



Full wwPDB NMR Structure Validation Report ⓘ

Feb 17, 2022 – 10:57 AM EST

PDB ID : 1P1E
Title : Structural Insights into the Inter-domain Chaperoning of Tandem PDZ Domains in Glutamate Receptor Interacting Proteins
Authors : Feng, W.; Shi, Y.; Li, M.; Zhang, M.
Deposited on : 2003-04-12

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:

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 : 2.26
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.26

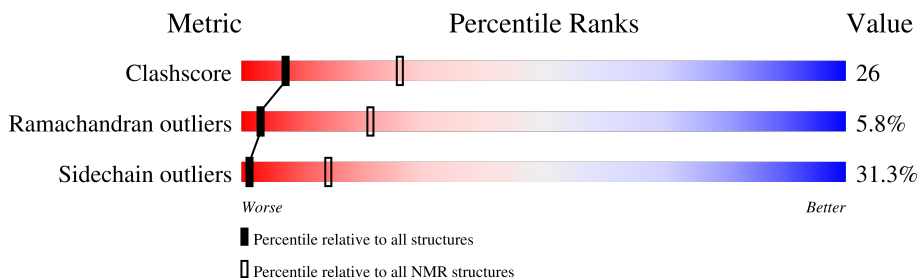
1 Overall quality at a glance

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	101	

2 Ensemble composition and analysis i

This entry contains 20 models. Model 9 is the overall representative, medoid model (most similar to other models). The authors have identified model 13 as representative, based on the following criterion: *minimized average structure*.

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:22-A:42, A:55-A:85, A:93-A:113 (73)	0.38	9

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. No single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1504 atoms, of which 744 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Glutamate receptor interacting protein.

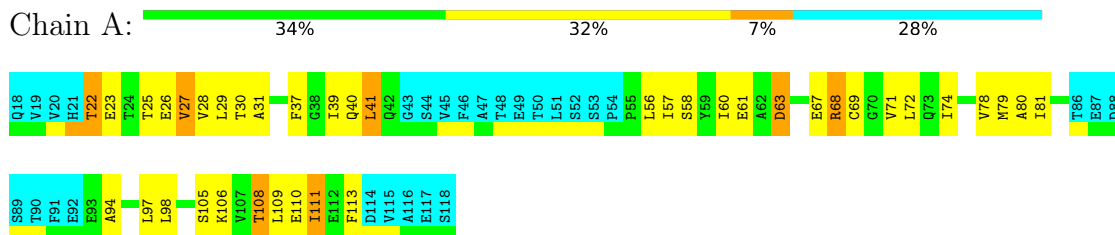
Mol	Chain	Residues	Atoms						Trace
			Total	C	H	N	O	S	
1	A	101	1504	473	744	120	165	2	0

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: Glutamate receptor interacting protein

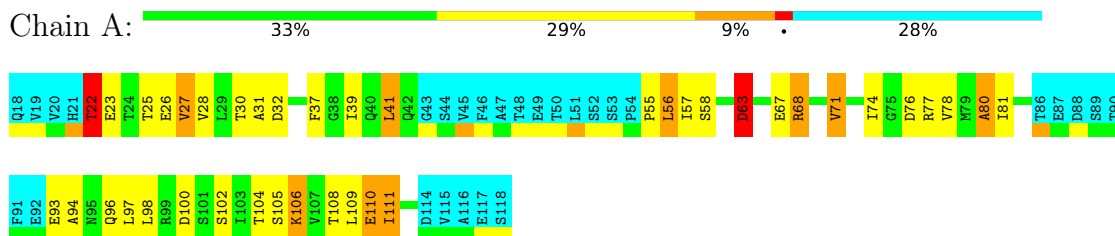


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

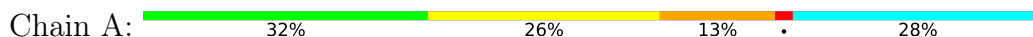
4.2.1 Score per residue for model 1

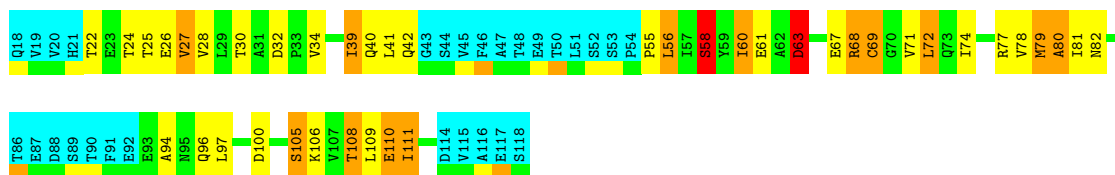
- Molecule 1: Glutamate receptor interacting protein



4.2.2 Score per residue for model 2

- Molecule 1: Glutamate receptor interacting protein

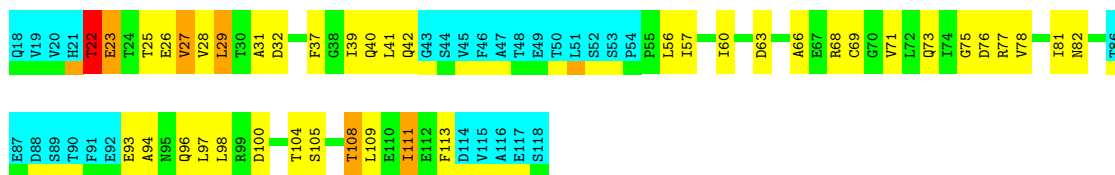




4.2.3 Score per residue for model 3

- Molecule 1: Glutamate receptor interacting protein

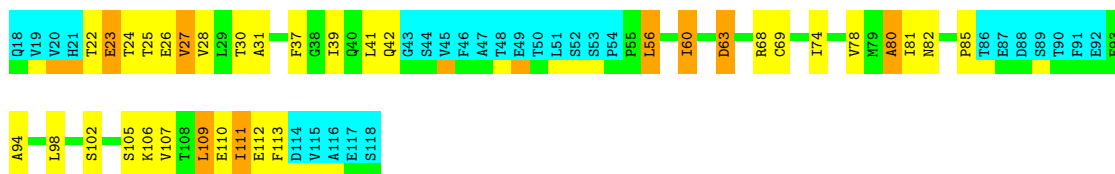
Chain A: 32% 35% 5% 28%



4.2.4 Score per residue for model 4

- Molecule 1: Glutamate receptor interacting protein

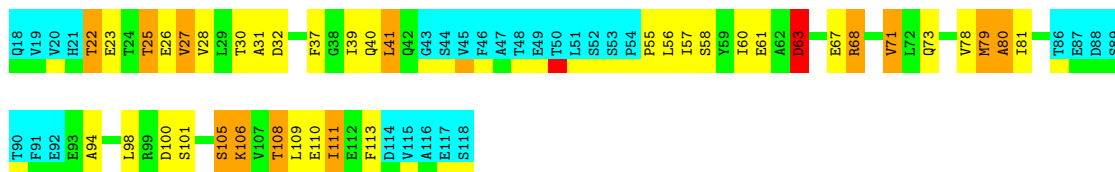
Chain A: 38% 27% 8% 28%



4.2.5 Score per residue for model 5

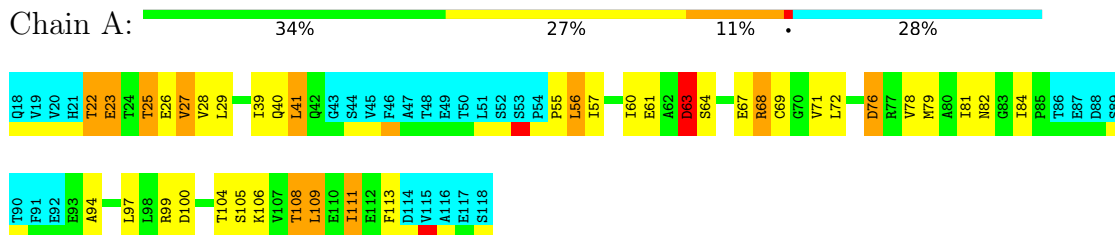
- Molecule 1: Glutamate receptor interacting protein

Chain A: 34% 26% 12% 28%



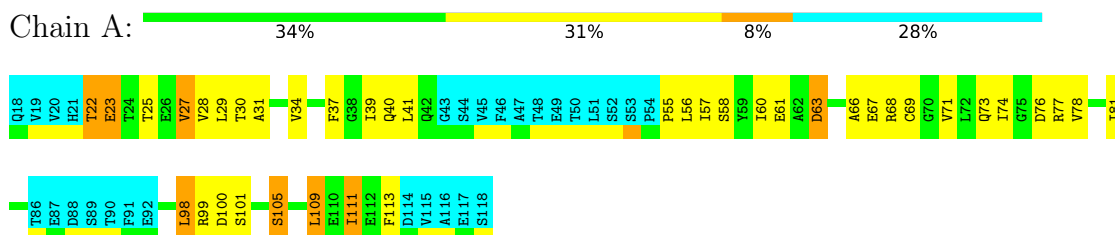
4.2.6 Score per residue for model 6

- Molecule 1: Glutamate receptor interacting protein



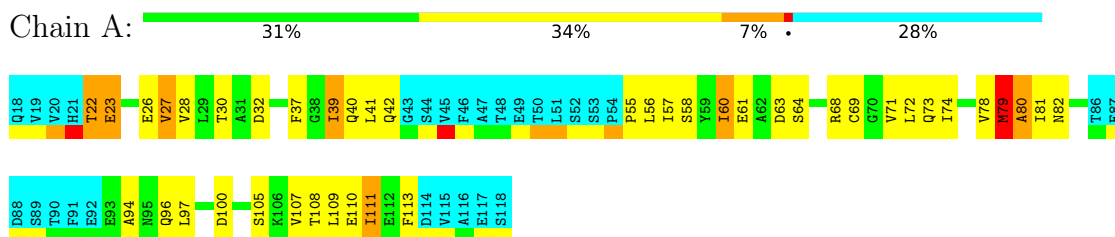
4.2.7 Score per residue for model 7

- Molecule 1: Glutamate receptor interacting protein



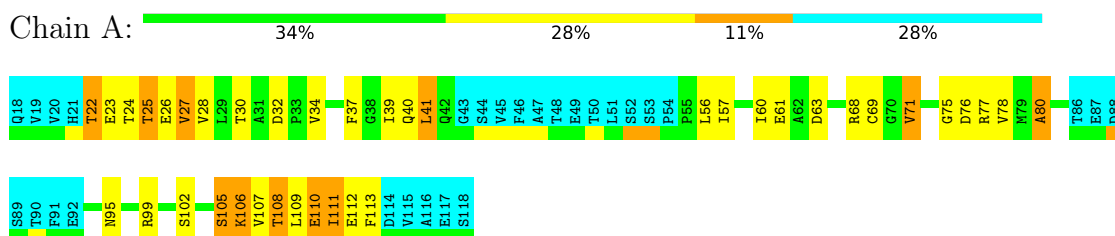
4.2.8 Score per residue for model 8

- Molecule 1: Glutamate receptor interacting protein



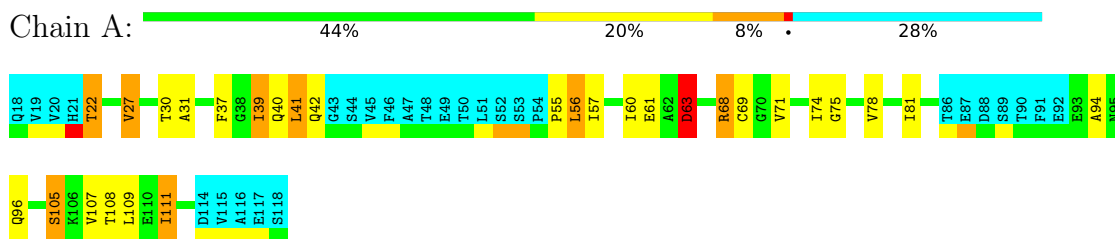
4.2.9 Score per residue for model 9 (medoid)

- Molecule 1: Glutamate receptor interacting protein



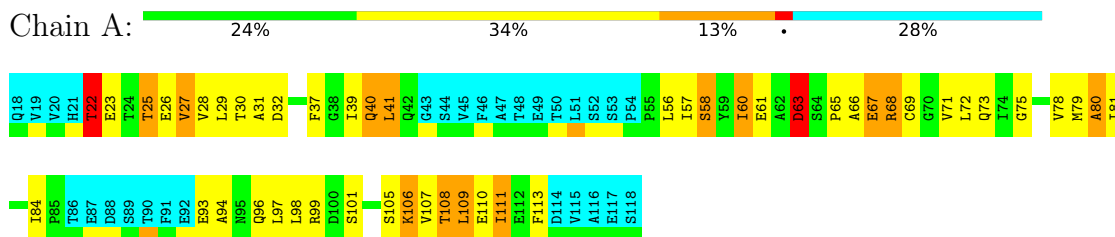
4.2.10 Score per residue for model 10

- Molecule 1: Glutamate receptor interacting protein



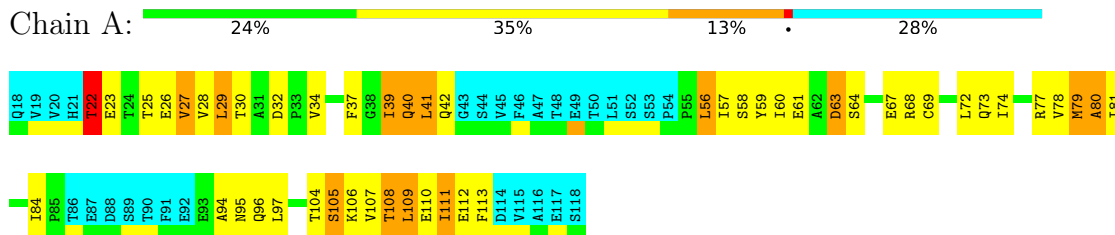
4.2.11 Score per residue for model 11

- Molecule 1: Glutamate receptor interacting protein



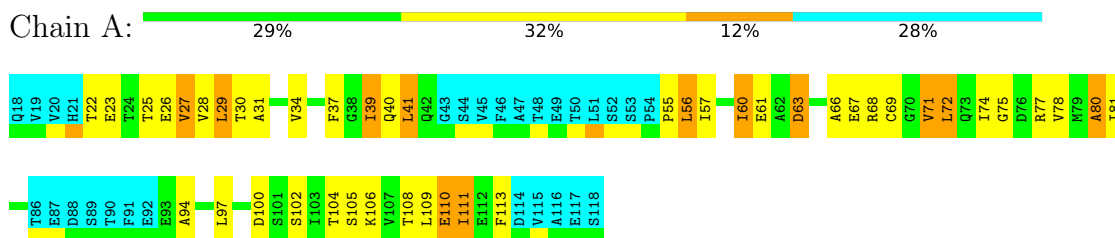
4.2.12 Score per residue for model 12

- Molecule 1: Glutamate receptor interacting protein



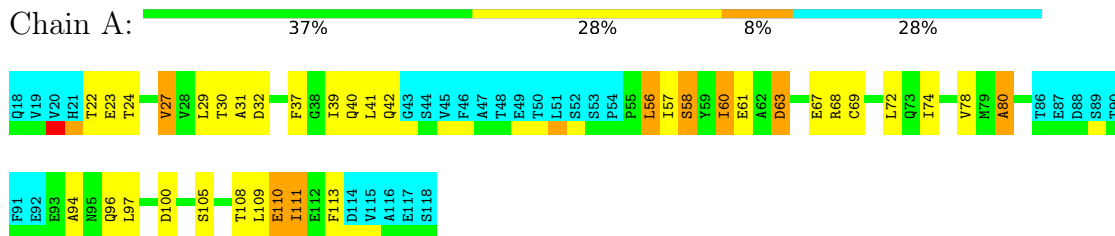
4.2.13 Score per residue for model 13

- Molecule 1: Glutamate receptor interacting protein



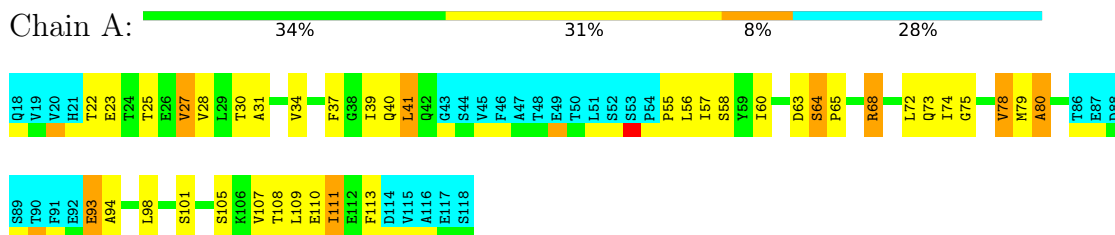
4.2.14 Score per residue for model 14

- Molecule 1: Glutamate receptor interacting protein



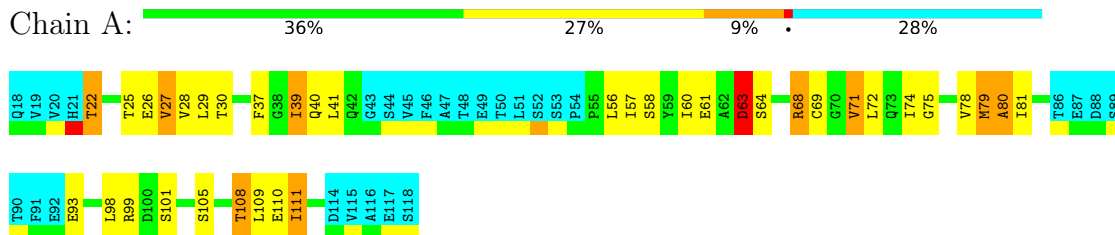
4.2.15 Score per residue for model 15

- Molecule 1: Glutamate receptor interacting protein



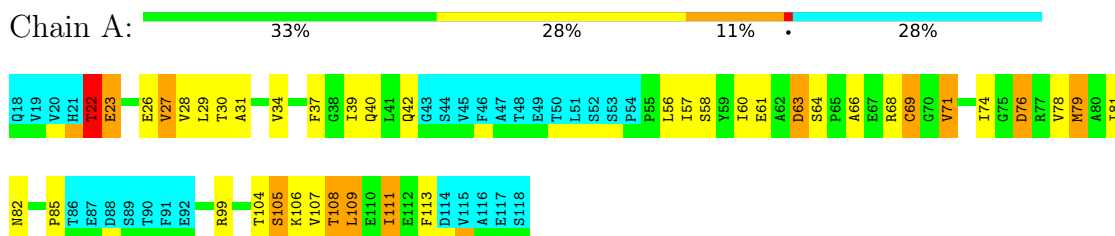
4.2.16 Score per residue for model 16

- Molecule 1: Glutamate receptor interacting protein



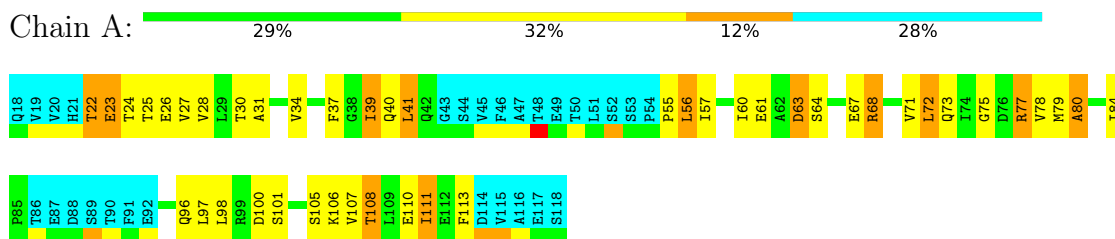
4.2.17 Score per residue for model 17

- Molecule 1: Glutamate receptor interacting protein



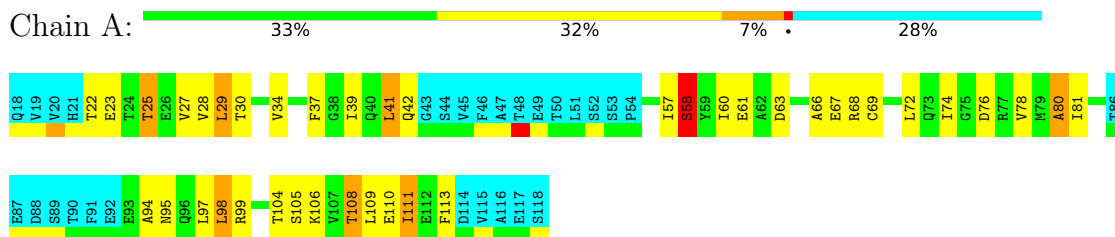
4.2.18 Score per residue for model 18

- Molecule 1: Glutamate receptor interacting protein



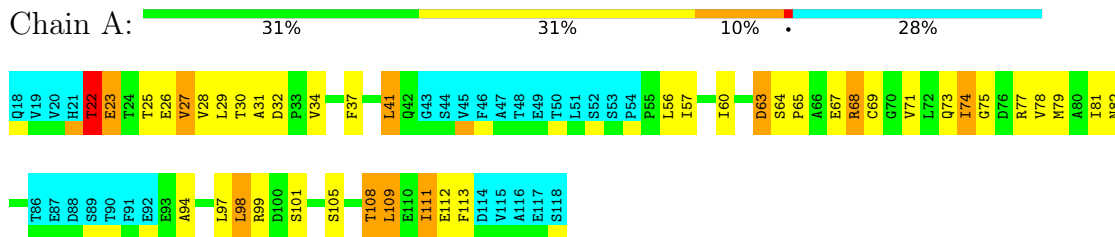
4.2.19 Score per residue for model 19

- Molecule 1: Glutamate receptor interacting protein



4.2.20 Score per residue for model 20

- Molecule 1: Glutamate receptor interacting protein



5 Refinement protocol and experimental data overview

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

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
CNS	structure solution	1.1
CNS	refinement	1.1

No chemical shift data was provided.

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	551	556	556	29±5
All	All	11020	11120	11120	584

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:78:VAL:HG11	1:A:81:ILE:HD11	0.95	1.35	12	12
1:A:78:VAL:HG21	1:A:81:ILE:HD11	0.92	1.38	17	1
1:A:28:VAL:HG22	1:A:108:THR:HG22	0.87	1.43	17	10
1:A:41:LEU:HD13	1:A:94:ALA:HB1	0.87	1.46	20	4
1:A:78:VAL:HG22	1:A:111:ILE:HG22	0.86	1.47	16	6
1:A:78:VAL:HG22	1:A:111:ILE:CG2	0.85	2.01	4	10
1:A:41:LEU:HD11	1:A:98:LEU:HD11	0.80	1.51	3	2
1:A:39:ILE:HD13	1:A:72:LEU:HD13	0.79	1.53	6	4
1:A:28:VAL:HG22	1:A:108:THR:CG2	0.79	2.07	16	10
1:A:39:ILE:HD11	1:A:57:ILE:CG2	0.79	2.08	14	9
1:A:97:LEU:HD12	1:A:98:LEU:N	0.78	1.93	20	1
1:A:28:VAL:HG23	1:A:108:THR:HG23	0.78	1.53	9	1
1:A:55:PRO:HB2	1:A:78:VAL:HG23	0.78	1.55	18	1
1:A:31:ALA:HB2	1:A:37:PHE:CE1	0.77	2.14	10	9

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:78:VAL:CG1	1:A:111:ILE:HG23	0.76	2.10	9	2
1:A:80:ALA:HB3	1:A:110:GLU:HB3	0.75	1.57	4	10
1:A:29:LEU:HD23	1:A:109:LEU:HD21	0.75	1.56	7	1
1:A:25:THR:HG23	1:A:111:ILE:CD1	0.75	2.11	5	7
1:A:78:VAL:HG13	1:A:111:ILE:HG23	0.75	1.59	9	1
1:A:30:THR:HG23	1:A:105:SER:O	0.74	1.83	9	18
1:A:78:VAL:CG2	1:A:81:ILE:HD11	0.73	2.13	17	1
1:A:41:LEU:HD12	1:A:41:LEU:O	0.73	1.84	8	1
1:A:55:PRO:O	1:A:56:LEU:HD12	0.72	1.83	7	3
1:A:22:THR:HG23	1:A:113:PHE:O	0.72	1.85	8	4
1:A:94:ALA:HA	1:A:97:LEU:HD12	0.72	1.61	14	10
1:A:31:ALA:HB2	1:A:37:PHE:CE2	0.72	2.20	18	3
1:A:58:SER:HA	1:A:74:ILE:HG23	0.72	1.59	7	7
1:A:78:VAL:HG11	1:A:81:ILE:CD1	0.71	2.15	12	9
1:A:41:LEU:CD1	1:A:98:LEU:HD11	0.71	2.15	1	2
1:A:39:ILE:HD11	1:A:57:ILE:HG21	0.71	1.62	1	11
1:A:56:LEU:HD13	1:A:56:LEU:N	0.70	2.01	14	6
1:A:82:ASN:OD1	1:A:97:LEU:HD22	0.69	1.86	6	1
1:A:25:THR:HG23	1:A:111:ILE:HD11	0.69	1.63	5	6
1:A:81:ILE:HG23	1:A:109:LEU:HB3	0.69	1.63	6	2
1:A:41:LEU:HD22	1:A:57:ILE:HG12	0.68	1.65	15	2
1:A:78:VAL:HG21	1:A:81:ILE:CD1	0.68	2.18	17	1
1:A:41:LEU:CD1	1:A:94:ALA:HB1	0.68	2.18	20	1
1:A:27:VAL:CG2	1:A:111:ILE:HD11	0.68	2.19	1	11
1:A:78:VAL:HB	1:A:81:ILE:HD11	0.68	1.65	10	2
1:A:39:ILE:HD13	1:A:72:LEU:CD1	0.67	2.18	2	1
1:A:80:ALA:HB3	1:A:110:GLU:CB	0.67	2.19	2	2
1:A:41:LEU:HD13	1:A:94:ALA:CB	0.67	2.20	10	3
1:A:41:LEU:HD11	1:A:94:ALA:CB	0.66	2.20	8	1
1:A:98:LEU:HD12	1:A:98:LEU:O	0.66	1.91	7	1
1:A:66:ALA:O	1:A:72:LEU:HD11	0.65	1.91	13	3
1:A:56:LEU:O	1:A:56:LEU:HD22	0.65	1.91	12	3
1:A:25:THR:HG23	1:A:111:ILE:HG13	0.65	1.69	19	4
1:A:27:VAL:HG21	1:A:111:ILE:HD11	0.64	1.69	1	7
1:A:28:VAL:HG22	1:A:106:LYS:HD3	0.63	1.70	5	1
1:A:37:PHE:CE1	1:A:107:VAL:HG12	0.62	2.29	8	3
1:A:66:ALA:HA	1:A:72:LEU:HD21	0.62	1.71	11	1
1:A:27:VAL:HG23	1:A:109:LEU:HD12	0.62	1.70	11	1
1:A:55:PRO:C	1:A:56:LEU:HD13	0.61	2.16	2	4
1:A:80:ALA:HB3	1:A:110:GLU:HB2	0.61	1.72	12	4
1:A:28:VAL:HG22	1:A:106:LYS:HD2	0.61	1.71	9	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:55:PRO:CB	1:A:78:VAL:HG23	0.61	2.26	18	1
1:A:41:LEU:HD11	1:A:94:ALA:HB3	0.61	1.71	8	2
1:A:55:PRO:C	1:A:56:LEU:HD12	0.61	2.16	15	2
1:A:41:LEU:HD11	1:A:98:LEU:CD2	0.60	2.27	18	1
1:A:41:LEU:HD12	1:A:94:ALA:HB1	0.60	1.73	15	1
1:A:40:GLN:C	1:A:41:LEU:HD13	0.60	2.17	12	1
1:A:41:LEU:HD23	1:A:95:ASN:N	0.60	2.12	12	1
1:A:78:VAL:HA	1:A:111:ILE:HG22	0.59	1.72	18	4
1:A:78:VAL:HG13	1:A:111:ILE:CG2	0.59	2.27	9	2
1:A:41:LEU:HD23	1:A:98:LEU:HD11	0.59	1.72	16	1
1:A:41:LEU:HD23	1:A:78:VAL:HG21	0.59	1.72	13	1
1:A:78:VAL:HG22	1:A:111:ILE:HG23	0.59	1.75	2	3
1:A:71:VAL:HG22	1:A:71:VAL:O	0.59	1.98	1	4
1:A:66:ALA:CA	1:A:72:LEU:HD21	0.59	2.26	11	1
1:A:69:CYS:O	1:A:71:VAL:HG23	0.58	1.98	2	2
1:A:84:ILE:HG21	1:A:97:LEU:HD21	0.58	1.74	11	2
1:A:23:GLU:HB2	1:A:113:PHE:CZ	0.58	2.32	7	4
1:A:31:ALA:HB3	1:A:105:SER:HA	0.58	1.75	7	4
1:A:81:ILE:HG23	1:A:109:LEU:HD22	0.58	1.75	20	1
1:A:57:ILE:CD1	1:A:111:ILE:HG21	0.58	2.28	12	5
1:A:27:VAL:HG11	1:A:71:VAL:HB	0.58	1.74	20	2
1:A:56:LEU:HD23	1:A:56:LEU:N	0.57	2.15	3	4
1:A:104:THR:HG23	1:A:106:LYS:O	0.57	1.99	6	6
1:A:60:ILE:HD11	1:A:74:ILE:HD13	0.57	1.76	4	1
1:A:74:ILE:HG23	1:A:74:ILE:O	0.56	2.00	15	2
1:A:84:ILE:HB	1:A:97:LEU:HD13	0.56	1.77	18	1
1:A:41:LEU:CD1	1:A:94:ALA:HB3	0.56	2.31	5	1
1:A:39:ILE:CD1	1:A:72:LEU:HD13	0.56	2.29	6	1
1:A:41:LEU:HD22	1:A:41:LEU:N	0.55	2.15	12	1
1:A:111:ILE:H	1:A:111:ILE:HD13	0.55	1.61	20	5
1:A:79:MET:O	1:A:85:PRO:HA	0.55	2.01	17	1
1:A:111:ILE:N	1:A:111:ILE:HD13	0.55	2.16	10	3
1:A:41:LEU:CD2	1:A:98:LEU:HD11	0.55	2.31	16	1
1:A:107:VAL:O	1:A:107:VAL:HG13	0.55	2.01	8	1
1:A:78:VAL:CA	1:A:111:ILE:HG22	0.55	2.32	18	4
1:A:31:ALA:HB2	1:A:37:PHE:CZ	0.55	2.37	11	4
1:A:41:LEU:N	1:A:41:LEU:HD12	0.54	2.18	1	2
1:A:29:LEU:HD11	1:A:66:ALA:N	0.54	2.16	3	1
1:A:27:VAL:CG1	1:A:71:VAL:HG23	0.54	2.31	7	1
1:A:41:LEU:HD22	1:A:78:VAL:HG21	0.54	1.78	8	1
1:A:57:ILE:HD12	1:A:111:ILE:HG21	0.54	1.79	16	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:27:VAL:HG12	1:A:71:VAL:CG2	0.54	2.31	5	1
1:A:29:LEU:HD11	1:A:72:LEU:CD1	0.54	2.32	16	1
1:A:78:VAL:HG12	1:A:111:ILE:CG2	0.53	2.34	10	2
1:A:29:LEU:HD11	1:A:66:ALA:CA	0.53	2.33	3	1
1:A:37:PHE:HB3	1:A:39:ILE:HG22	0.53	1.80	19	4
1:A:78:VAL:CB	1:A:111:ILE:HG22	0.53	2.34	17	1
1:A:78:VAL:HG12	1:A:81:ILE:HD11	0.53	1.80	13	1
1:A:25:THR:CG2	1:A:111:ILE:HD12	0.53	2.34	4	2
1:A:56:LEU:N	1:A:56:LEU:CD1	0.52	2.72	14	3
1:A:78:VAL:CG1	1:A:81:ILE:HD11	0.52	2.30	4	3
1:A:39:ILE:HD12	1:A:59:TYR:O	0.52	2.04	12	1
1:A:81:ILE:HG22	1:A:82:ASN:HD22	0.52	1.65	20	4
1:A:77:ARG:C	1:A:111:ILE:HG22	0.52	2.25	1	1
1:A:78:VAL:CG1	1:A:111:ILE:HG22	0.52	2.35	15	1
1:A:56:LEU:HD23	1:A:56:LEU:H	0.52	1.65	9	2
1:A:29:LEU:HD21	1:A:66:ALA:HA	0.52	1.82	17	1
1:A:41:LEU:HD23	1:A:98:LEU:CD1	0.52	2.35	16	1
1:A:57:ILE:HD12	1:A:78:VAL:CG2	0.51	2.35	3	1
1:A:58:SER:CA	1:A:74:ILE:HG23	0.51	2.36	2	2
1:A:95:ASN:HA	1:A:98:LEU:HD23	0.51	1.81	19	1
1:A:81:ILE:HB	1:A:97:LEU:HD22	0.51	1.80	13	1
1:A:66:ALA:O	1:A:72:LEU:HD21	0.51	2.05	11	1
1:A:22:THR:O	1:A:22:THR:HG23	0.51	2.05	7	2
1:A:57:ILE:HD12	1:A:76:ASP:HB2	0.51	1.82	19	3
1:A:76:ASP:HB3	1:A:111:ILE:HG21	0.51	1.82	1	1
1:A:41:LEU:HD23	1:A:57:ILE:CG1	0.51	2.36	19	2
1:A:28:VAL:O	1:A:28:VAL:HG13	0.51	2.06	9	8
1:A:41:LEU:HD23	1:A:57:ILE:HG13	0.51	1.83	9	2
1:A:84:ILE:HB	1:A:97:LEU:HD21	0.51	1.81	6	1
1:A:60:ILE:HG13	1:A:74:ILE:HD11	0.50	1.82	8	2
1:A:78:VAL:CG2	1:A:111:ILE:HG22	0.50	2.30	11	2
1:A:37:PHE:CE2	1:A:107:VAL:HG12	0.50	2.42	9	2
1:A:113:PHE:CD1	1:A:113:PHE:C	0.50	2.83	6	1
1:A:40:GLN:C	1:A:41:LEU:HD23	0.50	2.27	11	1
1:A:57:ILE:HD11	1:A:78:VAL:HG22	0.50	1.82	8	1
1:A:27:VAL:HG21	1:A:111:ILE:CD1	0.50	2.37	6	3
1:A:29:LEU:HD12	1:A:65:PRO:HB2	0.50	1.83	20	2
1:A:111:ILE:HD13	1:A:111:ILE:N	0.49	2.22	20	1
1:A:24:THR:HA	1:A:111:ILE:O	0.49	2.07	2	2
1:A:81:ILE:HG22	1:A:82:ASN:OD1	0.49	2.07	2	1
1:A:29:LEU:HB2	1:A:109:LEU:HD21	0.49	1.83	12	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:27:VAL:HG12	1:A:71:VAL:HG23	0.49	1.83	5	1
1:A:56:LEU:HA	1:A:78:VAL:HG22	0.49	1.84	1	1
1:A:78:VAL:HG22	1:A:111:ILE:HG21	0.49	1.83	4	1
1:A:56:LEU:HD13	1:A:56:LEU:H	0.49	1.65	12	2
1:A:57:ILE:HD13	1:A:76:ASP:HB2	0.49	1.85	3	1
1:A:55:PRO:C	1:A:56:LEU:HD23	0.49	2.28	6	1
1:A:39:ILE:HD12	1:A:66:ALA:HB1	0.49	1.85	7	2
1:A:27:VAL:CG1	1:A:71:VAL:HG13	0.48	2.38	9	2
1:A:31:ALA:HB2	1:A:37:PHE:HE1	0.48	1.63	7	3
1:A:27:VAL:HG13	1:A:71:VAL:HG11	0.48	1.84	6	1
1:A:78:VAL:HB	1:A:111:ILE:HG22	0.48	1.85	17	1
1:A:81:ILE:HG22	1:A:82:ASN:ND2	0.48	2.24	8	1
1:A:57:ILE:HD11	1:A:111:ILE:HG21	0.48	1.84	12	3
1:A:29:LEU:HD12	1:A:37:PHE:CD1	0.48	2.44	19	1
1:A:41:LEU:HD11	1:A:98:LEU:HD23	0.48	1.86	18	1
1:A:22:THR:HA	1:A:113:PHE:O	0.48	2.08	14	2
1:A:109:LEU:HD12	1:A:111:ILE:HG23	0.48	1.86	20	1
1:A:41:LEU:HD22	1:A:41:LEU:H	0.48	1.68	12	1
1:A:56:LEU:HD11	1:A:77:ARG:NE	0.48	2.24	18	1
1:A:41:LEU:HD11	1:A:98:LEU:CD1	0.47	2.34	3	1
1:A:60:ILE:HD12	1:A:67:GLU:OE1	0.47	2.09	14	1
1:A:78:VAL:HG11	1:A:81:ILE:CG1	0.47	2.39	20	6
1:A:97:LEU:HD12	1:A:97:LEU:C	0.47	2.29	20	1
1:A:111:ILE:HD13	1:A:111:ILE:H	0.47	1.69	18	3
1:A:98:LEU:HA	1:A:107:VAL:HG21	0.47	1.87	11	1
1:A:29:LEU:HD22	1:A:109:LEU:HD11	0.47	1.85	17	1
1:A:25:THR:HG22	1:A:111:ILE:HD12	0.47	1.86	4	1
1:A:37:PHE:CZ	1:A:101:SER:CB	0.47	2.98	7	2
1:A:28:VAL:CG2	1:A:108:THR:HG23	0.46	2.34	9	1
1:A:23:GLU:CB	1:A:113:PHE:CZ	0.46	2.97	7	1
1:A:27:VAL:HG12	1:A:71:VAL:HG13	0.46	1.87	13	2
1:A:27:VAL:CG1	1:A:71:VAL:HG21	0.46	2.39	8	4
1:A:27:VAL:HG12	1:A:71:VAL:CG1	0.46	2.40	13	1
1:A:29:LEU:HD23	1:A:109:LEU:HD11	0.46	1.88	11	1
1:A:81:ILE:HG23	1:A:108:THR:O	0.46	2.11	5	1
1:A:29:LEU:HD11	1:A:66:ALA:HA	0.46	1.87	13	1
1:A:74:ILE:O	1:A:74:ILE:HG22	0.46	2.11	20	1
1:A:78:VAL:N	1:A:111:ILE:HG22	0.46	2.26	14	1
1:A:98:LEU:HD12	1:A:99:ARG:N	0.46	2.25	11	1
1:A:78:VAL:HG13	1:A:111:ILE:HG22	0.46	1.86	15	1
1:A:60:ILE:HG22	1:A:67:GLU:HB2	0.46	1.88	13	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:79:MET:O	1:A:80:ALA:HB2	0.46	2.11	5	4
1:A:57:ILE:HD12	1:A:76:ASP:HB3	0.45	1.87	6	1
1:A:28:VAL:HG11	1:A:106:LYS:HD3	0.45	1.88	17	2
1:A:57:ILE:HD11	1:A:78:VAL:HG13	0.45	1.88	9	1
1:A:39:ILE:HG23	1:A:98:LEU:HD13	0.45	1.87	4	2
1:A:37:PHE:CE1	1:A:107:VAL:CG1	0.45	2.99	8	2
1:A:81:ILE:N	1:A:81:ILE:HD12	0.45	2.27	13	2
1:A:78:VAL:HG12	1:A:81:ILE:CD1	0.45	2.41	13	1
1:A:41:LEU:HD12	1:A:95:ASN:OD1	0.45	2.11	19	1
1:A:78:VAL:CG1	1:A:81:ILE:CG1	0.45	2.94	3	7
1:A:24:THR:HG22	1:A:112:GLU:HG3	0.45	1.89	9	1
1:A:23:GLU:CG	1:A:113:PHE:CZ	0.45	3.00	19	2
1:A:56:LEU:HA	1:A:78:VAL:HG23	0.45	1.89	3	1
1:A:39:ILE:HD12	1:A:60:ILE:HG23	0.44	1.89	8	1
1:A:37:PHE:CE1	1:A:101:SER:HB2	0.44	2.47	7	1
1:A:107:VAL:HG22	1:A:108:THR:N	0.44	2.28	12	1
1:A:78:VAL:CG1	1:A:111:ILE:CG2	0.44	2.96	18	4
1:A:22:THR:O	1:A:22:THR:CG2	0.44	2.66	7	2
1:A:27:VAL:CG2	1:A:111:ILE:CD1	0.44	2.96	12	8
1:A:27:VAL:HG22	1:A:111:ILE:HD11	0.44	1.90	14	1
1:A:29:LEU:HD11	1:A:69:CYS:SG	0.44	2.53	17	1
1:A:27:VAL:HG11	1:A:71:VAL:CB	0.43	2.42	10	1
1:A:63:ASP:N	1:A:68:ARG:NH2	0.43	2.67	2	7
1:A:104:THR:O	1:A:105:SER:CB	0.43	2.66	12	2
1:A:23:GLU:O	1:A:112:GLU:HA	0.43	2.13	4	1
1:A:71:VAL:HG13	1:A:72:LEU:N	0.43	2.28	11	2
1:A:41:LEU:HD13	1:A:41:LEU:N	0.43	2.27	12	1
1:A:72:LEU:HD12	1:A:72:LEU:O	0.43	2.13	18	1
1:A:31:ALA:CB	1:A:37:PHE:CE1	0.43	3.00	14	1
1:A:41:LEU:HD13	1:A:94:ALA:HB3	0.43	1.90	10	1
1:A:61:GLU:O	1:A:67:GLU:CB	0.43	2.67	11	1
1:A:22:THR:HG22	1:A:22:THR:O	0.43	2.13	6	1
1:A:78:VAL:HG22	1:A:79:MET:N	0.43	2.28	17	1
1:A:60:ILE:CG2	1:A:67:GLU:N	0.43	2.82	11	1
1:A:41:LEU:CD1	1:A:94:ALA:CB	0.42	2.97	14	1
1:A:57:ILE:CD1	1:A:78:VAL:HG22	0.42	2.44	8	2
1:A:57:ILE:HD13	1:A:72:LEU:CD2	0.42	2.44	16	1
1:A:81:ILE:HD12	1:A:81:ILE:H	0.42	1.74	13	1
1:A:22:THR:O	1:A:23:GLU:CG	0.42	2.68	12	1
1:A:67:GLU:CG	1:A:68:ARG:NH2	0.42	2.83	5	3
1:A:63:ASP:C	1:A:68:ARG:NH2	0.42	2.72	11	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:56:LEU:HD11	1:A:77:ARG:HE	0.42	1.74	18	1
1:A:23:GLU:O	1:A:113:PHE:CD1	0.42	2.73	15	5
1:A:37:PHE:CE2	1:A:98:LEU:O	0.42	2.72	11	1
1:A:29:LEU:N	1:A:29:LEU:HD22	0.42	2.30	7	1
1:A:23:GLU:O	1:A:113:PHE:CD2	0.42	2.73	20	2
1:A:56:LEU:N	1:A:56:LEU:HD12	0.42	2.29	17	1
1:A:27:VAL:CG2	1:A:111:ILE:CG1	0.41	2.99	7	1
1:A:23:GLU:HG2	1:A:113:PHE:CZ	0.41	2.50	13	1
1:A:78:VAL:HG12	1:A:111:ILE:HG23	0.41	1.92	1	2
1:A:27:VAL:HG23	1:A:109:LEU:CD1	0.41	2.45	4	1
1:A:28:VAL:O	1:A:28:VAL:CG1	0.41	2.68	9	1
1:A:41:LEU:HD21	1:A:57:ILE:HG12	0.41	1.92	16	1
1:A:69:CYS:SG	1:A:71:VAL:HG22	0.41	2.55	7	1
1:A:101:SER:O	1:A:105:SER:N	0.41	2.53	15	2
1:A:41:LEU:HD11	1:A:78:VAL:HG21	0.41	1.92	16	1
1:A:28:VAL:HG22	1:A:108:THR:HG23	0.41	1.90	2	1
1:A:37:PHE:CZ	1:A:101:SER:HB2	0.41	2.50	16	2
1:A:31:ALA:CB	1:A:37:PHE:CE2	0.41	2.99	11	1
1:A:81:ILE:HB	1:A:97:LEU:HD13	0.41	1.92	11	1
1:A:42:GLN:NE2	1:A:58:SER:CB	0.41	2.84	8	1
1:A:26:GLU:O	1:A:26:GLU:CG	0.41	2.68	9	2
1:A:82:ASN:HB3	1:A:107:VAL:HG23	0.41	1.90	17	1
1:A:107:VAL:O	1:A:107:VAL:CG1	0.41	2.69	8	1
1:A:41:LEU:HD23	1:A:57:ILE:HG12	0.41	1.93	10	1
1:A:41:LEU:CD2	1:A:57:ILE:CG1	0.41	2.98	14	1
1:A:78:VAL:CG2	1:A:81:ILE:CD1	0.41	2.92	17	1
1:A:111:ILE:N	1:A:111:ILE:CD1	0.41	2.84	10	1
1:A:71:VAL:O	1:A:71:VAL:CG2	0.40	2.69	9	1
1:A:25:THR:CG2	1:A:111:ILE:CD1	0.40	2.99	4	1
1:A:64:SER:O	1:A:68:ARG:NE	0.40	2.54	15	1
1:A:27:VAL:CG1	1:A:71:VAL:HG11	0.40	2.46	16	1
1:A:61:GLU:O	1:A:67:GLU:CG	0.40	2.70	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	73/101 (72%)	58±2 (80±3%)	11±2 (14±3%)	4±1 (6±2%)	3	21
All	All	1460/2020 (72%)	1164 (80%)	211 (14%)	85 (6%)	3	21

All 10 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	63	ASP	20
1	A	22	THR	15
1	A	80	ALA	14
1	A	34	VAL	9
1	A	75	GLY	9
1	A	71	VAL	7
1	A	58	SER	5
1	A	79	MET	3
1	A	74	ILE	2
1	A	55	PRO	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	63/88 (72%)	43±3 (69±5%)	20±3 (31±5%)	1	14
All	All	1260/1760 (72%)	866 (69%)	394 (31%)	1	14

All 45 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	27	VAL	20
1	A	68	ARG	20
1	A	111	ILE	20
1	A	109	LEU	19
1	A	60	ILE	19
1	A	108	THR	18
1	A	40	GLN	16
1	A	41	LEU	14

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Mol	Chain	Res	Type	Models (Total)
1	A	63	ASP	14
1	A	69	CYS	14
1	A	26	GLU	13
1	A	61	GLU	13
1	A	32	ASP	10
1	A	56	LEU	10
1	A	100	ASP	10
1	A	79	MET	10
1	A	23	GLU	10
1	A	25	THR	9
1	A	96	GLN	9
1	A	73	GLN	9
1	A	22	THR	8
1	A	42	GLN	8
1	A	77	ARG	8
1	A	105	SER	8
1	A	64	SER	8
1	A	106	LYS	7
1	A	39	ILE	7
1	A	99	ARG	7
1	A	67	GLU	6
1	A	29	LEU	6
1	A	93	GLU	5
1	A	110	GLU	5
1	A	58	SER	5
1	A	72	LEU	5
1	A	98	LEU	5
1	A	102	SER	4
1	A	101	SER	3
1	A	76	ASP	3
1	A	112	GLU	2
1	A	24	THR	2
1	A	34	VAL	1
1	A	95	ASN	1
1	A	107	VAL	1
1	A	65	PRO	1
1	A	78	VAL	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 monosaccharides 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