



Full wwPDB NMR Structure Validation Report ⓘ

Mar 12, 2026 – 02:22 PM UTC

PDB ID : 1P4Q / pdb_00001p4q
Title : Solution structure of the CITED2 transactivation domain in complex with the p300 CH1 domain
Authors : Freedman, S.J.; Sun, Z.-Y.J.; Kung, A.L.; France, D.S.; Wagner, G.; Eck, M.J.
Deposited on : 2003-04-23

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

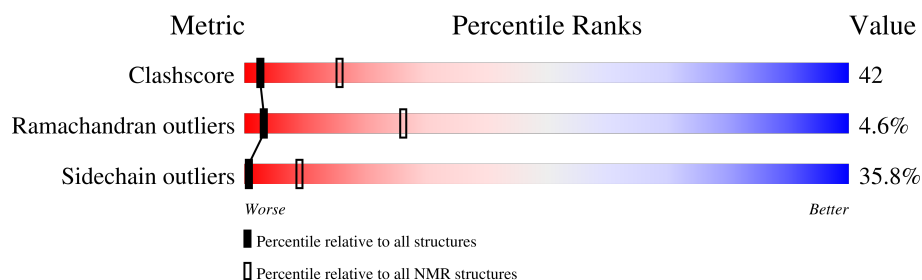
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

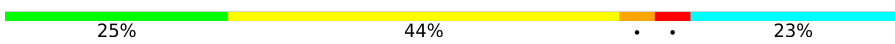
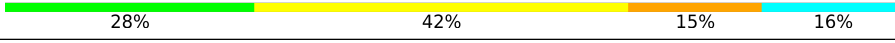
The overall completeness of chemical shifts assignment was not calculated.

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	52	 25% 44% . . 23%
2	B	101	 28% 42% 15% 16%

2 Ensemble composition and analysis

This entry contains 17 models. Model 12 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:10-A:49, B:109-B:184, B:189-B:197 (125)	0.77	12

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 5 clusters and 3 single-model clusters were found.

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

3 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2371 atoms, of which 1172 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Cbp/p300-interacting transactivator 2.

Mol	Chain	Residues	Atoms						Trace
1	A	52	Total	C	H	N	O	S	0
			788	256	382	60	87	3	

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	cloning artifact	UNP Q99967
A	2	SER	-	cloning artifact	UNP Q99967
A	3	GLY	-	cloning artifact	UNP Q99967
A	4	SER	-	cloning artifact	UNP Q99967
A	5	GLY	-	cloning artifact	UNP Q99967
A	6	SER	-	cloning artifact	UNP Q99967
A	7	GLY	-	cloning artifact	UNP Q99967
A	8	SER	-	cloning artifact	UNP Q99967

- Molecule 2 is a protein called E1A-associated protein p300.

Mol	Chain	Residues	Atoms						Trace
2	B	101	Total	C	H	N	O	S	0
			1580	475	790	164	139	12	

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

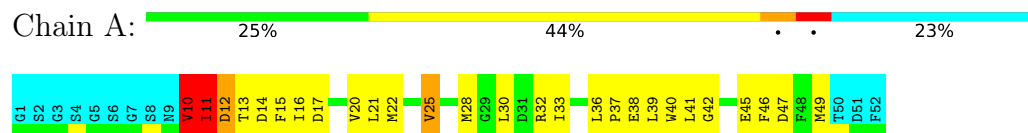
Mol	Chain	Residues	Atoms	
3	B	3	Total	Zn
			3	3

4 Residue-property plots

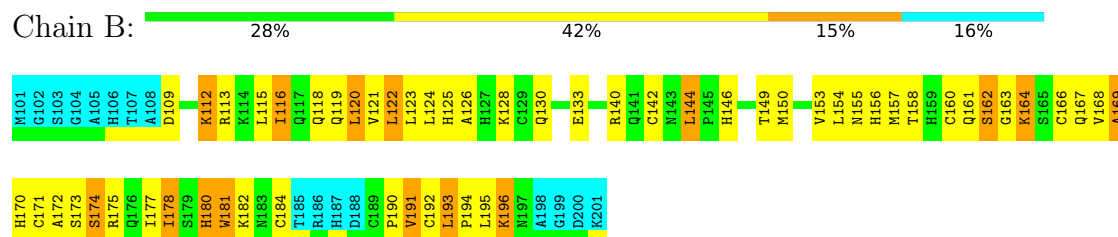
4.1 Average score per residue in the NMR ensemble

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: Cbp/p300-interacting transactivator 2



- Molecule 2: E1A-associated protein p300

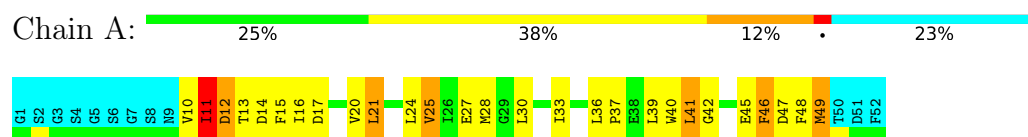


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

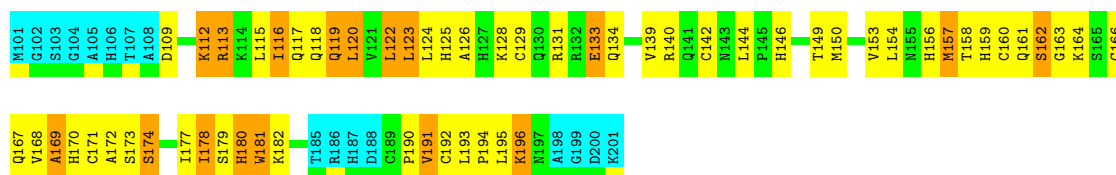
4.2.1 Score per residue for model 1

- Molecule 1: Cbp/p300-interacting transactivator 2



- Molecule 2: E1A-associated protein p300



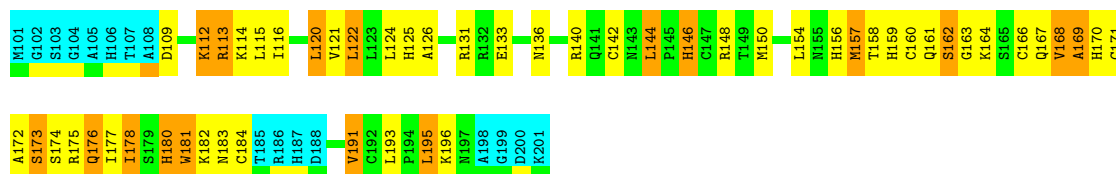


4.2.2 Score per residue for model 2

- Molecule 1: Cbp/p300-interacting transactivator 2



- Molecule 2: E1A-associated protein p300

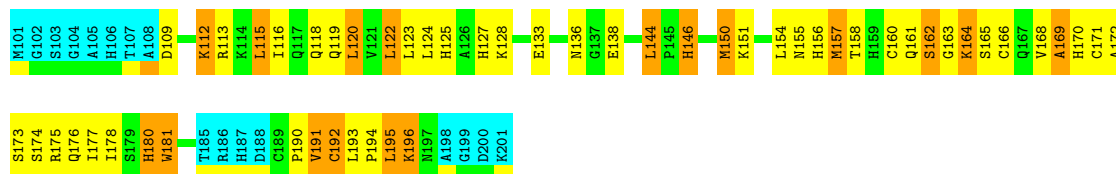
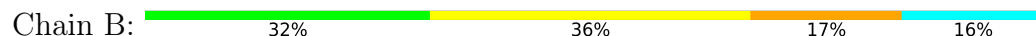


4.2.3 Score per residue for model 3

- Molecule 1: Cbp/p300-interacting transactivator 2

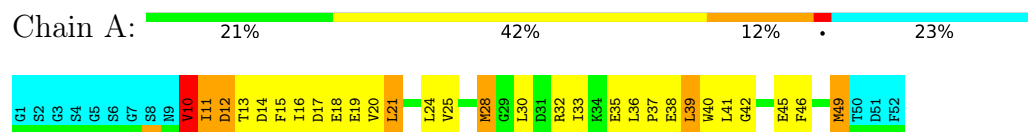


- Molecule 2: E1A-associated protein p300

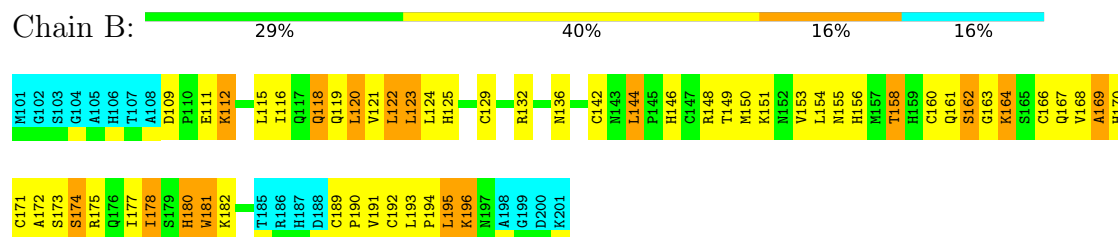


4.2.4 Score per residue for model 4

- Molecule 1: Cbp/p300-interacting transactivator 2

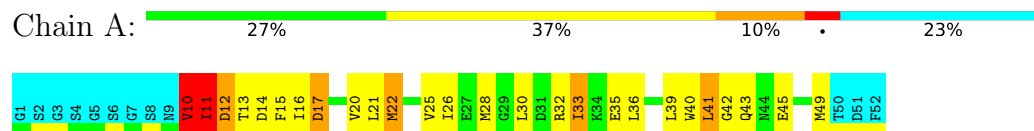


- Molecule 2: E1A-associated protein p300

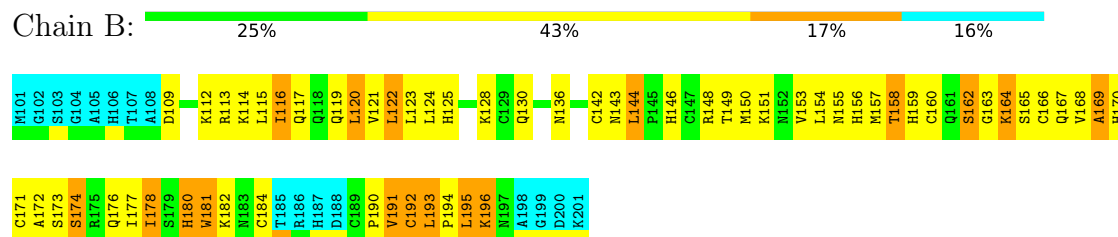


4.2.5 Score per residue for model 5

- Molecule 1: Cbp/p300-interacting transactivator 2

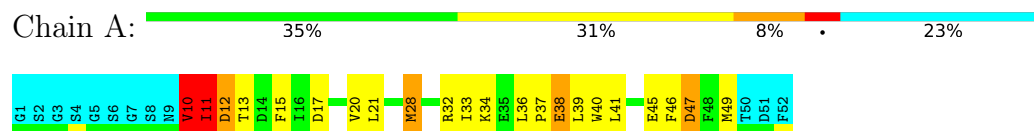


- Molecule 2: E1A-associated protein p300



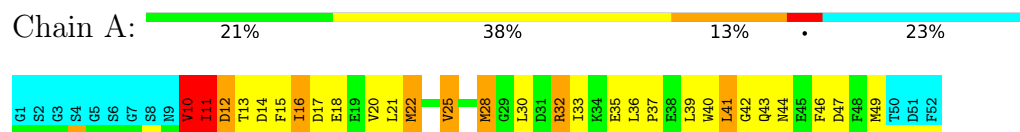
4.2.6 Score per residue for model 6

- Molecule 1: Cbp/p300-interacting transactivator 2

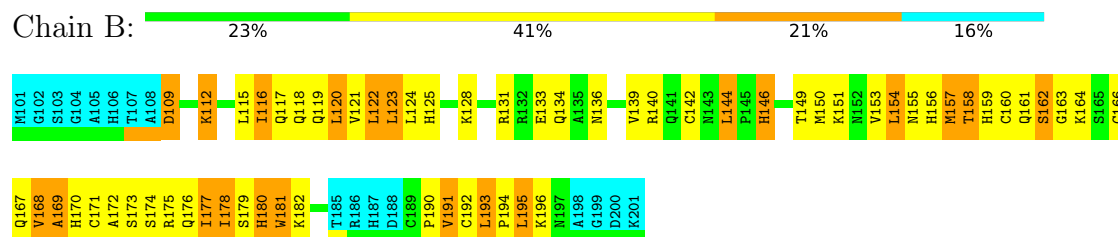


- Molecule 2: E1A-associated protein p300



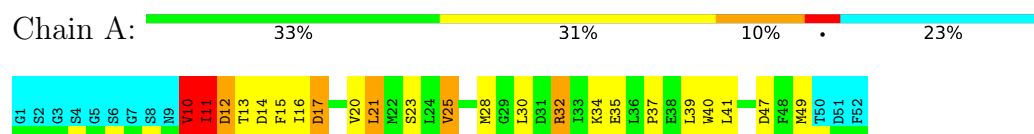


- Molecule 2: E1A-associated protein p300

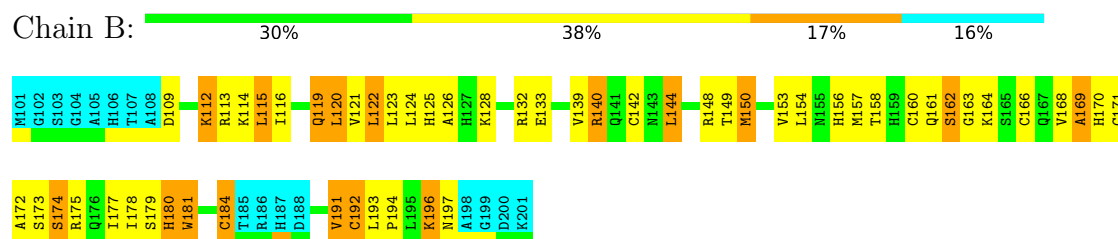


4.2.10 Score per residue for model 10

- Molecule 1: Cbp/p300-interacting transactivator 2

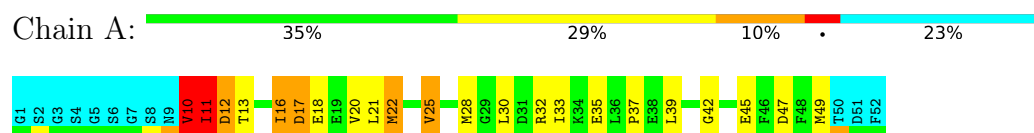


- Molecule 2: E1A-associated protein p300

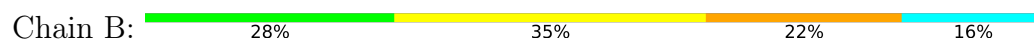


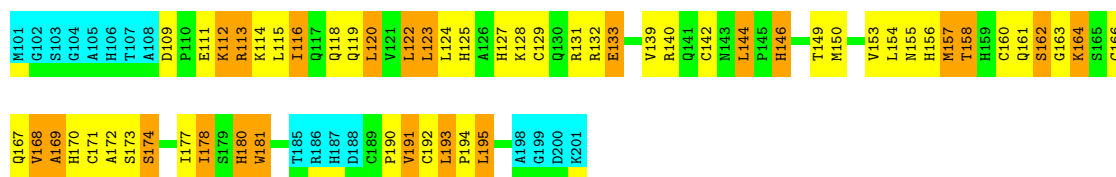
4.2.11 Score per residue for model 11

- Molecule 1: Cbp/p300-interacting transactivator 2



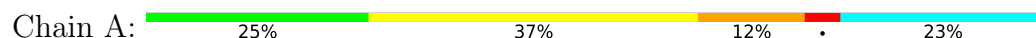
- Molecule 2: E1A-associated protein p300



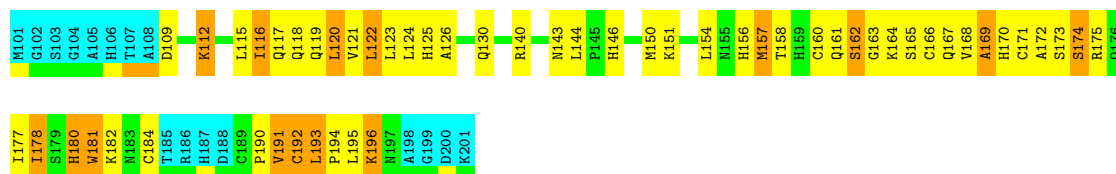


4.2.12 Score per residue for model 12 (medoid)

- Molecule 1: Cbp/p300-interacting transactivator 2

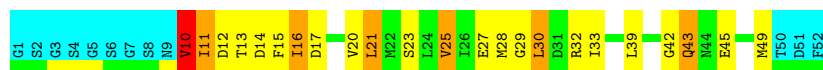


- Molecule 2: E1A-associated protein p300

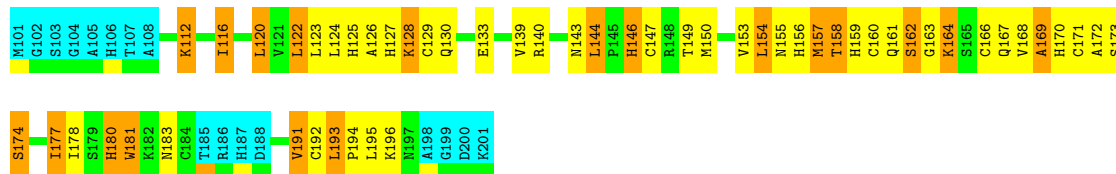
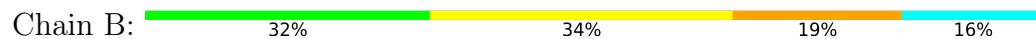


4.2.13 Score per residue for model 13

- Molecule 1: Cbp/p300-interacting transactivator 2

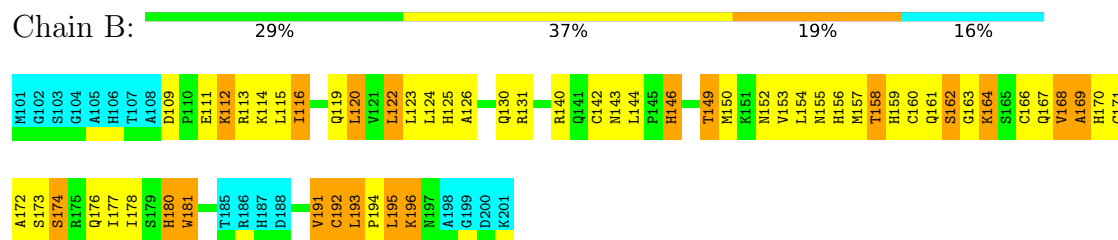
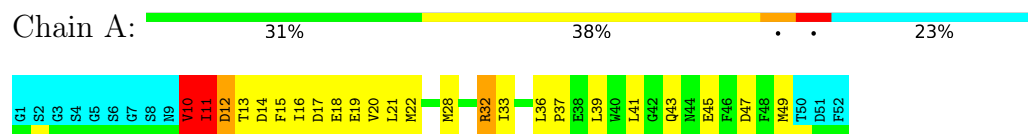


- Molecule 2: E1A-associated protein p300



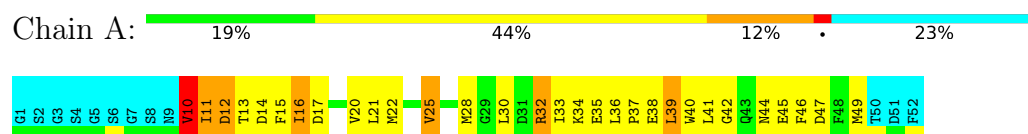
4.2.14 Score per residue for model 14

- Molecule 1: Cbp/p300-interacting transactivator 2

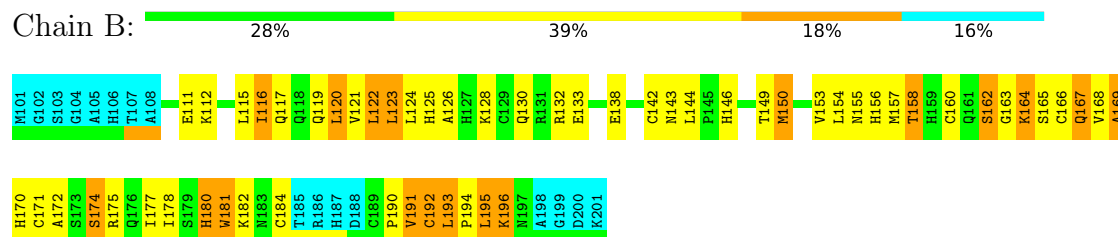


4.2.15 Score per residue for model 15

- Molecule 1: Cbp/p300-interacting transactivator 2

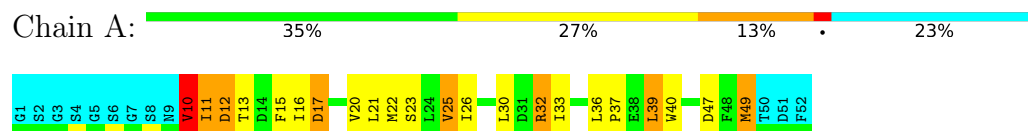


- Molecule 2: E1A-associated protein p300



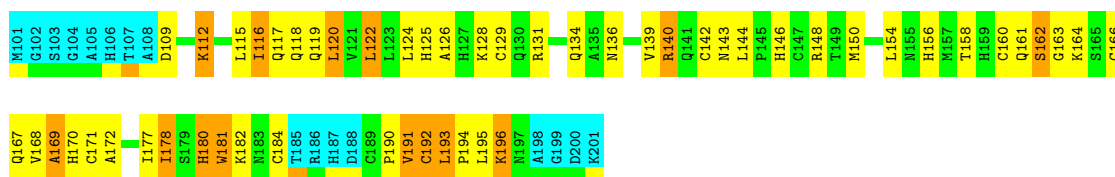
4.2.16 Score per residue for model 16

- Molecule 1: Cbp/p300-interacting transactivator 2



- Molecule 2: E1A-associated protein p300



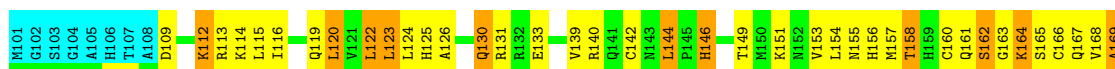


4.2.17 Score per residue for model 17

- Molecule 1: Cbp/p300-interacting transactivator 2



- Molecule 2: E1A-associated protein p300



5 Refinement protocol and experimental data overview

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

Of the 25 calculated structures, 17 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
CYANA	structure solution	1.0.4
X-PLOR	refinement	3.1

No chemical shift data was provided.

6 Model quality ⓘ

6.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section:
ZN

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	331	322	322	39±5
2	B	678	686	682	69±7
All	All	17204	17136	17068	1439

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 42.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:49:MET:HE1	2:B:116:ILE:HD13	1.09	1.19	4	11
1:A:33:ILE:HD12	2:B:146:HIS:CD2	0.87	2.03	14	13
2:B:166:CYS:O	2:B:172:ALA:HB2	0.85	1.70	2	17
1:A:11:ILE:CG2	1:A:13:THR:HG23	0.84	2.01	11	14
1:A:11:ILE:CG2	1:A:13:THR:HG22	0.83	2.03	2	2
1:A:25:VAL:HG11	2:B:144:LEU:HD22	0.82	1.49	5	3
1:A:11:ILE:HG21	1:A:13:THR:HG23	0.82	1.50	11	14
1:A:49:MET:CE	2:B:116:ILE:HD13	0.79	2.06	17	5
1:A:25:VAL:HG11	2:B:144:LEU:HD11	0.78	1.53	7	5
1:A:49:MET:HE1	2:B:116:ILE:CD1	0.78	2.09	11	8
2:B:157:MET:HE2	2:B:174:SER:OG	0.77	1.80	3	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:36:LEU:HD23	2:B:149:THR:HG21	0.77	1.55	7	1
2:B:149:THR:O	2:B:153:VAL:HG23	0.76	1.79	17	12
1:A:37:PRO:CG	2:B:115:LEU:HD13	0.75	2.10	15	6
2:B:154:LEU:HA	2:B:157:MET:HE2	0.75	1.58	17	1
2:B:124:LEU:HD21	2:B:181:TRP:CH2	0.75	2.17	15	2
1:A:17:ASP:HB3	1:A:20:VAL:HG23	0.74	1.60	2	17
2:B:153:VAL:HG12	2:B:157:MET:SD	0.73	2.23	8	1
2:B:123:LEU:HD11	2:B:174:SER:OG	0.73	1.84	7	11
2:B:112:LYS:CE	2:B:116:ILE:HD11	0.73	2.14	4	3
2:B:122:LEU:HD12	2:B:150:MET:SD	0.73	2.24	10	7
1:A:39:LEU:HD21	2:B:116:ILE:HD12	0.72	1.61	9	2
1:A:39:LEU:HD21	2:B:116:ILE:CD1	0.72	2.14	9	4
2:B:123:LEU:HD13	2:B:178:ILE:CD1	0.72	2.15	4	2
2:B:113:ARG:HA	2:B:116:ILE:HD12	0.71	1.62	17	3
2:B:112:LYS:HD3	2:B:116:ILE:HD11	0.71	1.62	17	2
1:A:39:LEU:HD22	2:B:112:LYS:HB2	0.70	1.62	9	3
2:B:123:LEU:HD22	2:B:178:ILE:HD11	0.70	1.63	15	3
1:A:33:ILE:HD12	2:B:146:HIS:NE2	0.70	2.02	4	11
2:B:120:LEU:HD22	2:B:177:ILE:CG2	0.70	2.17	7	16
1:A:49:MET:HE1	2:B:116:ILE:HG12	0.70	1.63	16	3
1:A:39:LEU:HD22	1:A:39:LEU:N	0.69	2.01	15	5
2:B:112:LYS:O	2:B:116:ILE:N	0.69	2.25	8	17
1:A:11:ILE:HG21	1:A:13:THR:HG22	0.69	1.65	1	2
2:B:180:HIS:CE1	2:B:191:VAL:HG21	0.69	2.23	11	12
2:B:112:LYS:O	2:B:116:ILE:HD12	0.68	1.89	6	4
1:A:21:LEU:HD13	2:B:125:HIS:HA	0.68	1.64	15	4
2:B:181:TRP:CZ2	2:B:195:LEU:HD22	0.68	2.22	15	5
1:A:37:PRO:CB	2:B:115:LEU:HD13	0.68	2.18	7	6
2:B:157:MET:HE2	2:B:174:SER:CB	0.67	2.18	2	2
1:A:28:MET:HE1	2:B:118:GLN:CB	0.67	2.18	6	1
2:B:120:LEU:O	2:B:124:LEU:N	0.67	2.27	1	17
1:A:25:VAL:HG11	2:B:144:LEU:HD21	0.67	1.66	3	2
2:B:122:LEU:HD12	2:B:150:MET:HG2	0.67	1.67	11	5
2:B:123:LEU:HD22	2:B:178:ILE:CD1	0.67	2.20	17	4
2:B:112:LYS:HE2	2:B:116:ILE:HD11	0.66	1.66	8	3
1:A:11:ILE:HD12	2:B:178:ILE:HG13	0.66	1.66	6	10
1:A:11:ILE:HG23	1:A:12:ASP:N	0.66	2.06	7	16
2:B:127:HIS:CE1	2:B:154:LEU:HD21	0.66	2.26	13	1
1:A:39:LEU:C	1:A:39:LEU:HD13	0.65	2.16	10	7
2:B:123:LEU:HD13	2:B:178:ILE:HD13	0.65	1.68	4	3
1:A:10:VAL:HG22	1:A:11:ILE:N	0.65	2.07	10	16

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:11:ILE:HD11	2:B:120:LEU:CD1	0.64	2.21	11	7
2:B:124:LEU:HD22	2:B:195:LEU:CD2	0.64	2.22	15	1
1:A:39:LEU:HD23	2:B:115:LEU:HB3	0.64	1.67	10	3
1:A:39:LEU:HD21	2:B:116:ILE:HG12	0.64	1.69	3	2
2:B:117:GLN:HG2	2:B:191:VAL:HG13	0.63	1.70	15	2
1:A:21:LEU:HD13	2:B:121:VAL:HG13	0.63	1.70	7	6
2:B:120:LEU:O	2:B:124:LEU:CD1	0.63	2.47	15	4
1:A:39:LEU:HD23	2:B:116:ILE:HD12	0.63	1.68	16	1
2:B:181:TRP:HE1	2:B:191:VAL:HG12	0.63	1.53	2	17
1:A:40:TRP:CH2	2:B:172:ALA:HB3	0.63	2.28	5	9
2:B:180:HIS:NE2	2:B:191:VAL:HG21	0.63	2.09	9	12
2:B:117:GLN:O	2:B:121:VAL:HG23	0.63	1.92	15	2
2:B:154:LEU:C	2:B:154:LEU:HD12	0.62	2.19	17	1
1:A:36:LEU:HD12	1:A:37:PRO:HD2	0.62	1.70	9	8
1:A:10:VAL:HG13	1:A:11:ILE:H	0.62	1.55	13	16
1:A:49:MET:HE1	2:B:116:ILE:HG13	0.62	1.72	10	2
2:B:112:LYS:CD	2:B:116:ILE:HD11	0.62	2.25	4	2
1:A:28:MET:HE1	2:B:118:GLN:HB2	0.62	1.70	4	3
2:B:120:LEU:O	2:B:124:LEU:HG	0.61	1.94	1	15
1:A:39:LEU:HD11	2:B:116:ILE:CD1	0.61	2.25	10	3
1:A:38:GLU:C	1:A:39:LEU:HD22	0.61	2.20	2	7
1:A:41:LEU:HD11	2:B:176:GLN:NE2	0.60	2.11	5	1
1:A:11:ILE:HD11	2:B:178:ILE:HG13	0.60	1.72	13	1
1:A:39:LEU:HD23	2:B:116:ILE:CD1	0.60	2.25	16	1
2:B:120:LEU:C	2:B:124:LEU:HD12	0.60	2.22	17	13
2:B:112:LYS:CE	2:B:116:ILE:HD13	0.60	2.26	1	1
2:B:115:LEU:HD23	2:B:118:GLN:CD	0.59	2.22	7	2
2:B:133:GLU:CD	2:B:139:VAL:HG22	0.59	2.23	10	1
1:A:21:LEU:HD11	2:B:125:HIS:HB2	0.59	1.75	8	10
1:A:39:LEU:CG	2:B:177:ILE:HD11	0.59	2.27	9	7
2:B:157:MET:HE2	2:B:174:SER:HB3	0.59	1.74	11	3
2:B:120:LEU:O	2:B:124:LEU:CG	0.59	2.51	13	15
2:B:120:LEU:O	2:B:124:LEU:HD13	0.58	1.97	15	2
1:A:37:PRO:HG2	2:B:115:LEU:HD21	0.58	1.72	3	4
1:A:39:LEU:HD12	2:B:177:ILE:HD11	0.58	1.75	5	7
1:A:37:PRO:HB2	2:B:115:LEU:HD13	0.58	1.73	7	6
2:B:124:LEU:HD21	2:B:181:TRP:CZ3	0.58	2.34	5	1
1:A:49:MET:HE1	2:B:116:ILE:CG1	0.58	2.29	12	2
1:A:39:LEU:HG	2:B:177:ILE:HD11	0.58	1.74	9	7
1:A:39:LEU:HD11	2:B:115:LEU:HB3	0.58	1.76	16	3
1:A:11:ILE:HD11	2:B:120:LEU:HD12	0.57	1.73	11	6

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:156:HIS:HE2	2:B:171:CYS:CB	0.57	2.10	16	15
2:B:180:HIS:CD2	2:B:181:TRP:N	0.57	2.72	2	17
1:A:39:LEU:HD11	2:B:116:ILE:HD11	0.57	1.76	10	4
1:A:39:LEU:CD1	2:B:177:ILE:HD11	0.57	2.29	14	7
1:A:30:LEU:N	1:A:30:LEU:HD23	0.57	2.14	13	8
1:A:39:LEU:HD21	2:B:112:LYS:O	0.57	2.00	14	3
1:A:25:VAL:HG13	1:A:30:LEU:HB2	0.56	1.76	13	9
2:B:191:VAL:O	2:B:195:LEU:HD23	0.56	2.00	12	2
1:A:39:LEU:HD11	2:B:115:LEU:CB	0.56	2.30	17	5
1:A:39:LEU:HD21	2:B:116:ILE:CG1	0.56	2.31	9	4
1:A:39:LEU:HB3	2:B:177:ILE:HD11	0.56	1.76	17	2
1:A:28:MET:HE1	2:B:118:GLN:HB3	0.56	1.78	6	3
2:B:120:LEU:CD2	2:B:181:TRP:CD1	0.55	2.89	11	16
2:B:126:ALA:HB3	2:B:154:LEU:CD1	0.55	2.31	13	1
2:B:115:LEU:O	2:B:119:GLN:N	0.55	2.36	16	10
1:A:37:PRO:HG2	2:B:115:LEU:HD11	0.55	1.76	3	4
1:A:21:LEU:HD13	2:B:125:HIS:CA	0.55	2.32	15	3
2:B:123:LEU:HD23	2:B:154:LEU:CD1	0.55	2.31	9	1
2:B:133:GLU:CD	2:B:139:VAL:HG13	0.55	2.27	1	1
2:B:123:LEU:HD11	2:B:174:SER:HB3	0.55	1.79	4	1
1:A:39:LEU:HD21	2:B:116:ILE:HG13	0.54	1.79	11	5
2:B:123:LEU:HD21	2:B:174:SER:OG	0.54	2.02	14	7
1:A:10:VAL:HG13	1:A:11:ILE:N	0.54	2.16	1	1
1:A:21:LEU:HD13	2:B:125:HIS:HB2	0.54	1.77	11	3
1:A:39:LEU:HD22	2:B:177:ILE:CD1	0.54	2.32	16	1
2:B:154:LEU:HD12	2:B:155:ASN:N	0.54	2.18	17	1
2:B:157:MET:HE2	2:B:174:SER:HB2	0.54	1.78	2	1
2:B:193:LEU:CB	2:B:194:PRO:HD3	0.54	2.33	3	5
2:B:112:LYS:C	2:B:116:ILE:HD12	0.54	2.27	14	4
2:B:133:GLU:OE2	2:B:139:VAL:HG13	0.54	2.03	1	1
2:B:133:GLU:HB3	2:B:139:VAL:HG22	0.54	1.80	1	1
1:A:39:LEU:HD11	2:B:116:ILE:CG1	0.54	2.33	13	2
1:A:25:VAL:HG22	1:A:30:LEU:HB2	0.54	1.80	16	5
2:B:117:GLN:CG	2:B:191:VAL:HG13	0.54	2.32	15	1
2:B:120:LEU:C	2:B:124:LEU:HD13	0.53	2.28	15	2
1:A:21:LEU:O	1:A:25:VAL:HG23	0.53	2.03	7	1
1:A:17:ASP:CB	1:A:20:VAL:HG23	0.53	2.33	15	8
2:B:122:LEU:HD21	2:B:144:LEU:HD13	0.53	1.79	5	1
1:A:13:THR:CG2	2:B:124:LEU:HD23	0.53	2.34	17	8
2:B:154:LEU:CA	2:B:157:MET:HE2	0.53	2.33	17	1
1:A:39:LEU:HD21	2:B:116:ILE:HA	0.53	1.80	16	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:122:LEU:HD11	2:B:146:HIS:C	0.53	2.29	6	2
1:A:39:LEU:HD11	2:B:116:ILE:HG13	0.52	1.81	14	3
2:B:162:SER:O	2:B:164:LYS:N	0.52	2.42	15	17
2:B:181:TRP:CZ2	2:B:192:CYS:HA	0.52	2.39	16	12
2:B:120:LEU:HG	2:B:181:TRP:CD1	0.52	2.40	1	12
1:A:37:PRO:HG3	2:B:115:LEU:HD22	0.52	1.81	15	2
1:A:39:LEU:HD21	2:B:116:ILE:CA	0.52	2.34	16	1
2:B:112:LYS:HE2	2:B:116:ILE:HD13	0.52	1.81	1	1
2:B:124:LEU:CD1	2:B:195:LEU:HD23	0.52	2.35	4	1
1:A:21:LEU:HD13	1:A:21:LEU:N	0.52	2.19	4	1
1:A:39:LEU:HD11	2:B:119:GLN:HB2	0.52	1.80	6	1
2:B:120:LEU:CD1	2:B:178:ILE:HD12	0.52	2.35	16	7
1:A:13:THR:HG22	2:B:124:LEU:HD23	0.52	1.82	5	1
2:B:156:HIS:CE1	2:B:168:VAL:HG11	0.51	2.39	3	11
1:A:29:GLY:C	1:A:30:LEU:HD23	0.51	2.30	13	2
2:B:120:LEU:HD22	2:B:177:ILE:HG23	0.51	1.81	11	8
1:A:39:LEU:HD11	2:B:119:GLN:CB	0.51	2.36	6	1
1:A:40:TRP:O	1:A:42:GLY:N	0.51	2.42	1	8
1:A:13:THR:CB	2:B:124:LEU:HD22	0.51	2.35	1	1
1:A:41:LEU:H	1:A:41:LEU:HD22	0.51	1.65	2	1
1:A:36:LEU:HD22	2:B:146:HIS:CE1	0.51	2.41	8	3
1:A:15:PHE:CD1	2:B:196:LYS:CG	0.50	2.94	16	9
2:B:130:GLN:NE2	2:B:154:LEU:HD11	0.50	2.22	17	1
2:B:156:HIS:NE2	2:B:171:CYS:HB2	0.50	2.22	5	15
1:A:40:TRP:O	2:B:116:ILE:HD11	0.50	2.06	2	1
1:A:11:ILE:HG12	2:B:181:TRP:CE3	0.50	2.41	11	15
1:A:39:LEU:HD21	2:B:119:GLN:HG2	0.50	1.83	1	1
1:A:39:LEU:HD22	2:B:112:LYS:CB	0.50	2.35	9	2
1:A:13:THR:HG22	2:B:124:LEU:CD2	0.50	2.35	17	6
2:B:112:LYS:CD	2:B:113:ARG:N	0.50	2.74	14	3
1:A:39:LEU:C	1:A:39:LEU:CD1	0.50	2.85	10	2
1:A:46:PHE:CZ	2:B:113:ARG:CD	0.50	2.95	7	1
2:B:127:HIS:CE1	2:B:154:LEU:CD2	0.50	2.94	13	1
1:A:15:PHE:CE1	2:B:196:LYS:CB	0.49	2.95	5	6
1:A:41:LEU:HD21	2:B:176:GLN:CD	0.49	2.32	3	1
2:B:130:GLN:OE1	2:B:154:LEU:HD22	0.49	2.07	13	1
2:B:122:LEU:HD21	2:B:144:LEU:CD2	0.49	2.37	2	2
2:B:115:LEU:HD12	2:B:118:GLN:CD	0.49	2.32	11	1
1:A:39:LEU:N	1:A:39:LEU:CD2	0.49	2.73	15	5
1:A:10:VAL:CG1	1:A:11:ILE:N	0.49	2.76	1	1
2:B:155:ASN:O	2:B:158:THR:HG22	0.49	2.06	17	8

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:36:LEU:CD2	2:B:149:THR:HG21	0.49	2.37	5	1
2:B:156:HIS:HE2	2:B:168:VAL:HB	0.49	1.65	17	2
1:A:11:ILE:CG2	1:A:12:ASP:N	0.49	2.75	16	9
1:A:40:TRP:CZ3	2:B:172:ALA:HB3	0.49	2.42	16	2
2:B:123:LEU:O	2:B:127:HIS:CD2	0.49	2.66	13	1
1:A:15:PHE:CZ	2:B:196:LYS:CG	0.49	2.95	1	3
2:B:156:HIS:O	2:B:160:CYS:N	0.49	2.45	17	17
2:B:193:LEU:CB	2:B:194:PRO:CD	0.49	2.90	1	13
2:B:120:LEU:O	2:B:124:LEU:HD12	0.48	2.08	17	8
2:B:112:LYS:HG3	2:B:113:ARG:N	0.48	2.23	7	3
1:A:10:VAL:CG2	1:A:11:ILE:N	0.48	2.76	3	7
1:A:38:GLU:O	1:A:39:LEU:HD22	0.48	2.08	6	1
1:A:46:PHE:CD2	2:B:113:ARG:CG	0.48	2.96	1	1
1:A:33:ILE:HD12	2:B:146:HIS:HB2	0.48	1.85	6	1
1:A:11:ILE:HG23	1:A:13:THR:H	0.48	1.67	5	13
1:A:40:TRP:O	2:B:112:LYS:HG3	0.48	2.09	9	1
2:B:173:SER:O	2:B:177:ILE:N	0.48	2.38	2	1
2:B:125:HIS:O	2:B:129:CYS:N	0.48	2.47	13	6
2:B:168:VAL:O	2:B:169:ALA:C	0.48	2.57	11	17
1:A:49:MET:HE3	2:B:116:ILE:HD13	0.48	1.84	13	1
2:B:162:SER:HB3	2:B:166:CYS:N	0.48	2.24	17	2
1:A:12:ASP:O	1:A:15:PHE:CZ	0.47	2.67	13	6
2:B:120:LEU:CD2	2:B:177:ILE:CG2	0.47	2.92	10	17
1:A:24:LEU:HD12	2:B:121:VAL:HG21	0.47	1.84	4	1
2:B:120:LEU:HD11	2:B:178:ILE:HD12	0.47	1.85	16	2
1:A:39:LEU:HD22	2:B:177:ILE:HD11	0.47	1.86	16	1
2:B:180:HIS:CD2	2:B:180:HIS:C	0.47	2.93	8	17
2:B:168:VAL:O	2:B:171:CYS:N	0.47	2.46	14	16
2:B:157:MET:HE1	2:B:174:SER:HB3	0.47	1.87	12	3
2:B:122:LEU:HD21	2:B:144:LEU:HD21	0.47	1.85	17	1
1:A:40:TRP:O	1:A:41:LEU:C	0.47	2.58	17	5
2:B:124:LEU:HD13	2:B:195:LEU:HD22	0.47	1.87	3	2
2:B:190:PRO:O	2:B:194:PRO:CG	0.47	2.62	1	10
1:A:30:LEU:HA	1:A:33:ILE:HD11	0.47	1.86	3	1
2:B:115:LEU:O	2:B:119:GLN:CB	0.47	2.62	11	8
2:B:190:PRO:O	2:B:194:PRO:HG3	0.47	2.09	8	3
1:A:33:ILE:CD1	2:B:146:HIS:CD2	0.47	2.98	5	1
2:B:156:HIS:NE2	2:B:171:CYS:CB	0.47	2.78	1	9
2:B:181:TRP:NE1	2:B:191:VAL:HG12	0.47	2.24	1	1
1:A:21:LEU:HD13	2:B:121:VAL:CG1	0.47	2.40	2	1
1:A:37:PRO:CG	2:B:115:LEU:HD11	0.47	2.40	3	2

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:11:ILE:CG1	2:B:181:TRP:CE3	0.47	2.98	17	5
1:A:12:ASP:O	1:A:15:PHE:CE2	0.46	2.69	1	1
1:A:36:LEU:HD21	2:B:149:THR:HG21	0.46	1.86	5	1
1:A:21:LEU:HD12	1:A:21:LEU:C	0.46	2.35	8	3
1:A:49:MET:HE2	2:B:116:ILE:HD12	0.46	1.85	3	1
2:B:193:LEU:HB3	2:B:194:PRO:CD	0.46	2.40	4	2
2:B:116:ILE:HG22	2:B:117:GLN:N	0.46	2.25	1	3
1:A:15:PHE:CE1	2:B:196:LYS:CG	0.46	2.99	6	6
2:B:189:CYS:O	2:B:193:LEU:CB	0.46	2.64	4	1
1:A:41:LEU:HD12	1:A:47:ASP:OD1	0.46	2.11	6	1
1:A:16:ILE:CD1	2:B:195:LEU:HD23	0.46	2.40	15	1
1:A:39:LEU:CD2	2:B:116:ILE:HD12	0.46	2.39	16	1
2:B:174:SER:O	2:B:178:ILE:HD13	0.46	2.10	15	1
1:A:10:VAL:O	1:A:11:ILE:O	0.46	2.34	1	1
1:A:21:LEU:HD13	2:B:125:HIS:CB	0.46	2.41	14	3
1:A:40:TRP:O	2:B:112:LYS:CD	0.46	2.64	8	2
1:A:11:ILE:HG23	1:A:13:THR:HG22	0.46	1.83	2	2
1:A:12:ASP:O	1:A:15:PHE:CE1	0.46	2.69	7	9
2:B:120:LEU:HD21	2:B:181:TRP:CD1	0.46	2.46	11	4
1:A:41:LEU:HD13	1:A:41:LEU:N	0.45	2.26	2	1
1:A:33:ILE:HD13	2:B:146:HIS:CD2	0.45	2.47	5	1
1:A:22:MET:HE1	2:B:125:HIS:CE1	0.45	2.47	9	3
2:B:120:LEU:HG	2:B:181:TRP:NE1	0.45	2.27	7	6
1:A:39:LEU:HD11	2:B:115:LEU:HB2	0.45	1.88	15	2
2:B:173:SER:O	2:B:177:ILE:HG13	0.45	2.12	10	2
1:A:41:LEU:HD23	1:A:41:LEU:N	0.45	2.27	6	1
1:A:37:PRO:CG	2:B:115:LEU:HD22	0.45	2.42	1	4
1:A:39:LEU:HD12	2:B:116:ILE:HD13	0.45	1.88	1	1
2:B:124:LEU:CD2	2:B:181:TRP:CH2	0.45	2.97	5	1
2:B:133:GLU:OE1	2:B:139:VAL:HG13	0.45	2.12	9	1
2:B:139:VAL:CG2	2:B:140:ARG:N	0.44	2.80	17	3
1:A:21:LEU:HA	1:A:24:LEU:HD12	0.44	1.89	1	1
1:A:30:LEU:HD22	2:B:118:GLN:CD	0.44	2.37	9	2
2:B:122:LEU:HD12	2:B:150:MET:CG	0.44	2.41	11	1
2:B:111:GLU:O	2:B:115:LEU:N	0.44	2.40	4	1
1:A:36:LEU:CD2	2:B:146:HIS:CE1	0.44	3.01	6	1
1:A:11:ILE:HG23	1:A:13:THR:HG23	0.44	1.89	7	3
1:A:15:PHE:CG	2:B:196:LYS:CG	0.44	3.01	15	1
2:B:170:HIS:O	2:B:173:SER:N	0.44	2.51	14	11
1:A:15:PHE:CE1	2:B:196:LYS:HA	0.44	2.47	7	1
2:B:139:VAL:HG22	2:B:140:ARG:N	0.44	2.27	17	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:112:LYS:HE2	2:B:116:ILE:CD1	0.44	2.43	12	1
1:A:46:PHE:CZ	2:B:113:ARG:HG3	0.44	2.47	7	1
2:B:156:HIS:NE2	2:B:168:VAL:HB	0.44	2.28	15	2
1:A:10:VAL:CG2	2:B:178:ILE:CG2	0.44	2.96	6	5
2:B:120:LEU:O	2:B:124:LEU:CB	0.44	2.65	13	1
2:B:156:HIS:HE2	2:B:171:CYS:HB2	0.43	1.67	3	6
2:B:190:PRO:C	2:B:194:PRO:CG	0.43	2.91	9	6
2:B:180:HIS:NE2	2:B:191:VAL:CG2	0.43	2.80	5	12
2:B:133:GLU:HB3	2:B:139:VAL:HB	0.43	1.90	11	1
2:B:170:HIS:O	2:B:171:CYS:C	0.43	2.61	4	17
1:A:33:ILE:HG22	1:A:34:LYS:N	0.43	2.28	15	1
1:A:39:LEU:HD23	2:B:119:GLN:OE1	0.43	2.13	8	1
2:B:116:ILE:CG2	2:B:117:GLN:N	0.43	2.81	7	3
2:B:123:LEU:HD13	2:B:178:ILE:HD11	0.43	1.89	4	1
2:B:124:LEU:O	2:B:128:LYS:HG3	0.43	2.13	13	1
2:B:154:LEU:C	2:B:154:LEU:CD1	0.43	2.90	17	1
1:A:41:LEU:HD11	2:B:176:GLN:OE1	0.43	2.14	2	1
1:A:11:ILE:CD1	2:B:120:LEU:HD12	0.43	2.43	11	1
1:A:38:GLU:O	2:B:112:LYS:HG2	0.43	2.14	2	1
2:B:156:HIS:CD2	2:B:171:CYS:SG	0.43	3.10	9	5
1:A:11:ILE:HD13	2:B:124:LEU:HD23	0.43	1.89	13	1
2:B:113:ARG:CA	2:B:116:ILE:HD12	0.43	2.39	17	1
2:B:126:ALA:CB	2:B:150:MET:CB	0.43	2.97	10	6
2:B:120:LEU:CD2	2:B:177:ILE:HG23	0.43	2.44	2	1
2:B:190:PRO:O	2:B:194:PRO:HG2	0.43	2.13	3	1
2:B:112:LYS:O	2:B:115:LEU:N	0.43	2.52	9	2
1:A:13:THR:CG2	2:B:124:LEU:CD2	0.43	2.97	17	7
1:A:21:LEU:HA	1:A:24:LEU:HD21	0.43	1.91	7	1
1:A:16:ILE:HD11	2:B:195:LEU:CD2	0.43	2.42	8	1
2:B:139:VAL:HG12	2:B:140:ARG:N	0.43	2.29	10	3
2:B:122:LEU:O	2:B:126:ALA:CB	0.43	2.67	17	4
2:B:167:GLN:O	2:B:167:GLN:CG	0.43	2.67	11	15
2:B:109:ASP:O	2:B:112:LYS:CG	0.43	2.67	6	1
2:B:156:HIS:CD2	2:B:171:CYS:HB2	0.43	2.49	17	1
1:A:12:ASP:HB3	1:A:15:PHE:CZ	0.42	2.49	6	2
2:B:121:VAL:HG22	2:B:195:LEU:HD21	0.42	1.90	8	2
1:A:40:TRP:CZ3	2:B:173:SER:N	0.42	2.87	12	1
1:A:36:LEU:HD11	2:B:146:HIS:CE1	0.42	2.49	5	1
1:A:26:ILE:HD12	1:A:26:ILE:N	0.42	2.29	12	2
1:A:37:PRO:HG2	2:B:115:LEU:CD1	0.42	2.44	11	2
2:B:121:VAL:CG2	2:B:195:LEU:HD21	0.42	2.44	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
2:B:125:HIS:CD2	2:B:125:HIS:C	0.42	2.98	4	1
1:A:38:GLU:O	2:B:112:LYS:CD	0.42	2.68	6	1
1:A:15:PHE:CD1	2:B:196:LYS:HG2	0.42	2.50	10	4
1:A:39:LEU:CD1	1:A:39:LEU:C	0.42	2.93	11	2
1:A:30:LEU:N	1:A:30:LEU:CD2	0.42	2.82	13	1
2:B:126:ALA:CB	2:B:154:LEU:CD1	0.42	2.97	13	1
1:A:15:PHE:C	1:A:15:PHE:CD1	0.42	2.97	1	1
1:A:49:MET:CE	2:B:116:ILE:CD1	0.42	2.97	12	2
2:B:126:ALA:CB	2:B:150:MET:HB2	0.42	2.44	16	1
2:B:123:LEU:O	2:B:127:HIS:HB2	0.42	2.15	3	1
2:B:156:HIS:O	2:B:160:CYS:HB2	0.42	2.15	17	1
1:A:11:ILE:HG21	2:B:124:LEU:HD23	0.42	1.92	13	1
2:B:166:CYS:O	2:B:172:ALA:CB	0.42	2.65	15	1
1:A:46:PHE:CD2	2:B:113:ARG:HG3	0.42	2.50	1	1
1:A:15:PHE:CE1	2:B:196:LYS:HG3	0.42	2.50	12	2
1:A:16:ILE:HD11	2:B:195:LEU:HD23	0.42	1.92	9	1
1:A:39:LEU:HD22	2:B:112:LYS:HA	0.42	1.92	14	2
1:A:15:PHE:CZ	2:B:196:LYS:HG2	0.41	2.50	1	1
1:A:39:LEU:O	1:A:40:TRP:CE3	0.41	2.73	1	1
2:B:122:LEU:CD2	2:B:144:LEU:CD2	0.41	2.98	2	1
1:A:39:LEU:CD2	2:B:112:LYS:O	0.41	2.67	3	2
2:B:174:SER:O	2:B:178:ILE:CD1	0.41	2.67	5	1
1:A:13:THR:HG22	2:B:124:LEU:HD22	0.41	1.92	17	3
2:B:193:LEU:O	2:B:196:LYS:CG	0.41	2.68	3	2
2:B:115:LEU:HD12	2:B:118:GLN:OE1	0.41	2.16	11	1
2:B:157:MET:HG3	2:B:158:THR:N	0.41	2.30	17	1
1:A:11:ILE:HG23	1:A:13:THR:N	0.41	2.30	5	3
1:A:22:MET:O	1:A:26:ILE:CD1	0.41	2.68	5	2
1:A:13:THR:HG22	1:A:16:ILE:HG21	0.41	1.90	11	1
1:A:15:PHE:CD1	1:A:16:ILE:N	0.41	2.88	13	1
2:B:119:GLN:O	2:B:123:LEU:N	0.41	2.53	15	1
1:A:30:LEU:HD11	2:B:118:GLN:HB2	0.41	1.92	16	1
1:A:15:PHE:CE1	2:B:196:LYS:HB2	0.41	2.50	3	4
2:B:180:HIS:O	2:B:184:CYS:N	0.41	2.54	10	1
2:B:119:GLN:O	2:B:122:LEU:N	0.41	2.53	16	3
2:B:122:LEU:HD22	2:B:122:LEU:HA	0.41	1.81	10	2
1:A:36:LEU:HD12	1:A:37:PRO:CD	0.41	2.46	2	2
2:B:113:ARG:O	2:B:116:ILE:HB	0.41	2.14	10	1
2:B:125:HIS:CD2	2:B:147:CYS:SG	0.41	3.14	13	1
1:A:49:MET:CE	2:B:116:ILE:HG12	0.41	2.46	1	1
1:A:40:TRP:CD1	2:B:112:LYS:HE2	0.41	2.51	2	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:37:PRO:CG	2:B:115:LEU:CD1	0.41	2.98	4	1
2:B:121:VAL:HG23	2:B:195:LEU:HD21	0.41	1.93	4	1
2:B:191:VAL:O	2:B:195:LEU:HD13	0.41	2.16	9	1
2:B:133:GLU:OE1	2:B:139:VAL:HG22	0.41	2.16	13	1
1:A:11:ILE:CG2	1:A:13:THR:CG2	0.41	2.94	16	1
2:B:127:HIS:CD2	2:B:127:HIS:C	0.41	2.98	11	1
1:A:15:PHE:CD1	2:B:196:LYS:HG3	0.41	2.51	16	1
2:B:120:LEU:CD2	2:B:177:ILE:HG22	0.41	2.46	17	1
1:A:26:ILE:N	1:A:26:ILE:HD12	0.40	2.31	8	1
1:A:15:PHE:CZ	2:B:196:LYS:HG3	0.40	2.51	13	1
2:B:120:LEU:O	2:B:124:LEU:HB2	0.40	2.16	5	1
2:B:109:ASP:O	2:B:112:LYS:CD	0.40	2.70	9	1
1:A:21:LEU:HD12	1:A:22:MET:N	0.40	2.30	11	1
2:B:124:LEU:O	2:B:127:HIS:HB3	0.40	2.17	11	1
2:B:175:ARG:O	2:B:179:SER:CB	0.40	2.70	9	1
1:A:11:ILE:HD13	2:B:124:LEU:CD2	0.40	2.46	13	1
2:B:120:LEU:HD13	2:B:120:LEU:HA	0.40	1.82	14	1
2:B:130:GLN:NE2	2:B:154:LEU:CB	0.40	2.85	6	1
1:A:25:VAL:HG11	2:B:144:LEU:CD1	0.40	2.38	7	1
2:B:124:LEU:CD1	2:B:195:LEU:HD12	0.40	2.47	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	40/52 (77%)	33±1 (82±3%)	4±1 (9±3%)	3±1 (9±2%)	1	12
2	B	85/101 (84%)	75±1 (88±1%)	8±1 (9±1%)	2±0 (3±1%)	6	41
All	All	2125/2601 (82%)	1838 (86%)	189 (9%)	98 (5%)	3	26

All 12 unique Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
2	B	163	GLY	17

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Mol	Chain	Res	Type	Models (Total)
2	B	169	ALA	17
1	A	10	VAL	16
1	A	32	ARG	16
1	A	11	ILE	15
1	A	41	LEU	8
2	B	177	ILE	2
1	A	42	GLY	2
1	A	43	GLN	2
2	B	145	PRO	1
2	B	197	ASN	1
2	B	143	ASN	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	38/46 (83%)	25±1 (66±4%)	13±1 (34±4%)	1	11
2	B	79/89 (89%)	50±4 (63±5%)	29±4 (37±5%)	1	8
All	All	1989/2295 (87%)	1276 (64%)	713 (36%)	1	9

All 89 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)
2	B	120	LEU	17
2	B	122	LEU	17
2	B	144	LEU	17
2	B	158	THR	17
2	B	162	SER	17
2	B	180	HIS	17
2	B	181	TRP	17
1	A	16	ILE	16
1	A	28	MET	16
2	B	154	LEU	16
1	A	10	VAL	16
1	A	11	ILE	15
1	A	12	ASP	15

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Mol	Chain	Res	Type	Models (Total)
2	B	161	GLN	15
2	B	178	ILE	15
2	B	191	VAL	15
2	B	109	ASP	14
1	A	25	VAL	13
2	B	142	CYS	13
2	B	157	MET	13
2	B	174	SER	13
2	B	192	CYS	13
2	B	195	LEU	13
1	A	14	ASP	12
2	B	112	LYS	12
2	B	128	LYS	12
1	A	45	GLU	11
1	A	47	ASP	11
2	B	116	ILE	11
2	B	196	LYS	11
2	B	193	LEU	11
2	B	182	LYS	10
2	B	164	LYS	10
2	B	146	HIS	9
2	B	175	ARG	9
2	B	184	CYS	9
1	A	21	LEU	8
2	B	131	ARG	8
2	B	140	ARG	8
1	A	22	MET	8
2	B	136	ASN	8
2	B	151	LYS	8
2	B	165	SER	8
1	A	46	PHE	7
2	B	123	LEU	7
2	B	133	GLU	7
2	B	159	HIS	7
1	A	41	LEU	7
2	B	114	LYS	7
1	A	35	GLU	7
1	A	32	ARG	7
1	A	49	MET	6
2	B	148	ARG	6
1	A	17	ASP	6
1	A	43	GLN	6

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Mol	Chain	Res	Type	Models (Total)
2	B	130	GLN	6
2	B	150	MET	5
1	A	18	GLU	5
2	B	143	ASN	5
2	B	113	ARG	4
2	B	134	GLN	4
2	B	168	VAL	4
1	A	39	LEU	4
2	B	132	ARG	4
1	A	27	GLU	3
2	B	119	GLN	3
2	B	176	GLN	3
2	B	183	ASN	3
1	A	44	ASN	3
1	A	23	SER	3
2	B	111	GLU	3
1	A	36	LEU	2
1	A	48	PHE	2
2	B	179	SER	2
2	B	115	LEU	2
2	B	138	GLU	2
1	A	19	GLU	2
2	B	118	GLN	2
1	A	34	LYS	2
2	B	149	THR	2
2	B	173	SER	1
2	B	155	ASN	1
1	A	33	ILE	1
1	A	38	GLU	1
2	B	117	GLN	1
1	A	13	THR	1
1	A	30	LEU	1
2	B	152	ASN	1
2	B	167	GLN	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 [i](#)

There are no oligosaccharides in this entry.

6.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

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