



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 04:25 AM UTC

PDB ID : 1Q5L / pdb_00001q5l
Title : NMR structure of the substrate binding domain of DnaK bound to the peptide
NRLLLTG
Authors : Stevens, S.Y.; Cai, S.; Pellecchia, M.; Zuiderweg, E.R.
Deposited on : 2003-08-08

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

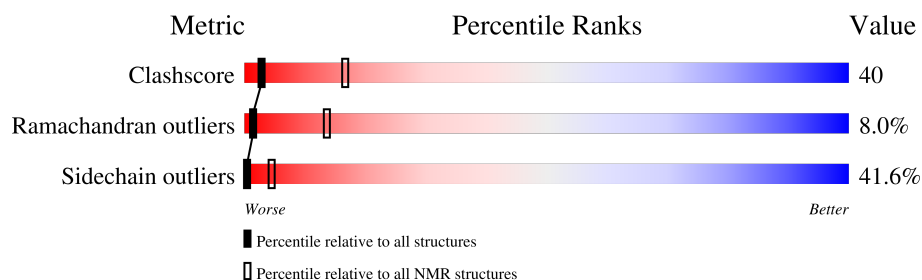
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

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

Mol	Chain	Length	Quality of chain
1	A	135	
2	B	7	

2 Ensemble composition and analysis

This entry contains 15 models. Model 2 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:394-A:503, B:902-B:907 (116)	1.25	2

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 3 clusters and 1 single-model cluster was found.

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

3 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1707 atoms, of which 795 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Chaperone protein dnaK.

Mol	Chain	Residues	Atoms						Trace
1	A	115	Total	C	H	N	O	S	0
			1595	529	738	151	174	3	

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	373	MET	-	cloning artifact	UNP P0A6Y8
A	374	GLY	-	cloning artifact	UNP P0A6Y8
A	375	SER	-	cloning artifact	UNP P0A6Y8
A	376	SER	-	cloning artifact	UNP P0A6Y8
A	377	HIS	-	cloning artifact	UNP P0A6Y8
A	378	HIS	-	cloning artifact	UNP P0A6Y8
A	379	HIS	-	cloning artifact	UNP P0A6Y8
A	380	HIS	-	cloning artifact	UNP P0A6Y8
A	381	HIS	-	cloning artifact	UNP P0A6Y8
A	382	HIS	-	cloning artifact	UNP P0A6Y8
A	383	GLY	-	cloning artifact	UNP P0A6Y8
A	384	LEU	-	cloning artifact	UNP P0A6Y8
A	385	VAL	-	cloning artifact	UNP P0A6Y8
A	386	PRO	-	cloning artifact	UNP P0A6Y8
A	387	ARG	-	cloning artifact	UNP P0A6Y8
A	388	GLY	-	cloning artifact	UNP P0A6Y8
A	389	SER	-	cloning artifact	UNP P0A6Y8
A	390	HIS	-	cloning artifact	UNP P0A6Y8
A	391	MET	-	cloning artifact	UNP P0A6Y8
A	392	VAL	-	cloning artifact	UNP P0A6Y8

- Molecule 2 is a protein called peptide NRLLLTG.

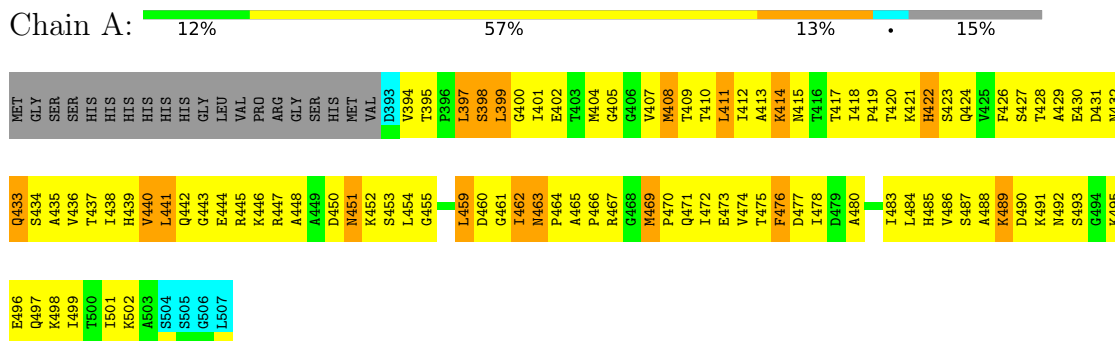
Mol	Chain	Residues	Atoms						Trace
2	B	7	Total	C	H	N	O		0
			112	34	57	11	10		

4 Residue-property plots

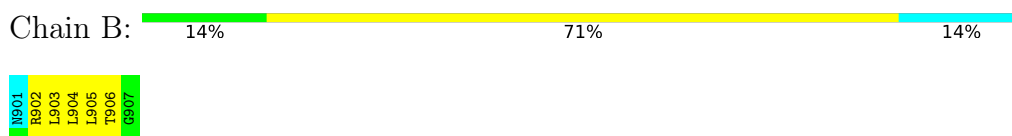
4.1 Average score per residue in the NMR ensemble

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: Chaperone protein dnaK



- Molecule 2: peptide NRLLLTG

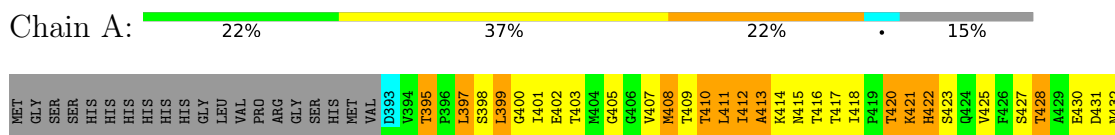


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: Chaperone protein dnaK

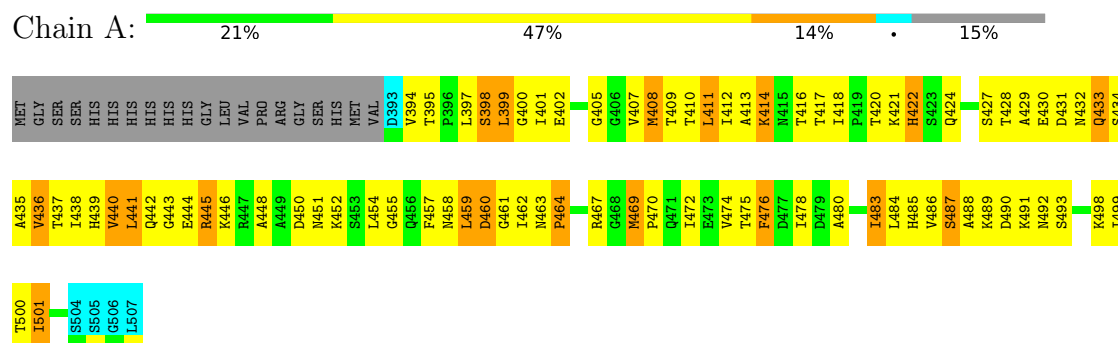


- Molecule 2: peptide NRLLTG



4.2.2 Score per residue for model 2 (medoid)

- Molecule 1: Chaperone protein dnaK

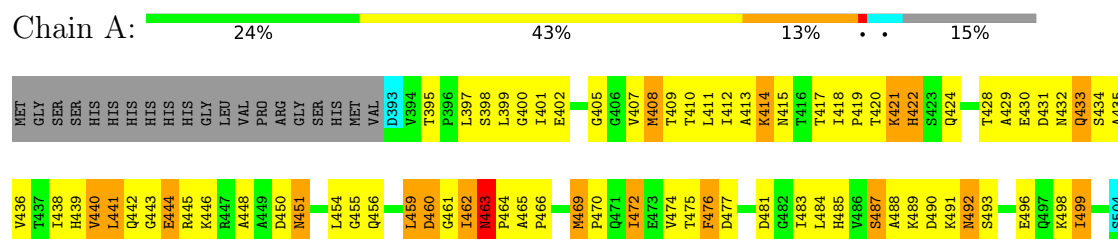


- Molecule 2: peptide NRLLTG



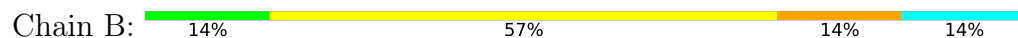
4.2.3 Score per residue for model 3

- Molecule 1: Chaperone protein dnaK



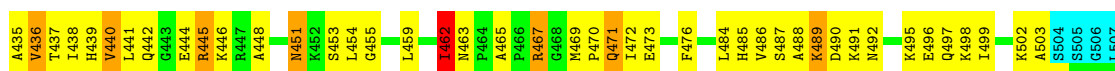


- Molecule 2: peptide NRLLLTG



4.2.4 Score per residue for model 4

- Molecule 1: Chaperone protein dnaK



- Molecule 2: peptide NRLLLTG



4.2.5 Score per residue for model 5

- Molecule 1: Chaperone protein dnaK



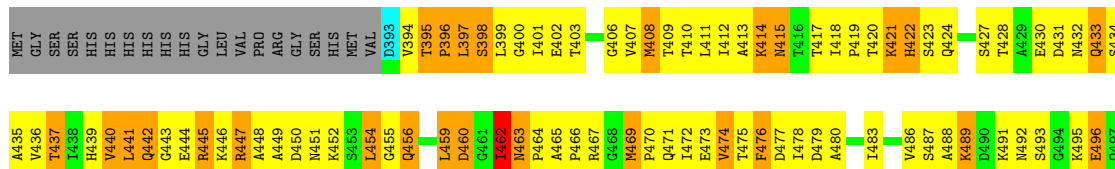
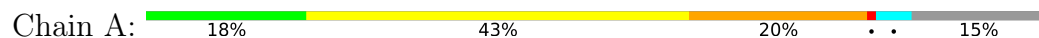
- Molecule 2: peptide NRLLLTG





4.2.6 Score per residue for model 6

- Molecule 1: Chaperone protein dnaK

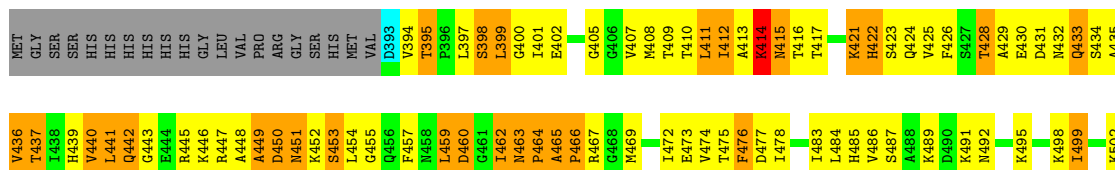
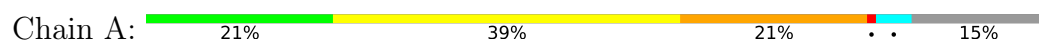


- Molecule 2: peptide NRLLLTG

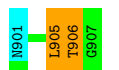


4.2.7 Score per residue for model 7

- Molecule 1: Chaperone protein dnaK

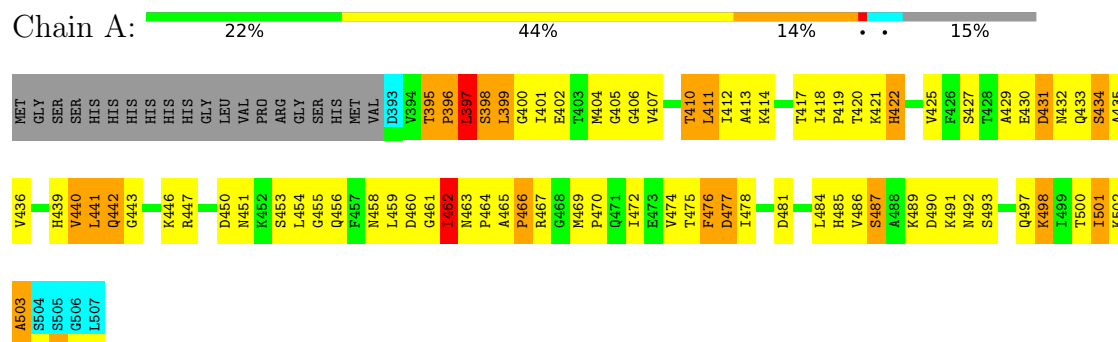


- Molecule 2: peptide NRLLLTG

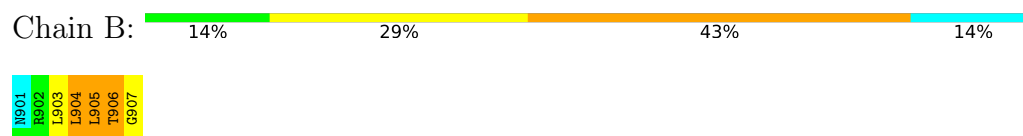


4.2.8 Score per residue for model 8

- Molecule 1: Chaperone protein dnaK

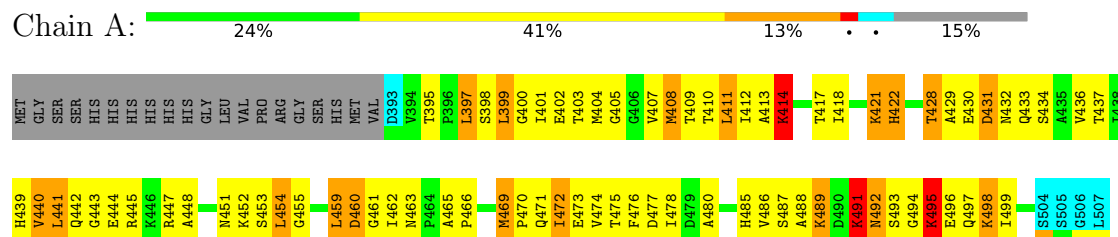


- Molecule 2: peptide NRLLLTG



4.2.9 Score per residue for model 9

- Molecule 1: Chaperone protein dnaK

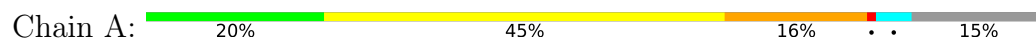


- Molecule 2: peptide NRLLLTG



4.2.10 Score per residue for model 10

- Molecule 1: Chaperone protein dnaK





- Molecule 2: peptide NRLLLTG

Chain B: 14% 57% 14% 14%



4.2.11 Score per residue for model 11

- Molecule 1: Chaperone protein dnaK

Chain A: 18% 45% 17% 15%



- Molecule 2: peptide NRLLLTG

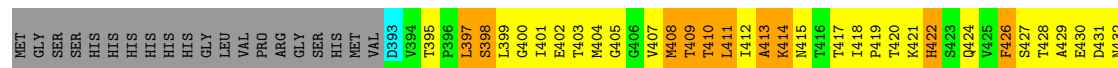
Chain B: 43% 14% 29% 14%

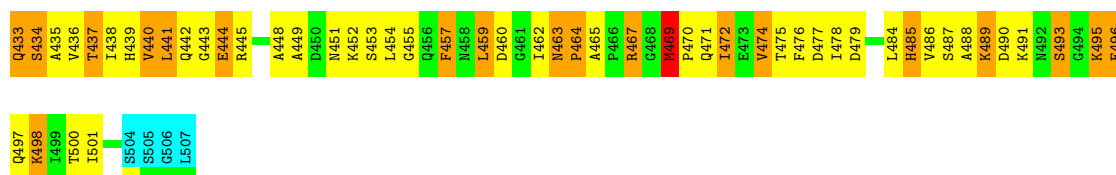


4.2.12 Score per residue for model 12

- Molecule 1: Chaperone protein dnaK

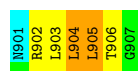
Chain A: 18% 41% 21% 15%





- Molecule 2: peptide NRLLLTG

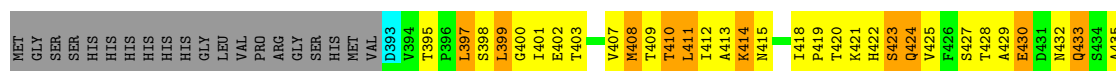
Chain B: 14% 43% 29% 14%



4.2.13 Score per residue for model 13

- Molecule 1: Chaperone protein dnaK

Chain A: 19% 41% 19% . . 15%



- Molecule 2: peptide NRLLLTG

Chain B: 14% 29% 43% 14%



4.2.14 Score per residue for model 14

- Molecule 1: Chaperone protein dnaK

Chain A: 18% 42% 21% . . 15%



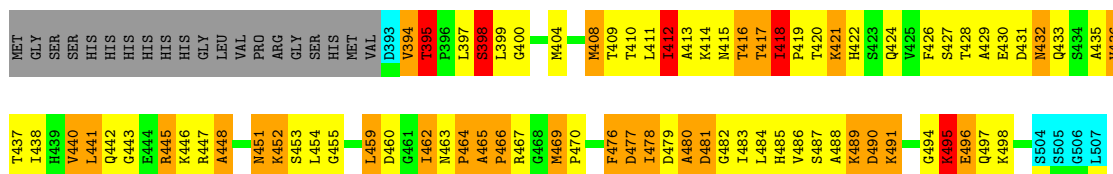
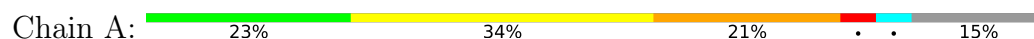


- Molecule 2: peptide NRLLTG

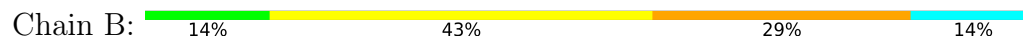


4.2.15 Score per residue for model 15

- Molecule 1: Chaperone protein dnaK



- Molecule 2: peptide NRLLTG



5 Refinement protocol and experimental data overview

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

Of the 60 calculated structures, 15 were deposited, based on the following criterion: *structures with acceptable covalent geometry, 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
ARIA/(CNS)	structure solution	1.0
ARIA/(CNS)	refinement	1.0

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	0.36±0.04	0±0/835 (0.0± 0.0%)	0.71±0.02	0±0/1130 (0.0± 0.0%)
2	B	0.31±0.03	0±0/46 (0.0± 0.0%)	0.58±0.11	0±0/59 (0.0± 0.0%)
All	All	0.36	0/13215 (0.0%)	0.70	2/17835 (0.0%)

There are no bond-length outliers.

All unique angle outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	493	SER	N-CA-C	-5.70	106.37	113.55	12	1
1	A	418	ILE	N-CA-CB	-5.22	106.13	111.87	15	1

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	824	714	842	71±17
2	B	47	50	56	6±3
All	All	13065	11460	13470	1069

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 40.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:489:LYS:HB3	1:A:496:GLU:HA	0.93	1.40	12	1
1:A:416:THR:HG21	1:A:478:ILE:HD12	0.83	1.50	15	1
1:A:440:VAL:HB	1:A:454:LEU:HB2	0.82	1.48	10	4
1:A:440:VAL:HG11	1:A:486:VAL:HG21	0.81	1.52	15	7
1:A:463:ASN:HB3	1:A:464:PRO:HD3	0.81	1.51	14	2
1:A:427:SER:HB2	2:B:904:LEU:HD12	0.80	1.52	14	1
1:A:400:GLY:HA3	1:A:410:THR:HA	0.80	1.50	15	3
1:A:489:LYS:HG2	1:A:496:GLU:HB3	0.80	1.54	15	1
1:A:402:GLU:HB3	1:A:439:HIS:HB3	0.80	1.51	4	12
1:A:398:SER:N	1:A:414:LYS:HG3	0.79	1.93	6	1
1:A:408:MET:HB3	1:A:448:ALA:HB2	0.78	1.55	4	7
1:A:397:LEU:CB	1:A:414:LYS:HB3	0.77	2.09	15	1
1:A:412:ILE:HD11	1:A:476:PHE:HB2	0.77	1.55	6	3
1:A:429:ALA:HB3	2:B:905:LEU:HA	0.77	1.56	8	2
1:A:397:LEU:HB2	1:A:414:LYS:HD3	0.77	1.56	15	1
1:A:440:VAL:HG23	1:A:455:GLY:H	0.77	1.39	11	13
1:A:401:ILE:HG22	1:A:440:VAL:HG13	0.77	1.57	10	1
1:A:417:THR:HA	1:A:482:GLY:H	0.76	1.38	15	1
1:A:435:ALA:O	2:B:906:THR:HG22	0.75	1.81	10	2
1:A:402:GLU:HB2	1:A:439:HIS:HB3	0.75	1.57	2	1
1:A:436:VAL:HG21	1:A:472:ILE:HG12	0.74	1.56	14	3
1:A:431:ASP:N	1:A:467:ARG:HB2	0.73	1.98	14	1
1:A:395:THR:HG21	1:A:412:ILE:HG22	0.73	1.61	15	1
1:A:428:THR:HB	1:A:467:ARG:C	0.73	2.08	14	1
1:A:420:THR:HG22	1:A:478:ILE:HB	0.73	1.61	15	5
1:A:399:LEU:HD13	1:A:413:ALA:C	0.72	2.08	15	1
1:A:417:THR:HA	1:A:482:GLY:N	0.72	2.00	15	1
1:A:397:LEU:HB3	1:A:414:LYS:HB3	0.71	1.60	15	1
1:A:395:THR:HB	1:A:414:LYS:HB3	0.71	1.59	6	1
1:A:442:GLN:HB2	1:A:454:LEU:HD22	0.71	1.61	13	1
1:A:429:ALA:HB2	2:B:903:LEU:HD23	0.71	1.61	15	3
1:A:427:SER:CB	2:B:904:LEU:HD12	0.70	2.15	14	1
1:A:395:THR:HA	1:A:414:LYS:O	0.70	1.85	15	1
1:A:440:VAL:HG23	1:A:455:GLY:N	0.70	2.02	8	12
1:A:399:LEU:HD12	1:A:442:GLN:HG3	0.70	1.64	15	3
1:A:436:VAL:HG12	1:A:459:LEU:HB3	0.70	1.62	13	12
1:A:498:LYS:O	1:A:499:ILE:HG23	0.70	1.87	13	1
1:A:397:LEU:HG	1:A:398:SER:H	0.70	1.45	6	1
1:A:394:VAL:O	1:A:395:THR:HG23	0.69	1.86	15	1
1:A:427:SER:HA	2:B:904:LEU:HD23	0.69	1.62	13	1
1:A:397:LEU:N	1:A:414:LYS:HB2	0.69	2.02	6	1
1:A:427:SER:HB2	1:A:472:ILE:HB	0.69	1.64	4	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:430:GLU:C	1:A:466:PRO:O	0.69	2.36	14	1
1:A:431:ASP:HA	1:A:466:PRO:HD2	0.68	1.64	9	2
1:A:404:MET:H	2:B:904:LEU:HA	0.68	1.46	14	5
1:A:489:LYS:HA	1:A:496:GLU:H	0.68	1.48	9	1
1:A:399:LEU:HA	1:A:442:GLN:HB2	0.68	1.64	11	1
1:A:397:LEU:HB3	1:A:414:LYS:H	0.68	1.49	15	1
1:A:491:LYS:HB3	1:A:494:GLY:H	0.68	1.49	5	2
1:A:408:MET:HG2	1:A:450:ASP:HB3	0.67	1.65	11	1
1:A:472:ILE:HA	1:A:489:LYS:O	0.67	1.90	5	14
1:A:412:ILE:HG21	1:A:420:THR:HG21	0.66	1.66	14	5
1:A:442:GLN:CB	1:A:454:LEU:HD22	0.66	2.19	13	1
1:A:442:GLN:HG2	1:A:454:LEU:HG	0.66	1.68	3	3
1:A:402:GLU:HB3	1:A:439:HIS:HB2	0.66	1.67	10	1
1:A:435:ALA:HA	1:A:459:LEU:O	0.66	1.91	8	4
1:A:436:VAL:HA	2:B:907:GLY:HA2	0.66	1.67	10	1
1:A:408:MET:HG3	1:A:449:ALA:HB3	0.65	1.69	11	2
1:A:404:MET:HG3	2:B:905:LEU:HD23	0.65	1.67	8	1
1:A:431:ASP:HB3	1:A:466:PRO:HA	0.65	1.69	5	1
1:A:397:LEU:HB3	1:A:414:LYS:HE2	0.65	1.67	6	1
1:A:433:GLN:HG3	2:B:906:THR:HA	0.65	1.66	8	1
1:A:473:GLU:O	1:A:488:ALA:HA	0.65	1.91	14	5
1:A:441:LEU:HB3	1:A:450:ASP:HB3	0.64	1.67	7	1
1:A:395:THR:HG22	1:A:413:ALA:O	0.64	1.93	15	1
1:A:433:GLN:O	1:A:462:ILE:HB	0.64	1.93	14	1
1:A:440:VAL:C	1:A:454:LEU:HD12	0.63	2.18	6	11
1:A:442:GLN:HB3	1:A:454:LEU:HD11	0.63	1.69	3	4
1:A:473:GLU:H	1:A:489:LYS:HB3	0.63	1.51	9	1
1:A:442:GLN:HB2	1:A:454:LEU:HD11	0.63	1.67	5	7
1:A:474:VAL:HA	1:A:487:SER:O	0.63	1.94	6	10
1:A:432:ASN:O	1:A:433:GLN:HB2	0.63	1.93	14	1
1:A:398:SER:O	1:A:442:GLN:HB2	0.63	1.93	4	2
1:A:432:ASN:H	1:A:467:ARG:HB3	0.63	1.53	14	1
1:A:445:ARG:HD2	1:A:451:ASN:HB2	0.63	1.71	2	1
1:A:427:SER:CB	2:B:904:LEU:HD11	0.63	2.23	4	1
1:A:397:LEU:HG	1:A:398:SER:N	0.63	2.09	6	2
1:A:398:SER:HA	1:A:414:LYS:N	0.62	2.08	11	9
1:A:394:VAL:HG23	1:A:416:THR:O	0.62	1.93	7	1
1:A:400:GLY:HA3	1:A:409:THR:O	0.62	1.93	12	13
1:A:429:ALA:N	1:A:467:ARG:N	0.62	2.47	14	1
1:A:397:LEU:H	1:A:414:LYS:HB3	0.62	1.54	8	1
1:A:430:GLU:N	1:A:466:PRO:O	0.62	2.32	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:404:MET:HG3	2:B:905:LEU:HD21	0.62	1.69	15	1
1:A:436:VAL:HB	1:A:462:ILE:HD11	0.62	1.70	13	3
1:A:399:LEU:HD11	1:A:484:LEU:HD22	0.62	1.71	5	2
1:A:469:MET:N	1:A:470:PRO:HD2	0.62	2.09	9	3
1:A:466:PRO:O	1:A:467:ARG:HB2	0.62	1.93	14	1
1:A:408:MET:HE1	1:A:445:ARG:HG3	0.61	1.72	4	1
1:A:445:ARG:CZ	1:A:451:ASN:HB2	0.61	2.25	15	1
1:A:476:PHE:HB3	1:A:484:LEU:HD11	0.61	1.71	8	8
1:A:432:ASN:HA	1:A:464:PRO:HA	0.61	1.71	13	3
1:A:471:GLN:O	1:A:490:ASP:HA	0.61	1.96	12	1
1:A:432:ASN:C	1:A:465:ALA:O	0.61	2.43	14	1
1:A:401:ILE:HG12	1:A:411:LEU:HD23	0.61	1.72	8	7
1:A:435:ALA:HB1	1:A:459:LEU:O	0.61	1.96	14	5
1:A:429:ALA:HB3	2:B:905:LEU:HB2	0.61	1.71	2	1
1:A:464:PRO:HD2	1:A:466:PRO:HD3	0.61	1.72	10	1
1:A:429:ALA:O	1:A:430:GLU:HB2	0.61	1.96	14	3
1:A:435:ALA:HA	1:A:460:ASP:HA	0.61	1.72	13	7
1:A:448:ALA:HA	1:A:451:ASN:ND2	0.61	2.10	3	1
1:A:418:ILE:HB	1:A:419:PRO:HD3	0.61	1.71	11	8
1:A:479:ASP:HB2	1:A:483:ILE:HD12	0.60	1.72	11	2
1:A:412:ILE:HG12	1:A:420:THR:HG21	0.60	1.72	1	2
1:A:429:ALA:HB3	2:B:905:LEU:HB3	0.60	1.71	13	1
1:A:398:SER:HA	1:A:413:ALA:HA	0.60	1.72	1	12
1:A:436:VAL:HG12	1:A:459:LEU:CB	0.60	2.27	15	4
1:A:491:LYS:HD3	1:A:493:SER:H	0.59	1.57	5	1
1:A:448:ALA:O	1:A:450:ASP:N	0.59	2.34	7	1
1:A:488:ALA:O	1:A:496:GLU:HB3	0.59	1.98	9	1
1:A:421:LYS:HA	1:A:476:PHE:O	0.59	1.97	15	12
1:A:404:MET:N	2:B:904:LEU:HA	0.59	2.12	4	2
1:A:404:MET:C	2:B:905:LEU:HB2	0.59	2.22	10	2
1:A:452:LYS:HG3	1:A:454:LEU:HB3	0.59	1.74	13	1
1:A:398:SER:HA	1:A:414:LYS:H	0.59	1.58	1	7
1:A:459:LEU:HD13	1:A:488:ALA:O	0.59	1.98	12	1
1:A:465:ALA:H	1:A:466:PRO:HD2	0.59	1.56	7	1
1:A:431:ASP:O	1:A:464:PRO:HB2	0.59	1.98	14	1
1:A:432:ASN:HB3	1:A:463:ASN:O	0.59	1.98	15	1
1:A:476:PHE:HA	1:A:485:HIS:O	0.59	1.97	10	11
1:A:475:THR:HB	1:A:487:SER:OG	0.58	1.98	14	3
1:A:463:ASN:HB2	1:A:464:PRO:HD3	0.58	1.75	5	2
1:A:489:LYS:HD3	1:A:495:LYS:HA	0.58	1.75	9	1
1:A:463:ASN:N	1:A:464:PRO:HD3	0.58	2.13	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:432:ASN:ND2	1:A:464:PRO:HD3	0.58	2.13	7	2
1:A:430:GLU:HG3	2:B:905:LEU:HG	0.58	1.75	10	1
1:A:428:THR:HG21	1:A:470:PRO:HG2	0.58	1.76	13	2
1:A:397:LEU:HA	1:A:414:LYS:HD2	0.58	1.74	11	3
1:A:395:THR:CG2	1:A:412:ILE:HG22	0.58	2.29	15	1
1:A:412:ILE:HD13	1:A:478:ILE:CG1	0.57	2.29	15	1
1:A:438:ILE:HG13	1:A:488:ALA:HB2	0.57	1.76	15	6
1:A:395:THR:HB	1:A:413:ALA:HB3	0.57	1.75	15	1
1:A:435:ALA:O	2:B:906:THR:HG21	0.57	1.99	14	1
1:A:399:LEU:HD11	1:A:503:ALA:HB1	0.57	1.76	8	1
1:A:416:THR:HG21	1:A:478:ILE:CD1	0.57	2.27	15	1
1:A:466:PRO:HA	1:A:470:PRO:HG2	0.57	1.74	3	1
1:A:397:LEU:H	1:A:414:LYS:HB2	0.57	1.59	6	1
1:A:442:GLN:HG3	1:A:443:GLY:N	0.57	2.14	6	1
1:A:402:GLU:HA	1:A:407:VAL:O	0.57	1.99	8	7
1:A:436:VAL:HA	2:B:906:THR:O	0.57	1.99	10	1
1:A:463:ASN:HB3	1:A:464:PRO:CD	0.57	2.26	14	1
1:A:431:ASP:C	1:A:465:ALA:O	0.57	2.47	14	1
1:A:394:VAL:HG21	1:A:417:THR:HG22	0.57	1.75	15	1
1:A:419:PRO:HA	1:A:478:ILE:O	0.57	2.00	15	1
1:A:399:LEU:HD13	1:A:478:ILE:HD11	0.57	1.77	5	5
1:A:440:VAL:O	1:A:454:LEU:HB2	0.57	1.99	12	10
1:A:476:PHE:N	1:A:476:PHE:CD1	0.57	2.73	15	6
1:A:472:ILE:HD13	1:A:489:LYS:O	0.57	2.00	13	1
1:A:411:LEU:HG	1:A:476:PHE:CZ	0.56	2.35	5	7
1:A:491:LYS:O	1:A:491:LYS:HG2	0.56	1.99	15	2
1:A:427:SER:O	1:A:472:ILE:HD13	0.56	2.00	14	1
1:A:412:ILE:HG23	1:A:416:THR:HG21	0.56	1.76	1	1
1:A:469:MET:HB3	1:A:470:PRO:HD3	0.56	1.77	3	7
1:A:397:LEU:CG	1:A:398:SER:N	0.56	2.66	8	1
1:A:466:PRO:HD2	1:A:470:PRO:HD3	0.56	1.77	8	1
1:A:465:ALA:O	1:A:466:PRO:O	0.56	2.22	14	1
1:A:418:ILE:HG21	1:A:480:ALA:HA	0.56	1.76	6	7
1:A:422:HIS:O	1:A:475:THR:HA	0.56	2.01	12	12
1:A:438:ILE:O	1:A:457:PHE:HB2	0.56	2.01	2	1
1:A:412:ILE:HD13	1:A:478:ILE:HG13	0.56	1.77	15	3
1:A:445:ARG:NH1	1:A:448:ALA:HA	0.56	2.14	15	1
1:A:444:GLU:C	1:A:445:ARG:HG2	0.56	2.25	10	4
1:A:416:THR:HG23	1:A:482:GLY:HA2	0.56	1.76	15	1
1:A:440:VAL:HB	1:A:454:LEU:CD1	0.56	2.30	11	7
1:A:451:ASN:CG	1:A:452:LYS:H	0.56	2.09	11	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:498:LYS:C	1:A:499:ILE:HD13	0.56	2.25	13	1
1:A:459:LEU:HD13	1:A:488:ALA:HB3	0.56	1.77	3	3
1:A:442:GLN:CB	1:A:454:LEU:HD11	0.56	2.30	11	1
1:A:478:ILE:HA	1:A:483:ILE:O	0.56	2.00	15	2
1:A:441:LEU:HD22	1:A:448:ALA:HB1	0.56	1.77	10	1
1:A:399:LEU:HA	1:A:442:GLN:CB	0.56	2.31	11	1
1:A:411:LEU:HG	1:A:476:PHE:CE1	0.55	2.37	10	5
1:A:491:LYS:HB3	1:A:494:GLY:N	0.55	2.15	5	1
1:A:487:SER:HB2	1:A:498:LYS:HB3	0.55	1.77	12	1
1:A:489:LYS:CB	1:A:496:GLU:HA	0.55	2.26	12	1
1:A:486:VAL:N	1:A:499:ILE:HG22	0.55	2.16	13	1
1:A:465:ALA:HB1	1:A:466:PRO:HD2	0.55	1.78	15	1
1:A:491:LYS:HE3	1:A:491:LYS:C	0.55	2.27	10	3
1:A:410:THR:C	1:A:411:LEU:HD22	0.55	2.27	15	1
1:A:479:ASP:C	1:A:481:ASP:H	0.55	2.10	15	1
1:A:394:VAL:HG22	1:A:395:THR:H	0.55	1.61	7	1
1:A:443:GLY:O	1:A:450:ASP:HB2	0.55	2.01	11	1
1:A:476:PHE:CE2	1:A:486:VAL:HG13	0.55	2.37	12	1
1:A:477:ASP:HB3	1:A:485:HIS:NE2	0.55	2.15	5	1
1:A:472:ILE:HD12	1:A:488:ALA:HB1	0.55	1.79	11	1
1:A:428:THR:HG22	1:A:467:ARG:H	0.55	1.61	14	1
1:A:442:GLN:O	1:A:442:GLN:HG2	0.54	2.02	11	3
1:A:433:GLN:HA	1:A:465:ALA:C	0.54	2.27	14	1
1:A:432:ASN:O	1:A:433:GLN:C	0.54	2.50	8	3
1:A:440:VAL:CG1	1:A:486:VAL:HG21	0.54	2.28	15	2
1:A:459:LEU:HD13	1:A:460:ASP:H	0.54	1.62	1	3
1:A:465:ALA:N	1:A:466:PRO:HD3	0.54	2.17	9	3
1:A:440:VAL:HG23	1:A:455:GLY:O	0.54	2.03	9	1
1:A:397:LEU:HD23	1:A:442:GLN:CD	0.54	2.28	11	1
1:A:491:LYS:HB2	1:A:495:LYS:H	0.54	1.63	5	1
1:A:426:PHE:CD1	2:B:904:LEU:HD11	0.54	2.38	12	2
1:A:442:GLN:HG2	1:A:452:LYS:N	0.54	2.17	6	1
1:A:433:GLN:HA	1:A:466:PRO:N	0.54	2.17	14	1
1:A:463:ASN:HB3	1:A:464:PRO:HD2	0.54	1.78	1	2
1:A:476:PHE:CD2	1:A:484:LEU:HD11	0.54	2.37	4	1
1:A:440:VAL:HG21	1:A:486:VAL:HG21	0.54	1.80	9	2
1:A:425:VAL:HG23	1:A:472:ILE:O	0.54	2.03	14	1
1:A:461:GLY:C	1:A:462:ILE:HG13	0.54	2.28	14	2
1:A:448:ALA:O	1:A:449:ALA:C	0.53	2.51	7	1
1:A:433:GLN:OE1	1:A:466:PRO:HD2	0.53	2.02	10	1
1:A:436:VAL:HG21	1:A:472:ILE:HG13	0.53	1.78	4	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:423:SER:HA	1:A:474:VAL:O	0.53	2.03	6	4
1:A:397:LEU:HD23	1:A:442:GLN:HG2	0.53	1.80	10	1
1:A:447:ARG:O	1:A:449:ALA:N	0.53	2.41	6	1
1:A:469:MET:H	1:A:470:PRO:HD2	0.53	1.62	14	2
1:A:443:GLY:HA3	1:A:451:ASN:CG	0.53	2.29	2	2
1:A:432:ASN:HA	1:A:463:ASN:O	0.53	2.04	4	3
1:A:445:ARG:O	1:A:449:ALA:HA	0.53	2.04	7	1
1:A:408:MET:HG2	1:A:451:ASN:HD21	0.53	1.64	5	1
1:A:441:LEU:CB	1:A:450:ASP:HB3	0.53	2.34	7	1
1:A:434:SER:HB2	1:A:463:ASN:O	0.53	2.04	14	1
1:A:463:ASN:CB	1:A:464:PRO:HD3	0.53	2.31	14	1
1:A:466:PRO:HG2	2:B:906:THR:H	0.53	1.64	14	1
1:A:398:SER:O	1:A:442:GLN:HG3	0.53	2.04	1	2
1:A:441:LEU:HB3	1:A:450:ASP:CB	0.53	2.33	7	1
1:A:448:ALA:HA	1:A:451:ASN:HD21	0.53	1.64	5	2
1:A:471:GLN:C	1:A:472:ILE:HD13	0.53	2.29	5	1
1:A:445:ARG:HG2	1:A:451:ASN:HD22	0.53	1.63	9	1
1:A:472:ILE:HD13	1:A:472:ILE:N	0.53	2.18	5	1
1:A:462:ILE:HD12	2:B:907:GLY:HA2	0.53	1.81	8	1
1:A:399:LEU:HD21	1:A:484:LEU:HD22	0.52	1.80	4	1
1:A:430:GLU:HG3	2:B:905:LEU:HD21	0.52	1.81	12	1
1:A:491:LYS:HE3	1:A:492:ASN:N	0.52	2.19	13	1
1:A:433:GLN:HG3	2:B:906:THR:HG23	0.52	1.80	13	1
1:A:427:SER:HB3	1:A:436:VAL:HG21	0.52	1.81	13	1
1:A:403:THR:HB	2:B:903:LEU:O	0.52	2.05	1	4
1:A:426:PHE:HA	1:A:471:GLN:HG2	0.52	1.81	4	1
1:A:397:LEU:HD23	1:A:398:SER:N	0.52	2.19	15	2
1:A:463:ASN:ND2	1:A:463:ASN:H	0.52	2.02	5	1
1:A:401:ILE:O	1:A:409:THR:N	0.52	2.43	10	3
1:A:443:GLY:HA3	1:A:451:ASN:ND2	0.52	2.20	2	1
1:A:431:ASP:O	1:A:432:ASN:HB2	0.52	2.04	4	6
1:A:408:MET:HG2	1:A:451:ASN:ND2	0.52	2.20	5	1
1:A:408:MET:CG	1:A:449:ALA:HB3	0.52	2.34	7	1
1:A:489:LYS:HE2	1:A:495:LYS:HB3	0.52	1.81	9	1
1:A:400:GLY:O	1:A:441:LEU:O	0.52	2.28	7	7
1:A:427:SER:HB3	2:B:904:LEU:HD11	0.52	1.81	4	1
1:A:489:LYS:HG3	1:A:490:ASP:N	0.52	2.20	13	2
1:A:429:ALA:HB3	2:B:905:LEU:CA	0.51	2.36	3	1
1:A:441:LEU:HD22	1:A:451:ASN:HD21	0.51	1.64	4	1
1:A:407:VAL:O	1:A:441:LEU:HD23	0.51	2.05	6	1
1:A:408:MET:HB3	1:A:448:ALA:N	0.51	2.20	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:491:LYS:HB2	1:A:495:LYS:N	0.51	2.20	5	1
1:A:397:LEU:O	1:A:414:LYS:HD2	0.51	2.05	8	1
1:A:462:ILE:HG22	1:A:465:ALA:HB2	0.51	1.81	8	1
1:A:395:THR:HB	1:A:414:LYS:CB	0.51	2.33	6	1
1:A:397:LEU:HD21	1:A:442:GLN:NE2	0.51	2.19	6	1
1:A:429:ALA:HA	1:A:467:ARG:HA	0.51	1.81	14	1
1:A:496:GLU:HG2	1:A:497:GLN:N	0.51	2.21	5	1
1:A:499:ILE:HG13	1:A:500:THR:N	0.51	2.18	5	1
1:A:489:LYS:HD3	1:A:495:LYS:O	0.51	2.04	12	1
1:A:466:PRO:HG3	2:B:906:THR:HG23	0.51	1.81	14	1
1:A:472:ILE:HD12	1:A:490:ASP:OD1	0.51	2.06	5	1
1:A:420:THR:O	1:A:477:ASP:HA	0.51	2.05	5	5
1:A:434:SER:HA	1:A:462:ILE:C	0.51	2.31	4	3
1:A:477:ASP:N	1:A:485:HIS:HB2	0.51	2.20	13	1
1:A:399:LEU:HD22	1:A:478:ILE:HD11	0.51	1.81	11	1
1:A:408:MET:CG	1:A:450:ASP:HB3	0.51	2.35	11	1
1:A:479:ASP:HB2	1:A:483:ILE:HB	0.51	1.82	15	1
1:A:496:GLU:HG3	1:A:497:GLN:N	0.50	2.20	9	1
1:A:440:VAL:HB	1:A:454:LEU:HD13	0.50	1.82	3	8
1:A:459:LEU:HD13	1:A:460:ASP:N	0.50	2.21	2	2
1:A:442:GLN:O	1:A:452:LYS:HB3	0.50	2.05	11	3
1:A:462:ILE:O	1:A:464:PRO:HD2	0.50	2.06	14	1
1:A:459:LEU:HD11	1:A:496:GLU:HG2	0.50	1.82	9	1
1:A:486:VAL:O	1:A:498:LYS:O	0.50	2.30	13	1
1:A:397:LEU:C	1:A:414:LYS:HB2	0.50	2.32	6	1
1:A:442:GLN:HB2	1:A:454:LEU:HD21	0.50	1.82	10	1
1:A:500:THR:O	1:A:501:ILE:HD13	0.50	2.07	13	1
1:A:433:GLN:HG3	1:A:466:PRO:CA	0.50	2.36	14	1
1:A:398:SER:O	1:A:443:GLY:N	0.50	2.45	10	8
1:A:437:THR:OG1	2:B:906:THR:HA	0.50	2.07	15	4
1:A:489:LYS:HD3	1:A:490:ASP:N	0.50	2.22	15	1
1:A:396:PRO:O	1:A:397:LEU:HB2	0.50	2.05	6	1
1:A:473:GLU:HB3	1:A:489:LYS:HB3	0.50	1.83	10	1
1:A:478:ILE:HD13	1:A:484:LEU:HB2	0.50	1.83	8	4
1:A:462:ILE:HB	2:B:907:GLY:O	0.50	2.07	8	1
1:A:441:LEU:HD22	1:A:451:ASN:HB2	0.50	1.82	12	1
1:A:394:VAL:O	1:A:416:THR:HA	0.50	2.07	15	1
1:A:394:VAL:CG2	1:A:417:THR:HG22	0.50	2.37	15	1
1:A:412:ILE:HD11	1:A:476:PHE:CB	0.49	2.33	6	1
1:A:473:GLU:HB2	1:A:489:LYS:HG2	0.49	1.83	7	1
1:A:418:ILE:HG23	1:A:478:ILE:HG22	0.49	1.85	11	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:395:THR:CG2	1:A:416:THR:HB	0.49	2.36	15	1
1:A:481:ASP:HB3	1:A:483:ILE:HD11	0.49	1.85	11	1
1:A:471:GLN:HB3	1:A:491:LYS:HB2	0.49	1.83	1	1
1:A:489:LYS:HA	1:A:496:GLU:N	0.49	2.20	9	1
1:A:434:SER:HA	1:A:462:ILE:H	0.49	1.67	2	1
1:A:490:ASP:HB2	1:A:494:GLY:O	0.49	2.07	15	1
1:A:440:VAL:HB	1:A:454:LEU:CD2	0.49	2.38	10	1
1:A:417:THR:O	1:A:480:ALA:HA	0.49	2.07	15	1
1:A:440:VAL:CB	1:A:454:LEU:HB2	0.49	2.30	10	1
1:A:459:LEU:HD21	1:A:496:GLU:CD	0.49	2.33	9	1
1:A:399:LEU:CD1	1:A:442:GLN:HG3	0.49	2.35	11	1
1:A:408:MET:HE3	1:A:450:ASP:CB	0.49	2.37	11	1
2:B:903:LEU:O	2:B:904:LEU:HD22	0.49	2.08	13	2
1:A:400:GLY:O	1:A:441:LEU:N	0.49	2.46	10	8
1:A:472:ILE:HG21	1:A:488:ALA:HB1	0.49	1.85	4	1
1:A:440:VAL:O	1:A:454:LEU:HD12	0.49	2.07	6	1
1:A:431:ASP:HA	1:A:466:PRO:HB2	0.49	1.83	7	1
1:A:404:MET:HA	2:B:905:LEU:H	0.49	1.68	11	2
1:A:433:GLN:O	1:A:434:SER:C	0.49	2.55	8	2
1:A:433:GLN:C	1:A:435:ALA:H	0.49	2.16	5	3
1:A:434:SER:HB2	1:A:463:ASN:H	0.49	1.68	12	1
1:A:429:ALA:O	1:A:430:GLU:CB	0.49	2.61	14	1
1:A:451:ASN:CG	1:A:452:LYS:N	0.48	2.71	11	2
1:A:438:ILE:HD13	2:B:904:LEU:HD21	0.48	1.85	3	1
1:A:462:ILE:C	1:A:464:PRO:HD3	0.48	2.33	10	1
1:A:489:LYS:HD2	1:A:494:GLY:HA2	0.48	1.85	10	1
1:A:465:ALA:N	1:A:466:PRO:CD	0.48	2.76	7	1
1:A:500:THR:C	1:A:501:ILE:HD13	0.48	2.32	13	2
1:A:447:ARG:O	1:A:448:ALA:HB3	0.48	2.09	11	1
1:A:397:LEU:HB3	1:A:414:LYS:HG3	0.48	1.84	6	1
1:A:429:ALA:HB3	2:B:905:LEU:HG	0.48	1.85	7	1
1:A:401:ILE:CG2	1:A:440:VAL:HG13	0.48	2.37	10	1
1:A:401:ILE:HG23	1:A:411:LEU:HD22	0.48	1.84	6	1
1:A:443:GLY:N	1:A:450:ASP:HA	0.48	2.22	7	1
1:A:500:THR:C	1:A:501:ILE:HG13	0.48	2.32	10	2
1:A:442:GLN:O	1:A:442:GLN:HG3	0.48	2.09	3	2
1:A:466:PRO:HG2	1:A:469:MET:HB2	0.48	1.85	8	1
1:A:433:GLN:C	1:A:435:ALA:N	0.48	2.70	12	12
1:A:433:GLN:OE1	1:A:462:ILE:HB	0.48	2.09	11	2
1:A:476:PHE:CB	1:A:484:LEU:HD11	0.48	2.37	8	2
1:A:465:ALA:N	1:A:466:PRO:HD2	0.48	2.23	7	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:440:VAL:HG22	1:A:454:LEU:HG	0.48	1.86	13	1
1:A:434:SER:O	1:A:462:ILE:HG13	0.48	2.09	4	1
1:A:443:GLY:HA3	1:A:450:ASP:O	0.48	2.07	7	1
1:A:429:ALA:H	1:A:466:PRO:HB3	0.48	1.68	14	1
1:A:416:THR:OG1	1:A:478:ILE:HG21	0.48	2.08	15	1
1:A:481:ASP:HB2	1:A:483:ILE:HD11	0.47	1.85	1	1
1:A:433:GLN:O	1:A:435:ALA:N	0.47	2.47	8	2
1:A:431:ASP:HA	1:A:465:ALA:O	0.47	2.09	6	4
1:A:425:VAL:HG21	1:A:471:GLN:HG2	0.47	1.86	14	1
1:A:445:ARG:O	1:A:446:LYS:HB2	0.47	2.09	6	1
1:A:464:PRO:O	1:A:466:PRO:HD3	0.47	2.09	13	2
1:A:395:THR:HG21	1:A:412:ILE:CG2	0.47	2.35	15	1
1:A:457:PHE:HZ	1:A:501:ILE:HD11	0.47	1.68	5	1
1:A:428:THR:HB	1:A:468:GLY:N	0.47	2.23	14	1
1:A:436:VAL:HG22	2:B:904:LEU:HD13	0.47	1.84	4	1
1:A:395:THR:O	1:A:414:LYS:HA	0.47	2.08	8	1
1:A:483:ILE:HG22	1:A:484:LEU:N	0.47	2.25	15	1
1:A:408:MET:O	1:A:408:MET:HG3	0.47	2.09	14	1
1:A:431:ASP:N	1:A:466:PRO:O	0.47	2.47	14	1
1:A:442:GLN:CG	1:A:443:GLY:N	0.47	2.78	6	1
1:A:476:PHE:HB3	1:A:485:HIS:O	0.47	2.09	13	1
1:A:430:GLU:O	1:A:433:GLN:HG2	0.47	2.09	15	2
1:A:436:VAL:HG23	2:B:907:GLY:HA2	0.47	1.86	10	1
1:A:467:ARG:O	1:A:469:MET:N	0.47	2.47	14	1
1:A:401:ILE:O	1:A:409:THR:HB	0.47	2.09	12	2
1:A:471:GLN:O	1:A:489:LYS:O	0.47	2.32	5	1
1:A:463:ASN:CB	1:A:464:PRO:HD2	0.47	2.40	13	1
1:A:398:SER:HB3	1:A:443:GLY:O	0.47	2.10	15	1
1:A:408:MET:HB2	1:A:448:ALA:HB2	0.47	1.86	2	1
1:A:432:ASN:H	1:A:467:ARG:CB	0.47	2.23	14	1
2:B:904:LEU:C	2:B:906:THR:H	0.46	2.18	3	1
1:A:397:LEU:C	1:A:414:LYS:HB3	0.46	2.34	13	2
1:A:399:LEU:CD1	1:A:478:ILE:HD11	0.46	2.40	12	1
1:A:424:GLN:O	1:A:474:VAL:HG23	0.46	2.10	14	2
1:A:457:PHE:CE1	1:A:486:VAL:HB	0.46	2.44	14	1
1:A:439:HIS:CD2	1:A:441:LEU:HD13	0.46	2.45	11	1
1:A:491:LYS:HB2	1:A:496:GLU:HB3	0.46	1.87	12	1
1:A:399:LEU:HD23	1:A:400:GLY:N	0.46	2.25	12	3
1:A:408:MET:SD	1:A:447:ARG:HA	0.46	2.51	6	1
1:A:394:VAL:HG21	1:A:415:ASN:N	0.46	2.25	7	1
1:A:440:VAL:O	1:A:454:LEU:N	0.46	2.48	7	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:429:ALA:H	1:A:466:PRO:CB	0.46	2.22	14	1
1:A:470:PRO:HA	1:A:491:LYS:HD2	0.46	1.88	13	1
1:A:428:THR:HG22	1:A:470:PRO:HG2	0.46	1.87	1	1
1:A:429:ALA:HB2	2:B:903:LEU:HB3	0.46	1.85	4	2
1:A:476:PHE:CD1	1:A:486:VAL:HA	0.46	2.46	9	1
1:A:430:GLU:HB2	2:B:905:LEU:HG	0.46	1.85	11	1
1:A:441:LEU:HB3	1:A:450:ASP:OD1	0.46	2.11	11	1
1:A:426:PHE:CE1	2:B:904:LEU:HD11	0.46	2.46	12	1
1:A:440:VAL:CG2	1:A:454:LEU:HG	0.46	2.40	13	1
1:A:420:THR:CG2	1:A:478:ILE:HG13	0.46	2.40	15	1
1:A:412:ILE:CG2	1:A:420:THR:HG21	0.46	2.40	5	2
1:A:474:VAL:HG12	1:A:476:PHE:CE2	0.46	2.46	14	1
1:A:484:LEU:O	1:A:501:ILE:HD12	0.46	2.10	5	1
1:A:489:LYS:HD3	1:A:494:GLY:HA2	0.46	1.87	5	1
1:A:394:VAL:HG11	1:A:414:LYS:HG3	0.46	1.88	7	1
1:A:443:GLY:O	1:A:450:ASP:HA	0.46	2.10	7	1
1:A:434:SER:O	1:A:460:ASP:HA	0.46	2.10	9	1
1:A:408:MET:HE3	1:A:450:ASP:CG	0.46	2.36	11	1
1:A:433:GLN:HG3	1:A:466:PRO:HA	0.46	1.87	14	1
1:A:434:SER:HB2	1:A:464:PRO:O	0.46	2.10	14	1
1:A:440:VAL:O	1:A:454:LEU:HD22	0.46	2.10	10	1
1:A:463:ASN:O	1:A:465:ALA:N	0.46	2.49	13	3
1:A:440:VAL:HG23	1:A:454:LEU:HD21	0.46	1.87	13	1
1:A:469:MET:N	1:A:470:PRO:CD	0.46	2.79	9	2
1:A:430:GLU:H	1:A:433:GLN:HE22	0.46	1.53	10	1
1:A:463:ASN:O	1:A:464:PRO:C	0.45	2.58	13	4
1:A:416:THR:HG23	1:A:417:THR:N	0.45	2.26	7	1
1:A:438:ILE:HB	1:A:457:PHE:CE2	0.45	2.45	12	1
1:A:491:LYS:HD3	1:A:496:GLU:OE1	0.45	2.11	12	1
1:A:471:GLN:H	1:A:491:LYS:HD3	0.45	1.71	13	1
1:A:398:SER:CA	1:A:413:ALA:HA	0.45	2.42	12	7
1:A:439:HIS:CE1	1:A:441:LEU:HD13	0.45	2.46	7	1
1:A:427:SER:OG	2:B:904:LEU:HG	0.45	2.11	13	1
1:A:484:LEU:O	1:A:501:ILE:HG12	0.45	2.11	13	1
1:A:433:GLN:C	1:A:462:ILE:HB	0.45	2.36	14	1
1:A:435:ALA:CB	1:A:459:LEU:O	0.45	2.65	14	1
1:A:395:THR:C	1:A:413:ALA:HB3	0.45	2.37	15	1
2:B:906:THR:OG1	2:B:907:GLY:N	0.45	2.49	3	1
1:A:428:THR:O	2:B:903:LEU:HB3	0.45	2.12	4	1
1:A:417:THR:OG1	1:A:419:PRO:HD2	0.45	2.12	11	1
1:A:453:SER:O	1:A:454:LEU:C	0.45	2.59	7	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:407:VAL:HG12	1:A:447:ARG:HA	0.45	1.87	10	1
1:A:399:LEU:HD23	1:A:484:LEU:HD22	0.45	1.87	3	1
1:A:433:GLN:HG3	1:A:466:PRO:CD	0.45	2.41	14	1
1:A:433:GLN:OE1	2:B:906:THR:HG23	0.45	2.11	15	1
1:A:491:LYS:C	1:A:493:SER:H	0.45	2.20	12	1
1:A:472:ILE:HG23	1:A:489:LYS:O	0.45	2.11	4	1
1:A:397:LEU:N	1:A:414:LYS:HD2	0.45	2.26	8	1
1:A:433:GLN:HB3	2:B:907:GLY:N	0.45	2.27	8	1
1:A:399:LEU:CD2	1:A:478:ILE:HD11	0.45	2.42	11	1
1:A:399:LEU:HG	1:A:442:GLN:HB2	0.44	1.88	11	1
1:A:409:THR:CG2	1:A:410:THR:N	0.44	2.80	12	1
1:A:429:ALA:O	1:A:466:PRO:HB2	0.44	2.12	14	1
1:A:430:GLU:CA	1:A:466:PRO:O	0.44	2.65	14	1
1:A:465:ALA:H	1:A:466:PRO:CD	0.44	2.25	7	1
1:A:402:GLU:HG2	1:A:406:GLY:H	0.44	1.72	8	1
1:A:433:GLN:HB3	1:A:462:ILE:CD1	0.44	2.42	14	1
1:A:397:LEU:CA	1:A:414:LYS:HD2	0.44	2.43	8	1
1:A:408:MET:C	1:A:408:MET:SD	0.44	3.00	11	1
1:A:433:GLN:CG	1:A:466:PRO:HA	0.44	2.43	14	1
1:A:408:MET:N	1:A:448:ALA:HB2	0.44	2.28	12	2
1:A:425:VAL:HA	1:A:472:ILE:O	0.44	2.12	13	1
1:A:441:LEU:HG	1:A:450:ASP:HB3	0.44	1.89	7	1
1:A:429:ALA:CB	2:B:905:LEU:HB3	0.44	2.41	13	1
1:A:417:THR:H	1:A:418:ILE:HG23	0.44	1.72	15	1
1:A:408:MET:SD	1:A:441:LEU:HD12	0.44	2.53	2	1
1:A:418:ILE:N	1:A:419:PRO:HD2	0.44	2.27	12	2
1:A:462:ILE:O	1:A:463:ASN:O	0.44	2.36	3	1
1:A:496:GLU:CG	1:A:497:GLN:N	0.44	2.81	9	1
1:A:428:THR:HG22	1:A:466:PRO:HA	0.44	1.89	15	1
1:A:428:THR:HG22	1:A:469:MET:HE3	0.44	1.90	5	1
1:A:397:LEU:CG	1:A:398:SER:H	0.44	2.13	6	1
1:A:397:LEU:HB2	1:A:414:LYS:CD	0.44	2.37	15	1
1:A:399:LEU:N	1:A:413:ALA:HA	0.44	2.28	15	1
1:A:395:THR:HG21	1:A:412:ILE:HG21	0.44	1.89	1	1
1:A:399:LEU:HB2	1:A:412:ILE:O	0.44	2.13	11	2
1:A:439:HIS:ND1	1:A:456:GLN:HB2	0.44	2.27	6	1
1:A:486:VAL:O	1:A:498:LYS:HA	0.44	2.12	6	3
1:A:441:LEU:HD22	1:A:451:ASN:ND2	0.44	2.27	4	1
1:A:430:GLU:O	1:A:433:GLN:HG3	0.44	2.13	5	1
1:A:408:MET:HG3	1:A:448:ALA:CA	0.44	2.43	10	1
1:A:432:ASN:HB2	1:A:464:PRO:HB2	0.44	1.89	14	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:442:GLN:O	1:A:442:GLN:CG	0.43	2.65	11	1
1:A:450:ASP:O	1:A:451:ASN:HB3	0.43	2.13	11	1
1:A:417:THR:OG1	1:A:480:ALA:C	0.43	2.61	15	1
1:A:397:LEU:CB	1:A:414:LYS:HE2	0.43	2.39	6	1
1:A:408:MET:CE	1:A:441:LEU:HD12	0.43	2.43	13	1
1:A:414:LYS:C	1:A:416:THR:H	0.43	2.20	15	1
1:A:442:GLN:HG2	1:A:451:ASN:C	0.43	2.39	6	1
1:A:396:PRO:HA	1:A:414:LYS:HB2	0.43	1.89	8	1
1:A:487:SER:CA	1:A:498:LYS:O	0.43	2.66	13	1
1:A:435:ALA:CA	1:A:459:LEU:O	0.43	2.64	8	1
1:A:411:LEU:HG	1:A:476:PHE:CE2	0.43	2.49	9	1
1:A:428:THR:HG21	1:A:466:PRO:HB3	0.43	1.91	7	1
1:A:454:LEU:O	1:A:501:ILE:HG21	0.43	2.14	10	1
1:A:394:VAL:HG12	1:A:417:THR:HA	0.43	1.91	11	1
1:A:398:SER:HB3	1:A:414:LYS:HD3	0.43	1.89	13	1
1:A:394:VAL:HG12	1:A:416:THR:O	0.43	2.13	2	1
1:A:442:GLN:HB3	1:A:452:LYS:O	0.43	2.14	6	1
1:A:394:VAL:HG13	1:A:395:THR:N	0.43	2.28	7	1
1:A:402:GLU:HG2	1:A:439:HIS:ND1	0.43	2.27	10	1
1:A:448:ALA:O	1:A:449:ALA:HB3	0.43	2.13	11	1
1:A:467:ARG:HD3	2:B:903:LEU:HD21	0.43	1.91	13	1
1:A:479:ASP:C	1:A:481:ASP:N	0.43	2.77	15	1
1:A:408:MET:HB2	1:A:441:LEU:HG	0.43	1.90	1	1
2:B:905:LEU:HG	2:B:906:THR:N	0.43	2.29	14	1
1:A:462:ILE:HG21	1:A:470:PRO:HB3	0.43	1.91	1	1
1:A:485:HIS:HA	1:A:499:ILE:O	0.43	2.13	9	1
1:A:434:SER:HA	1:A:462:ILE:HB	0.43	1.91	12	1
1:A:459:LEU:HG	1:A:488:ALA:HB3	0.43	1.89	15	1
1:A:489:LYS:CE	1:A:495:LYS:HA	0.43	2.43	15	1
1:A:442:GLN:C	1:A:451:ASN:OD1	0.42	2.61	2	1
1:A:429:ALA:HB3	2:B:904:LEU:C	0.42	2.39	4	1
1:A:476:PHE:CD1	1:A:476:PHE:N	0.42	2.85	8	3
1:A:474:VAL:HB	1:A:476:PHE:HE1	0.42	1.74	6	1
1:A:428:THR:CG2	1:A:466:PRO:HB3	0.42	2.44	7	1
1:A:437:THR:OG1	2:B:906:THR:HG23	0.42	2.14	7	1
1:A:430:GLU:N	1:A:433:GLN:NE2	0.42	2.67	10	1
1:A:441:LEU:HD21	1:A:448:ALA:HB1	0.42	1.90	11	1
1:A:434:SER:HA	1:A:463:ASN:N	0.42	2.28	3	1
1:A:402:GLU:HG3	1:A:407:VAL:O	0.42	2.14	8	1
1:A:457:PHE:HD2	1:A:498:LYS:HB3	0.42	1.74	13	1
1:A:402:GLU:N	1:A:439:HIS:O	0.42	2.51	14	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:444:GLU:HG3	1:A:450:ASP:HB3	0.42	1.89	6	1
1:A:472:ILE:HG13	1:A:489:LYS:O	0.42	2.14	9	1
1:A:436:VAL:CA	2:B:907:GLY:HA2	0.42	2.43	10	1
1:A:457:PHE:CD2	1:A:498:LYS:HB3	0.42	2.49	13	1
1:A:401:ILE:HG12	1:A:411:LEU:CD2	0.42	2.44	5	2
1:A:451:ASN:OD1	1:A:451:ASN:N	0.42	2.51	4	1
1:A:436:VAL:HG12	1:A:459:LEU:HB2	0.42	1.92	9	1
1:A:487:SER:HA	1:A:498:LYS:O	0.42	2.13	13	1
1:A:428:THR:HG22	1:A:467:ARG:N	0.42	2.29	14	1
1:A:428:THR:HG23	1:A:465:ALA:HB1	0.42	1.91	13	1
1:A:428:THR:CG2	1:A:429:ALA:N	0.42	2.81	15	1
1:A:431:ASP:O	1:A:432:ASN:CB	0.42	2.68	5	1
1:A:398:SER:N	1:A:414:LYS:CG	0.42	2.77	6	1
1:A:426:PHE:HE1	2:B:902:ARG:HB2	0.42	1.74	15	1
1:A:483:ILE:HG22	1:A:484:LEU:H	0.42	1.73	15	1
1:A:401:ILE:CG2	1:A:411:LEU:HD22	0.42	2.44	6	1
1:A:476:PHE:CE1	1:A:486:VAL:HG13	0.42	2.50	9	1
1:A:428:THR:HG23	1:A:429:ALA:N	0.42	2.29	11	1
1:A:411:LEU:HD22	1:A:476:PHE:CD1	0.42	2.50	12	1
1:A:432:ASN:CB	1:A:464:PRO:HD2	0.42	2.44	14	1
1:A:462:ILE:HG12	1:A:490:ASP:OD2	0.42	2.14	15	1
1:A:437:THR:O	2:B:904:LEU:HD22	0.42	2.15	4	1
1:A:429:ALA:C	1:A:466:PRO:HB2	0.42	2.40	14	1
1:A:433:GLN:HB3	1:A:462:ILE:HD12	0.42	1.91	14	1
1:A:427:SER:O	2:B:903:LEU:HA	0.42	2.14	2	2
1:A:408:MET:CB	1:A:448:ALA:HB2	0.42	2.45	3	2
1:A:463:ASN:HB2	1:A:464:PRO:HD2	0.42	1.92	3	1
1:A:442:GLN:HB3	1:A:452:LYS:H	0.42	1.75	6	1
1:A:414:LYS:CG	1:A:415:ASN:N	0.42	2.82	15	1
1:A:426:PHE:O	1:A:426:PHE:CG	0.42	2.73	4	1
1:A:475:THR:O	1:A:486:VAL:HA	0.42	2.15	5	1
1:A:432:ASN:O	1:A:466:PRO:C	0.42	2.63	14	1
1:A:489:LYS:HD3	1:A:489:LYS:C	0.42	2.40	15	1
1:A:412:ILE:HG22	1:A:416:THR:HG21	0.41	1.92	7	1
1:A:397:LEU:N	1:A:414:LYS:HG3	0.41	2.30	11	1
1:A:408:MET:HE1	1:A:410:THR:HG23	0.41	1.92	11	1
1:A:484:LEU:HB2	1:A:501:ILE:O	0.41	2.15	11	1
1:A:499:ILE:HD13	1:A:499:ILE:N	0.41	2.29	3	1
1:A:397:LEU:HB3	1:A:442:GLN:CD	0.41	2.39	13	1
1:A:432:ASN:HB3	1:A:462:ILE:O	0.41	2.15	14	1
1:A:400:GLY:CA	1:A:410:THR:HA	0.41	2.32	15	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:420:THR:HG22	1:A:478:ILE:CB	0.41	2.40	15	1
1:A:471:GLN:HG3	1:A:489:LYS:HE3	0.41	1.91	5	1
1:A:472:ILE:HG23	1:A:490:ASP:OD1	0.41	2.14	5	1
1:A:404:MET:HA	2:B:905:LEU:N	0.41	2.30	10	1
1:A:440:VAL:HB	1:A:454:LEU:HD12	0.41	1.91	11	1
1:A:470:PRO:HG2	1:A:490:ASP:OD1	0.41	2.16	11	1
1:A:397:LEU:HB2	1:A:414:LYS:HB3	0.41	1.90	15	1
1:A:443:GLY:HA3	1:A:445:ARG:HE	0.41	1.75	15	1
1:A:418:ILE:O	1:A:478:ILE:HB	0.41	2.16	2	2
1:A:469:MET:O	1:A:470:PRO:C	0.41	2.63	5	1
1:A:489:LYS:HB2	1:A:495:LYS:CB	0.41	2.44	9	1
1:A:491:LYS:O	1:A:492:ASN:HB2	0.41	2.15	9	1
1:A:412:ILE:HD13	1:A:478:ILE:HG12	0.41	1.92	11	1
1:A:427:SER:O	2:B:904:LEU:HD12	0.41	2.16	12	1
1:A:472:ILE:HG22	1:A:474:VAL:HG22	0.41	1.92	12	1
1:A:485:HIS:HB3	1:A:499:ILE:CG2	0.41	2.45	13	1
1:A:445:ARG:HG3	1:A:447:ARG:O	0.41	2.16	15	1
1:A:402:GLU:HG3	1:A:406:GLY:H	0.41	1.74	6	1
1:A:442:GLN:HB3	1:A:452:LYS:N	0.41	2.30	6	1
1:A:403:THR:HA	2:B:904:LEU:CD1	0.41	2.45	13	1
1:A:451:ASN:HD22	1:A:452:LYS:H	0.41	1.59	1	1
2:B:905:LEU:HG	2:B:907:GLY:H	0.41	1.74	2	1
1:A:414:LYS:H	1:A:414:LYS:HG3	0.41	1.52	4	1
1:A:489:LYS:HB2	1:A:496:GLU:HB2	0.41	1.92	13	2
1:A:502:LYS:H	1:A:502:LYS:HG3	0.41	1.55	5	1
1:A:429:ALA:O	2:B:905:LEU:HA	0.41	2.16	10	1
1:A:399:LEU:CD1	1:A:484:LEU:HD22	0.41	2.46	13	1
1:A:427:SER:HB3	1:A:436:VAL:CG2	0.41	2.45	13	1
1:A:397:LEU:CA	1:A:414:LYS:HB2	0.41	2.46	6	1
1:A:457:PHE:CE1	1:A:499:ILE:HD12	0.41	2.51	7	1
1:A:397:LEU:HG	1:A:398:SER:OG	0.41	2.16	8	1
1:A:483:ILE:HA	1:A:502:LYS:HA	0.41	1.92	10	1
1:A:440:VAL:CG2	1:A:455:GLY:O	0.41	2.69	11	1
1:A:428:THR:N	2:B:904:LEU:HB2	0.41	2.31	13	1
1:A:431:ASP:O	1:A:465:ALA:O	0.41	2.39	14	1
1:A:489:LYS:HB2	1:A:496:GLU:HG3	0.41	1.92	6	1
1:A:439:HIS:CE1	1:A:453:SER:HB3	0.41	2.51	8	1
1:A:474:VAL:HB	1:A:476:PHE:CZ	0.41	2.51	8	1
1:A:397:LEU:HD23	1:A:442:GLN:NE2	0.41	2.31	11	1
2:B:903:LEU:HB3	2:B:904:LEU:H	0.41	1.61	14	1
1:A:418:ILE:HD11	1:A:420:THR:HB	0.41	1.92	15	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:408:MET:HE1	1:A:443:GLY:N	0.40	2.31	3	1
1:A:410:THR:O	1:A:411:LEU:HD22	0.40	2.16	8	1
1:A:432:ASN:H	1:A:466:PRO:HD3	0.40	1.75	9	1
1:A:500:THR:O	1:A:501:ILE:HG13	0.40	2.15	12	1
1:A:463:ASN:N	1:A:464:PRO:CD	0.40	2.84	15	1
1:A:395:THR:OG1	1:A:416:THR:HG22	0.40	2.16	1	1
1:A:436:VAL:HA	2:B:906:THR:HG22	0.40	1.92	3	1
1:A:408:MET:HG3	1:A:449:ALA:CB	0.40	2.46	7	1
1:A:428:THR:CG2	1:A:466:PRO:HB2	0.40	2.46	9	1
2:B:902:ARG:HB3	2:B:903:LEU:H	0.40	1.53	9	1
1:A:466:PRO:HG2	2:B:906:THR:N	0.40	2.29	14	1
1:A:417:THR:C	1:A:418:ILE:CG2	0.40	2.89	15	1
1:A:397:LEU:C	1:A:414:LYS:HD2	0.40	2.41	8	1
1:A:427:SER:HB2	1:A:472:ILE:CB	0.40	2.41	13	1
1:A:427:SER:O	1:A:427:SER:OG	0.40	2.32	14	1
1:A:477:ASP:O	1:A:484:LEU:HA	0.40	2.17	5	1
1:A:489:LYS:HD3	1:A:495:LYS:C	0.40	2.40	12	1
1:A:427:SER:HA	2:B:904:LEU:CD2	0.40	2.41	13	1
1:A:427:SER:OG	2:B:904:LEU:HD12	0.40	2.16	14	1
1:A:430:GLU:HB3	1:A:433:GLN:HG2	0.40	1.92	4	1
1:A:408:MET:HG3	1:A:449:ALA:CA	0.40	2.46	7	1
1:A:442:GLN:C	1:A:450:ASP:O	0.40	2.64	7	1
1:A:395:THR:CB	1:A:413:ALA:HB3	0.40	2.44	15	1
1:A:418:ILE:O	1:A:478:ILE:O	0.40	2.40	15	1

6.3 Torsion angles ⓘ

6.3.1 Protein backbone ⓘ

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	110/135 (81%)	81±4 (74±4%)	20±4 (18±3%)	9±2 (8±2%)	1	13
2	B	5/7 (71%)	3±1 (67±21%)	1±1 (28±18%)	0±1 (5±11%)	2	22
All	All	1725/2130 (81%)	1272 (74%)	315 (18%)	138 (8%)	1	13

All 47 unique Ramachandran outliers are listed below. They are sorted by the frequency of

occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	405	GLY	11
1	A	492	ASN	10
1	A	461	GLY	9
1	A	467	ARG	8
1	A	464	PRO	7
1	A	450	ASP	7
1	A	430	GLU	6
1	A	466	PRO	6
1	A	414	LYS	5
1	A	463	ASN	5
1	A	462	ILE	5
1	A	413	ALA	3
1	A	448	ALA	3
1	A	503	ALA	3
1	A	470	PRO	3
1	A	469	MET	3
1	A	451	ASN	3
1	A	432	ASN	2
2	B	903	LEU	2
1	A	434	SER	2
1	A	396	PRO	2
1	A	397	LEU	2
1	A	449	ALA	2
1	A	465	ALA	2
1	A	495	LYS	2
1	A	452	LYS	2
1	A	447	ARG	2
1	A	468	GLY	2
1	A	394	VAL	1
1	A	494	GLY	1
1	A	415	ASN	1
1	A	431	ASP	1
1	A	454	LEU	1
1	A	491	LYS	1
1	A	493	SER	1
2	B	904	LEU	1
2	B	906	THR	1
1	A	404	MET	1
1	A	499	ILE	1
1	A	433	GLN	1
1	A	453	SER	1
1	A	395	THR	1

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Mol	Chain	Res	Type	Models (Total)
1	A	398	SER	1
1	A	412	ILE	1
1	A	416	THR	1
1	A	417	THR	1
1	A	480	ALA	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	92/113 (81%)	54±3 (58±3%)	38±3 (42±3%)	0 4
2	B	5/6 (83%)	3±1 (59±24%)	2±1 (41±24%)	0 4
All	All	1455/1785 (82%)	850 (58%)	605 (42%)	0 4

All 86 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	395	THR	15
1	A	440	VAL	15
1	A	422	HIS	14
1	A	491	LYS	14
1	A	498	LYS	14
1	A	410	THR	13
1	A	441	LEU	13
1	A	459	LEU	13
1	A	399	LEU	12
1	A	460	ASP	12
1	A	462	ILE	12
1	A	476	PHE	12
1	A	414	LYS	12
1	A	397	LEU	11
1	A	469	MET	11
1	A	487	SER	11
1	A	408	MET	10
1	A	411	LEU	10
1	A	428	THR	10

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Mol	Chain	Res	Type	Models (Total)
1	A	433	GLN	10
1	A	437	THR	10
1	A	495	LYS	10
1	A	424	GLN	10
1	A	446	LYS	10
1	A	477	ASP	10
1	A	489	LYS	10
1	A	421	LYS	9
1	A	431	ASP	9
1	A	467	ARG	9
2	B	902	ARG	9
1	A	417	THR	8
1	A	427	SER	8
1	A	444	GLU	8
1	A	445	ARG	8
1	A	451	ASN	8
1	A	496	GLU	8
1	A	502	LYS	8
1	A	398	SER	8
1	A	493	SER	8
1	A	471	GLN	8
1	A	415	ASN	7
2	B	905	LEU	7
1	A	499	ILE	7
1	A	497	GLN	7
1	A	474	VAL	7
1	A	472	ILE	6
1	A	407	VAL	6
1	A	458	ASN	6
1	A	490	ASP	6
2	B	904	LEU	6
1	A	453	SER	6
1	A	412	ILE	5
1	A	423	SER	5
1	A	430	GLU	5
1	A	436	VAL	5
1	A	483	ILE	5
2	B	903	LEU	5
1	A	501	ILE	5
1	A	456	GLN	5
1	A	426	PHE	5
1	A	447	ARG	5

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Mol	Chain	Res	Type	Models (Total)
1	A	452	LYS	5
1	A	425	VAL	4
1	A	434	SER	4
1	A	481	ASP	4
1	A	463	ASN	4
1	A	485	HIS	4
1	A	450	ASP	4
1	A	442	GLN	4
2	B	906	THR	4
1	A	492	ASN	3
1	A	394	VAL	3
1	A	454	LEU	3
1	A	420	THR	2
1	A	473	GLU	2
1	A	478	ILE	2
1	A	484	LEU	2
1	A	479	ASP	2
1	A	486	VAL	1
1	A	404	MET	1
1	A	402	GLU	1
1	A	403	THR	1
1	A	409	THR	1
1	A	457	PHE	1
1	A	418	ILE	1
1	A	432	ASN	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

6.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.6 Ligand geometry

There are no ligands in this entry.

6.7 Other polymers

There are no such molecules in this entry.

6.8 Polymer linkage issues

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided