEU:HD23	1:A:383:LEU:C	0.43	2.34	5	1
2:B:1160:MET:HA	2:B:1160:MET:CE	0.43	2.43	5	3
2:B:1107:GLU:C	2:B:1181:LEU:HD11	0.43	2.33	14	1
2:B:1193:MET:HB3	2:B:1196:LEU:HD23	0.43	1.91	14	1
1:A:383:LEU:O	1:A:384:MET:CG	0.43	2.66	18	1
2:B:1142:LEU:CD1	2:B:1153:TYR:CD1	0.43	3.01	16	1
2:B:1133:MET:HE1	2:B:1159:LEU:HD22	0.43	1.90	4	1
2:B:1115:VAL:O	2:B:1115:VAL:CG1	0.43	2.67	8	1
2:B:1158:TRP:O	2:B:1162:ASN:N	0.43	2.51	8	1
2:B:1092:GLN:O	2:B:1095:MET:N	0.43	2.50	7	8
2:B:1150:PRO:O	2:B:1154:VAL:CG2	0.43	2.66	12	12
2:B:1102:TYR:HB2	2:B:1139:LYS:CE	0.43	2.43	18	3
2:B:1105:ASP:OD2	2:B:1108:SER:N	0.43	2.51	13	1
2:B:1112:ARG:O	2:B:1112:ARG:CG	0.43	2.66	7	1
2:B:1099:GLU:O	2:B:1103:ARG:CG	0.43	2.66	2	3
2:B:1122:ILE:HD12	2:B:1123:PRO:CD	0.43	2.44	15	1
2:B:1102:TYR:CE2	2:B:1135:LEU:O	0.43	2.71	4	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1131:ASN:O	2:B:1159:LEU:CD2	0.43	2.67	12	5
2:B:1167:TYR:CD2	2:B:1167:TYR:O	0.43	2.71	13	1
2:B:1101:LEU:O	2:B:1181:LEU:CD1	0.43	2.67	12	2
2:B:1098:LEU:CD1	2:B:1139:LYS:CE	0.43	2.96	1	1
2:B:1190:ASP:CB	2:B:1191:PRO:CD	0.43	2.97	12	2
2:B:1119:LEU:O	2:B:1119:LEU:CG	0.43	2.66	19	1
2:B:1164:ALA:C	2:B:1167:TYR:CE1	0.43	2.91	8	2
2:B:1185:PHE:O	2:B:1189:ILE:CG1	0.43	2.67	20	2
2:B:1155:ASP:O	2:B:1159:LEU:CB	0.43	2.66	5	3
2:B:1131:ASN:O	2:B:1133:MET:CE	0.43	2.67	20	2
2:B:1163:ASN:ND2	2:B:1167:TYR:OH	0.43	2.51	10	1
2:B:1125:TYR:HA	2:B:1128:ILE:CG1	0.43	2.43	20	2
2:B:1138:ILE:O	2:B:1142:LEU:HB2	0.43	2.14	16	2
2:B:1164:ALA:O	2:B:1168:ASN:ND2	0.43	2.52	1	3
2:B:1105:ASP:CB	2:B:1106:PRO:CD	0.43	2.96	14	3
2:B:1154:VAL:O	2:B:1158:TRP:CB	0.43	2.67	8	2
2:B:1169:ARG:O	2:B:1175:TYR:CB	0.43	2.67	6	1
2:B:1129:VAL:HG22	2:B:1166:LEU:HG	0.43	1.90	13	1
2:B:1111:PHE:O	2:B:1113:GLN:N	0.43	2.52	15	3
2:B:1107:GLU:O	2:B:1177:PHE:CA	0.43	2.67	17	1
2:B:1161:PHE:CE1	2:B:1181:LEU:C	0.43	2.92	2	1
2:B:1169:ARG:CG	2:B:1169:ARG:O	0.43	2.67	20	1
2:B:1118:GLN:O	2:B:1119:LEU:HB2	0.43	2.14	14	1
2:B:1147:TYR:O	2:B:1149:GLU:N	0.43	2.51	6	1
2:B:1180:LYS:O	2:B:1184:VAL:CG1	0.43	2.66	7	3
2:B:1128:ILE:HG13	2:B:1129:VAL:HG23	0.43	1.91	5	1
2:B:1106:PRO:O	2:B:1110:PRO:CD	0.43	2.67	18	3
2:B:1111:PHE:CD2	2:B:1135:LEU:CD1	0.43	3.01	9	1
2:B:1116:ASP:OD1	2:B:1118:GLN:N	0.43	2.51	18	1
2:B:1177:PHE:CE1	2:B:1178:CYS:SG	0.43	3.09	16	1
2:B:1136:SER:O	2:B:1140:ARG:CG	0.43	2.67	4	1
2:B:1128:ILE:O	2:B:1129:VAL:CG2	0.43	2.67	8	2
2:B:1168:ASN:ND2	2:B:1175:TYR:CD2	0.43	2.87	11	1
2:B:1107:GLU:O	2:B:1181:LEU:CD1	0.43	2.67	2	3
2:B:1156:ASP:O	2:B:1160:MET:CB	0.43	2.67	1	1
2:B:1125:TYR:CE2	2:B:1129:VAL:HG21	0.43	2.48	5	1
2:B:1138:ILE:O	2:B:1142:LEU:N	0.42	2.52	11	1
2:B:1114:PRO:O	2:B:1119:LEU:CD2	0.42	2.67	13	1
1:A:383:LEU:CD2	1:A:383:LEU:O	0.42	2.67	1	2
2:B:1094:LEU:O	2:B:1097:THR:CG2	0.42	2.67	1	1
1:A:384:MET:O	2:B:1173:ARG:NH2	0.42	2.52	20	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1102:TYR:CD1	2:B:1139:LYS:CD	0.42	3.02	20	1
2:B:1172:SER:OG	2:B:1173:ARG:NH2	0.42	2.52	20	1
2:B:1168:ASN:ND2	2:B:1174:VAL:O	0.42	2.52	3	1
2:B:1095:MET:CE	2:B:1095:MET:O	0.42	2.67	16	1
2:B:1190:ASP:CB	2:B:1191:PRO:HD3	0.42	2.44	1	8
2:B:1163:ASN:O	2:B:1164:ALA:CB	0.42	2.67	10	1
2:B:1168:ASN:N	2:B:1168:ASN:ND2	0.42	2.67	14	1
2:B:1192:VAL:HG12	2:B:1192:VAL:O	0.42	2.14	14	1
2:B:1116:ASP:N	2:B:1125:TYR:OH	0.42	2.52	18	1
2:B:1147:TYR:N	2:B:1147:TYR:CD1	0.42	2.87	18	1
2:B:1194:GLN:O	2:B:1194:GLN:NE2	0.42	2.52	15	1
2:B:1112:ARG:O	2:B:1112:ARG:NH1	0.42	2.52	8	1
2:B:1119:LEU:O	2:B:1120:LEU:CB	0.42	2.67	19	2
2:B:1115:VAL:CG2	2:B:1115:VAL:O	0.42	2.67	7	1
2:B:1142:LEU:CD2	2:B:1147:TYR:O	0.42	2.67	9	1
1:A:384:MET:CE	2:B:1167:TYR:O	0.42	2.67	16	1
2:B:1100:ALA:O	2:B:1184:VAL:CG1	0.42	2.67	16	1
2:B:1150:PRO:O	2:B:1154:VAL:HB	0.42	2.14	11	4
2:B:1192:VAL:O	2:B:1193:MET:HB2	0.42	2.13	11	4
2:B:1092:GLN:O	2:B:1096:PRO:HD3	0.42	2.14	6	2
2:B:1131:ASN:O	2:B:1159:LEU:CD1	0.42	2.67	6	2
2:B:1141:LYS:HB3	2:B:1147:TYR:CD2	0.42	2.50	6	2
1:A:384:MET:HA	2:B:1174:VAL:HG21	0.42	1.92	6	1
2:B:1189:ILE:CD1	2:B:1195:SER:OG	0.42	2.66	12	2
2:B:1177:PHE:C	2:B:1177:PHE:CD1	0.42	2.92	12	1
2:B:1154:VAL:CG1	2:B:1194:GLN:NE2	0.42	2.83	11	1
2:B:1122:ILE:HD11	2:B:1125:TYR:CE1	0.42	2.49	7	1
2:B:1102:TYR:CD1	2:B:1139:LYS:HE3	0.42	2.49	7	1
2:B:1081:ARG:NE	2:B:1196:LEU:CB	0.42	2.83	12	1
2:B:1081:ARG:CB	2:B:1196:LEU:HD23	0.42	2.45	18	1
2:B:1095:MET:HE2	2:B:1095:MET:O	0.42	2.15	16	1
2:B:1101:LEU:HD23	2:B:1139:LYS:CE	0.42	2.45	4	1
2:B:1102:TYR:CD2	2:B:1139:LYS:HG2	0.42	2.50	19	1
2:B:1102:TYR:CD2	2:B:1139:LYS:HE2	0.42	2.50	19	1
2:B:1165:TRP:HA	2:B:1168:ASN:OD1	0.42	2.15	19	1
2:B:1156:ASP:O	2:B:1159:LEU:CB	0.42	2.68	8	1
2:B:1170:LYS:O	2:B:1172:SER:N	0.42	2.52	8	2
1:A:384:MET:O	2:B:1173:ARG:NE	0.42	2.53	11	2
2:B:1101:LEU:O	2:B:1181:LEU:CD2	0.42	2.66	2	5
2:B:1196:LEU:HG	2:B:1197:GLY:N	0.42	2.30	6	1
2:B:1154:VAL:CG2	2:B:1194:GLN:OE1	0.42	2.67	5	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1181:LEU:HA	2:B:1184:VAL:CG1	0.42	2.45	1	2
2:B:1102:TYR:CE1	2:B:1139:LYS:HG3	0.42	2.49	20	1
2:B:1118:GLN:O	2:B:1122:ILE:HG23	0.42	2.14	14	1
2:B:1179:SER:O	2:B:1183:GLU:CB	0.42	2.67	9	1
2:B:1129:VAL:CG2	2:B:1163:ASN:HA	0.42	2.44	16	1
2:B:1087:PRO:CB	2:B:1148:GLN:OE1	0.42	2.67	6	1
2:B:1196:LEU:HD23	2:B:1196:LEU:C	0.42	2.34	13	1
2:B:1129:VAL:O	2:B:1131:ASN:N	0.42	2.52	5	2
2:B:1115:VAL:O	2:B:1115:VAL:CG2	0.42	2.66	17	2
2:B:1108:SER:N	2:B:1181:LEU:HD11	0.42	2.29	20	1
2:B:1178:CYS:O	2:B:1182:ALA:CB	0.42	2.67	15	1
2:B:1111:PHE:CD1	2:B:1160:MET:SD	0.42	3.13	4	1
2:B:1184:VAL:O	2:B:1188:GLU:CG	0.42	2.67	11	1
2:B:1094:LEU:O	2:B:1098:LEU:CD1	0.42	2.67	10	1
2:B:1122:ILE:HG12	2:B:1125:TYR:CE1	0.42	2.50	1	2
2:B:1132:PRO:HB3	2:B:1163:ASN:ND2	0.42	2.30	12	2
2:B:1090:LEU:CD1	2:B:1149:GLU:OE2	0.42	2.67	17	1
2:B:1170:LYS:HA	2:B:1175:TYR:CD2	0.42	2.50	2	1
1:A:384:MET:O	1:A:384:MET:CG	0.42	2.67	15	1
2:B:1115:VAL:O	2:B:1116:ASP:O	0.42	2.38	18	2
2:B:1101:LEU:HD11	2:B:1161:PHE:CZ	0.42	2.50	1	1
2:B:1090:LEU:O	2:B:1094:LEU:CG	0.42	2.68	17	2
2:B:1086:LYS:CB	2:B:1089:GLU:OE1	0.42	2.67	2	1
2:B:1189:ILE:HB	2:B:1194:GLN:HG2	0.42	1.91	2	1
2:B:1141:LYS:HB2	2:B:1147:TYR:CE2	0.42	2.50	20	1
2:B:1105:ASP:O	2:B:1109:LEU:CD1	0.42	2.67	9	1
2:B:1095:MET:HB3	2:B:1096:PRO:CD	0.42	2.45	16	1
2:B:1139:LYS:HA	2:B:1142:LEU:HB2	0.42	1.92	15	1
2:B:1098:LEU:HD23	2:B:1139:LYS:NZ	0.42	2.30	10	1
1:A:384:MET:CE	2:B:1169:ARG:CD	0.42	2.97	1	1
2:B:1163:ASN:ND2	2:B:1163:ASN:O	0.42	2.52	5	1
2:B:1111:PHE:CD2	2:B:1181:LEU:HD12	0.42	2.49	16	2
2:B:1139:LYS:HD2	2:B:1139:LYS:N	0.42	2.29	16	1
2:B:1110:PRO:HD2	2:B:1177:PHE:CD2	0.42	2.50	4	1
2:B:1081:ARG:CD	2:B:1081:ARG:C	0.41	2.88	11	1
2:B:1138:ILE:CG2	2:B:1139:LYS:N	0.41	2.83	13	1
2:B:1159:LEU:O	2:B:1163:ASN:N	0.41	2.49	16	2
2:B:1143:ASP:O	2:B:1144:THR:CG2	0.41	2.68	10	1
2:B:1173:ARG:HG3	2:B:1174:VAL:N	0.41	2.30	14	2
2:B:1090:LEU:HB3	2:B:1148:GLN:O	0.41	2.15	9	1
1:A:383:LEU:O	1:A:384:MET:CB	0.41	2.67	18	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1081:ARG:CZ	2:B:1195:SER:O	0.41	2.68	4	1
2:B:1165:TRP:O	2:B:1167:TYR:N	0.41	2.52	4	1
2:B:1128:ILE:O	2:B:1129:VAL:HG23	0.41	2.15	8	1
2:B:1102:TYR:CG	2:B:1139:LYS:HG2	0.41	2.50	6	1
2:B:1116:ASP:HB2	2:B:1117:PRO:CD	0.41	2.44	13	1
2:B:1098:LEU:O	2:B:1102:TYR:CB	0.41	2.67	15	2
2:B:1125:TYR:CD2	2:B:1129:VAL:HB	0.41	2.50	17	1
2:B:1163:ASN:OD1	2:B:1164:ALA:N	0.41	2.53	2	1
2:B:1108:SER:HG	2:B:1135:LEU:HD23	0.41	1.73	20	1
2:B:1161:PHE:O	2:B:1165:TRP:CD1	0.41	2.73	8	1
2:B:1128:ILE:HD13	2:B:1167:TYR:CG	0.41	2.50	6	1
2:B:1110:PRO:O	2:B:1113:GLN:CG	0.41	2.68	13	1
1:A:384:MET:HG3	1:A:384:MET:O	0.41	2.15	3	1
2:B:1091:ARG:NH1	2:B:1143:ASP:O	0.41	2.53	16	1
2:B:1148:GLN:CD	2:B:1149:GLU:N	0.41	2.73	16	1
2:B:1161:PHE:O	2:B:1165:TRP:CB	0.41	2.69	6	1
2:B:1116:ASP:OD1	2:B:1117:PRO:N	0.41	2.53	13	1
2:B:1157:VAL:O	2:B:1160:MET:N	0.41	2.54	5	1
2:B:1168:ASN:O	2:B:1169:ARG:HG2	0.41	2.15	20	1
2:B:1189:ILE:O	2:B:1192:VAL:N	0.41	2.53	20	1
2:B:1173:ARG:CG	2:B:1174:VAL:N	0.41	2.83	14	3
2:B:1116:ASP:OD1	2:B:1118:GLN:CB	0.41	2.69	3	1
2:B:1112:ARG:O	2:B:1113:GLN:CG	0.41	2.69	16	1
2:B:1097:THR:CG2	2:B:1098:LEU:N	0.41	2.83	6	1
2:B:1182:ALA:O	2:B:1185:PHE:CD2	0.41	2.74	10	1
2:B:1156:ASP:O	2:B:1160:MET:N	0.41	2.52	12	1
2:B:1094:LEU:HD23	2:B:1153:TYR:CD2	0.41	2.50	14	1
2:B:1089:GLU:N	2:B:1089:GLU:OE1	0.41	2.53	16	1
2:B:1112:ARG:CZ	2:B:1112:ARG:HB3	0.41	2.45	16	1
2:B:1084:ILE:HG12	2:B:1090:LEU:CD2	0.41	2.46	15	1
2:B:1173:ARG:HG2	2:B:1174:VAL:HG23	0.41	1.92	15	1
2:B:1116:ASP:CB	2:B:1117:PRO:C	0.41	2.89	8	1
2:B:1158:TRP:CD1	2:B:1158:TRP:O	0.41	2.74	6	1
2:B:1161:PHE:CZ	2:B:1182:ALA:HA	0.41	2.51	10	1
2:B:1187:GLN:O	2:B:1191:PRO:CG	0.41	2.68	5	2
2:B:1133:MET:O	2:B:1134:ASP:OD1	0.41	2.39	17	1
2:B:1133:MET:SD	2:B:1156:ASP:CB	0.41	3.09	9	2
2:B:1194:GLN:NE2	2:B:1194:GLN:O	0.41	2.53	9	1
2:B:1111:PHE:CE1	2:B:1160:MET:HG3	0.41	2.51	18	1
2:B:1107:GLU:OE1	2:B:1107:GLU:CA	0.41	2.67	13	1
2:B:1154:VAL:HG21	2:B:1194:GLN:CG	0.41	2.46	13	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1119:LEU:HD12	2:B:1119:LEU:C	0.41	2.35	17	1
2:B:1189:ILE:O	2:B:1192:VAL:HB	0.41	2.15	2	1
2:B:1102:TYR:CG	2:B:1139:LYS:HG3	0.41	2.51	9	1
2:B:1110:PRO:HB2	2:B:1177:PHE:CE1	0.41	2.51	16	1
2:B:1138:ILE:HD11	2:B:1153:TYR:CD1	0.41	2.51	11	1
2:B:1159:LEU:O	2:B:1163:ASN:CB	0.41	2.69	11	1
2:B:1134:ASP:O	2:B:1137:THR:N	0.41	2.53	3	4
2:B:1094:LEU:O	2:B:1098:LEU:CG	0.41	2.69	13	1
2:B:1133:MET:O	2:B:1160:MET:HE3	0.41	2.16	13	1
1:A:383:LEU:HD22	1:A:383:LEU:N	0.41	2.31	16	1
2:B:1095:MET:HB3	2:B:1096:PRO:HD3	0.41	1.92	3	2
2:B:1095:MET:N	2:B:1096:PRO:HD2	0.41	2.31	5	1
2:B:1169:ARG:HB2	2:B:1175:TYR:CD2	0.41	2.50	20	1
2:B:1159:LEU:CD1	2:B:1159:LEU:C	0.41	2.89	14	1
2:B:1098:LEU:HD13	2:B:1102:TYR:HB2	0.41	1.93	3	1
2:B:1147:TYR:CD1	2:B:1152:GLN:CD	0.41	2.94	3	1
1:A:383:LEU:CD2	1:A:383:LEU:N	0.41	2.84	16	1
2:B:1168:ASN:ND2	2:B:1175:TYR:CG	0.41	2.88	4	1
1:A:383:LEU:CD1	1:A:383:LEU:O	0.41	2.65	19	1
2:B:1142:LEU:CD2	2:B:1153:TYR:CE1	0.41	3.04	11	1
2:B:1094:LEU:HD21	2:B:1153:TYR:CD2	0.41	2.50	10	1
2:B:1129:VAL:HG11	2:B:1163:ASN:HD22	0.41	1.75	10	1
2:B:1177:PHE:CD1	2:B:1177:PHE:C	0.41	2.94	17	1
2:B:1141:LYS:CG	2:B:1146:GLN:HB3	0.41	2.45	14	1
2:B:1189:ILE:HD12	2:B:1194:GLN:HE21	0.41	1.76	14	1
2:B:1111:PHE:CD1	2:B:1135:LEU:HG	0.41	2.51	15	1
2:B:1131:ASN:O	2:B:1131:ASN:ND2	0.40	2.53	8	1
2:B:1157:VAL:HG12	2:B:1161:PHE:CD2	0.40	2.51	11	1
2:B:1095:MET:CE	2:B:1099:GLU:OE1	0.40	2.69	3	1
2:B:1161:PHE:CE1	2:B:1182:ALA:HA	0.40	2.52	4	1
2:B:1165:TRP:CD1	2:B:1178:CYS:SG	0.40	3.13	6	2
2:B:1116:ASP:CB	2:B:1117:PRO:HD2	0.40	2.47	13	1
2:B:1081:ARG:HG2	2:B:1196:LEU:CD2	0.40	2.46	1	1
2:B:1157:VAL:O	2:B:1161:PHE:CD2	0.40	2.75	2	1
2:B:1194:GLN:NE2	2:B:1194:GLN:C	0.40	2.75	2	2
2:B:1189:ILE:HB	2:B:1194:GLN:CG	0.40	2.46	2	1
2:B:1149:GLU:CG	2:B:1149:GLU:O	0.40	2.69	14	1
2:B:1115:VAL:CA	2:B:1119:LEU:HD21	0.40	2.47	16	1
2:B:1111:PHE:CD1	2:B:1135:LEU:CG	0.40	3.04	15	1
2:B:1081:ARG:NH2	2:B:1197:GLY:OXT	0.40	2.54	4	1
2:B:1110:PRO:CG	2:B:1177:PHE:CG	0.40	3.04	19	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:1185:PHE:O	2:B:1189:ILE:HG12	0.40	2.16	11	1
2:B:1133:MET:SD	2:B:1133:MET:N	0.40	2.95	6	1
2:B:1124:ASP:OD1	2:B:1124:ASP:N	0.40	2.54	13	1
2:B:1163:ASN:CG	2:B:1163:ASN:O	0.40	2.60	10	1
1:A:384:MET:CE	2:B:1169:ARG:HD2	0.40	2.45	1	1
2:B:1125:TYR:O	2:B:1127:ASP:N	0.40	2.52	2	1
2:B:1141:LYS:HB3	2:B:1147:TYR:CD1	0.40	2.52	19	2
1:A:383:LEU:CD2	2:B:1119:LEU:CD1	0.40	2.99	9	1
2:B:1173:ARG:CG	2:B:1173:ARG:O	0.40	2.67	5	1
2:B:1107:GLU:OE1	2:B:1107:GLU:N	0.40	2.55	17	1
2:B:1168:ASN:N	2:B:1168:ASN:OD1	0.40	2.53	17	1
2:B:1117:PRO:O	2:B:1118:GLN:HB3	0.40	2.16	14	1
2:B:1141:LYS:CD	2:B:1147:TYR:CZ	0.40	3.04	14	1
2:B:1102:TYR:HD2	2:B:1135:LEU:HD22	0.40	1.76	18	1
2:B:1101:LEU:HD12	2:B:1101:LEU:O	0.40	2.17	13	1
2:B:1185:PHE:O	2:B:1188:GLU:N	0.40	2.55	5	1
2:B:1101:LEU:HD23	2:B:1102:TYR:N	0.40	2.32	2	1
2:B:1151:TRP:CD1	2:B:1151:TRP:C	0.40	2.94	16	1
2:B:1102:TYR:CB	2:B:1139:LYS:HE2	0.40	2.46	19	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	2/20 (10%)	1±1 (50±27%)	1±1 (45±31%)	0±0 (5±15%)	4	26
2	B	116/121 (96%)	87±4 (75±3%)	21±3 (18±3%)	8±2 (7±2%)	2	16
All	All	2360/2820 (84%)	1750 (74%)	439 (19%)	171 (7%)	2	16

All 42 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
2	B	1085	PHE	20
2	B	1104	GLN	12

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Mol	Chain	Res	Type	Models (Total)
2	B	1196	LEU	10
2	B	1114	PRO	9
2	B	1176	LYS	8
2	B	1171	THR	7
2	B	1115	VAL	7
2	B	1123	PRO	7
2	B	1129	VAL	6
2	B	1120	LEU	6
2	B	1084	ILE	6
2	B	1116	ASP	6
2	B	1170	LYS	6
2	B	1121	GLY	6
2	B	1124	ASP	5
2	B	1169	ARG	4
2	B	1111	PHE	4
2	B	1168	ASN	4
2	B	1118	GLN	3
2	B	1081	ARG	3
2	B	1122	ILE	3
2	B	1157	VAL	2
2	B	1126	PHE	2
2	B	1119	LEU	2
2	B	1173	ARG	2
2	B	1125	TYR	2
2	B	1117	PRO	2
1	A	383	LEU	2
2	B	1105	ASP	2
2	B	1175	TYR	1
2	B	1150	PRO	1
2	B	1112	ARG	1
2	B	1130	LYS	1
2	B	1164	ALA	1
2	B	1134	ASP	1
2	B	1127	ASP	1
2	B	1165	TRP	1
2	B	1174	VAL	1
2	B	1132	PRO	1
2	B	1133	MET	1
2	B	1148	GLN	1
2	B	1167	TYR	1

6.3.2 Protein sidechains ⓘ

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	2/18 (11%)	1±1 (52±33%)	1±1 (48±33%)	0	1
2	B	110/113 (97%)	79±4 (72±3%)	31±4 (28±3%)	2	20
All	All	2240/2620 (85%)	1604 (72%)	636 (28%)	2	19

All 92 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
2	B	1185	PHE	19
2	B	1085	PHE	18
2	B	1095	MET	17
2	B	1190	ASP	16
2	B	1111	PHE	15
2	B	1105	ASP	15
2	B	1194	GLN	14
2	B	1125	TYR	14
1	A	383	LEU	13
2	B	1167	TYR	13
2	B	1098	LEU	13
2	B	1118	GLN	13
2	B	1120	LEU	12
2	B	1081	ARG	11
2	B	1193	MET	11
2	B	1195	SER	11
2	B	1162	ASN	11
2	B	1184	VAL	10
2	B	1160	MET	10
2	B	1112	ARG	10
2	B	1165	TRP	10
2	B	1083	LYS	10
2	B	1173	ARG	10
2	B	1166	LEU	10
2	B	1133	MET	10
2	B	1130	LYS	9
2	B	1104	GLN	9
2	B	1143	ASP	9

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Mol	Chain	Res	Type	Models (Total)
2	B	1196	LEU	9
2	B	1169	ARG	9
2	B	1180	LYS	9
2	B	1176	LYS	9
2	B	1168	ASN	8
2	B	1144	THR	8
2	B	1091	ARG	8
2	B	1170	LYS	7
2	B	1086	LYS	7
2	B	1101	LEU	7
2	B	1103	ARG	7
2	B	1140	ARG	7
2	B	1139	LYS	7
2	B	1082	LYS	6
1	A	384	MET	6
2	B	1149	GLU	6
2	B	1142	LEU	6
2	B	1155	ASP	6
2	B	1148	GLN	6
2	B	1119	LEU	6
2	B	1131	ASN	6
2	B	1159	LEU	6
2	B	1135	LEU	6
2	B	1172	SER	6
2	B	1088	GLU	6
2	B	1136	SER	6
2	B	1102	TYR	6
2	B	1128	ILE	5
2	B	1137	THR	5
2	B	1163	ASN	5
2	B	1113	GLN	5
2	B	1161	PHE	5
2	B	1156	ASP	5
2	B	1124	ASP	5
2	B	1116	ASP	5
2	B	1092	GLN	5
2	B	1084	ILE	4
2	B	1122	ILE	4
2	B	1152	GLN	4
2	B	1126	PHE	4
2	B	1127	ASP	4
2	B	1186	GLU	3

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Mol	Chain	Res	Type	Models (Total)
2	B	1178	CYS	3
2	B	1146	GLN	3
2	B	1181	LEU	3
2	B	1141	LYS	3
2	B	1099	GLU	3
2	B	1189	ILE	3
2	B	1107	GLU	2
2	B	1108	SER	2
2	B	1183	GLU	2
2	B	1187	GLN	2
2	B	1134	ASP	2
2	B	1138	ILE	1
2	B	1171	THR	1
2	B	1188	GLU	1
2	B	1117	PRO	1
2	B	1110	PRO	1
2	B	1089	GLU	1
2	B	1109	LEU	1
2	B	1097	THR	1
2	B	1179	SER	1
2	B	1115	VAL	1
2	B	1175	TYR	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length is the number of standard deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
1	ALY	A	382	1	11,11,12	1.20±0.08	0±0 (0±0%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
1	ALY	A	382	1	9,12,14	1.19±0.09	0±0 (0±0%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	ALY	A	382	1	-	0±0,8,10,12	0±0,0,0,0

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided