Full wwPDB X-ray Structure Validation Report (i)

Nov 15, 2023 – 03:17 PM JST

PDB ID	:	6JDQ
Title	:	Crystal structure of Nme1Cas9 in complex with sgRNA
Authors	:	Sun, W.; Yang, J.; Cheng, Z.; Liu, C.; Wang, K.; Huang, X.; Wang, Y.
Deposited on	:	2019-02-02
Resolution	:	2.95 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*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

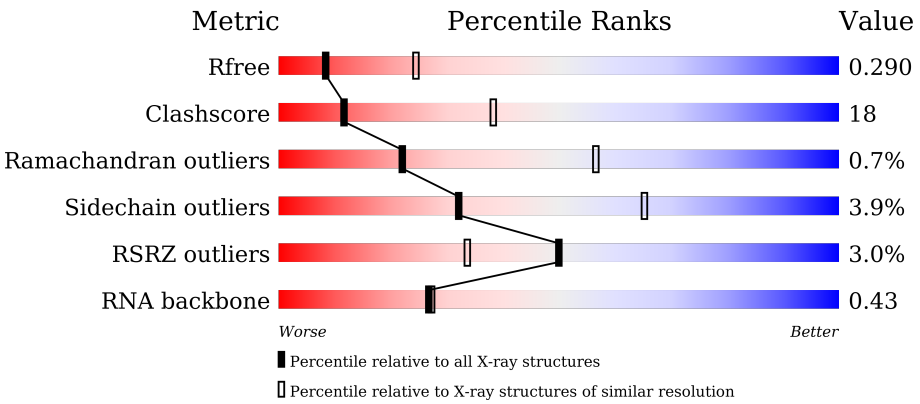
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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}	130704	3104 (3.00-2.92)
Clashscore	141614	3462 (3.00-2.92)
Ramachandran outliers	138981	3340 (3.00-2.92)
Sidechain outliers	138945	3343 (3.00-2.92)
RSRZ outliers	127900	2986 (3.00-2.92)
RNA backbone	3102	1065 (3.22-2.70)

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 <=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	1092	2%, 69%, 27%, • •
2	B	135	13%, 18%, 45%, 18%, 6%, 13%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 10857 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 CRISPR-associated endonuclease Cas9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1069	8365	5292	1514	1535	24	0	3	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1083	SER	-	expression tag	UNP C9X1G5
A	1084	GLU	-	expression tag	UNP C9X1G5
A	1085	HIS	-	expression tag	UNP C9X1G5
A	1086	HIS	-	expression tag	UNP C9X1G5
A	1087	HIS	-	expression tag	UNP C9X1G5
A	1088	HIS	-	expression tag	UNP C9X1G5
A	1089	HIS	-	expression tag	UNP C9X1G5
A	1090	HIS	-	expression tag	UNP C9X1G5
A	1091	HIS	-	expression tag	UNP C9X1G5
A	1092	HIS	-	expression tag	UNP C9X1G5

- Molecule 2 is a RNA chain called sgRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	B	117	2473	1107	426	823	117	0	0	0

- Molecule 3 is water.

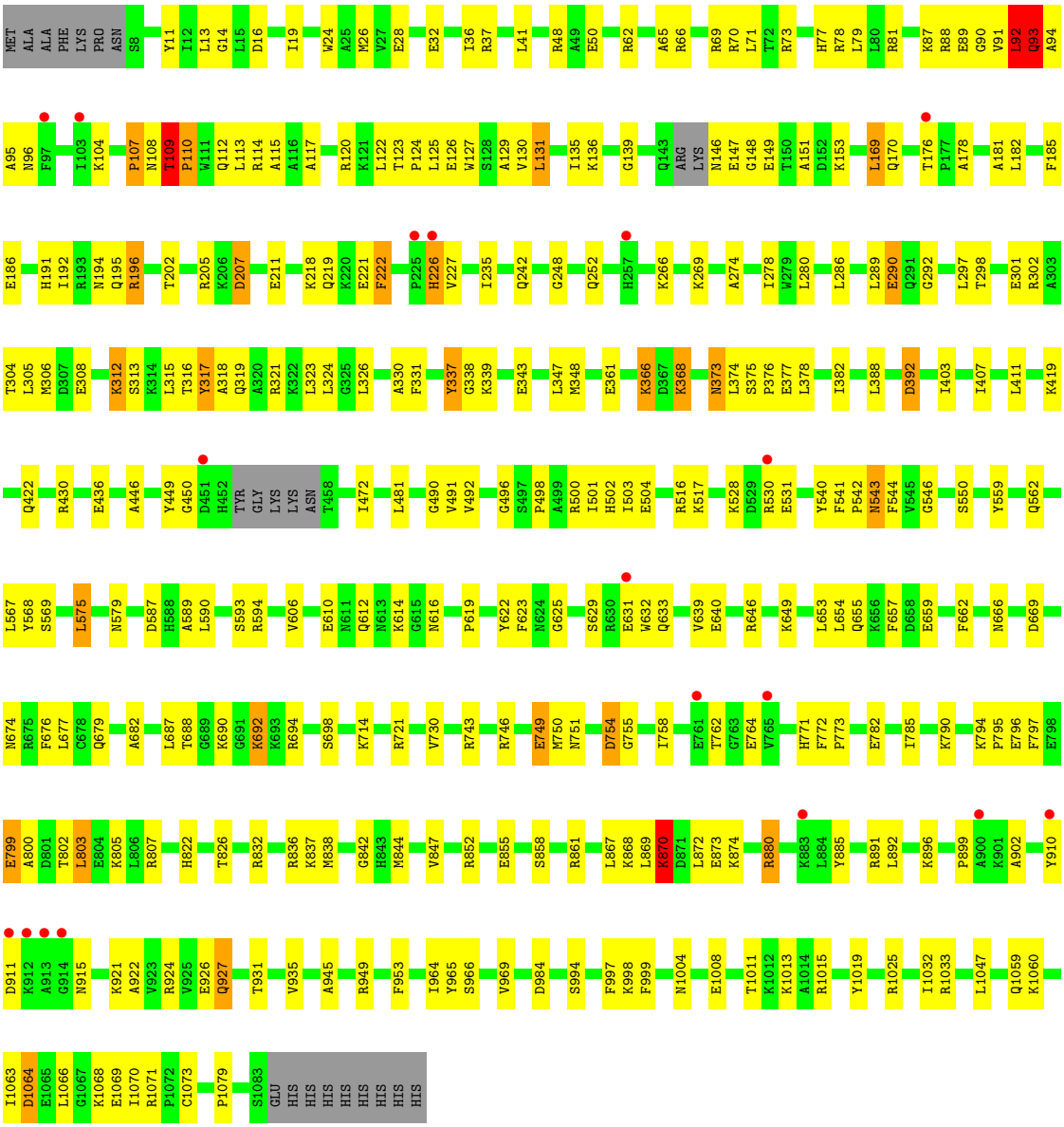
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	16	16	16	0	0
3	B	3	3	3	0	0

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.

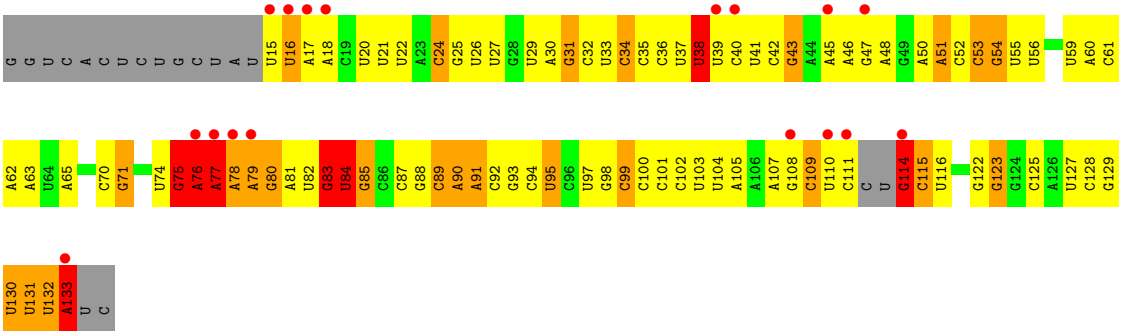
- Molecule 1: CRISPR-associated endonuclease Cas9

Chain A:



• Molecule 2: sgRNA

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	72.08Å 122.45Å 238.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.91 – 2.95 48.91 – 2.95	Depositor EDS
% Data completeness (in resolution range)	98.7 (48.91-2.95) 98.7 (48.91-2.95)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.27 (at 2.96Å)	Xtriage
Refinement program	PHENIX (1.14_3247: ???)	Depositor
R, R_{free}	0.255 , 0.290 0.255 , 0.290	Depositor DCC
R_{free} test set	2241 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	50.8	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 50.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.87	EDS
Total number of atoms	10857	wwPDB-VP
Average B, all atoms (Å^2)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the 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	0.81	11/8529 (0.1%)	0.72	17/11518 (0.1%)
2	B	1.21	40/2759 (1.4%)	1.25	28/4290 (0.7%)
All	All	0.92	51/11288 (0.5%)	0.90	45/15808 (0.3%)

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
2	B	0	1

All (51) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	91	A	O3'-P	-13.29	1.45	1.61
2	B	130	U	O3'-P	-12.66	1.46	1.61
2	B	85	G	O3'-P	-11.89	1.46	1.61
2	B	92	C	O3'-P	-10.24	1.48	1.61
1	A	984	ASP	C-N	10.21	1.57	1.34
2	B	36	C	O3'-P	-9.40	1.49	1.61
2	B	122	G	O3'-P	-9.26	1.50	1.61
2	B	81	A	O3'-P	-9.25	1.50	1.61
2	B	89	C	O3'-P	-8.96	1.50	1.61
2	B	123	G	O3'-P	-8.83	1.50	1.61
2	B	35	C	O3'-P	-8.74	1.50	1.61
2	B	90	A	O3'-P	-8.73	1.50	1.61
2	B	53	C	O3'-P	-7.95	1.51	1.61
2	B	92	C	P-OP1	-7.75	1.35	1.49
2	B	93	G	O3'-P	-7.65	1.51	1.61
2	B	90	A	P-OP2	-7.53	1.36	1.49
2	B	37	U	O3'-P	-7.50	1.52	1.61
2	B	84	U	O3'-P	-7.49	1.52	1.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	91	A	P-OP1	-7.31	1.36	1.49
2	B	83	G	O3'-P	-7.18	1.52	1.61
2	B	92	C	P-OP2	-7.17	1.36	1.49
2	B	80	G	O3'-P	-7.00	1.52	1.61
2	B	91	A	P-OP2	-6.96	1.37	1.49
1	A	317	TYR	CE1-CZ	-6.94	1.29	1.38
2	B	90	A	P-OP1	-6.83	1.37	1.49
2	B	99	C	O3'-P	-6.78	1.53	1.61
2	B	131	U	O3'-P	-6.62	1.53	1.61
2	B	34	C	O3'-P	-6.37	1.53	1.61
2	B	89	C	P-OP1	-6.36	1.38	1.49
2	B	89	C	P-OP2	-6.27	1.38	1.49
2	B	88	G	O3'-P	-6.07	1.53	1.61
2	B	100	C	O3'-P	-5.99	1.53	1.61
1	A	337	TYR	CE1-CZ	-5.95	1.30	1.38
1	A	313	SER	CB-OG	-5.91	1.34	1.42
1	A	337	TYR	CG-CD2	-5.85	1.31	1.39
2	B	54	G	O3'-P	-5.80	1.54	1.61
1	A	337	TYR	CE2-CZ	-5.78	1.31	1.38
1	A	799	GLU	CD-OE1	-5.63	1.19	1.25
1	A	348	MET	C-O	-5.44	1.13	1.23
2	B	101	C	O3'-P	-5.43	1.54	1.61
2	B	71	G	O3'-P	-5.39	1.54	1.61
2	B	51	A	O3'-P	-5.30	1.54	1.61
1	A	782	GLU	CD-OE2	-5.19	1.20	1.25
1	A	337	TYR	CB-CG	-5.19	1.43	1.51
2	B	94	C	P-OP1	-5.15	1.40	1.49
2	B	131	U	P-OP2	-5.15	1.40	1.49
2	B	95	U	O3'-P	-5.11	1.55	1.61
2	B	94	C	P-OP2	-5.05	1.40	1.49
2	B	93	G	P-OP2	-5.04	1.40	1.49
1	A	749	GLU	CD-OE2	-5.03	1.20	1.25
2	B	50	A	O3'-P	-5.02	1.55	1.61

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	984	ASP	O-C-N	-15.09	98.56	122.70
2	B	77	A	N9-C1'-C2'	13.94	132.12	114.00
2	B	85	G	O5'-P-OP1	-11.42	95.42	105.70
2	B	76	A	O5'-P-OP2	-10.90	95.89	105.70
2	B	78	A	C4'-C3'-O3'	10.30	133.59	113.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	984	ASP	CA-C-N	10.19	139.62	117.20
1	A	546	GLY	C-N-CA	8.61	143.22	121.70
2	B	75	G	C2'-C3'-O3'	8.42	128.03	109.50
2	B	85	G	O5'-P-OP2	8.31	120.68	110.70
1	A	546	GLY	O-C-N	8.30	135.98	122.70
2	B	34	C	O5'-P-OP1	-8.08	98.43	105.70
2	B	77	A	C8-N9-C1'	-7.93	113.42	127.70
1	A	92	LEU	N-CA-C	-7.77	90.03	111.00
2	B	78	A	N9-C1'-C2'	-7.74	103.49	112.00
2	B	91	A	O5'-P-OP2	-7.50	98.95	105.70
1	A	984	ASP	C-N-CA	7.33	140.03	121.70
2	B	90	A	N9-C1'-C2'	7.12	123.26	114.00
2	B	77	A	C4-N9-C1'	6.61	138.20	126.30
1	A	546	GLY	CA-C-N	-6.59	102.69	117.20
2	B	76	A	N9-C1'-C2'	6.48	122.43	114.00
2	B	77	A	C4'-C3'-O3'	6.35	125.70	113.00
2	B	99	C	O5'-P-OP1	-6.08	100.23	105.70
2	B	133	A	O5'-P-OP2	5.92	117.81	110.70
2	B	77	A	C1'-O4'-C4'	-5.92	105.17	109.90
1	A	870	LYS	CD-CE-NZ	5.88	125.21	111.70
1	A	169	LEU	CA-CB-CG	5.83	128.72	115.30
1	A	93	GLN	N-CA-C	-5.76	95.44	111.00
1	A	107	PRO	N-CA-C	5.72	126.97	112.10
2	B	78	A	C2'-C3'-O3'	-5.52	97.35	109.50
2	B	38	U	C5-C6-N1	5.51	125.45	122.70
1	A	368	LYS	N-CA-CB	5.44	120.40	110.60
2	B	101	C	C1'-C2'-O2'	-5.39	94.42	110.60
1	A	750	MET	N-CA-C	-5.32	96.64	111.00
1	A	543	ASN	N-CA-C	-5.31	96.66	111.00
1	A	803	LEU	CB-CG-CD2	-5.30	101.99	111.00
2	B	114	G	C8-N9-C1'	-5.23	120.21	127.00
2	B	114	G	C4-N9-C1'	5.22	133.29	126.50
2	B	16	U	C5-C6-N1	5.20	125.30	122.70
2	B	133	A	C5'-C4'-O4'	5.20	115.34	109.10
2	B	93	G	N9-C1'-C2'	5.18	120.73	114.00
1	A	339	LYS	N-CA-C	5.16	124.93	111.00
2	B	38	U	C5-C4-O4	-5.07	122.86	125.90
2	B	78	A	C4-N9-C1'	-5.04	117.23	126.30
1	A	338	GLY	N-CA-C	5.02	125.66	113.10
2	B	77	A	O4'-C1'-N9	5.01	112.21	108.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	77	A	Sidechain

5.2 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 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	8365	0	8161	276	1
2	B	2473	0	1256	89	0
3	A	16	0	0	2	0
3	B	3	0	0	0	0
All	All	10857	0	9417	355	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 18.

All (355) 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:109:THR:H	1:A:110:PRO:HD3	1.11	1.15
1:A:218:LYS:O	1:A:222:PHE:CE2	2.05	1.10
2:B:74:U:H5"	2:B:74:U:H6	1.14	1.07
1:A:148:GLY:O	1:A:149:GLU:HG2	1.53	1.07
1:A:218:LYS:O	1:A:222:PHE:CD2	2.13	1.01
1:A:266:LYS:NZ	1:A:422:GLN:OE1	1.94	1.00
1:A:109:THR:N	1:A:110:PRO:HD3	1.73	0.99
2:B:38:U:H5"	2:B:38:U:H6	1.27	0.95
1:A:298:THR:OG1	1:A:301:GLU:OE1	1.84	0.95
1:A:91:VAL:CG2	1:A:124:PRO:HB3	1.96	0.95
1:A:91:VAL:HG23	1:A:124:PRO:HB3	1.49	0.94
2:B:79:A:H8	2:B:79:A:H5"	1.31	0.93
1:A:146:ASN:N	1:A:149:GLU:OE2	2.03	0.91
1:A:593:SER:OG	1:A:749:GLU:OE1	1.89	0.90
2:B:74:U:H5"	2:B:74:U:C6	2.07	0.89
1:A:319:GLN:HE21	1:A:323:LEU:CD1	1.87	0.87
1:A:109:THR:N	1:A:110:PRO:CD	2.38	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:LEU:O	1:A:135:ILE:HD12	1.76	0.85
1:A:319:GLN:NE2	1:A:323:LEU:CD1	2.40	0.85
1:A:131:LEU:HG	1:A:235:ILE:HD13	1.59	0.84
1:A:373:ASN:O	1:A:373:ASN:ND2	2.10	0.84
1:A:880:ARG:HH11	1:A:880:ARG:HG2	1.43	0.84
1:A:205:ARG:NH1	1:A:242:GLN:OE1	2.11	0.83
2:B:38:U:H5"	2:B:38:U:C6	2.13	0.83
1:A:500:ARG:NH2	1:A:690:LYS:HB2	1.95	0.81
1:A:298:THR:CB	1:A:301:GLU:OE1	2.30	0.80
1:A:308:GLU:OE1	1:A:312:LYS:HD3	1.81	0.79
1:A:131:LEU:HG	1:A:235:ILE:CD1	2.13	0.79
1:A:148:GLY:O	1:A:149:GLU:CG	2.31	0.79
1:A:562:GLN:HE22	1:A:567:LEU:HB2	1.48	0.78
2:B:15:U:H2'	2:B:16:U:H6	1.47	0.78
1:A:298:THR:N	1:A:301:GLU:OE1	2.18	0.77
2:B:74:U:H6	2:B:74:U:C5'	1.95	0.76
2:B:79:A:H5"	2:B:79:A:C8	2.18	0.76
1:A:109:THR:H	1:A:110:PRO:CD	1.95	0.76
1:A:688:THR:HG23	1:A:688:THR:O	1.85	0.75
1:A:692:LYS:HD3	1:A:692:LYS:H	1.52	0.75
1:A:880:ARG:HH11	1:A:880:ARG:CG	2.00	0.75
1:A:803:LEU:H	1:A:803:LEU:HD12	1.52	0.74
1:A:319:GLN:HE21	1:A:323:LEU:HD13	1.52	0.74
2:B:79:A:H8	2:B:79:A:C5'	2.00	0.73
1:A:743:ARG:NH2	3:A:1101:HOH:O	2.19	0.73
1:A:562:GLN:NE2	1:A:567:LEU:HB2	2.03	0.72
1:A:306:MET:O	1:A:306:MET:HG3	1.88	0.72
1:A:196:ARG:O	2:B:24:C:O2'	2.08	0.72
1:A:368:LYS:O	1:A:368:LYS:HG2	1.89	0.72
1:A:298:THR:HG23	1:A:301:GLU:OE1	1.90	0.71
1:A:568:TYR:OH	1:A:614:LYS:NZ	2.22	0.71
1:A:181:ALA:O	1:A:185:PHE:HB2	1.91	0.71
1:A:375:SER:OG	1:A:376:PRO:HD2	1.90	0.71
1:A:289:LEU:HD12	1:A:331:PHE:CE1	2.26	0.71
1:A:266:LYS:HZ3	1:A:422:GLN:CD	1.95	0.70
2:B:17:A:H2'	2:B:18:A:H8	1.56	0.70
2:B:60:A:H2'	2:B:61:C:C6	2.27	0.70
2:B:76:A:C2	2:B:77:A:N6	2.59	0.70
2:B:133:A:OP2	2:B:133:A:O3'	2.05	0.70
1:A:308:GLU:OE1	1:A:312:LYS:CD	2.39	0.70
1:A:147:GLU:CB	1:A:392:ASP:OD2	2.41	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:594:ARG:NH2	1:A:659:GLU:OE2	2.22	0.69
2:B:15:U:H2'	2:B:16:U:C6	2.28	0.69
1:A:91:VAL:H	1:A:226:HIS:CD2	2.10	0.69
1:A:218:LYS:C	1:A:222:PHE:HE2	1.97	0.69
1:A:559:TYR:CZ	1:A:575:LEU:HG	2.28	0.69
1:A:91:VAL:HG21	1:A:124:PRO:HB3	1.74	0.68
1:A:148:GLY:C	1:A:149:GLU:CG	2.62	0.67
1:A:218:LYS:C	1:A:222:PHE:CE2	2.68	0.67
1:A:319:GLN:NE2	1:A:323:LEU:HD11	2.09	0.67
1:A:96:ASN:O	1:A:104:LYS:N	2.20	0.66
1:A:682:ALA:O	1:A:694:ARG:NH2	2.28	0.66
1:A:807:ARG:NH2	1:A:822:HIS:O	2.24	0.66
1:A:298:THR:CG2	1:A:301:GLU:OE1	2.43	0.66
1:A:692:LYS:HD3	1:A:692:LYS:N	2.10	0.66
2:B:42:C:H2'	2:B:43:G:H5'	1.78	0.65
1:A:191:HIS:CE1	1:A:195:GLN:HG3	2.31	0.65
1:A:227:VAL:HG12	1:A:227:VAL:O	1.96	0.65
2:B:16:U:H3'	2:B:17:A:H8	1.63	0.64
2:B:59:U:H2'	2:B:60:A:H8	1.62	0.64
1:A:178:ALA:O	1:A:182:LEU:HD12	1.97	0.64
2:B:76:A:C2	2:B:77:A:C6	2.86	0.64
1:A:503:ILE:HD13	1:A:677:LEU:HD23	1.80	0.64
1:A:331:PHE:HB2	1:A:337:TYR:CE2	2.34	0.63
1:A:587:ASP:OD2	1:A:614:LYS:NZ	2.21	0.63
1:A:869:LEU:HD11	1:A:892:LEU:HB3	1.80	0.62
2:B:54:G:H2'	2:B:55:U:O4'	1.99	0.62
1:A:502:HIS:HB3	1:A:730:VAL:HG22	1.81	0.62
2:B:17:A:H2'	2:B:18:A:C8	2.35	0.62
1:A:361:GLU:OE2	1:A:366:LYS:NZ	2.32	0.62
1:A:375:SER:OG	1:A:377:GLU:OE1	2.17	0.62
2:B:60:A:H2'	2:B:61:C:H6	1.64	0.62
1:A:148:GLY:C	1:A:149:GLU:HG2	2.19	0.61
1:A:290:GLU:OE1	1:A:330:ALA:HB2	2.00	0.61
2:B:41:U:H2'	2:B:42:C:C6	2.34	0.61
2:B:132:U:H4'	2:B:133:A:O5'	2.01	0.61
1:A:867:LEU:HD21	1:A:872:LEU:HB2	1.82	0.61
2:B:76:A:H2'	2:B:77:A:C8	2.35	0.60
1:A:880:ARG:NH2	2:B:53:C:O2	2.34	0.60
1:A:147:GLU:CB	1:A:392:ASP:CG	2.70	0.60
1:A:762:THR:HB	1:A:764:GLU:OE1	2.02	0.60
1:A:91:VAL:N	1:A:226:HIS:CD2	2.70	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:861:ARG:HH21	1:A:922:ALA:HB3	1.67	0.60
1:A:66:ARG:HD2	1:A:70:ARG:HH21	1.67	0.59
1:A:324:LEU:HB2	1:A:326:LEU:HG	1.84	0.59
1:A:266:LYS:NZ	1:A:422:GLN:CD	2.51	0.59
1:A:321:ARG:NH2	1:A:343:GLU:OE2	2.33	0.59
1:A:290:GLU:OE1	1:A:330:ALA:N	2.35	0.59
2:B:76:A:H8	2:B:76:A:P	2.26	0.59
2:B:76:A:P	2:B:76:A:C8	2.96	0.59
1:A:79:LEU:HD11	1:A:136:LYS:HG3	1.85	0.58
1:A:758:ILE:CB	1:A:762:THR:OG1	2.52	0.58
1:A:297:LEU:HD22	1:A:301:GLU:HB3	1.84	0.58
1:A:500:ARG:HH22	1:A:690:LYS:HB2	1.67	0.58
1:A:569:SER:HB3	1:A:610[A]:GLU:HG2	1.86	0.58
1:A:870:LYS:H	1:A:870:LYS:HD2	1.68	0.58
1:A:113:LEU:HD13	1:A:126:GLU:HB3	1.86	0.57
1:A:319:GLN:HE21	1:A:323:LEU:HD11	1.64	0.57
1:A:222:PHE:CD2	1:A:222:PHE:N	2.72	0.57
1:A:73:ARG:NH2	2:B:87:C:OP2	2.31	0.57
1:A:899:PRO:HA	1:A:902:ALA:HB3	1.87	0.57
1:A:698:SER:HB2	1:A:730:VAL:HG23	1.85	0.56
2:B:59:U:H2'	2:B:60:A:C8	2.40	0.56
2:B:75:G:C8	2:B:75:G:H3'	2.41	0.56
1:A:794:LYS:N	1:A:795:PRO:CD	2.64	0.56
1:A:1063:ILE:HG13	1:A:1070:ILE:HG12	1.87	0.56
2:B:38:U:C6	2:B:38:U:C5'	2.87	0.56
2:B:130:U:H2'	2:B:131:U:C6	2.40	0.56
1:A:112:GLN:HG3	1:A:182:LEU:HD21	1.88	0.56
1:A:290:GLU:OE1	1:A:330:ALA:CA	2.53	0.56
1:A:873:GLU:OE2	1:A:885:TYR:OH	2.24	0.56
1:A:541:PHE:O	1:A:544:PHE:HB2	2.06	0.56
1:A:640:GLU:HA	1:A:649:LYS:HE3	1.89	0.55
1:A:802:THR:OG1	1:A:805:LYS:HG3	2.07	0.55
2:B:51:A:H2'	2:B:52:C:C6	2.42	0.55
1:A:91:VAL:HG13	1:A:91:VAL:O	2.06	0.55
1:A:302:ARG:HG3	1:A:302:ARG:HH11	1.72	0.55
2:B:82:U:O2'	2:B:83:G:H5'	2.07	0.55
1:A:375:SER:OG	1:A:376:PRO:CD	2.56	0.54
2:B:90:A:H2'	2:B:91:A:C8	2.41	0.54
1:A:248:GLY:HA2	1:A:388:LEU:HD22	1.90	0.54
1:A:868:LYS:HG2	1:A:899:PRO:HG3	1.90	0.54
1:A:568:TYR:CD2	1:A:606:VAL:HG11	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:ALA:HB2	1:A:122:LEU:HD11	1.90	0.53
1:A:374:LEU:N	1:A:374:LEU:HD23	2.22	0.53
1:A:88:ARG:O	1:A:89:GLU:HG2	2.09	0.53
1:A:266:LYS:HZ1	1:A:422:GLN:HE22	1.57	0.53
1:A:880:ARG:HG2	1:A:880:ARG:NH1	2.18	0.53
1:A:147:GLU:CB	1:A:392:ASP:OD1	2.56	0.53
1:A:292:GLY:HA3	1:A:653:LEU:O	2.09	0.53
1:A:317:TYR:CE1	1:A:347:LEU:HD13	2.43	0.53
1:A:181:ALA:HA	1:A:185:PHE:CD2	2.44	0.52
1:A:754:ASP:OD1	1:A:755:GLY:N	2.27	0.52
2:B:97:U:H2'	2:B:98:G:H8	1.74	0.52
2:B:79:A:C8	2:B:79:A:C5'	2.85	0.52
1:A:169:LEU:HD12	1:A:170:GLN:N	2.24	0.52
1:A:90:GLY:CA	1:A:226:HIS:HD2	2.23	0.52
1:A:500:ARG:HH11	1:A:694:ARG:HA	1.75	0.52
1:A:931:THR:HG21	1:A:1025:ARG:NH1	2.24	0.52
2:B:75:G:O5'	2:B:76:A:OP2	2.27	0.52
1:A:1011:THR:HG23	1:A:1013:LYS:H	1.75	0.52
2:B:70:C:H3'	2:B:71:G:H8	1.75	0.52
1:A:81:ARG:HH12	2:B:84:U:H4'	1.76	0.51
1:A:289:LEU:HD12	1:A:331:PHE:CZ	2.44	0.51
1:A:378:LEU:O	1:A:382:ILE:HG13	2.10	0.51
1:A:616:ASN:CB	1:A:749:GLU:OE2	2.58	0.51
2:B:38:U:H6	2:B:38:U:C5'	2.09	0.51
2:B:41:U:H2'	2:B:42:C:H6	1.73	0.51
1:A:88:ARG:C	1:A:89:GLU:HG2	2.31	0.51
1:A:115:ALA:HB2	1:A:182:LEU:HD13	1.92	0.51
1:A:304:THR:HG22	1:A:323:LEU:HD23	1.92	0.51
1:A:501:ILE:HG13	1:A:687:LEU:HD21	1.93	0.51
1:A:629:SER:OG	1:A:631:GLU:HG3	2.10	0.51
2:B:29:U:H2'	2:B:30:A:H8	1.75	0.51
2:B:115:C:H2'	2:B:116:U:H6	1.74	0.51
1:A:127:TRP:CZ3	1:A:219:GLN:HG3	2.44	0.51
2:B:76:A:N1	2:B:77:A:N6	2.58	0.51
1:A:407:ILE:O	1:A:411:LEU:HG	2.10	0.51
1:A:500:ARG:NH1	1:A:694:ARG:HA	2.25	0.51
2:B:16:U:H3'	2:B:17:A:C8	2.43	0.51
1:A:280:LEU:HD11	1:A:436:GLU:HG3	1.93	0.51
2:B:74:U:C6	2:B:74:U:C5'	2.83	0.51
1:A:500:ARG:NH2	1:A:690:LYS:CB	2.73	0.50
2:B:42:C:C2'	2:B:43:G:H5'	2.39	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:114:G:HO2'	2:B:115:C:P	2.34	0.50
1:A:803:LEU:HG	1:A:826:THR:HG22	1.93	0.50
2:B:76:A:H8	2:B:76:A:O5'	1.93	0.50
1:A:28:GLU:OE1	1:A:37:ARG:NE	2.35	0.50
1:A:248:GLY:O	1:A:252:GLN:HG3	2.11	0.50
1:A:1060:LYS:HG2	1:A:1073:CYS:SG	2.51	0.50
2:B:76:A:C8	2:B:76:A:O5'	2.65	0.50
2:B:98:G:H2'	2:B:99:C:H6	1.76	0.50
2:B:62:A:H2'	2:B:63:A:C8	2.47	0.50
1:A:195:GLN:OE1	1:A:921:LYS:NZ	2.38	0.50
1:A:803:LEU:HD12	1:A:803:LEU:N	2.20	0.50
2:B:30:A:H2'	2:B:31:G:C8	2.47	0.50
1:A:500:ARG:HH11	1:A:500:ARG:HG2	1.77	0.50
1:A:266:LYS:HZ1	1:A:422:GLN:NE2	2.09	0.49
1:A:998:LYS:HE2	1:A:998:LYS:HA	1.94	0.49
1:A:373:ASN:HD22	1:A:373:ASN:C	2.04	0.49
1:A:629:SER:O	1:A:633:GLN:HG2	2.12	0.49
1:A:869:LEU:HD12	1:A:892:LEU:HD13	1.92	0.49
1:A:800:ALA:HB1	1:A:805:LYS:HB2	1.95	0.49
1:A:999:PHE:CE1	1:A:1059:GLN:HA	2.46	0.49
1:A:176:THR:HB	1:A:211:GLU:HG3	1.94	0.49
1:A:589:ALA:O	1:A:619:PRO:HD3	2.12	0.49
2:B:26:U:H2'	2:B:27:U:C6	2.47	0.49
2:B:104:U:H2'	2:B:105:A:H8	1.77	0.49
2:B:115:C:H2'	2:B:116:U:C6	2.48	0.49
1:A:528:LYS:O	1:A:531:GLU:HG2	2.12	0.48
2:B:76:A:H8	2:B:76:A:C5'	2.24	0.48
1:A:92:LEU:HD12	1:A:93:GLN:H	1.77	0.48
1:A:120:ARG:HD2	1:A:122:LEU:HD23	1.95	0.48
2:B:102:C:H2'	2:B:103:U:H6	1.78	0.48
1:A:50:GLU:OE2	1:A:516:ARG:NH2	2.46	0.48
1:A:120:ARG:NH2	1:A:126:GLU:OE2	2.46	0.48
1:A:838:MET:HG3	2:B:127:U:C5	2.49	0.48
1:A:542:PRO:HD2	1:A:543:ASN:H	1.79	0.48
1:A:92:LEU:HD13	1:A:94:ALA:HB2	1.95	0.47
1:A:139:GLY:O	2:B:22:U:H4'	2.13	0.47
1:A:541:PHE:CE1	1:A:579:ASN:ND2	2.82	0.47
1:A:794:LYS:HA	1:A:795:PRO:HD2	1.40	0.47
1:A:218:LYS:O	1:A:221:GLU:HB2	2.15	0.47
1:A:692:LYS:N	1:A:692:LYS:CD	2.73	0.47
1:A:870:LYS:H	1:A:870:LYS:CD	2.24	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:836:ARG:HH21	2:B:125:C:H4'	1.79	0.47
1:A:107:PRO:HD2	1:A:129:ALA:HB2	1.97	0.47
1:A:653:LEU:O	1:A:655:GLN:HG3	2.14	0.47
1:A:924:ARG:HG3	2:B:56:U:OP1	2.14	0.47
1:A:87:LYS:HA	1:A:92:LEU:HB3	1.95	0.47
1:A:302:ARG:HG3	1:A:302:ARG:NH1	2.30	0.47
1:A:114:ARG:HG3	1:A:130:VAL:HB	1.97	0.47
1:A:361:GLU:HG3	1:A:366:LYS:HD3	1.95	0.47
1:A:949:ARG:CZ	1:A:1079:PRO:HG2	2.44	0.47
1:A:91:VAL:HG21	1:A:124:PRO:CB	2.45	0.47
1:A:852:ARG:NH2	1:A:926:GLU:OE2	2.45	0.47
1:A:1032:ILE:HD11	1:A:1047:LEU:HD12	1.97	0.47
2:B:109:C:H6	2:B:109:C:O5'	1.98	0.46
1:A:186:GLU:OE1	1:A:192:ILE:HD11	2.14	0.46
1:A:91:VAL:HB	1:A:226:HIS:CE1	2.51	0.46
1:A:688:THR:O	1:A:688:THR:CG2	2.57	0.46
1:A:844:MET:HE3	1:A:945:ALA:HA	1.97	0.46
2:B:74:U:C6	2:B:74:U:C4'	2.98	0.46
1:A:298:THR:CA	1:A:301:GLU:OE1	2.64	0.46
1:A:785:ILE:HD13	1:A:790:LYS:C	2.35	0.46
1:A:965:TYR:O	1:A:969:VAL:HG23	2.16	0.46
1:A:71:LEU:HD23	1:A:71:LEU:HA	1.81	0.46
1:A:207:ASP:N	1:A:207:ASP:OD1	2.49	0.46
1:A:832:ARG:HD2	1:A:1019:TYR:CD1	2.51	0.46
1:A:90:GLY:CA	1:A:226:HIS:CD2	2.99	0.45
1:A:331:PHE:HB2	1:A:337:TYR:CZ	2.50	0.45
1:A:274:ALA:O	1:A:278:ILE:HG13	2.17	0.45
2:B:114:G:O2'	2:B:115:C:OP1	2.31	0.45
2:B:43:G:O2'	2:B:45:A:N7	2.43	0.45
2:B:76:A:C8	2:B:76:A:OP1	2.70	0.45
1:A:269:LYS:HE2	1:A:419:LYS:CB	2.47	0.45
1:A:286:LEU:HD23	1:A:297:LEU:HD11	1.97	0.45
1:A:542:PRO:C	1:A:544:PHE:N	2.69	0.45
2:B:104:U:H2'	2:B:105:A:C8	2.51	0.45
1:A:290:GLU:OE1	1:A:330:ALA:CB	2.63	0.45
1:A:319:GLN:NE2	1:A:323:LEU:HD13	2.20	0.45
2:B:31:G:H2'	2:B:32:C:H6	1.82	0.45
1:A:541:PHE:HE1	1:A:579:ASN:HD21	1.64	0.45
1:A:868:LYS:HG2	1:A:899:PRO:CG	2.46	0.45
1:A:874:LYS:O	1:A:926:GLU:HG2	2.17	0.45
1:A:910:TYR:HA	1:A:915:ASN:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:378:LEU:HD13	1:A:403:ILE:HG22	1.98	0.44
1:A:517:LYS:HA	1:A:517:LYS:HD3	1.84	0.44
1:A:77:HIS:CE1	1:A:81:ARG:HD2	2.52	0.44
1:A:472:ILE:HD11	1:A:676:PHE:CD2	2.52	0.44
2:B:75:G:C8	2:B:75:G:C3'	3.00	0.44
1:A:122:LEU:HB3	1:A:126:GLU:HB2	1.99	0.44
1:A:337:TYR:OH	1:A:343:GLU:OE2	2.26	0.44
1:A:90:GLY:HA3	1:A:226:HIS:CD2	2.52	0.44
1:A:194:ASN:ND2	1:A:202:THR:OG1	2.50	0.44
2:B:90:A:H2'	2:B:91:A:H8	1.82	0.44
1:A:16:ASP:HA	1:A:504:GLU:O	2.18	0.44
1:A:616:ASN:HB2	1:A:749:GLU:OE2	2.17	0.44
2:B:133:A:P	2:B:133:A:H3'	2.58	0.44
1:A:301:GLU:HG2	1:A:324:LEU:HD22	1.99	0.44
1:A:896:LYS:HB3	1:A:896:LYS:HE2	1.73	0.44
1:A:847:VAL:O	2:B:26:U:O2'	2.35	0.43
1:A:616:ASN:HB3	1:A:749:GLU:OE2	2.19	0.43
1:A:842:GLY:HA2	1:A:966:SER:OG	2.18	0.43
1:A:911:ASP:OD1	1:A:915:ASN:HB2	2.18	0.43
1:A:1064:ASP:HB2	1:A:1068:LYS:O	2.19	0.43
1:A:772:PHE:HA	1:A:773:PRO:HD3	1.74	0.43
1:A:891:ARG:HD3	1:A:902:ALA:O	2.19	0.43
1:A:953:PHE:CE2	1:A:997:PHE:HB2	2.54	0.43
1:A:562:GLN:NE2	1:A:567:LEU:HD13	2.34	0.43
1:A:625:GLY:HA2	1:A:632:TRP:CG	2.54	0.43
1:A:743:ARG:HG2	1:A:746:ARG:HH22	1.83	0.43
2:B:114:G:O4'	2:B:114:G:P	2.76	0.43
1:A:147:GLU:HA	1:A:392:ASP:OD1	2.19	0.43
2:B:62:A:H2'	2:B:63:A:H8	1.84	0.43
1:A:269:LYS:HE2	1:A:269:LYS:HB3	1.83	0.43
2:B:98:G:H2'	2:B:99:C:C6	2.52	0.43
1:A:855:GLU:H	1:A:855:GLU:HG3	1.69	0.43
1:A:191:HIS:CE1	1:A:195:GLN:CG	3.00	0.42
1:A:446:ALA:O	1:A:450:GLY:HA2	2.19	0.42
1:A:500:ARG:NH1	1:A:500:ARG:HG2	2.32	0.42
1:A:490:GLY:C	1:A:1066:LEU:HD11	2.40	0.42
1:A:858:SER:O	1:A:927:GLN:N	2.41	0.42
1:A:13:LEU:HD13	1:A:491:VAL:HG11	2.01	0.42
1:A:26:MET:HB3	1:A:26:MET:HE2	1.90	0.42
1:A:266:LYS:NZ	1:A:422:GLN:NE2	2.67	0.42
1:A:837:LYS:HB2	1:A:837:LYS:HE3	1.74	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:949:ARG:HB3	1:A:964:ILE:HB	2.01	0.42
1:A:541:PHE:HA	1:A:542:PRO:HD3	1.69	0.42
2:B:45:A:OP2	2:B:45:A:H8	2.03	0.42
2:B:70:C:H2'	2:B:71:G:C8	2.55	0.42
2:B:102:C:H2'	2:B:103:U:C6	2.54	0.42
2:B:51:A:H2'	2:B:52:C:H6	1.85	0.42
1:A:14:GLY:O	1:A:24:TRP:HA	2.20	0.42
1:A:78:ARG:NE	3:A:1103:HOH:O	2.47	0.42
1:A:148:GLY:C	1:A:149:GLU:HG3	2.40	0.42
2:B:133:A:OP2	2:B:133:A:C3'	2.67	0.42
1:A:530:ARG:NH1	1:A:550:SER:OG	2.53	0.42
1:A:62:ARG:HD2	2:B:90:A:O4'	2.20	0.42
1:A:50:GLU:CD	1:A:516:ARG:HH22	2.24	0.42
1:A:492:VAL:HA	1:A:496:GLY:O	2.20	0.41
2:B:31:G:H2'	2:B:32:C:C6	2.54	0.41
1:A:227:VAL:O	1:A:227:VAL:CG1	2.66	0.41
2:B:25:G:H2'	2:B:26:U:H6	1.84	0.41
1:A:218:LYS:HB3	1:A:222:PHE:HE2	1.86	0.41
1:A:590:LEU:HD11	1:A:654:LEU:HD12	2.01	0.41
1:A:19:ILE:O	1:A:48:ARG:HB3	2.20	0.41
1:A:153:LYS:H	1:A:153:LYS:HG3	1.44	0.41
1:A:316:THR:HG22	1:A:318:ALA:H	1.86	0.41
1:A:623:PHE:O	1:A:632:TRP:HB2	2.20	0.41
1:A:847:VAL:HG22	1:A:935:VAL:HG21	2.02	0.41
1:A:109:THR:O	1:A:112:GLN:N	2.54	0.41
1:A:123:THR:HG22	1:A:125:LEU:H	1.85	0.41
1:A:662:PHE:O	1:A:666:ASN:ND2	2.53	0.41
1:A:869:LEU:HD11	1:A:892:LEU:CB	2.49	0.41
1:A:91:VAL:O	1:A:91:VAL:CG1	2.69	0.41
1:A:135:ILE:HD11	1:A:235:ILE:HG23	2.03	0.41
1:A:721:ARG:HH22	1:A:799:GLU:CD	2.24	0.41
2:B:133:A:P	2:B:133:A:C3'	3.09	0.41
1:A:87:LYS:O	1:A:92:LEU:HD23	2.21	0.41
1:A:88:ARG:C	1:A:89:GLU:CG	2.88	0.41
1:A:669:ASP:OD1	1:A:674:ASN:ND2	2.53	0.41
2:B:97:U:H2'	2:B:98:G:C8	2.54	0.41
1:A:639:VAL:HG21	1:A:653:LEU:HD11	2.03	0.41
1:A:785:ILE:HD12	1:A:785:ILE:HG23	1.61	0.41
1:A:481:LEU:HD22	1:A:677:LEU:HD11	2.02	0.41
1:A:542:PRO:CD	1:A:543:ASN:H	2.34	0.41
1:A:610[B]:GLU:OE2	1:A:622:TYR:OH	2.15	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1004:ASN:OD1	1:A:1033:ARG:NH1	2.54	0.40
1:A:1008:GLU:OE2	1:A:1015:ARG:NH2	2.44	0.40
1:A:11:TYR:CE2	1:A:498:PRO:HB3	2.56	0.40
1:A:633:GLN:HG2	1:A:633:GLN:H	1.74	0.40
1:A:65:ALA:O	1:A:69:ARG:HG3	2.21	0.40
1:A:305:LEU:HA	1:A:305:LEU:HD23	1.83	0.40
1:A:315:LEU:HD23	1:A:347:LEU:HD23	2.02	0.40
1:A:654:LEU:HB3	1:A:657:PHE:HD2	1.86	0.40
1:A:1069:GLU:OE2	1:A:1071:ARG:NH2	2.41	0.40
2:B:128:C:H2'	2:B:129:G:H8	1.86	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)
1:A:319:GLN:OE1	1:A:915:ASN:ND2[3_455]	1.99	0.21

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
1	A	1066/1092 (98%)	1003 (94%)	56 (5%)	7 (1%)	22 56

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	93	GLN
1	A	109	THR
1	A	196	ARG
1	A	95	ALA
1	A	151	ALA
1	A	751	ASN

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Mol	Chain	Res	Type
1	A	110	PRO

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	840/942 (89%)	807 (96%)	33 (4%)	32 65

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	GLU
1	A	36	ILE
1	A	41	LEU
1	A	92	LEU
1	A	108	ASN
1	A	109	THR
1	A	131	LEU
1	A	207	ASP
1	A	222	PHE
1	A	226	HIS
1	A	290	GLU
1	A	312	LYS
1	A	366	LYS
1	A	373	ASN
1	A	392	ASP
1	A	430	ARG
1	A	449	TYR
1	A	540	TYR
1	A	575	LEU
1	A	612	GLN
1	A	646	ARG
1	A	679	GLN
1	A	692	LYS
1	A	714	LYS
1	A	754	ASP

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Mol	Chain	Res	Type
1	A	771	HIS
1	A	796	GLU
1	A	797	PHE
1	A	870	LYS
1	A	880	ARG
1	A	927	GLN
1	A	994	SER
1	A	1064	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	191	HIS
1	A	226	HIS
1	A	319	GLN
1	A	354	HIS
1	A	373	ASN
1	A	739	GLN
1	A	1059	GLN

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	116/135 (85%)	34 (29%)	7 (6%)

All (34) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	20	U
2	B	21	U
2	B	24	C
2	B	31	G
2	B	33	U
2	B	34	C
2	B	38	U
2	B	39	U
2	B	40	C
2	B	43	G
2	B	46	A
2	B	47	G

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Mol	Chain	Res	Type
2	B	48	A
2	B	65	A
2	B	75	G
2	B	76	A
2	B	77	A
2	B	78	A
2	B	79	A
2	B	80	G
2	B	83	G
2	B	84	U
2	B	85	G
2	B	89	C
2	B	95	U
2	B	107	A
2	B	108	G
2	B	109	C
2	B	110	U
2	B	111	C
2	B	115	C
2	B	123	G
2	B	132	U
2	B	133	A

All (7) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	75	G
2	B	76	A
2	B	77	A
2	B	79	A
2	B	83	G
2	B	114	G
2	B	132	U

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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	A	1069/1092 (97%)	0.17	18 (1%) 70 53	11, 47, 76, 95	0
2	B	117/135 (86%)	0.83	17 (14%) 2 1	17, 57, 126, 156	0
All	All	1186/1227 (96%)	0.23	35 (2%) 50 34	11, 48, 80, 156	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	111	C	5.6
2	B	114	G	5.3
1	A	914	GLY	4.8
2	B	17	A	4.0
2	B	77	A	3.2
2	B	78	A	3.1
2	B	16	U	3.0
1	A	913	ALA	2.8
2	B	76	A	2.8
2	B	110	U	2.8
1	A	225	PRO	2.8
2	B	18	A	2.8
2	B	40	C	2.8
1	A	911	ASP	2.7
1	A	176	THR	2.7
2	B	133	A	2.7
2	B	47	G	2.6
1	A	900	ALA	2.5
2	B	39	U	2.4
1	A	451	ASP	2.4
1	A	226	HIS	2.4
2	B	108	G	2.4
2	B	79	A	2.4
2	B	15	U	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	883	LYS	2.3
1	A	761	GLU	2.3
1	A	631	GLU	2.2
1	A	103	ILE	2.2
1	A	257	HIS	2.2
1	A	530	ARG	2.1
1	A	912	LYS	2.1
1	A	97	PHE	2.1
1	A	910	TYR	2.1
2	B	45	A	2.0
1	A	765	VAL	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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.