



Full wwPDB NMR Structure Validation Report ⓘ

Apr 27, 2016 – 03:55 AM BST

PDB ID : 2N3T
Title : Solution structure of the Rpn1 substrate receptor site toroid 1 (T1)
Authors : Chen, X.; Walters, K.J.
Deposited on : 2015-06-10

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

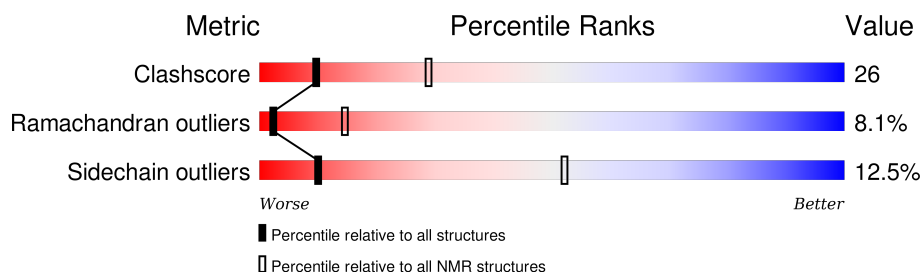
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 61%.

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	131	

2 Ensemble composition and analysis

This entry contains 10 models. Model 3 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

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:490-A:607 (118)	0.84	3

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

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

The models can be grouped into 3 clusters. No single-model clusters were found.

Cluster number	Models
1	1, 6, 7, 9, 10
2	2, 3, 4
3	5, 8

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 1921 atoms, of which 967 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called 26S proteasome regulatory subunit RPN1.

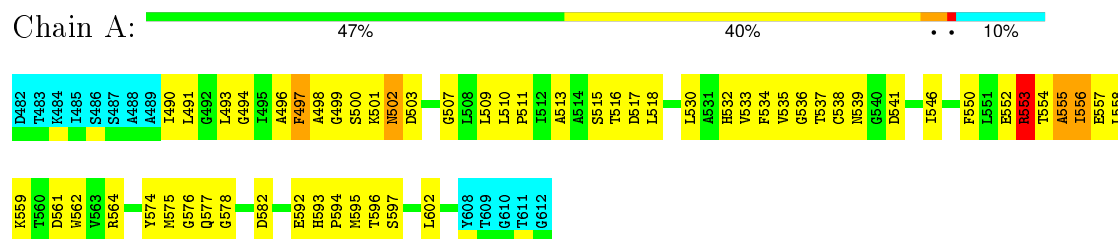
Mol	Chain	Residues	Atoms						Trace
1	A	131	Total	C	H	N	O	S	0
			1921	606	967	150	192	6	

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: 26S proteasome regulatory subunit RPN1

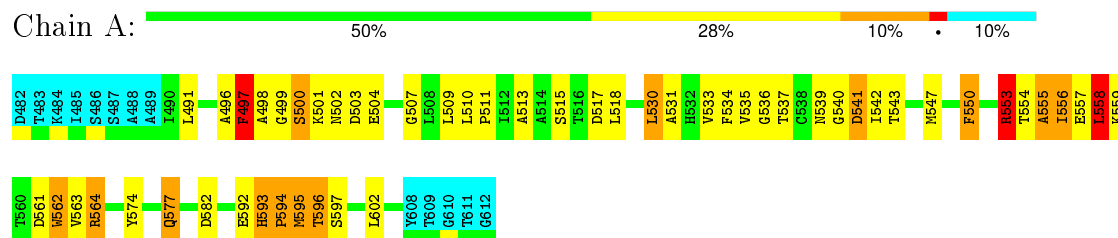


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

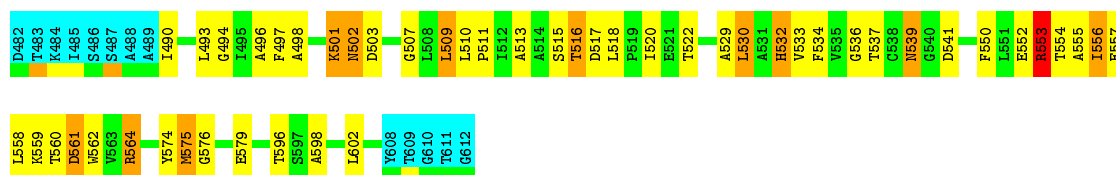
- Molecule 1: 26S proteasome regulatory subunit RPN1



4.2.2 Score per residue for model 2

- Molecule 1: 26S proteasome regulatory subunit RPN1

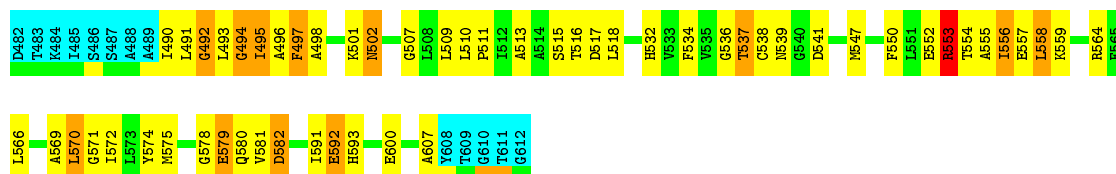




4.2.3 Score per residue for model 3 (medoid)

- Molecule 1: 26S proteasome regulatory subunit RPN1

Chain A: 48% 32% 9% 10%



4.2.4 Score per residue for model 4

- Molecule 1: 26S proteasome regulatory subunit RPN1

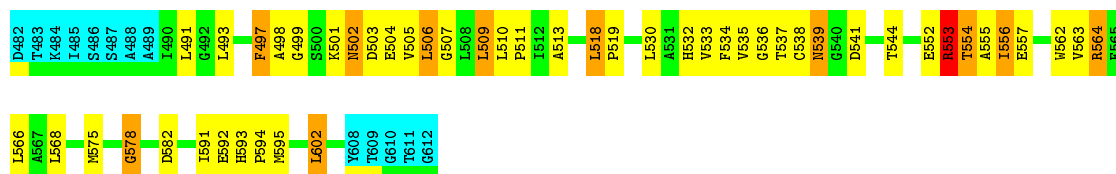
Chain A: 53% 28% 8% 10%



4.2.5 Score per residue for model 5

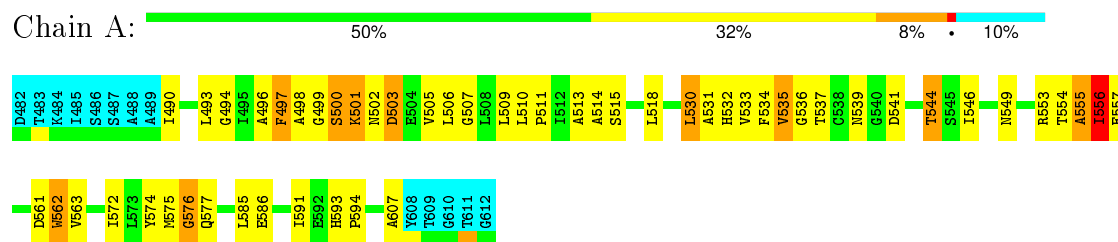
- Molecule 1: 26S proteasome regulatory subunit RPN1

Chain A: 53% 28% 8% 10%



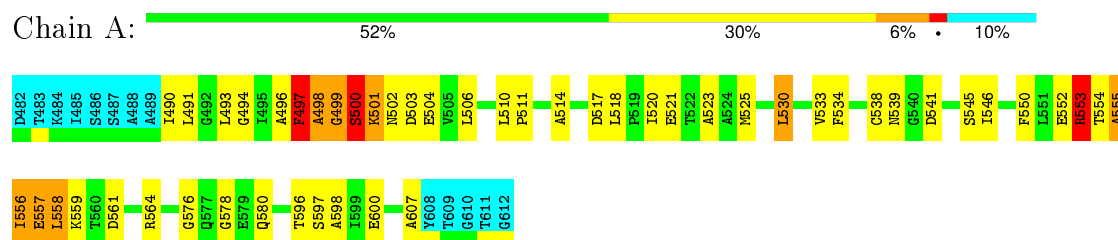
4.2.6 Score per residue for model 6

- Molecule 1: 26S proteasome regulatory subunit RPN1



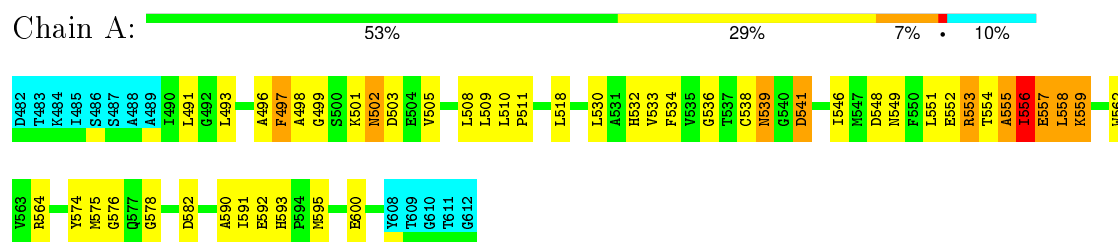
4.2.7 Score per residue for model 7

- Molecule 1: 26S proteasome regulatory subunit RPN1



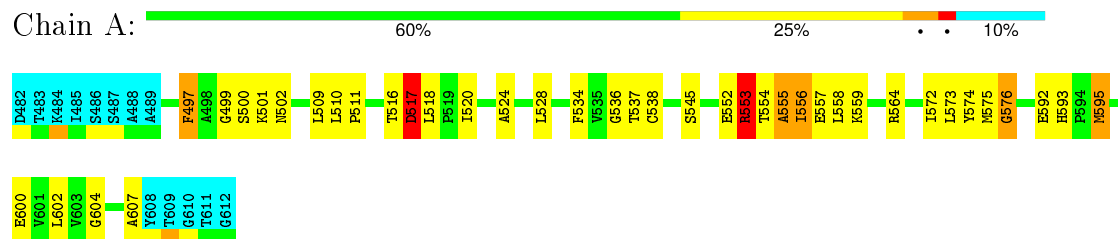
4.2.8 Score per residue for model 8

- Molecule 1: 26S proteasome regulatory subunit RPN1



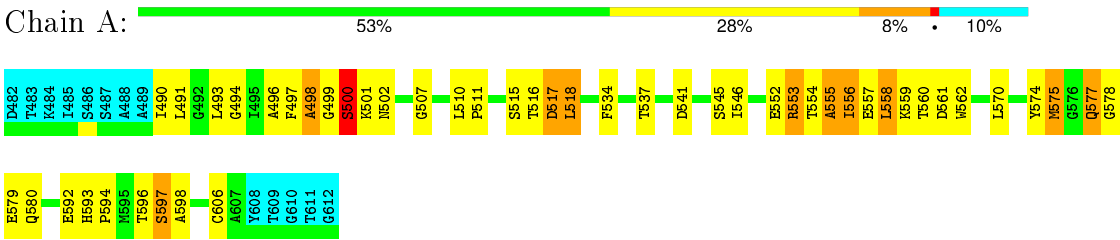
4.2.9 Score per residue for model 9

- Molecule 1: 26S proteasome regulatory subunit RPN1



4.2.10 Score per residue for model 10

- Molecule 1: 26S proteasome regulatory subunit RPN1



5 Refinement protocol and experimental data overview

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

Of the 50 calculated structures, 10 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	structure solution	
X-PLOR NIH	refinement	

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

Chemical shift file(s)	2n3t_cs.str
Number of chemical shift lists	1
Total number of shifts	1034
Number of shifts mapped to atoms	1034
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	61%

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.

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

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	1.6±0.8
All	All	0	16

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
1	A	564	ARG	Sidechain	8
1	A	553	ARG	Sidechain	8

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	865	881	881	45±11
All	All	8650	8810	8810	450

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 26.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:497:PHE:CD1	1:A:498:ALA:N	0.71	2.58	2	1
1:A:494:GLY:O	1:A:497:PHE:CD1	0.71	2.44	3	1
1:A:494:GLY:C	1:A:497:PHE:CZ	0.69	2.66	3	1
1:A:494:GLY:O	1:A:497:PHE:CG	0.68	2.47	3	2
1:A:494:GLY:O	1:A:497:PHE:CE1	0.67	2.46	3	1
1:A:499:GLY:O	1:A:501:LYS:N	0.67	2.27	7	5
1:A:577:GLN:NE2	1:A:577:GLN:H	0.65	1.89	10	1
1:A:494:GLY:O	1:A:497:PHE:CD2	0.64	2.51	4	3
1:A:506:LEU:HD23	1:A:534:PHE:CE1	0.63	2.28	5	1
1:A:497:PHE:CZ	1:A:533:VAL:HG22	0.62	2.29	2	1
1:A:501:LYS:O	1:A:502:ASN:ND2	0.62	2.33	3	1
1:A:591:ILE:N	1:A:591:ILE:HD12	0.61	2.11	3	1
1:A:494:GLY:O	1:A:497:PHE:CZ	0.61	2.53	3	1
1:A:497:PHE:CD1	1:A:497:PHE:N	0.61	2.65	5	1
1:A:572:ILE:N	1:A:572:ILE:HD12	0.61	2.10	6	1
1:A:497:PHE:O	1:A:499:GLY:N	0.59	2.35	7	4
1:A:525:MET:N	1:A:525:MET:SD	0.59	2.76	7	1
1:A:494:GLY:O	1:A:496:ALA:N	0.59	2.36	4	2
1:A:497:PHE:N	1:A:497:PHE:CD1	0.59	2.68	8	1
1:A:494:GLY:HA2	1:A:497:PHE:CE2	0.59	2.32	3	1
1:A:494:GLY:O	1:A:497:PHE:CE2	0.58	2.57	3	3
1:A:508:LEU:HD12	1:A:508:LEU:N	0.58	2.14	8	1
1:A:497:PHE:CD1	1:A:505:VAL:HG21	0.57	2.33	8	1
1:A:554:THR:O	1:A:556:ILE:N	0.57	2.38	7	9
1:A:502:ASN:OD1	1:A:502:ASN:N	0.56	2.38	8	1
1:A:497:PHE:CE1	1:A:533:VAL:HG22	0.56	2.35	2	1
1:A:593:HIS:O	1:A:594:PRO:O	0.56	2.23	1	1
1:A:490:ILE:N	1:A:490:ILE:HD12	0.56	2.16	3	1
1:A:591:ILE:N	1:A:591:ILE:CD1	0.55	2.69	3	1
1:A:572:ILE:HG23	1:A:573:LEU:N	0.55	2.17	9	1
1:A:501:LYS:O	1:A:502:ASN:CB	0.55	2.55	3	3
1:A:493:LEU:O	1:A:497:PHE:CE1	0.55	2.60	8	2
1:A:562:TRP:CD2	1:A:563:VAL:N	0.55	2.75	1	2
1:A:498:ALA:C	1:A:500:SER:H	0.55	2.05	7	1
1:A:574:TYR:O	1:A:576:GLY:N	0.54	2.40	6	5
1:A:595:MET:N	1:A:595:MET:SD	0.54	2.80	8	1
1:A:579:GLU:H	1:A:579:GLU:CD	0.54	2.06	3	2
1:A:508:LEU:N	1:A:508:LEU:CD1	0.54	2.71	8	1
1:A:557:GLU:O	1:A:559:LYS:N	0.54	2.40	7	5
1:A:496:ALA:C	1:A:498:ALA:H	0.53	2.06	7	6
1:A:493:LEU:O	1:A:497:PHE:CD1	0.53	2.61	6	2
1:A:497:PHE:CD1	1:A:533:VAL:HG21	0.53	2.38	4	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:502:ASN:N	1:A:502:ASN:OD1	0.53	2.37	5	2
1:A:558:LEU:N	1:A:558:LEU:HD23	0.53	2.17	3	2
1:A:501:LYS:O	1:A:502:ASN:CG	0.53	2.47	4	2
1:A:516:THR:O	1:A:518:LEU:N	0.53	2.42	10	1
1:A:497:PHE:C	1:A:499:GLY:N	0.53	2.62	7	5
1:A:494:GLY:C	1:A:496:ALA:N	0.53	2.62	4	3
1:A:502:ASN:OD1	1:A:534:PHE:CZ	0.53	2.61	3	1
1:A:530:LEU:O	1:A:533:VAL:N	0.53	2.42	8	1
1:A:539:ASN:ND2	1:A:541:ASP:OD2	0.53	2.41	3	1
1:A:502:ASN:HD22	1:A:533:VAL:CG1	0.53	2.17	5	1
1:A:554:THR:C	1:A:556:ILE:N	0.52	2.62	7	10
1:A:498:ALA:O	1:A:500:SER:N	0.52	2.43	7	1
1:A:497:PHE:CD1	1:A:533:VAL:CG2	0.52	2.93	4	1
1:A:499:GLY:O	1:A:502:ASN:ND2	0.52	2.43	10	1
1:A:491:LEU:O	1:A:494:GLY:N	0.52	2.42	7	2
1:A:574:TYR:N	1:A:574:TYR:CD1	0.52	2.78	3	1
1:A:551:LEU:N	1:A:551:LEU:HD12	0.52	2.19	8	1
1:A:569:ALA:O	1:A:571:GLY:N	0.52	2.42	3	1
1:A:593:HIS:CG	1:A:593:HIS:O	0.52	2.62	5	1
1:A:497:PHE:C	1:A:499:GLY:H	0.52	2.08	8	3
1:A:499:GLY:C	1:A:501:LYS:N	0.51	2.62	7	5
1:A:507:GLY:O	1:A:511:PRO:CD	0.51	2.57	10	7
1:A:554:THR:C	1:A:556:ILE:H	0.51	2.08	5	7
1:A:593:HIS:ND1	1:A:597:SER:OG	0.51	2.43	10	1
1:A:557:GLU:C	1:A:559:LYS:H	0.51	2.09	1	6
1:A:538:CYS:O	1:A:539:ASN:ND2	0.51	2.41	5	1
1:A:502:ASN:O	1:A:504:GLU:N	0.51	2.43	1	2
1:A:550:PHE:O	1:A:553:ARG:NE	0.51	2.43	1	1
1:A:534:PHE:CD2	1:A:537:THR:OG1	0.51	2.64	2	2
1:A:498:ALA:C	1:A:500:SER:N	0.51	2.63	7	1
1:A:520:ILE:O	1:A:524:ALA:N	0.51	2.44	9	1
1:A:497:PHE:CG	1:A:498:ALA:N	0.51	2.78	2	1
1:A:501:LYS:O	1:A:502:ASN:OD1	0.51	2.29	4	1
1:A:520:ILE:HD11	1:A:556:ILE:CG2	0.51	2.35	2	1
1:A:537:THR:OG1	1:A:538:CYS:N	0.50	2.44	3	2
1:A:578:GLY:O	1:A:582:ASP:N	0.50	2.44	8	1
1:A:494:GLY:C	1:A:497:PHE:CE2	0.50	2.85	3	2
1:A:541:ASP:OD1	1:A:541:ASP:N	0.50	2.43	1	1
1:A:560:THR:O	1:A:561:ASP:CB	0.50	2.59	10	1
1:A:494:GLY:C	1:A:496:ALA:H	0.50	2.10	4	3
1:A:510:LEU:CB	1:A:511:PRO:CD	0.50	2.90	2	10

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:507:GLY:O	1:A:511:PRO:CG	0.50	2.60	3	5
1:A:515:SER:O	1:A:517:ASP:N	0.50	2.44	2	3
1:A:503:ASP:CG	1:A:504:GLU:N	0.50	2.66	5	1
1:A:510:LEU:O	1:A:514:ALA:N	0.50	2.43	7	1
1:A:494:GLY:CA	1:A:497:PHE:CE2	0.49	2.94	3	1
1:A:572:ILE:N	1:A:572:ILE:CD1	0.49	2.75	6	1
1:A:559:LYS:O	1:A:562:TRP:CD1	0.49	2.65	10	1
1:A:557:GLU:C	1:A:559:LYS:N	0.49	2.66	7	7
1:A:541:ASP:O	1:A:544:THR:N	0.49	2.44	6	1
1:A:502:ASN:C	1:A:504:GLU:N	0.49	2.65	1	2
1:A:492:GLY:O	1:A:494:GLY:N	0.49	2.46	3	2
1:A:569:ALA:C	1:A:571:GLY:N	0.49	2.66	3	1
1:A:502:ASN:OD1	1:A:502:ASN:C	0.49	2.51	3	1
1:A:558:LEU:CD2	1:A:558:LEU:N	0.49	2.75	3	2
1:A:571:GLY:O	1:A:575:MET:N	0.49	2.46	3	1
1:A:578:GLY:O	1:A:580:GLN:N	0.49	2.46	3	2
1:A:493:LEU:O	1:A:497:PHE:CD2	0.49	2.65	10	1
1:A:497:PHE:O	1:A:502:ASN:OD1	0.49	2.31	5	3
1:A:530:LEU:O	1:A:530:LEU:HD23	0.49	2.08	2	1
1:A:585:LEU:HD12	1:A:585:LEU:N	0.49	2.23	4	1
1:A:530:LEU:HD23	1:A:530:LEU:O	0.49	2.08	7	2
1:A:500:SER:HB2	1:A:533:VAL:HG11	0.49	1.83	1	1
1:A:515:SER:C	1:A:517:ASP:N	0.48	2.66	2	4
1:A:606:CYS:O	1:A:606:CYS:SG	0.48	2.71	10	1
1:A:568:LEU:CD2	1:A:568:LEU:N	0.48	2.76	5	1
1:A:579:GLU:O	1:A:582:ASP:OD2	0.48	2.31	3	1
1:A:552:GLU:O	1:A:553:ARG:C	0.48	2.51	5	5
1:A:592:GLU:O	1:A:593:HIS:ND1	0.48	2.45	3	1
1:A:534:PHE:C	1:A:536:GLY:N	0.48	2.67	2	8
1:A:541:ASP:N	1:A:541:ASP:OD1	0.48	2.42	8	1
1:A:509:LEU:HD23	1:A:530:LEU:HD23	0.48	1.84	6	1
1:A:594:PRO:C	1:A:596:THR:H	0.48	2.11	1	1
1:A:497:PHE:CD1	1:A:533:VAL:HG13	0.48	2.43	2	1
1:A:521:GLU:O	1:A:525:MET:SD	0.48	2.72	7	1
1:A:560:THR:OG1	1:A:561:ASP:N	0.48	2.46	10	1
1:A:502:ASN:C	1:A:504:GLU:H	0.48	2.12	7	2
1:A:561:ASP:OD1	1:A:562:TRP:N	0.48	2.47	2	1
1:A:496:ALA:C	1:A:498:ALA:N	0.48	2.67	1	6
1:A:550:PHE:CE1	1:A:566:LEU:HD21	0.48	2.44	3	1
1:A:496:ALA:O	1:A:498:ALA:N	0.48	2.47	6	4
1:A:492:GLY:C	1:A:494:GLY:N	0.47	2.66	3	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:555:ALA:C	1:A:557:GLU:H	0.47	2.12	8	5
1:A:575:MET:SD	1:A:575:MET:O	0.47	2.72	10	1
1:A:531:ALA:O	1:A:535:VAL:N	0.47	2.48	1	2
1:A:506:LEU:O	1:A:510:LEU:HD13	0.47	2.09	7	1
1:A:509:LEU:O	1:A:513:ALA:N	0.47	2.48	6	6
1:A:555:ALA:C	1:A:557:GLU:N	0.47	2.67	8	8
1:A:552:GLU:O	1:A:554:THR:N	0.47	2.47	8	1
1:A:520:ILE:O	1:A:523:ALA:N	0.47	2.48	7	1
1:A:520:ILE:HD12	1:A:556:ILE:CG2	0.47	2.39	9	1
1:A:595:MET:O	1:A:595:MET:SD	0.47	2.73	9	1
1:A:502:ASN:C	1:A:502:ASN:OD1	0.47	2.53	4	1
1:A:509:LEU:CD2	1:A:530:LEU:HD23	0.47	2.40	6	1
1:A:534:PHE:O	1:A:537:THR:O	0.47	2.33	3	8
1:A:504:GLU:CD	1:A:504:GLU:N	0.47	2.68	4	1
1:A:575:MET:C	1:A:575:MET:SD	0.47	2.93	10	2
1:A:502:ASN:ND2	1:A:534:PHE:CD1	0.47	2.83	4	1
1:A:503:ASP:OD1	1:A:504:GLU:N	0.47	2.47	5	1
1:A:568:LEU:N	1:A:568:LEU:HD22	0.47	2.25	5	1
1:A:490:ILE:O	1:A:494:GLY:N	0.46	2.49	3	5
1:A:516:THR:C	1:A:518:LEU:N	0.46	2.67	10	1
1:A:562:TRP:CE3	1:A:563:VAL:N	0.46	2.82	6	1
1:A:594:PRO:C	1:A:596:THR:N	0.46	2.68	1	1
1:A:578:GLY:C	1:A:580:GLN:N	0.46	2.68	3	3
1:A:577:GLN:CD	1:A:577:GLN:H	0.46	2.11	10	1
1:A:593:HIS:O	1:A:593:HIS:CG	0.46	2.68	6	1
1:A:581:VAL:CG1	1:A:582:ASP:N	0.46	2.78	3	1
1:A:554:THR:HG22	1:A:555:ALA:N	0.46	2.25	3	1
1:A:558:LEU:N	1:A:558:LEU:CD2	0.46	2.78	10	1
1:A:538:CYS:O	1:A:539:ASN:OD1	0.46	2.34	8	1
1:A:579:GLU:N	1:A:579:GLU:OE1	0.46	2.49	3	1
1:A:499:GLY:C	1:A:501:LYS:H	0.46	2.14	7	5
1:A:595:MET:SD	1:A:595:MET:C	0.46	2.94	9	1
1:A:532:HIS:O	1:A:532:HIS:ND1	0.46	2.49	6	1
1:A:574:TYR:O	1:A:575:MET:C	0.46	2.54	2	1
1:A:532:HIS:O	1:A:535:VAL:HG12	0.46	2.11	4	1
1:A:593:HIS:N	1:A:593:HIS:CD2	0.45	2.83	10	1
1:A:503:ASP:O	1:A:506:LEU:CB	0.45	2.64	5	1
1:A:502:ASN:ND2	1:A:534:PHE:CE1	0.45	2.83	4	1
1:A:534:PHE:O	1:A:536:GLY:N	0.45	2.49	2	7
1:A:569:ALA:O	1:A:572:ILE:N	0.45	2.49	3	1
1:A:555:ALA:O	1:A:557:GLU:N	0.45	2.49	8	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:585:LEU:N	1:A:585:LEU:HD12	0.45	2.26	6	1
1:A:574:TYR:CD1	1:A:574:TYR:N	0.45	2.85	1	1
1:A:596:THR:O	1:A:598:ALA:N	0.45	2.50	10	4
1:A:602:LEU:HD23	1:A:602:LEU:C	0.45	2.32	1	1
1:A:502:ASN:OD1	1:A:502:ASN:O	0.45	2.35	3	1
1:A:575:MET:CG	1:A:575:MET:O	0.45	2.64	3	1
1:A:572:ILE:CG2	1:A:573:LEU:N	0.45	2.80	9	1
1:A:586:GLU:OE2	1:A:586:GLU:O	0.45	2.35	6	1
1:A:556:ILE:HG23	1:A:556:ILE:O	0.45	2.12	3	1
1:A:552:GLU:O	1:A:553:ARG:O	0.45	2.35	2	3
1:A:500:SER:HB3	1:A:533:VAL:HG11	0.45	1.88	6	1
1:A:490:ILE:N	1:A:490:ILE:CD1	0.45	2.80	3	1
1:A:520:ILE:HD11	1:A:556:ILE:HG21	0.45	1.89	2	1
1:A:497:PHE:O	1:A:500:SER:OG	0.45	2.35	1	1
1:A:570:LEU:O	1:A:574:TYR:CE2	0.44	2.71	10	1
1:A:551:LEU:CD1	1:A:551:LEU:N	0.44	2.80	8	1
1:A:582:ASP:OD1	1:A:582:ASP:C	0.44	2.55	3	1
1:A:558:LEU:HD23	1:A:558:LEU:N	0.44	2.26	1	3
1:A:561:ASP:N	1:A:561:ASP:OD1	0.44	2.47	2	1
1:A:593:HIS:O	1:A:595:MET:N	0.44	2.50	9	1
1:A:506:LEU:O	1:A:506:LEU:HD13	0.44	2.13	5	1
1:A:491:LEU:HD12	1:A:491:LEU:N	0.44	2.28	8	1
1:A:493:LEU:O	1:A:496:ALA:HB3	0.44	2.13	8	1
1:A:550:PHE:CE2	1:A:566:LEU:HD21	0.44	2.48	4	1
1:A:497:PHE:C	1:A:497:PHE:CD1	0.44	2.87	7	1
1:A:592:GLU:CD	1:A:592:GLU:H	0.44	2.15	10	1
1:A:497:PHE:O	1:A:500:SER:N	0.44	2.50	10	1
1:A:500:SER:OG	1:A:505:VAL:HG11	0.44	2.13	6	1
1:A:561:ASP:O	1:A:561:ASP:OD1	0.44	2.36	7	1
1:A:596:THR:C	1:A:598:ALA:N	0.43	2.71	7	4
1:A:518:LEU:CB	1:A:519:PRO:CD	0.43	2.96	5	1
1:A:504:GLU:H	1:A:504:GLU:CD	0.43	2.16	4	1
1:A:595:MET:O	1:A:597:SER:N	0.43	2.51	1	2
1:A:516:THR:O	1:A:517:ASP:OD1	0.43	2.37	9	1
1:A:539:ASN:ND2	1:A:541:ASP:CB	0.43	2.82	2	1
1:A:503:ASP:OD1	1:A:503:ASP:N	0.43	2.51	6	2
1:A:592:GLU:OE2	1:A:600:GLU:OE1	0.43	2.36	8	1
1:A:594:PRO:O	1:A:596:THR:N	0.43	2.51	1	1
1:A:550:PHE:CE1	1:A:570:LEU:HD11	0.43	2.49	3	1
1:A:592:GLU:O	1:A:592:GLU:OE2	0.43	2.37	9	1
1:A:513:ALA:O	1:A:549:ASN:ND2	0.43	2.49	6	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:556:ILE:O	1:A:556:ILE:CG2	0.43	2.66	10	1
1:A:515:SER:C	1:A:517:ASP:H	0.43	2.17	3	3
1:A:590:ALA:HB3	1:A:591:ILE:HD12	0.43	1.91	8	1
1:A:581:VAL:HG13	1:A:582:ASP:N	0.43	2.29	3	1
1:A:490:ILE:O	1:A:494:GLY:CA	0.43	2.67	10	1
1:A:514:ALA:O	1:A:549:ASN:OD1	0.43	2.37	6	1
1:A:506:LEU:HD13	1:A:534:PHE:CZ	0.43	2.49	6	1
1:A:585:LEU:CD1	1:A:585:LEU:N	0.43	2.82	4	2
1:A:538:CYS:SG	1:A:539:ASN:N	0.43	2.92	7	1
1:A:600:GLU:OE2	1:A:600:GLU:O	0.42	2.37	3	1
1:A:600:GLU:O	1:A:604:GLY:N	0.42	2.52	9	1
1:A:532:HIS:O	1:A:535:VAL:CG1	0.42	2.67	5	1
1:A:539:ASN:OD1	1:A:540:GLY:N	0.42	2.52	1	1
1:A:563:VAL:HG13	1:A:564:ARG:N	0.42	2.29	1	2
1:A:501:LYS:C	1:A:502:ASN:OD1	0.42	2.58	2	1
1:A:534:PHE:C	1:A:536:GLY:H	0.42	2.18	2	4
1:A:592:GLU:CD	1:A:592:GLU:N	0.42	2.72	10	1
1:A:541:ASP:O	1:A:544:THR:OG1	0.42	2.32	5	1
1:A:500:SER:C	1:A:502:ASN:H	0.42	2.18	7	3
1:A:491:LEU:HD23	1:A:491:LEU:C	0.42	2.35	1	1
1:A:595:MET:C	1:A:597:SER:N	0.42	2.73	4	2
1:A:591:ILE:HG22	1:A:592:GLU:N	0.42	2.30	5	1
1:A:592:GLU:OE2	1:A:593:HIS:O	0.42	2.37	1	1
1:A:582:ASP:OD1	1:A:582:ASP:O	0.42	2.38	1	1
1:A:529:ALA:O	1:A:532:HIS:N	0.41	2.53	2	1
1:A:528:LEU:HD23	1:A:528:LEU:O	0.41	2.16	9	1
1:A:562:TRP:CD1	1:A:563:VAL:N	0.41	2.88	5	1
1:A:602:LEU:CD1	1:A:602:LEU:C	0.41	2.89	5	1
1:A:543:THR:O	1:A:547:MET:N	0.41	2.53	1	1
1:A:499:GLY:O	1:A:502:ASN:OD1	0.41	2.38	6	1
1:A:502:ASN:ND2	1:A:533:VAL:CG1	0.41	2.84	5	1
1:A:495:ILE:O	1:A:498:ALA:HB3	0.41	2.16	3	1
1:A:594:PRO:O	1:A:597:SER:CB	0.41	2.68	4	1
1:A:501:LYS:CD	1:A:501:LYS:N	0.41	2.84	7	1
1:A:531:ALA:O	1:A:533:VAL:N	0.40	2.54	1	1
1:A:539:ASN:OD1	1:A:541:ASP:HB2	0.40	2.15	6	1
1:A:578:GLY:O	1:A:582:ASP:OD2	0.40	2.39	5	1
1:A:491:LEU:CD1	1:A:491:LEU:N	0.40	2.85	8	1
1:A:539:ASN:OD1	1:A:541:ASP:N	0.40	2.55	7	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	118/131 (90%)	91±4 (77±4%)	18±4 (15±4%)	10±3 (8±3%)	2	14
All	All	1180/1310 (90%)	910 (77%)	175 (15%)	95 (8%)	2	14

All 35 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	553	ARG	10
1	A	555	ALA	7
1	A	558	LEU	5
1	A	500	SER	5
1	A	594	PRO	5
1	A	575	MET	5
1	A	517	ASP	4
1	A	498	ALA	4
1	A	607	ALA	4
1	A	516	THR	3
1	A	576	GLY	3
1	A	497	PHE	3
1	A	597	SER	3
1	A	561	ASP	3
1	A	503	ASP	2
1	A	579	GLU	2
1	A	556	ILE	2
1	A	494	GLY	2
1	A	559	LYS	2
1	A	577	GLN	2
1	A	492	GLY	2
1	A	515	SER	2
1	A	495	ILE	2
1	A	493	LEU	2
1	A	591	ILE	1
1	A	539	ASN	1
1	A	578	GLY	1

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Mol	Chain	Res	Type	Models (Total)
1	A	595	MET	1
1	A	491	LEU	1
1	A	499	GLY	1
1	A	535	VAL	1
1	A	570	LEU	1
1	A	596	THR	1
1	A	502	ASN	1
1	A	592	GLU	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	93/102 (91%)	81±3 (88±4%)	12±3 (12±4%)	10	52
All	All	930/1020 (91%)	814 (88%)	116 (12%)	10	52

All 44 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	556	ILE	10
1	A	518	LEU	10
1	A	497	PHE	7
1	A	530	LEU	6
1	A	501	LYS	6
1	A	546	ILE	5
1	A	502	ASN	5
1	A	558	LEU	4
1	A	509	LEU	4
1	A	550	PHE	4
1	A	503	ASP	3
1	A	541	ASP	3
1	A	532	HIS	3
1	A	562	TRP	3
1	A	602	LEU	3
1	A	557	GLU	3
1	A	545	SER	3

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Mol	Chain	Res	Type	Models (Total)
1	A	539	ASN	2
1	A	534	PHE	2
1	A	505	VAL	2
1	A	595	MET	2
1	A	593	HIS	2
1	A	577	GLN	2
1	A	500	SER	2
1	A	548	ASP	1
1	A	566	LEU	1
1	A	564	ARG	1
1	A	600	GLU	1
1	A	517	ASP	1
1	A	560	THR	1
1	A	537	THR	1
1	A	549	ASN	1
1	A	516	THR	1
1	A	491	LEU	1
1	A	542	ILE	1
1	A	506	LEU	1
1	A	533	VAL	1
1	A	575	MET	1
1	A	522	THR	1
1	A	493	LEU	1
1	A	582	ASP	1
1	A	544	THR	1
1	A	547	MET	1
1	A	554	THR	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 carbohydrates 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 [i](#)

The completeness of assignment taking into account all chemical shift lists is 61% for the well-defined parts and 58% for the entire structure.

7.1 Chemical shift list 1

File name: 2n3t_cs.str

Chemical shift list name: *assigned_chem_shift_list_1*

7.1.1 Bookkeeping [i](#)

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

Total number of shifts	1034
Number of shifts mapped to atoms	1034
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	96	-0.46 ± 0.19	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	92	0.22 ± 0.07	None needed (< 0.5 ppm)
$^{13}\text{C}'$	0	—	—
^{15}N	63	0.80 ± 0.70	None needed (imprecise)

7.1.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 61%, i.e. 807 atoms were assigned a chemical shift out of a possible 1317. 21 out of 25 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	313/584 (54%)	157/233 (67%)	93/236 (39%)	63/115 (55%)
Sidechain	474/663 (71%)	280/377 (74%)	194/273 (71%)	0/13 (0%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	20/70 (29%)	20/38 (53%)	0/29 (0%)	0/3 (0%)
Overall	807/1317 (61%)	457/648 (71%)	287/538 (53%)	63/131 (48%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 58%, i.e. 833 atoms were assigned a chemical shift out of a possible 1441. 21 out of 25 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	319/649 (49%)	160/259 (62%)	96/262 (37%)	63/128 (49%)
Sidechain	490/714 (69%)	290/406 (71%)	200/294 (68%)	0/14 (0%)
Aromatic	24/78 (31%)	24/42 (57%)	0/33 (0%)	0/3 (0%)
Overall	833/1441 (58%)	474/707 (67%)	296/589 (50%)	63/145 (43%)

7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

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

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition.

Random coil index (RCI) for chain A:

