



Full wwPDB NMR Structure Validation Report ⓘ

Mar 7, 2026 – 02:13 AM UTC

PDB ID : 1IHW / pdb_00001ihw
Title : SOLUTION STRUCTURE OF THE DNA BINDING DOMAIN OF HIV-1
INTEGRASE, NMR, 40 STRUCTURES
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Deposited on : 1995-05-12

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 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

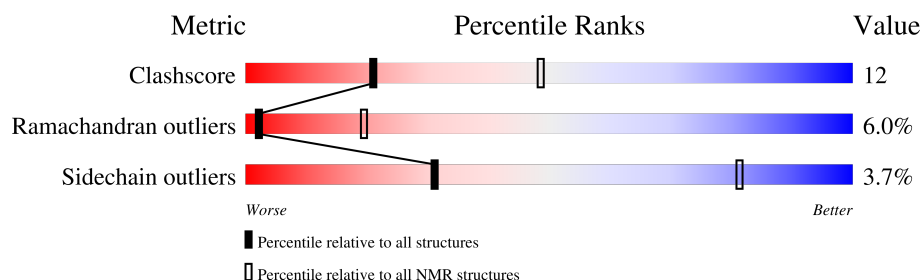
1 Overall quality at a glance

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	
1	B	52	

2 Ensemble composition and analysis

This entry contains 40 models. Model 16 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:222-A:269, B:222-B:270 (97)	0.45	16

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 2 single-model clusters were found.

Cluster number	Models
1	5, 8, 9, 13, 14, 16, 25, 29, 37, 40
2	1, 3, 12, 18, 20, 26, 35, 36
3	2, 7, 15, 19, 24, 32, 39
4	4, 6, 10, 11, 21, 27, 38
5	22, 23, 28, 31, 33, 34
Single-model clusters	17; 30

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1770 atoms, of which 904 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called HIV-1 INTEGRASE.

Mol	Chain	Residues	Atoms							Trace
1	A	52	Total	C	H	N	O	S	0	
			885	277	452	82	73	1		
1	B	52	Total	C	H	N	O	S	0	
			885	277	452	82	73	1		

There are 4 discrepancies between the modelled and reference sequences:

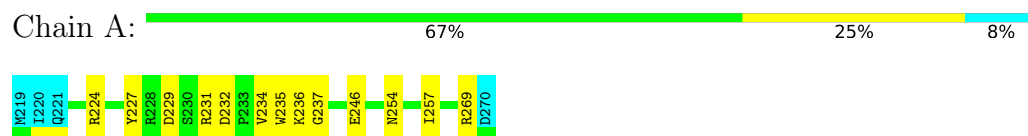
Chain	Residue	Modelled	Actual	Comment	Reference
A	234	VAL	ILE	conflict	UNP P04586
A	265	ALA	VAL	conflict	UNP P04586
B	234	VAL	ILE	conflict	UNP P04586
B	265	ALA	VAL	conflict	UNP P04586

4 Residue-property plots [i](#)

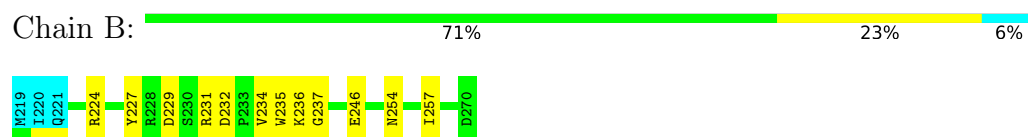
4.1 Average score per residue in the NMR ensemble

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: HIV-1 INTEGRASE



• Molecule 1: HIV-1 INTEGRASE

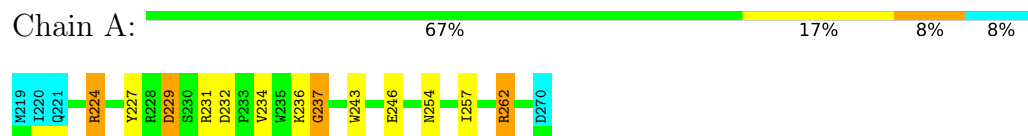


4.2 Scores per residue for each member of the ensemble

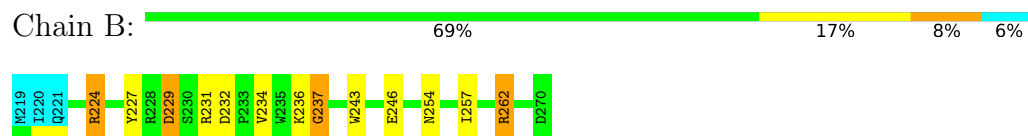
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

• Molecule 1: HIV-1 INTEGRASE

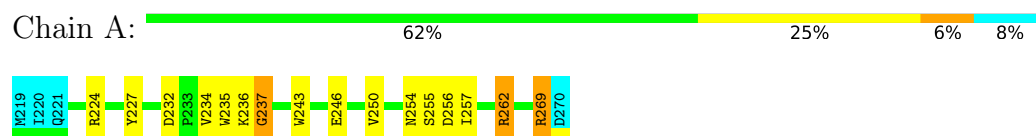


• Molecule 1: HIV-1 INTEGRASE

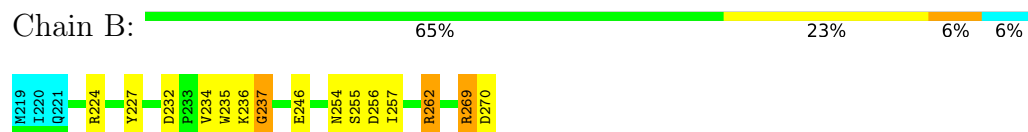


4.2.2 Score per residue for model 2

- Molecule 1: HIV-1 INTEGRASE

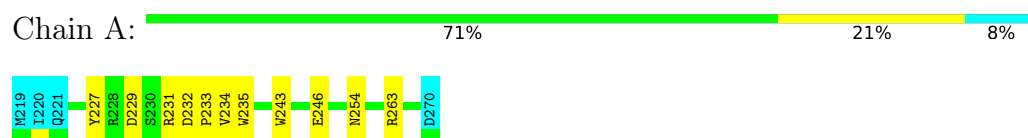


- Molecule 1: HIV-1 INTEGRASE

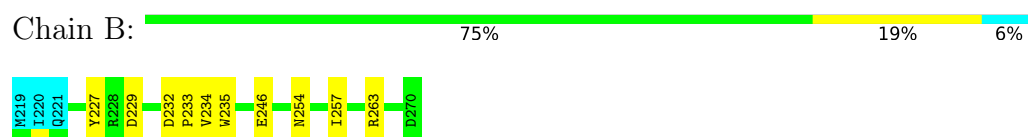


4.2.3 Score per residue for model 3

- Molecule 1: HIV-1 INTEGRASE

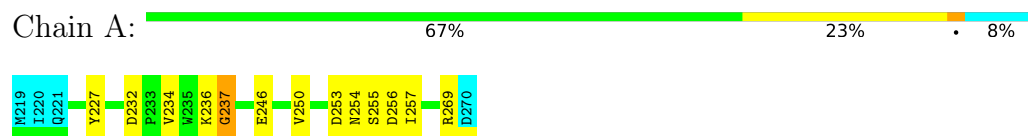


- Molecule 1: HIV-1 INTEGRASE

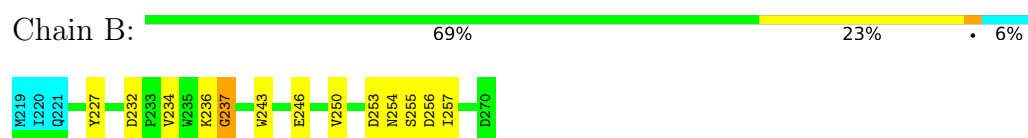


4.2.4 Score per residue for model 4

- Molecule 1: HIV-1 INTEGRASE



- Molecule 1: HIV-1 INTEGRASE



4.2.5 Score per residue for model 5

- Molecule 1: HIV-1 INTEGRASE

Chain A:  67% 23% 8%



- Molecule 1: HIV-1 INTEGRASE

Chain B:  69% 23% 6%



4.2.6 Score per residue for model 6

- Molecule 1: HIV-1 INTEGRASE

Chain A:  62% 23% 8% 8%



- Molecule 1: HIV-1 INTEGRASE

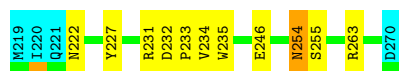
Chain B:  62% 25% 6% 6%



4.2.7 Score per residue for model 7

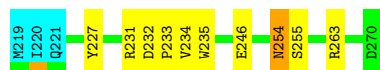
- Molecule 1: HIV-1 INTEGRASE

Chain A:  71% 19% 8%



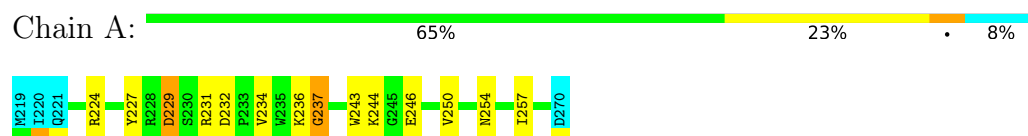
- Molecule 1: HIV-1 INTEGRASE

Chain B:  75% 17% 6%

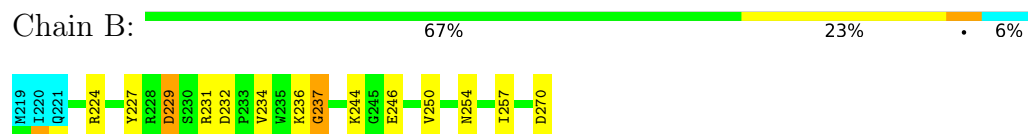


4.2.8 Score per residue for model 8

- Molecule 1: HIV-1 INTEGRASE

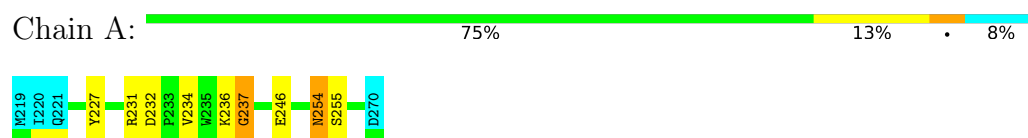


- Molecule 1: HIV-1 INTEGRASE

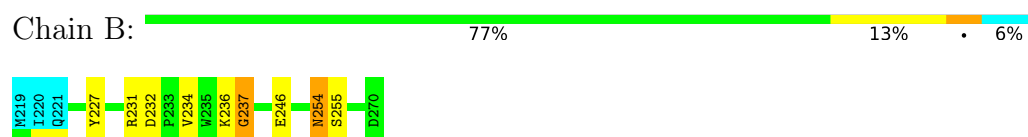


4.2.9 Score per residue for model 9

- Molecule 1: HIV-1 INTEGRASE

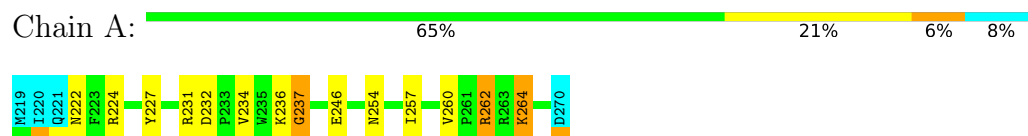


- Molecule 1: HIV-1 INTEGRASE

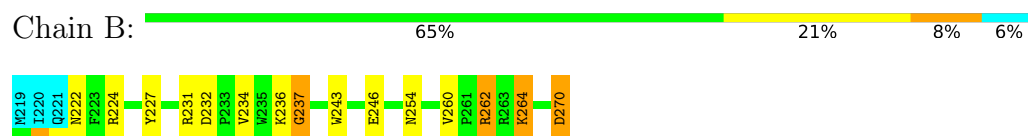


4.2.10 Score per residue for model 10

- Molecule 1: HIV-1 INTEGRASE

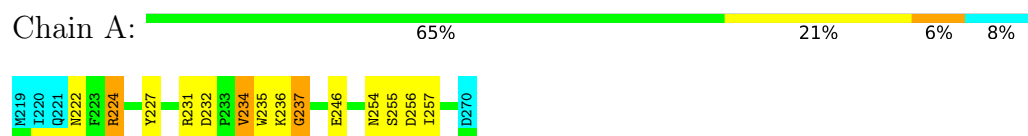


- Molecule 1: HIV-1 INTEGRASE

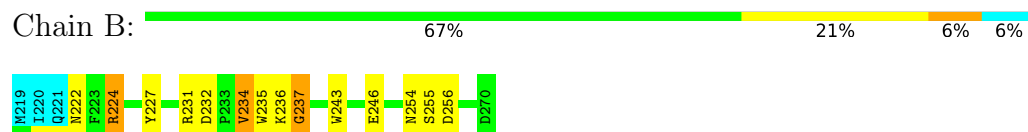


4.2.11 Score per residue for model 11

- Molecule 1: HIV-1 INTEGRASE

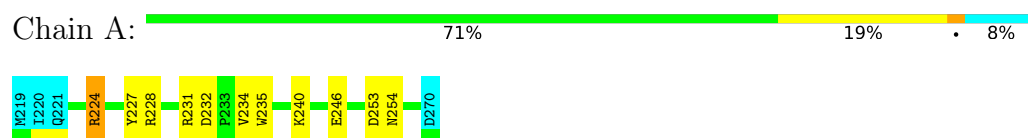


- Molecule 1: HIV-1 INTEGRASE

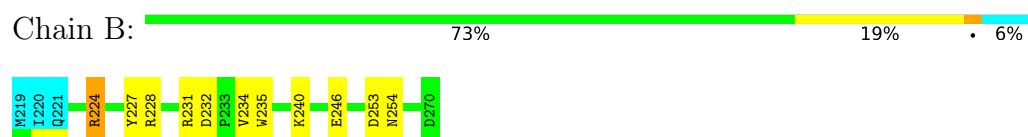


4.2.12 Score per residue for model 12

- Molecule 1: HIV-1 INTEGRASE

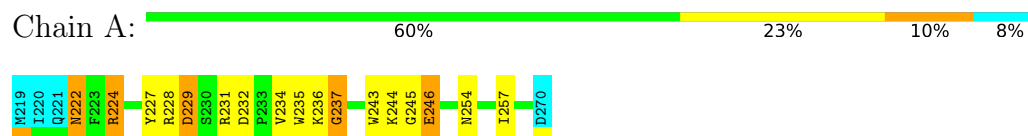


- Molecule 1: HIV-1 INTEGRASE

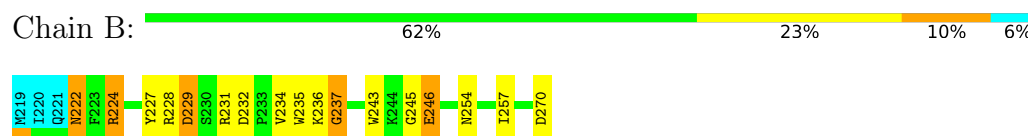


4.2.13 Score per residue for model 13

- Molecule 1: HIV-1 INTEGRASE

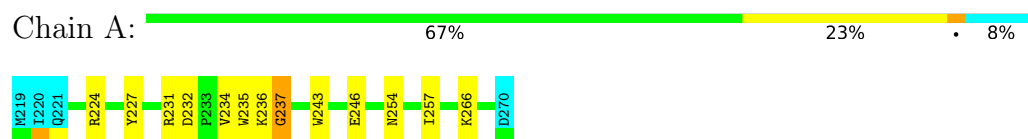


- Molecule 1: HIV-1 INTEGRASE

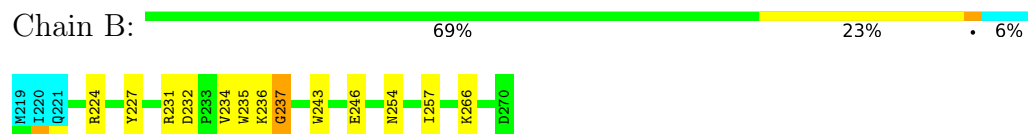


4.2.14 Score per residue for model 14

- Molecule 1: HIV-1 INTEGRASE

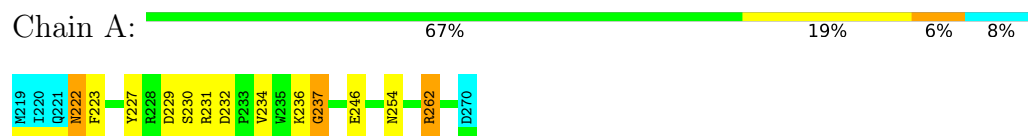


- Molecule 1: HIV-1 INTEGRASE

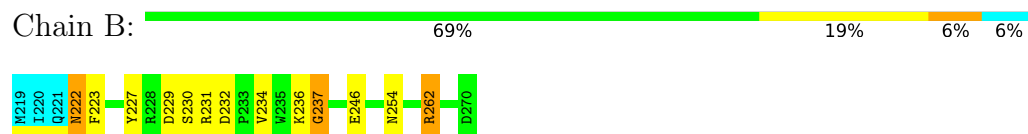


4.2.15 Score per residue for model 15

- Molecule 1: HIV-1 INTEGRASE

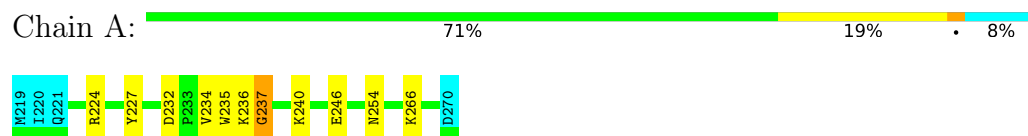


- Molecule 1: HIV-1 INTEGRASE

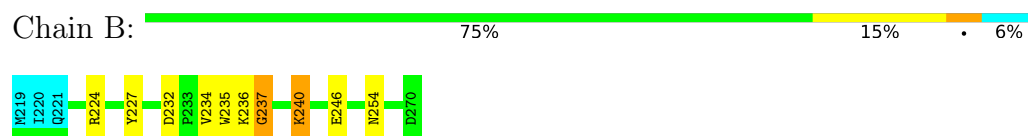


4.2.16 Score per residue for model 16 (medoid)

- Molecule 1: HIV-1 INTEGRASE

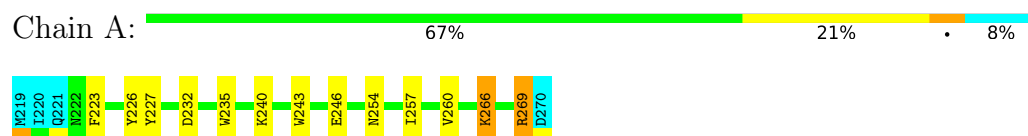


- Molecule 1: HIV-1 INTEGRASE

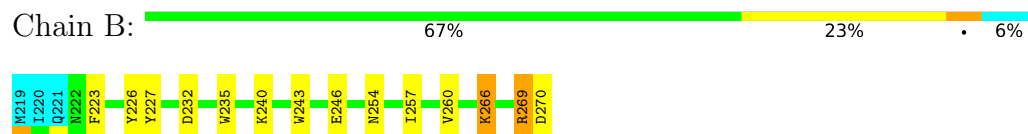


4.2.17 Score per residue for model 17

- Molecule 1: HIV-1 INTEGRASE

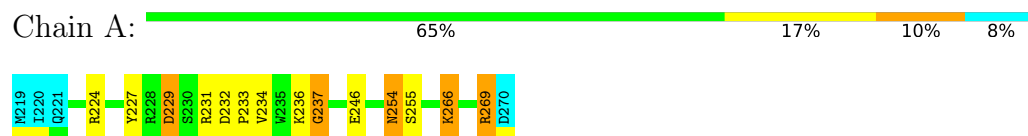


- Molecule 1: HIV-1 INTEGRASE

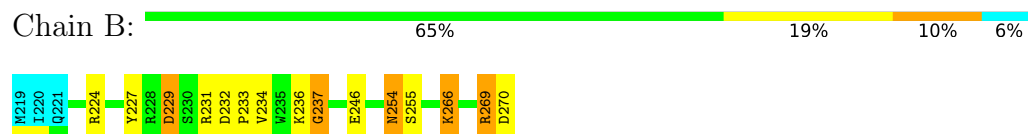


4.2.18 Score per residue for model 18

- Molecule 1: HIV-1 INTEGRASE

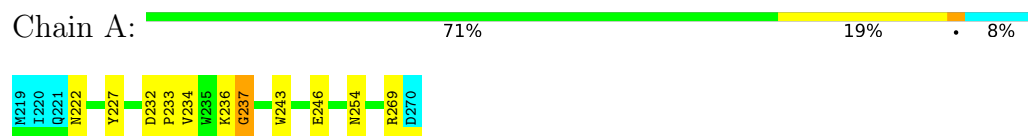


- Molecule 1: HIV-1 INTEGRASE

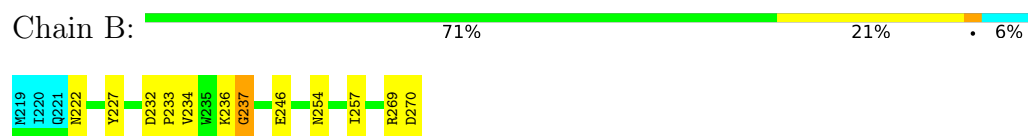


4.2.19 Score per residue for model 19

- Molecule 1: HIV-1 INTEGRASE



- Molecule 1: HIV-1 INTEGRASE



4.2.20 Score per residue for model 20

- Molecule 1: HIV-1 INTEGRASE

Chain A:  73% 17% 8%



- Molecule 1: HIV-1 INTEGRASE

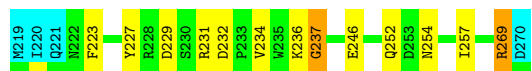
Chain B:  73% 19% 6%



4.2.21 Score per residue for model 21

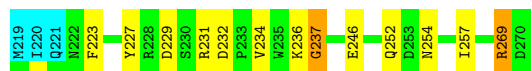
- Molecule 1: HIV-1 INTEGRASE

Chain A:  67% 21% 8%



- Molecule 1: HIV-1 INTEGRASE

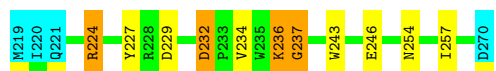
Chain B:  69% 21% 6%



4.2.22 Score per residue for model 22

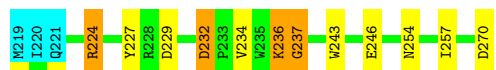
- Molecule 1: HIV-1 INTEGRASE

Chain A:  71% 13% 8% 8%



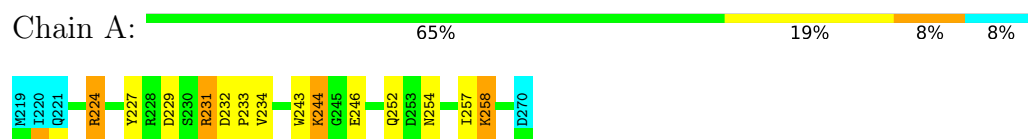
- Molecule 1: HIV-1 INTEGRASE

Chain B:  71% 15% 8% 6%

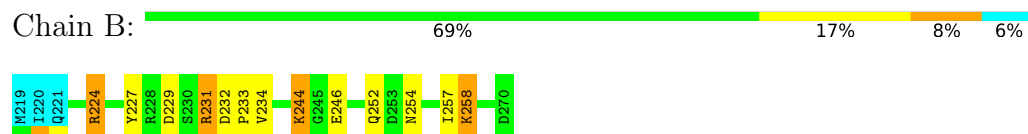


4.2.23 Score per residue for model 23

- Molecule 1: HIV-1 INTEGRASE

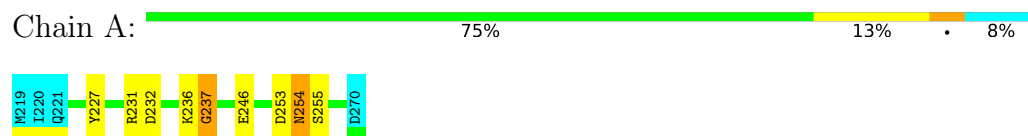


- Molecule 1: HIV-1 INTEGRASE

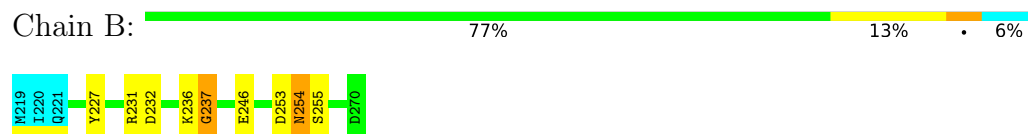


4.2.24 Score per residue for model 24

- Molecule 1: HIV-1 INTEGRASE

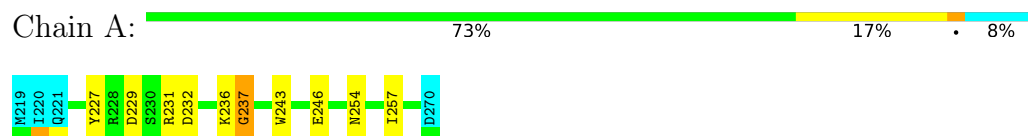


- Molecule 1: HIV-1 INTEGRASE

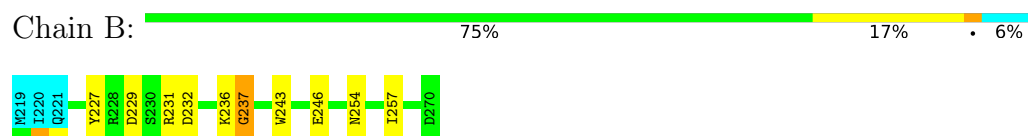


4.2.25 Score per residue for model 25

- Molecule 1: HIV-1 INTEGRASE

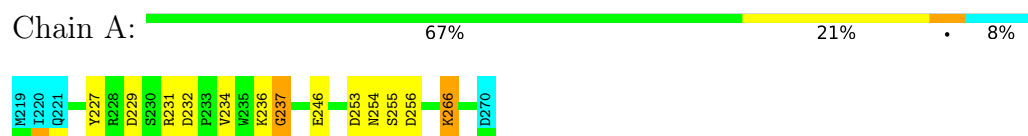


- Molecule 1: HIV-1 INTEGRASE

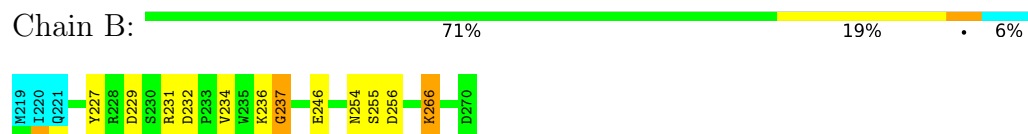


4.2.26 Score per residue for model 26

- Molecule 1: HIV-1 INTEGRASE

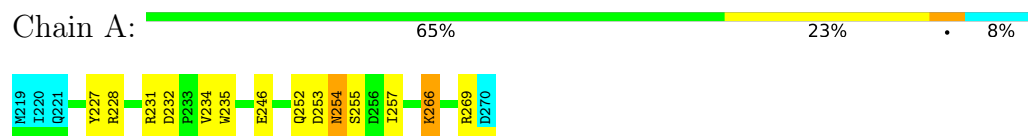


- Molecule 1: HIV-1 INTEGRASE

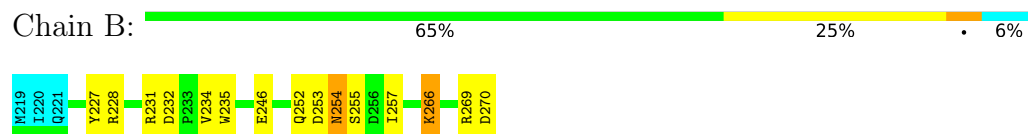


4.2.27 Score per residue for model 27

- Molecule 1: HIV-1 INTEGRASE

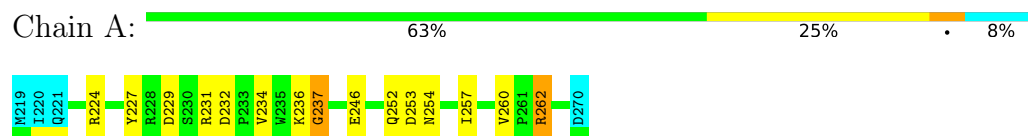


- Molecule 1: HIV-1 INTEGRASE

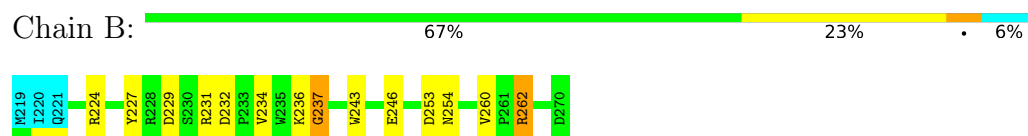


4.2.28 Score per residue for model 28

- Molecule 1: HIV-1 INTEGRASE

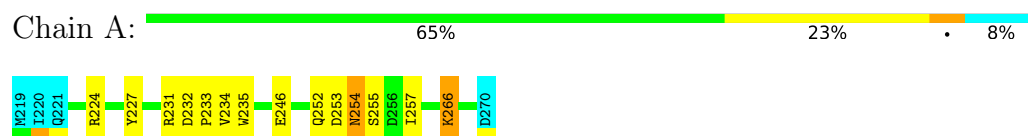


- Molecule 1: HIV-1 INTEGRASE

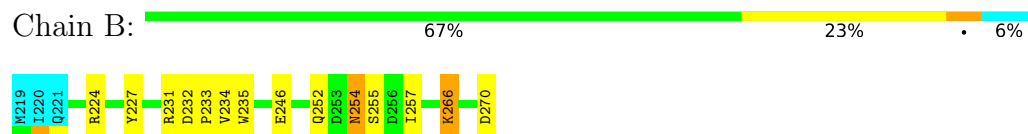


4.2.29 Score per residue for model 29

- Molecule 1: HIV-1 INTEGRASE

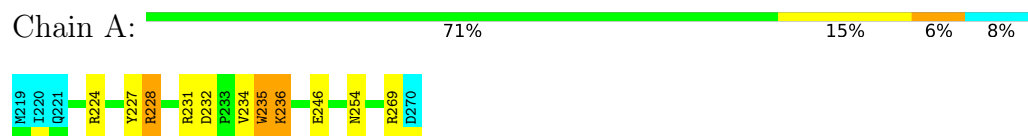


- Molecule 1: HIV-1 INTEGRASE

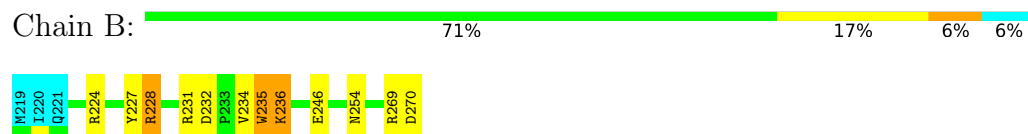


4.2.30 Score per residue for model 30

- Molecule 1: HIV-1 INTEGRASE

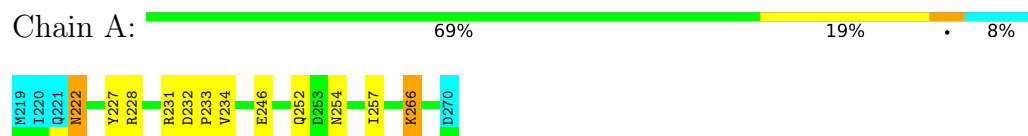


- Molecule 1: HIV-1 INTEGRASE

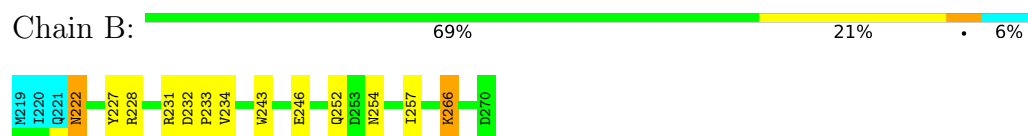


4.2.31 Score per residue for model 31

- Molecule 1: HIV-1 INTEGRASE

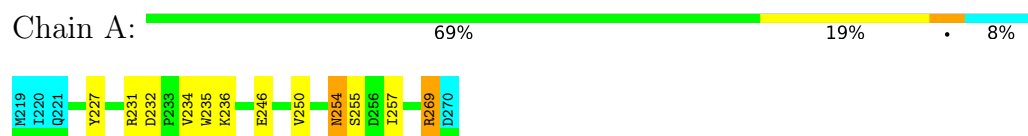


- Molecule 1: HIV-1 INTEGRASE

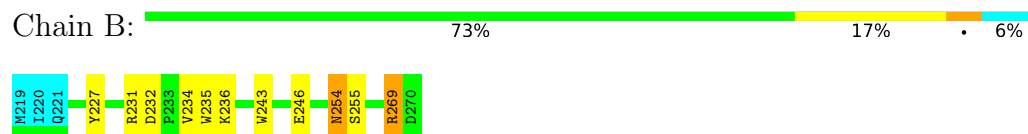


4.2.32 Score per residue for model 32

- Molecule 1: HIV-1 INTEGRASE

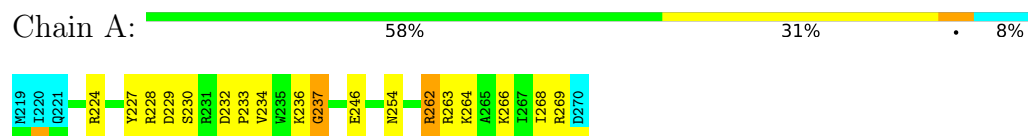


- Molecule 1: HIV-1 INTEGRASE

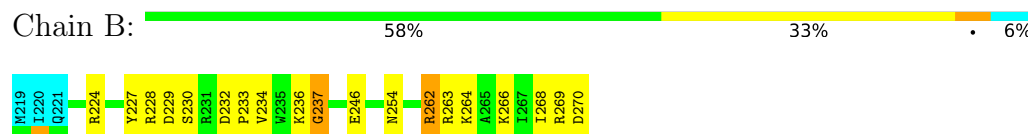


4.2.33 Score per residue for model 33

- Molecule 1: HIV-1 INTEGRASE

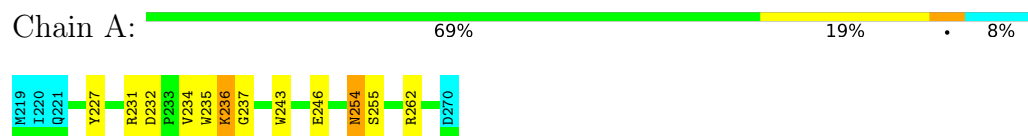


- Molecule 1: HIV-1 INTEGRASE

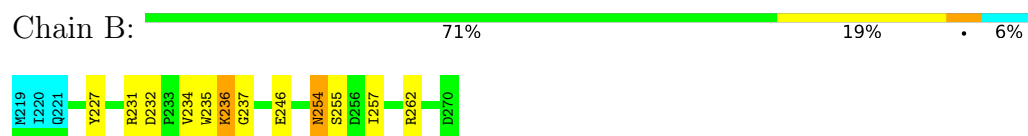


4.2.34 Score per residue for model 34

- Molecule 1: HIV-1 INTEGRASE

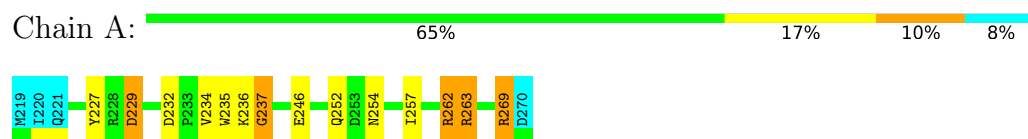


- Molecule 1: HIV-1 INTEGRASE

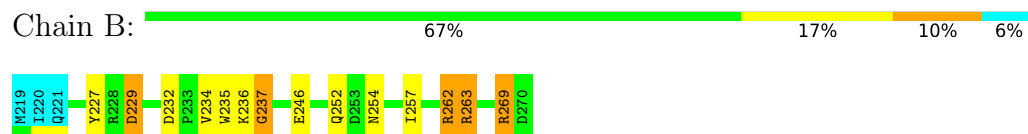


4.2.35 Score per residue for model 35

- Molecule 1: HIV-1 INTEGRASE

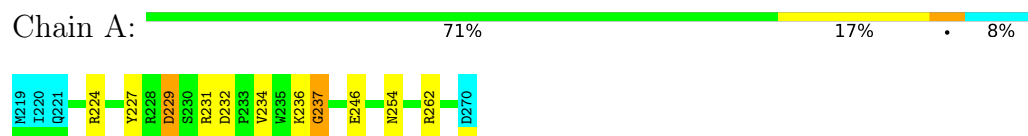


- Molecule 1: HIV-1 INTEGRASE

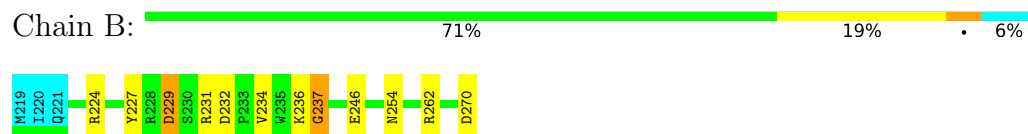


4.2.36 Score per residue for model 36

- Molecule 1: HIV-1 INTEGRASE

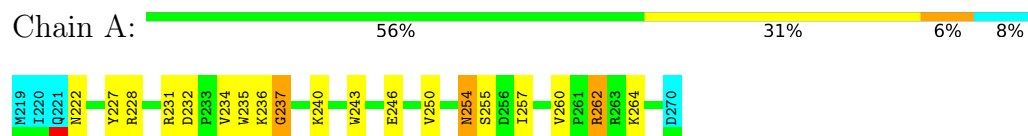


- Molecule 1: HIV-1 INTEGRASE

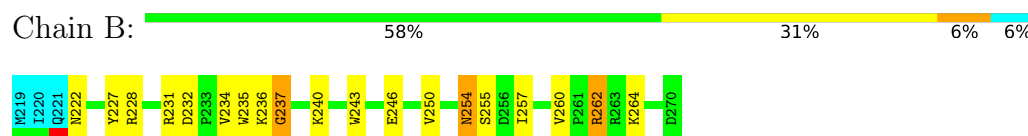


4.2.37 Score per residue for model 37

- Molecule 1: HIV-1 INTEGRASE



- Molecule 1: HIV-1 INTEGRASE



4.2.38 Score per residue for model 38

- Molecule 1: HIV-1 INTEGRASE

Chain A:  65% 25% 8%



- Molecule 1: HIV-1 INTEGRASE

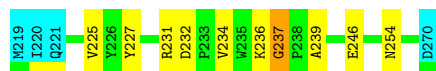
Chain B:  67% 25% 6%



4.2.39 Score per residue for model 39

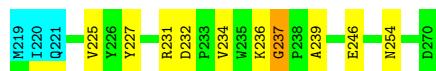
- Molecule 1: HIV-1 INTEGRASE

Chain A:  73% 17% 8%



- Molecule 1: HIV-1 INTEGRASE

Chain B:  75% 17% 6%



4.2.40 Score per residue for model 40

- Molecule 1: HIV-1 INTEGRASE

Chain A:  71% 17% 8%



- Molecule 1: HIV-1 INTEGRASE

Chain B:  73% 17% 6%



5 Refinement protocol and experimental data overview ⓘ

The models were refined using the following method: ?.

Of the ? calculated structures, 40 were deposited, based on the following criterion: ?.

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

Software name	Classification	Version
X-PLOR	refinement	3.1

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#Z>5	RMSZ	#Z>5
1	A	1.49±0.01	0±0/409 (0.0± 0.1%)	1.18±0.01	0±0/552 (0.0± 0.0%)
1	B	1.48±0.01	0±0/417 (0.0± 0.1%)	1.18±0.01	0±0/560 (0.0± 0.0%)
All	All	1.48	8/33040 (0.0%)	1.18	0/44480 (0.0%)

All unique bond outliers are listed below. They are sorted according to the Z-score of the worst occurrence in the ensemble.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	B	260	VAL	N-CA	5.55	1.50	1.45	17	4
1	A	260	VAL	N-CA	5.45	1.50	1.45	17	4

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	399	420	420	10±3
1	B	408	424	424	10±3
All	All	32280	33760	33760	784

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 12.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:224:ARG:NH2	1:B:226:TYR:CD2	0.64	2.66	38	1
1:A:224:ARG:NH2	1:A:226:TYR:CD2	0.63	2.66	38	1
1:A:254:ASN:HD22	1:A:255:SER:H	0.61	1.36	32	3
1:A:224:ARG:CZ	1:A:226:TYR:CE2	0.60	2.85	38	1
1:B:224:ARG:CZ	1:B:226:TYR:CE2	0.60	2.85	38	1
1:A:224:ARG:NH2	1:A:226:TYR:CG	0.60	2.70	38	1
1:B:254:ASN:HD22	1:B:255:SER:H	0.60	1.37	32	3
1:A:254:ASN:ND2	1:A:255:SER:H	0.59	1.96	32	4
1:B:224:ARG:NH2	1:B:226:TYR:CG	0.59	2.70	38	1
1:B:254:ASN:ND2	1:B:255:SER:H	0.58	1.96	32	4
1:A:228:ARG:NH1	1:A:235:TRP:CZ2	0.58	2.72	30	1
1:A:254:ASN:HD22	1:A:255:SER:N	0.58	1.97	29	2
1:B:254:ASN:HD22	1:B:255:SER:N	0.58	1.97	29	2
1:B:228:ARG:NH1	1:B:235:TRP:CZ2	0.58	2.71	30	1
1:B:234:VAL:HG22	1:B:235:TRP:N	0.56	2.16	2	13
1:A:234:VAL:HG22	1:A:235:TRP:N	0.56	2.16	14	13
1:A:224:ARG:HH11	1:A:224:ARG:CG	0.56	2.14	13	5
1:B:224:ARG:N	1:B:270:ASP:OD1	0.55	2.40	6	9
1:A:229:ASP:OD1	1:A:229:ASP:N	0.55	2.40	8	6
1:B:229:ASP:OD1	1:B:229:ASP:N	0.55	2.40	8	5
1:A:262:ARG:NH1	1:A:262:ARG:CG	0.54	2.71	35	5
1:B:224:ARG:CG	1:B:224:ARG:HH11	0.54	2.15	1	8
1:A:231:ARG:HH11	1:A:231:ARG:CG	0.54	2.15	39	1
1:B:224:ARG:CG	1:B:224:ARG:NH1	0.54	2.70	13	5
1:A:269:ARG:CG	1:A:269:ARG:NH1	0.54	2.71	35	6
1:B:269:ARG:HH11	1:B:269:ARG:CG	0.54	2.15	17	2
1:B:229:ASP:N	1:B:229:ASP:OD1	0.54	2.41	13	6
1:A:262:ARG:CG	1:A:262:ARG:HH11	0.54	2.16	35	5
1:B:269:ARG:CG	1:B:269:ARG:HH11	0.54	2.15	35	7
1:B:262:ARG:HH11	1:B:262:ARG:CG	0.54	2.16	35	3
1:B:231:ARG:HH11	1:B:231:ARG:CG	0.54	2.15	39	1
1:A:263:ARG:HH11	1:A:263:ARG:CG	0.54	2.16	35	2
1:B:231:ARG:NH1	1:B:231:ARG:CG	0.54	2.71	13	2
1:A:262:ARG:HH11	1:A:262:ARG:CG	0.54	2.16	10	3
1:B:263:ARG:NH1	1:B:263:ARG:CG	0.54	2.71	35	2
1:B:263:ARG:CG	1:B:263:ARG:HH11	0.54	2.16	35	3
1:B:224:ARG:NH1	1:B:224:ARG:CG	0.54	2.71	28	5
1:A:229:ASP:N	1:A:229:ASP:OD1	0.53	2.41	18	5
1:B:254:ASN:ND2	1:B:255:SER:N	0.53	2.56	29	4
1:A:224:ARG:CG	1:A:224:ARG:NH1	0.53	2.72	28	10
1:A:243:TRP:CZ3	1:B:257:ILE:CD1	0.53	2.91	5	9
1:A:269:ARG:CG	1:A:269:ARG:HH11	0.53	2.15	17	9

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:254:ASN:ND2	1:A:255:SER:N	0.53	2.56	29	4
1:B:227:TYR:OH	1:B:236:LYS:NZ	0.53	2.41	38	2
1:B:224:ARG:HH11	1:B:224:ARG:CG	0.53	2.16	22	1
1:B:269:ARG:CG	1:B:269:ARG:NH1	0.53	2.70	17	7
1:A:227:TYR:OH	1:A:236:LYS:NZ	0.53	2.41	38	2
1:A:228:ARG:HH11	1:A:228:ARG:CG	0.53	2.16	30	1
1:B:231:ARG:CG	1:B:231:ARG:HH11	0.53	2.17	13	2
1:B:262:ARG:CG	1:B:262:ARG:NH1	0.53	2.72	37	9
1:B:228:ARG:HH11	1:B:228:ARG:CG	0.52	2.17	30	1
1:B:258:LYS:CB	1:B:258:LYS:NZ	0.52	2.73	23	1
1:B:263:ARG:CG	1:B:263:ARG:NH1	0.52	2.71	7	1
1:A:269:ARG:NH1	1:A:269:ARG:CG	0.52	2.72	20	3
1:A:258:LYS:CB	1:A:258:LYS:NZ	0.52	2.73	23	1
1:B:269:ARG:NH1	1:B:269:ARG:CG	0.52	2.73	40	2
1:B:227:TYR:N	1:B:227:TYR:CD1	0.52	2.77	29	18
1:A:262:ARG:CG	1:A:262:ARG:NH1	0.52	2.71	10	4
1:A:263:ARG:CG	1:A:263:ARG:NH1	0.52	2.71	7	3
1:A:231:ARG:CG	1:A:231:ARG:HH11	0.52	2.18	13	2
1:B:231:ARG:CG	1:B:231:ARG:NH1	0.52	2.71	39	1
1:B:227:TYR:CD1	1:B:227:TYR:N	0.51	2.79	4	21
1:A:224:ARG:CG	1:A:224:ARG:HH11	0.51	2.18	28	4
1:A:227:TYR:N	1:A:227:TYR:CD1	0.51	2.77	29	18
1:B:228:ARG:HH21	1:B:233:PRO:CB	0.51	2.18	31	1
1:A:227:TYR:CD1	1:A:227:TYR:N	0.51	2.79	4	21
1:A:225:VAL:O	1:A:239:ALA:N	0.51	2.44	39	1
1:A:228:ARG:HH21	1:A:233:PRO:CB	0.51	2.18	31	1
1:B:225:VAL:O	1:B:239:ALA:N	0.51	2.44	39	1
1:B:262:ARG:CG	1:B:262:ARG:HH11	0.51	2.16	10	5
1:A:222:ASN:HD22	1:A:222:ASN:N	0.51	2.04	13	1
1:A:231:ARG:CG	1:A:231:ARG:NH1	0.51	2.71	39	3
1:A:257:ILE:CD1	1:B:243:TRP:CZ3	0.51	2.94	13	10
1:B:228:ARG:NH1	1:B:228:ARG:CG	0.51	2.72	30	1
1:A:228:ARG:NH1	1:A:228:ARG:CG	0.50	2.71	30	1
1:B:222:ASN:HD22	1:B:222:ASN:N	0.50	2.04	13	1
1:B:229:ASP:OD2	1:B:232:ASP:N	0.50	2.44	25	1
1:A:229:ASP:OD1	1:A:230:SER:N	0.50	2.45	15	2
1:A:254:ASN:OD1	1:A:255:SER:N	0.50	2.44	24	3
1:A:229:ASP:OD2	1:A:232:ASP:N	0.50	2.44	25	1
1:A:236:LYS:HZ3	1:A:237:GLY:H	0.50	1.50	34	1
1:B:264:LYS:CB	1:B:264:LYS:NZ	0.50	2.75	10	1
1:B:270:ASP:OD1	1:B:270:ASP:N	0.50	2.42	18	3

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:264:LYS:CB	1:A:264:LYS:NZ	0.49	2.75	10	1
1:B:254:ASN:OD1	1:B:255:SER:N	0.49	2.44	24	3
1:B:229:ASP:OD1	1:B:230:SER:N	0.49	2.45	15	2
1:A:234:VAL:CG2	1:A:235:TRP:N	0.49	2.76	2	8
1:B:234:VAL:CG2	1:B:235:TRP:N	0.49	2.75	2	8
1:B:252:GLN:HE21	1:B:257:ILE:CG1	0.49	2.21	21	7
1:B:228:ARG:HH21	1:B:233:PRO:CA	0.49	2.21	31	1
1:B:232:ASP:C	1:B:234:VAL:H	0.48	2.16	32	23
1:B:229:ASP:OD2	1:B:231:ARG:NH1	0.48	2.45	26	1
1:B:236:LYS:HZ3	1:B:237:GLY:H	0.48	1.50	34	1
1:A:222:ASN:OD1	1:A:222:ASN:N	0.48	2.46	37	5
1:A:229:ASP:OD2	1:A:231:ARG:NH1	0.48	2.45	26	1
1:A:250:VAL:CG1	1:A:257:ILE:CG2	0.48	2.91	37	5
1:A:232:ASP:C	1:A:234:VAL:H	0.48	2.16	32	22
1:A:232:ASP:O	1:A:232:ASP:OD1	0.48	2.31	2	7
1:A:252:GLN:HE21	1:A:257:ILE:CG1	0.48	2.21	21	8
1:A:228:ARG:HH21	1:A:233:PRO:CA	0.48	2.21	31	1
1:B:250:VAL:CG1	1:B:257:ILE:CG2	0.47	2.91	37	3
1:A:266:LYS:NZ	1:A:266:LYS:CB	0.47	2.77	29	7
1:B:266:LYS:CB	1:B:266:LYS:NZ	0.47	2.77	29	2
1:B:232:ASP:O	1:B:232:ASP:OD1	0.47	2.31	2	7
1:B:254:ASN:HD22	1:B:254:ASN:C	0.47	2.18	29	2
1:A:254:ASN:HD22	1:A:254:ASN:C	0.47	2.18	29	1
1:B:228:ARG:NH1	1:B:235:TRP:CH2	0.47	2.83	37	1
1:A:228:ARG:NH1	1:A:235:TRP:CH2	0.46	2.83	37	1
1:B:222:ASN:OD1	1:B:222:ASN:N	0.46	2.49	15	3
1:B:266:LYS:NZ	1:B:266:LYS:CB	0.46	2.78	26	5
1:B:236:LYS:O	1:B:237:GLY:O	0.46	2.33	13	25
1:A:224:ARG:CZ	1:A:226:TYR:CD2	0.46	2.99	38	1
1:B:222:ASN:N	1:B:222:ASN:OD1	0.46	2.46	37	1
1:A:224:ARG:HH22	1:A:240:LYS:NZ	0.46	2.09	12	1
1:A:236:LYS:O	1:A:237:GLY:O	0.46	2.33	13	26
1:B:258:LYS:HZ2	1:B:258:LYS:HB2	0.46	1.71	23	1
1:B:264:LYS:CB	1:B:264:LYS:HZ2	0.45	2.24	10	1
1:B:224:ARG:HH22	1:B:240:LYS:NZ	0.45	2.09	12	1
1:B:224:ARG:CZ	1:B:226:TYR:CD2	0.45	2.99	38	1
1:A:231:ARG:O	1:A:232:ASP:CG	0.45	2.60	24	29
1:B:227:TYR:OH	1:B:236:LYS:CE	0.45	2.65	30	3
1:A:224:ARG:NH2	1:A:226:TYR:CE2	0.45	2.85	38	1
1:B:231:ARG:O	1:B:232:ASP:CG	0.45	2.60	24	28
1:A:257:ILE:CD1	1:B:243:TRP:CH2	0.45	3.00	17	6

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:229:ASP:OD1	1:A:232:ASP:O	0.45	2.35	22	3
1:B:229:ASP:OD1	1:B:232:ASP:O	0.45	2.35	22	3
1:A:227:TYR:OH	1:A:236:LYS:CE	0.45	2.64	40	3
1:A:236:LYS:O	1:A:237:GLY:C	0.45	2.60	8	24
1:B:236:LYS:O	1:B:237:GLY:C	0.45	2.60	8	24
1:B:234:VAL:O	1:B:234:VAL:HG23	0.45	2.12	10	2
1:A:234:VAL:HG23	1:A:234:VAL:O	0.45	2.12	10	2
1:A:264:LYS:CB	1:A:264:LYS:HZ2	0.45	2.24	10	1
1:B:270:ASP:N	1:B:270:ASP:OD1	0.45	2.49	20	1
1:A:234:VAL:HG12	1:A:235:TRP:N	0.44	2.27	16	3
1:A:228:ARG:HH21	1:A:235:TRP:NE1	0.44	2.10	12	1
1:B:228:ARG:HH21	1:B:235:TRP:NE1	0.44	2.10	12	1
1:A:231:ARG:CD	1:A:231:ARG:N	0.44	2.81	6	1
1:B:231:ARG:N	1:B:231:ARG:CD	0.44	2.81	6	1
1:B:234:VAL:HG12	1:B:235:TRP:N	0.44	2.27	16	3
1:B:224:ARG:NH2	1:B:226:TYR:CE2	0.44	2.85	38	1
1:A:222:ASN:N	1:A:222:ASN:ND2	0.44	2.65	13	1
1:A:255:SER:O	1:A:256:ASP:OD1	0.43	2.36	2	7
1:B:231:ARG:O	1:B:232:ASP:OD1	0.43	2.37	39	5
1:B:234:VAL:HG23	1:B:234:VAL:O	0.43	2.12	15	1
1:B:269:ARG:O	1:B:270:ASP:OD1	0.43	2.36	27	2
1:A:234:VAL:O	1:A:234:VAL:HG23	0.43	2.13	26	1
1:A:240:LYS:N	1:A:240:LYS:CD	0.43	2.81	37	1
1:A:231:ARG:O	1:A:232:ASP:OD1	0.43	2.37	5	5
1:A:228:ARG:CZ	1:A:228:ARG:CB	0.43	2.97	27	2
1:B:222:ASN:N	1:B:222:ASN:ND2	0.43	2.65	13	1
1:B:224:ARG:HH22	1:B:225:VAL:C	0.43	2.21	38	1
1:B:255:SER:O	1:B:256:ASP:OD1	0.43	2.37	2	7
1:A:227:TYR:OH	1:A:253:ASP:CG	0.43	2.62	27	7
1:B:227:TYR:OH	1:B:253:ASP:CG	0.43	2.62	4	5
1:B:240:LYS:N	1:B:240:LYS:CD	0.43	2.81	37	1
1:A:243:TRP:CH2	1:B:257:ILE:CD1	0.43	3.02	34	7
1:B:232:ASP:C	1:B:234:VAL:N	0.43	2.77	32	3
1:A:264:LYS:HZ2	1:A:264:LYS:HB2	0.43	1.74	10	1
1:B:228:ARG:HH21	1:B:264:LYS:NZ	0.43	2.12	33	1
1:A:228:ARG:NH2	1:A:264:LYS:CD	0.43	2.82	37	1
1:A:224:ARG:HH22	1:A:225:VAL:C	0.43	2.21	38	1
1:B:222:ASN:C	1:B:270:ASP:OD2	0.42	2.62	10	1
1:A:245:GLY:O	1:A:246:GLU:O	0.42	2.37	6	2
1:B:245:GLY:O	1:B:246:GLU:O	0.42	2.37	6	2
1:B:264:LYS:HZ2	1:B:264:LYS:HB2	0.42	1.74	10	1

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:227:TYR:O	1:B:236:LYS:N	0.42	2.53	38	1
1:B:222:ASN:O	1:B:270:ASP:OXT	0.42	2.38	19	1
1:A:263:ARG:CG	1:A:263:ARG:HH11	0.42	2.26	33	1
1:A:232:ASP:C	1:A:234:VAL:N	0.42	2.77	32	3
1:B:231:ARG:C	1:B:232:ASP:CG	0.42	2.87	14	1
1:A:258:LYS:HZ2	1:A:258:LYS:HB2	0.42	1.73	23	1
1:A:227:TYR:O	1:A:236:LYS:N	0.42	2.53	38	1
1:B:228:ARG:CZ	1:B:228:ARG:CB	0.42	2.97	27	2
1:B:228:ARG:NH2	1:B:264:LYS:CD	0.42	2.82	37	1
1:B:258:LYS:CB	1:B:258:LYS:HZ2	0.42	2.28	23	1
1:B:234:VAL:O	1:B:235:TRP:O	0.42	2.38	30	1
1:A:228:ARG:HH21	1:A:264:LYS:NZ	0.42	2.12	33	1
1:A:231:ARG:C	1:A:232:ASP:CG	0.41	2.87	14	1
1:A:232:ASP:OD2	1:A:234:VAL:CG2	0.41	2.68	16	1
1:B:232:ASP:OD2	1:B:234:VAL:CG2	0.41	2.68	16	1
1:A:234:VAL:O	1:A:235:TRP:O	0.41	2.39	30	1
1:A:244:LYS:CB	1:A:244:LYS:NZ	0.41	2.82	23	1
1:B:240:LYS:HZ2	1:B:240:LYS:HB2	0.41	1.76	16	1
1:B:223:PHE:CA	1:B:270:ASP:OD1	0.41	2.68	17	1
1:B:254:ASN:OD1	1:B:254:ASN:N	0.41	2.53	6	1
1:A:226:TYR:CD1	1:A:226:TYR:N	0.41	2.89	17	1
1:A:228:ARG:NH2	1:A:233:PRO:O	0.41	2.54	31	1
1:B:228:ARG:NH2	1:B:233:PRO:CB	0.41	2.84	31	1
1:B:229:ASP:CG	1:B:231:ARG:NH1	0.41	2.79	15	1
1:B:228:ARG:NH2	1:B:233:PRO:O	0.41	2.54	31	1
1:B:244:LYS:NZ	1:B:244:LYS:CB	0.40	2.83	23	1
1:A:224:ARG:NH2	1:A:268:ILE:HG21	0.40	2.31	33	1
1:B:226:TYR:CD1	1:B:226:TYR:N	0.40	2.89	17	1
1:B:224:ARG:NH2	1:B:268:ILE:HG21	0.40	2.31	33	1
1:A:244:LYS:C	1:A:244:LYS:CD	0.40	2.94	13	1
1:A:235:TRP:CZ3	1:A:266:LYS:NZ	0.40	2.89	16	1
1:A:228:ARG:NH2	1:A:233:PRO:CB	0.40	2.84	31	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	48/52 (92%)	40±1 (82±2%)	6±1 (11±2%)	3±0 (6±1%)	2	19
1	B	48/52 (92%)	40±1 (82±2%)	6±1 (11±2%)	3±0 (6±1%)	2	19
All	All	3840/4160 (92%)	3168 (82%)	440 (11%)	232 (6%)	2	19

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	246	GLU	40
1	B	246	GLU	40
1	A	254	ASN	39
1	B	254	ASN	39
1	A	237	GLY	27
1	B	237	GLY	27
1	A	233	PRO	7
1	B	233	PRO	7
1	A	235	TRP	3
1	B	235	TRP	3

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	42/46 (91%)	40±1 (96±3%)	2±1 (4±3%)	30	81
1	B	43/46 (93%)	41±1 (96±3%)	2±1 (4±3%)	31	81
All	All	3400/3680 (92%)	3273 (96%)	127 (4%)	31	81

All 35 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	262	ARG	9
1	B	262	ARG	9
1	A	269	ARG	9
1	B	269	ARG	9
1	A	266	LYS	9
1	B	266	LYS	9

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Mol	Chain	Res	Type	Models (Total)
1	A	224	ARG	8
1	B	224	ARG	8
1	A	229	ASP	6
1	B	229	ASP	6
1	A	236	LYS	4
1	B	236	LYS	4
1	A	222	ASN	3
1	B	222	ASN	3
1	A	263	ARG	2
1	B	263	ARG	2
1	A	254	ASN	2
1	B	254	ASN	2
1	A	244	LYS	2
1	B	244	LYS	2
1	A	234	VAL	2
1	B	234	VAL	2
1	A	240	LYS	2
1	B	240	LYS	2
1	A	264	LYS	1
1	B	264	LYS	1
1	B	270	ASP	1
1	A	232	ASP	1
1	B	232	ASP	1
1	A	231	ARG	1
1	A	258	LYS	1
1	B	231	ARG	1
1	B	258	LYS	1
1	A	228	ARG	1
1	B	228	ARG	1

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.6 Ligand geometry

There are no ligands in this entry.

6.7 Other polymers

There are no such molecules in this entry.

6.8 Polymer linkage issues

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided