



Full wwPDB NMR Structure Validation Report ⓘ

Mar 8, 2026 – 06:02 AM UTC

PDB ID : 6TPH / pdb_00006tph
BMRB ID : 34465
Title : Structure of a protein-RNA complex by ssNMR
Authors : Mumdooh, A.; Marchanka, A.; Carlomagno, T.
Deposited on : 2019-12-13

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

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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
BMRB Restraints Analysis : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

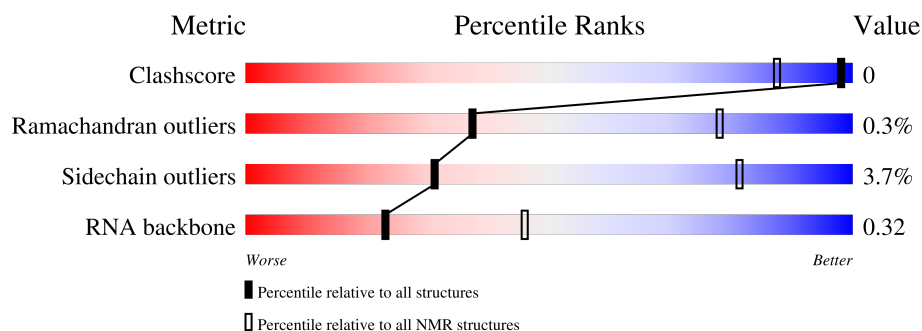
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLID-STATE NMR

The overall completeness of chemical shifts assignment is 20%.

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
RNA backbone	8273	777

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	123	
2	B	26	

2 Ensemble composition and analysis

This entry contains 10 models. Model 7 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *closest to the average*.

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:5-A:124 (120)	0.53	7

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

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

The models can be grouped into 2 clusters and 4 single-model clusters were found.

Cluster number	Models
1	3, 4, 7, 10
2	2, 6
Single-model clusters	1; 5; 8; 9

3 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 2647 atoms, of which 1227 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called 50S ribosomal protein L7Ae.

Mol	Chain	Residues	Atoms						Trace
1	A	121	Total	C	H	N	O	S	0
			1904	591	978	153	179	3	

- Molecule 2 is a RNA chain called RNA (26-MER).

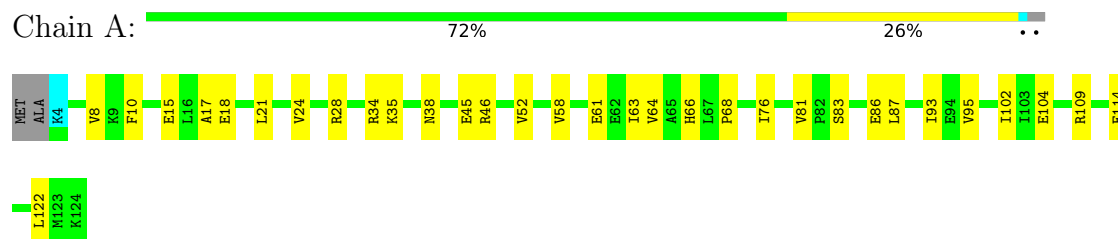
Mol	Chain	Residues	Atoms						Trace
2	B	23	Total	C	H	N	O	P	0
			743	222	249	95	155	22	

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: 50S ribosomal protein L7Ae



- Molecule 2: RNA (26-MER)

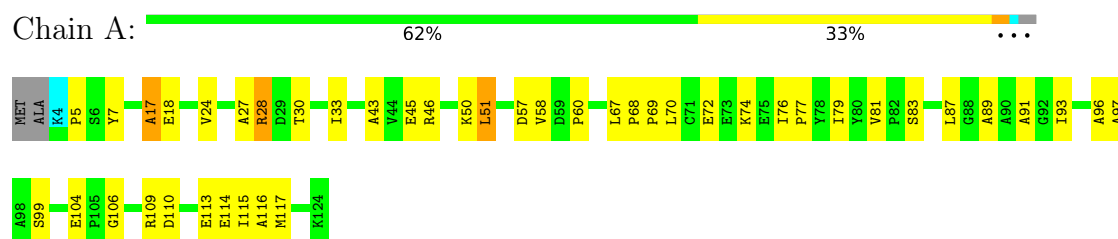


4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: 50S ribosomal protein L7Ae

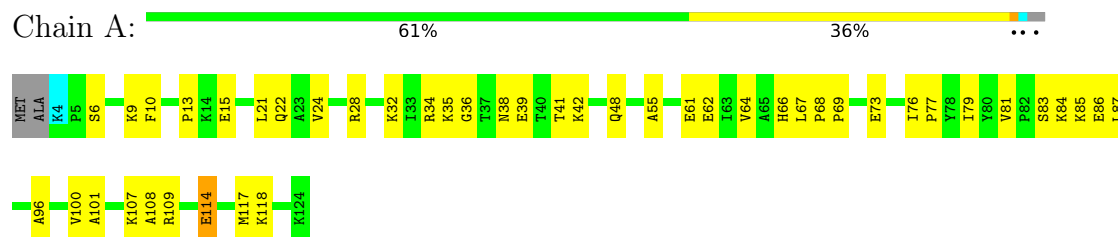


- Molecule 2: RNA (26-MER)

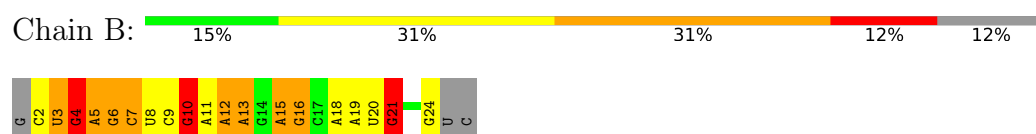


4.2.2 Score per residue for model 2

- Molecule 1: 50S ribosomal protein L7Ae

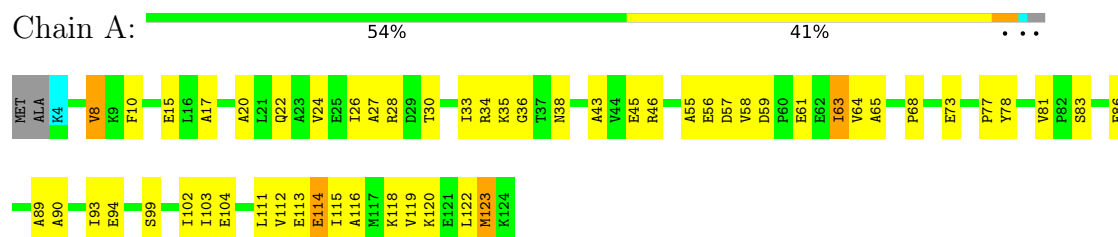


- Molecule 2: RNA (26-MER)



4.2.3 Score per residue for model 3

- Molecule 1: 50S ribosomal protein L7Ae

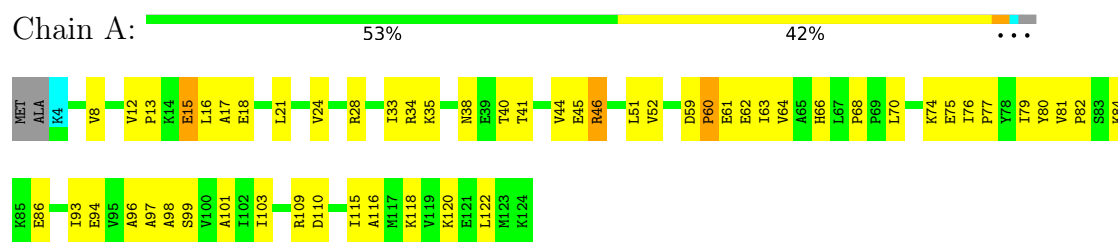


- Molecule 2: RNA (26-MER)



4.2.4 Score per residue for model 4

- Molecule 1: 50S ribosomal protein L7Ae

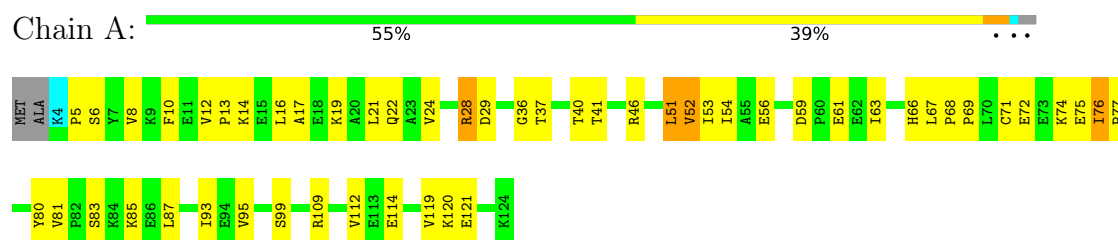


- Molecule 2: RNA (26-MER)



4.2.5 Score per residue for model 5

- Molecule 1: 50S ribosomal protein L7Ae

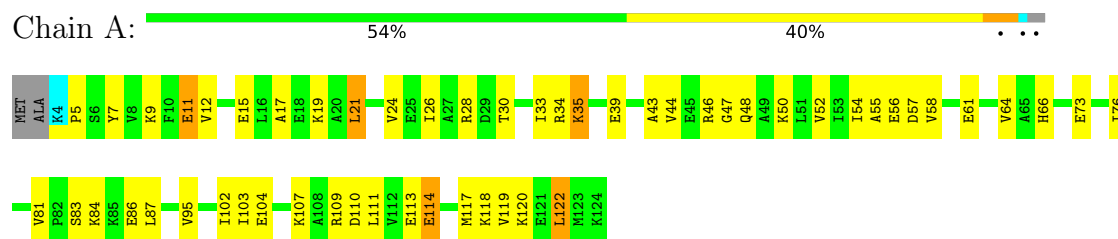


- Molecule 2: RNA (26-MER)



4.2.6 Score per residue for model 6

- Molecule 1: 50S ribosomal protein L7Ae

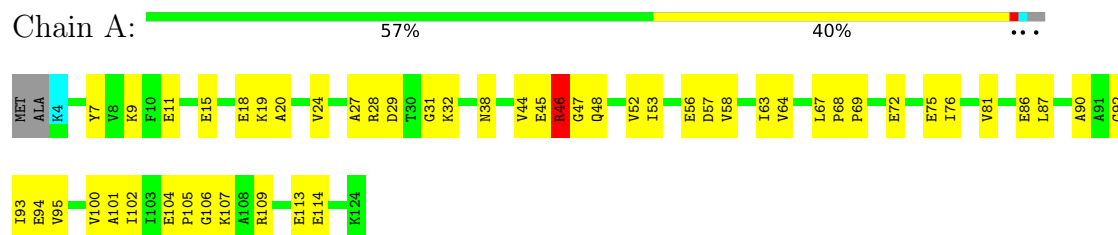


- Molecule 2: RNA (26-MER)



4.2.7 Score per residue for model 7 (medoid)

- Molecule 1: 50S ribosomal protein L7Ae

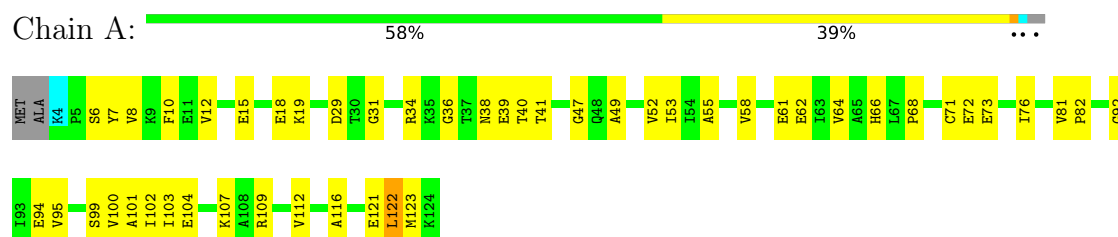


- Molecule 2: RNA (26-MER)

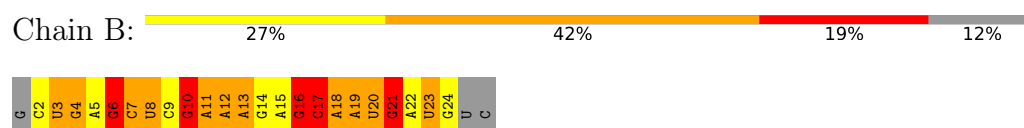


4.2.8 Score per residue for model 8

- Molecule 1: 50S ribosomal protein L7Ae

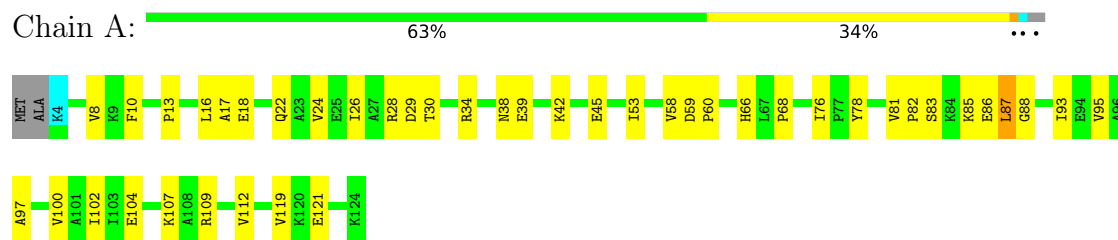


- Molecule 2: RNA (26-MER)



4.2.9 Score per residue for model 9

- Molecule 1: 50S ribosomal protein L7Ae

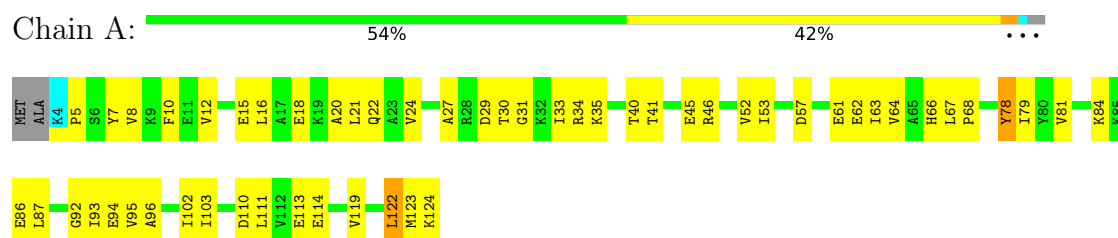


- Molecule 2: RNA (26-MER)

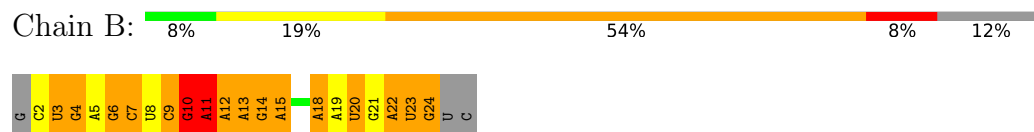


4.2.10 Score per residue for model 10

- Molecule 1: 50S ribosomal protein L7Ae



- Molecule 2: RNA (26-MER)



5 Refinement protocol and experimental data overview

The models were refined using the following method: *na*.

Of the 200 calculated structures, 10 were deposited, based on the following criterion: *target function*.

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

Software name	Classification	Version
Amber	refinement	
HADDOCK	structure calculation	

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	425
Number of shifts mapped to atoms	425
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	20%

Note: This is a solid-state NMR structure, where hydrogen atoms are typically not assigned a chemical shift value, which may lead to lower completeness of assignment measure.

6 Model quality

6.1 Standard geometry

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	2.03±0.04	22±4/929 (2.3± 0.4%)	2.28±0.03	47±6/1252 (3.7± 0.5%)
2	B	1.92±0.04	8±2/554 (1.4± 0.4%)	2.04±0.11	31±8/863 (3.6± 0.9%)
All	All	1.99	294/14830 (2.0%)	2.18	778/21150 (3.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0.0±0.0	1.7±1.1
2	B	0.0±0.0	14.0±2.4
All	All	0	157

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	A	31	GLY	N-CA	11.91	1.54	1.44	10	2
1	A	63	ILE	CA-C	10.34	1.64	1.52	7	2
1	A	32	LYS	N-CA	9.30	1.57	1.46	2	1
1	A	31	GLY	CA-C	9.10	1.64	1.51	8	2
1	A	28	ARG	N-CA	8.65	1.56	1.46	2	2
1	A	75	GLU	C-N	8.34	1.39	1.33	5	2
1	A	57	ASP	CA-C	8.24	1.63	1.53	10	1
1	A	67	LEU	N-CA	8.23	1.52	1.46	10	1
1	A	33	ILE	CA-C	8.09	1.62	1.52	1	2
1	A	106	GLY	CA-C	7.98	1.63	1.51	7	1
1	A	66	HIS	CG-ND1	7.97	1.47	1.38	10	1
1	A	66	HIS	CB-CG	7.88	1.61	1.50	8	1
2	B	7	C	C1'-N1	7.80	1.60	1.48	3	1
1	A	45	GLU	CA-C	7.79	1.63	1.52	4	2

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	40	THR	N-CA	7.71	1.55	1.46	8	2
1	A	58	VAL	CA-C	7.63	1.62	1.52	1	1
1	A	100	VAL	N-CA	7.56	1.54	1.46	7	1
1	A	59	ASP	CA-C	7.55	1.62	1.52	9	1
2	B	22	A	P-O5'	7.54	1.71	1.59	3	1
1	A	68	PRO	CA-C	7.52	1.59	1.52	2	4
2	B	20	U	C5'-C4'	7.50	1.62	1.51	4	1
2	B	4	G	C1'-N9	7.49	1.58	1.47	2	2
1	A	76	ILE	CA-C	7.46	1.59	1.52	7	2
2	B	18	A	C3'-O3'	7.44	1.53	1.42	8	1
1	A	114	GLU	N-CA	7.40	1.55	1.46	10	3
1	A	17	ALA	N-CA	7.37	1.55	1.46	1	2
2	B	19	A	C2'-O2'	7.29	1.52	1.41	2	1
1	A	94	GLU	N-CA	7.29	1.55	1.46	10	3
1	A	64	VAL	N-CA	7.26	1.52	1.46	8	4
1	A	17	ALA	CA-C	7.25	1.62	1.52	5	1
1	A	77	PRO	C-O	7.25	1.32	1.23	5	2
1	A	48	GLN	N-CA	7.21	1.55	1.46	6	1
1	A	97	ALA	N-CA	7.18	1.54	1.46	4	1
1	A	67	LEU	CA-C	7.16	1.60	1.52	5	1
1	A	116	ALA	N-CA	-7.10	1.38	1.46	4	1
2	B	10	G	C2-N2	-7.07	1.20	1.34	10	1
1	A	36	GLY	N-CA	7.02	1.53	1.45	5	2
1	A	44	VAL	CA-C	6.92	1.62	1.52	7	1
2	B	17	C	C4'-O4'	-6.91	1.35	1.45	9	1
2	B	14	G	O5'-C5'	-6.85	1.32	1.42	3	1
1	A	6	SER	C-O	6.82	1.32	1.24	5	1
2	B	22	A	C3'-C2'	6.81	1.62	1.52	8	2
1	A	117	MET	N-CA	6.81	1.54	1.46	1	1
2	B	13	A	C4'-O4'	-6.78	1.35	1.45	8	1
2	B	9	C	O3'-P	6.76	1.71	1.61	10	3
1	A	107	LYS	C-O	-6.71	1.15	1.24	7	1
1	A	27	ALA	N-CA	6.68	1.54	1.46	7	1
1	A	39	GLU	CA-C	6.67	1.61	1.52	2	2
1	A	49	ALA	CA-C	6.67	1.61	1.52	8	1
1	A	109	ARG	CZ-NH2	-6.65	1.24	1.33	1	1
1	A	77	PRO	CA-C	6.65	1.59	1.52	1	2
1	A	39	GLU	C-O	-6.61	1.16	1.24	2	1
1	A	96	ALA	CA-C	6.60	1.60	1.52	1	1
2	B	18	A	O5'-C5'	6.59	1.52	1.42	5	2
1	A	8	VAL	CA-C	6.59	1.60	1.52	9	2

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	51	LEU	N-CA	6.59	1.54	1.46	4	1
1	A	20	ALA	N-CA	6.58	1.54	1.46	3	1
2	B	15	A	C4'-O4'	-6.56	1.35	1.45	10	1
1	A	10	PHE	CA-C	6.51	1.60	1.52	2	2
1	A	79	ILE	N-CA	6.49	1.54	1.46	1	1
2	B	23	U	C5'-C4'	6.48	1.60	1.51	10	2
1	A	43	ALA	C-O	-6.47	1.16	1.24	3	1
1	A	101	ALA	CA-C	6.46	1.60	1.52	8	1
1	A	85	LYS	C-O	-6.45	1.16	1.24	9	1
2	B	2	C	O3'-P	-6.44	1.51	1.61	10	1
2	B	7	C	P-O5'	-6.44	1.50	1.59	6	1
1	A	81	VAL	CA-C	6.42	1.59	1.52	5	3
1	A	62	GLU	N-CA	6.40	1.54	1.46	4	1
1	A	119	VAL	C-O	-6.40	1.16	1.24	3	1
1	A	82	PRO	CA-C	6.39	1.61	1.52	4	3
2	B	21	G	C1'-N9	6.37	1.56	1.47	7	1
2	B	9	C	C2'-O2'	-6.37	1.32	1.42	9	1
2	B	14	G	C5'-C4'	6.36	1.60	1.51	5	1
1	A	52	VAL	N-CA	6.33	1.53	1.46	7	2
1	A	13	PRO	N-CA	-6.33	1.39	1.47	5	1
2	B	19	A	O3'-P	-6.32	1.51	1.61	1	1
2	B	10	G	C5'-C4'	6.32	1.60	1.51	3	1
2	B	19	A	C1'-N9	6.26	1.56	1.47	6	1
1	A	63	ILE	C-N	6.26	1.39	1.33	7	1
2	B	4	G	C5'-C4'	6.25	1.60	1.51	8	1
1	A	18	GLU	N-CA	6.24	1.54	1.46	1	1
1	A	8	VAL	N-CA	6.22	1.53	1.46	10	2
1	A	34	ARG	CZ-NH2	-6.20	1.25	1.33	8	2
1	A	122	LEU	N-CA	6.20	1.54	1.45	6	1
1	A	123	MET	CA-C	6.19	1.61	1.52	8	1
1	A	74	LYS	C-O	-6.19	1.15	1.24	4	1
2	B	17	C	C2'-O2'	6.18	1.51	1.42	5	1
1	A	12	VAL	CA-C	6.13	1.59	1.52	10	2
1	A	107	LYS	CA-C	6.13	1.61	1.52	9	2
1	A	24	VAL	N-CA	6.12	1.54	1.46	10	1
2	B	13	A	C3'-O3'	6.12	1.51	1.42	3	1
2	B	12	A	C4'-O4'	-6.08	1.36	1.45	10	1
1	A	103	ILE	CA-C	6.08	1.60	1.52	6	2
1	A	81	VAL	N-CA	6.07	1.54	1.46	6	4
1	A	104	GLU	C-N	6.05	1.41	1.33	6	2
2	B	21	G	O3'-P	-6.04	1.52	1.61	4	2

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	61	GLU	CA-C	6.04	1.60	1.52	4	1
1	A	95	VAL	N-CA	6.02	1.54	1.46	7	1
1	A	63	ILE	N-CA	6.02	1.53	1.46	10	1
2	B	16	G	C5'-C4'	6.02	1.60	1.51	1	1
2	B	6	G	C3'-O3'	-6.01	1.33	1.42	3	1
2	B	7	C	C4'-O4'	-6.01	1.36	1.45	1	2
1	A	122	LEU	CA-C	5.99	1.60	1.52	10	3
1	A	109	ARG	N-CA	5.99	1.53	1.46	7	2
2	B	6	G	P-O5'	5.97	1.68	1.59	10	2
1	A	42	LYS	CA-C	5.95	1.60	1.52	9	1
1	A	29	ASP	CA-C	5.90	1.59	1.52	10	1
2	B	5	A	C3'-C2'	5.89	1.61	1.52	6	1
1	A	58	VAL	C-O	5.88	1.29	1.24	6	1
1	A	41	THR	CA-C	5.87	1.60	1.52	2	1
2	B	11	A	O4'-C1'	5.87	1.50	1.41	7	2
1	A	99	SER	CA-C	5.86	1.59	1.52	3	1
1	A	30	THR	CB-OG1	-5.86	1.34	1.43	6	2
1	A	28	ARG	CD-NE	-5.86	1.38	1.46	6	1
2	B	10	G	C2'-O2'	5.85	1.50	1.41	2	1
1	A	72	GLU	N-CA	5.85	1.53	1.46	5	1
1	A	58	VAL	N-CA	5.82	1.53	1.46	3	1
1	A	34	ARG	N-CA	5.81	1.53	1.45	6	1
1	A	121	GLU	CA-C	5.81	1.60	1.52	9	2
1	A	53	ILE	N-CA	5.80	1.53	1.46	8	1
1	A	35	LYS	N-CA	5.77	1.53	1.45	6	1
1	A	111	LEU	CA-C	5.77	1.60	1.52	3	1
2	B	17	C	C5'-C4'	5.76	1.59	1.51	1	3
1	A	70	LEU	N-CA	5.75	1.53	1.46	4	1
1	A	62	GLU	C-O	-5.74	1.16	1.24	10	1
1	A	75	GLU	CA-C	5.72	1.60	1.53	7	1
1	A	19	LYS	CA-C	5.71	1.60	1.52	6	1
1	A	117	MET	CA-C	5.71	1.60	1.52	6	2
1	A	61	GLU	C-O	-5.69	1.16	1.24	6	1
1	A	103	ILE	N-CA	5.67	1.53	1.46	10	1
2	B	8	U	C1'-N1	5.66	1.56	1.48	9	1
2	B	8	U	O3'-P	5.65	1.69	1.61	7	1
1	A	58	VAL	CA-CB	5.65	1.60	1.54	1	1
1	A	53	ILE	CA-C	5.64	1.59	1.52	10	1
1	A	30	THR	N-CA	5.62	1.53	1.46	3	2
1	A	92	GLY	N-CA	5.62	1.53	1.45	7	1
1	A	67	LEU	C-N	5.61	1.40	1.33	2	2

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	36	GLY	CA-C	5.61	1.58	1.52	8	1
1	A	97	ALA	CA-C	5.59	1.59	1.52	1	1
1	A	48	GLN	C-O	-5.59	1.17	1.23	7	1
1	A	78	TYR	C-O	-5.58	1.16	1.23	9	1
1	A	86	GLU	CA-C	-5.58	1.45	1.52	10	2
1	A	15	GLU	N-CA	5.58	1.52	1.45	3	1
1	A	93	ILE	CA-C	5.57	1.59	1.52	1	2
1	A	84	LYS	CA-CB	5.56	1.62	1.53	6	1
1	A	24	VAL	C-O	5.55	1.30	1.24	6	2
1	A	111	LEU	C-N	5.55	1.40	1.33	6	1
1	A	104	GLU	CA-C	5.54	1.59	1.52	8	2
1	A	123	MET	N-CA	-5.53	1.39	1.46	10	2
1	A	21	LEU	N-CA	5.53	1.52	1.46	6	1
1	A	38	ASN	N-CA	5.52	1.53	1.46	2	1
1	A	66	HIS	CE1-NE2	5.51	1.38	1.32	6	2
1	A	42	LYS	N-CA	5.51	1.53	1.46	9	1
2	B	20	U	C4'-C3'	5.50	1.61	1.53	7	1
1	A	56	GLU	CA-C	5.48	1.60	1.52	6	1
1	A	76	ILE	N-CA	-5.47	1.41	1.46	1	1
1	A	85	LYS	N-CA	5.44	1.53	1.46	5	2
1	A	79	ILE	CA-C	5.43	1.59	1.52	4	1
2	B	3	U	C5'-C4'	5.43	1.59	1.51	9	1
1	A	41	THR	N-CA	5.43	1.52	1.46	4	1
1	A	32	LYS	CA-C	5.41	1.58	1.52	7	1
2	B	13	A	C4'-C3'	5.39	1.61	1.53	8	1
1	A	52	VAL	CA-C	-5.38	1.45	1.52	6	1
1	A	51	LEU	CA-C	5.35	1.59	1.52	1	1
2	B	13	A	O4'-C1'	5.33	1.49	1.41	3	1
2	B	22	A	O3'-P	-5.32	1.53	1.61	1	1
1	A	64	VAL	CA-CB	5.32	1.60	1.54	4	1
1	A	62	GLU	CA-CB	5.32	1.63	1.53	8	1
1	A	24	VAL	CA-C	5.31	1.59	1.52	3	1
1	A	54	ILE	CA-C	5.31	1.59	1.52	6	1
1	A	20	ALA	C-N	-5.30	1.27	1.33	7	1
2	B	7	C	C4-N4	-5.30	1.23	1.33	2	1
1	A	87	LEU	CA-C	5.29	1.59	1.52	9	1
1	A	61	GLU	N-CA	5.29	1.52	1.46	2	1
1	A	107	LYS	N-CA	5.28	1.53	1.46	2	1
2	B	24	G	P-O5'	-5.28	1.51	1.59	2	1
1	A	16	LEU	CA-C	5.26	1.59	1.52	5	1
2	B	11	A	O3'-P	-5.26	1.53	1.61	7	2

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
2	B	2	C	C1'-N1	5.25	1.56	1.48	8	1
1	A	28	ARG	CA-C	5.25	1.58	1.53	9	1
1	A	18	GLU	CA-C	5.25	1.59	1.52	1	1
2	B	22	A	C1'-N9	5.22	1.55	1.48	1	1
2	B	3	U	C4'-C3'	5.22	1.60	1.52	6	1
1	A	102	ILE	CA-CB	5.22	1.60	1.54	10	1
1	A	8	VAL	CA-CB	5.21	1.61	1.54	8	1
1	A	97	ALA	C-O	-5.21	1.18	1.24	1	1
1	A	73	GLU	C-O	5.20	1.30	1.24	6	1
1	A	10	PHE	C-N	5.20	1.40	1.33	9	1
1	A	47	GLY	N-CA	5.19	1.52	1.45	7	1
1	A	93	ILE	N-CA	5.19	1.52	1.46	4	1
1	A	16	LEU	N-CA	5.18	1.52	1.46	9	1
2	B	13	A	C5'-C4'	-5.17	1.43	1.51	6	1
2	B	8	U	C4'-O4'	-5.17	1.37	1.45	8	1
1	A	5	PRO	CA-C	-5.16	1.47	1.53	1	1
1	A	120	LYS	N-CA	5.16	1.52	1.46	6	1
1	A	45	GLU	N-CA	5.15	1.53	1.46	7	1
1	A	46	ARG	CA-C	5.14	1.59	1.52	10	1
2	B	24	G	C4'-O4'	-5.14	1.38	1.45	3	1
1	A	35	LYS	CA-C	5.14	1.58	1.52	10	1
1	A	92	GLY	CA-C	5.13	1.59	1.51	8	1
1	A	62	GLU	C-N	5.13	1.39	1.34	2	2
1	A	96	ALA	C-O	-5.12	1.17	1.23	4	1
2	B	19	A	P-O5'	5.12	1.67	1.59	2	1
2	B	18	A	C5'-C4'	5.12	1.58	1.51	4	1
2	B	6	G	C4'-O4'	-5.12	1.37	1.45	7	1
1	A	10	PHE	C-O	5.11	1.30	1.23	5	1
1	A	7	TYR	CA-CB	5.10	1.62	1.53	6	1
2	B	14	G	O4'-C1'	5.10	1.49	1.41	6	1
2	B	19	A	C5'-C4'	5.10	1.59	1.51	4	1
2	B	4	G	P-O5'	-5.09	1.52	1.59	8	1
2	B	20	U	O4'-C1'	5.09	1.49	1.42	10	1
1	A	78	TYR	C-N	5.08	1.39	1.33	3	1
1	A	19	LYS	N-CA	5.08	1.52	1.46	8	1
2	B	20	U	C4'-O4'	-5.08	1.38	1.45	3	1
1	A	27	ALA	CA-CB	5.07	1.62	1.53	3	1
1	A	18	GLU	C-N	5.07	1.41	1.33	9	1
1	A	120	LYS	C-N	-5.06	1.27	1.33	4	1
1	A	21	LEU	C-O	-5.04	1.18	1.24	5	1
1	A	12	VAL	N-CA	5.04	1.52	1.46	8	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)	Models	
								Worst	Total
1	A	9	LYS	CA-CB	5.02	1.62	1.53	2	1
1	A	48	GLN	CA-C	-5.01	1.46	1.52	7	1
2	B	4	G	C4'-C3'	5.01	1.60	1.53	10	1
2	B	24	G	C2'-O2'	-5.00	1.34	1.42	7	1
2	B	22	A	C4'-O4'	-5.00	1.38	1.45	8	1
2	B	9	C	C4'-O4'	-5.00	1.38	1.45	4	1

All unique angle 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(°)	Ideal(°)	Models	
								Worst	Total
2	B	10	G	C3'-C2'-C1'	12.04	113.34	101.30	10	5
2	B	24	G	O4'-C1'-C2'	-11.35	96.25	107.60	4	1
2	B	4	G	O4'-C4'-C3'	11.18	117.28	106.10	9	2
2	B	22	A	O4'-C4'-C3'	10.27	114.27	104.00	8	2
1	A	38	ASN	OD1-CG-ND2	-10.26	112.34	122.60	3	3
2	B	10	G	O4'-C4'-C3'	10.21	114.21	104.00	9	2
1	A	48	GLN	OE1-CD-NE2	-9.94	112.66	122.60	7	3
2	B	4	G	C3'-C2'-C1'	9.75	111.25	101.50	6	3
2	B	15	A	O4'-C4'-C3'	9.61	113.61	104.00	1	1
1	A	76	ILE	CB-CA-C	9.54	119.38	110.13	5	1
2	B	22	A	C4'-C3'-C2'	-9.48	93.12	102.60	6	3
1	A	100	VAL	O-C-N	-9.46	114.74	122.97	9	2
2	B	8	U	C1'-O4'-C4'	-9.45	100.45	109.90	8	1
2	B	18	A	O4'-C1'-C2'	-9.42	98.18	107.60	10	2
2	B	14	G	C4'-C3'-O3'	-9.34	99.00	113.00	10	1
1	A	54	ILE	O-C-N	-9.25	113.20	123.10	5	1
1	A	96	ALA	O-C-N	9.05	133.17	122.85	10	2
2	B	24	G	C3'-C2'-C1'	9.05	110.36	101.30	4	2
2	B	11	A	O4'-C1'-C2'	-9.01	96.79	105.80	5	2
1	A	81	VAL	O-C-N	-8.97	115.44	121.72	1	2
2	B	24	G	C1'-O4'-C4'	8.78	118.68	109.90	4	1
1	A	99	SER	CA-C-N	8.70	135.46	122.58	1	3
1	A	99	SER	C-N-CA	8.70	135.46	122.58	1	3
2	B	16	G	O4'-C1'-N9	8.68	121.51	108.50	8	1
2	B	23	U	C4'-C3'-C2'	-8.54	94.06	102.60	10	2
2	B	10	G	O4'-C1'-C2'	-8.45	99.15	107.60	10	2
2	B	21	G	O4'-C1'-C2'	8.45	114.25	105.80	4	4
1	A	57	ASP	CA-CB-CG	8.42	121.02	112.60	7	2
2	B	3	U	O4'-C4'-C3'	8.37	112.37	104.00	4	6
1	A	123	MET	N-CA-C	8.35	120.38	111.28	10	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	36	GLY	O-C-N	-8.31	114.74	122.47	2	1
1	A	83	SER	N-CA-C	8.31	121.70	109.07	6	5
1	A	24	VAL	CA-CB-CG1	8.27	124.46	110.40	9	4
2	B	23	U	O3'-P-O5'	8.19	116.29	104.00	7	1
2	B	4	G	O4'-C1'-C2'	-8.17	97.63	105.80	7	1
1	A	118	LYS	CA-C-O	8.14	129.05	120.42	2	2
1	A	12	VAL	N-CA-CB	8.13	122.59	111.21	6	1
2	B	22	A	C3'-C2'-O2'	8.13	122.89	110.70	10	2
2	B	3	U	C3'-C2'-C1'	8.12	109.42	101.30	3	1
2	B	11	A	C1'-C2'-O2'	8.03	123.85	111.80	3	1
2	B	12	A	C1'-C2'-O2'	-8.03	99.75	111.80	4	1
2	B	6	G	C1'-O4'-C4'	8.02	117.92	109.90	7	1
2	B	6	G	C3'-C2'-C1'	7.99	109.29	101.30	9	1
1	A	118	LYS	CA-C-N	7.97	130.89	120.60	6	2
1	A	118	LYS	C-N-CA	7.97	130.89	120.60	6	2
1	A	58	VAL	CA-C-O	-7.93	112.75	121.23	9	2
1	A	28	ARG	NE-CZ-NH2	7.88	126.29	119.20	3	3
2	B	23	U	O4'-C1'-C2'	-7.87	99.73	107.60	3	3
2	B	18	A	C1'-O4'-C4'	-7.87	102.03	109.90	3	1
1	A	109	ARG	NE-CZ-NH1	7.83	129.33	121.50	8	1
1	A	17	ALA	O-C-N	7.82	130.12	122.07	6	1
2	B	16	G	C4'-C3'-C2'	-7.79	94.81	102.60	4	1
1	A	119	VAL	CA-CB-CG1	7.75	123.58	110.40	9	2
1	A	110	ASP	CA-CB-CG	7.72	120.32	112.60	6	1
1	A	53	ILE	CA-C-O	-7.66	112.24	120.36	9	1
2	B	15	A	C4'-C3'-C2'	-7.61	94.99	102.60	6	3
1	A	76	ILE	CA-C-N	7.60	127.88	119.90	4	3
1	A	76	ILE	C-N-CA	7.60	127.88	119.90	4	3
1	A	97	ALA	CA-C-O	-7.60	112.17	120.30	9	2
1	A	103	ILE	CB-CA-C	7.59	117.76	110.63	3	2
1	A	77	PRO	N-CA-CB	7.58	110.03	103.36	2	3
1	A	28	ARG	CA-C-O	-7.57	112.44	120.92	2	1
1	A	29	ASP	CA-CB-CG	7.56	120.16	112.60	9	5
2	B	7	C	C5'-C4'-C3'	-7.54	104.68	116.00	7	2
2	B	18	A	C3'-C2'-C1'	-7.52	93.78	101.30	9	2
1	A	109	ARG	N-CA-C	7.52	120.43	111.33	6	2
2	B	19	A	C2'-C3'-O3'	-7.51	98.24	109.50	1	1
2	B	7	C	C3'-C2'-C1'	7.49	108.79	101.30	1	1
2	B	4	G	O3'-P-O5'	-7.48	92.78	104.00	9	1
1	A	90	ALA	CA-C-O	-7.45	112.88	121.07	3	2
1	A	104	GLU	CA-C-N	7.44	127.15	119.56	7	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	104	GLU	C-N-CA	7.44	127.15	119.56	7	1
1	A	106	GLY	CA-C-N	7.44	130.57	120.38	1	1
1	A	106	GLY	C-N-CA	7.44	130.57	120.38	1	1
2	B	6	G	O4'-C1'-C2'	-7.44	100.16	107.60	7	1
1	A	103	ILE	N-CA-C	-7.41	103.72	111.58	6	1
2	B	15	A	C3'-C2'-O2'	7.38	121.77	110.70	7	1
2	B	24	G	C4'-C3'-O3'	7.38	120.47	109.40	3	1
2	B	18	A	C4'-C3'-C2'	-7.37	95.23	102.60	8	2
2	B	11	A	C3'-C2'-C1'	7.37	108.67	101.30	1	2
2	B	13	A	C3'-C2'-C1'	7.34	108.84	101.50	7	3
2	B	18	A	O4'-C4'-C3'	7.32	111.32	104.00	3	1
2	B	8	U	C4'-C3'-C2'	-7.32	95.28	102.60	6	3
1	A	68	PRO	N-CA-CB	7.29	110.16	103.08	1	4
2	B	3	U	C2'-C3'-O3'	-7.26	102.81	113.70	3	1
1	A	52	VAL	CA-C-N	7.25	132.18	123.19	5	1
1	A	52	VAL	C-N-CA	7.25	132.18	123.19	5	1
1	A	57	ASP	CA-C-N	7.24	130.49	120.49	6	2
1	A	57	ASP	C-N-CA	7.24	130.49	120.49	6	2
2	B	15	A	O5'-C5'-C4'	7.23	122.35	111.50	4	1
1	A	87	LEU	CA-C-N	7.20	128.12	119.99	5	3
1	A	87	LEU	C-N-CA	7.20	128.12	119.99	5	3
2	B	3	U	P-O3'-C3'	7.19	130.98	120.20	4	1
2	B	10	G	C3'-C2'-O2'	-7.17	103.84	114.60	6	3
2	B	23	U	O4'-C1'-N1	7.17	119.25	108.50	1	3
2	B	2	C	O5'-C5'-C4'	7.16	122.24	111.50	2	1
1	A	7	TYR	N-CA-C	7.16	121.86	113.20	6	2
1	A	76	ILE	N-CA-C	7.15	116.48	108.05	2	2
2	B	10	G	C4'-C3'-C2'	-7.14	95.46	102.60	4	4
1	A	56	GLU	O-C-N	-7.14	112.27	122.41	7	1
1	A	16	LEU	CA-C-N	7.13	130.42	120.29	10	1
1	A	16	LEU	C-N-CA	7.13	130.42	120.29	10	1
2	B	8	U	C3'-C2'-C1'	7.12	108.42	101.30	10	1
1	A	41	THR	OG1-CB-CG2	-7.10	95.09	109.30	4	1
1	A	36	GLY	CA-C-N	7.08	129.76	120.28	2	2
1	A	36	GLY	C-N-CA	7.08	129.76	120.28	2	2
2	B	5	A	O4'-C1'-N9	7.06	118.80	108.20	5	4
1	A	52	VAL	O-C-N	-7.06	115.62	123.03	8	2
2	B	9	C	O4'-C4'-C3'	7.05	111.06	104.00	5	3
1	A	84	LYS	O-C-N	-7.04	113.00	122.43	10	1
2	B	7	C	O3'-P-O5'	7.03	114.55	104.00	10	1
2	B	7	C	O4'-C1'-N1	7.03	119.04	108.50	8	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	22	GLN	N-CA-C	7.02	119.96	111.82	9	1
1	A	68	PRO	N-CA-C	7.00	119.24	110.70	5	2
2	B	24	G	O4'-C4'-C3'	6.99	110.99	104.00	5	2
1	A	64	VAL	CA-C-O	-6.98	114.02	119.46	7	1
2	B	16	G	C3'-C2'-O2'	6.97	121.16	110.70	3	2
1	A	55	ALA	O-C-N	-6.97	115.09	122.96	3	1
1	A	47	GLY	CA-C-O	-6.96	113.29	121.08	8	1
2	B	16	G	O4'-C4'-C3'	6.95	110.95	104.00	4	1
1	A	115	ILE	CA-CB-CG1	6.94	122.20	110.40	4	1
1	A	38	ASN	O-C-N	-6.93	114.78	122.12	3	2
1	A	46	ARG	N-CA-C	6.91	120.81	112.38	5	1
1	A	68	PRO	CA-C-O	-6.89	110.57	120.56	3	2
1	A	74	LYS	CA-C-O	6.88	127.61	119.28	1	1
1	A	6	SER	N-CA-C	6.86	120.88	112.23	8	1
1	A	22	GLN	OE1-CD-NE2	-6.86	115.74	122.60	5	4
1	A	40	THR	CA-C-N	6.84	129.34	120.44	10	1
1	A	40	THR	C-N-CA	6.84	129.34	120.44	10	1
2	B	21	G	C3'-C2'-C1'	-6.81	94.69	101.50	4	4
2	B	9	C	O4'-C1'-C2'	-6.77	100.83	107.60	4	1
1	A	19	LYS	CA-C-O	-6.76	113.26	120.42	7	1
1	A	96	ALA	N-CA-C	6.74	119.74	110.24	1	1
2	B	11	A	O5'-C5'-C4'	6.73	121.79	111.70	8	1
1	A	6	SER	O-C-N	-6.70	113.45	122.43	2	1
1	A	50	LYS	N-CA-C	6.70	119.59	111.82	6	1
2	B	8	U	O5'-C5'-C4'	-6.70	101.45	111.50	8	1
1	A	21	LEU	CB-CA-C	6.69	121.39	110.88	4	1
1	A	113	GLU	O-C-N	-6.69	115.18	122.07	7	1
2	B	3	U	C4'-C3'-C2'	-6.69	95.91	102.60	7	1
2	B	20	U	C1'-O4'-C4'	-6.68	103.02	109.70	4	1
1	A	32	LYS	CA-C-N	6.67	133.30	122.69	7	1
1	A	32	LYS	C-N-CA	6.67	133.30	122.69	7	1
2	B	5	A	C4'-C3'-O3'	6.67	119.40	109.40	8	3
2	B	2	C	O4'-C4'-C3'	6.66	110.66	104.00	4	2
2	B	19	A	C5'-C4'-O4'	6.66	119.08	109.10	7	2
2	B	14	G	C3'-C2'-O2'	6.64	120.66	110.70	10	1
1	A	65	ALA	CA-C-O	-6.63	112.77	120.20	3	1
1	A	46	ARG	CD-NE-CZ	6.62	133.67	124.40	6	1
1	A	112	VAL	CA-C-N	6.62	129.04	120.44	5	1
1	A	112	VAL	C-N-CA	6.62	129.04	120.44	5	1
1	A	76	ILE	N-CA-CB	-6.61	105.36	112.37	5	1
1	A	15	GLU	CA-C-O	-6.60	114.46	121.99	8	2

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	100	VAL	CA-C-N	6.59	132.28	122.99	2	1
1	A	100	VAL	C-N-CA	6.59	132.28	122.99	2	1
1	A	74	LYS	CA-C-N	6.59	131.48	122.19	1	2
1	A	74	LYS	C-N-CA	6.59	131.48	122.19	1	2
1	A	107	LYS	N-CA-C	6.58	120.41	112.38	9	2
2	B	14	G	C4'-C3'-C2'	-6.58	96.02	102.60	5	1
2	B	23	U	O4'-C4'-C3'	6.57	110.57	104.00	7	1
1	A	40	THR	N-CA-C	6.57	118.99	111.11	8	1
2	B	12	A	O4'-C4'-C3'	-6.55	97.45	104.00	10	4
2	B	12	A	C3'-C2'-C1'	6.55	107.85	101.30	9	4
2	B	20	U	P-O3'-C3'	6.55	130.02	120.20	2	1
1	A	38	ASN	CA-CB-CG	6.55	119.15	112.60	4	1
1	A	91	ALA	CA-C-N	6.54	132.11	122.10	1	1
1	A	91	ALA	C-N-CA	6.54	132.11	122.10	1	1
1	A	59	ASP	CA-C-O	-6.54	113.58	119.86	4	2
1	A	119	VAL	CA-C-N	6.52	129.90	120.38	3	2
1	A	119	VAL	C-N-CA	6.52	129.90	120.38	3	2
1	A	99	SER	CA-C-O	-6.52	114.47	121.45	1	2
2	B	12	A	C2'-C3'-O3'	6.51	119.26	109.50	4	1
2	B	19	A	C4'-C3'-C2'	-6.50	96.10	102.60	10	1
1	A	93	ILE	CB-CA-C	6.49	120.47	111.38	7	2
2	B	5	A	O4'-C1'-C2'	-6.49	99.31	105.80	5	1
2	B	4	G	C5'-C4'-C3'	-6.47	105.50	115.20	5	2
1	A	66	HIS	CB-CG-CD2	-6.46	122.81	131.20	4	4
1	A	12	VAL	CB-CA-C	6.45	116.68	110.16	4	2
2	B	2	C	C1'-C2'-O2'	-6.45	98.72	108.40	1	1
2	B	18	A	C1'-C2'-O2'	-6.45	98.72	108.40	6	1
2	B	23	U	C3'-C2'-C1'	6.43	107.73	101.30	5	4
1	A	5	PRO	CA-C-N	6.42	129.76	120.38	5	3
1	A	5	PRO	C-N-CA	6.42	129.76	120.38	5	3
2	B	5	A	O3'-P-O5'	6.41	113.62	104.00	3	2
2	B	4	G	C4'-C3'-C2'	-6.41	96.19	102.60	9	2
1	A	60	PRO	N-CD-CG	6.39	111.47	103.80	1	2
2	B	9	C	C4'-C3'-O3'	-6.39	103.42	113.00	1	1
1	A	110	ASP	CA-C-N	6.37	128.73	120.44	4	2
1	A	110	ASP	C-N-CA	6.37	128.73	120.44	4	2
1	A	56	GLU	CA-C-O	6.36	126.74	119.18	7	1
1	A	114	GLU	N-CA-C	6.35	117.87	111.07	3	1
2	B	19	A	C1'-C2'-O2'	-6.35	102.27	111.80	4	4
2	B	2	C	C4'-C3'-C2'	-6.35	96.25	102.60	3	1
1	A	58	VAL	CA-C-N	6.34	129.23	122.26	1	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	58	VAL	C-N-CA	6.34	129.23	122.26	1	1
2	B	22	A	O3'-P-O5'	6.34	113.51	104.00	10	2
2	B	19	A	O4'-C4'-C3'	6.34	112.44	106.10	2	1
1	A	52	VAL	CA-CB-CG2	6.31	121.13	110.40	6	1
2	B	17	C	C4'-C3'-O3'	6.31	122.47	113.00	8	1
1	A	109	ARG	NH1-CZ-NH2	-6.30	111.11	119.30	8	1
1	A	80	TYR	CA-C-O	-6.30	113.80	121.36	4	1
1	A	69	PRO	CA-C-N	6.29	128.71	120.28	5	2
1	A	69	PRO	C-N-CA	6.29	128.71	120.28	5	2
1	A	17	ALA	CA-C-N	6.28	128.94	120.65	5	2
1	A	17	ALA	C-N-CA	6.28	128.94	120.65	5	2
2	B	16	G	O3'-P-O5'	6.26	113.39	104.00	9	1
1	A	43	ALA	CA-C-N	6.25	128.35	120.72	1	1
1	A	43	ALA	C-N-CA	6.25	128.35	120.72	1	1
2	B	17	C	C5'-C4'-C3'	-6.25	106.63	116.00	7	2
1	A	102	ILE	CA-C-N	6.24	130.06	122.63	8	1
1	A	102	ILE	C-N-CA	6.24	130.06	122.63	8	1
1	A	27	ALA	CA-C-N	6.22	128.94	120.54	1	2
1	A	27	ALA	C-N-CA	6.22	128.94	120.54	1	2
1	A	93	ILE	O-C-N	-6.22	116.34	122.93	5	2
1	A	109	ARG	NE-CZ-NH2	6.22	124.80	119.20	2	1
1	A	46	ARG	NE-CZ-NH2	6.21	124.79	119.20	7	2
2	B	13	A	O4'-C4'-C3'	6.20	112.30	106.10	7	1
2	B	21	G	C4'-C3'-C2'	-6.20	96.40	102.60	6	1
1	A	38	ASN	CA-C-N	6.18	129.07	120.29	3	2
1	A	38	ASN	C-N-CA	6.18	129.07	120.29	3	2
1	A	13	PRO	CA-C-N	6.17	128.55	120.28	4	3
1	A	13	PRO	C-N-CA	6.17	128.55	120.28	4	3
1	A	120	LYS	N-CA-CB	-6.16	100.76	109.94	5	1
2	B	18	A	C2'-C3'-O3'	6.16	122.94	113.70	5	1
1	A	105	PRO	CB-CA-C	-6.13	103.14	111.85	7	1
2	B	7	C	C5'-C4'-O4'	6.13	119.00	109.80	5	4
2	B	20	U	C3'-C2'-C1'	-6.13	95.37	101.50	10	1
1	A	75	GLU	CA-C-N	6.12	130.17	123.43	7	1
1	A	75	GLU	C-N-CA	6.12	130.17	123.43	7	1
1	A	46	ARG	NE-CZ-NH1	6.12	127.62	121.50	1	2
1	A	89	ALA	CA-C-N	6.12	130.00	120.82	1	1
1	A	89	ALA	C-N-CA	6.12	130.00	120.82	1	1
1	A	121	GLU	CA-C-N	6.12	128.98	120.29	9	1
1	A	121	GLU	C-N-CA	6.12	128.98	120.29	9	1
2	B	14	G	O4'-C4'-C3'	6.11	110.11	104.00	5	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	41	THR	N-CA-C	6.11	118.47	111.02	8	1
2	B	13	A	C1'-C2'-O2'	-6.10	102.66	111.80	9	2
2	B	13	A	O4'-C1'-C2'	-6.07	99.73	105.80	8	1
1	A	17	ALA	N-CA-C	6.07	117.57	111.07	6	1
2	B	4	G	C1'-O4'-C4'	-6.04	103.66	109.70	9	2
1	A	81	VAL	CA-C-N	6.04	125.25	118.97	6	5
1	A	81	VAL	C-N-CA	6.04	125.25	118.97	6	5
1	A	45	GLU	CA-C-N	6.04	131.14	122.08	3	1
1	A	45	GLU	C-N-CA	6.04	131.14	122.08	3	1
2	B	20	U	C5'-C4'-C3'	-6.03	106.15	115.20	5	2
1	A	109	ARG	CD-NE-CZ	6.02	132.83	124.40	2	2
2	B	14	G	C3'-C2'-C1'	6.02	107.32	101.30	8	3
1	A	45	GLU	CB-CG-CD	6.02	122.83	112.60	10	1
2	B	3	U	O4'-C1'-N1	6.01	117.52	108.50	8	1
2	B	2	C	C2'-C3'-O3'	-6.01	104.68	113.70	7	1
1	A	33	ILE	CA-CB-CG1	6.00	120.61	110.40	6	1
1	A	121	GLU	CA-C-O	-6.00	112.68	120.00	8	1
2	B	18	A	O4'-C1'-N9	5.99	117.49	108.50	7	1
2	B	2	C	O4'-C1'-N1	5.99	117.49	108.50	10	1
1	A	72	GLU	CA-C-N	5.98	128.57	120.38	8	2
1	A	72	GLU	C-N-CA	5.98	128.57	120.38	8	2
1	A	112	VAL	CA-CB-CG1	5.97	120.56	110.40	3	3
2	B	19	A	C5'-C4'-C3'	-5.97	106.24	115.20	9	2
2	B	6	G	C3'-C2'-O2'	5.97	119.65	110.70	2	2
2	B	10	G	C5'-C4'-O4'	5.97	118.76	109.80	3	1
2	B	10	G	C1'-O4'-C4'	-5.97	103.93	109.90	4	2
2	B	21	G	C1'-O4'-C4'	-5.96	103.74	109.70	4	2
1	A	86	GLU	CA-C-N	5.96	128.26	120.28	2	2
1	A	86	GLU	C-N-CA	5.96	128.26	120.28	2	2
1	A	34	ARG	CD-NE-CZ	5.95	132.74	124.40	2	2
1	A	116	ALA	O-C-N	-5.95	115.36	122.15	3	1
2	B	24	G	O5'-C5'-C4'	5.93	120.60	111.70	3	1
2	B	8	U	C5'-C4'-C3'	-5.93	107.10	116.00	4	2
1	A	83	SER	CA-C-N	5.92	128.81	120.28	5	2
1	A	83	SER	C-N-CA	5.92	128.81	120.28	5	2
1	A	87	LEU	CD1-CG-CD2	-5.92	97.77	110.80	1	1
2	B	20	U	C5'-C4'-O4'	5.91	118.67	109.80	3	2
2	B	7	C	N1-C1'-C2'	-5.91	103.13	112.00	3	2
2	B	19	A	O5'-C5'-C4'	-5.91	102.83	111.70	1	2
1	A	53	ILE	O-C-N	-5.90	116.68	123.00	7	1
1	A	66	HIS	CA-CB-CG	-5.89	107.91	113.80	2	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
2	B	22	A	O5'-C5'-C4'	5.88	120.33	111.50	3	2
1	A	61	GLU	CA-C-O	5.88	126.55	119.61	6	1
1	A	77	PRO	CB-CA-C	-5.88	100.98	110.21	1	1
1	A	84	LYS	CG-CD-CE	-5.88	97.78	111.30	2	1
1	A	26	ILE	CA-C-O	-5.87	112.57	119.85	9	1
1	A	115	ILE	CA-C-N	5.87	128.63	120.29	3	2
1	A	115	ILE	C-N-CA	5.87	128.63	120.29	3	2
2	B	7	C	O4'-C4'-C3'	5.87	109.87	104.00	8	2
2	B	12	A	C4'-C3'-C2'	5.87	108.47	102.60	10	3
2	B	5	A	C5'-C4'-C3'	-5.87	106.40	115.20	2	1
2	B	18	A	C5'-C4'-O4'	5.86	118.60	109.80	4	1
2	B	20	U	O4'-C1'-C2'	-5.86	99.94	105.80	6	3
1	A	69	PRO	O-C-N	-5.86	115.50	122.24	7	1
2	B	3	U	P-O5'-C5'	5.86	129.69	120.90	8	1
1	A	28	ARG	NE-CZ-NH1	5.86	127.36	121.50	4	3
2	B	13	A	O3'-P-O5'	5.85	112.78	104.00	5	1
1	A	39	GLU	CA-C-N	5.84	128.04	120.44	9	1
1	A	39	GLU	C-N-CA	5.84	128.04	120.44	9	1
2	B	10	G	O5'-C5'-C4'	-5.84	102.73	111.50	10	1
1	A	73	GLU	N-CA-C	5.84	118.39	111.33	8	3
1	A	30	THR	OG1-CB-CG2	-5.84	97.62	109.30	10	1
2	B	14	G	O3'-P-O5'	5.83	112.75	104.00	1	1
2	B	8	U	O4'-C4'-C3'	5.82	109.82	104.00	9	2
2	B	22	A	C4'-C3'-O3'	5.82	121.73	113.00	10	1
2	B	11	A	O3'-P-O5'	-5.82	95.28	104.00	4	1
2	B	17	C	O5'-C5'-C4'	5.80	120.21	111.50	5	2
1	A	71	CYS	O-C-N	-5.80	116.10	122.07	5	1
2	B	10	G	C5'-C4'-C3'	-5.79	107.31	116.00	10	1
1	A	95	VAL	CA-CB-CG2	5.79	120.24	110.40	6	2
2	B	6	G	C5'-C4'-O4'	5.79	118.48	109.80	3	1
1	A	113	GLU	CB-CG-CD	5.79	122.44	112.60	10	1
2	B	5	A	C3'-C2'-O2'	-5.78	105.93	114.60	10	1
2	B	7	C	C4'-C3'-C2'	-5.77	96.83	102.60	8	1
2	B	16	G	C1'-O4'-C4'	-5.77	104.13	109.90	3	1
1	A	69	PRO	N-CA-CB	5.77	109.83	103.26	7	2
1	A	54	ILE	N-CA-C	5.76	117.05	108.46	5	1
1	A	38	ASN	CB-CA-C	5.75	120.34	110.79	8	1
1	A	30	THR	CA-CB-OG1	5.75	118.23	109.60	3	2
1	A	111	LEU	N-CA-C	5.75	117.63	111.36	3	1
1	A	101	ALA	N-CA-CB	-5.75	101.25	111.08	4	1
1	A	89	ALA	O-C-N	-5.75	116.15	122.07	1	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	116	ALA	CA-C-O	5.74	126.50	120.42	3	1
2	B	3	U	O3'-P-O5'	5.74	112.61	104.00	3	1
1	A	81	VAL	CA-CB-CG2	5.74	120.15	110.40	1	2
1	A	120	LYS	CA-C-O	-5.72	113.79	120.20	3	2
1	A	18	GLU	N-CA-C	5.72	117.51	111.28	4	1
1	A	79	ILE	CA-C-N	5.72	131.17	122.65	10	1
1	A	79	ILE	C-N-CA	5.72	131.17	122.65	10	1
1	A	117	MET	N-CA-C	5.72	117.19	111.07	6	1
1	A	45	GLU	N-CA-C	5.71	118.45	111.82	1	1
1	A	20	ALA	N-CA-C	5.71	117.50	111.28	7	1
2	B	17	C	O4'-C1'-N1	5.70	117.05	108.50	7	1
2	B	2	C	O4'-C1'-C2'	-5.70	101.91	107.60	3	2
2	B	12	A	C5'-C4'-O4'	5.69	117.64	109.10	2	1
2	B	21	G	O4'-C1'-N9	-5.68	99.68	108.20	6	1
1	A	39	GLU	CB-CA-C	5.67	119.87	110.90	8	1
2	B	23	U	C2'-C3'-O3'	-5.67	105.19	113.70	3	1
1	A	24	VAL	CA-C-N	5.67	128.14	120.38	2	1
1	A	24	VAL	C-N-CA	5.67	128.14	120.38	2	1
1	A	15	GLU	O-C-N	-5.66	116.05	122.68	2	1
1	A	116	ALA	N-CA-C	5.66	117.13	111.07	8	1
2	B	17	C	C3'-C2'-C1'	-5.66	95.64	101.30	9	1
2	B	24	G	C1'-C2'-O2'	5.66	116.89	108.40	6	1
2	B	24	G	C5'-C4'-C3'	-5.65	107.52	116.00	8	1
2	B	9	C	C4'-C3'-C2'	-5.64	96.96	102.60	4	1
1	A	24	VAL	CA-C-O	-5.63	114.69	120.71	7	2
1	A	119	VAL	N-CA-C	-5.62	105.36	110.82	5	1
1	A	39	GLU	N-CA-C	-5.62	105.30	111.82	6	1
1	A	31	GLY	CA-C-N	5.62	131.40	122.74	8	1
1	A	31	GLY	C-N-CA	5.62	131.40	122.74	8	1
1	A	62	GLU	N-CA-C	5.62	119.61	112.87	8	1
2	B	17	C	O3'-P-O5'	-5.61	95.59	104.00	1	1
2	B	15	A	C5'-C4'-O4'	5.61	118.21	109.80	9	1
1	A	118	LYS	O-C-N	-5.60	115.77	122.15	2	1
1	A	28	ARG	N-CA-CB	-5.60	101.67	110.06	7	2
2	B	8	U	O4'-C1'-C2'	-5.59	102.00	107.60	6	1
2	B	23	U	C5'-C4'-C3'	-5.59	107.61	116.00	10	1
2	B	11	A	N9-C1'-C2'	5.58	120.38	112.00	7	1
1	A	41	THR	O-C-N	-5.58	116.21	122.12	5	1
2	B	11	A	O4'-C4'-C3'	5.58	109.58	104.00	4	2
2	B	17	C	N1-C1'-C2'	-5.57	103.64	112.00	4	1
1	A	53	ILE	N-CA-C	5.57	115.97	108.17	5	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	51	LEU	CA-C-N	5.57	130.82	122.58	4	1
1	A	51	LEU	C-N-CA	5.57	130.82	122.58	4	1
2	B	11	A	C2'-C3'-O3'	5.57	117.85	109.50	8	1
2	B	12	A	O4'-C1'-C2'	5.57	113.17	107.60	8	1
2	B	21	G	C2'-C3'-O3'	5.56	117.84	109.50	4	1
1	A	44	VAL	O-C-N	-5.56	114.21	121.82	4	1
2	B	22	A	C3'-C2'-C1'	5.56	106.86	101.30	7	1
1	A	19	LYS	CA-C-N	5.54	127.64	120.44	5	1
1	A	19	LYS	C-N-CA	5.54	127.64	120.44	5	1
1	A	34	ARG	O-C-N	-5.53	116.80	123.27	3	1
1	A	14	LYS	CA-C-O	-5.53	114.56	120.42	5	1
2	B	9	C	C3'-C2'-C1'	5.53	106.83	101.30	4	2
1	A	41	THR	N-CA-CB	5.52	118.01	110.01	10	1
1	A	87	LEU	N-CA-C	5.52	117.37	111.36	7	3
1	A	59	ASP	N-CA-CB	-5.52	100.55	110.37	5	1
2	B	17	C	O4'-C1'-C2'	-5.52	102.08	107.60	5	1
1	A	82	PRO	CA-C-N	5.50	131.59	121.85	9	1
1	A	82	PRO	C-N-CA	5.50	131.59	121.85	9	1
1	A	44	VAL	N-CA-C	5.50	116.14	110.36	6	1
2	B	14	G	O5'-C5'-C4'	-5.50	103.25	111.50	8	1
1	A	10	PHE	O-C-N	-5.48	117.00	123.25	2	1
1	A	59	ASP	CA-CB-CG	5.47	118.07	112.60	5	1
1	A	99	SER	N-CA-C	5.47	117.16	108.79	4	1
1	A	45	GLU	O-C-N	-5.47	115.32	122.59	9	1
1	A	64	VAL	CA-C-N	5.46	129.84	120.72	6	1
1	A	64	VAL	C-N-CA	5.46	129.84	120.72	6	1
1	A	64	VAL	O-C-N	-5.46	115.04	122.03	8	1
1	A	68	PRO	CA-C-N	5.46	125.54	119.32	7	1
1	A	68	PRO	C-N-CA	5.46	125.54	119.32	7	1
1	A	20	ALA	CA-C-N	5.45	128.37	120.79	10	1
1	A	20	ALA	C-N-CA	5.45	128.37	120.79	10	1
1	A	101	ALA	N-CA-C	5.45	118.33	109.24	7	2
2	B	23	U	C3'-C2'-O2'	-5.44	102.54	110.70	8	1
2	B	19	A	C1'-O4'-C4'	5.44	115.14	109.70	4	1
2	B	21	G	O4'-C4'-C3'	5.44	111.54	106.10	6	1
1	A	40	THR	CA-CB-CG2	5.43	119.73	110.50	4	1
2	B	3	U	C3'-C2'-O2'	-5.43	102.56	110.70	6	1
2	B	9	C	O4'-C1'-N1	5.42	116.64	108.50	8	3
2	B	2	C	C5'-C4'-C3'	-5.42	107.86	116.00	2	1
1	A	102	ILE	O-C-N	-5.42	116.33	122.72	3	2
1	A	9	LYS	CA-C-N	5.42	132.03	121.63	6	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	9	LYS	C-N-CA	5.42	132.03	121.63	6	1
1	A	43	ALA	N-CA-C	5.42	117.04	111.03	6	1
1	A	8	VAL	CA-C-N	5.42	129.26	120.60	4	1
1	A	8	VAL	C-N-CA	5.42	129.26	120.60	4	1
1	A	63	ILE	N-CA-C	5.41	120.03	113.00	3	1
1	A	94	GLU	CA-C-N	5.41	128.59	122.48	3	1
1	A	94	GLU	C-N-CA	5.41	128.59	122.48	3	1
1	A	89	ALA	N-CA-C	5.40	117.17	111.28	3	1
1	A	26	ILE	CA-C-N	5.39	128.50	120.31	6	1
1	A	26	ILE	C-N-CA	5.39	128.50	120.31	6	1
1	A	93	ILE	CA-C-O	-5.39	115.61	121.92	4	1
1	A	86	GLU	CA-C-O	-5.38	114.72	120.42	6	1
1	A	102	ILE	CB-CG1-CD1	5.37	125.08	113.80	9	1
1	A	34	ARG	NH1-CZ-NH2	-5.37	112.32	119.30	9	1
2	B	5	A	O4'-C4'-C3'	5.37	111.47	106.10	1	1
1	A	60	PRO	N-CA-CB	5.36	108.50	102.60	4	2
1	A	64	VAL	N-CA-C	-5.36	108.06	113.47	8	2
2	B	12	A	O4'-C1'-N9	5.36	116.23	108.20	5	1
2	B	14	G	O4'-C1'-C2'	-5.35	102.25	107.60	4	3
2	B	14	G	C2'-C3'-O3'	5.33	121.70	113.70	10	1
1	A	122	LEU	CA-C-N	5.33	131.98	121.58	8	1
1	A	122	LEU	C-N-CA	5.33	131.98	121.58	8	1
1	A	93	ILE	N-CA-CB	-5.33	106.00	112.07	3	1
1	A	88	GLY	CA-C-O	5.32	126.15	120.40	9	1
1	A	78	TYR	CA-C-N	5.32	129.88	122.23	9	1
1	A	78	TYR	C-N-CA	5.32	129.88	122.23	9	1
1	A	31	GLY	O-C-N	-5.31	116.37	123.64	8	1
2	B	23	U	C5'-C4'-O4'	5.30	117.74	109.80	4	1
1	A	116	ALA	CA-C-N	5.29	127.63	120.38	1	2
1	A	116	ALA	C-N-CA	5.29	127.63	120.38	1	2
1	A	52	VAL	CA-C-O	5.29	126.89	121.28	4	1
2	B	9	C	N1-C1'-C2'	-5.29	104.06	112.00	8	2
2	B	5	A	C5'-C4'-O4'	5.29	117.03	109.10	8	2
2	B	23	U	P-O5'-C5'	5.28	128.83	120.90	8	1
1	A	33	ILE	N-CA-C	5.28	117.66	108.90	3	1
1	A	12	VAL	CA-C-N	5.28	126.44	119.84	6	2
1	A	12	VAL	C-N-CA	5.28	126.44	119.84	6	2
1	A	18	GLU	CA-C-N	5.27	127.61	120.44	8	1
1	A	18	GLU	C-N-CA	5.27	127.61	120.44	8	1
1	A	61	GLU	CA-C-N	5.27	129.52	120.72	3	2
1	A	61	GLU	C-N-CA	5.27	129.52	120.72	3	2

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
2	B	3	U	C1'-O4'-C4'	-5.27	104.63	109.90	10	1
1	A	8	VAL	N-CA-CB	5.26	117.78	111.31	3	1
1	A	89	ALA	CA-C-O	5.26	126.13	120.55	3	1
1	A	79	ILE	N-CA-C	5.26	115.29	108.35	4	1
1	A	124	LYS	N-CA-CB	-5.26	101.56	110.50	10	1
1	A	28	ARG	N-CA-C	5.25	117.00	111.28	5	1
2	B	15	A	C5'-C4'-C3'	-5.25	108.13	116.00	4	1
1	A	48	GLN	N-CA-C	5.25	116.62	107.49	6	1
1	A	70	LEU	CA-C-N	5.24	128.27	120.31	1	1
1	A	70	LEU	C-N-CA	5.24	128.27	120.31	1	1
1	A	67	LEU	CA-C-O	-5.23	112.99	120.16	7	1
1	A	95	VAL	O-C-N	-5.23	115.62	122.47	7	1
2	B	2	C	C4'-C3'-O3'	5.21	120.82	113.00	1	1
1	A	111	LEU	O-C-N	-5.20	116.68	122.09	10	1
2	B	22	A	O4'-C1'-C2'	-5.20	102.40	107.60	6	1
1	A	26	ILE	CA-CB-CG1	5.20	119.24	110.40	3	1
1	A	113	GLU	CB-CA-C	5.19	119.03	110.88	3	1
1	A	98	ALA	N-CA-CB	-5.19	102.34	110.44	4	1
2	B	5	A	C1'-O4'-C4'	5.18	114.89	109.70	6	1
1	A	116	ALA	N-CA-CB	-5.18	102.50	110.01	8	1
2	B	6	G	C5'-C4'-C3'	-5.18	108.23	116.00	8	1
2	B	4	G	O4'-C1'-N9	5.18	115.97	108.20	10	1
1	A	8	VAL	CG1-CB-CG2	-5.17	99.43	110.80	5	1
1	A	34	ARG	NE-CZ-NH1	5.17	126.67	121.50	8	1
2	B	9	C	O3'-P-O5'	5.17	111.75	104.00	8	1
1	A	104	GLU	CB-CA-C	5.17	117.06	110.34	9	1
1	A	34	ARG	NE-CZ-NH2	5.16	123.85	119.20	10	1
2	B	23	U	C1'-C2'-O2'	5.16	116.14	108.40	8	1
2	B	20	U	O5'-C5'-C4'	5.16	119.24	111.50	3	1
2	B	13	A	C5'-C4'-O4'	5.16	117.54	109.80	5	1
2	B	16	G	P-O5'-C5'	-5.16	113.17	120.90	3	1
2	B	4	G	C2'-C3'-O3'	5.15	117.23	109.50	9	1
2	B	3	U	C5'-C4'-C3'	-5.15	108.27	116.00	2	1
2	B	6	G	O4'-C4'-C3'	5.15	109.15	104.00	5	1
1	A	92	GLY	CA-C-O	-5.15	113.30	119.06	10	1
1	A	102	ILE	CA-CB-CG1	5.15	119.15	110.40	7	1
1	A	35	LYS	N-CA-C	5.14	117.63	109.50	3	2
1	A	58	VAL	N-CA-C	5.13	117.02	109.63	1	1
1	A	10	PHE	CA-CB-CG	-5.13	108.67	113.80	10	2
1	A	95	VAL	N-CA-C	5.13	117.07	109.17	9	1
1	A	108	ALA	N-CA-C	5.12	119.10	113.21	2	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
1	A	112	VAL	O-C-N	-5.12	116.01	121.80	5	1
1	A	109	ARG	CA-C-O	-5.12	115.45	120.82	5	1
1	A	117	MET	CA-C-N	5.12	127.55	120.29	6	1
1	A	117	MET	C-N-CA	5.12	127.55	120.29	6	1
2	B	16	G	C5'-C4'-O4'	5.12	117.47	109.80	7	1
2	B	3	U	O5'-C5'-C4'	5.11	119.16	111.50	8	1
1	A	90	ALA	O-C-N	5.11	127.54	122.08	3	1
2	B	9	C	C1'-O4'-C4'	-5.10	104.80	109.90	5	1
1	A	55	ALA	N-CA-C	5.09	118.25	110.36	2	1
1	A	12	VAL	CA-C-O	-5.09	117.30	119.94	4	1
1	A	11	GLU	CA-CB-CG	5.08	124.25	114.10	6	1
1	A	81	VAL	N-CA-C	5.07	119.84	108.88	4	1
1	A	9	LYS	O-C-N	-5.07	116.00	122.34	6	1
2	B	2	C	O3'-P-O5'	5.07	111.61	104.00	2	1
1	A	55	ALA	CA-C-O	-5.07	116.03	121.56	8	2
1	A	42	LYS	N-CA-C	-5.07	105.67	111.14	2	1
1	A	5	PRO	O-C-N	-5.07	116.82	123.00	10	1
1	A	84	LYS	CB-CG-CD	-5.06	99.66	111.30	4	1
2	B	11	A	O4'-C1'-N9	5.06	115.80	108.20	6	1
1	A	71	CYS	CA-C-N	5.06	129.23	120.58	8	1
1	A	71	CYS	C-N-CA	5.06	129.23	120.58	8	1
2	B	20	U	O3'-P-O5'	5.05	111.58	104.00	8	1
1	A	94	GLU	CA-C-O	-5.05	114.00	120.31	8	1
2	B	11	A	C5'-C4'-O4'	5.04	116.67	109.10	6	1
1	A	71	CYS	N-CA-C	5.04	116.46	111.07	5	1
1	A	113	GLU	N-CA-C	-5.04	105.49	111.69	6	1
1	A	94	GLU	N-CA-CB	-5.04	102.58	110.44	7	1
1	A	33	ILE	CA-C-O	-5.03	116.38	121.67	10	1
1	A	93	ILE	CA-C-N	5.03	127.44	120.29	10	1
1	A	93	ILE	C-N-CA	5.03	127.44	120.29	10	1
1	A	60	PRO	CA-N-CD	-5.03	104.46	111.50	9	1
1	A	7	TYR	CA-C-N	5.03	127.32	120.63	7	1
1	A	7	TYR	C-N-CA	5.03	127.32	120.63	7	1
2	B	8	U	O4'-C1'-N1	5.03	116.04	108.50	1	1
2	B	2	C	C5'-C4'-O4'	5.03	117.34	109.80	6	1
1	A	78	TYR	CA-C-O	-5.03	115.23	121.11	10	1
2	B	7	C	C1'-O4'-C4'	-5.02	104.88	109.90	2	1
2	B	6	G	C4'-C3'-C2'	-5.02	97.58	102.60	6	1
1	A	30	THR	O-C-N	-5.02	115.27	122.04	9	1
1	A	113	GLU	N-CA-CB	-5.01	102.80	110.07	1	1
2	B	15	A	C1'-C2'-O2'	-5.01	100.88	108.40	2	1

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)	Models	
								Worst	Total
2	B	23	U	C1'-O4'-C4'	5.01	114.91	109.90	3	1
1	A	80	TYR	CA-C-N	5.01	128.75	121.68	5	1
1	A	80	TYR	C-N-CA	5.01	128.75	121.68	5	1
1	A	79	ILE	CA-C-O	5.01	126.35	121.19	2	1
2	B	10	G	C1'-C2'-O2'	-5.01	100.89	108.40	1	1
1	A	22	GLN	CG-CD-NE2	5.00	123.90	116.40	2	1
1	A	85	LYS	O-C-N	5.00	127.22	122.07	2	1

There are no chirality outliers.

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

Mol	Chain	Res	Type	Group	Models (Total)
2	B	3	U	Sidechain	10
2	B	4	G	Sidechain	9
2	B	14	G	Sidechain	8
2	B	18	A	Sidechain	8
2	B	20	U	Sidechain	8
2	B	16	G	Sidechain	7
2	B	24	G	Sidechain	7
2	B	9	C	Sidechain	6
2	B	11	A	Sidechain	6
2	B	22	A	Sidechain	6
2	B	23	U	Sidechain	6
2	B	12	A	Sidechain	6
2	B	15	A	Sidechain	6
2	B	6	G	Sidechain	6
2	B	10	G	Sidechain	6
2	B	19	A	Sidechain	6
2	B	7	C	Sidechain	5
2	B	8	U	Sidechain	5
2	B	13	A	Sidechain	5
2	B	21	G	Sidechain	5
2	B	5	A	Sidechain	4
2	B	17	C	Sidechain	3
1	A	7	TYR	Sidechain	2
1	A	17	ALA	Mainchain	2
1	A	28	ARG	Sidechain	2
2	B	2	C	Sidechain	2
1	A	46	ARG	Peptide,Sidechain	2
1	A	109	ARG	Sidechain	2

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Mol	Chain	Res	Type	Group	Models (Total)
1	A	8	VAL	Mainchain	1
1	A	10	PHE	Sidechain	1
1	A	123	MET	Peptide	1
1	A	37	THR	Mainchain	1
1	A	87	LEU	Mainchain	1
1	A	18	GLU	Mainchain	1
1	A	78	TYR	Sidechain	1

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	917	965	965	0±0
2	B	494	249	241	1±1
All	All	14110	12140	12073	11

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 0.

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

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:60:PRO:O	1:A:63:ILE:HG22	0.60	1.97	4	1
1:A:46:ARG:HH22	2:B:3:U:P	0.49	2.31	7	1
1:A:51:LEU:HD23	1:A:52:VAL:H	0.47	1.68	5	1
1:A:21:LEU:C	1:A:21:LEU:HD13	0.46	2.36	10	1
2:B:8:U:H2'	2:B:9:C:C6	0.45	2.47	7	1
2:B:22:A:C5	2:B:23:U:C5	0.45	3.05	1	1
2:B:10:G:H1	2:B:14:G:N2	0.44	2.10	1	1
1:A:58:VAL:HG23	1:A:61:GLU:HA	0.44	1.90	8	1
2:B:4:G:N2	2:B:21:G:H1'	0.44	2.28	2	1
2:B:16:G:C5	2:B:17:C:C5	0.42	3.07	8	1
2:B:10:G:H3'	2:B:11:A:C8	0.41	2.50	2	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	119/123 (97%)	111±2 (93±2%)	8±2 (6±2%)	0±0 (0±0%)	37 78
All	All	1190/1230 (97%)	1110 (93%)	76 (6%)	4 (0%)	37 78

All 4 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	16	LEU	1
1	A	95	VAL	1
1	A	47	GLY	1
1	A	13	PRO	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	97/99 (98%)	93±2 (96±2%)	4±2 (4±2%)	31 81
All	All	970/990 (98%)	934 (96%)	36 (4%)	31 81

All 19 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	114	GLU	6
1	A	122	LEU	4
1	A	15	GLU	4
1	A	51	LEU	2
1	A	21	LEU	2
1	A	56	GLU	2

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Mol	Chain	Res	Type	Models (Total)
1	A	35	LYS	2
1	A	11	GLU	2
1	A	86	GLU	2
1	A	50	LYS	1
1	A	72	GLU	1
1	A	63	ILE	1
1	A	103	ILE	1
1	A	76	ILE	1
1	A	9	LYS	1
1	A	18	GLU	1
1	A	95	VAL	1
1	A	61	GLU	1
1	A	110	ASP	1

6.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers	Suiteness
2	B	22/26 (85%)	2±0 (9±0%)	2±1 (7±4%)	0.41±0.00
All	All	220/260 (85%)	20 (9%)	16 (7%)	0.41

The overall RNA backbone suiteness is 0.32.

All unique RNA backbone outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
2	B	6	G	10
2	B	10	G	10

All unique RNA pucker outliers are listed below:

Mol	Chain	Res	Type	Models (Total)
2	B	13	A	3
2	B	20	U	2
2	B	12	A	2
2	B	11	A	2
2	B	19	A	1
2	B	6	G	1
2	B	7	C	1
2	B	5	A	1
2	B	18	A	1
2	B	21	G	1
2	B	10	G	1

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

The completeness of assignment taking into account all chemical shift lists is 20% for the well-defined parts and 20% for the entire structure.

7.1 Chemical shift list 1

File name: `working_cs.cif`

Chemical shift list name: *starch_output*

7.1.1 Bookkeeping

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	425
Number of shifts mapped to atoms	425
Number of unparsed shifts	0
Number of shifts with mapping errors	0
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	1

7.1.2 Chemical shift referencing

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction $\pm$ precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	97	-0.65 ± 0.26	Should be checked
$^{13}\text{C}_\beta$	70	0.75 ± 0.15	Should be checked
$^{13}\text{C}'$	94	-0.25 ± 0.19	None needed (< 0.5 ppm)
^{15}N	98	1.71 ± 0.57	Should be applied

7.1.3 Completeness of resonance assignments

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 20%, i.e. 414 atoms were assigned a chemical shift out of a possible 2071. 0 out of 20 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	^1H	^{13}C	^{15}N
Backbone	280/590 (47%)	0/238 (0%)	189/240 (79%)	91/112 (81%)
Sidechain	134/1001 (13%)	0/654 (0%)	134/318 (42%)	0/29 (0%)

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	Total	¹ H	¹³ C	¹⁵ N
Aromatic	0/44 (0%)	0/21 (0%)	0/22 (0%)	0/1 (0%)
Sugar	0/253 (0%)	0/138 (0%)	0/115 (0%)	0/0 (—%)
Base	0/183 (0%)	0/114 (0%)	0/39 (0%)	0/30 (0%)
Overall	414/2071 (20%)	0/1165 (0%)	323/734 (44%)	91/172 (53%)

Note: This is a solid-state NMR structure, where hydrogen atoms are typically not assigned a chemical shift value, which may lead to lower completeness of assignment measure.

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 20%, i.e. 418 atoms were assigned a chemical shift out of a possible 2089. 0 out of 20 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹ H	¹³ C	¹⁵ N
Backbone	283/595 (48%)	0/240 (0%)	191/242 (79%)	92/113 (81%)
Sidechain	135/1014 (13%)	0/662 (0%)	135/322 (42%)	0/30 (0%)
Aromatic	0/44 (0%)	0/21 (0%)	0/22 (0%)	0/1 (0%)
Sugar	0/253 (0%)	0/138 (0%)	0/115 (0%)	0/0 (—%)
Base	0/183 (0%)	0/114 (0%)	0/39 (0%)	0/30 (0%)
Overall	418/2089 (20%)	0/1175 (0%)	326/740 (44%)	92/174 (53%)

Note: This is a solid-state NMR structure, where hydrogen atoms are typically not assigned a chemical shift value, which may lead to lower completeness of assignment measure.

7.1.4 Statistically unusual chemical shifts [i](#)

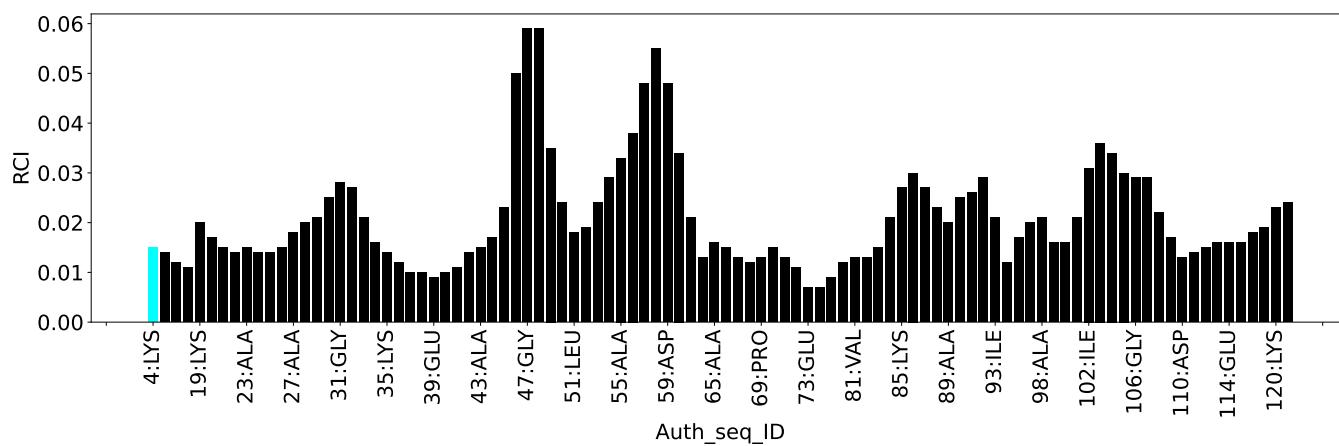
The following table lists the statistically unusual chemical shifts. These are statistical measures, and large deviations from the mean do not necessarily imply incorrect assignments. Molecules containing paramagnetic centres or hemes are expected to give rise to anomalous chemical shifts.

List Id	Chain	Res	Type	Atom	Shift, ppm	Expected range, ppm	Z-score
1	A	53	ILE	CG2	25.20	10.93 – 24.12	5.8

7.1.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:



8 NMR restraints analysis

8.1 Conformationally restricting restraints

The following table provides the summary of experimentally observed NMR restraints in different categories. Restraints are classified into different categories based on the sequence separation of the atoms involved.

Description	Value
Total distance restraints	5
Intra-residue ($ i-j =0$)	0
Sequential ($ i-j =1$)	0
Medium range ($ i-j >1$ and $ i-j <5$)	0
Long range ($ i-j \geq 5$)	0
Inter-chain	5
Hydrogen bond restraints	0
Disulfide bond restraints	0
Total dihedral-angle restraints	0
Number of unmapped restraints	0
Number of restraints per residue	0.0
Number of long range restraints per residue ¹	0.0

¹Long range hydrogen bonds and disulfide bonds are counted as long range restraints while calculating the number of long range restraints per residue

8.2 Residual restraint violations

This section provides the overview of the restraint violations analysis. The violations are binned as small, medium and large violations based on its absolute value. Average number of violations per model is calculated by dividing the total number of violations in each bin by the size of the ensemble.

8.2.1 Average number of distance violations per model

Distance violations less than 0.1 Å are not included in the calculation.

Bins (Å)	Average number of violations per model	Max (Å)
0.1-0.2 (Small)	0.2	0.2
0.2-0.5 (Medium)	1.4	0.47
>0.5 (Large)	1.4	4.29

8.2.2 Average number of dihedral-angle violations per model [i](#)

Dihedral-angle violations less than 1° are not included in the calculation. There are no dihedral-angle violations

9 Distance violation analysis ⓘ

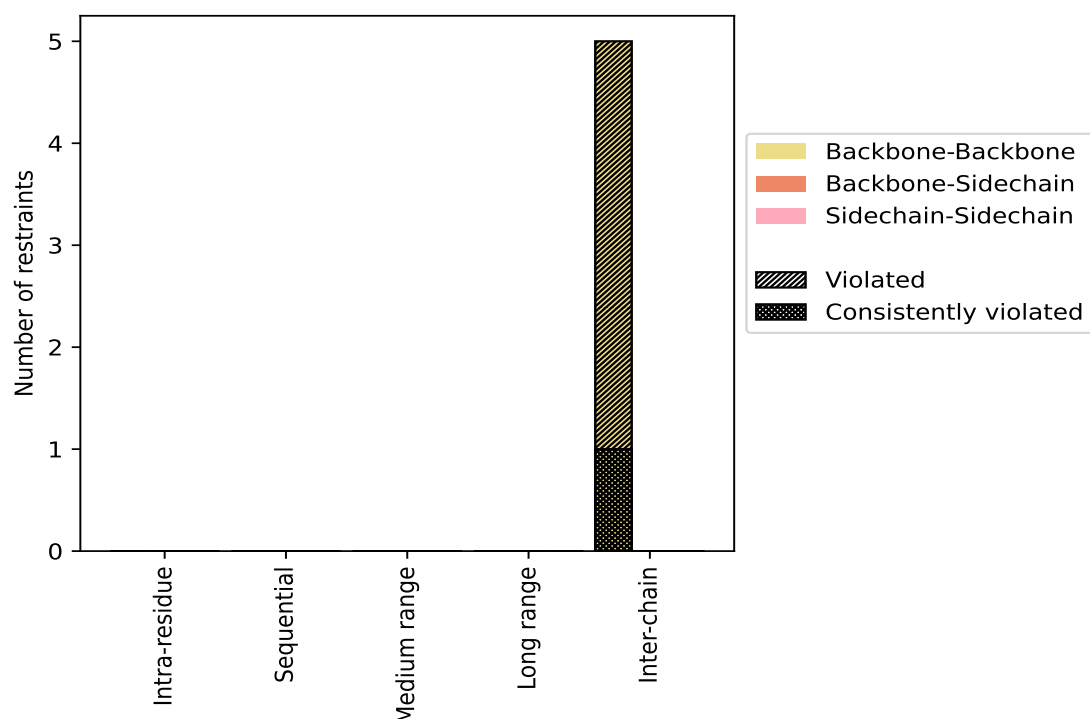
9.1 Summary of distance violations ⓘ

The following table shows the summary of distance violations in different restraint categories based on the sequence separation of the atoms involved. Each category is further sub-divided into three sub-categories based on the atoms involved. Violations less than 0.1 Å are not included in the statistics.

Restrains type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue (i-j =0)	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sequential (i-j =1)	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Medium range (i-j >1 & i-j <5)	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Long range (i-j ≥5)	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Inter-chain	5	100.0	5	100.0	100.0	1	20.0	20.0
Backbone-Backbone	5	100.0	5	100.0	100.0	1	20.0	20.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Hydrogen bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	5	100.0	5	100.0	100.0	1	20.0	20.0
Backbone-Backbone	5	100.0	5	100.0	100.0	1	20.0	20.0
Backbone-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	0	0.0	0	0.0	0.0	0	0.0	0.0

¹ percentage calculated with respect to the total number of distance restraints, ² percentage calculated with respect to the number of restraints in a particular restraint category, ³ violated in at least one model, ⁴ violated in all the models

9.1.1 Bar chart : Distribution of distance restraints and violations [i](#)



Violated and consistently violated restraints are shown using different hatch patterns in their respective categories. The hydrogen bonds and disulfied bonds are counted in their appropriate category on the x-axis

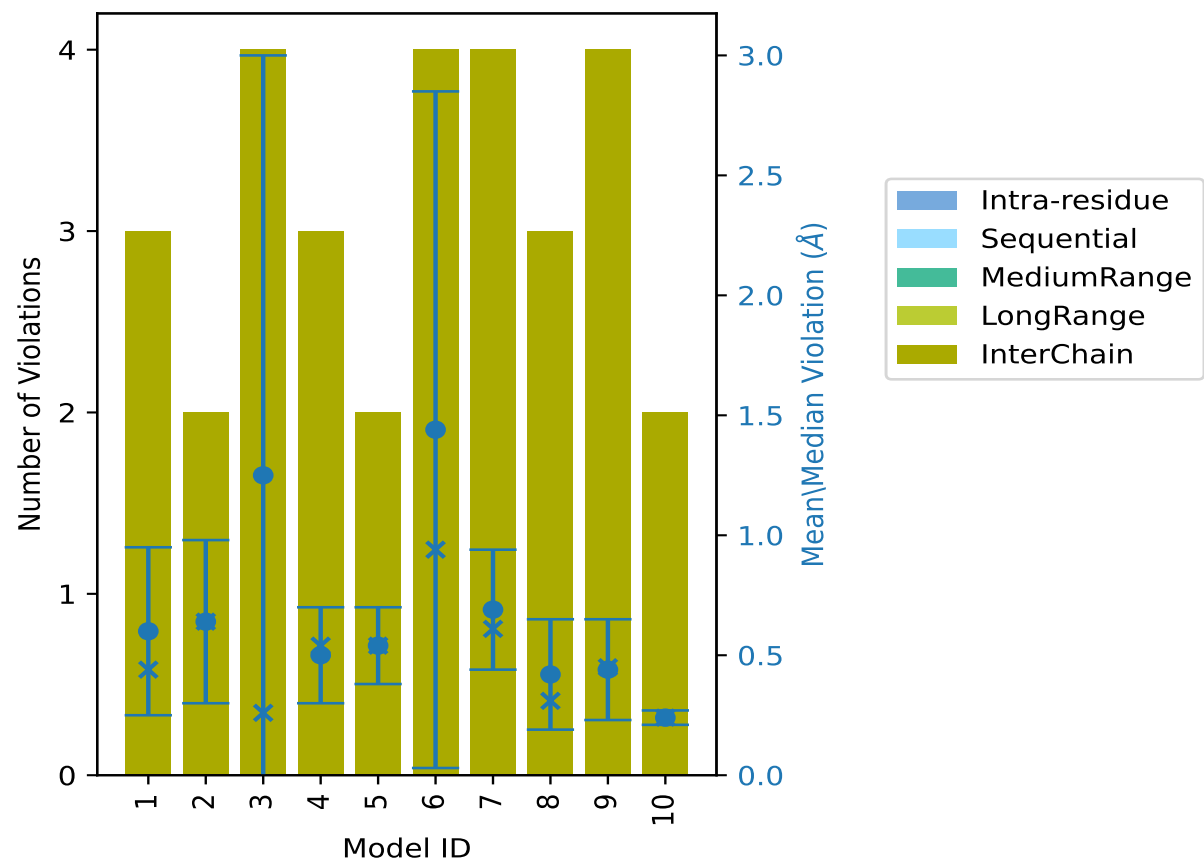
9.2 Distance violation statistics for each model [i](#)

The following table provides the distance violation statistics for each model in the ensemble. Violations less than 0.1 Å are not included in the statistics.

Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
1	0	0	0	0	3	3	0.6	1.09	0.35	0.44
2	0	0	0	0	2	2	0.64	0.98	0.34	0.64
3	0	0	0	0	4	4	1.25	4.29	1.75	0.26
4	0	0	0	0	3	3	0.5	0.73	0.2	0.54
5	0	0	0	0	2	2	0.54	0.7	0.16	0.54
6	0	0	0	0	4	4	1.44	3.8	1.41	0.94
7	0	0	0	0	4	4	0.69	1.09	0.25	0.61
8	0	0	0	0	3	3	0.42	0.74	0.23	0.31
9	0	0	0	0	4	4	0.44	0.74	0.21	0.45
10	0	0	0	0	2	2	0.24	0.27	0.03	0.24

¹Intra-residue restraints, ²Sequential restraints, ³Medium range restraints, ⁴Long range restraints, ⁵Inter-chain restraints, ⁶Standard deviation

9.2.1 Bar graph : Distance Violation statistics for each model ⓘ



The mean(dot),median(x) and the standard deviation are shown in blue with respect to the y axis on the right

9.3 Distance violation statistics for the ensemble ⓘ

Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 0(IR:0, SQ:0, MR:0, LR:0, IC:0) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total	Count ⁶	%
0	0	0	0	0	0	1	10.0
0	0	0	0	0	0	2	20.0
0	0	0	0	1	1	3	30.0

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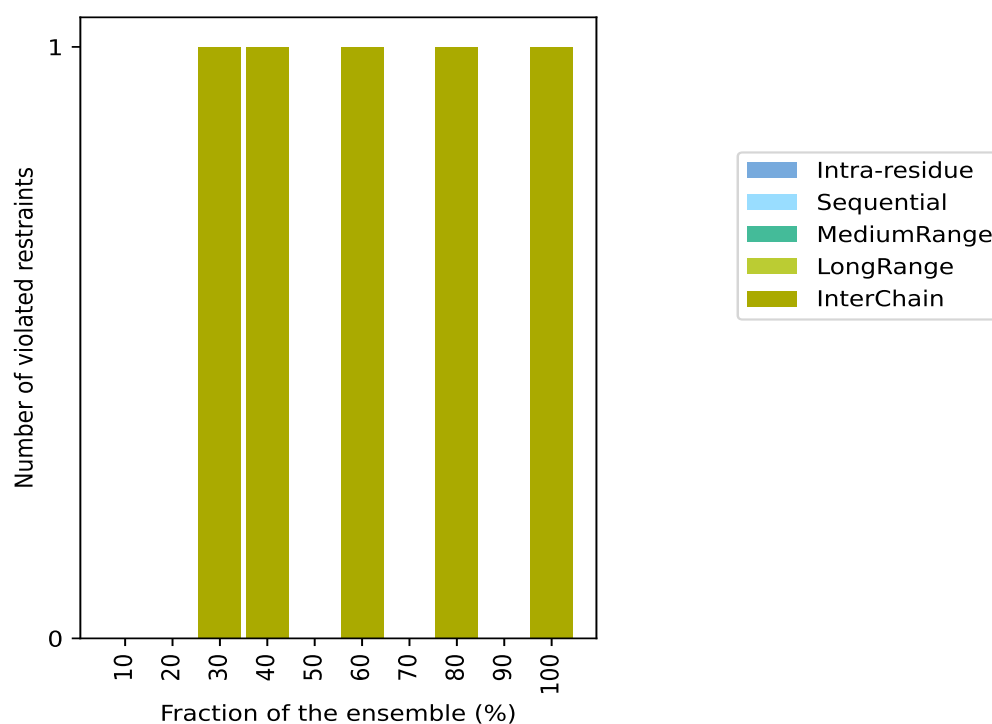
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Number of violated restraints						Fraction of the ensemble	
IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total	Count ⁶	%
0	0	0	0	1	1	4	40.0
0	0	0	0	0	0	5	50.0
0	0	0	0	1	1	6	60.0
0	0	0	0	0	0	7	70.0
0	0	0	0	1	1	8	80.0
0	0	0	0	0	0	9	90.0
0	0	0	0	1	1	10	100.0

¹Intra-residue restraints, ²Sequential restraints, ³Medium range restraints, ⁴Long range restraints,

⁵Inter-chain restraints, ⁶ Number of models with violations

9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)

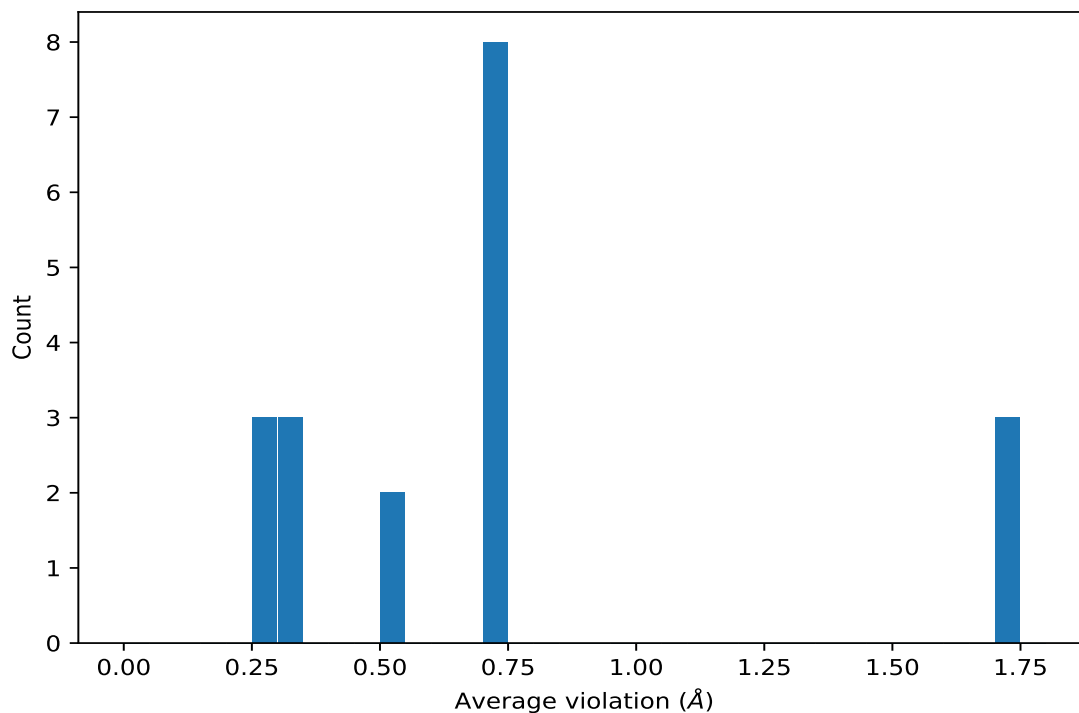


9.4 Most violated distance restraints in the ensemble [i](#)

9.4.1 Histogram : Distribution of mean distance violations [i](#)

The following histogram shows the distribution of the average value of the violation. The average is calculated for each restraint that is violated in more than one model over all the violated models

in the ensemble



9.4.2 Table: Most violated distance restraints [i](#)

The following table provides the mean and the standard deviation of the violation for each restraint sorted by number of violated models and the mean value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,4)	1:93:A:ILE:HG22	2:19:B:A:H61	10	0.74	0.28	0.74
(1,4)	1:93:A:ILE:HD13	2:19:B:A:HO2'	10	0.74	0.28	0.74
(1,4)	1:93:A:ILE:HG23	2:19:B:A:H61	10	0.74	0.28	0.74
(1,4)	1:93:A:ILE:HD13	2:19:B:A:H2'	10	0.74	0.28	0.74
(1,4)	1:93:A:ILE:HG22	2:19:B:A:C6	10	0.74	0.28	0.74
(1,4)	1:93:A:ILE:HD11	2:19:B:A:HO2'	10	0.74	0.28	0.74
(1,4)	1:93:A:ILE:HG22	2:19:B:A:N6	10	0.74	0.28	0.74
(1,4)	1:93:A:ILE:HD12	2:19:B:A:H2'	10	0.74	0.28	0.74
(1,1)	1:36:A:GLY:HA3	2:19:B:A:H1'	8	0.31	0.11	0.26
(1,1)	1:36:A:GLY:HA2	2:19:B:A:HO2'	8	0.31	0.11	0.26
(1,1)	1:36:A:GLY:HA2	2:19:B:A:O2'	8	0.31	0.11	0.26
(1,3)	1:84:A:LYS:HD2	2:20:B:U:H5	6	1.72	1.68	0.82
(1,3)	1:84:A:LYS:HB2	2:20:B:U:OP1	6	1.72	1.68	0.82
(1,3)	1:84:A:LYS:HZ1	2:20:B:U:H5	6	1.72	1.68	0.82
(1,2)	1:38:A:ASN:H	2:21:B:G:O6	4	0.28	0.13	0.29

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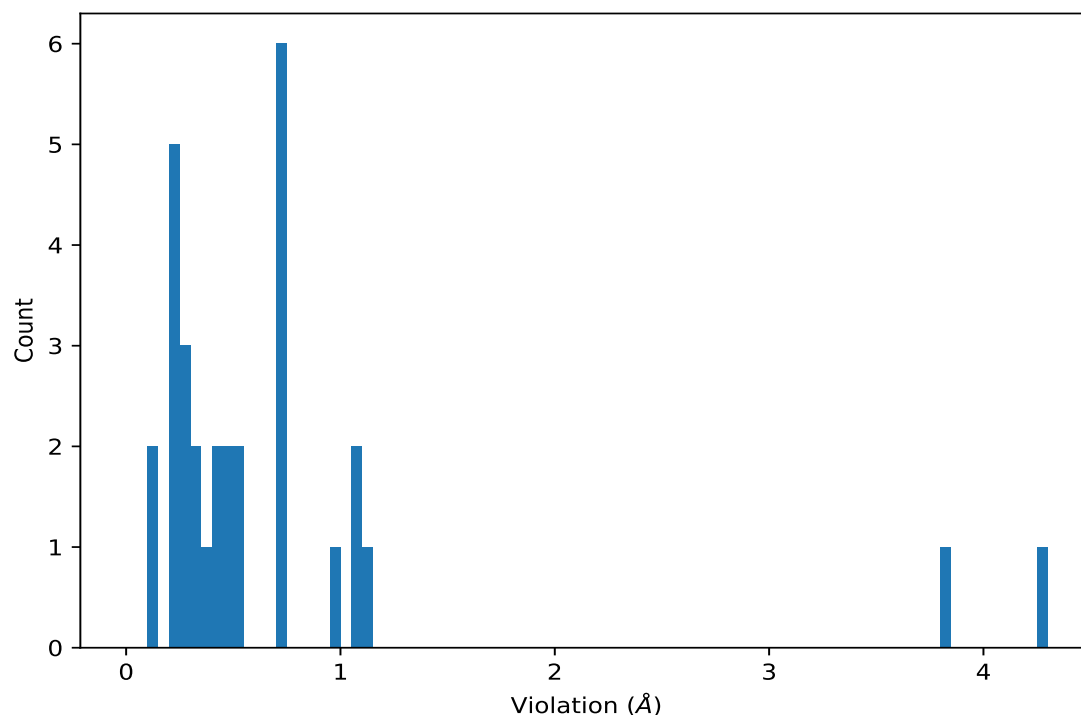
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,2)	1:38:A:ASN:HD21	2:21:B:G:N7	4	0.28	0.13	0.29
(1,2)	1:38:A:ASN:HB2	2:21:B:G:O6	4	0.28	0.13	0.29
(1,5)	1:98:A:ALA:H	2:20:B:U:OP2	3	0.5	0.17	0.45
(1,5)	1:98:A:ALA:HB2	2:20:B:U:H5	3	0.5	0.17	0.45

¹Number of violated models, ²Standard deviation

9.5 All violated distance restraints [i](#)

9.5.1 Histogram : Distribution of distance violations [i](#)

The following histogram shows the distribution of the absolute value of the violation for all violated restraints in the ensemble.



9.5.2 Table : All distance violations [i](#)

The following table lists the absolute value of the violation for each restraint in the ensemble sorted by its value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,3)	1:84:A:LYS:HB2	2:20:B:U:OP1	3	4.29
(1,3)	1:84:A:LYS:HB2	2:20:B:U:OP1	6	3.8
(1,4)	1:93:A:ILE:HD13	2:19:B:A:HO2'	6	1.14
(1,4)	1:93:A:ILE:HG22	2:19:B:A:H61	1	1.09
(1,3)	1:84:A:LYS:HD2	2:20:B:U:H5	7	1.09
(1,4)	1:93:A:ILE:HD13	2:19:B:A:HO2'	2	0.98
(1,4)	1:93:A:ILE:HD11	2:19:B:A:HO2'	8	0.74
(1,4)	1:93:A:ILE:HG22	2:19:B:A:N6	9	0.74
(1,5)	1:98:A:ALA:H	2:20:B:U:OP2	6	0.73
(1,4)	1:93:A:ILE:HG22	2:19:B:A:H61	4	0.73
(1,4)	1:93:A:ILE:HG22	2:19:B:A:C6	7	0.71
(1,4)	1:93:A:ILE:HD13	2:19:B:A:H2'	5	0.7
(1,3)	1:84:A:LYS:HZ1	2:20:B:U:H5	4	0.54
(1,1)	1:36:A:GLY:HA2	2:19:B:A:O2'	7	0.51
(1,1)	1:36:A:GLY:HA3	2:19:B:A:H1'	9	0.47
(1,5)	1:98:A:ALA:HB2	2:20:B:U:H5	7	0.45
(1,3)	1:84:A:LYS:HD2	2:20:B:U:H5	1	0.44
(1,2)	1:38:A:ASN:H	2:21:B:G:O6	9	0.43
(1,2)	1:38:A:ASN:HB2	2:21:B:G:O6	5	0.38
(1,5)	1:98:A:ALA:H	2:20:B:U:OP2	8	0.31
(1,4)	1:93:A:ILE:HG23	2:19:B:A:H61	3	0.3
(1,1)	1:36:A:GLY:HA3	2:19:B:A:H1'	2	0.29
(1,1)	1:36:A:GLY:HA3	2:19:B:A:H1'	1	0.28
(1,4)	1:93:A:ILE:HD12	2:19:B:A:H2'	10	0.27
(1,1)	1:36:A:GLY:HA3	2:19:B:A:H1'	4	0.24
(1,1)	1:36:A:GLY:HA2	2:19:B:A:HO2'	3	0.23
(1,1)	1:36:A:GLY:HA3	2:19:B:A:H1'	10	0.22
(1,1)	1:36:A:GLY:HA3	2:19:B:A:H1'	8	0.21
(1,2)	1:38:A:ASN:HD21	2:21:B:G:N7	3	0.2
(1,3)	1:84:A:LYS:HD2	2:20:B:U:H5	9	0.14
(1,2)	1:38:A:ASN:HD21	2:21:B:G:N7	6	0.1

10 Dihedral-angle violation analysis ⓘ

No dihedral-angle restraints found