



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:21 PM GMT

PDB ID : 4RI9  
Title : FAN1 Nuclease bound to 5' phosphorylated p(dT)/3'(dT-dT-dT-dT-dT-dT-dT-dT) double flap DNA  
Authors : Pavletich, N.P.; Wang, R.  
Deposited on : 2014-10-05  
Resolution : 2.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

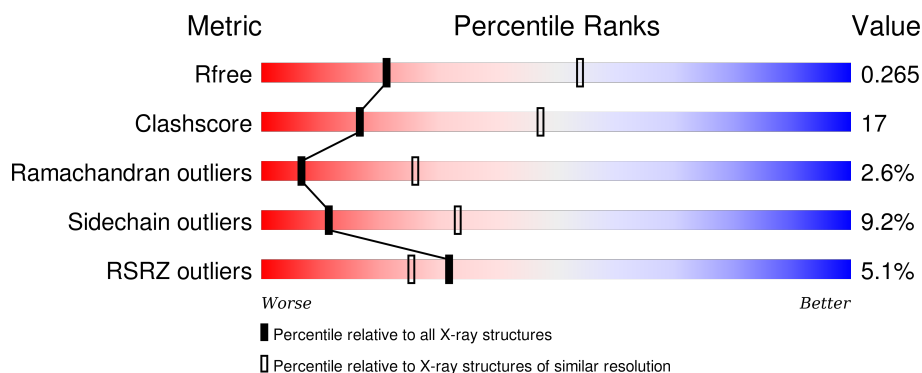
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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.



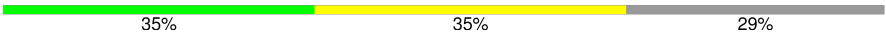
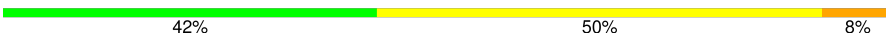



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	A	652	<div> <div>4%</div> <div>58%</div> <div>31%</div> <div>5%</div> <div>6%</div> </div>
1	B	652	<div> <div>6%</div> <div>60%</div> <div>30%</div> <div>•</div> <div>6%</div> </div>
2	S	11	<div> <div>9%</div> <div>91%</div> </div>
2	W	11	<div> <div>18%</div> <div>82%</div> </div>
3	U	17	<div> <div>53%</div> <div>18%</div> <div>29%</div> </div>

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Mol	Chain	Length	Quality of chain
3	Z	17	
4	V	12	
4	X	12	
5	T	14	
5	Y	14	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	BA	B	1101	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11606 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 protein called Fanconi-associated nuclease 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	615	Total	C	N	O	S	0	0	0
			4942	3145	880	890	27			
1	B	615	Total	C	N	O	S	0	0	0
			4942	3145	880	890	27			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	357	SER	-	EXPRESSION TAG	UNP Q9Y2M0
A	358	GLY	-	EXPRESSION TAG	UNP Q9Y2M0
A	359	ALA	-	EXPRESSION TAG	UNP Q9Y2M0
A	360	HIS	-	EXPRESSION TAG	UNP Q9Y2M0
A	361	MET	-	EXPRESSION TAG	UNP Q9Y2M0
A	362	THR	-	EXPRESSION TAG	UNP Q9Y2M0
A	363	ARG	-	EXPRESSION TAG	UNP Q9Y2M0
A	364	ASN	-	EXPRESSION TAG	UNP Q9Y2M0
A	365	GLY	-	EXPRESSION TAG	UNP Q9Y2M0
A	366	PRO	-	EXPRESSION TAG	UNP Q9Y2M0
A	367	GLY	-	EXPRESSION TAG	UNP Q9Y2M0
A	368	GLN	-	EXPRESSION TAG	UNP Q9Y2M0
A	369	THR	-	EXPRESSION TAG	UNP Q9Y2M0
A	487	ALA	VAL	ENGINEERED MUTATION	UNP Q9Y2M0
A	?	-	CYS	DELETION	UNP Q9Y2M0
A	?	-	THR	DELETION	UNP Q9Y2M0
A	?	-	TRP	DELETION	UNP Q9Y2M0
A	?	-	GLY	DELETION	UNP Q9Y2M0
A	?	-	LYS	DELETION	UNP Q9Y2M0
A	?	-	ASN	DELETION	UNP Q9Y2M0
A	?	-	LYS	DELETION	UNP Q9Y2M0
A	?	-	PRO	DELETION	UNP Q9Y2M0
A	?	-	GLY	DELETION	UNP Q9Y2M0
B	357	SER	-	EXPRESSION TAG	UNP Q9Y2M0
B	358	GLY	-	EXPRESSION TAG	UNP Q9Y2M0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	359	ALA	-	EXPRESSION TAG	UNP Q9Y2M0
B	360	HIS	-	EXPRESSION TAG	UNP Q9Y2M0
B	361	MET	-	EXPRESSION TAG	UNP Q9Y2M0
B	362	THR	-	EXPRESSION TAG	UNP Q9Y2M0
B	363	ARG	-	EXPRESSION TAG	UNP Q9Y2M0
B	364	ASN	-	EXPRESSION TAG	UNP Q9Y2M0
B	365	GLY	-	EXPRESSION TAG	UNP Q9Y2M0
B	366	PRO	-	EXPRESSION TAG	UNP Q9Y2M0
B	367	GLY	-	EXPRESSION TAG	UNP Q9Y2M0
B	368	GLN	-	EXPRESSION TAG	UNP Q9Y2M0
B	369	THR	-	EXPRESSION TAG	UNP Q9Y2M0
B	487	ALA	VAL	ENGINEERED MUTATION	UNP Q9Y2M0
B	?	-	CYS	DELETION	UNP Q9Y2M0
B	?	-	THR	DELETION	UNP Q9Y2M0
B	?	-	TRP	DELETION	UNP Q9Y2M0
B	?	-	GLY	DELETION	UNP Q9Y2M0
B	?	-	LYS	DELETION	UNP Q9Y2M0
B	?	-	ASN	DELETION	UNP Q9Y2M0
B	?	-	LYS	DELETION	UNP Q9Y2M0
B	?	-	PRO	DELETION	UNP Q9Y2M0
B	?	-	GLY	DELETION	UNP Q9Y2M0

- Molecule 2 is a DNA chain called DNA (5'-D(P\*TP\*AP\*GP\*CP\*CP\*AP\*CP\*GP\*CP\*CP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	S	11	Total	C	N	O	P	0	0	0
			222	105	39	67	11			
2	W	11	Total	C	N	O	P	0	0	0
			222	105	39	67	11			

- Molecule 3 is a DNA chain called DNA (5'-D(P\*AP\*GP\*AP\*CP\*TP\*CP\*CP\*TP\*CP\*TP\*TP\*TP\*TP\*TP\*TP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	U	12	Total	C	N	O	P	0	0	0
			240	116	37	75	12			
3	Z	12	Total	C	N	O	P	0	0	0
			240	116	37	75	12			

- Molecule 4 is a DNA chain called DNA (5'-D(P\*GP\*CP\*TP\*GP\*AP\*GP\*GP\*AP\*GP\*TP\*P\*CP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	V	12	Total	C	N	O	P	3	0	0
			250	118	47	73	12			
4	X	12	Total	C	N	O	P	3	0	0
			250	118	47	73	12			

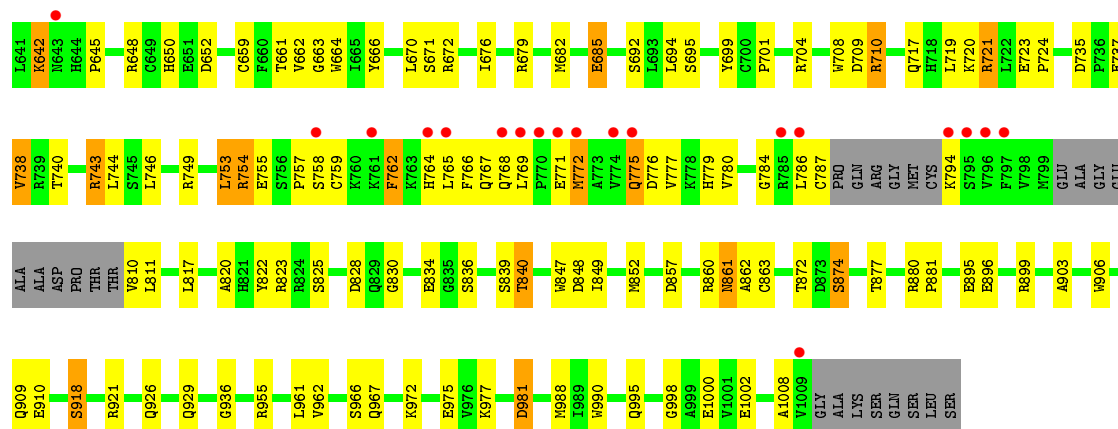
- Molecule 5 is a DNA chain called DNA (5'-D(\*TP\*TP\*TP\*TP\*TP\*TP\*GP\*AP\*GP\*GP\*CP\*GP\*TP\*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	Y	7	Total	C	N	O	P	0	0	0
			148	69	30	42	7			
5	T	7	Total	C	N	O	P	0	0	0
			148	69	30	42	7			

- Molecule 6 is BARIUM ION (three-letter code: BA) (formula: Ba).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ba	0	0
			1	1		
6	A	1	Total	Ba	0	0
			1	1		

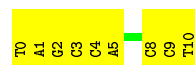




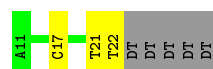
- Molecule 2: DNA (5'-D(P\*TP\*AP\*GP\*CP\*CP\*AP\*CP\*GP\*CP\*CP\*T)-3')



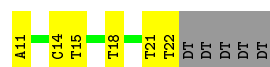
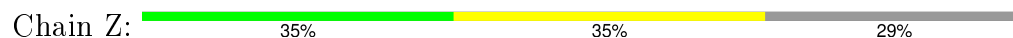
- Molecule 2: DNA (5'-D(P\*TP\*AP\*GP\*CP\*CP\*AP\*CP\*GP\*CP\*CP\*T)-3')



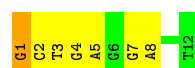
- Molecule 3: DNA (5'-D(P<sup>\*</sup>AP<sup>\*</sup>GP<sup>\*</sup>AP<sup>\*</sup>CP<sup>\*</sup>TP<sup>\*</sup>CP<sup>\*</sup>CP<sup>\*</sup>TP<sup>\*</sup>CP<sup>\*</sup>TP<sup>\*</sup>TP<sup>\*</sup>TP<sup>\*</sup>TP<sup>\*</sup>TP<sup>\*</sup>T  
P<sup>\*</sup>T)-3')



- Molecule 3: DNA (5'-D(P<sup>\*</sup>AP<sup>\*</sup>GP<sup>\*</sup>AP<sup>\*</sup>CP<sup>\*</sup>TP<sup>\*</sup>CP<sup>\*</sup>CP<sup>\*</sup>TP<sup>\*</sup>CP<sup>\*</sup>TP<sup>\*</sup>TP<sup>\*</sup>TP<sup>\*</sup>TP<sup>\*</sup>TP<sup>\*</sup>T  
P<sup>\*</sup>T)-3')



- Molecule 4: DNA (5'-D(P\*GP\*CP\*TP\*GP\*AP\*GP\*GP\*AP\*GP\*TP\*CP\*T)-3')

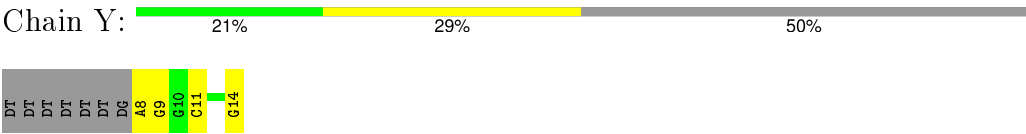


- Molecule 4: DNA (5'-D(P\*GP\*CP\*TP\*GP\*AP\*GP\*GP\*AP\*GP\*TP\*CP\*T)-3')

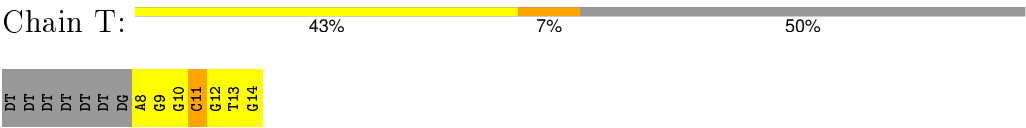




• Molecule 5: DNA (5'-D(\*TP\*TP\*TP\*TP\*TP\*TP\*GP\*AP\*GP\*GP\*CP\*GP\*TP\*G)-3')



• Molecule 5: DNA (5'-D(\*TP\*TP\*TP\*TP\*TP\*TP\*GP\*AP\*GP\*GP\*CP\*GP\*TP\*G)-3')



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.86Å 211.57Å 69.27Å 90.00° 99.15° 90.00°	Depositor
Resolution (Å)	50.00 – 2.90 68.39 – 2.90	Depositor EDS
% Data completeness (in resolution range)	91.6 (50.00-2.90) 91.6 (68.39-2.90)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.69 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, $R_{free}$	0.227 , 0.266 0.227 , 0.265	Depositor DCC
$R_{free}$ test set	1626 reflections (4.41%)	DCC
Wilson B-factor (Å <sup>2</sup> )	62.0	Xtriage
Anisotropy	0.387	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 26.8	EDS
Estimated twinning fraction	0.078 for l,-k,h	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 40067 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	11606	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: BA

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	A	0.51	0/5044	0.71	1/6818 (0.0%)
1	B	0.49	0/5044	0.71	1/6818 (0.0%)
2	S	0.71	1/247 (0.4%)	0.87	1/376 (0.3%)
2	W	0.75	1/247 (0.4%)	0.87	0/376
3	U	0.50	0/266	0.82	1/407 (0.2%)
3	Z	0.41	0/266	0.82	0/407
4	V	0.83	1/280 (0.4%)	0.97	2/431 (0.5%)
4	X	2.18	1/280 (0.4%)	1.39	4/431 (0.9%)
5	T	0.58	0/166	0.93	2/255 (0.8%)
5	Y	0.56	0/166	0.83	1/255 (0.4%)
All	All	0.61	4/12006 (0.0%)	0.76	13/16574 (0.1%)

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	A	0	2
1	B	0	2
All	All	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	X	1	DG	P-O5'	35.95	1.95	1.59
4	V	1	DG	P-O5'	12.60	1.72	1.59
2	W	0	DT	OP3-P	-10.29	1.48	1.61
2	S	0	DT	OP3-P	-9.46	1.49	1.61

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	X	1	DG	P-O5'-C5'	-20.55	88.01	120.90
4	X	1	DG	O5'-P-OP1	10.82	123.68	110.70
4	V	1	DG	P-O5'-C5'	-10.05	104.81	120.90
4	X	1	DG	O5'-P-OP2	-6.59	99.77	105.70
5	T	11	DC	C1'-O4'-C4'	-6.30	103.80	110.10
4	X	1	DG	OP1-P-OP2	-6.28	110.19	119.60
5	Y	11	DC	C1'-O4'-C4'	-6.27	103.83	110.10
4	V	1	DG	OP1-P-OP2	-6.13	110.40	119.60
5	T	10	DG	C1'-O4'-C4'	-5.83	104.27	110.10
1	B	558	MET	N-CA-C	-5.53	96.07	111.00
1	A	558	MET	N-CA-C	-5.46	96.25	111.00
2	S	6	DC	C1'-O4'-C4'	-5.33	104.77	110.10
3	U	17	DC	C1'-O4'-C4'	-5.10	105.00	110.10

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	567	GLY	Peptide
1	A	861	ASN	Peptide
1	B	567	GLY	Peptide
1	B	861	ASN	Peptide

## 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	A	4942	0	4978	157	0
1	B	4942	0	4978	156	0
2	S	222	0	124	6	0
2	W	222	0	124	11	1
3	U	240	0	138	4	0
3	Z	240	0	138	11	1
4	V	250	0	136	9	0
4	X	250	0	136	12	0
5	T	148	0	79	8	0
5	Y	148	0	79	5	0
6	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	1	0	0	0	0
All	All	11606	0	10910	367	1

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 17.

All (367) 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)
1:A:679:ARG:NH1	5:Y:14:DG:OP2	1.87	1.05
1:B:836:SER:O	1:B:840:THR:HG23	1.60	1.01
1:B:565:CYS:HB3	1:B:609:THR:HG23	1.49	0.94
1:A:836:SER:O	1:A:840:THR:HG23	1.68	0.94
2:W:1:DA:H2''	2:W:2:DG:OP2	1.69	0.93
3:Z:21:DT:H4'	3:Z:22:DT:H5'	1.51	0.92
3:Z:21:DT:H2''	3:Z:22:DT:OP2	1.72	0.90
1:B:591:ARG:HD3	1:B:860:ARG:O	1.72	0.89
5:T:8:DA:H2''	5:T:9:DG:OP2	1.73	0.86
1:B:370:THR:HG22	1:B:377:ARG:HD3	1.59	0.85
2:W:5:DA:N6	5:T:13:DT:O4	2.08	0.84
1:A:370:THR:HG22	1:A:377:ARG:HD3	1.58	0.83
1:B:738:VAL:O	1:B:743:ARG:NH1	2.12	0.83
1:B:565:CYS:CB	1:B:609:THR:HG23	2.08	0.83
1:A:389:ASN:HB3	1:A:392:ASP:HB2	1.61	0.83
1:A:591:ARG:HD3	1:A:860:ARG:O	1.83	0.79
1:B:918:SER:HB2	1:B:921:ARG:HB2	1.65	0.78
1:A:828:ASP:OD2	1:A:966:SER:HB2	1.84	0.78
1:B:640:ARG:HH11	1:B:640:ARG:HG2	1.50	0.77
1:A:723:GLU:HB3	1:A:724:PRO:HD3	1.67	0.76
1:B:389:ASN:HB3	1:B:392:ASP:HB2	1.68	0.76
1:B:723:GLU:HB3	1:B:724:PRO:HD3	1.67	0.75
3:U:22:DT:H2'	3:U:22:DT:O2	1.86	0.75
4:X:8:DA:H2''	4:X:9:DG:OP2	1.85	0.75
1:A:385:THR:HG23	1:A:862:ALA:HB3	1.68	0.75
3:U:21:DT:H4'	3:U:22:DT:H5'	1.67	0.74
4:V:1:DG:H2''	4:V:2:DC:H5'	1.67	0.74
1:A:431:MET:HG3	1:A:533:GLN:HB3	1.70	0.74
1:B:640:ARG:CG	1:B:640:ARG:HH11	2.02	0.73
1:A:421:LEU:HB3	1:A:537:ILE:HD13	1.70	0.73
1:B:385:THR:HG23	1:B:862:ALA:HB3	1.68	0.73
1:B:672:ARG:O	1:B:676:ILE:HG13	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:V:4:DG:H2"	4:V:5:DA:OP2	1.89	0.72
1:A:775:GLN:HE22	1:A:996:LYS:HG3	1.54	0.72
1:A:520:GLY:HA2	1:A:523:ILE:HD12	1.72	0.72
1:B:757:PRO:C	1:B:759:CYS:H	1.93	0.72
2:W:2:DG:H2"	2:W:3:DC:C5'	2.20	0.71
1:A:757:PRO:C	1:A:759:CYS:H	1.93	0.71
1:A:640:ARG:HG2	1:A:640:ARG:HH11	1.56	0.71
1:B:794:LYS:HA	2:W:5:DA:H5"	1.72	0.71
1:B:520:GLY:HA2	1:B:523:ILE:HD12	1.73	0.71
1:A:982:ARG:NH2	5:Y:9:DG:H2"	2.06	0.70
1:B:565:CYS:HB3	1:B:609:THR:CG2	2.21	0.70
1:A:458:THR:HG22	1:A:460:SER:H	1.57	0.70
1:A:754:ARG:HG3	1:A:755:GLU:N	2.07	0.69
1:B:872:THR:HG22	1:B:874:SER:H	1.57	0.69
1:B:568:GLN:HE22	1:B:610:HIS:HD2	1.41	0.69
1:A:738:VAL:O	1:A:743:ARG:NH1	2.24	0.69
1:A:444:THR:N	1:A:445:PRO:HD2	2.07	0.69
4:X:4:DG:H2"	4:X:5:DA:OP2	1.92	0.69
1:A:639:ASN:HA	1:A:642:LYS:HG3	1.75	0.69
1:B:719:LEU:O	1:B:721:ARG:N	2.27	0.68
1:A:565:CYS:HB3	1:A:609:THR:HG23	1.75	0.68
4:V:2:DC:H2'	4:V:3:DT:C6	2.28	0.67
4:X:6:DG:H2"	4:X:7:DG:OP2	1.94	0.67
1:A:744:LEU:HA	1:A:772:MET:HE1	1.76	0.67
1:B:444:THR:N	1:B:445:PRO:HD2	2.09	0.67
1:B:568:GLN:NE2	1:B:610:HIS:HD2	1.92	0.66
1:A:694:LEU:O	1:A:704:ARG:NH2	2.29	0.65
1:A:640:ARG:CG	1:A:640:ARG:HH11	2.09	0.65
1:B:391:ASP:CG	1:B:591:ARG:HH12	2.01	0.64
1:B:639:ASN:HA	1:B:642:LYS:HG3	1.80	0.64
1:A:872:THR:HG22	1:A:874:SER:H	1.62	0.64
1:B:638:TRP:HB2	1:B:666:TYR:CD1	2.32	0.64
1:A:995:GLN:O	1:A:998:GLY:N	2.26	0.64
1:B:754:ARG:HG3	1:B:755:GLU:N	2.12	0.64
1:B:482:LYS:HG3	1:B:489:PRO:HB3	1.80	0.63
1:B:559:GLU:O	1:B:560:ASP:HB2	1.98	0.63
1:A:899:ARG:NH2	1:A:926:GLN:HE22	1.96	0.63
1:B:391:ASP:HB2	1:B:591:ARG:NH1	2.13	0.63
1:B:565:CYS:SG	1:B:609:THR:HG23	2.39	0.63
1:A:638:TRP:HB2	1:A:666:TYR:CD1	2.34	0.62
4:X:4:DG:H8	4:X:4:DG:OP1	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:774:VAL:HG13	1:A:992:ALA:HB2	1.81	0.62
4:X:7:DG:H2''	4:X:8:DA:OP2	1.99	0.62
1:B:426:LEU:HD13	1:B:537:ILE:HG22	1.80	0.62
1:A:975:GLU:OE2	1:A:977:LYS:NZ	2.27	0.62
1:B:459:GLU:O	1:B:462:LEU:HB3	2.00	0.62
1:B:458:THR:HG22	1:B:460:SER:H	1.64	0.61
2:S:7:DG:H2''	2:S:8:DC:OP2	1.99	0.61
1:B:421:LEU:HD21	1:B:434:LEU:HD21	1.81	0.61
1:A:559:GLU:O	1:A:560:ASP:HB2	2.00	0.61
1:A:672:ARG:O	1:A:676:ILE:HG13	2.00	0.61
1:B:568:GLN:NE2	1:B:610:HIS:CD2	2.69	0.61
1:B:762:PHE:HB2	1:B:764:HIS:HD2	1.66	0.61
1:A:421:LEU:HB3	1:A:537:ILE:CD1	2.31	0.61
1:B:679:ARG:NH1	5:T:14:DG:OP2	2.31	0.60
1:B:765:LEU:HD22	1:B:769:LEU:HD11	1.83	0.60
1:A:918:SER:HB2	1:A:921:ARG:HB2	1.84	0.60
1:A:426:LEU:HD22	1:A:542:ARG:HD2	1.84	0.60
1:B:584:PHE:CZ	1:B:863:CYS:HB3	2.37	0.60
3:Z:22:DT:O2	3:Z:22:DT:C2'	2.50	0.60
1:A:765:LEU:HD22	1:A:769:LEU:HD11	1.82	0.60
1:A:584:PHE:CZ	1:A:863:CYS:HB3	2.37	0.60
1:A:719:LEU:O	1:A:721:ARG:N	2.35	0.60
1:A:701:PRO:O	1:A:704:ARG:HG3	2.01	0.59
1:A:762:PHE:HB2	1:A:764:HIS:HD2	1.66	0.59
1:B:694:LEU:O	1:B:704:ARG:NH2	2.36	0.59
5:Y:8:DA:H2'	5:Y:8:DA:OP2	2.02	0.59
1:A:692:SER:O	1:A:695:SER:OG	2.20	0.59
1:B:387:LEU:HD11	1:B:404:VAL:HG11	1.84	0.59
3:U:22:DT:C2'	3:U:22:DT:O2	2.49	0.59
1:B:757:PRO:O	1:B:759:CYS:N	2.35	0.58
1:B:431:MET:HG3	1:B:533:GLN:HB3	1.85	0.58
1:B:744:LEU:HA	1:B:772:MET:HE3	1.84	0.58
2:W:2:DG:H2''	2:W:3:DC:H5'	1.84	0.58
1:A:895:GLU:O	1:A:899:ARG:HG3	2.03	0.58
1:A:482:LYS:HG3	1:A:489:PRO:HB3	1.86	0.58
1:A:786:LEU:HB2	1:A:1008:ALA:HB1	1.84	0.58
2:S:2:DG:H2''	2:S:3:DC:OP2	2.03	0.58
4:V:1:DG:H5'	5:T:14:DG:O3'	2.02	0.58
3:Z:22:DT:O2	3:Z:22:DT:H2'	2.02	0.58
1:A:558:MET:C	1:A:560:ASP:H	2.05	0.58
1:B:975:GLU:OE2	1:B:977:LYS:NZ	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:X:8:DA:H1'	4:X:9:DG:H5'	1.86	0.58
1:B:640:ARG:CG	1:B:640:ARG:NH1	2.67	0.57
1:A:558:MET:C	1:A:560:ASP:N	2.58	0.57
1:A:591:ARG:HG2	1:A:860:ARG:HA	1.87	0.57
2:W:2:DG:H2''	2:W:3:DC:O4'	2.05	0.56
1:A:659:CYS:HA	1:A:664:TRP:CD2	2.40	0.56
1:B:561:GLU:O	1:B:565:CYS:HB2	2.06	0.56
1:A:757:PRO:O	1:A:759:CYS:N	2.38	0.56
1:B:635:LYS:HG3	1:B:670:LEU:HD21	1.86	0.56
1:A:387:LEU:HD11	1:A:404:VAL:HG11	1.88	0.56
4:V:1:DG:H2''	4:V:2:DC:C5'	2.35	0.56
4:X:1:DG:H5'	5:Y:14:DG:O3'	2.05	0.56
3:U:21:DT:H4'	3:U:22:DT:C5'	2.34	0.56
1:B:492:GLN:HB2	1:B:495:GLN:HG2	1.88	0.56
1:B:899:ARG:NH2	1:B:926:GLN:HE22	2.04	0.56
1:A:566:GLY:HA3	1:A:613:SER:OG	2.07	0.55
1:A:635:LYS:HG3	1:A:670:LEU:HD21	1.88	0.55
1:B:465:LEU:HD23	1:B:501:LEU:HD23	1.89	0.55
1:B:682:MET:HB3	1:B:685:GLU:HG3	1.88	0.55
1:A:840:THR:CG2	1:A:918:SER:H	2.20	0.55
1:B:744:LEU:HD21	1:B:988:MET:HG3	1.89	0.55
1:B:577:VAL:O	1:B:581:ARG:NH2	2.40	0.54
2:S:1:DA:H2''	2:S:2:DG:OP2	2.07	0.54
1:B:822:TYR:OH	1:B:1002:GLU:OE2	2.23	0.54
1:B:487:ALA:O	1:B:489:PRO:HD3	2.08	0.54
1:B:421:LEU:HB3	1:B:537:ILE:HD13	1.89	0.54
1:A:481:ALA:O	1:A:487:ALA:HB3	2.07	0.54
1:A:822:TYR:OH	1:A:1002:GLU:OE2	2.21	0.53
1:B:481:ALA:O	1:B:487:ALA:HB3	2.08	0.53
1:B:828:ASP:OD2	1:B:966:SER:HB2	2.09	0.53
1:B:659:CYS:HA	1:B:664:TRP:CD2	2.43	0.53
1:B:558:MET:C	1:B:560:ASP:H	2.12	0.53
1:B:744:LEU:CD2	1:B:988:MET:HG3	2.39	0.53
1:A:487:ALA:O	1:A:489:PRO:HD3	2.08	0.53
1:A:383:LEU:O	1:A:387:LEU:HG	2.07	0.53
1:B:553:SER:HB2	1:B:559:GLU:HG2	1.90	0.53
1:B:560:ASP:HB3	1:B:563:ALA:H	1.73	0.53
1:A:836:SER:OG	1:A:921:ARG:HD2	2.08	0.53
2:S:8:DC:H2''	2:S:9:DC:O5'	2.08	0.53
1:B:573:THR:O	1:B:577:VAL:HG23	2.09	0.53
1:B:426:LEU:HD22	1:B:542:ARG:HD2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:830:GLY:HA2	1:A:962:VAL:O	2.09	0.53
1:A:942:GLY:HA3	1:A:997:LEU:HD21	1.91	0.52
1:A:477:LEU:HD12	1:A:493:LYS:HG3	1.91	0.52
1:B:723:GLU:CB	1:B:724:PRO:HD3	2.39	0.52
1:B:704:ARG:NH1	1:B:735:ASP:OD2	2.40	0.52
1:A:972:LYS:HG3	1:A:1000:GLU:HB3	1.90	0.52
1:B:565:CYS:CB	1:B:609:THR:CG2	2.83	0.52
1:B:895:GLU:O	1:B:899:ARG:HG3	2.09	0.52
1:A:568:GLN:NE2	1:A:610:HIS:HD2	2.07	0.52
1:B:766:PHE:O	1:B:768:GLN:N	2.42	0.52
1:B:786:LEU:HD12	1:B:787:CYS:H	1.74	0.51
1:A:644:HIS:CG	1:A:645:PRO:HD2	2.45	0.51
1:A:757:PRO:C	1:A:759:CYS:N	2.62	0.51
1:B:972:LYS:HG3	1:B:1000:GLU:HB3	1.91	0.51
1:A:481:ALA:HB1	1:A:489:PRO:HA	1.92	0.51
1:A:682:MET:HB3	1:A:685:GLU:HG3	1.92	0.51
1:B:977:LYS:NZ	1:B:981:ASP:OD2	2.43	0.51
1:B:391:ASP:HB2	1:B:591:ARG:HH12	1.75	0.51
1:A:709:ASP:HA	1:A:746:LEU:HD21	1.91	0.51
1:A:659:CYS:HA	1:A:664:TRP:CG	2.45	0.51
2:W:1:DA:O5'	2:W:1:DA:H8	1.93	0.50
1:A:860:ARG:O	1:A:861:ASN:HB3	2.11	0.50
1:B:477:LEU:HD12	1:B:493:LYS:HG3	1.93	0.50
1:B:391:ASP:CB	1:B:591:ARG:HH12	2.24	0.50
1:B:757:PRO:C	1:B:759:CYS:N	2.62	0.50
1:A:645:PRO:O	1:A:648:ARG:HG2	2.11	0.50
1:B:420:ARG:O	1:B:424:ARG:HG2	2.10	0.50
1:A:577:VAL:O	1:A:581:ARG:NH2	2.45	0.50
1:B:591:ARG:HG2	1:B:860:ARG:HA	1.93	0.50
1:B:645:PRO:O	1:B:648:ARG:HG2	2.12	0.50
1:B:558:MET:C	1:B:560:ASP:N	2.63	0.50
1:A:694:LEU:HD22	1:A:708:TRP:CE2	2.46	0.50
1:B:383:LEU:O	1:B:387:LEU:HG	2.12	0.50
1:A:766:PHE:O	1:A:768:GLN:N	2.44	0.50
1:A:723:GLU:CB	1:A:724:PRO:HD3	2.40	0.50
1:A:431:MET:HG3	1:A:533:GLN:CB	2.40	0.50
1:A:568:GLN:HA	1:A:568:GLN:NE2	2.27	0.50
1:B:557:SER:O	1:B:558:MET:HB2	2.12	0.49
1:B:481:ALA:HB1	1:B:489:PRO:HA	1.94	0.49
1:B:836:SER:OG	1:B:921:ARG:HD2	2.12	0.49
1:B:694:LEU:HD22	1:B:708:TRP:CE2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:560:ASP:HB3	1:A:563:ALA:H	1.77	0.49
1:B:370:THR:C	1:B:372:HIS:H	2.15	0.49
4:X:5:DA:H2''	4:X:6:DG:OP2	2.12	0.49
1:A:625:GLU:HA	1:A:625:GLU:OE1	2.12	0.49
1:B:840:THR:CG2	1:B:918:SER:H	2.26	0.49
1:A:640:ARG:CG	1:A:640:ARG:NH1	2.74	0.49
2:W:3:DC:H1'	2:W:4:DC:H5'	1.94	0.49
1:B:482:LYS:C	1:B:484:PHE:H	2.16	0.49
1:A:786:LEU:HD12	1:A:787:CYS:H	1.76	0.48
1:A:568:GLN:HE22	1:A:610:HIS:HD2	1.60	0.48
1:B:820:ALA:HA	1:B:823:ARG:NH1	2.28	0.48
1:B:737:GLU:OE1	1:B:737:GLU:HA	2.13	0.48
1:B:880:ARG:HB3	1:B:881:PRO:HD3	1.94	0.48
1:A:663:GLY:HA3	1:A:699:TYR:CZ	2.48	0.48
1:A:421:LEU:CB	1:A:537:ILE:CD1	2.90	0.48
1:B:903:ALA:HB2	1:B:929:GLN:NE2	2.28	0.48
2:W:1:DA:O5'	2:W:1:DA:C8	2.67	0.48
1:A:872:THR:HG22	1:A:873:ASP:N	2.29	0.48
1:A:820:ALA:HA	1:A:823:ARG:NH1	2.29	0.48
4:X:1:DG:H2'	4:X:2:DC:C6	2.49	0.48
1:A:710:ARG:HA	1:A:710:ARG:HD2	1.58	0.48
1:A:557:SER:O	1:A:558:MET:HB2	2.14	0.48
1:A:849:ILE:HG23	1:A:852:MET:CE	2.44	0.48
1:A:840:THR:HG21	1:A:918:SER:H	1.78	0.48
2:W:8:DC:H2''	2:W:9:DC:H6	1.78	0.48
1:B:977:LYS:HA	1:B:977:LYS:HD3	1.65	0.47
1:B:507:ARG:O	1:B:519:ILE:HG22	2.14	0.47
1:A:465:LEU:HD23	1:A:501:LEU:HD23	1.95	0.47
4:V:2:DC:H2'	4:V:3:DT:H6	1.74	0.47
1:B:496:LEU:O	1:B:500:PHE:HD2	1.96	0.47
1:B:571:LEU:HD22	1:B:576:LEU:HD12	1.96	0.47
1:B:556:ASP:C	1:B:558:MET:H	2.16	0.47
1:A:558:MET:O	1:A:560:ASP:N	2.47	0.47
1:B:849:ILE:HG23	1:B:852:MET:CE	2.45	0.47
1:A:507:ARG:O	1:A:519:ILE:HG22	2.15	0.47
1:B:692:SER:O	1:B:695:SER:OG	2.28	0.47
1:B:995:GLN:O	1:B:998:GLY:N	2.44	0.47
3:Z:21:DT:C4'	3:Z:22:DT:H5'	2.36	0.47
1:A:982:ARG:HH21	5:Y:9:DG:H2''	1.80	0.47
1:A:482:LYS:C	1:A:484:PHE:H	2.18	0.47
1:B:663:GLY:HA3	1:B:699:TYR:CZ	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:709:ASP:HA	1:B:746:LEU:HD21	1.97	0.47
2:S:5:DA:OP2	2:S:5:DA:H8	1.98	0.47
1:B:558:MET:HG2	1:B:561:GLU:OE2	2.14	0.47
1:A:565:CYS:CB	1:A:609:THR:HG23	2.44	0.47
1:A:876:PHE:CD2	1:A:883:LEU:HD13	2.49	0.47
1:A:977:LYS:NZ	1:A:981:ASP:OD2	2.48	0.47
4:X:4:DG:C2'	4:X:5:DA:OP2	2.59	0.46
1:A:848:ASP:OD1	1:A:848:ASP:N	2.49	0.46
1:A:977:LYS:HD3	1:A:977:LYS:HA	1.74	0.46
1:A:744:LEU:HD23	1:A:985:HIS:HB3	1.97	0.46
1:A:622:GLY:O	1:A:624:TRP:CD1	2.69	0.46
1:A:840:THR:HB	1:A:917:VAL:HA	1.98	0.46
1:A:706:ARG:O	1:A:710:ARG:HB2	2.16	0.46
5:T:11:DC:H2''	5:T:12:DG:O5'	2.15	0.46
1:B:857:ASP:OD1	1:B:860:ARG:NH1	2.48	0.46
1:A:749:ARG:O	1:A:753:LEU:HB2	2.16	0.46
1:A:739:ARG:HG3	1:A:953:HIS:CD2	2.51	0.46
3:Z:21:DT:H6	3:Z:21:DT:OP1	1.98	0.46
1:A:492:GLN:HB2	1:A:495:GLN:HG2	1.98	0.46
1:B:710:ARG:HD2	1:B:710:ARG:HA	1.60	0.46
1:B:822:TYR:HA	1:B:825:SER:HB2	1.98	0.45
1:B:786:LEU:HB2	1:B:1008:ALA:HB1	1.98	0.45
1:B:810:VAL:HG12	1:B:811:LEU:N	2.31	0.45
1:A:835:GLY:O	1:A:839:SER:HB3	2.16	0.45
1:B:601:ASP:OD2	1:B:650:HIS:NE2	2.48	0.45
1:A:424:ARG:HD3	3:Z:18:DT:H3'	1.99	0.45
1:B:494:GLN:OE1	1:B:494:GLN:HA	2.16	0.45
2:S:10:DT:OP2	2:S:10:DT:H3'	2.15	0.45
1:A:494:GLN:OE1	1:A:494:GLN:HA	2.17	0.45
1:B:830:GLY:HA2	1:B:962:VAL:O	2.16	0.45
1:A:745:SER:HB2	1:A:985:HIS:CE1	2.52	0.45
1:B:458:THR:HG23	1:B:459:GLU:CD	2.37	0.45
1:B:906:TRP:O	1:B:910:GLU:HB2	2.16	0.45
1:B:568:GLN:NE2	1:B:568:GLN:HA	2.26	0.45
1:A:458:THR:HG23	1:A:459:GLU:CD	2.38	0.45
1:B:426:LEU:CD1	1:B:537:ILE:HG22	2.47	0.45
1:B:749:ARG:O	1:B:753:LEU:HB2	2.17	0.45
1:B:565:CYS:O	1:B:610:HIS:HA	2.17	0.44
4:X:4:DG:OP1	4:X:4:DG:C8	2.67	0.44
1:A:973:LEU:HB2	1:A:1001:VAL:HG12	1.97	0.44
1:A:426:LEU:HD13	1:A:537:ILE:HG22	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:650:HIS:HB3	1:A:661:THR:HG21	1.99	0.44
1:A:700:CYS:N	1:A:701:PRO:HD3	2.32	0.44
1:B:425:LYS:HG2	1:B:426:LEU:N	2.32	0.44
1:A:733:LEU:O	1:A:743:ARG:NH2	2.50	0.44
4:V:7:DG:H2"	4:V:8:DA:C8	2.53	0.44
1:A:549:LEU:HD11	1:A:571:LEU:HD12	1.99	0.44
1:A:556:ASP:C	1:A:558:MET:H	2.17	0.44
1:B:721:ARG:HD2	1:B:721:ARG:HA	1.84	0.44
1:B:421:LEU:HB3	1:B:537:ILE:CD1	2.48	0.44
4:V:2:DC:H2"	4:V:3:DT:O5'	2.18	0.44
1:A:444:THR:N	1:A:445:PRO:CD	2.78	0.44
1:B:701:PRO:O	1:B:704:ARG:HG3	2.17	0.44
1:B:659:CYS:HA	1:B:664:TRP:CG	2.52	0.44
1:B:877:THR:HG22	1:B:880:ARG:NH2	2.32	0.44
1:A:421:LEU:HD21	1:A:434:LEU:HD21	1.99	0.43
1:A:370:THR:C	1:A:372:HIS:H	2.20	0.43
1:B:650:HIS:HB3	1:B:661:THR:HG21	2.00	0.43
1:B:777:VAL:O	1:B:779:HIS:HD2	2.01	0.43
1:A:810:VAL:HG12	1:A:811:LEU:N	2.32	0.43
1:A:880:ARG:HB3	1:A:881:PRO:HD3	1.99	0.43
1:A:558:MET:HG2	1:A:561:GLU:OE2	2.19	0.43
1:A:898:LEU:O	1:A:899:ARG:C	2.57	0.43
1:B:820:ALA:HA	1:B:823:ARG:CZ	2.49	0.43
1:A:506:GLN:CG	1:A:507:ARG:H	2.32	0.43
3:Z:14:DC:H2"	3:Z:15:DT:OP2	2.19	0.43
1:A:465:LEU:HD13	1:A:524:LEU:HB2	1.99	0.43
1:A:737:GLU:HA	1:A:737:GLU:OE1	2.19	0.43
1:A:496:LEU:O	1:A:500:PHE:HD2	2.01	0.43
1:A:877:THR:HG22	1:A:880:ARG:NH2	2.33	0.43
1:A:784:GLY:O	1:A:1008:ALA:HA	2.19	0.43
1:B:458:THR:CG2	1:B:459:GLU:N	2.81	0.43
1:B:877:THR:HG22	1:B:880:ARG:HH21	1.84	0.43
1:A:549:LEU:HD11	1:A:571:LEU:CD1	2.48	0.43
1:B:482:LYS:O	1:B:484:PHE:N	2.52	0.43
1:B:566:GLY:HA3	1:B:613:SER:OG	2.19	0.43
1:A:421:LEU:HB2	1:A:537:ILE:HD11	2.01	0.43
1:B:848:ASP:N	1:B:848:ASP:OD1	2.51	0.43
1:B:775:GLN:HG3	1:B:775:GLN:H	1.54	0.43
1:B:555:THR:HG21	1:B:861:ASN:ND2	2.34	0.42
1:A:391:ASP:CG	1:A:591:ARG:HH12	2.23	0.42
1:B:370:THR:C	1:B:372:HIS:N	2.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:465:LEU:O	1:A:468:VAL:N	2.52	0.42
1:A:903:ALA:HA	1:A:925:LEU:HD21	2.01	0.42
5:T:9:DG:H8	5:T:9:DG:OP2	2.02	0.42
1:B:444:THR:N	1:B:445:PRO:CD	2.79	0.42
1:A:899:ARG:HH21	1:A:926:GLN:HE22	1.65	0.42
1:B:428:TRP:CE2	1:B:462:LEU:HD13	2.55	0.42
1:B:784:GLY:HA3	1:B:817:LEU:HD21	2.01	0.42
1:A:426:LEU:O	1:A:427:SER:HB3	2.20	0.42
1:A:704:ARG:NH1	1:A:735:ASP:OD2	2.45	0.42
1:A:677:LEU:HD13	1:A:685:GLU:HB2	2.01	0.42
3:Z:14:DC:C2'	3:Z:15:DT:H72	2.50	0.42
1:A:786:LEU:CB	1:A:1008:ALA:HB1	2.49	0.42
1:B:441:LEU:HA	1:B:441:LEU:HD12	1.93	0.41
1:A:847:TRP:CE2	1:A:909:GLN:HG3	2.55	0.41
1:B:396:PHE:O	1:B:401:LYS:HE3	2.20	0.41
1:B:847:TRP:CE2	1:B:909:GLN:HG3	2.55	0.41
4:X:2:DC:H2'	4:X:3:DT:C6	2.55	0.41
1:B:637:ASP:O	1:B:640:ARG:HB2	2.20	0.41
1:A:425:LYS:HG2	1:A:426:LEU:N	2.34	0.41
1:B:717:GLN:HG3	5:T:11:DC:H5''	2.01	0.41
4:V:7:DG:H2''	4:V:8:DA:H8	1.84	0.41
1:A:908:GLU:HG2	1:B:896:GLU:OE1	2.20	0.41
1:A:822:TYR:HA	1:A:825:SER:HB2	2.01	0.41
1:B:506:GLN:CG	1:B:507:ARG:H	2.33	0.41
1:A:903:ALA:HB2	1:A:929:GLN:NE2	2.35	0.41
1:B:415:GLN:O	1:B:416:LYS:C	2.59	0.41
1:B:556:ASP:O	1:B:558:MET:N	2.33	0.41
1:B:906:TRP:NE1	1:B:910:GLU:HG3	2.35	0.41
1:B:721:ARG:HH21	1:B:723:GLU:CB	2.34	0.41
1:B:759:CYS:O	1:B:762:PHE:CD1	2.73	0.41
1:B:611:MET:O	1:B:615:ILE:HG13	2.20	0.41
1:B:524:LEU:O	1:B:528:LYS:HG3	2.20	0.41
1:A:577:VAL:HG11	3:Z:21:DT:C4	2.56	0.41
1:A:784:GLY:HA3	1:A:817:LEU:HD21	2.03	0.41
3:Z:14:DC:H2''	3:Z:15:DT:H72	2.03	0.41
1:A:965:ASN:HB3	1:A:968:SER:OG	2.21	0.41
1:B:423:GLN:HA	1:B:570:GLN:HG2	2.01	0.41
1:A:431:MET:HB2	1:A:533:GLN:HB2	2.03	0.41
1:A:458:THR:CG2	1:A:459:GLU:N	2.83	0.40
1:A:744:LEU:CD2	1:A:988:MET:HG3	2.51	0.40
5:T:8:DA:C2'	5:T:9:DG:OP2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:370:THR:CG2	1:B:377:ARG:HD3	2.40	0.40
1:A:721:ARG:HH21	1:A:723:GLU:CB	2.33	0.40
1:A:561:GLU:O	1:A:565:CYS:HB2	2.22	0.40
1:B:465:LEU:O	1:B:468:VAL:N	2.55	0.40
1:A:799:MET:HE1	1:A:820:ALA:CB	2.52	0.40
1:A:601:ASP:OD2	1:A:650:HIS:NE2	2.53	0.40
2:W:1:DA:C2'	2:W:2:DG:OP2	2.50	0.40

All (1) 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)
2:W:10:DT:O3'	3:Z:11:DA:P[2_645]	1.91	0.29

## 5.3 Torsion angles

### 5.3.1 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 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	
1	A	609/652 (93%)	530 (87%)	60 (10%)	19 (3%)	5	21
1	B	609/652 (93%)	525 (86%)	71 (12%)	13 (2%)	9	32
All	All	1218/1304 (93%)	1055 (87%)	131 (11%)	32 (3%)	7	26

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	427	SER
1	A	558	MET
1	A	560	ASP
1	A	720	LYS
1	A	767	GLN
1	B	427	SER

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Mol	Chain	Res	Type
1	B	558	MET
1	B	560	ASP
1	B	720	LYS
1	B	758	SER
1	B	767	GLN
1	A	483	THR
1	A	559	GLU
1	A	642	LYS
1	A	758	SER
1	A	937	GLY
1	B	483	THR
1	B	559	GLU
1	B	642	LYS
1	A	557	SER
1	B	557	SER
1	B	740	THR
1	A	459	GLU
1	A	637	ASP
1	B	772	MET
1	A	761	LYS
1	A	772	MET
1	A	895	GLU
1	A	740	THR
1	A	814	VAL
1	B	936	GLY
1	A	936	GLY

### 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	
1	A	538/564 (95%)	489 (91%)	49 (9%)	12	34
1	B	538/564 (95%)	488 (91%)	50 (9%)	11	32
All	All	1076/1128 (95%)	977 (91%)	99 (9%)	11	33

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	370	THR
1	A	389	ASN
1	A	390	GLU
1	A	399	GLN
1	A	405	THR
1	A	406	LYS
1	A	424	ARG
1	A	435	GLU
1	A	441	LEU
1	A	444	THR
1	A	458	THR
1	A	486	LEU
1	A	488	ASN
1	A	525	LYS
1	A	535	VAL
1	A	537	ILE
1	A	539	LYS
1	A	542	ARG
1	A	547	ARG
1	A	549	LEU
1	A	559	GLU
1	A	568	GLN
1	A	573	THR
1	A	586	SER
1	A	593	THR
1	A	633	CYS
1	A	652	ASP
1	A	662	VAL
1	A	671	SER
1	A	685	GLU
1	A	710	ARG
1	A	721	ARG
1	A	738	VAL
1	A	753	LEU
1	A	754	ARG
1	A	762	PHE
1	A	771	GLU
1	A	775	GLN
1	A	776	ASP
1	A	780	VAL
1	A	834	GLU
1	A	839	SER
1	A	840	THR

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Mol	Chain	Res	Type
1	A	874	SER
1	A	918	SER
1	A	955	ARG
1	A	961	LEU
1	A	967	GLN
1	A	981	ASP
1	B	370	THR
1	B	389	ASN
1	B	390	GLU
1	B	395	LEU
1	B	399	GLN
1	B	405	THR
1	B	406	LYS
1	B	424	ARG
1	B	435	GLU
1	B	441	LEU
1	B	444	THR
1	B	458	THR
1	B	486	LEU
1	B	488	ASN
1	B	525	LYS
1	B	535	VAL
1	B	537	ILE
1	B	539	LYS
1	B	542	ARG
1	B	549	LEU
1	B	559	GLU
1	B	568	GLN
1	B	573	THR
1	B	586	SER
1	B	633	CYS
1	B	652	ASP
1	B	662	VAL
1	B	671	SER
1	B	685	GLU
1	B	710	ARG
1	B	721	ARG
1	B	738	VAL
1	B	743	ARG
1	B	753	LEU
1	B	754	ARG
1	B	762	PHE

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Mol	Chain	Res	Type
1	B	771	GLU
1	B	775	GLN
1	B	776	ASP
1	B	780	VAL
1	B	834	GLU
1	B	839	SER
1	B	840	THR
1	B	874	SER
1	B	918	SER
1	B	955	ARG
1	B	961	LEU
1	B	967	GLN
1	B	981	ASP
1	B	990	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	488	ASN
1	A	568	GLN
1	A	610	HIS
1	A	623	ASN
1	A	764	HIS
1	A	775	GLN
1	A	832	HIS
1	A	926	GLN
1	A	953	HIS
1	B	488	ASN
1	B	568	GLN
1	B	610	HIS
1	B	623	ASN
1	B	764	HIS
1	B	832	HIS
1	B	926	GLN
1	B	953	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	615/652 (94%)	0.08	26 (4%) 40 33	39, 66, 143, 171	0
1	B	615/652 (94%)	0.20	41 (6%) 21 15	39, 71, 150, 171	0
2	S	11/11 (100%)	-0.82	0 100 100	81, 87, 94, 98	0
2	W	11/11 (100%)	-0.79	0 100 100	89, 108, 126, 129	0
3	U	12/17 (70%)	-0.72	0 100 100	68, 88, 130, 148	0
3	Z	12/17 (70%)	-0.59	0 100 100	61, 85, 128, 139	0
4	V	12/12 (100%)	-0.96	0 100 100	74, 89, 114, 136	1 (8%)
4	X	12/12 (100%)	-0.92	0 100 100	80, 90, 105, 112	1 (8%)
5	T	7/14 (50%)	-0.81	0 100 100	63, 74, 93, 101	0
5	Y	7/14 (50%)	-0.72	0 100 100	55, 66, 81, 94	0
All	All	1314/1412 (93%)	0.08	67 (5%) 32 25	39, 70, 147, 171	2 (0%)

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	506	GLN	8.1
1	B	770	PRO	7.4
1	B	519	ILE	7.0
1	B	567	GLY	6.7
1	B	774	VAL	6.3
1	A	567	GLY	6.3
1	A	775	GLN	5.9
1	A	486	LEU	5.2
1	A	772	MET	5.1
1	B	775	GLN	4.9
1	A	1009	VAL	4.7
1	A	765	LEU	4.7
1	B	794	LYS	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	486	LEU	4.1
1	B	497	VAL	4.0
1	A	796	VAL	3.9
1	B	507	ARG	3.9
1	A	757	PRO	3.9
1	B	503	LEU	3.8
1	B	785	ARG	3.8
1	B	761	LYS	3.7
1	B	508	SER	3.6
1	B	524	LEU	3.6
1	B	764	HIS	3.5
1	B	768	GLN	3.5
1	A	785	ARG	3.4
1	B	565	CYS	3.4
1	B	526	ARG	3.2
1	A	505	LYS	3.1
1	B	523	ILE	3.1
1	B	1009	VAL	3.1
1	B	795	SER	3.0
1	B	772	MET	3.0
1	B	796	VAL	3.0
1	A	503	LEU	3.0
1	A	526	ARG	2.9
1	A	764	HIS	2.9
1	A	761	LYS	2.9
1	B	568	GLN	2.8
1	B	566	GLY	2.8
1	B	797	PHE	2.7
1	A	812	CYS	2.7
1	B	504	ALA	2.6
1	A	566	GLY	2.6
1	A	508	SER	2.6
1	A	758	SER	2.6
1	B	505	LYS	2.5
1	B	765	LEU	2.5
1	A	519	ILE	2.5
1	B	769	LEU	2.4
1	B	484	PHE	2.3
1	A	1008	ALA	2.3
1	A	509	VAL	2.3
1	B	771	GLU	2.2
1	B	453	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	501	LEU	2.2
1	A	565	CYS	2.1
1	A	484	PHE	2.1
1	B	758	SER	2.1
1	A	768	GLN	2.1
1	B	491	GLY	2.1
1	B	643	ASN	2.1
1	B	485	HIS	2.1
1	B	480	LEU	2.0
1	A	504	ALA	2.0
1	B	786	LEU	2.0
1	B	501	LEU	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
6	BA	B	1101	1/1	1.00	0.20	2.82	73,73,73,73	0
6	BA	A	1101	1/1	1.00	0.21	1.54	71,71,71,71	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.