



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:42 AM UTC

PDB ID : 1R4O / pdb_00001r4o
Title : Crystallographic analysis of the interaction of the glucocorticoid receptor with DNA
Authors : Luisi, B.F.; Xu, W.X.; Otwinowski, Z.; Freedman, L.P.; Yamamoto, K.R.; Sigler, P.B.
Deposited on : 2003-10-07
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray 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/XrayValidationReportHelp>

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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

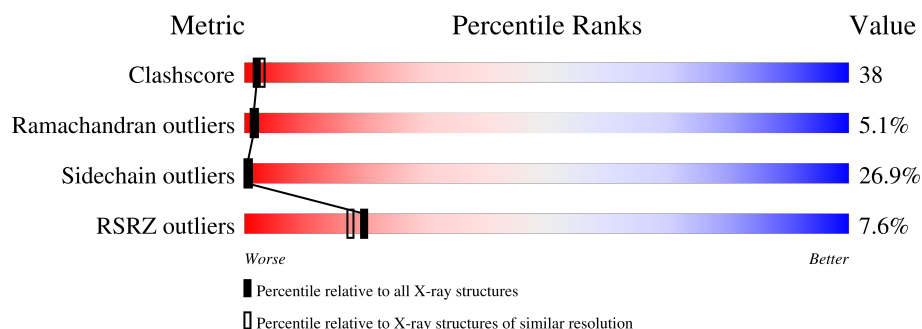
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	C	19	
1	D	19	
2	A	92	
2	B	92	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2121 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a DNA chain called 5'-D(*CP*CP*AP*GP*AP*AP*CP*AP*TP*CP*GP*AP*TP*GP*TP*TP*CP*TP*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	19	Total	C	N	O	P	0	0	0
			385	185	70	112	18			
1	D	19	Total	C	N	O	P	0	0	0
			385	185	70	112	18			

- Molecule 2 is a protein called Glucocorticoid receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	80	Total	C	N	O	S	0	0	0
			619	379	122	106	12			
2	B	80	Total	C	N	O	S	0	0	0
			619	379	122	106	12			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	434	MET	-	cloning artifact	UNP P06536
A	435	LYS	-	cloning artifact	UNP P06536
A	436	PRO	-	cloning artifact	UNP P06536
A	437	ALA	-	cloning artifact	UNP P06536
A	438	ARG	-	cloning artifact	UNP P06536
A	439	PRO	-	cloning artifact	UNP P06536
B	434	MET	-	cloning artifact	UNP P06536
B	435	LYS	-	cloning artifact	UNP P06536
B	436	PRO	-	cloning artifact	UNP P06536
B	437	ALA	-	cloning artifact	UNP P06536
B	438	ARG	-	cloning artifact	UNP P06536
B	439	PRO	-	cloning artifact	UNP P06536

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total 2	Zn 2	0	0
3	B	2	Total 2	Zn 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	23	Total 23	O 23	0	0
4	D	13	Total 13	O 13	0	0
4	A	32	Total 32	O 32	0	0
4	B	41	Total 41	O 41	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-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. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

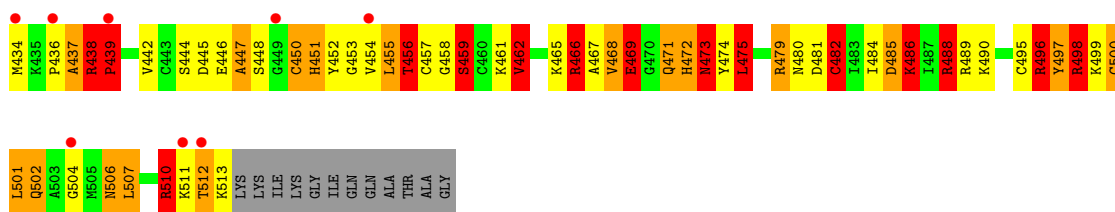
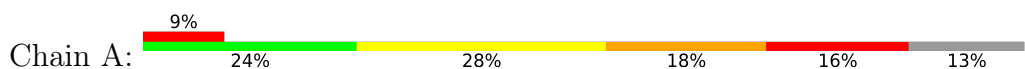
- Molecule 1: 5'-D(*CP*CP*AP*GP*AP*AP*CP*AP*TP*CP*GP*AP*TP*GP*TP*TP*CP*TP*G)-3'



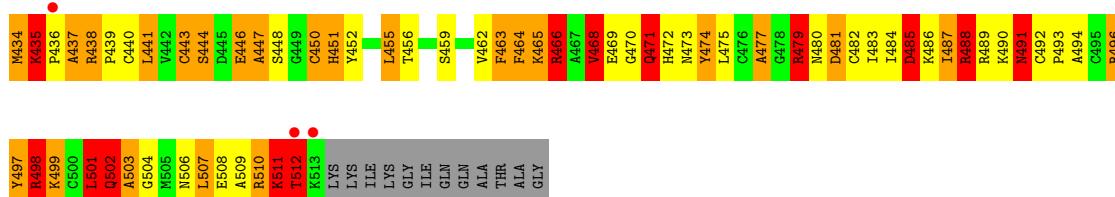
- Molecule 1: 5'-D(*CP*CP*AP*GP*AP*AP*CP*AP*TP*CP*GP*AP*TP*GP*TP*TP*CP*TP*G)-3'



- Molecule 2: Glucocorticoid receptor



- Molecule 2: Glucocorticoid receptor



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	38.50Å 97.50Å 120.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 2.50 6.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (6.00-2.50) 79.1 (6.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.78 (at 2.46Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.196 , (Not available) 0.271 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	44.8	Xtriage
Anisotropy	0.528	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.63 , 138.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	2121	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.88% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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 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	C	1.49	6/431 (1.4%)	3.04	57/663 (8.6%)
1	D	1.60	8/431 (1.9%)	2.88	58/663 (8.7%)
2	A	1.32	0/628	3.01	71/837 (8.5%)
2	B	1.40	1/628 (0.2%)	3.46	92/837 (11.0%)
All	All	1.44	15/2118 (0.7%)	3.12	278/3000 (9.3%)

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 outliers	#Planarity outliers
1	C	0	2
1	D	0	4
2	A	0	8
2	B	0	13
All	All	0	27

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	2	DC	O3'-P	8.36	1.73	1.61
1	C	17	DC	O3'-P	8.12	1.73	1.61
1	C	6	DA	O3'-P	7.66	1.72	1.61
1	D	10	DC	O3'-P	6.97	1.71	1.61
1	D	17	DC	O3'-P	6.70	1.71	1.61
1	C	2	DC	P-OP2	6.35	1.61	1.48
1	D	14	DG	O3'-P	-6.21	1.51	1.61
1	D	17	DC	P-OP1	6.17	1.60	1.48
1	C	8	DA	O3'-P	5.83	1.69	1.61
1	D	8	DA	O3'-P	-5.74	1.52	1.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	5	DA	P-O5'	5.65	1.71	1.60
1	C	9	DT	P-OP2	5.64	1.59	1.48
2	B	498	ARG	NE-CZ	5.49	1.39	1.33
1	D	4	DG	O3'-P	5.15	1.68	1.61
1	D	17	DC	O4'-C1'	5.08	1.52	1.41

All (278) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	479	ARG	CD-NE-CZ	39.51	179.71	124.40
1	C	16	DT	P-O3'-C3'	15.12	142.88	120.20
2	B	498	ARG	CD-NE-CZ	-15.11	103.24	124.40
1	C	1	DC	P-O3'-C3'	14.88	142.51	120.20
1	D	12	DA	P-O5'-C5'	-13.27	100.09	120.00
1	C	18	DT	P-O5'-C5'	-13.24	100.14	120.00
2	B	481	ASP	CA-C-O	-12.74	105.61	122.31
2	B	510	ARG	N-CA-C	-11.95	98.65	113.23
2	B	464	PHE	CA-C-O	11.79	133.51	120.90
2	B	498	ARG	NE-CZ-NH2	-11.70	108.67	119.20
1	D	9	DT	P-O5'-C5'	-11.68	102.48	120.00
1	C	5	DA	P-O3'-C3'	11.28	137.12	120.20
1	C	2	DC	O5'-C5'-C4'	-11.07	94.19	110.80
2	A	467	ALA	CA-C-N	10.65	134.33	120.60
2	A	467	ALA	C-N-CA	10.65	134.33	120.60
2	A	499	LYS	CA-C-N	10.63	135.56	120.79
2	A	499	LYS	C-N-CA	10.63	135.56	120.79
1	C	3	DA	P-O5'-C5'	-10.37	104.44	120.00
2	B	481	ASP	CA-CB-CG	10.36	122.96	112.60
2	A	458	GLY	CA-C-N	10.23	134.54	120.63
2	A	458	GLY	C-N-CA	10.23	134.54	120.63
2	A	498	ARG	CD-NE-CZ	-10.08	110.29	124.40
2	B	463	PHE	CA-C-O	-10.06	110.53	120.90
2	B	444	SER	N-CA-CB	-10.03	102.74	114.17
1	D	19	DG	C1'-O4'-C4'	-9.88	94.89	109.70
1	D	17	DC	O5'-P-OP2	9.84	137.51	108.00
1	C	5	DA	P-O5'-C5'	-9.83	105.25	120.00
1	C	6	DA	P-O5'-C5'	-9.82	105.27	120.00
2	A	466	ARG	O-C-N	-9.78	110.78	122.22
1	C	13	DT	O4'-C1'-C2'	-9.76	91.76	106.40
1	C	8	DA	P-O5'-C5'	-9.74	105.40	120.00
2	B	464	PHE	O-C-N	-9.73	111.97	122.09
2	B	482	CYS	CA-C-O	-9.68	109.85	121.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	473	ASN	O-C-N	9.48	134.28	123.28
2	A	473	ASN	CB-CA-C	-9.48	99.28	111.70
2	A	510	ARG	N-CA-C	9.40	124.42	109.86
2	B	488	ARG	NE-CZ-NH1	-9.37	112.13	121.50
2	B	485	ASP	CB-CA-C	9.28	128.88	110.42
2	B	463	PHE	N-CA-C	-9.14	100.88	111.03
2	B	507	LEU	N-CA-C	9.05	120.75	111.07
1	C	18	DT	O4'-C1'-C2'	-9.02	92.87	106.40
1	C	14	DG	P-O3'-C3'	9.01	133.71	120.20
2	B	510	ARG	CA-C-O	8.80	130.24	120.12
1	D	11	DG	N1-C2-N2	-8.79	103.12	116.30
1	D	13	DT	O3'-P-O5'	-8.74	90.89	104.00
1	D	11	DG	P-O5'-C5'	-8.69	106.96	120.00
2	A	471	GLN	OE1-CD-NE2	-8.64	113.96	122.60
1	D	10	DC	O3'-P-O5'	-8.54	91.19	104.00
1	D	10	DC	C2'-C3'-O3'	8.51	124.27	111.50
1	D	11	DG	O5'-C5'-C4'	-8.47	98.09	110.80
2	B	501	LEU	O-C-N	-8.46	111.34	122.59
2	B	496	ARG	NE-CZ-NH2	8.46	126.81	119.20
1	C	9	DT	O5'-C5'-C4'	-8.42	98.17	110.80
2	A	502	GLN	OE1-CD-NE2	-8.36	114.24	122.60
1	C	17	DC	P-O3'-C3'	-8.32	107.72	120.20
1	D	2	DC	O4'-C1'-C2'	-8.28	93.99	106.40
2	B	511	LYS	CA-C-O	8.27	129.48	120.80
2	A	462	VAL	CB-CA-C	-8.26	101.39	111.97
1	D	17	DC	O4'-C4'-C3'	-8.15	93.17	105.40
2	A	474	TYR	CA-C-O	-8.08	111.65	120.30
2	A	455	LEU	CA-C-O	-8.07	111.72	120.92
1	D	12	DA	O4'-C1'-C2'	-7.98	94.43	106.40
2	B	450	CYS	CA-C-O	-7.95	111.67	120.81
1	D	19	DG	O4'-C4'-C3'	-7.89	93.56	105.40
1	C	12	DA	N9-C1'-C2'	-7.87	101.69	113.50
2	A	445	ASP	CA-CB-CG	7.87	120.47	112.60
2	B	468	VAL	O-C-N	7.87	129.51	121.87
1	C	11	DG	P-O3'-C3'	7.85	131.97	120.20
1	C	9	DT	P-O3'-C3'	7.83	131.95	120.20
2	A	501	LEU	O-C-N	-7.79	113.86	122.12
2	A	453	GLY	CA-C-O	-7.78	110.25	119.04
1	D	12	DA	O5'-C5'-C4'	-7.73	99.20	110.80
2	B	511	LYS	CA-C-N	7.70	136.25	121.54
2	B	511	LYS	C-N-CA	7.70	136.25	121.54
1	D	1	DC	P-O3'-C3'	7.62	131.63	120.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	468	VAL	CA-C-O	-7.61	113.03	120.95
2	A	451	HIS	CA-C-N	7.59	136.03	121.54
2	A	451	HIS	C-N-CA	7.59	136.03	121.54
2	B	444	SER	O-C-N	-7.57	114.36	121.56
2	B	488	ARG	NH1-CZ-NH2	7.53	129.09	119.30
1	D	5	DA	O5'-P-OP1	7.52	131.56	109.00
2	A	473	ASN	CA-C-O	-7.51	111.35	120.89
1	C	9	DT	O4'-C1'-N1	7.50	119.64	108.40
2	B	440	CYS	CA-C-O	-7.47	112.73	120.80
2	A	472	HIS	CA-C-N	7.43	135.47	121.87
2	A	472	HIS	C-N-CA	7.43	135.47	121.87
2	A	459	SER	CB-CA-C	-7.41	99.47	110.95
1	D	13	DT	O4'-C1'-C2'	-7.41	95.29	106.40
2	B	447	ALA	CA-C-O	-7.30	113.40	121.87
2	B	468	VAL	N-CA-CB	7.28	120.44	110.54
1	D	4	DG	O5'-C5'-C4'	-7.23	99.96	110.80
2	A	485	ASP	CA-CB-CG	7.22	119.83	112.60
2	B	492	CYS	CA-C-N	7.17	127.50	119.32
2	B	492	CYS	C-N-CA	7.17	127.50	119.32
1	C	18	DT	O5'-C5'-C4'	-7.16	100.06	110.80
2	B	436	PRO	N-CA-C	-7.16	98.88	110.55
2	A	459	SER	O-C-N	7.13	129.73	122.03
1	D	11	DG	N3-C2-N2	7.10	130.36	119.70
2	B	463	PHE	N-CA-CB	7.07	120.23	109.98
1	D	18	DT	O4'-C1'-N1	7.03	118.95	108.40
2	A	504	GLY	CA-C-O	-6.98	111.81	119.27
1	C	12	DA	O5'-P-OP1	6.97	129.92	109.00
1	D	19	DG	C5'-C4'-O4'	6.91	119.77	109.40
1	C	6	DA	C5'-C4'-C3'	-6.91	104.54	114.90
1	C	19	DG	P-O5'-C5'	-6.90	109.66	120.00
2	A	482	CYS	CA-C-N	6.87	131.76	122.01
2	A	482	CYS	C-N-CA	6.87	131.76	122.01
1	C	15	DT	O4'-C1'-C2'	-6.86	96.11	106.40
1	D	14	DG	C4'-C3'-O3'	-6.84	99.73	110.00
1	C	12	DA	P-O3'-C3'	6.84	130.46	120.20
1	D	16	DT	C3'-C2'-C1'	-6.84	91.34	101.60
1	C	17	DC	O3'-P-O5'	-6.83	93.75	104.00
1	D	7	DC	O4'-C1'-N1	6.80	118.61	108.40
1	D	17	DC	C1'-O4'-C4'	-6.79	99.51	109.70
2	B	482	CYS	O-C-N	6.75	130.86	123.10
1	C	3	DA	P-O3'-C3'	-6.73	110.10	120.20
1	C	2	DC	C1'-O4'-C4'	-6.72	99.62	109.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	2	DC	O5'-C5'-C4'	-6.68	100.77	110.80
2	B	447	ALA	O-C-N	6.67	130.40	122.79
2	B	486	LYS	O-C-N	-6.67	114.59	122.20
1	C	19	DG	OP1-P-OP2	-6.63	100.10	120.00
2	B	510	ARG	CB-CA-C	6.63	122.34	111.26
1	C	19	DG	C4'-C3'-C2'	-6.62	92.46	102.40
2	A	459	SER	N-CA-C	6.60	118.35	111.03
1	C	6	DA	O4'-C1'-N9	6.58	118.27	108.40
1	C	14	DG	C2'-C3'-O3'	6.55	121.32	111.50
1	C	6	DA	C4'-C3'-C2'	-6.51	92.63	102.40
2	B	470	GLY	N-CA-C	6.51	120.25	111.20
2	B	459	SER	CA-CB-OG	6.48	124.06	111.10
2	B	498	ARG	CA-C-N	6.47	131.53	120.72
2	B	498	ARG	C-N-CA	6.47	131.53	120.72
2	B	443	CYS	N-CA-C	6.46	121.96	113.30
2	A	444	SER	CA-C-O	-6.45	112.70	120.32
2	B	477	ALA	CA-C-O	-6.42	111.54	119.18
1	C	2	DC	O5'-P-OP1	6.40	128.20	109.00
2	B	465	LYS	CA-C-N	6.39	128.75	120.44
2	B	465	LYS	C-N-CA	6.39	128.75	120.44
2	B	498	ARG	CA-C-O	-6.38	112.63	119.97
2	B	468	VAL	CB-CA-C	-6.37	103.54	112.14
2	B	437	ALA	CA-C-O	-6.33	114.07	121.16
1	D	19	DG	C2'-C3'-O3'	6.30	120.96	111.50
1	C	16	DT	O4'-C1'-N1	6.30	117.85	108.40
1	C	11	DG	N9-C1'-C2'	-6.27	104.09	113.50
1	D	10	DC	C5'-C4'-C3'	-6.27	105.50	114.90
2	B	494	ALA	N-CA-CB	6.25	119.07	110.01
2	B	435	LYS	CB-CA-C	-6.24	99.76	109.11
1	C	19	DG	O5'-C5'-C4'	-6.23	101.46	110.80
2	B	508	GLU	CB-CA-C	-6.22	102.86	112.07
2	B	508	GLU	N-CA-CB	6.22	117.32	110.35
1	D	9	DT	C5'-C4'-O4'	6.22	118.73	109.40
1	D	8	DA	P-O5'-C5'	-6.21	110.68	120.00
1	C	1	DC	C2'-C3'-O3'	6.21	120.81	111.50
2	B	474	TYR	CA-C-O	-6.18	114.12	121.66
1	C	18	DT	N3-C4-O4	-6.16	113.36	122.60
1	D	3	DA	O5'-C5'-C4'	-6.12	101.62	110.80
1	D	1	DC	O4'-C1'-C2'	-6.11	97.23	106.40
2	A	458	GLY	CA-C-O	-6.09	114.78	121.05
1	C	18	DT	P-O3'-C3'	-6.08	111.08	120.20
2	B	438	ARG	CB-CA-C	6.06	116.51	110.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	474	TYR	CA-CB-CG	6.04	124.78	113.90
2	B	503	ALA	O-C-N	-6.04	114.60	122.39
1	D	2	DC	O3'-P-O5'	-6.03	94.96	104.00
2	A	474	TYR	N-CA-CB	-6.00	101.30	110.65
1	C	6	DA	O4'-C4'-C3'	-5.97	96.44	105.40
1	C	9	DT	C6-N1-C1'	5.96	128.29	119.35
1	D	4	DG	O4'-C1'-C2'	-5.96	97.46	106.40
2	A	459	SER	CA-C-N	-5.96	111.70	120.28
2	A	459	SER	C-N-CA	-5.96	111.70	120.28
2	A	488	ARG	NH1-CZ-NH2	-5.95	111.56	119.30
2	A	442	VAL	O-C-N	-5.94	115.63	122.17
2	B	497	TYR	CA-C-O	5.92	126.69	120.42
2	A	479	ARG	NH1-CZ-NH2	-5.92	111.61	119.30
2	A	475	LEU	CA-C-O	-5.90	113.99	120.24
2	A	462	VAL	O-C-N	5.89	127.68	121.91
1	C	2	DC	O4'-C4'-C3'	-5.89	96.57	105.40
2	B	496	ARG	CA-C-O	-5.89	114.18	120.42
2	A	450	CYS	CA-C-O	-5.88	114.99	121.58
1	C	6	DA	C2'-C3'-O3'	5.85	120.28	111.50
1	D	16	DT	N3-C4-O4	-5.85	113.82	122.60
2	B	488	ARG	CA-C-O	-5.85	110.22	119.23
1	C	7	DC	O5'-C5'-C4'	-5.82	102.08	110.80
2	B	497	TYR	O-C-N	-5.81	115.53	122.15
2	A	459	SER	CA-C-O	-5.80	114.93	120.90
2	B	491	ASN	O-C-N	-5.79	115.19	122.20
2	B	443	CYS	CA-C-N	-5.79	116.20	126.45
2	B	443	CYS	C-N-CA	-5.79	116.20	126.45
1	D	17	DC	O4'-C1'-C2'	-5.78	97.74	106.40
2	A	451	HIS	CA-CB-CG	-5.76	108.04	113.80
1	C	17	DC	OP1-P-O3'	-5.76	90.73	108.00
1	D	15	DT	C2-N3-C4	-5.75	118.38	127.00
1	D	5	DA	O4'-C1'-C2'	-5.74	97.79	106.40
1	D	10	DC	P-O3'-C3'	-5.70	111.66	120.20
1	D	10	DC	O4'-C4'-C3'	-5.69	96.86	105.40
2	A	468	VAL	CA-C-O	-5.69	115.35	121.27
1	C	5	DA	C4'-C3'-C2'	-5.66	93.92	102.40
1	C	16	DT	P-O5'-C5'	-5.65	111.53	120.00
2	A	466	ARG	CA-C-O	5.65	126.48	120.10
1	C	12	DA	O5'-C5'-C4'	5.64	119.26	110.80
2	B	446	GLU	O-C-N	5.64	130.09	122.59
1	C	4	DG	O3'-P-O5'	-5.63	95.55	104.00
2	A	471	GLN	CG-CD-NE2	5.62	124.83	116.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	438	ARG	NE-CZ-NH2	5.61	124.25	119.20
1	C	8	DA	O4'-C1'-C2'	-5.60	98.00	106.40
1	D	18	DT	C3'-C2'-C1'	-5.59	93.22	101.60
2	B	480	ASN	CB-CG-ND2	-5.58	108.03	116.40
1	C	10	DC	O5'-C5'-C4'	-5.57	102.45	110.80
2	B	488	ARG	N-CA-C	5.56	119.77	113.15
2	B	480	ASN	CB-CG-OD1	5.56	131.92	120.80
1	C	11	DG	OP2-P-O3'	-5.55	91.34	108.00
2	A	495	CYS	N-CA-C	-5.55	105.63	112.90
2	A	506	ASN	CA-CB-CG	5.55	118.15	112.60
1	D	3	DA	OP1-P-OP2	5.53	136.58	120.00
1	D	6	DA	O5'-P-OP1	5.52	125.56	109.00
2	B	496	ARG	NE-CZ-NH1	-5.52	115.98	121.50
2	B	434	MET	O-C-N	5.51	131.82	123.00
2	B	512	THR	CA-C-N	5.50	131.61	121.70
2	B	512	THR	C-N-CA	5.50	131.61	121.70
1	C	1	DC	O4'-C4'-C3'	-5.50	97.16	105.40
1	C	9	DT	O3'-P-O5'	-5.49	95.76	104.00
2	A	442	VAL	N-CA-CB	-5.49	104.78	112.35
1	D	2	DC	P-O3'-C3'	-5.45	112.02	120.20
2	A	496	ARG	N-CA-C	-5.45	104.98	111.03
1	D	16	DT	O4'-C1'-N1	5.43	116.55	108.40
2	B	446	GLU	CG-CD-OE1	5.43	130.89	118.40
1	D	13	DT	C2-N3-C4	-5.43	118.86	127.00
1	D	16	DT	C4'-C3'-C2'	-5.42	94.27	102.40
2	A	486	LYS	N-CA-CB	5.38	118.64	110.14
2	A	468	VAL	O-C-N	5.36	127.30	121.94
2	A	500	CYS	N-CA-C	-5.34	104.51	111.02
2	B	498	ARG	NH1-CZ-NH2	5.32	126.22	119.30
2	B	502	GLN	CB-CG-CD	-5.30	103.59	112.60
2	A	482	CYS	CA-C-O	-5.29	115.73	121.87
1	D	1	DC	O3'-P-O5'	-5.28	96.07	104.00
2	A	436	PRO	N-CA-CB	5.25	108.76	103.25
2	A	484	ILE	CA-C-O	-5.24	114.88	120.59
2	A	447	ALA	N-CA-CB	-5.23	102.78	110.26
2	B	471	GLN	N-CA-C	5.22	117.18	107.99
2	B	489	ARG	CA-C-O	5.22	125.80	119.38
2	B	466	ARG	CA-CB-CG	5.20	124.50	114.10
2	A	498	ARG	O-C-N	-5.19	115.68	122.59
2	A	474	TYR	CA-C-N	-5.19	115.45	122.77
2	A	474	TYR	C-N-CA	-5.19	115.45	122.77
2	B	506	ASN	CA-C-O	-5.18	114.17	120.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	10	DC	C1'-O4'-C4'	-5.17	101.94	109.70
2	A	507	LEU	CB-CA-C	5.17	118.96	110.95
1	D	7	DC	C4'-C3'-C2'	-5.16	94.66	102.40
2	A	500	CYS	CA-C-O	-5.15	115.41	121.07
2	A	444	SER	CB-CA-C	-5.12	103.02	111.36
2	A	472	HIS	CA-C-O	5.11	127.35	121.32
2	A	496	ARG	CD-NE-CZ	-5.11	117.24	124.40
2	B	466	ARG	CG-CD-NE	-5.11	100.77	112.00
1	D	7	DC	C3'-C2'-C1'	-5.10	93.95	101.60
1	C	19	DG	O4'-C4'-C3'	-5.09	97.76	105.40
1	D	6	DA	C1'-O4'-C4'	-5.09	102.06	109.70
2	B	512	THR	O-C-N	5.09	129.37	122.59
2	B	499	LYS	N-CA-CB	5.09	118.84	110.39
1	D	9	DT	P-O3'-C3'	-5.08	112.57	120.20
2	B	456	THR	OG1-CB-CG2	5.08	119.47	109.30
2	B	491	ASN	CA-CB-CG	5.08	117.68	112.60
2	B	435	LYS	O-C-N	5.08	128.09	121.64
2	A	480	ASN	OD1-CG-ND2	-5.07	117.53	122.60
1	D	14	DG	O5'-P-OP2	5.07	123.21	108.00
2	B	506	ASN	OD1-CG-ND2	-5.07	117.53	122.60
1	C	13	DT	O5'-P-OP2	5.07	123.20	108.00
1	C	11	DG	O4'-C1'-N9	5.06	115.98	108.40
2	B	473	ASN	CA-C-O	-5.05	115.67	122.44
1	D	2	DC	P-O5'-C5'	5.05	127.58	120.00
2	A	472	HIS	CA-CB-CG	5.05	118.85	113.80
2	B	504	GLY	O-C-N	-5.05	116.10	122.41
2	A	469	GLU	CB-CG-CD	5.04	121.17	112.60
1	D	15	DT	P-O3'-C3'	5.04	127.76	120.20
1	D	11	DG	P-O3'-C3'	-5.04	112.64	120.20
2	B	503	ALA	CA-C-N	5.04	131.46	121.53
2	B	503	ALA	C-N-CA	5.04	131.46	121.53
2	B	481	ASP	CB-CA-C	-5.01	103.81	112.44
2	A	456	THR	N-CA-C	5.01	115.14	108.38
2	B	436	PRO	CA-C-N	-5.00	113.77	120.87
2	B	436	PRO	C-N-CA	-5.00	113.77	120.87

There are no chirality outliers.

All (27) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	437	ALA	Mainchain
2	A	450	CYS	Mainchain

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Mol	Chain	Res	Type	Group
2	A	466	ARG	Mainchain
2	A	482	CYS	Mainchain
2	A	488	ARG	Sidechain
2	A	496	ARG	Sidechain,Mainchain
2	A	498	ARG	Sidechain
2	B	441	LEU	Mainchain
2	B	451	HIS	Mainchain
2	B	452	TYR	Sidechain
2	B	463	PHE	Mainchain
2	B	466	ARG	Sidechain
2	B	468	VAL	Mainchain
2	B	471	GLN	Mainchain
2	B	485	ASP	Mainchain
2	B	491	ASN	Mainchain
2	B	497	TYR	Sidechain
2	B	498	ARG	Sidechain
2	B	501	LEU	Mainchain
2	B	502	GLN	Sidechain
1	C	19	DG	Sidechain
1	C	2	DC	Sidechain
1	D	12	DA	Sidechain
1	D	17	DC	Sidechain
1	D	19	DG	Sidechain
1	D	8	DA	Sidechain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the 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 within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	385	0	216	20	1
1	D	385	0	216	23	0
2	A	619	0	623	54	0
2	B	619	0	623	44	1
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	32	0	0	3	0
4	B	41	0	0	7	0
4	C	23	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	13	0	0	6	0
All	All	2121	0	1678	128	2

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 38.

All (128) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:484:ILE:O	2:B:485:ASP:HB3	1.42	1.16
1:D:6:DA:H2''	1:D:7:DC:H5'	1.28	1.12
1:D:8:DA:H2''	1:D:9:DT:O5'	1.62	1.00
1:D:16:DT:O4	4:D:32:HOH:O	1.78	0.99
2:B:443:CYS:O	2:B:444:SER:HB2	1.63	0.95
1:D:16:DT:H73	4:D:32:HOH:O	1.67	0.94
2:B:471:GLN:HA	4:B:13:HOH:O	1.67	0.94
2:A:486:LYS:HG2	2:A:486:LYS:O	1.69	0.90
2:A:475:LEU:HD12	2:B:488:ARG:NH2	1.87	0.90
2:B:484:ILE:O	2:B:485:ASP:CB	2.20	0.90
2:A:511:LYS:O	2:A:512:THR:HB	1.75	0.87
1:D:6:DA:C2'	1:D:7:DC:H5'	2.05	0.86
1:D:16:DT:O5'	1:D:16:DT:H2'	1.76	0.83
2:A:456:THR:CG2	2:A:457:CYS:N	2.41	0.83
1:C:7:DC:H2''	1:C:8:DA:O5'	1.79	0.82
1:C:5:DA:OP2	4:C:31:HOH:O	1.98	0.81
1:D:18:DT:O5'	1:D:18:DT:H2'	1.81	0.80
1:C:5:DA:N6	1:D:16:DT:O4	2.15	0.80
1:D:16:DT:C7	4:D:32:HOH:O	2.29	0.78
1:C:18:DT:H2''	1:C:19:DG:H5''	1.67	0.76
2:B:484:ILE:HD12	4:B:55:HOH:O	1.84	0.76
1:D:10:DC:H2'	1:D:11:DG:C8	2.21	0.76
2:A:456:THR:CG2	2:A:457:CYS:H	2.02	0.72
2:A:456:THR:HG22	2:A:457:CYS:N	2.03	0.72
2:A:465:LYS:HZ2	2:A:469:GLU:CD	1.97	0.72
2:A:456:THR:HG23	2:A:457:CYS:H	1.53	0.72
2:A:511:LYS:O	2:A:512:THR:CB	2.38	0.72
1:C:14:DG:O5'	1:C:14:DG:H2'	1.91	0.69
1:C:10:DC:H2''	1:C:11:DG:H5'	1.76	0.68
2:A:437:ALA:O	2:A:438:ARG:HB2	1.94	0.67
2:A:446:GLU:O	4:A:105:HOH:O	2.11	0.67
2:A:452:TYR:HB3	2:A:506:ASN:O	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:439:PRO:O	2:A:439:PRO:HG2	1.96	0.66
2:A:475:LEU:CD1	2:B:488:ARG:NH2	2.58	0.66
2:B:437:ALA:HB1	4:B:3:HOH:O	1.95	0.66
2:B:501:LEU:O	2:B:503:ALA:N	2.29	0.66
1:D:11:DG:O5'	1:D:11:DG:H2'	1.96	0.65
2:B:437:ALA:CB	4:B:3:HOH:O	2.44	0.64
2:B:481:ASP:CG	2:B:481:ASP:O	2.41	0.64
2:B:509:ALA:C	2:B:511:LYS:N	2.52	0.63
2:B:435:LYS:HD2	2:B:446:GLU:OE2	1.99	0.63
1:D:9:DT:H2''	1:D:10:DC:OP2	1.97	0.63
2:A:498:ARG:HH21	2:A:498:ARG:HB3	1.64	0.61
1:C:18:DT:H2''	1:C:19:DG:C5'	2.31	0.59
1:D:16:DT:O5'	1:D:16:DT:C2'	2.49	0.59
1:D:14:DG:O6	2:A:466:ARG:NH1	2.29	0.59
2:B:448:SER:OG	2:B:451:HIS:HE1	1.86	0.59
2:A:437:ALA:O	2:A:438:ARG:CB	2.51	0.58
2:B:462:VAL:O	2:B:466:ARG:HB2	2.04	0.58
2:A:496:ARG:O	2:A:497:TYR:C	2.46	0.58
2:B:483:ILE:HD13	2:B:483:ILE:N	2.20	0.57
1:C:10:DC:H2''	1:C:11:DG:C5'	2.34	0.57
2:B:484:ILE:HB	4:B:55:HOH:O	2.04	0.57
2:B:443:CYS:O	2:B:444:SER:CB	2.38	0.57
1:C:2:DC:H1'	2:A:511:LYS:HD3	1.86	0.57
2:A:465:LYS:NZ	2:A:469:GLU:CD	2.62	0.57
2:A:497:TYR:O	2:A:500:CYS:HB2	2.05	0.57
2:A:488:ARG:HD2	2:B:477:ALA:HA	1.86	0.56
1:C:2:DC:H2'	1:C:3:DA:C8	2.40	0.56
2:A:488:ARG:NH2	2:B:475:LEU:HD13	2.20	0.56
1:C:7:DC:C2'	1:C:8:DA:O5'	2.51	0.56
2:A:465:LYS:O	2:A:469:GLU:HG3	2.06	0.56
2:A:468:VAL:HG23	2:A:469:GLU:N	2.19	0.56
4:D:26:HOH:O	2:B:511:LYS:HE3	2.06	0.55
2:B:501:LEU:O	2:B:502:GLN:C	2.48	0.55
2:A:439:PRO:O	2:A:439:PRO:CG	2.55	0.55
1:D:16:DT:C4	4:D:32:HOH:O	2.48	0.54
2:A:447:ALA:HB1	2:A:455:LEU:CD1	2.38	0.53
4:D:24:HOH:O	2:A:451:HIS:HE1	1.90	0.53
1:D:9:DT:C2'	1:D:10:DC:OP2	2.53	0.52
2:B:472:HIS:N	4:B:13:HOH:O	2.41	0.52
1:D:11:DG:O5'	1:D:11:DG:C2'	2.58	0.52
2:A:501:LEU:O	2:A:502:GLN:C	2.52	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2:DC:P	4:C:28:HOH:O	2.68	0.51
1:C:2:DC:H2''	2:A:511:LYS:HE3	1.92	0.51
1:D:8:DA:C2'	1:D:9:DT:O5'	2.47	0.51
2:A:468:VAL:HG23	2:A:469:GLU:H	1.75	0.51
2:B:468:VAL:HG23	2:B:469:GLU:N	2.26	0.50
2:A:475:LEU:HD12	2:B:488:ARG:HH21	1.75	0.49
2:A:456:THR:HG22	2:A:457:CYS:O	2.12	0.49
2:A:498:ARG:NH2	4:A:72:HOH:O	2.28	0.49
2:B:511:LYS:NZ	2:B:511:LYS:HB2	2.27	0.49
2:A:485:ASP:O	2:A:486:LYS:C	2.53	0.49
1:D:18:DT:O5'	1:D:18:DT:C2'	2.52	0.49
2:A:485:ASP:O	2:A:489:ARG:HB3	2.12	0.48
2:A:465:LYS:NZ	2:A:469:GLU:OE2	2.47	0.48
2:A:481:ASP:OD1	2:A:481:ASP:C	2.56	0.48
1:C:14:DG:O5'	1:C:14:DG:C2'	2.49	0.48
1:C:18:DT:C2'	1:C:19:DG:H5''	2.42	0.48
2:A:438:ARG:N	2:A:439:PRO:CD	2.77	0.48
2:A:488:ARG:HH22	2:B:475:LEU:HD13	1.79	0.48
2:B:464:PHE:O	2:B:465:LYS:C	2.56	0.48
2:B:479:ARG:NH1	4:B:76:HOH:O	2.48	0.47
2:A:447:ALA:CB	2:A:455:LEU:HD11	2.45	0.46
2:A:462:VAL:O	2:A:466:ARG:HB2	2.15	0.46
2:A:498:ARG:NH1	4:A:72:HOH:O	2.43	0.46
2:B:487:ILE:H	2:B:487:ILE:HG12	1.18	0.46
2:A:447:ALA:HB1	2:A:455:LEU:HD11	1.97	0.46
2:A:447:ALA:HA	2:A:456:THR:O	2.16	0.46
1:D:13:DT:H2''	1:D:14:DG:O5'	2.16	0.45
2:A:475:LEU:HB3	2:B:488:ARG:HD3	1.98	0.45
2:B:483:ILE:HG22	2:B:484:ILE:H	1.82	0.45
2:B:465:LYS:NZ	2:B:469:GLU:CD	2.75	0.45
2:B:484:ILE:HD11	2:B:496:ARG:HG3	2.00	0.44
2:B:501:LEU:HA	2:B:501:LEU:HD12	1.58	0.44
1:C:11:DG:H1	1:D:10:DC:N4	2.15	0.43
1:C:6:DA:H2'	1:C:6:DA:H5'	1.22	0.43
2:B:483:ILE:HG22	2:B:484:ILE:N	2.33	0.43
2:B:465:LYS:O	2:B:469:GLU:HG3	2.18	0.43
2:B:465:LYS:HZ1	2:B:469:GLU:CD	2.27	0.43
2:B:498:ARG:HH21	2:B:498:ARG:HD2	1.10	0.43
2:A:459:SER:HB2	2:A:496:ARG:NH1	2.33	0.43
2:B:488:ARG:HH21	2:B:488:ARG:HD3	1.65	0.42
2:A:510:ARG:HH21	2:A:510:ARG:HD2	1.68	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:438:ARG:H	2:A:455:LEU:HD21	1.85	0.41
1:C:15:DT:H2''	1:C:16:DT:H5'	2.01	0.41
2:A:507:LEU:HD12	2:A:507:LEU:HA	1.87	0.41
2:B:447:ALA:CB	2:B:455:LEU:HD23	2.51	0.41
2:B:507:LEU:HD23	2:B:507:LEU:HA	1.85	0.41
2:B:509:ALA:C	2:B:511:LYS:H	2.09	0.41
1:C:12:DA:N3	4:C:34:HOH:O	2.37	0.41
1:D:17:DC:H2'	1:D:18:DT:C6	2.56	0.41
1:C:19:DG:H5'	1:C:19:DG:H2'	1.20	0.40
2:A:448:SER:N	2:A:456:THR:O	2.49	0.40
2:B:511:LYS:HB3	2:B:512:THR:H	1.21	0.40
1:D:13:DT:H4'	2:A:490:LYS:HE3	2.03	0.40
2:A:438:ARG:HA	2:A:438:ARG:HD3	1.84	0.40
2:A:472:HIS:HB3	2:A:473:ASN:H	1.68	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1:DC:N3	1:C:19:DG:N7[2_455]	1.98	0.22
2:B:488:ARG:NH1	2:B:498:ARG:NE[4_555]	2.02	0.18

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	
2	A	78/92 (85%)	63 (81%)	10 (13%)	5 (6%)	1	1
2	B	78/92 (85%)	66 (85%)	9 (12%)	3 (4%)	2	3
All	All	156/184 (85%)	129 (83%)	19 (12%)	8 (5%)	1	2

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	438	ARG
2	B	512	THR
2	A	439	PRO
2	B	501	LEU
2	A	497	TYR
2	A	512	THR
2	B	485	ASP
2	A	498	ARG

5.3.2 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 X-ray entries followed by that with respect to entries of similar resolution.

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	
2	A	67/75 (89%)	49 (73%)	18 (27%)	0	1
2	B	67/75 (89%)	49 (73%)	18 (27%)	0	1
All	All	134/150 (89%)	98 (73%)	36 (27%)	0	1

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	434	MET
2	A	438	ARG
2	A	439	PRO
2	A	454	VAL
2	A	456	THR
2	A	459	SER
2	A	461	LYS
2	A	462	VAL
2	A	469	GLU
2	A	471	GLN
2	A	473	ASN
2	A	475	LEU
2	A	479	ARG
2	A	482	CYS
2	A	486	LYS
2	A	498	ARG

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Mol	Chain	Res	Type
2	A	510	ARG
2	A	513	LYS
2	B	434	MET
2	B	435	LYS
2	B	438	ARG
2	B	439	PRO
2	B	441	LEU
2	B	450	CYS
2	B	455	LEU
2	B	474	TYR
2	B	479	ARG
2	B	487	ILE
2	B	488	ARG
2	B	490	LYS
2	B	491	ASN
2	B	493	PRO
2	B	499	LYS
2	B	501	LEU
2	B	510	ARG
2	B	511	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	473	ASN
2	A	491	ASN
2	A	506	ASN
2	B	451	HIS
2	B	480	ASN
2	B	491	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	C	19/19 (100%)	0.90	3 (15%) 5 4	22, 27, 36, 44	0
1	D	19/19 (100%)	0.60	1 (5%) 32 28	22, 27, 33, 36	0
2	A	80/92 (86%)	0.74	8 (10%) 12 10	14, 26, 52, 57	0
2	B	80/92 (86%)	0.41	3 (3%) 44 39	13, 24, 45, 50	0
All	All	198/222 (89%)	0.61	15 (7%) 20 17	13, 26, 48, 57	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	436	PRO	3.9
2	A	504	GLY	2.8
1	C	1	DC	2.7
1	D	17	DC	2.7
2	B	512	THR	2.7
2	A	434	MET	2.6
2	A	454	VAL	2.6
1	C	2	DC	2.6
2	A	512	THR	2.5
2	A	439	PRO	2.3
2	B	436	PRO	2.3
2	A	511	LYS	2.3
1	C	7	DC	2.1
2	B	513	LYS	2.0
2	A	449	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	ZN	A	526	1/1	0.64	0.07	26,26,26,26	0
3	ZN	B	527	1/1	0.71	0.07	17,17,17,17	0
3	ZN	A	527	1/1	0.83	0.06	16,16,16,16	0
3	ZN	B	526	1/1	0.92	0.07	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.