



Full wwPDB NMR Structure Validation Report ⓘ

May 28, 2020 – 10:08 pm BST

PDB ID : 2JGW
Title : Structure of CCP module 7 of complement factor H - The AMD at risk variant (402H)
Authors : Herbert, A.P.; Deakin, J.A.; Schmidt, C.Q.; Blaum, B.S.; Egan, C.; Ferreira, V.P.; Pangburn, M.K.; Lyon, M.; Uhrin, D.; Barlow, P.N.
Deposited on : 2007-02-16

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

Cyrange : Kirchner and Güntert (2011)
NmrClust : Kelley et al. (1996)
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
ShiftChecker : 2.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

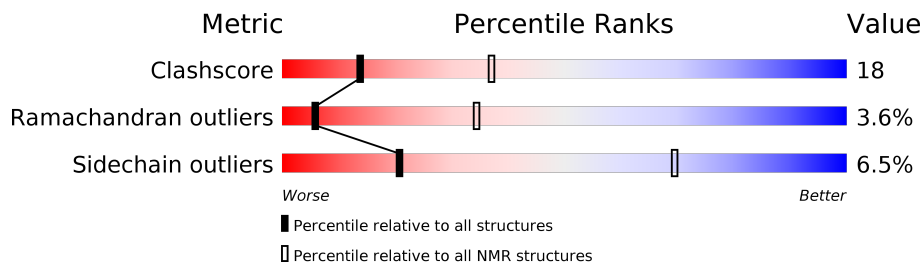
1 Overall quality at a glance

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	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 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	61	

2 Ensemble composition and analysis i

This entry contains 20 models. Model 6 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative.

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:387-A:446 (60)	0.46	6

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 2 clusters and 16 single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 969 atoms, of which 480 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called COMPLEMENT FACTOR H.

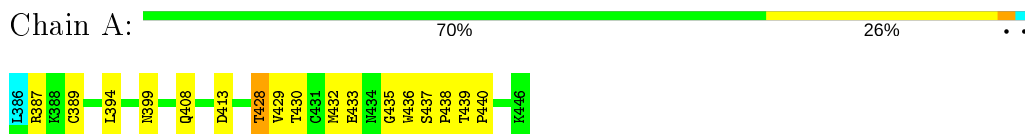
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	61	969	309	480	90	85	5	0

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: COMPLEMENT FACTOR H

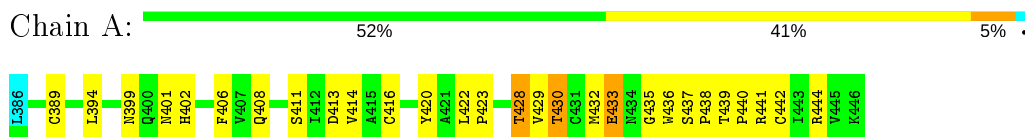


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

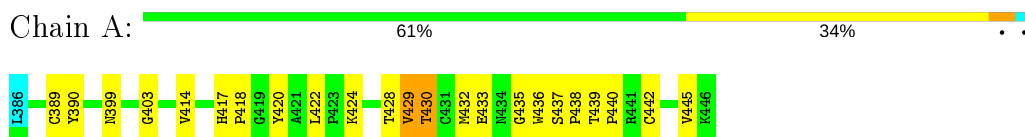
4.2.1 Score per residue for model 1

- Molecule 1: COMPLEMENT FACTOR H



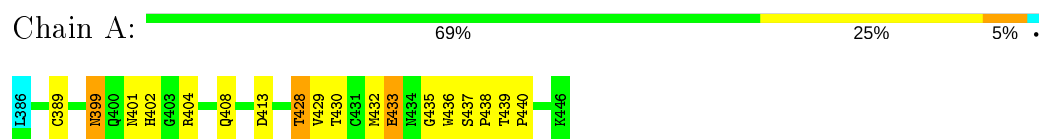
4.2.2 Score per residue for model 2

- Molecule 1: COMPLEMENT FACTOR H



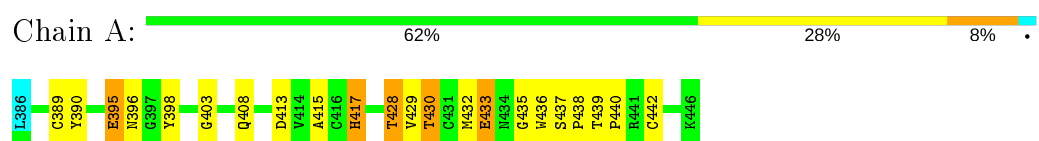
4.2.3 Score per residue for model 3

- Molecule 1: COMPLEMENT FACTOR H



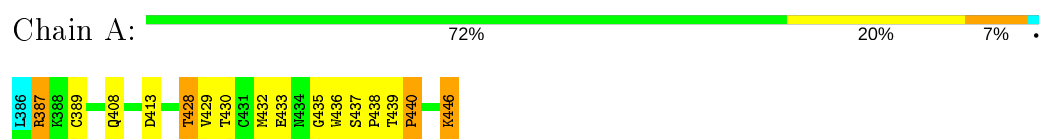
4.2.4 Score per residue for model 4

- Molecule 1: COMPLEMENT FACTOR H



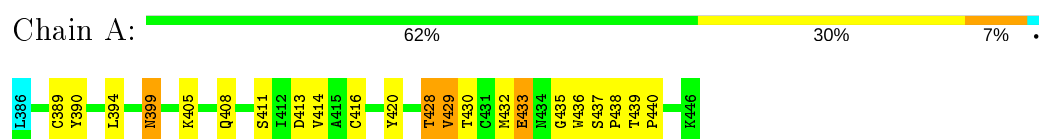
4.2.5 Score per residue for model 5

- Molecule 1: COMPLEMENT FACTOR H



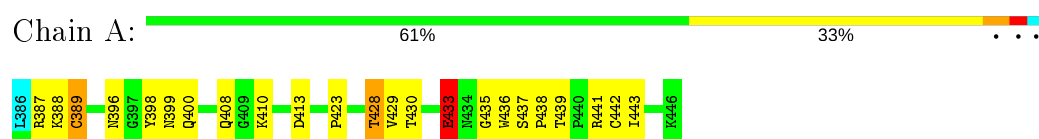
4.2.6 Score per residue for model 6 (medoid)

- Molecule 1: COMPLEMENT FACTOR H



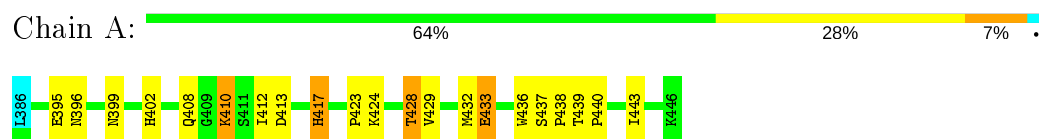
4.2.7 Score per residue for model 7

- Molecule 1: COMPLEMENT FACTOR H



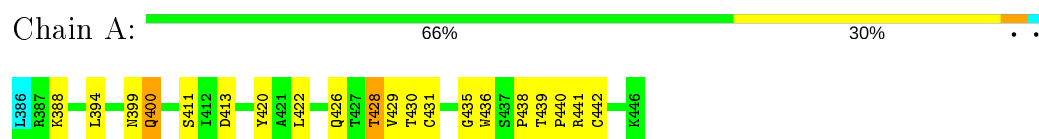
4.2.8 Score per residue for model 8

- Molecule 1: COMPLEMENT FACTOR H



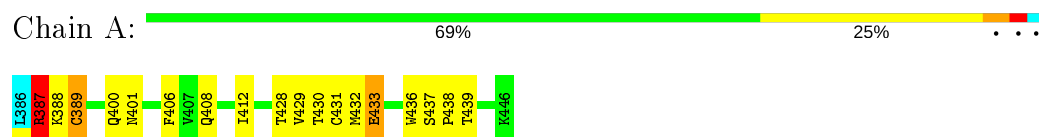
4.2.9 Score per residue for model 9

- Molecule 1: COMPLEMENT FACTOR H



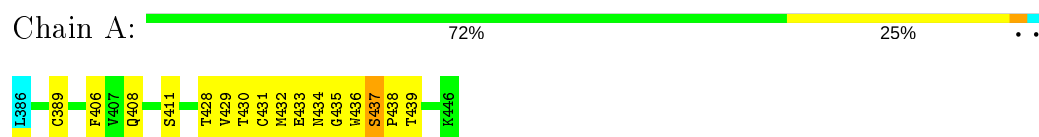
4.2.10 Score per residue for model 10

- Molecule 1: COMPLEMENT FACTOR H



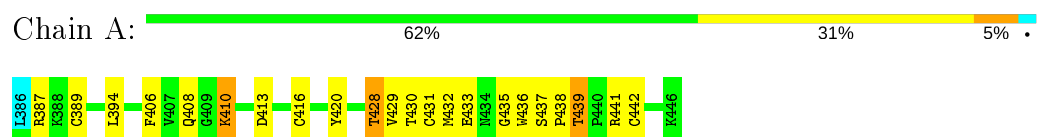
4.2.11 Score per residue for model 11

- Molecule 1: COMPLEMENT FACTOR H



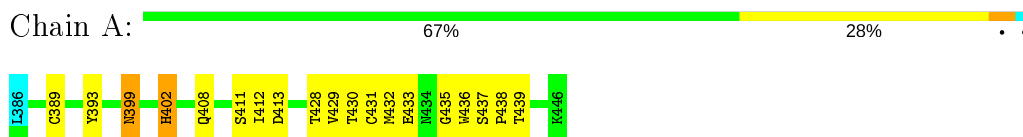
4.2.12 Score per residue for model 12

- Molecule 1: COMPLEMENT FACTOR H



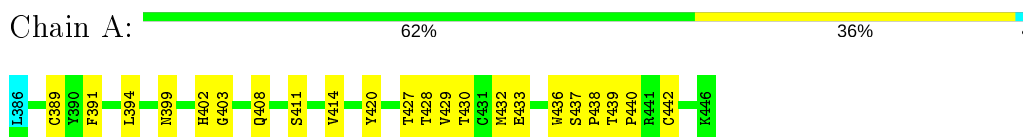
4.2.13 Score per residue for model 13

- Molecule 1: COMPLEMENT FACTOR H



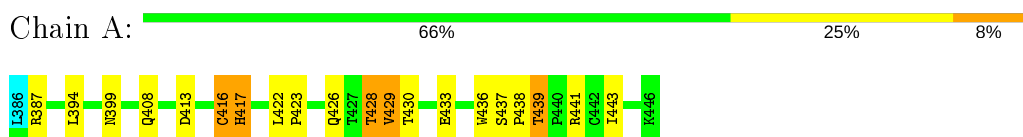
4.2.14 Score per residue for model 14

- Molecule 1: COMPLEMENT FACTOR H



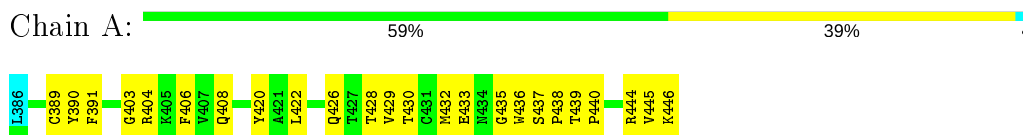
4.2.15 Score per residue for model 15

- Molecule 1: COMPLEMENT FACTOR H



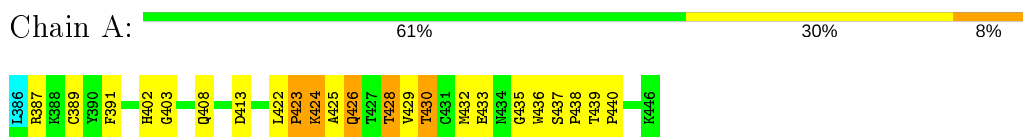
4.2.16 Score per residue for model 16

- Molecule 1: COMPLEMENT FACTOR H



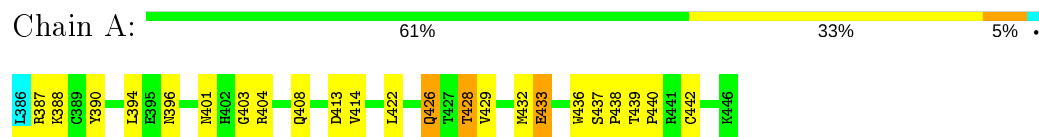
4.2.17 Score per residue for model 17

- Molecule 1: COMPLEMENT FACTOR H



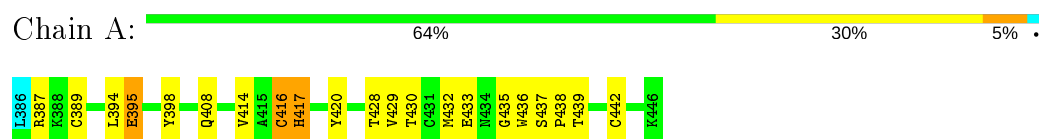
4.2.18 Score per residue for model 18

- Molecule 1: COMPLEMENT FACTOR H



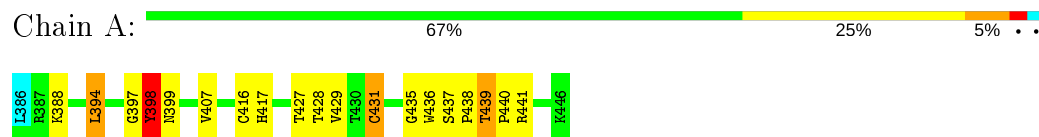
4.2.19 Score per residue for model 19

- Molecule 1: COMPLEMENT FACTOR H



4.2.20 Score per residue for model 20

- Molecule 1: COMPLEMENT FACTOR H



5 Refinement protocol and experimental data overview

The models were refined using the following method: *ARIA 2.0*.

Of the 50 calculated structures, 20 were deposited, based on the following criterion: *LOWEST OVERALL ENERGY*.

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

Software name	Classification	Version
CNS	refinement	1.1
CCPNMR ANALYSIS	structure solution	ANALYSIS

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

6 Model quality i

6.1 Standard geometry i

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 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	481	469	467	17±5
All	All	9620	9380	9340	336

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 18.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:399:ASN:HB3	1:A:402:HIS:HB2	0.90	1.43	13	2
1:A:389:CYS:SG	1:A:406:PHE:HB2	0.83	2.13	11	2
1:A:406:PHE:HB2	1:A:410:LYS:HE3	0.83	1.50	12	1
1:A:429:VAL:HA	1:A:438:PRO:HG2	0.82	1.52	8	19
1:A:406:PHE:HB3	1:A:410:LYS:HG2	0.77	1.55	12	1
1:A:398:TYR:CD1	1:A:416:CYS:HA	0.76	2.14	20	1
1:A:387:ARG:HA	1:A:408:GLN:HB2	0.74	1.58	5	2
1:A:410:LYS:HA	1:A:410:LYS:HE2	0.74	1.60	7	1
1:A:390:TYR:HA	1:A:405:LYS:HB3	0.74	1.60	6	1
1:A:422:LEU:HB2	1:A:426:GLN:O	0.73	1.84	17	3
1:A:391:PHE:CZ	1:A:406:PHE:CD2	0.73	2.77	16	1
1:A:391:PHE:HZ	1:A:406:PHE:N	0.71	1.83	16	1
1:A:423:PRO:HD2	1:A:425:ALA:HB3	0.70	1.63	17	1
1:A:410:LYS:HE2	1:A:436:TRP:CH2	0.69	2.23	12	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:391:PHE:CD1	1:A:436:TRP:HZ2	0.69	2.05	16	1
1:A:413:ASP:HA	1:A:428:THR:HB	0.68	1.65	13	1
1:A:406:PHE:HB2	1:A:410:LYS:CE	0.67	2.19	12	1
1:A:408:GLN:NE2	1:A:433:GLU:HA	0.66	2.06	16	14
1:A:424:LYS:HE2	1:A:424:LYS:HA	0.66	1.67	2	1
1:A:406:PHE:CB	1:A:410:LYS:HE3	0.66	2.20	12	1
1:A:389:CYS:SG	1:A:435:GLY:HA2	0.65	2.31	1	11
1:A:391:PHE:CD2	1:A:404:ARG:HB2	0.65	2.27	16	1
1:A:391:PHE:HZ	1:A:406:PHE:H	0.65	1.33	16	1
1:A:429:VAL:CG2	1:A:436:TRP:HB3	0.65	2.22	4	19
1:A:401:ASN:HD22	1:A:406:PHE:HZ	0.65	1.35	1	1
1:A:410:LYS:HE3	1:A:410:LYS:HA	0.64	1.70	8	1
1:A:431:CYS:SG	1:A:436:TRP:CZ3	0.64	2.90	10	2
1:A:417:HIS:CG	1:A:418:PRO:HD2	0.63	2.28	2	1
1:A:417:HIS:H	1:A:417:HIS:CD2	0.63	2.10	15	2
1:A:391:PHE:CD1	1:A:436:TRP:CZ2	0.62	2.86	16	1
1:A:437:SER:HB3	1:A:438:PRO:HD3	0.62	1.69	10	10
1:A:437:SER:OG	1:A:438:PRO:HD3	0.62	1.94	5	9
1:A:387:ARG:HB2	1:A:408:GLN:OE1	0.62	1.95	17	1
1:A:394:LEU:HB2	1:A:397:GLY:O	0.61	1.95	20	1
1:A:414:VAL:HB	1:A:422:LEU:HD21	0.61	1.71	18	1
1:A:394:LEU:HD21	1:A:441:ARG:HA	0.60	1.73	9	1
1:A:394:LEU:CD2	1:A:441:ARG:HA	0.58	2.28	9	1
1:A:413:ASP:HA	1:A:428:THR:HA	0.58	1.76	1	12
1:A:410:LYS:HE2	1:A:436:TRP:CZ3	0.58	2.32	12	1
1:A:391:PHE:CZ	1:A:406:PHE:N	0.58	2.71	16	1
1:A:417:HIS:CD2	1:A:417:HIS:H	0.58	2.14	8	1
1:A:398:TYR:HE2	1:A:416:CYS:CA	0.58	2.11	19	1
1:A:420:TYR:HB3	1:A:442:CYS:HB3	0.58	1.75	19	4
1:A:413:ASP:HA	1:A:428:THR:CB	0.58	2.28	13	1
1:A:391:PHE:CE2	1:A:406:PHE:CD2	0.58	2.92	16	1
1:A:422:LEU:HB2	1:A:426:GLN:HB2	0.57	1.75	9	1
1:A:429:VAL:HG23	1:A:438:PRO:HD2	0.57	1.75	16	17
1:A:391:PHE:N	1:A:391:PHE:CD1	0.57	2.64	16	1
1:A:429:VAL:HG11	1:A:440:PRO:HB3	0.57	1.76	16	1
1:A:424:LYS:O	1:A:426:GLN:HG3	0.57	2.00	17	1
1:A:387:ARG:CA	1:A:408:GLN:HB2	0.56	2.30	5	1
1:A:430:THR:O	1:A:436:TRP:HA	0.56	2.00	14	16
1:A:398:TYR:OH	1:A:414:VAL:HG13	0.56	2.01	19	1
1:A:408:GLN:HE21	1:A:433:GLU:HA	0.55	1.62	11	3
1:A:417:HIS:CD2	1:A:418:PRO:HD2	0.55	2.36	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:391:PHE:HZ	1:A:414:VAL:HG22	0.55	1.62	14	1
1:A:414:VAL:HB	1:A:422:LEU:HD11	0.55	1.79	1	2
1:A:432:MET:HG2	1:A:435:GLY:O	0.55	2.01	16	6
1:A:399:ASN:O	1:A:400:GLN:HG2	0.55	2.01	9	1
1:A:390:TYR:CE1	1:A:403:GLY:HA2	0.54	2.36	16	3
1:A:420:TYR:CE2	1:A:444:ARG:HB2	0.54	2.37	16	1
1:A:429:VAL:HG21	1:A:440:PRO:HG3	0.54	1.80	2	1
1:A:390:TYR:HA	1:A:405:LYS:CB	0.54	2.30	6	1
1:A:429:VAL:HG22	1:A:436:TRP:CE3	0.54	2.38	12	5
1:A:436:TRP:CB	1:A:440:PRO:HD3	0.53	2.33	4	12
1:A:389:CYS:SG	1:A:406:PHE:CB	0.53	2.95	11	2
1:A:398:TYR:N	1:A:398:TYR:CD1	0.53	2.77	20	1
1:A:388:LYS:HG2	1:A:407:VAL:HG23	0.53	1.80	20	1
1:A:408:GLN:HA	1:A:431:CYS:SG	0.53	2.44	13	2
1:A:398:TYR:CE2	1:A:416:CYS:N	0.53	2.77	19	1
1:A:398:TYR:CE2	1:A:416:CYS:CA	0.52	2.91	19	1
1:A:399:ASN:CB	1:A:402:HIS:HB2	0.52	2.34	1	1
1:A:398:TYR:N	1:A:398:TYR:CD2	0.52	2.71	19	1
1:A:425:ALA:O	1:A:426:GLN:HB2	0.52	2.05	17	1
1:A:391:PHE:CE2	1:A:404:ARG:C	0.51	2.84	16	1
1:A:398:TYR:HE2	1:A:416:CYS:HA	0.51	1.66	19	1
1:A:398:TYR:CE2	1:A:416:CYS:HA	0.51	2.40	19	1
1:A:394:LEU:HD23	1:A:416:CYS:SG	0.50	2.46	15	1
1:A:445:VAL:O	1:A:446:LYS:HB2	0.50	2.06	16	1
1:A:432:MET:SD	1:A:437:SER:HB3	0.50	2.47	1	6
1:A:432:MET:HB3	1:A:437:SER:HB3	0.50	1.83	11	8
1:A:394:LEU:HD23	1:A:442:CYS:SG	0.49	2.47	18	3
1:A:399:ASN:HB2	1:A:402:HIS:HB2	0.49	1.84	8	2
1:A:423:PRO:HA	1:A:443:ILE:HD11	0.49	1.84	7	3
1:A:424:LYS:HE2	1:A:438:PRO:HA	0.49	1.84	8	1
1:A:399:ASN:HB3	1:A:402:HIS:CB	0.49	2.29	13	1
1:A:410:LYS:HZ3	1:A:431:CYS:HB2	0.49	1.68	12	1
1:A:389:CYS:HB3	1:A:435:GLY:HA2	0.49	1.85	7	1
1:A:396:ASN:OD1	1:A:442:CYS:HB2	0.49	2.08	7	1
1:A:400:GLN:HG3	1:A:401:ASN:OD1	0.49	2.08	10	1
1:A:432:MET:HG3	1:A:433:GLU:N	0.48	2.23	5	9
1:A:446:LYS:HB3	1:A:446:LYS:NZ	0.48	2.23	5	1
1:A:387:ARG:HB2	1:A:408:GLN:NE2	0.48	2.24	12	1
1:A:396:ASN:HB2	1:A:417:HIS:ND1	0.48	2.24	4	2
1:A:432:MET:HB3	1:A:437:SER:HB2	0.48	1.84	10	1
1:A:387:ARG:O	1:A:387:ARG:HG3	0.48	2.09	18	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:439:THR:O	1:A:441:ARG:HG2	0.48	2.09	12	2
1:A:416:CYS:SG	1:A:420:TYR:HB2	0.48	2.49	6	3
1:A:410:LYS:HD2	1:A:431:CYS:CB	0.48	2.39	12	1
1:A:394:LEU:HD13	1:A:442:CYS:H	0.47	1.69	9	1
1:A:441:ARG:HG3	1:A:441:ARG:O	0.47	2.09	9	1
1:A:423:PRO:HD3	1:A:441:ARG:O	0.46	2.10	7	2
1:A:394:LEU:HG	1:A:398:TYR:CE2	0.46	2.45	20	1
1:A:391:PHE:O	1:A:402:HIS:ND1	0.45	2.49	14	2
1:A:439:THR:HG22	1:A:441:ARG:HG2	0.45	1.88	20	1
1:A:410:LYS:CE	1:A:431:CYS:HB2	0.45	2.40	12	1
1:A:389:CYS:HB2	1:A:410:LYS:NZ	0.45	2.26	12	1
1:A:429:VAL:HG21	1:A:436:TRP:HB3	0.45	1.87	13	2
1:A:420:TYR:HA	1:A:444:ARG:HA	0.45	1.89	1	1
1:A:422:LEU:HB2	1:A:426:GLN:HB3	0.45	1.87	18	1
1:A:398:TYR:HB2	1:A:415:ALA:HB3	0.45	1.88	4	1
1:A:394:LEU:HB3	1:A:417:HIS:CE1	0.44	2.47	19	1
1:A:398:TYR:HB3	1:A:400:GLN:OE1	0.44	2.12	7	1
1:A:417:HIS:N	1:A:417:HIS:CD2	0.44	2.85	8	2
1:A:394:LEU:HD11	1:A:440:PRO:HB2	0.44	1.89	14	1
1:A:422:LEU:HD13	1:A:425:ALA:O	0.44	2.12	17	1
1:A:394:LEU:HB3	1:A:417:HIS:HE1	0.44	1.72	19	1
1:A:394:LEU:H	1:A:399:ASN:ND2	0.44	2.10	6	1
1:A:431:CYS:SG	1:A:435:GLY:HA2	0.44	2.53	20	2
1:A:391:PHE:CE2	1:A:404:ARG:HB2	0.43	2.47	16	1
1:A:393:TYR:HA	1:A:402:HIS:CD2	0.43	2.48	13	1
1:A:414:VAL:HG21	1:A:440:PRO:HB3	0.43	1.90	6	1
1:A:390:TYR:CA	1:A:405:LYS:HB3	0.43	2.40	6	1
1:A:411:SER:OG	1:A:430:THR:HG22	0.43	2.14	14	4
1:A:391:PHE:HD2	1:A:404:ARG:HB2	0.42	1.68	16	1
1:A:411:SER:HA	1:A:430:THR:HB	0.42	1.90	1	1
1:A:423:PRO:HD2	1:A:425:ALA:CB	0.42	2.42	17	1
1:A:429:VAL:HG23	1:A:436:TRP:CE3	0.42	2.50	2	1
1:A:401:ASN:HA	1:A:404:ARG:HD2	0.42	1.91	3	1
1:A:406:PHE:CB	1:A:410:LYS:HG2	0.41	2.39	12	1
1:A:401:ASN:CB	1:A:404:ARG:HD2	0.41	2.46	18	1
1:A:402:HIS:CG	1:A:403:GLY:N	0.41	2.88	14	2
1:A:417:HIS:CG	1:A:418:PRO:CD	0.41	3.03	2	1
1:A:406:PHE:HB2	1:A:410:LYS:NZ	0.41	2.29	12	1
1:A:389:CYS:HB2	1:A:410:LYS:HZ3	0.41	1.76	12	1
1:A:410:LYS:NZ	1:A:431:CYS:HB2	0.41	2.30	12	1
1:A:422:LEU:HB3	1:A:425:ALA:HB3	0.41	1.92	17	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:436:TRP:HB2	1:A:440:PRO:HD3	0.41	1.92	4	1
1:A:436:TRP:HB3	1:A:440:PRO:HD3	0.41	1.91	20	2
1:A:411:SER:CB	1:A:430:THR:HG22	0.40	2.46	13	1
1:A:390:TYR:CZ	1:A:403:GLY:HA2	0.40	2.51	2	1
1:A:410:LYS:CE	1:A:436:TRP:CH2	0.40	3.02	12	1

6.3 Torsion angles [i](#)

6.3.1 Protein backbone [i](#)

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	59/61 (97%)	50±2 (84±3%)	7±2 (12±3%)	2±1 (4±1%)	6	34
All	All	1180/1220 (97%)	993 (84%)	144 (12%)	43 (4%)	6	34

All 14 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	439	THR	20
1	A	433	GLU	9
1	A	395	GLU	3
1	A	387	ARG	1
1	A	423	PRO	1
1	A	426	GLN	1
1	A	445	VAL	1
1	A	388	LYS	1
1	A	442	CYS	1
1	A	398	TYR	1
1	A	400	GLN	1
1	A	416	CYS	1
1	A	440	PRO	1
1	A	412	ILE	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	52/53 (98%)	49±2 (93±3%)	3±2 (7±3%)	21 69
All	All	1040/1060 (98%)	972 (93%)	68 (7%)	21 69

All 23 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	428	THR	19
1	A	399	ASN	8
1	A	417	HIS	5
1	A	387	ARG	5
1	A	430	THR	4
1	A	389	CYS	3
1	A	429	VAL	3
1	A	427	THR	2
1	A	388	LYS	2
1	A	410	LYS	2
1	A	412	ILE	2
1	A	395	GLU	2
1	A	446	LYS	1
1	A	424	LYS	1
1	A	426	GLN	1
1	A	398	TYR	1
1	A	416	CYS	1
1	A	394	LEU	1
1	A	396	ASN	1
1	A	433	GLU	1
1	A	437	SER	1
1	A	431	CYS	1
1	A	402	HIS	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

No chemical shift data were provided