



Full wwPDB NMR Structure Validation Report ⓘ

Apr 27, 2016 – 03:50 AM BST

PDB ID : 2MZ7
Title : Structure of Tau(267-312) bound to Microtubules
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Deposited on : 2015-02-06

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/NMRValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

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

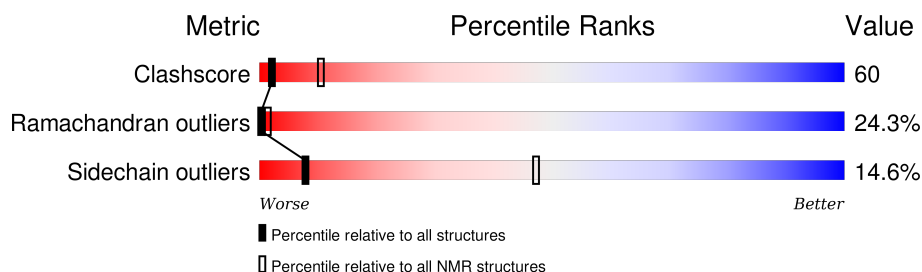
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 8%.

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



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	114402	11133
Ramachandran outliers	111179	9975
Sidechain outliers	111093	9958

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

Mol	Chain	Length	Quality of chain
1	A	46	<div> <div></div> <div>20%</div> <div>67%</div> <div>13%</div> </div>

2 Ensemble composition and analysis ⓘ

This entry contains 20 models.

Cyrange was unable to find well-defined residues.

Error message: The number of core atoms (5) was below the domain threshold value (8).

The models can be grouped into 2 clusters. No single-model clusters were found.

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

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 708 atoms, of which 365 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Microtubule-associated protein tau.

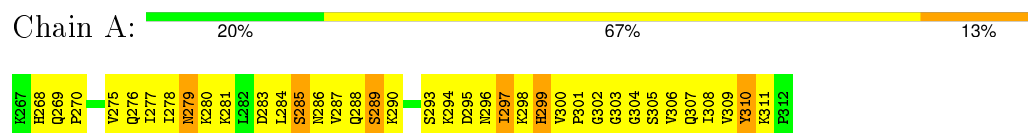
Mol	Chain	Residues	Atoms						Trace
1	A	46	Total	C	H	N	O	S	0
			708	214	365	65	63	1	

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: Microtubule-associated protein tau

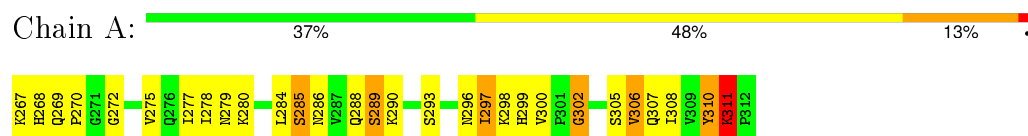


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

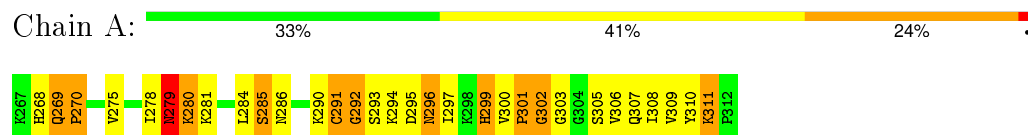
4.2.1 Score per residue for model 1

- Molecule 1: Microtubule-associated protein tau



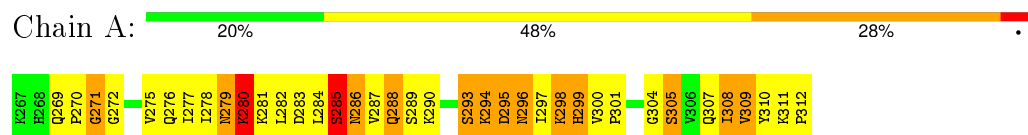
4.2.2 Score per residue for model 2

- Molecule 1: Microtubule-associated protein tau



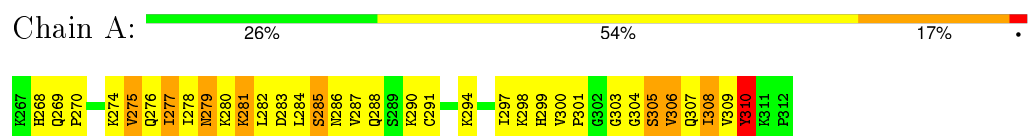
4.2.3 Score per residue for model 3

- Molecule 1: Microtubule-associated protein tau



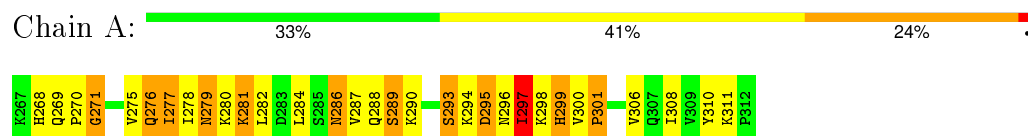
4.2.4 Score per residue for model 4

- Molecule 1: Microtubule-associated protein tau



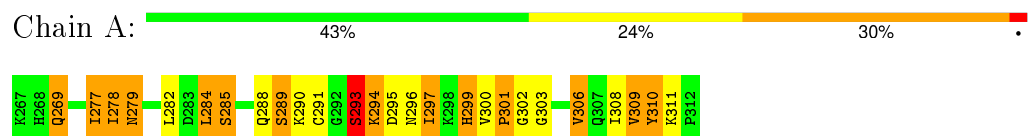
4.2.5 Score per residue for model 5

- Molecule 1: Microtubule-associated protein tau



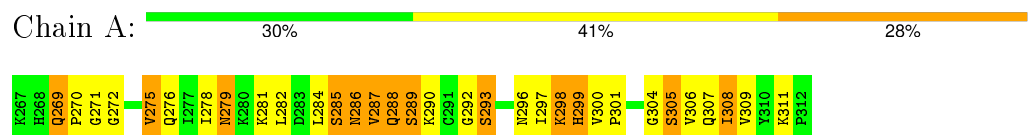
4.2.6 Score per residue for model 6

- Molecule 1: Microtubule-associated protein tau



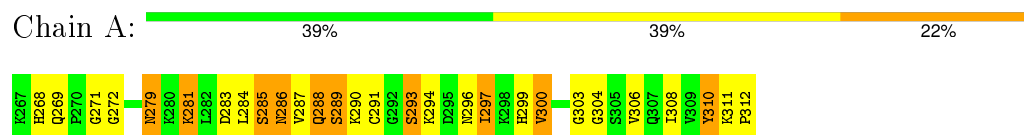
4.2.7 Score per residue for model 7

- Molecule 1: Microtubule-associated protein tau



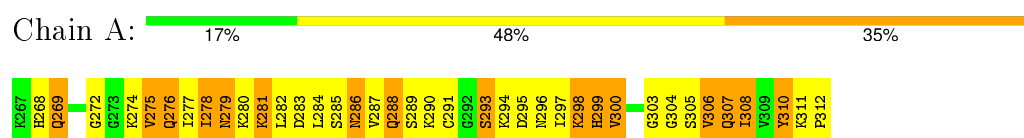
4.2.8 Score per residue for model 8

- Molecule 1: Microtubule-associated protein tau



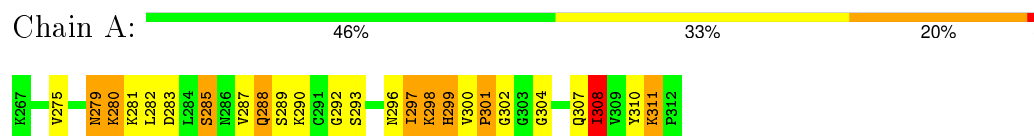
4.2.9 Score per residue for model 9

- Molecule 1: Microtubule-associated protein tau



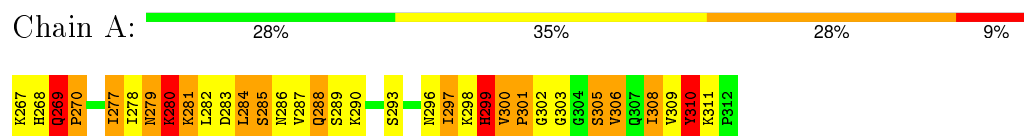
4.2.10 Score per residue for model 10

- Molecule 1: Microtubule-associated protein tau



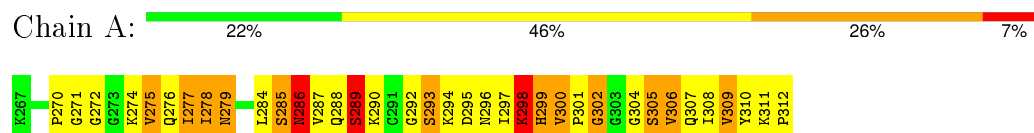
4.2.11 Score per residue for model 11

- Molecule 1: Microtubule-associated protein tau



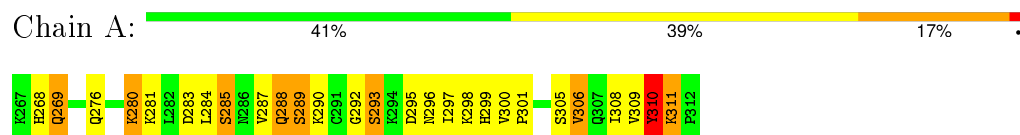
4.2.12 Score per residue for model 12

- Molecule 1: Microtubule-associated protein tau



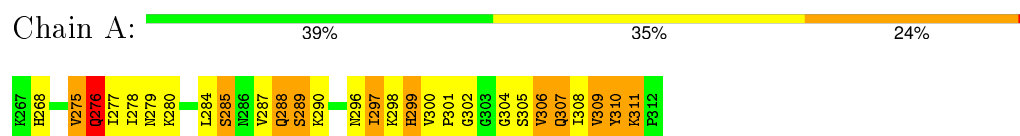
4.2.13 Score per residue for model 13

- Molecule 1: Microtubule-associated protein tau



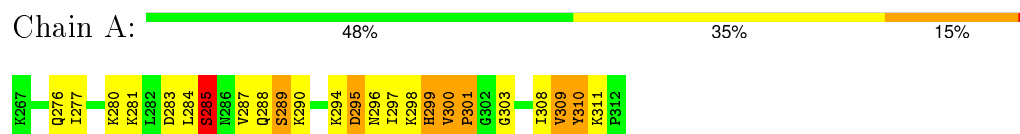
4.2.14 Score per residue for model 14

- Molecule 1: Microtubule-associated protein tau



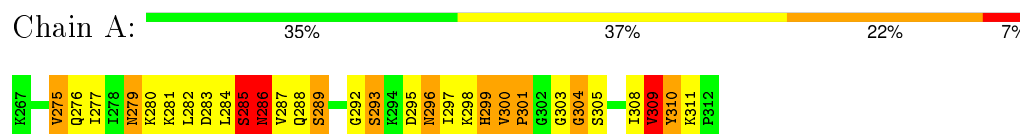
4.2.15 Score per residue for model 15

- Molecule 1: Microtubule-associated protein tau



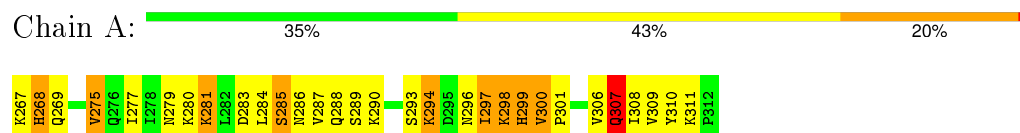
4.2.16 Score per residue for model 16

- Molecule 1: Microtubule-associated protein tau



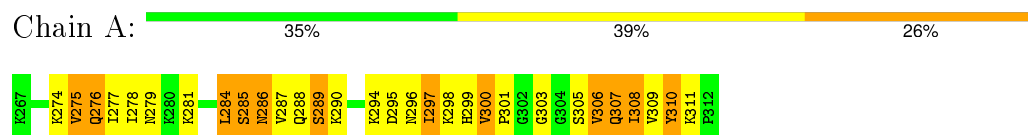
4.2.17 Score per residue for model 17

- Molecule 1: Microtubule-associated protein tau



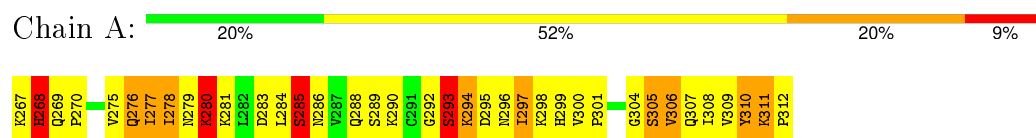
4.2.18 Score per residue for model 18

- Molecule 1: Microtubule-associated protein tau



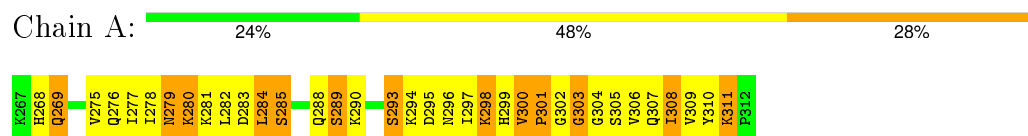
4.2.19 Score per residue for model 19

- Molecule 1: Microtubule-associated protein tau



4.2.20 Score per residue for model 20

- Molecule 1: Microtubule-associated protein tau



5 Refinement protocol and experimental data overview

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

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

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

Software name	Classification	Version
X-PLOR NIH	refinement	2.33
CYANA	structure solution	3.0

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

Chemical shift file(s)	2mz7_cs.str
Number of chemical shift lists	1
Total number of shifts	43
Number of shifts mapped to atoms	43
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	8%

No validations of the models with respect to experimental NMR restraints is performed at this time.

6 Model quality

6.1 Standard geometry

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

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	343	365	362	42±8
All	All	6860	7300	7240	840

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 60.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:308:ILE:HD12	1:A:309:VAL:N	0.95	1.76	14	5
1:A:284:LEU:H	1:A:284:LEU:HD22	0.93	1.23	11	3
1:A:308:ILE:HD12	1:A:308:ILE:O	0.88	1.67	19	2
1:A:299:HIS:O	1:A:300:VAL:HG13	0.82	1.73	4	1
1:A:308:ILE:HD13	1:A:308:ILE:O	0.82	1.75	7	2
1:A:284:LEU:N	1:A:284:LEU:HD13	0.81	1.90	11	2
1:A:284:LEU:HD22	1:A:284:LEU:N	0.80	1.90	18	3
1:A:294:LYS:O	1:A:297:ILE:HG23	0.78	1.79	17	1
1:A:284:LEU:N	1:A:284:LEU:HD12	0.76	1.93	4	2
1:A:296:ASN:O	1:A:297:ILE:HG22	0.74	1.82	10	1
1:A:275:VAL:HG22	1:A:276:GLN:N	0.74	1.96	7	1
1:A:300:VAL:H	1:A:301:PRO:CD	0.72	1.97	17	1
1:A:278:ILE:N	1:A:278:ILE:HD12	0.71	2.00	7	1
1:A:279:ASN:ND2	1:A:279:ASN:N	0.70	2.39	8	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:308:ILE:HD12	1:A:308:ILE:C	0.70	2.07	6	4
1:A:297:ILE:N	1:A:297:ILE:HD12	0.70	2.02	20	1
1:A:297:ILE:O	1:A:297:ILE:HG23	0.70	1.84	10	1
1:A:297:ILE:HD12	1:A:297:ILE:C	0.69	2.07	15	1
1:A:297:ILE:HG22	1:A:297:ILE:O	0.69	1.86	2	1
1:A:288:GLN:O	1:A:290:LYS:N	0.68	2.25	15	16
1:A:300:VAL:N	1:A:301:PRO:CD	0.68	2.55	5	9
1:A:299:HIS:O	1:A:300:VAL:HG23	0.68	1.89	12	1
1:A:305:SER:H	1:A:307:GLN:HE22	0.68	1.28	9	1
1:A:308:ILE:HD13	1:A:310:TYR:O	0.68	1.87	5	1
1:A:283:ASP:C	1:A:284:LEU:HD13	0.68	2.08	11	1
1:A:297:ILE:O	1:A:299:HIS:N	0.68	2.27	12	9
1:A:300:VAL:HG13	1:A:305:SER:O	0.68	1.87	3	1
1:A:297:ILE:HD13	1:A:297:ILE:O	0.67	1.89	6	2
1:A:278:ILE:H	1:A:278:ILE:HD12	0.67	1.48	7	1
1:A:299:HIS:NE2	1:A:302:GLY:N	0.66	2.41	12	1
1:A:297:ILE:H	1:A:297:ILE:HD12	0.66	1.49	20	1
1:A:299:HIS:CD2	1:A:300:VAL:H	0.66	2.07	12	1
1:A:279:ASN:H	1:A:279:ASN:HD22	0.66	1.33	11	1
1:A:269:GLN:NE2	1:A:286:ASN:ND2	0.65	2.44	5	1
1:A:306:VAL:O	1:A:306:VAL:HG22	0.65	1.89	6	2
1:A:299:HIS:C	1:A:308:ILE:HD11	0.65	2.11	1	1
1:A:279:ASN:O	1:A:281:LYS:N	0.65	2.29	3	5
1:A:299:HIS:N	1:A:299:HIS:ND1	0.65	2.44	11	1
1:A:281:LYS:CD	1:A:281:LYS:H	0.65	2.04	8	3
1:A:279:ASN:HD22	1:A:279:ASN:N	0.65	1.87	6	2
1:A:286:ASN:N	1:A:286:ASN:HD22	0.65	1.90	5	1
1:A:300:VAL:O	1:A:304:GLY:N	0.64	2.29	16	8
1:A:296:ASN:O	1:A:299:HIS:CG	0.64	2.50	6	4
1:A:284:LEU:N	1:A:288:GLN:HE21	0.64	1.90	15	1
1:A:308:ILE:O	1:A:310:TYR:N	0.64	2.31	14	2
1:A:306:VAL:N	1:A:307:GLN:NE2	0.64	2.46	18	2
1:A:279:ASN:N	1:A:279:ASN:HD22	0.64	1.90	16	3
1:A:296:ASN:O	1:A:299:HIS:ND1	0.64	2.31	6	1
1:A:308:ILE:C	1:A:308:ILE:HD13	0.64	2.13	10	1
1:A:284:LEU:H	1:A:288:GLN:HE21	0.63	1.36	15	1
1:A:267:LYS:N	1:A:280:LYS:CD	0.63	2.61	1	1
1:A:310:TYR:CG	1:A:310:TYR:O	0.63	2.50	1	3
1:A:275:VAL:HG13	1:A:276:GLN:H	0.63	1.53	7	1
1:A:308:ILE:HD13	1:A:308:ILE:H	0.63	1.53	18	1
1:A:299:HIS:HD1	1:A:299:HIS:C	0.63	1.97	16	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:299:HIS:H	1:A:299:HIS:CD2	0.63	2.10	17	2
1:A:300:VAL:O	1:A:300:VAL:HG22	0.63	1.93	20	2
1:A:286:ASN:ND2	1:A:287:VAL:H	0.62	1.92	9	2
1:A:268:HIS:O	1:A:270:PRO:N	0.62	2.32	11	2
1:A:296:ASN:O	1:A:299:HIS:CD2	0.62	2.53	20	3
1:A:299:HIS:C	1:A:299:HIS:ND1	0.62	2.52	8	3
1:A:308:ILE:N	1:A:308:ILE:HD13	0.62	2.10	3	1
1:A:287:VAL:O	1:A:289:SER:N	0.62	2.33	3	5
1:A:311:LYS:N	1:A:312:PRO:CD	0.62	2.63	3	1
1:A:305:SER:O	1:A:307:GLN:N	0.62	2.33	2	2
1:A:310:TYR:N	1:A:310:TYR:CD1	0.62	2.67	8	4
1:A:296:ASN:O	1:A:297:ILE:CG2	0.62	2.47	10	2
1:A:299:HIS:ND1	1:A:299:HIS:C	0.62	2.53	4	3
1:A:300:VAL:O	1:A:302:GLY:N	0.62	2.33	20	5
1:A:308:ILE:HD12	1:A:309:VAL:H	0.62	1.54	11	2
1:A:310:TYR:O	1:A:310:TYR:CD2	0.61	2.53	3	1
1:A:310:TYR:CD1	1:A:310:TYR:O	0.61	2.54	12	4
1:A:276:GLN:NE2	1:A:277:ILE:O	0.61	2.33	15	2
1:A:284:LEU:N	1:A:284:LEU:CD2	0.61	2.63	18	1
1:A:310:TYR:CD2	1:A:310:TYR:O	0.61	2.54	11	1
1:A:305:SER:N	1:A:307:GLN:HE22	0.61	1.94	14	2
1:A:275:VAL:O	1:A:275:VAL:HG12	0.61	1.95	17	1
1:A:295:ASP:OD1	1:A:296:ASN:N	0.61	2.33	6	1
1:A:285:SER:O	1:A:289:SER:N	0.61	2.33	9	1
1:A:284:LEU:O	1:A:286:ASN:N	0.61	2.34	16	4
1:A:294:LYS:O	1:A:296:ASN:N	0.61	2.33	9	5
1:A:277:ILE:O	1:A:279:ASN:N	0.61	2.33	11	5
1:A:284:LEU:H	1:A:288:GLN:NE2	0.61	1.93	15	1
1:A:286:ASN:ND2	1:A:287:VAL:N	0.61	2.47	9	3
1:A:284:LEU:CD1	1:A:284:LEU:N	0.61	2.60	11	2
1:A:275:VAL:O	1:A:277:ILE:N	0.61	2.33	19	2
1:A:285:SER:O	1:A:287:VAL:N	0.60	2.34	11	10
1:A:299:HIS:O	1:A:299:HIS:CD2	0.60	2.54	10	2
1:A:296:ASN:O	1:A:299:HIS:CE1	0.60	2.55	6	1
1:A:300:VAL:O	1:A:300:VAL:HG12	0.60	1.96	17	1
1:A:310:TYR:O	1:A:310:TYR:CD1	0.60	2.54	13	2
1:A:280:LYS:O	1:A:282:LEU:N	0.60	2.35	10	1
1:A:295:ASP:O	1:A:297:ILE:N	0.60	2.35	16	1
1:A:270:PRO:O	1:A:272:GLY:N	0.60	2.35	3	3
1:A:297:ILE:O	1:A:299:HIS:CG	0.60	2.54	13	1
1:A:268:HIS:ND1	1:A:269:GLN:OE1	0.60	2.35	13	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:298:LYS:O	1:A:307:GLN:NE2	0.59	2.35	18	2
1:A:277:ILE:HG22	1:A:277:ILE:O	0.59	1.97	1	2
1:A:295:ASP:O	1:A:296:ASN:ND2	0.59	2.36	13	1
1:A:306:VAL:H	1:A:307:GLN:NE2	0.59	1.95	14	2
1:A:300:VAL:O	1:A:300:VAL:HG13	0.59	1.97	11	1
1:A:283:ASP:O	1:A:286:ASN:ND2	0.59	2.36	9	1
1:A:299:HIS:O	1:A:299:HIS:ND1	0.59	2.36	6	2
1:A:286:ASN:C	1:A:286:ASN:ND2	0.59	2.55	16	1
1:A:299:HIS:ND1	1:A:299:HIS:O	0.59	2.36	14	2
1:A:299:HIS:ND1	1:A:300:VAL:N	0.59	2.50	20	4
1:A:268:HIS:CD2	1:A:268:HIS:O	0.59	2.56	5	3
1:A:290:LYS:O	1:A:292:GLY:N	0.59	2.36	2	2
1:A:293:SER:OG	1:A:294:LYS:N	0.58	2.36	20	1
1:A:276:GLN:O	1:A:278:ILE:N	0.58	2.36	12	4
1:A:284:LEU:O	1:A:285:SER:CB	0.58	2.51	1	7
1:A:279:ASN:ND2	1:A:280:LYS:O	0.58	2.36	10	1
1:A:286:ASN:HD22	1:A:287:VAL:H	0.58	1.41	9	1
1:A:305:SER:OG	1:A:306:VAL:N	0.58	2.35	11	2
1:A:296:ASN:O	1:A:298:LYS:N	0.58	2.36	5	4
1:A:306:VAL:N	1:A:307:GLN:HE22	0.58	1.96	18	2
1:A:309:VAL:O	1:A:311:LYS:N	0.58	2.37	18	3
1:A:278:ILE:O	1:A:279:ASN:CB	0.58	2.51	12	1
1:A:282:LEU:N	1:A:282:LEU:HD22	0.58	2.12	16	1
1:A:284:LEU:H	1:A:284:LEU:CD2	0.58	2.06	6	2
1:A:269:GLN:OE1	1:A:286:ASN:ND2	0.58	2.37	9	1
1:A:280:LYS:O	1:A:283:ASP:N	0.58	2.36	3	3
1:A:300:VAL:O	1:A:303:GLY:N	0.58	2.36	20	1
1:A:291:CYS:CB	1:A:296:ASN:HD22	0.58	2.12	9	1
1:A:281:LYS:H	1:A:281:LYS:CD	0.58	2.12	17	2
1:A:308:ILE:CD1	1:A:308:ILE:H	0.58	2.12	18	3
1:A:284:LEU:N	1:A:284:LEU:CD1	0.58	2.65	4	1
1:A:283:ASP:OD1	1:A:288:GLN:NE2	0.58	2.36	19	1
1:A:297:ILE:O	1:A:299:HIS:CD2	0.58	2.56	13	1
1:A:308:ILE:CD1	1:A:309:VAL:N	0.57	2.63	14	2
1:A:282:LEU:CD2	1:A:282:LEU:N	0.57	2.67	16	1
1:A:293:SER:O	1:A:295:ASP:N	0.57	2.38	2	3
1:A:288:GLN:O	1:A:291:CYS:N	0.57	2.38	6	2
1:A:308:ILE:O	1:A:310:TYR:CD1	0.57	2.57	17	1
1:A:299:HIS:O	1:A:300:VAL:CG2	0.57	2.51	12	1
1:A:280:LYS:C	1:A:282:LEU:H	0.57	2.03	10	1
1:A:274:LYS:O	1:A:275:VAL:HG23	0.57	1.99	18	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:306:VAL:H	1:A:307:GLN:HE22	0.57	1.43	14	1
1:A:297:ILE:HG23	1:A:299:HIS:NE2	0.57	2.15	2	1
1:A:308:ILE:C	1:A:309:VAL:HG12	0.57	2.20	16	1
1:A:292:GLY:O	1:A:293:SER:CB	0.56	2.53	13	3
1:A:309:VAL:O	1:A:309:VAL:HG12	0.56	1.99	11	2
1:A:278:ILE:C	1:A:280:LYS:H	0.56	2.04	3	3
1:A:268:HIS:O	1:A:269:GLN:CB	0.56	2.53	9	2
1:A:286:ASN:HD22	1:A:286:ASN:N	0.56	1.99	9	1
1:A:307:GLN:CD	1:A:307:GLN:N	0.56	2.59	18	2
1:A:290:LYS:NZ	1:A:310:TYR:OH	0.56	2.39	2	1
1:A:293:SER:O	1:A:294:LYS:CB	0.56	2.53	6	1
1:A:297:ILE:H	1:A:297:ILE:CD1	0.56	2.13	20	1
1:A:294:LYS:C	1:A:296:ASN:N	0.56	2.59	19	3
1:A:284:LEU:O	1:A:288:GLN:NE2	0.56	2.37	9	1
1:A:294:LYS:O	1:A:297:ILE:N	0.56	2.38	18	1
1:A:300:VAL:CG2	1:A:307:GLN:N	0.56	2.69	7	2
1:A:297:ILE:O	1:A:307:GLN:NE2	0.56	2.39	10	1
1:A:305:SER:O	1:A:306:VAL:HG13	0.56	2.00	11	2
1:A:307:GLN:O	1:A:307:GLN:CG	0.56	2.53	17	1
1:A:299:HIS:N	1:A:303:GLY:O	0.56	2.39	6	1
1:A:299:HIS:CE1	1:A:301:PRO:O	0.56	2.59	18	2
1:A:276:GLN:C	1:A:278:ILE:H	0.56	2.05	12	2
1:A:299:HIS:CE1	1:A:308:ILE:HD11	0.56	2.36	11	2
1:A:279:ASN:ND2	1:A:282:LEU:H	0.55	2.00	7	1
1:A:306:VAL:HG22	1:A:306:VAL:O	0.55	2.02	18	1
1:A:275:VAL:C	1:A:277:ILE:H	0.55	2.05	5	5
1:A:288:GLN:C	1:A:290:LYS:N	0.55	2.60	10	10
1:A:291:CYS:O	1:A:296:ASN:N	0.55	2.40	8	1
1:A:299:HIS:O	1:A:300:VAL:CB	0.55	2.54	12	1
1:A:290:LYS:NZ	1:A:295:ASP:OD2	0.55	2.36	13	1
1:A:296:ASN:O	1:A:297:ILE:CB	0.55	2.55	10	1
1:A:275:VAL:HG13	1:A:276:GLN:NE2	0.55	2.17	14	1
1:A:301:PRO:C	1:A:303:GLY:N	0.54	2.58	11	2
1:A:293:SER:H	1:A:299:HIS:CD2	0.54	2.20	7	1
1:A:297:ILE:O	1:A:297:ILE:CG2	0.54	2.56	10	1
1:A:296:ASN:C	1:A:298:LYS:H	0.54	2.05	5	3
1:A:276:GLN:NE2	1:A:277:ILE:C	0.54	2.61	15	2
1:A:297:ILE:O	1:A:298:LYS:C	0.54	2.45	10	6
1:A:278:ILE:H	1:A:278:ILE:CD1	0.54	2.15	7	1
1:A:282:LEU:N	1:A:282:LEU:HD12	0.54	2.17	20	3
1:A:278:ILE:O	1:A:280:LYS:N	0.54	2.41	3	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:269:GLN:NE2	1:A:286:ASN:CG	0.54	2.61	5	1
1:A:301:PRO:O	1:A:303:GLY:N	0.54	2.40	11	1
1:A:308:ILE:HD13	1:A:308:ILE:N	0.54	2.18	20	2
1:A:293:SER:C	1:A:295:ASP:H	0.54	2.06	12	3
1:A:293:SER:H	1:A:297:ILE:CG1	0.54	2.15	1	1
1:A:297:ILE:C	1:A:299:HIS:N	0.54	2.60	4	6
1:A:278:ILE:O	1:A:279:ASN:C	0.54	2.46	2	1
1:A:275:VAL:O	1:A:275:VAL:HG22	0.54	2.02	2	1
1:A:285:SER:C	1:A:287:VAL:H	0.54	2.05	16	10
1:A:285:SER:C	1:A:287:VAL:N	0.54	2.61	17	8
1:A:269:GLN:CD	1:A:269:GLN:N	0.54	2.61	6	2
1:A:284:LEU:O	1:A:287:VAL:HG22	0.54	2.03	18	2
1:A:269:GLN:N	1:A:270:PRO:CD	0.53	2.71	1	5
1:A:287:VAL:C	1:A:289:SER:N	0.53	2.61	7	5
1:A:308:ILE:CG2	1:A:311:LYS:H	0.53	2.16	2	1
1:A:285:SER:OG	1:A:286:ASN:N	0.53	2.38	3	2
1:A:309:VAL:C	1:A:311:LYS:H	0.53	2.05	13	4
1:A:306:VAL:O	1:A:306:VAL:HG12	0.53	2.03	9	1
1:A:306:VAL:O	1:A:306:VAL:CG2	0.53	2.56	6	2
1:A:279:ASN:H	1:A:279:ASN:ND2	0.53	2.01	11	1
1:A:278:ILE:HG22	1:A:278:ILE:O	0.53	2.01	9	1
1:A:300:VAL:O	1:A:308:ILE:CG1	0.53	2.56	1	1
1:A:279:ASN:C	1:A:281:LYS:H	0.53	2.06	19	3
1:A:293:SER:O	1:A:294:LYS:C	0.53	2.47	3	1
1:A:310:TYR:O	1:A:310:TYR:CG	0.53	2.62	13	2
1:A:308:ILE:C	1:A:309:VAL:HG23	0.53	2.23	17	2
1:A:310:TYR:O	1:A:311:LYS:C	0.53	2.47	1	9
1:A:293:SER:C	1:A:295:ASP:N	0.53	2.59	19	3
1:A:297:ILE:CD1	1:A:297:ILE:N	0.53	2.72	20	1
1:A:268:HIS:O	1:A:271:GLY:N	0.53	2.40	8	1
1:A:309:VAL:O	1:A:309:VAL:HG13	0.53	2.03	4	2
1:A:286:ASN:HD22	1:A:286:ASN:C	0.53	2.08	18	2
1:A:305:SER:N	1:A:307:GLN:NE2	0.53	2.56	9	1
1:A:310:TYR:O	1:A:311:LYS:O	0.53	2.28	10	4
1:A:295:ASP:C	1:A:297:ILE:N	0.52	2.62	16	1
1:A:296:ASN:C	1:A:298:LYS:N	0.52	2.62	1	4
1:A:308:ILE:CD1	1:A:310:TYR:O	0.52	2.57	12	1
1:A:310:TYR:CD1	1:A:310:TYR:C	0.52	2.82	1	2
1:A:305:SER:C	1:A:306:VAL:HG23	0.52	2.24	12	3
1:A:299:HIS:ND1	1:A:301:PRO:CD	0.52	2.73	16	2
1:A:282:LEU:HD12	1:A:282:LEU:N	0.52	2.19	4	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:288:GLN:O	1:A:289:SER:C	0.52	2.47	6	8
1:A:268:HIS:CE1	1:A:269:GLN:OE1	0.52	2.63	13	1
1:A:297:ILE:O	1:A:299:HIS:CE1	0.52	2.62	11	1
1:A:277:ILE:C	1:A:279:ASN:N	0.52	2.61	12	1
1:A:284:LEU:O	1:A:285:SER:O	0.52	2.28	15	6
1:A:276:GLN:C	1:A:276:GLN:NE2	0.52	2.64	3	2
1:A:305:SER:C	1:A:307:GLN:N	0.52	2.63	2	2
1:A:300:VAL:O	1:A:301:PRO:C	0.52	2.49	11	2
1:A:272:GLY:O	1:A:275:VAL:O	0.52	2.28	1	1
1:A:286:ASN:ND2	1:A:286:ASN:C	0.52	2.63	18	1
1:A:275:VAL:C	1:A:277:ILE:N	0.51	2.63	19	2
1:A:279:ASN:ND2	1:A:279:ASN:H	0.51	2.03	8	1
1:A:296:ASN:O	1:A:297:ILE:C	0.51	2.48	17	2
1:A:277:ILE:HG22	1:A:278:ILE:N	0.51	2.20	6	1
1:A:288:GLN:C	1:A:290:LYS:H	0.51	2.08	18	6
1:A:300:VAL:CG1	1:A:300:VAL:O	0.51	2.57	17	1
1:A:306:VAL:HG12	1:A:309:VAL:HG22	0.51	1.82	20	1
1:A:279:ASN:C	1:A:279:ASN:HD22	0.51	2.08	10	2
1:A:278:ILE:C	1:A:280:LYS:N	0.51	2.63	3	3
1:A:280:LYS:O	1:A:283:ASP:O	0.51	2.29	15	4
1:A:298:LYS:C	1:A:299:HIS:CG	0.51	2.83	15	1
1:A:279:ASN:C	1:A:281:LYS:N	0.51	2.64	2	5
1:A:308:ILE:N	1:A:308:ILE:CD1	0.51	2.73	20	3
1:A:308:ILE:C	1:A:310:TYR:N	0.51	2.64	14	1
1:A:276:GLN:C	1:A:278:ILE:N	0.51	2.63	19	4
1:A:300:VAL:N	1:A:301:PRO:HD2	0.51	2.21	10	6
1:A:306:VAL:O	1:A:307:GLN:O	0.51	2.29	17	1
1:A:305:SER:H	1:A:307:GLN:NE2	0.51	2.02	9	1
1:A:290:LYS:C	1:A:291:CYS:SG	0.50	2.89	4	1
1:A:276:GLN:O	1:A:276:GLN:NE2	0.50	2.43	3	2
1:A:300:VAL:C	1:A:302:GLY:N	0.50	2.65	10	4
1:A:309:VAL:C	1:A:311:LYS:N	0.50	2.65	13	3
1:A:277:ILE:C	1:A:279:ASN:H	0.50	2.10	12	5
1:A:287:VAL:O	1:A:288:GLN:O	0.50	2.28	8	1
1:A:304:GLY:O	1:A:305:SER:O	0.50	2.29	19	2
1:A:293:SER:O	1:A:293:SER:OG	0.50	2.28	8	2
1:A:300:VAL:HG22	1:A:300:VAL:O	0.50	2.04	8	1
1:A:292:GLY:O	1:A:293:SER:O	0.50	2.30	7	1
1:A:280:LYS:H	1:A:280:LYS:CD	0.50	2.19	13	1
1:A:299:HIS:NE2	1:A:308:ILE:HD12	0.50	2.22	20	1
1:A:275:VAL:O	1:A:276:GLN:O	0.50	2.30	14	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:280:LYS:O	1:A:284:LEU:N	0.50	2.44	9	1
1:A:269:GLN:O	1:A:283:ASP:OD2	0.50	2.29	8	2
1:A:276:GLN:OE1	1:A:277:ILE:O	0.50	2.30	3	2
1:A:269:GLN:NE2	1:A:288:GLN:HE22	0.50	2.05	9	1
1:A:268:HIS:O	1:A:268:HIS:CD2	0.50	2.65	8	1
1:A:270:PRO:C	1:A:272:GLY:N	0.50	2.63	3	1
1:A:297:ILE:CG2	1:A:297:ILE:O	0.50	2.59	2	1
1:A:286:ASN:C	1:A:288:GLN:N	0.50	2.64	7	3
1:A:299:HIS:O	1:A:303:GLY:O	0.49	2.31	6	2
1:A:309:VAL:O	1:A:309:VAL:CG2	0.49	2.60	16	1
1:A:288:GLN:O	1:A:295:ASP:OD2	0.49	2.29	12	1
1:A:279:ASN:HD22	1:A:279:ASN:C	0.49	2.09	5	1
1:A:286:ASN:O	1:A:288:GLN:N	0.49	2.46	7	2
1:A:297:ILE:O	1:A:297:ILE:HG22	0.49	2.06	1	1
1:A:298:LYS:O	1:A:299:HIS:O	0.49	2.30	11	1
1:A:308:ILE:H	1:A:308:ILE:HD13	0.49	1.67	9	1
1:A:281:LYS:CD	1:A:281:LYS:N	0.49	2.76	8	3
1:A:308:ILE:CD1	1:A:308:ILE:O	0.49	2.54	7	1
1:A:269:GLN:O	1:A:283:ASP:O	0.49	2.30	11	1
1:A:284:LEU:CD2	1:A:284:LEU:N	0.49	2.68	11	1
1:A:299:HIS:CD2	1:A:308:ILE:HD12	0.49	2.42	20	3
1:A:308:ILE:CD1	1:A:308:ILE:C	0.49	2.80	5	2
1:A:294:LYS:C	1:A:296:ASN:H	0.49	2.11	19	3
1:A:275:VAL:HG22	1:A:275:VAL:O	0.49	2.08	10	1
1:A:267:LYS:O	1:A:268:HIS:O	0.49	2.31	19	2
1:A:287:VAL:C	1:A:289:SER:H	0.49	2.11	13	5
1:A:299:HIS:HB2	1:A:308:ILE:HD11	0.49	1.84	6	1
1:A:282:LEU:O	1:A:283:ASP:OD1	0.49	2.31	11	1
1:A:283:ASP:O	1:A:283:ASP:OD1	0.49	2.31	16	1
1:A:288:GLN:OE1	1:A:296:ASN:OD1	0.48	2.31	5	1
1:A:276:GLN:HE22	1:A:277:ILE:C	0.48	2.12	3	2
1:A:295:ASP:O	1:A:295:ASP:OD1	0.48	2.31	15	1
1:A:268:HIS:C	1:A:270:PRO:CD	0.48	2.81	11	1
1:A:311:LYS:O	1:A:312:PRO:OXT	0.48	2.31	19	3
1:A:297:ILE:C	1:A:299:HIS:H	0.48	2.10	4	4
1:A:268:HIS:CD2	1:A:268:HIS:H	0.48	2.19	19	1
1:A:302:GLY:O	1:A:306:VAL:N	0.48	2.45	1	1
1:A:301:PRO:O	1:A:302:GLY:O	0.48	2.32	12	1
1:A:289:SER:O	1:A:293:SER:OG	0.48	2.29	8	1
1:A:293:SER:O	1:A:294:LYS:O	0.48	2.31	3	1
1:A:310:TYR:C	1:A:310:TYR:CD1	0.48	2.87	12	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:290:LYS:O	1:A:295:ASP:OD2	0.48	2.31	12	1
1:A:269:GLN:H	1:A:269:GLN:CD	0.48	2.10	13	1
1:A:297:ILE:HD12	1:A:297:ILE:O	0.48	2.09	15	1
1:A:287:VAL:O	1:A:287:VAL:HG12	0.48	2.06	3	1
1:A:280:LYS:O	1:A:283:ASP:OD1	0.48	2.31	16	1
1:A:279:ASN:N	1:A:279:ASN:ND2	0.48	2.62	11	1
1:A:285:SER:O	1:A:285:SER:OG	0.48	2.30	14	1
1:A:292:GLY:O	1:A:293:SER:OG	0.47	2.30	13	2
1:A:309:VAL:O	1:A:309:VAL:HG22	0.47	2.10	7	4
1:A:288:GLN:O	1:A:290:LYS:O	0.47	2.32	12	1
1:A:284:LEU:O	1:A:285:SER:OG	0.47	2.30	14	4
1:A:279:ASN:ND2	1:A:279:ASN:C	0.47	2.67	10	2
1:A:269:GLN:OE1	1:A:284:LEU:O	0.47	2.32	9	1
1:A:299:HIS:ND1	1:A:310:TYR:OH	0.47	2.46	10	1
1:A:269:GLN:NE2	1:A:283:ASP:OD2	0.47	2.48	13	1
1:A:290:LYS:O	1:A:291:CYS:C	0.47	2.53	2	3
1:A:286:ASN:ND2	1:A:286:ASN:N	0.47	2.60	5	2
1:A:308:ILE:O	1:A:309:VAL:C	0.47	2.52	3	2
1:A:293:SER:OG	1:A:309:VAL:HG23	0.47	2.08	2	1
1:A:293:SER:O	1:A:297:ILE:CG1	0.47	2.63	11	1
1:A:295:ASP:O	1:A:296:ASN:OD1	0.47	2.33	19	1
1:A:292:GLY:O	1:A:293:SER:C	0.47	2.53	7	1
1:A:279:ASN:HD22	1:A:282:LEU:H	0.47	1.52	7	1
1:A:291:CYS:SG	1:A:296:ASN:ND2	0.47	2.88	9	1
1:A:276:GLN:CD	1:A:276:GLN:H	0.47	2.10	9	1
1:A:297:ILE:O	1:A:299:HIS:O	0.47	2.33	20	1
1:A:299:HIS:CD2	1:A:308:ILE:CD1	0.46	2.99	4	2
1:A:299:HIS:O	1:A:300:VAL:C	0.46	2.54	6	2
1:A:280:LYS:C	1:A:282:LEU:N	0.46	2.68	10	1
1:A:300:VAL:C	1:A:302:GLY:H	0.46	2.13	2	3
1:A:286:ASN:C	1:A:288:GLN:H	0.46	2.12	7	2
1:A:305:SER:C	1:A:307:GLN:H	0.46	2.14	2	1
1:A:301:PRO:C	1:A:303:GLY:H	0.46	2.13	20	2
1:A:295:ASP:O	1:A:296:ASN:CG	0.46	2.53	3	1
1:A:308:ILE:C	1:A:310:TYR:H	0.46	2.14	12	2
1:A:299:HIS:CD2	1:A:300:VAL:N	0.46	2.82	12	1
1:A:279:ASN:ND2	1:A:282:LEU:N	0.46	2.62	7	1
1:A:274:LYS:C	1:A:275:VAL:HG23	0.46	2.30	12	1
1:A:304:GLY:C	1:A:306:VAL:H	0.46	2.13	4	1
1:A:275:VAL:HG13	1:A:276:GLN:N	0.46	2.24	7	2
1:A:287:VAL:O	1:A:288:GLN:C	0.46	2.54	11	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:274:LYS:O	1:A:275:VAL:CG2	0.46	2.63	18	2
1:A:298:LYS:O	1:A:308:ILE:HD13	0.46	2.10	1	1
1:A:283:ASP:OD1	1:A:283:ASP:C	0.46	2.55	15	2
1:A:268:HIS:O	1:A:269:GLN:C	0.46	2.54	11	4
1:A:270:PRO:O	1:A:271:GLY:C	0.46	2.54	5	2
1:A:276:GLN:CA	1:A:276:GLN:HE21	0.46	2.24	14	1
1:A:299:HIS:O	1:A:300:VAL:CG1	0.45	2.55	4	1
1:A:305:SER:O	1:A:306:VAL:CG1	0.45	2.65	19	3
1:A:284:LEU:HD12	1:A:284:LEU:H	0.45	1.65	14	2
1:A:296:ASN:OD1	1:A:299:HIS:CD2	0.45	2.70	6	1
1:A:300:VAL:N	1:A:301:PRO:HD3	0.45	2.26	5	1
1:A:300:VAL:CG1	1:A:305:SER:O	0.45	2.62	3	1
1:A:269:GLN:CD	1:A:283:ASP:OD2	0.45	2.55	13	1
1:A:300:VAL:O	1:A:300:VAL:CG2	0.45	2.63	20	1
1:A:298:LYS:O	1:A:299:HIS:CB	0.45	2.64	15	1
1:A:286:ASN:H	1:A:286:ASN:HD22	0.45	1.51	5	1
1:A:280:LYS:CG	1:A:281:LYS:N	0.45	2.79	17	1
1:A:300:VAL:HG11	1:A:307:GLN:H	0.45	1.70	12	1
1:A:299:HIS:O	1:A:303:GLY:C	0.45	2.55	2	2
1:A:288:GLN:OE1	1:A:296:ASN:CG	0.45	2.54	5	1
1:A:281:LYS:NZ	1:A:282:LEU:HD21	0.45	2.27	5	1
1:A:275:VAL:HG13	1:A:276:GLN:HE21	0.45	1.72	7	1
1:A:309:VAL:O	1:A:310:TYR:CB	0.44	2.64	4	1
1:A:277:ILE:O	1:A:280:LYS:CB	0.44	2.66	4	1
1:A:293:SER:O	1:A:294:LYS:CG	0.44	2.65	9	1
1:A:287:VAL:CG1	1:A:295:ASP:OD2	0.44	2.65	3	1
1:A:284:LEU:CD1	1:A:284:LEU:H	0.44	2.24	14	1
1:A:294:LYS:O	1:A:295:ASP:C	0.44	2.55	18	1
1:A:285:SER:O	1:A:286:ASN:CG	0.44	2.55	16	1
1:A:279:ASN:O	1:A:280:LYS:C	0.44	2.56	10	3
1:A:311:LYS:O	1:A:312:PRO:C	0.44	2.55	9	1
1:A:296:ASN:O	1:A:297:ILE:HG23	0.44	2.12	19	1
1:A:296:ASN:OD1	1:A:296:ASN:C	0.44	2.55	20	1
1:A:288:GLN:CG	1:A:289:SER:N	0.44	2.80	18	1
1:A:293:SER:O	1:A:297:ILE:N	0.44	2.46	2	1
1:A:290:LYS:C	1:A:292:GLY:N	0.44	2.71	2	1
1:A:278:ILE:CD1	1:A:278:ILE:N	0.44	2.71	7	1
1:A:293:SER:OG	1:A:299:HIS:O	0.44	2.35	7	1
1:A:287:VAL:HG22	1:A:287:VAL:O	0.44	2.12	7	1
1:A:308:ILE:C	1:A:308:ILE:CD1	0.44	2.79	10	1
1:A:289:SER:O	1:A:293:SER:CB	0.44	2.66	8	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:269:GLN:CD	1:A:269:GLN:H	0.44	2.16	1	1
1:A:287:VAL:O	1:A:287:VAL:CG1	0.43	2.66	3	1
1:A:286:ASN:O	1:A:286:ASN:ND2	0.43	2.51	12	1
1:A:275:VAL:O	1:A:276:GLN:C	0.43	2.55	7	1
1:A:268:HIS:C	1:A:268:HIS:CD2	0.43	2.90	8	1
1:A:275:VAL:CG2	1:A:276:GLN:N	0.43	2.67	7	1
1:A:286:ASN:HD22	1:A:287:VAL:N	0.43	2.07	9	1
1:A:295:ASP:C	1:A:297:ILE:H	0.43	2.15	5	1
1:A:305:SER:OG	1:A:307:GLN:CG	0.43	2.66	20	1
1:A:297:ILE:CD1	1:A:297:ILE:C	0.43	2.79	15	1
1:A:284:LEU:C	1:A:285:SER:O	0.43	2.56	17	2
1:A:275:VAL:CG1	1:A:276:GLN:N	0.43	2.81	16	1
1:A:299:HIS:ND1	1:A:301:PRO:CG	0.43	2.81	5	1
1:A:300:VAL:HG21	1:A:307:GLN:HB2	0.43	1.90	12	1
1:A:280:LYS:N	1:A:280:LYS:CD	0.43	2.81	13	1
1:A:300:VAL:HG21	1:A:307:GLN:HE21	0.43	1.73	1	1
1:A:299:HIS:N	1:A:304:GLY:HA3	0.43	2.29	10	1
1:A:283:ASP:C	1:A:283:ASP:OD1	0.43	2.56	16	1
1:A:276:GLN:CD	1:A:277:ILE:O	0.43	2.57	3	2
1:A:283:ASP:O	1:A:286:ASN:CG	0.43	2.56	9	1
1:A:308:ILE:O	1:A:310:TYR:CE1	0.43	2.72	17	1
1:A:296:ASN:C	1:A:297:ILE:HG23	0.43	2.34	19	1
1:A:305:SER:O	1:A:308:ILE:O	0.43	2.37	16	1
1:A:279:ASN:O	1:A:282:LEU:N	0.43	2.45	9	1
1:A:296:ASN:C	1:A:296:ASN:OD1	0.43	2.56	2	1
1:A:307:GLN:O	1:A:309:VAL:HG23	0.43	2.14	20	1
1:A:290:LYS:NZ	1:A:297:ILE:CG1	0.42	2.82	4	1
1:A:308:ILE:O	1:A:309:VAL:CB	0.42	2.67	16	1
1:A:300:VAL:HG13	1:A:304:GLY:N	0.42	2.28	9	1
1:A:283:ASP:OD1	1:A:284:LEU:N	0.42	2.52	17	1
1:A:299:HIS:ND1	1:A:301:PRO:HD3	0.42	2.29	16	2
1:A:301:PRO:O	1:A:302:GLY:C	0.42	2.57	11	1
1:A:277:ILE:O	1:A:277:ILE:CG2	0.42	2.66	1	1
1:A:269:GLN:HE22	1:A:288:GLN:HE22	0.42	1.57	9	1
1:A:269:GLN:CB	1:A:270:PRO:CD	0.42	2.97	19	1
1:A:268:HIS:ND1	1:A:288:GLN:NE2	0.42	2.68	14	1
1:A:274:LYS:C	1:A:276:GLN:H	0.42	2.18	4	1
1:A:268:HIS:C	1:A:270:PRO:N	0.42	2.73	11	1
1:A:279:ASN:OD1	1:A:279:ASN:O	0.42	2.38	3	1
1:A:299:HIS:N	1:A:299:HIS:CD2	0.42	2.86	12	2
1:A:298:LYS:CE	1:A:307:GLN:NE2	0.42	2.83	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:305:SER:O	1:A:305:SER:OG	0.42	2.38	4	1
1:A:294:LYS:N	1:A:297:ILE:HG23	0.42	2.30	9	1
1:A:300:VAL:HG22	1:A:306:VAL:HA	0.41	1.92	13	1
1:A:308:ILE:N	1:A:308:ILE:HD12	0.41	2.30	11	2
1:A:293:SER:HB2	1:A:297:ILE:HG22	0.41	1.92	17	1
1:A:303:GLY:O	1:A:304:GLY:O	0.41	2.38	16	1
1:A:308:ILE:C	1:A:309:VAL:CG1	0.41	2.89	16	1
1:A:294:LYS:N	1:A:298:LYS:HB2	0.41	2.30	19	1
1:A:269:GLN:OE1	1:A:287:VAL:CG2	0.41	2.68	9	1
1:A:298:LYS:O	1:A:304:GLY:CA	0.41	2.69	14	1
1:A:281:LYS:N	1:A:281:LYS:CD	0.41	2.84	4	1
1:A:282:LEU:N	1:A:282:LEU:CD1	0.41	2.83	6	2
1:A:282:LEU:H	1:A:282:LEU:HD12	0.41	1.75	3	1
1:A:267:LYS:O	1:A:270:PRO:CD	0.41	2.69	11	1
1:A:274:LYS:C	1:A:276:GLN:N	0.40	2.74	4	1
1:A:299:HIS:CE1	1:A:301:PRO:CD	0.40	3.04	16	1
1:A:284:LEU:N	1:A:288:GLN:OE1	0.40	2.53	16	1
1:A:299:HIS:CE1	1:A:301:PRO:HD3	0.40	2.52	11	1
1:A:281:LYS:CB	1:A:281:LYS:NZ	0.40	2.84	9	1
1:A:310:TYR:O	1:A:311:LYS:CB	0.40	2.69	14	1
1:A:269:GLN:H	1:A:269:GLN:NE2	0.40	2.14	6	1
1:A:297:ILE:O	1:A:297:ILE:HD13	0.40	2.17	11	1
1:A:275:VAL:CG1	1:A:275:VAL:O	0.40	2.66	17	1
1:A:305:SER:O	1:A:306:VAL:CB	0.40	2.69	14	1
1:A:276:GLN:CD	1:A:276:GLN:N	0.40	2.73	20	1
1:A:280:LYS:O	1:A:281:LYS:C	0.40	2.59	3	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	44/46 (96%)	26±2 (58±6%)	8±2 (17±5%)	11±3 (24±7%)	0	1
All	All	880/920 (96%)	513 (58%)	153 (17%)	214 (24%)	0	1

All 38 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	285	SER	18
1	A	289	SER	15
1	A	306	VAL	13
1	A	293	SER	11
1	A	275	VAL	9
1	A	278	ILE	8
1	A	298	LYS	8
1	A	294	LYS	8
1	A	311	LYS	8
1	A	305	SER	7
1	A	301	PRO	7
1	A	286	ASN	7
1	A	288	GLN	7
1	A	280	LYS	6
1	A	309	VAL	6
1	A	300	VAL	6
1	A	297	ILE	6
1	A	277	ILE	6
1	A	279	ASN	5
1	A	310	TYR	5
1	A	296	ASN	5
1	A	295	ASP	5
1	A	303	GLY	4
1	A	271	GLY	4
1	A	276	GLN	4
1	A	269	GLN	4
1	A	302	GLY	3
1	A	299	HIS	3
1	A	292	GLY	3
1	A	268	HIS	2
1	A	270	PRO	2
1	A	307	GLN	2
1	A	272	GLY	2
1	A	287	VAL	1
1	A	304	GLY	1
1	A	291	CYS	1
1	A	308	ILE	1
1	A	281	LYS	1

6.3.2 Protein sidechains [i](#)

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	39/39 (100%)	33±2 (85±6%)	6±2 (15±6%)	8 47
All	All	780/780 (100%)	666 (85%)	114 (15%)	8 47

All 25 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	299	HIS	12
1	A	310	TYR	10
1	A	279	ASN	9
1	A	281	LYS	9
1	A	308	ILE	8
1	A	286	ASN	7
1	A	269	GLN	7
1	A	280	LYS	6
1	A	284	LEU	5
1	A	297	ILE	5
1	A	293	SER	4
1	A	285	SER	4
1	A	307	GLN	4
1	A	311	LYS	4
1	A	300	VAL	3
1	A	276	GLN	3
1	A	298	LYS	3
1	A	289	SER	3
1	A	301	PRO	2
1	A	268	HIS	1
1	A	305	SER	1
1	A	309	VAL	1
1	A	288	GLN	1
1	A	275	VAL	1
1	A	277	ILE	1

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

6.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

The completeness of assignment taking into account all chemical shift lists is 8% for the well-defined parts and 8% for the entire structure.

7.1 Chemical shift list 1

File name: 2mz7_cs.str

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping

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

Total number of shifts	43
Number of shifts mapped to atoms	43
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing

No chemical shift referencing corrections were calculated (not enough data).

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 8%, i.e. 42 atoms were assigned a chemical shift out of a possible 556. 0 out of 7 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	42/224 (19%)	42/89 (47%)	0/92 (0%)	0/43 (0%)
Sidechain	0/308 (0%)	0/183 (0%)	0/110 (0%)	0/15 (0%)
Aromatic	0/24 (0%)	0/12 (0%)	0/8 (0%)	0/4 (0%)
Overall	42/556 (8%)	42/284 (15%)	0/210 (0%)	0/62 (0%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 8%, i.e. 42 atoms were assigned a chemical shift out of a possible 556. 0 out of 7 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	42/224 (19%)	42/89 (47%)	0/92 (0%)	0/43 (0%)
Sidechain	0/308 (0%)	0/183 (0%)	0/110 (0%)	0/15 (0%)
Aromatic	0/24 (0%)	0/12 (0%)	0/8 (0%)	0/4 (0%)
Overall	42/556 (8%)	42/284 (15%)	0/210 (0%)	0/62 (0%)

7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

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

No *random coil index* (RCI) plot could be generated from the current chemical shift list (assigned_chem_shift_list_1). RCI is only applicable to proteins.