



Full wwPDB NMR Structure Validation Report ⓘ

Mar 5, 2026 – 07:47 AM UTC

PDB ID : 2E4J / pdb_00002e4j
Title : Solution Structure of mouse Lipocalin-type Prostaglandin D Synthase
Authors : Shimamoto, S.; Ohkubo, T.
Deposited on : 2006-12-11

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

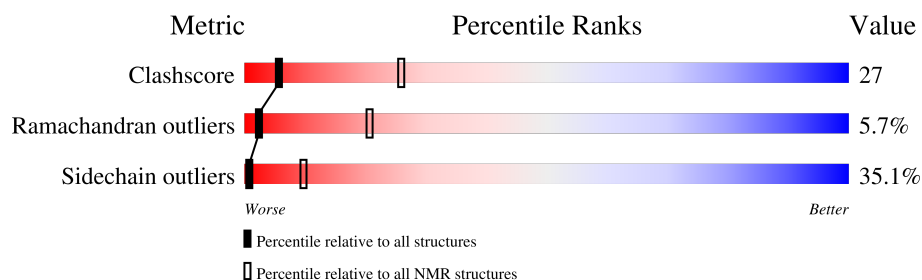
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

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



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

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

Mol	Chain	Length	Quality of chain
1	A	167	29% 46% 10% 15%

2 Ensemble composition and analysis

This entry contains 15 models. Model 1 is the overall representative, medoid model (most similar to other models).

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:30-A:84, A:88-A:108, A:114-A:137, A:142-A:183 (142)	0.92	1

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

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

The models can be grouped into 3 clusters and 5 single-model clusters were found.

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

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2592 atoms, of which 1282 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Prostaglandin-H2 D-isomerase.

Mol	Chain	Residues	Atoms						Trace
1	A	167	Total	C	H	N	O	S	0
			2592	828	1282	223	255	4	

There are 4 discrepancies between the modelled and reference sequences:

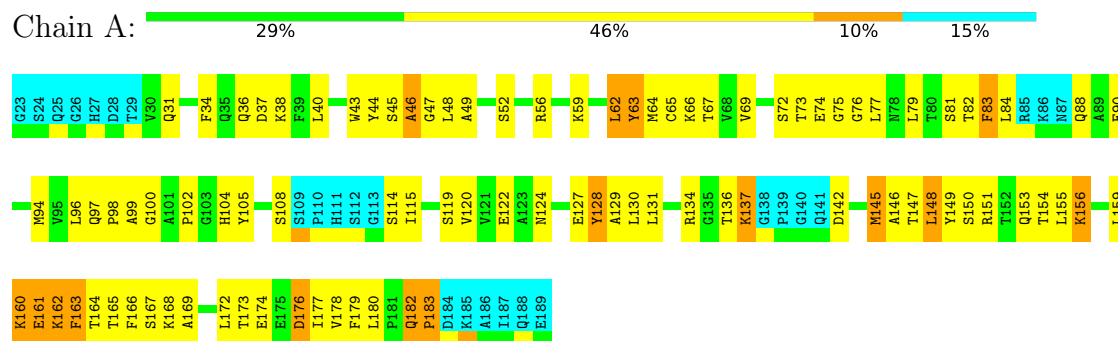
Chain	Residue	Modelled	Actual	Comment	Reference
A	23	GLY	-	expression tag	UNP O09114
A	24	SER	-	expression tag	UNP O09114
A	89	ALA	CYS	engineered mutation	UNP O09114
A	186	ALA	CYS	engineered mutation	UNP O09114

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: Prostaglandin-H2 D-isomerase

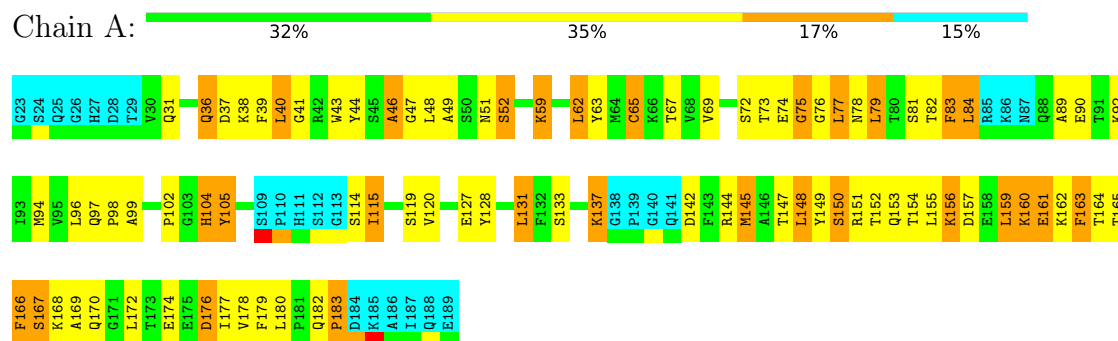


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

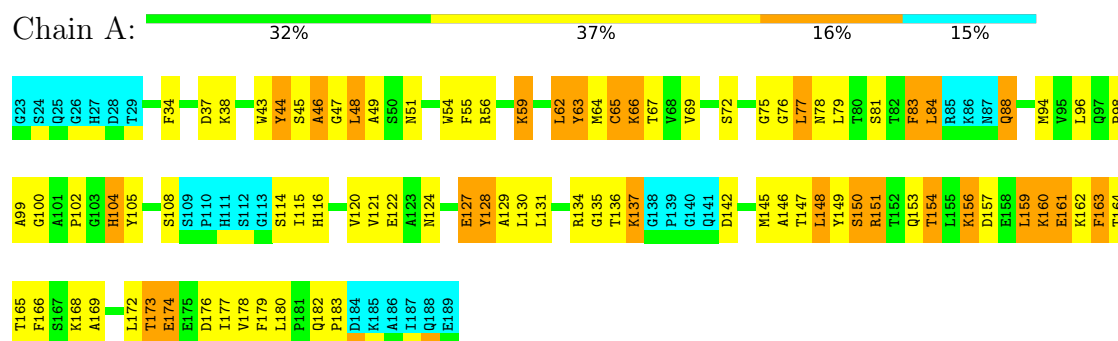
4.2.1 Score per residue for model 1 (medoid)

- Molecule 1: Prostaglandin-H2 D-isomerase



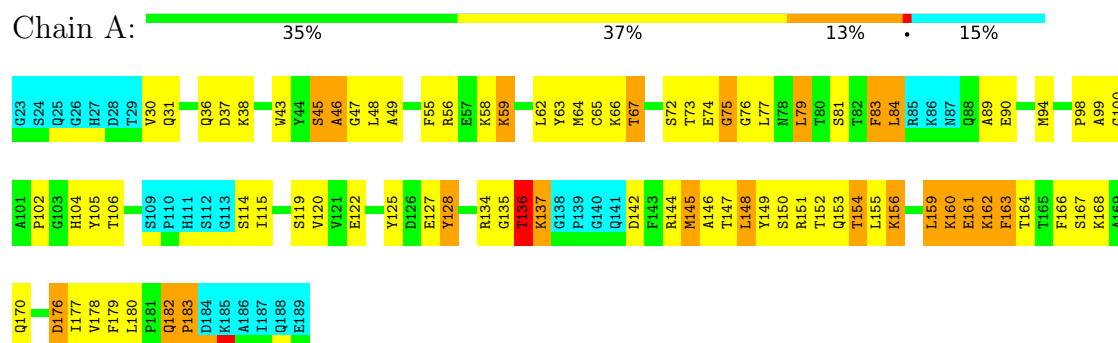
4.2.2 Score per residue for model 2

• Molecule 1: Prostaglandin-H2 D-isomerase



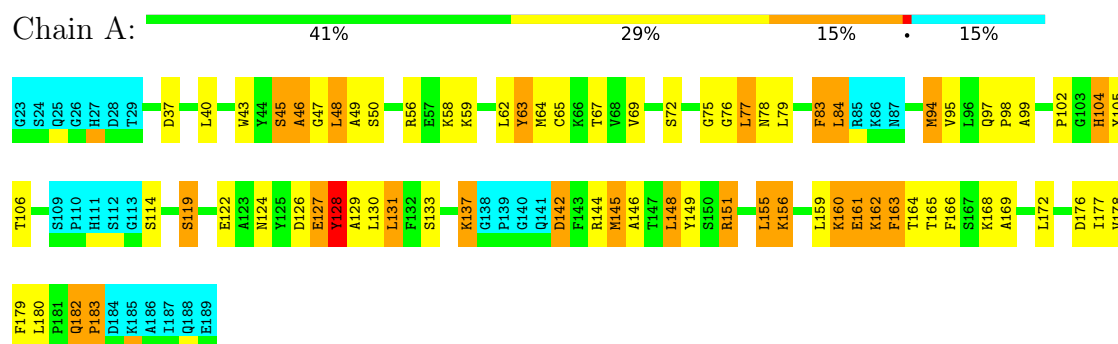
4.2.3 Score per residue for model 3

• Molecule 1: Prostaglandin-H2 D-isomerase



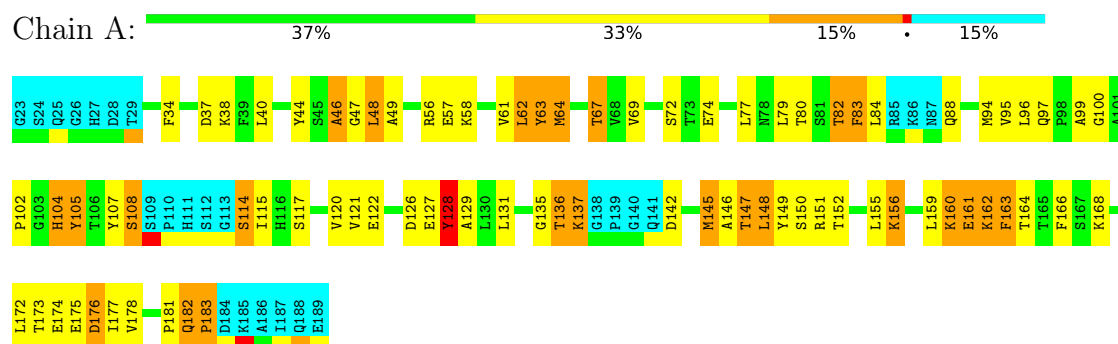
4.2.4 Score per residue for model 4

• Molecule 1: Prostaglandin-H2 D-isomerase



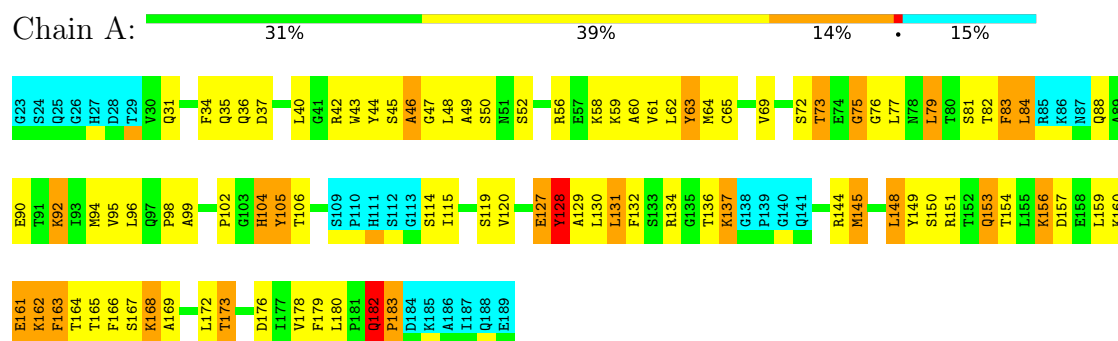
4.2.5 Score per residue for model 5

- Molecule 1: Prostaglandin-H2 D-isomerase



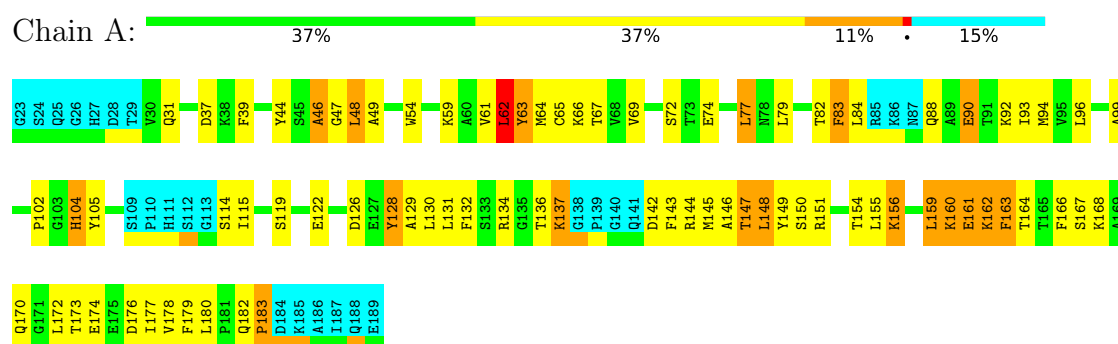
4.2.6 Score per residue for model 6

- Molecule 1: Prostaglandin-H2 D-isomerase



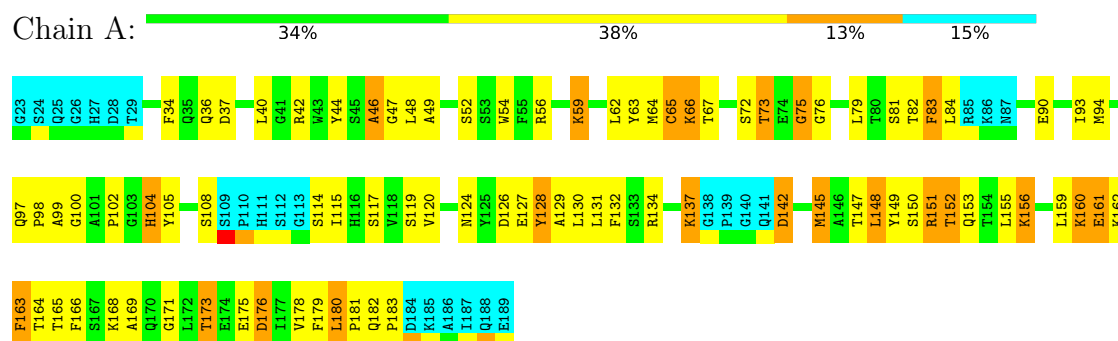
4.2.7 Score per residue for model 7

- Molecule 1: Prostaglandin-H2 D-isomerase



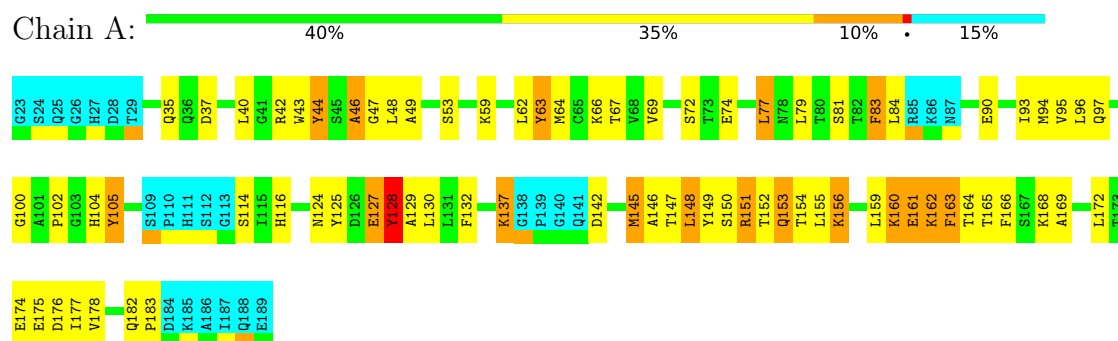
4.2.8 Score per residue for model 8

- Molecule 1: Prostaglandin-H2 D-isomerase



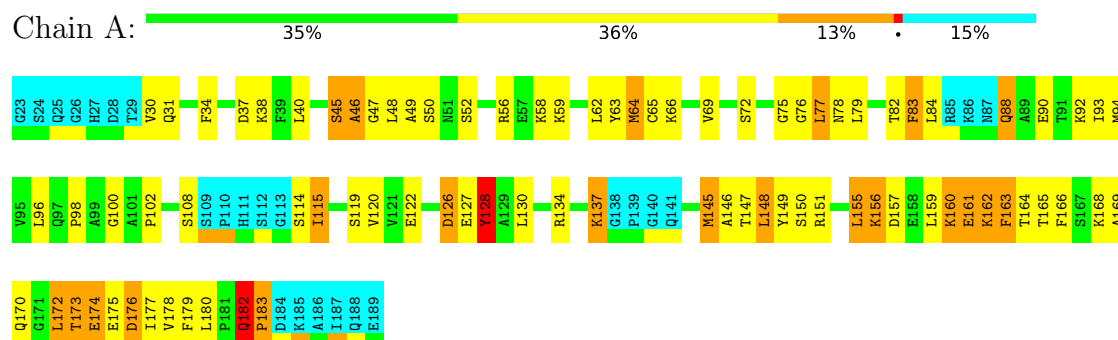
4.2.11 Score per residue for model 11

- Molecule 1: Prostaglandin-H2 D-isomerase



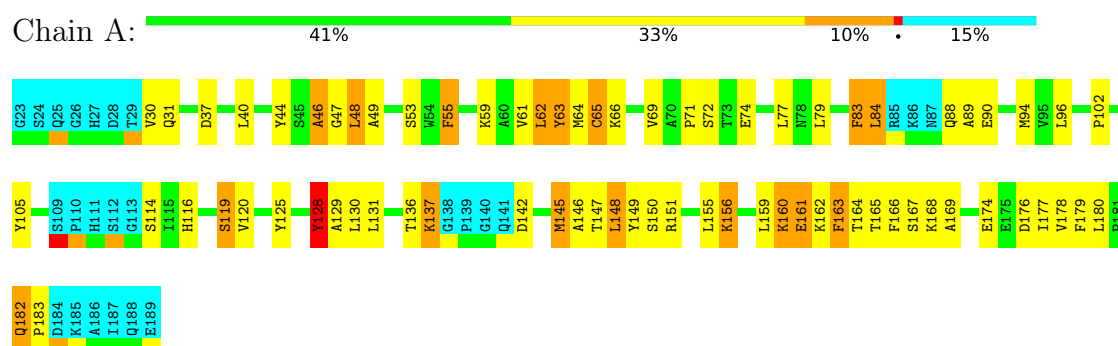
4.2.12 Score per residue for model 12

- Molecule 1: Prostaglandin-H2 D-isomerase



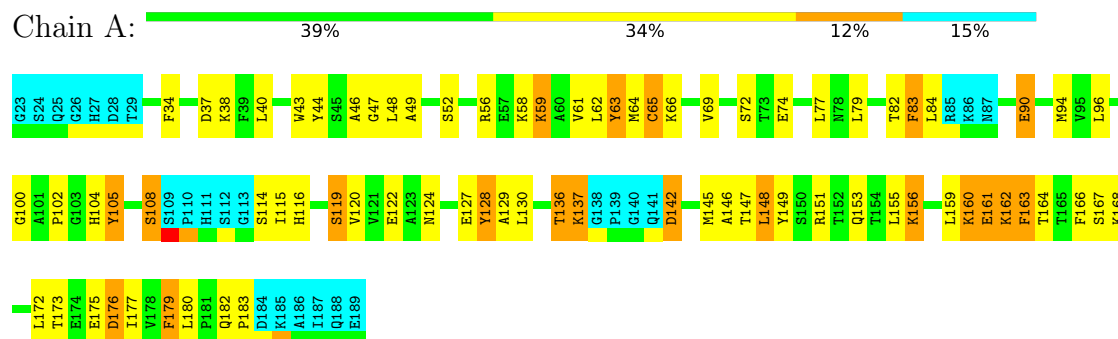
4.2.13 Score per residue for model 13

- Molecule 1: Prostaglandin-H2 D-isomerase



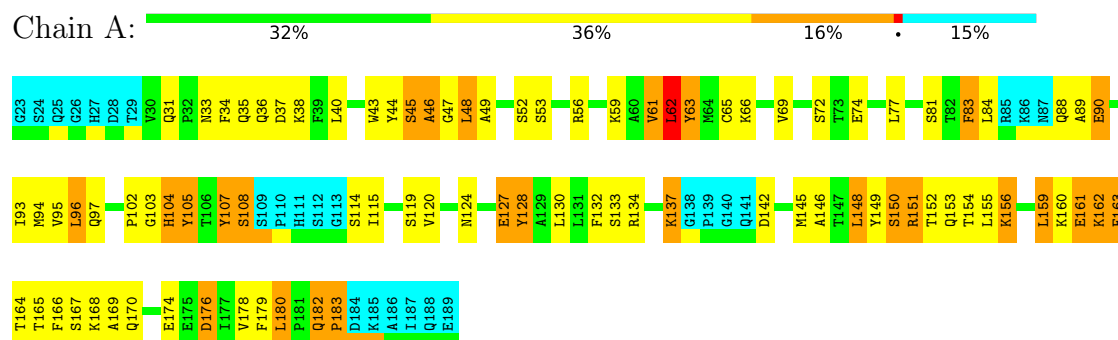
4.2.14 Score per residue for model 14

- Molecule 1: Prostaglandin-H2 D-isomerase



4.2.15 Score per residue for model 15

- Molecule 1: Prostaglandin-H2 D-isomerase



5 Refinement protocol and experimental data overview

The models were refined using the following method: *DISTANCE GEOMETRY, SIMULATED ANNEALING, TORSION ANGLE DYNAMICS*.

Of the 30 calculated structures, 15 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	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	1128	1114	1112	61±4
All	All	16920	16710	16680	917

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 27.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:148:LEU:HD21	1:A:159:LEU:HD13	1.09	1.20	5	13
1:A:148:LEU:HD11	1:A:159:LEU:HD22	0.94	1.36	12	8
1:A:159:LEU:HD12	1:A:160:LYS:N	0.93	1.79	13	10
1:A:84:LEU:HD13	1:A:89:ALA:HB2	0.85	1.48	15	3
1:A:164:THR:HG22	1:A:177:ILE:HD13	0.84	1.49	13	1
1:A:148:LEU:HD11	1:A:159:LEU:HD13	0.82	1.47	14	2
1:A:148:LEU:HD12	1:A:149:TYR:N	0.80	1.91	13	15
1:A:69:VAL:HG13	1:A:77:LEU:HD21	0.80	1.52	4	7
1:A:46:ALA:HB3	1:A:150:SER:OG	0.79	1.78	3	12
1:A:64:MET:HE1	1:A:66:LYS:CB	0.79	2.07	13	1
1:A:163:PHE:CE1	1:A:177:ILE:HD12	0.79	2.13	9	3
1:A:128:TYR:HB3	1:A:159:LEU:HD23	0.78	1.54	12	9
1:A:30:VAL:HG11	1:A:119:SER:OG	0.76	1.80	13	1
1:A:62:LEU:CD2	1:A:180:LEU:HD23	0.76	2.11	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:129:ALA:C	1:A:130:LEU:HD22	0.75	2.06	8	9
1:A:59:LYS:O	1:A:62:LEU:HD23	0.74	1.82	13	6
1:A:65:CYS:SG	1:A:180:LEU:HD13	0.73	2.23	12	2
1:A:108:SER:OG	1:A:115:ILE:HG23	0.73	1.82	14	2
1:A:148:LEU:CD2	1:A:159:LEU:HD13	0.72	2.12	12	6
1:A:129:ALA:C	1:A:130:LEU:HD12	0.72	2.10	6	1
1:A:59:LYS:HB2	1:A:62:LEU:HD12	0.71	1.61	3	3
1:A:62:LEU:HD12	1:A:62:LEU:O	0.70	1.84	8	5
1:A:148:LEU:HD21	1:A:159:LEU:CD1	0.70	2.17	8	7
1:A:108:SER:CB	1:A:115:ILE:HG23	0.70	2.16	14	4
1:A:148:LEU:CD1	1:A:159:LEU:HD13	0.69	2.17	14	2
1:A:179:PHE:CD1	1:A:179:PHE:N	0.69	2.58	14	1
1:A:34:PHE:CE2	1:A:120:VAL:HG11	0.69	2.22	14	6
1:A:82:THR:HG23	1:A:90:GLU:O	0.69	1.87	14	3
1:A:69:VAL:CG1	1:A:77:LEU:HD11	0.69	2.18	15	2
1:A:48:LEU:HD23	1:A:146:ALA:O	0.68	1.88	5	4
1:A:156:LYS:HD2	1:A:159:LEU:HD21	0.68	1.66	3	5
1:A:164:THR:HG23	1:A:177:ILE:HD11	0.68	1.65	10	2
1:A:168:LYS:NZ	1:A:173:THR:HG23	0.68	2.03	6	1
1:A:65:CYS:HB2	1:A:180:LEU:HD22	0.68	1.64	2	8
1:A:83:PHE:CD1	1:A:83:PHE:N	0.67	2.62	6	15
1:A:114:SER:HA	1:A:137:LYS:CB	0.67	2.19	13	3
1:A:64:MET:HE1	1:A:66:LYS:HB3	0.67	1.64	13	1
1:A:120:VAL:HA	1:A:131:LEU:HD23	0.67	1.67	8	2
1:A:108:SER:HA	1:A:115:ILE:HG23	0.67	1.64	9	1
1:A:67:THR:CG2	1:A:79:LEU:HD21	0.67	2.19	5	4
1:A:148:LEU:C	1:A:148:LEU:HD12	0.67	2.14	4	1
1:A:79:LEU:HD21	1:A:96:LEU:HD23	0.66	1.67	6	1
1:A:36:GLN:OE1	1:A:77:LEU:HD13	0.66	1.90	10	1
1:A:84:LEU:HD22	1:A:89:ALA:HB2	0.66	1.64	13	1
1:A:96:LEU:HD12	1:A:105:TYR:CD2	0.66	2.25	1	1
1:A:108:SER:HB2	1:A:115:ILE:HG22	0.66	1.67	2	1
1:A:49:ALA:HB1	1:A:176:ASP:O	0.66	1.91	11	15
1:A:131:LEU:HD22	1:A:131:LEU:N	0.66	2.06	6	2
1:A:176:ASP:O	1:A:177:ILE:HD13	0.66	1.90	3	4
1:A:148:LEU:CD1	1:A:159:LEU:HD22	0.66	2.18	12	1
1:A:130:LEU:HD11	1:A:163:PHE:HB2	0.65	1.67	10	3
1:A:114:SER:HA	1:A:137:LYS:HB2	0.65	1.69	13	3
1:A:67:THR:CG2	1:A:79:LEU:HD22	0.64	2.22	11	1
1:A:166:PHE:CD1	1:A:166:PHE:C	0.64	2.76	1	15
1:A:160:LYS:O	1:A:164:THR:HG22	0.64	1.92	12	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:135:GLY:O	1:A:136:THR:HG23	0.64	1.93	3	3
1:A:182:GLN:HB2	1:A:183:PRO:HD2	0.64	1.68	6	6
1:A:48:LEU:HD23	1:A:49:ALA:N	0.64	2.08	9	1
1:A:69:VAL:HG12	1:A:77:LEU:HD11	0.63	1.68	15	1
1:A:48:LEU:O	1:A:178:VAL:N	0.63	2.32	11	5
1:A:65:CYS:SG	1:A:180:LEU:HD22	0.62	2.34	8	2
1:A:84:LEU:CD1	1:A:89:ALA:HB2	0.62	2.22	15	2
1:A:159:LEU:O	1:A:163:PHE:HB3	0.62	1.95	7	15
1:A:62:LEU:HD13	1:A:181:PRO:HD3	0.62	1.70	8	1
1:A:64:MET:HE1	1:A:66:LYS:HG3	0.62	1.71	2	2
1:A:151:ARG:HD3	1:A:152:THR:HG23	0.62	1.71	8	1
1:A:65:CYS:HB2	1:A:180:LEU:HD13	0.62	1.72	14	1
1:A:67:THR:HG21	1:A:79:LEU:HD11	0.61	1.72	7	2
1:A:172:LEU:HD21	1:A:176:ASP:CG	0.61	2.21	2	7
1:A:168:LYS:HZ2	1:A:173:THR:HG23	0.61	1.56	6	1
1:A:64:MET:HE3	1:A:84:LEU:HD23	0.61	1.72	10	1
1:A:129:ALA:O	1:A:130:LEU:HD22	0.61	1.95	10	3
1:A:67:THR:CG2	1:A:79:LEU:HD11	0.61	2.25	7	2
1:A:166:PHE:CD1	1:A:167:SER:N	0.61	2.68	1	1
1:A:163:PHE:C	1:A:163:PHE:CD1	0.61	2.79	13	14
1:A:67:THR:HG22	1:A:79:LEU:HD22	0.61	1.73	11	2
1:A:143:PHE:CD2	1:A:145:MET:HE1	0.60	2.32	7	1
1:A:84:LEU:HD13	1:A:89:ALA:CB	0.60	2.25	15	1
1:A:148:LEU:HD11	1:A:159:LEU:CD1	0.59	2.26	14	1
1:A:76:GLY:HA2	1:A:98:PRO:CD	0.59	2.27	8	9
1:A:156:LYS:CE	1:A:159:LEU:HD21	0.59	2.28	2	2
1:A:55:PHE:CE1	1:A:178:VAL:HG23	0.58	2.34	13	1
1:A:64:MET:HE1	1:A:66:LYS:HB2	0.58	1.73	13	1
1:A:115:ILE:HD13	1:A:115:ILE:N	0.57	2.15	1	2
1:A:114:SER:HA	1:A:137:LYS:HA	0.57	1.76	15	10
1:A:159:LEU:HD12	1:A:159:LEU:C	0.57	2.24	14	7
1:A:96:LEU:C	1:A:96:LEU:HD23	0.57	2.24	15	1
1:A:108:SER:HB3	1:A:115:ILE:HG23	0.57	1.76	14	2
1:A:163:PHE:HE1	1:A:177:ILE:HD12	0.57	1.58	3	3
1:A:131:LEU:HB2	1:A:147:THR:HG23	0.57	1.77	7	2
1:A:173:THR:O	1:A:177:ILE:HD12	0.56	1.99	2	2
1:A:84:LEU:HD23	1:A:84:LEU:C	0.56	2.25	5	1
1:A:45:SER:OG	1:A:180:LEU:HD13	0.56	2.00	15	2
1:A:45:SER:CB	1:A:180:LEU:HD13	0.56	2.29	15	1
1:A:128:TYR:CB	1:A:159:LEU:HD23	0.56	2.30	4	3
1:A:114:SER:C	1:A:137:LYS:HG3	0.56	2.26	1	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:64:MET:HE3	1:A:84:LEU:CB	0.56	2.30	3	1
1:A:108:SER:HB2	1:A:115:ILE:HG23	0.56	1.77	10	1
1:A:47:GLY:HA3	1:A:148:LEU:HD23	0.56	1.78	14	13
1:A:154:THR:O	1:A:156:LYS:NZ	0.55	2.39	15	5
1:A:64:MET:HE1	1:A:66:LYS:CG	0.55	2.30	2	2
1:A:65:CYS:CB	1:A:180:LEU:HD22	0.55	2.31	15	1
1:A:84:LEU:HD22	1:A:89:ALA:CB	0.55	2.32	13	1
1:A:93:ILE:HD12	1:A:93:ILE:N	0.54	2.16	7	2
1:A:182:GLN:HB3	1:A:183:PRO:HD2	0.54	1.79	12	3
1:A:96:LEU:HD23	1:A:96:LEU:O	0.54	2.03	15	1
1:A:67:THR:HG23	1:A:79:LEU:HD21	0.54	1.78	7	2
1:A:161:GLU:HG2	1:A:162:LYS:N	0.54	2.18	1	15
1:A:30:VAL:HG11	1:A:119:SER:CB	0.54	2.32	13	1
1:A:84:LEU:CD2	1:A:89:ALA:HB2	0.54	2.31	13	1
1:A:69:VAL:HG22	1:A:79:LEU:HD21	0.54	1.78	11	1
1:A:96:LEU:HD12	1:A:105:TYR:HB2	0.54	1.79	6	2
1:A:150:SER:HA	1:A:156:LYS:HE3	0.53	1.78	10	3
1:A:67:THR:HG21	1:A:79:LEU:HD21	0.53	1.78	4	3
1:A:155:LEU:HD13	1:A:179:PHE:CE2	0.53	2.38	12	2
1:A:154:THR:O	1:A:156:LYS:HE2	0.53	2.03	3	1
1:A:64:MET:HB3	1:A:182:GLN:O	0.53	2.04	12	4
1:A:49:ALA:O	1:A:145:MET:HA	0.53	2.03	14	15
1:A:148:LEU:HD12	1:A:148:LEU:C	0.53	2.29	12	8
1:A:47:GLY:HA3	1:A:148:LEU:CD2	0.53	2.34	4	5
1:A:77:LEU:O	1:A:95:VAL:HG23	0.53	2.04	6	4
1:A:56:ARG:O	1:A:60:ALA:HB2	0.52	2.04	6	1
1:A:128:TYR:N	1:A:128:TYR:CD1	0.52	2.78	14	2
1:A:62:LEU:C	1:A:62:LEU:HD12	0.52	2.30	12	1
1:A:164:THR:CG2	1:A:177:ILE:HD13	0.52	2.29	13	1
1:A:96:LEU:HD11	1:A:105:TYR:CD2	0.52	2.40	15	1
1:A:115:ILE:HD13	1:A:115:ILE:H	0.52	1.63	1	2
1:A:65:CYS:CB	1:A:180:LEU:HD13	0.52	2.33	14	1
1:A:120:VAL:HA	1:A:131:LEU:HD12	0.52	1.81	6	2
1:A:49:ALA:HB2	1:A:163:PHE:CZ	0.52	2.40	9	12
1:A:49:ALA:HB3	1:A:146:ALA:HB3	0.52	1.82	10	9
1:A:182:GLN:O	1:A:183:PRO:O	0.51	2.27	15	1
1:A:163:PHE:CD1	1:A:163:PHE:C	0.51	2.88	1	1
1:A:163:PHE:HA	1:A:166:PHE:CD2	0.51	2.40	1	1
1:A:163:PHE:CD1	1:A:164:THR:N	0.51	2.78	1	15
1:A:182:GLN:CB	1:A:183:PRO:HD2	0.51	2.35	15	3
1:A:79:LEU:HD22	1:A:96:LEU:HD12	0.51	1.82	12	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:45:SER:HB3	1:A:180:LEU:HD13	0.51	1.82	15	2
1:A:151:ARG:H	1:A:156:LYS:CD	0.51	2.18	1	2
1:A:61:VAL:HG21	1:A:63:TYR:CE2	0.51	2.41	15	1
1:A:150:SER:CA	1:A:156:LYS:HE3	0.50	2.36	10	3
1:A:149:TYR:O	1:A:159:LEU:HD22	0.50	2.06	14	1
1:A:129:ALA:HB3	1:A:149:TYR:HB2	0.50	1.83	6	5
1:A:96:LEU:CD1	1:A:105:TYR:CD2	0.50	2.93	15	1
1:A:182:GLN:CB	1:A:183:PRO:CD	0.50	2.88	5	4
1:A:156:LYS:HZ1	1:A:156:LYS:H	0.50	1.49	9	1
1:A:156:LYS:CE	1:A:159:LEU:HD23	0.50	2.37	11	2
1:A:95:VAL:O	1:A:96:LEU:HD22	0.50	2.06	5	1
1:A:164:THR:HG23	1:A:177:ILE:CD1	0.50	2.36	10	3
1:A:43:TRP:CZ3	1:A:127:GLU:O	0.50	2.65	14	3
1:A:59:LYS:CB	1:A:62:LEU:HD12	0.50	2.34	3	1
1:A:114:SER:HA	1:A:137:LYS:CA	0.49	2.37	11	5
1:A:63:TYR:CD1	1:A:83:PHE:CB	0.49	2.96	4	3
1:A:152:THR:OG1	1:A:156:LYS:HG3	0.49	2.07	15	2
1:A:156:LYS:N	1:A:156:LYS:HE2	0.49	2.23	1	2
1:A:128:TYR:CD1	1:A:156:LYS:HD2	0.49	2.43	2	3
1:A:61:VAL:O	1:A:63:TYR:CD2	0.49	2.65	13	4
1:A:148:LEU:HD11	1:A:159:LEU:CD2	0.49	2.25	12	1
1:A:67:THR:HG23	1:A:79:LEU:HD12	0.49	1.84	2	1
1:A:59:LYS:HE3	1:A:178:VAL:HG23	0.49	1.84	3	1
1:A:69:VAL:CG1	1:A:77:LEU:HD21	0.48	2.33	4	1
1:A:167:SER:OG	1:A:172:LEU:HD23	0.48	2.08	14	1
1:A:148:LEU:HD21	1:A:159:LEU:HB2	0.48	1.85	3	2
1:A:156:LYS:HE2	1:A:156:LYS:H	0.48	1.67	10	3
1:A:49:ALA:CB	1:A:176:ASP:O	0.48	2.62	1	13
1:A:48:LEU:HA	1:A:146:ALA:O	0.48	2.09	10	7
1:A:156:LYS:O	1:A:159:LEU:CD1	0.48	2.62	4	8
1:A:30:VAL:HG12	1:A:120:VAL:O	0.48	2.09	3	2
1:A:64:MET:HE3	1:A:84:LEU:HB2	0.48	1.84	3	1
1:A:61:VAL:CG2	1:A:63:TYR:CE2	0.47	2.97	15	1
1:A:99:ALA:HB3	1:A:104:HIS:O	0.47	2.09	3	8
1:A:155:LEU:HD12	1:A:155:LEU:C	0.47	2.35	7	1
1:A:121:VAL:HG23	1:A:131:LEU:C	0.47	2.34	2	1
1:A:114:SER:HA	1:A:137:LYS:HG3	0.47	1.85	14	8
1:A:114:SER:CA	1:A:137:LYS:HA	0.47	2.40	5	5
1:A:34:PHE:CE2	1:A:120:VAL:HG21	0.47	2.44	12	1
1:A:128:TYR:CD1	1:A:128:TYR:N	0.47	2.82	15	3
1:A:76:GLY:HA2	1:A:98:PRO:HD2	0.47	1.86	8	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:83:PHE:CD2	1:A:90:GLU:CB	0.47	2.98	15	2
1:A:128:TYR:CD1	1:A:156:LYS:NZ	0.47	2.83	13	2
1:A:181:PRO:O	1:A:182:GLN:O	0.46	2.33	5	1
1:A:174:GLU:HA	1:A:177:ILE:HD12	0.46	1.87	12	2
1:A:108:SER:CA	1:A:115:ILE:HG23	0.46	2.37	9	1
1:A:156:LYS:HE2	1:A:156:LYS:N	0.46	2.25	3	1
1:A:64:MET:HB2	1:A:84:LEU:CB	0.46	2.40	6	2
1:A:156:LYS:HE2	1:A:159:LEU:HD21	0.46	1.87	10	1
1:A:46:ALA:O	1:A:179:PHE:CD2	0.46	2.68	3	12
1:A:47:GLY:O	1:A:148:LEU:N	0.46	2.48	15	12
1:A:64:MET:HE3	1:A:183:PRO:HA	0.46	1.87	14	1
1:A:43:TRP:CH2	1:A:127:GLU:HA	0.46	2.46	6	5
1:A:156:LYS:HZ1	1:A:159:LEU:HD23	0.46	1.70	13	2
1:A:34:PHE:CE2	1:A:120:VAL:CG1	0.46	2.99	14	1
1:A:44:TYR:CZ	1:A:183:PRO:HG3	0.46	2.45	10	2
1:A:77:LEU:N	1:A:96:LEU:O	0.46	2.49	5	7
1:A:45:SER:O	1:A:46:ALA:HB2	0.46	2.11	12	3
1:A:156:LYS:N	1:A:156:LYS:CE	0.45	2.79	1	2
1:A:178:VAL:HG13	1:A:178:VAL:O	0.45	2.11	10	9
1:A:150:SER:HB3	1:A:156:LYS:CE	0.45	2.40	3	1
1:A:150:SER:CB	1:A:156:LYS:NZ	0.45	2.79	5	1
1:A:114:SER:OG	1:A:137:LYS:HD3	0.45	2.11	11	3
1:A:151:ARG:H	1:A:156:LYS:HD3	0.45	1.71	1	2
1:A:130:LEU:HD22	1:A:130:LEU:N	0.45	2.25	14	2
1:A:44:TYR:CZ	1:A:183:PRO:CG	0.45	3.00	10	1
1:A:128:TYR:CD2	1:A:151:ARG:CB	0.45	3.00	4	2
1:A:48:LEU:HB3	1:A:178:VAL:HG13	0.45	1.87	9	1
1:A:151:ARG:CB	1:A:156:LYS:HD3	0.45	2.41	6	3
1:A:44:TYR:CD1	1:A:44:TYR:N	0.45	2.84	11	5
1:A:40:LEU:HD23	1:A:41:GLY:N	0.45	2.27	1	1
1:A:44:TYR:CE2	1:A:183:PRO:CG	0.45	2.99	9	1
1:A:156:LYS:NZ	1:A:159:LEU:HD23	0.45	2.27	13	2
1:A:157:ASP:HA	1:A:160:LYS:HB3	0.45	1.87	12	4
1:A:64:MET:O	1:A:83:PHE:HA	0.45	2.11	7	2
1:A:62:LEU:HD12	1:A:62:LEU:C	0.45	2.36	8	1
1:A:124:ASN:O	1:A:128:TYR:CD1	0.45	2.70	4	1
1:A:67:THR:HG23	1:A:79:LEU:CD2	0.45	2.41	7	2
1:A:108:SER:CB	1:A:115:ILE:HG22	0.45	2.42	5	2
1:A:173:THR:HG22	1:A:174:GLU:H	0.44	1.72	2	2
1:A:49:ALA:O	1:A:145:MET:CA	0.44	2.65	13	3
1:A:63:TYR:CE1	1:A:83:PHE:CD2	0.44	3.05	4	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:159:LEU:HD12	1:A:160:LYS:H	0.44	1.67	13	2
1:A:121:VAL:HG23	1:A:131:LEU:CA	0.44	2.42	2	1
1:A:119:SER:O	1:A:131:LEU:HD23	0.44	2.12	4	1
1:A:154:THR:O	1:A:156:LYS:CE	0.44	2.65	3	1
1:A:128:TYR:CD2	1:A:156:LYS:HG2	0.44	2.46	4	1
1:A:148:LEU:HD13	1:A:159:LEU:HD13	0.44	1.90	4	1
1:A:64:MET:HG3	1:A:182:GLN:O	0.44	2.13	11	2
1:A:64:MET:HB2	1:A:84:LEU:HB3	0.44	1.88	10	1
1:A:47:GLY:HA3	1:A:148:LEU:CG	0.44	2.42	9	4
1:A:83:PHE:CD2	1:A:90:GLU:HB3	0.44	2.48	15	2
1:A:152:THR:OG1	1:A:156:LYS:CE	0.44	2.66	1	2
1:A:152:THR:O	1:A:153:GLN:C	0.44	2.60	11	1
1:A:159:LEU:CD1	1:A:159:LEU:C	0.44	2.90	14	1
1:A:42:ARG:NH1	1:A:44:TYR:CE1	0.44	2.85	10	1
1:A:103:GLY:CA	1:A:120:VAL:HG12	0.44	2.43	15	1
1:A:83:PHE:O	1:A:89:ALA:HA	0.43	2.13	10	2
1:A:124:ASN:O	1:A:128:TYR:CD2	0.43	2.71	14	1
1:A:152:THR:CG2	1:A:156:LYS:CG	0.43	2.96	1	1
1:A:67:THR:CG2	1:A:79:LEU:HD12	0.43	2.42	2	1
1:A:99:ALA:CB	1:A:104:HIS:O	0.43	2.67	3	5
1:A:64:MET:HG3	1:A:84:LEU:HD23	0.43	1.89	9	1
1:A:96:LEU:HD11	1:A:105:TYR:CE2	0.43	2.49	15	1
1:A:64:MET:O	1:A:83:PHE:HB3	0.43	2.14	12	2
1:A:69:VAL:HG13	1:A:77:LEU:CD2	0.43	2.42	15	1
1:A:127:GLU:O	1:A:129:ALA:N	0.43	2.52	5	1
1:A:156:LYS:HB2	1:A:159:LEU:CD2	0.43	2.44	8	3
1:A:44:TYR:CD2	1:A:151:ARG:O	0.43	2.71	8	1
1:A:155:LEU:HD12	1:A:155:LEU:O	0.43	2.13	11	1
1:A:73:THR:O	1:A:75:GLY:N	0.43	2.52	6	6
1:A:128:TYR:CE1	1:A:151:ARG:HB3	0.43	2.48	1	2
1:A:121:VAL:HG12	1:A:122:GLU:HG2	0.43	1.91	5	1
1:A:127:GLU:O	1:A:128:TYR:C	0.43	2.61	5	1
1:A:177:ILE:HG22	1:A:179:PHE:CE1	0.43	2.49	14	1
1:A:79:LEU:N	1:A:94:MET:O	0.43	2.51	4	1
1:A:44:TYR:CE2	1:A:183:PRO:HG3	0.43	2.49	6	2
1:A:155:LEU:HD13	1:A:179:PHE:HE2	0.43	1.74	12	1
1:A:130:LEU:HD12	1:A:163:PHE:HB2	0.43	1.89	14	1
1:A:182:GLN:O	1:A:183:PRO:C	0.43	2.62	1	1
1:A:44:TYR:CD2	1:A:183:PRO:HG3	0.43	2.48	14	1
1:A:150:SER:OG	1:A:156:LYS:HE3	0.42	2.14	1	2
1:A:84:LEU:C	1:A:84:LEU:CD2	0.42	2.92	5	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:46:ALA:CB	1:A:150:SER:OG	0.42	2.67	7	1
1:A:165:THR:O	1:A:169:ALA:CB	0.42	2.67	4	11
1:A:44:TYR:O	1:A:150:SER:N	0.42	2.48	2	3
1:A:63:TYR:CD1	1:A:83:PHE:HB3	0.42	2.49	2	2
1:A:115:ILE:O	1:A:115:ILE:HD12	0.42	2.14	2	1
1:A:161:GLU:CG	1:A:162:LYS:N	0.42	2.82	1	2
1:A:69:VAL:HG22	1:A:79:LEU:CD1	0.42	2.45	4	1
1:A:156:LYS:HB2	1:A:159:LEU:HD21	0.42	1.91	9	2
1:A:64:MET:CG	1:A:182:GLN:O	0.42	2.68	4	2
1:A:178:VAL:O	1:A:178:VAL:HG13	0.42	2.14	8	3
1:A:115:ILE:N	1:A:115:ILE:CD1	0.42	2.82	1	2
1:A:173:THR:HG23	1:A:175:GLU:H	0.42	1.73	8	3
1:A:34:PHE:CD2	1:A:120:VAL:HG21	0.42	2.48	14	1
1:A:83:PHE:CD2	1:A:90:GLU:HB2	0.42	2.50	7	1
1:A:67:THR:HG22	1:A:79:LEU:CD2	0.42	2.44	10	1
1:A:128:TYR:CE1	1:A:156:LYS:CD	0.42	3.02	15	1
1:A:128:TYR:CG	1:A:159:LEU:HD23	0.42	2.50	15	2
1:A:150:SER:OG	1:A:156:LYS:NZ	0.42	2.52	5	1
1:A:46:ALA:O	1:A:179:PHE:CE2	0.42	2.73	8	1
1:A:54:TRP:CD1	1:A:57:GLU:OE2	0.42	2.73	10	1
1:A:104:HIS:C	1:A:105:TYR:CD1	0.42	2.98	10	4
1:A:182:GLN:HB2	1:A:183:PRO:CD	0.42	2.44	10	1
1:A:152:THR:HG22	1:A:153:GLN:N	0.42	2.30	11	2
1:A:160:LYS:HG2	1:A:179:PHE:CE2	0.42	2.50	3	1
1:A:59:LYS:HD2	1:A:62:LEU:HD23	0.42	1.92	8	1
1:A:95:VAL:O	1:A:107:TYR:CE1	0.42	2.72	15	1
1:A:54:TRP:CD1	1:A:54:TRP:C	0.41	2.98	9	1
1:A:34:PHE:CD1	1:A:35:GLN:N	0.41	2.88	15	1
1:A:62:LEU:CD1	1:A:181:PRO:CG	0.41	2.97	8	1
1:A:156:LYS:O	1:A:159:LEU:HG	0.41	2.16	11	1
1:A:34:PHE:HE2	1:A:120:VAL:HG21	0.41	1.75	12	1
1:A:173:THR:O	1:A:177:ILE:CD1	0.41	2.68	5	2
1:A:39:PHE:CZ	1:A:125:TYR:O	0.41	2.74	9	1
1:A:71:PRO:HA	1:A:77:LEU:HD12	0.41	1.91	13	1
1:A:130:LEU:C	1:A:131:LEU:HD13	0.41	2.40	6	1
1:A:83:PHE:CE2	1:A:90:GLU:HB3	0.41	2.51	7	2
1:A:83:PHE:CE1	1:A:92:LYS:HB2	0.41	2.51	6	1
1:A:128:TYR:CE1	1:A:156:LYS:HD2	0.41	2.50	11	1
1:A:120:VAL:O	1:A:120:VAL:HG23	0.41	2.14	13	1
1:A:114:SER:N	1:A:137:LYS:HA	0.41	2.31	11	2
1:A:133:SER:OG	1:A:145:MET:HB2	0.41	2.16	1	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:46:ALA:HB3	1:A:150:SER:HG	0.41	1.74	3	1
1:A:67:THR:HG23	1:A:79:LEU:HD13	0.41	1.93	8	1
1:A:67:THR:CG2	1:A:79:LEU:CD1	0.41	2.99	1	1
1:A:151:ARG:O	1:A:151:ARG:HG3	0.41	2.15	5	1
1:A:115:ILE:HG22	1:A:136:THR:O	0.41	2.16	6	1
1:A:130:LEU:HD12	1:A:130:LEU:N	0.41	2.30	6	1
1:A:127:GLU:C	1:A:128:TYR:CG	0.41	2.98	8	1
1:A:69:VAL:HA	1:A:79:LEU:HD23	0.41	1.92	11	1
1:A:172:LEU:CD2	1:A:176:ASP:CB	0.41	2.99	12	1
1:A:128:TYR:O	1:A:130:LEU:HD12	0.41	2.16	15	1
1:A:126:ASP:O	1:A:128:TYR:CZ	0.41	2.74	8	1
1:A:36:GLN:NE2	1:A:77:LEU:CD1	0.40	2.84	1	1
1:A:163:PHE:CE1	1:A:177:ILE:HG23	0.40	2.51	3	2
1:A:61:VAL:O	1:A:62:LEU:C	0.40	2.63	5	2
1:A:39:PHE:CD1	1:A:39:PHE:N	0.40	2.89	7	2
1:A:104:HIS:CD2	1:A:119:SER:OG	0.40	2.74	14	1
1:A:63:TYR:HA	1:A:84:LEU:O	0.40	2.16	5	1
1:A:46:ALA:HB3	1:A:155:LEU:HB3	0.40	1.93	4	1
1:A:67:THR:HG22	1:A:79:LEU:HD11	0.40	1.93	4	1
1:A:128:TYR:CD2	1:A:151:ARG:HB2	0.40	2.51	4	1
1:A:172:LEU:C	1:A:173:THR:OG1	0.40	2.65	6	2
1:A:64:MET:CB	1:A:84:LEU:O	0.40	2.70	8	1
1:A:49:ALA:O	1:A:146:ALA:N	0.40	2.53	13	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	142/167 (85%)	115±3 (81±2%)	18±3 (13±2%)	8±2 (6±1%)	2	21
All	All	2130/2505 (85%)	1732 (81%)	276 (13%)	122 (6%)	2	21

All 24 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	46	ALA	15
1	A	102	PRO	15
1	A	183	PRO	11
1	A	128	TYR	10
1	A	74	GLU	9
1	A	75	GLY	9
1	A	100	GLY	8
1	A	151	ARG	6
1	A	182	GLN	5
1	A	52	SER	4
1	A	136	THR	4
1	A	142	ASP	4
1	A	88	GLN	3
1	A	62	LEU	3
1	A	126	ASP	3
1	A	127	GLU	2
1	A	64	MET	2
1	A	153	GLN	2
1	A	173	THR	2
1	A	144	ARG	1
1	A	171	GLY	1
1	A	124	ASN	1
1	A	53	SER	1
1	A	33	ASN	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	123/142 (87%)	80±6 (65±5%)	43±6 (35±5%)	1	10
All	All	1845/2130 (87%)	1198 (65%)	647 (35%)	1	10

All 97 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	37	ASP	15
1	A	63	TYR	15

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Mol	Chain	Res	Type	Models (Total)
1	A	72	SER	15
1	A	83	PHE	15
1	A	137	LYS	15
1	A	148	LEU	15
1	A	156	LYS	15
1	A	161	GLU	15
1	A	163	PHE	15
1	A	168	LYS	15
1	A	94	MET	14
1	A	160	LYS	14
1	A	105	TYR	13
1	A	40	LEU	12
1	A	48	LEU	12
1	A	142	ASP	12
1	A	147	THR	12
1	A	128	TYR	12
1	A	90	GLU	11
1	A	119	SER	11
1	A	155	LEU	10
1	A	56	ARG	10
1	A	162	LYS	10
1	A	81	SER	9
1	A	145	MET	9
1	A	66	LYS	9
1	A	88	GLN	9
1	A	31	GLN	8
1	A	38	LYS	8
1	A	59	LYS	8
1	A	97	GLN	8
1	A	104	HIS	8
1	A	153	GLN	8
1	A	174	GLU	8
1	A	176	ASP	8
1	A	134	ARG	8
1	A	62	LEU	7
1	A	77	LEU	7
1	A	167	SER	7
1	A	122	GLU	7
1	A	151	ARG	7
1	A	58	LYS	7
1	A	182	GLN	7
1	A	36	GLN	6

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Mol	Chain	Res	Type	Models (Total)
1	A	65	CYS	6
1	A	79	LEU	6
1	A	84	LEU	6
1	A	144	ARG	6
1	A	159	LEU	6
1	A	154	THR	6
1	A	45	SER	6
1	A	52	SER	5
1	A	92	LYS	5
1	A	170	GLN	5
1	A	124	ASN	5
1	A	127	GLU	5
1	A	44	TYR	4
1	A	78	ASN	4
1	A	82	THR	4
1	A	54	TRP	4
1	A	116	HIS	4
1	A	173	THR	4
1	A	106	THR	4
1	A	136	THR	4
1	A	50	SER	4
1	A	117	SER	4
1	A	132	PHE	4
1	A	93	ILE	4
1	A	51	ASN	3
1	A	115	ILE	3
1	A	131	LEU	3
1	A	150	SER	3
1	A	55	PHE	3
1	A	67	THR	3
1	A	125	TYR	3
1	A	126	ASP	3
1	A	108	SER	3
1	A	172	LEU	3
1	A	175	GLU	3
1	A	35	GLN	3
1	A	42	ARG	3
1	A	73	THR	3
1	A	180	LEU	3
1	A	53	SER	3
1	A	133	SER	2
1	A	57	GLU	2

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Mol	Chain	Res	Type	Models (Total)
1	A	107	TYR	2
1	A	69	VAL	2
1	A	64	MET	2
1	A	166	PHE	1
1	A	80	THR	1
1	A	114	SER	1
1	A	152	THR	1
1	A	157	ASP	1
1	A	179	PHE	1
1	A	61	VAL	1
1	A	96	LEU	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 oligosaccharides 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