



Full wwPDB/EMDatabank EM Map/Model Validation Report ⓘ

Dec 11, 2019 – 03:36 AM EST

PDB ID : 6DBI
EMDB ID: : EMD-7843
Title : Cryo-EM structure of RAG in complex with 12-RSS and 23-RSS nicked DNA intermediates
Authors : Wu, H.; Liao, M.; Ru, H.; Mi, W.
Deposited on : 2018-05-03
Resolution : 3.36 Å(reported)

This is a Full wwPDB/EMDatabank EM Map/Model Validation Report
for a publicly released PDB/EMDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

MolProbity : 4.02b-467
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : 2.4

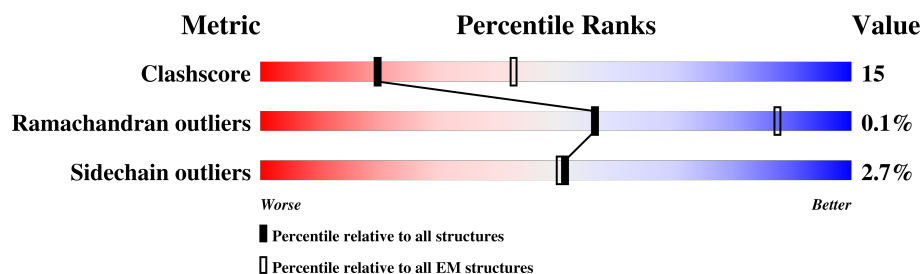
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.36 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)	EM structures (#Entries)
Clashscore	136327	1886
Ramachandran outliers	132723	1663
Sidechain outliers	132532	1531

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for ≥ 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	1159	38% 15% . 46%
1	C	1159	39% 14% . 46%
2	B	533	40% 25% . 34%
2	D	533	45% 20% . 34%
3	E	34	29% 68% .
4	F	50	40% 60%
5	G	61	46% 54%
6	H	45	58% 42%
7	I	16	44% 56%

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Mol	Chain	Length	Quality of chain
7	J	16	<div><div></div><div>50%</div><div></div><div>50%</div></div>

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 20001 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Recombination activating gene 1 - MBP chimera.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	C	622	Total	C	N	O	S	0	0
			5008	3134	901	936	37		
1	A	622	Total	C	N	O	S	0	0
			5008	3134	901	936	37		

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-127	MET	-	initiating methionine	UNP P0AEX9
C	-126	GLY	-	expression tag	UNP P0AEX9
C	-125	SER	-	expression tag	UNP P0AEX9
C	-124	SER	-	expression tag	UNP P0AEX9
C	-123	HIS	-	expression tag	UNP P0AEX9
C	-122	HIS	-	expression tag	UNP P0AEX9
C	-121	HIS	-	expression tag	UNP P0AEX9
C	-120	HIS	-	expression tag	UNP P0AEX9
C	-119	HIS	-	expression tag	UNP P0AEX9
C	-118	HIS	-	expression tag	UNP P0AEX9
C	-117	GLY	-	expression tag	UNP P0AEX9
C	-116	THR	-	expression tag	UNP P0AEX9
C	-115	LYS	-	expression tag	UNP P0AEX9
C	-114	THR	-	expression tag	UNP P0AEX9
C	251	GLY	-	linker	UNP P0AEX9
C	252	THR	-	linker	UNP P0AEX9
C	253	ASP	-	linker	UNP P0AEX9
C	254	TYR	-	linker	UNP P0AEX9
C	255	ASP	-	linker	UNP P0AEX9
C	256	ILE	-	linker	UNP P0AEX9
C	257	PRO	-	linker	UNP P0AEX9
C	258	THR	-	linker	UNP P0AEX9
C	259	THR	-	linker	UNP P0AEX9
C	260	LEU	-	linker	UNP P0AEX9
C	261	GLU	-	linker	UNP P0AEX9
C	262	VAL	-	linker	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	263	LEU	-	linker	UNP P0AEX9
C	264	PHE	-	linker	UNP P0AEX9
C	265	GLN	-	linker	UNP P0AEX9
C	266	GLY	-	linker	UNP P0AEX9
C	267	PRO	-	linker	UNP P0AEX9
C	268	LEU	-	linker	UNP P0AEX9
C	269	GLY	-	linker	UNP P0AEX9
C	270	SER	-	linker	UNP P0AEX9
A	-127	MET	-	initiating methionine	UNP P0AEX9
A	-126	GLY	-	expression tag	UNP P0AEX9
A	-125	SER	-	expression tag	UNP P0AEX9
A	-124	SER	-	expression tag	UNP P0AEX9
A	-123	HIS	-	expression tag	UNP P0AEX9
A	-122	HIS	-	expression tag	UNP P0AEX9
A	-121	HIS	-	expression tag	UNP P0AEX9
A	-120	HIS	-	expression tag	UNP P0AEX9
A	-119	HIS	-	expression tag	UNP P0AEX9
A	-118	HIS	-	expression tag	UNP P0AEX9
A	-117	GLY	-	expression tag	UNP P0AEX9
A	-116	THR	-	expression tag	UNP P0AEX9
A	-115	LYS	-	expression tag	UNP P0AEX9
A	-114	THR	-	expression tag	UNP P0AEX9
A	251	GLY	-	linker	UNP P0AEX9
A	252	THR	-	linker	UNP P0AEX9
A	253	ASP	-	linker	UNP P0AEX9
A	254	TYR	-	linker	UNP P0AEX9
A	255	ASP	-	linker	UNP P0AEX9
A	256	ILE	-	linker	UNP P0AEX9
A	257	PRO	-	linker	UNP P0AEX9
A	258	THR	-	linker	UNP P0AEX9
A	259	THR	-	linker	UNP P0AEX9
A	260	LEU	-	linker	UNP P0AEX9
A	261	GLU	-	linker	UNP P0AEX9
A	262	VAL	-	linker	UNP P0AEX9
A	263	LEU	-	linker	UNP P0AEX9
A	264	PHE	-	linker	UNP P0AEX9
A	265	GLN	-	linker	UNP P0AEX9
A	266	GLY	-	linker	UNP P0AEX9
A	267	PRO	-	linker	UNP P0AEX9
A	268	LEU	-	linker	UNP P0AEX9
A	269	GLY	-	linker	UNP P0AEX9
A	270	SER	-	linker	UNP P0AEX9

- Molecule 2 is a protein called Recombination activating gene 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	351	Total	C	N	O	S	0	0
			2714	1716	470	509	19		
2	B	351	Total	C	N	O	S	0	0
			2714	1716	470	509	19		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	-2	GLY	-	expression tag	UNP Q1RLW7
D	-1	GLY	-	expression tag	UNP Q1RLW7
D	0	SER	-	expression tag	UNP Q1RLW7
B	-2	GLY	-	expression tag	UNP Q1RLW7
B	-1	GLY	-	expression tag	UNP Q1RLW7
B	0	SER	-	expression tag	UNP Q1RLW7

- Molecule 3 is a DNA chain called Forward strand of 12-RSS signal end.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	34	Total	C	N	O	P	0	0
			697	330	138	195	34		

- Molecule 4 is a DNA chain called Reverse strand of 12-RSS.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	F	50	Total	C	N	O	P	0	0
			1027	489	183	305	50		

- Molecule 5 is a DNA chain called Reverse strand of 23-RSS.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	G	61	Total	C	N	O	P	0	0
			1256	596	235	364	61		

- Molecule 6 is a DNA chain called Forward strand of 23-RSS signal end.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	H	45	Total	C	N	O	P	0	0
			919	437	169	268	45		

- Molecule 7 is a DNA chain called Forward strand of coding flank.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	I	16	Total	C	N	O	P	0	0
			326	156	54	100	16		
7	J	16	Total	C	N	O	P	0	0
			326	156	54	100	16		

- Molecule 8 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
8	A	1	Total	Zn	0
			1	1	
8	C	1	Total	Zn	0
			1	1	

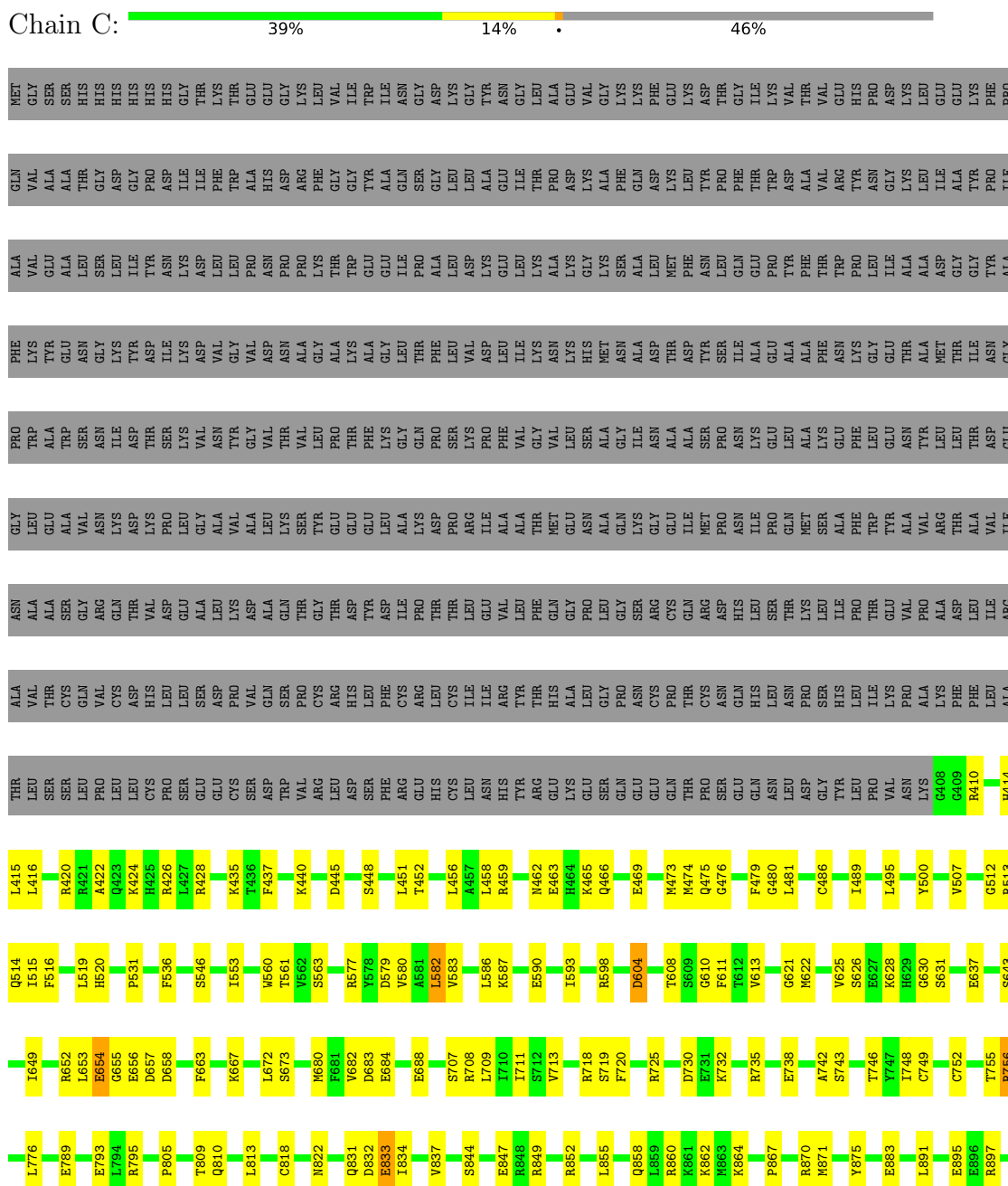
- Molecule 9 is CALCIUM ION (three-letter code: CA) (formula: Ca).

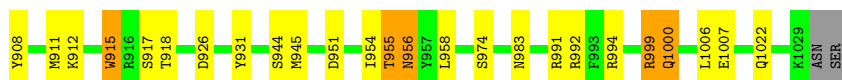
Mol	Chain	Residues	Atoms		AltConf
9	A	2	Total	Ca	0
			2	2	
9	C	2	Total	Ca	0
			2	2	

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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. 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. 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.

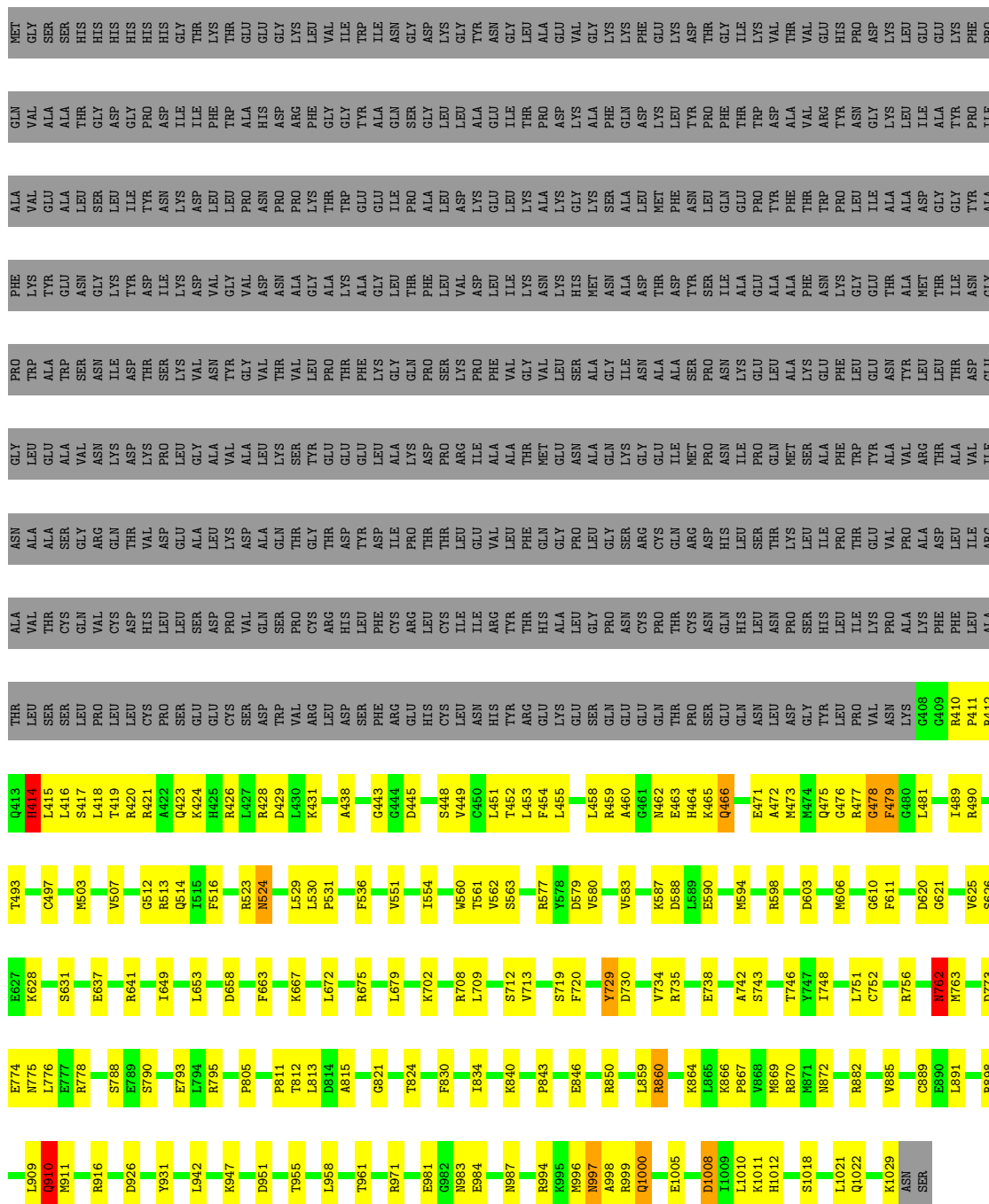
- Molecule 1: Recombination activating gene 1 - MBP chimera





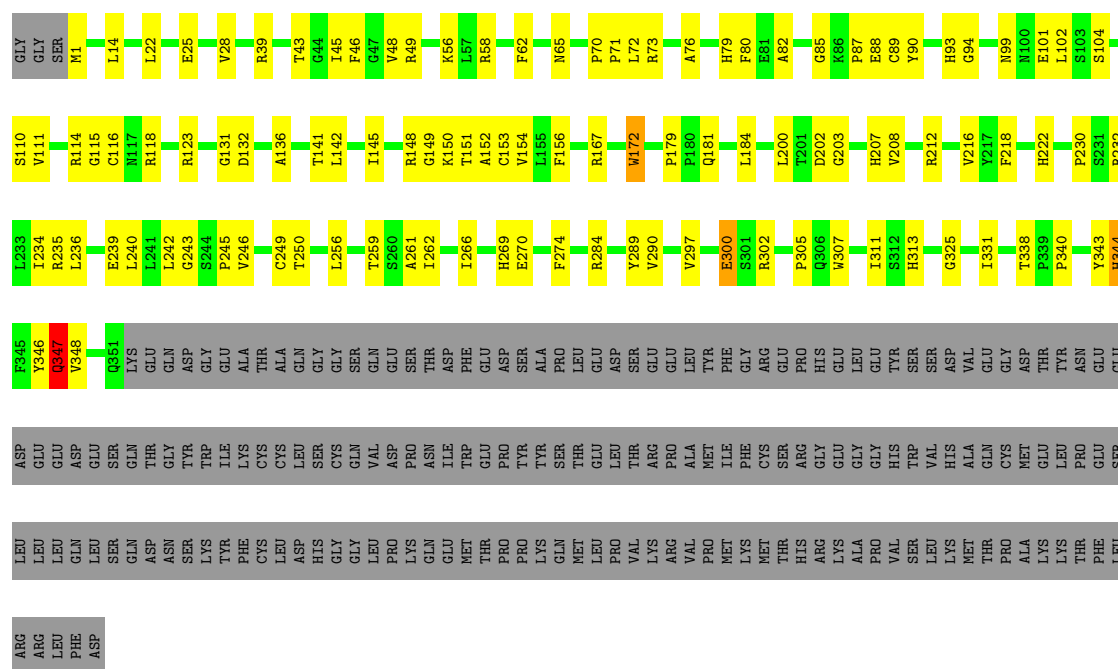
- Molecule 1: Recombination activating gene 1 - MBP chimera

Chain A: 38% 15% 46%



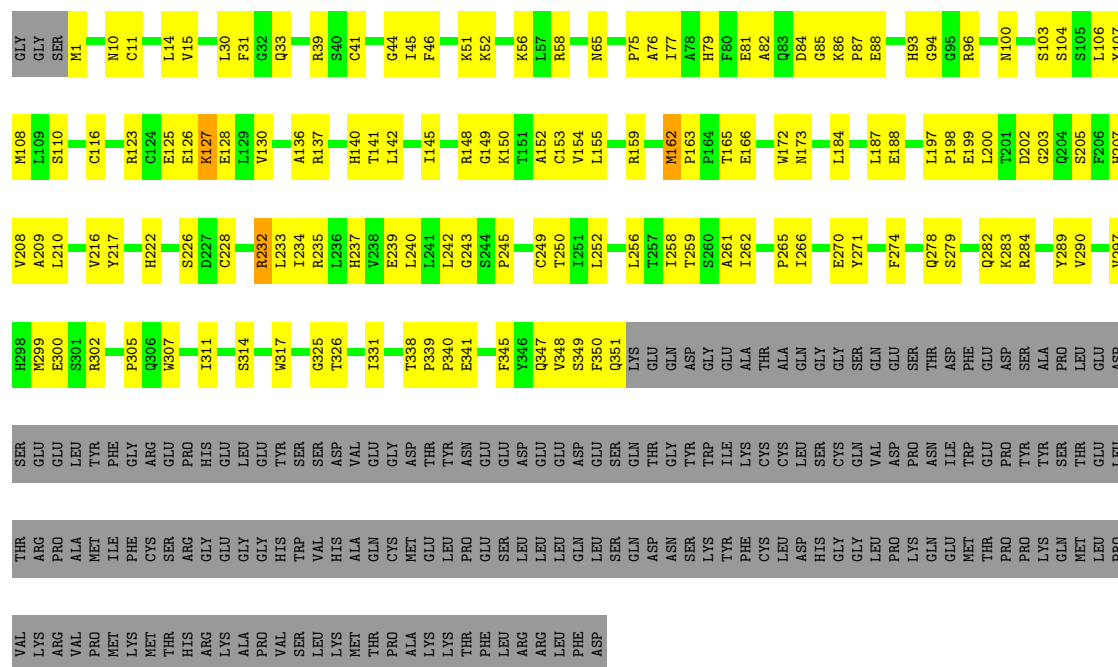
- Molecule 2: Recombination activating gene 2

Chain D:  45% 20% . 34%



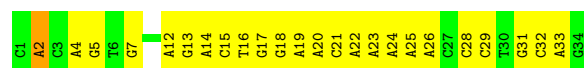
• Molecule 2: Recombination activating gene 2

Chain B: 40% 25% 34%



• Molecule 3: Forward strand of 12-RSS signal end

Chain E: 29% 68%



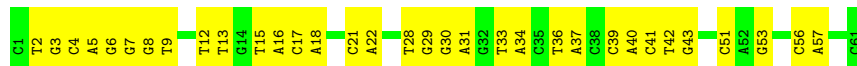
- Molecule 4: Reverse strand of 12-RSS

Chain F:  40% 60%



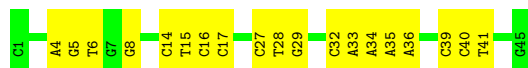
- Molecule 5: Reverse strand of 23-RSS

Chain G:  46% 54%



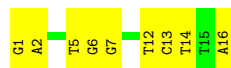
- Molecule 6: Forward strand of 23-RSS signal end

Chain H:  58% 42%



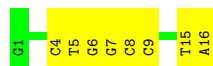
- Molecule 7: Forward strand of coding flank

Chain I:  44% 56%



- Molecule 7: Forward strand of coding flank

Chain J:  50% 50%



4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	53109	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	Not provided	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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 > 2$	RMSZ	# $ Z > 2$
1	A	0.60	6/5106 (0.1%)	0.87	17/6867 (0.2%)
1	C	0.38	2/5106 (0.0%)	0.65	9/6867 (0.1%)
2	B	0.38	1/2784 (0.0%)	0.77	6/3784 (0.2%)
2	D	0.52	4/2784 (0.1%)	0.81	14/3784 (0.4%)
3	E	0.70	0/784	0.90	1/1206 (0.1%)
4	F	0.77	0/1150	1.02	0/1774
5	G	0.66	0/1410	0.94	1/2175 (0.0%)
6	H	0.61	0/1030	0.92	0/1586
7	I	0.75	0/363	1.07	0/558
7	J	0.65	0/363	1.03	0/558
All	All	0.54	13/20880 (0.1%)	0.83	48/29159 (0.2%)

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	5
2	D	0	2
All	All	0	7

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	910	GLN	CD-OE1	-21.19	0.77	1.24
1	A	910	GLN	CD-NE2	17.51	1.76	1.32
2	D	347	GLN	CD-OE1	-14.75	0.91	1.24
1	A	762	ASN	CB-CG	11.73	1.78	1.51
2	D	347	GLN	CD-NE2	9.09	1.55	1.32
1	A	910	GLN	CG-CD	9.07	1.72	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	956	ASN	CG-ND2	8.71	1.54	1.32
1	A	762	ASN	CG-OD1	-7.97	1.06	1.24
1	C	956	ASN	CG-OD1	-6.79	1.09	1.24
1	A	762	ASN	CG-ND2	6.32	1.48	1.32
2	D	344	HIS	CG-CD2	6.17	1.46	1.35
2	D	347	GLN	CG-CD	5.42	1.63	1.51
2	B	173	ASN	CG-ND2	5.04	1.45	1.32

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	524	ASN	CB-CG-OD1	29.19	179.98	121.60
2	B	173	ASN	CB-CG-OD1	27.53	176.65	121.60
1	A	910	GLN	CA-CB-CG	26.34	171.34	113.40
1	C	956	ASN	CB-CG-OD1	24.11	169.82	121.60
1	A	762	ASN	CB-CA-C	-21.01	68.39	110.40
2	D	344	HIS	ND1-CG-CD2	-19.11	79.25	106.00
1	A	910	GLN	OE1-CD-NE2	-15.06	87.27	121.90
2	D	344	HIS	CG-ND1-CE1	14.56	128.59	108.20
1	A	524	ASN	CB-CG-ND2	-14.29	82.40	116.70
2	B	173	ASN	CB-CG-ND2	-14.28	82.44	116.70
1	C	956	ASN	CB-CG-ND2	-14.07	82.92	116.70
1	C	956	ASN	OD1-CG-ND2	-13.94	89.84	121.90
2	D	344	HIS	N-CA-CB	12.89	133.79	110.60
2	D	344	HIS	CB-CG-CD2	-12.79	91.14	130.80
2	D	344	HIS	CB-CG-ND1	12.54	154.56	123.20
1	A	762	ASN	CB-CG-ND2	-11.86	88.23	116.70
2	B	173	ASN	OD1-CG-ND2	-11.51	95.42	121.90
1	A	524	ASN	OD1-CG-ND2	-10.56	97.62	121.90
1	A	762	ASN	CB-CG-OD1	10.32	142.24	121.60
2	D	347	GLN	OE1-CD-NE2	-9.06	101.06	121.90
1	A	910	GLN	CG-CD-NE2	-8.12	97.21	116.70
1	A	762	ASN	N-CA-CB	8.01	125.02	110.60
2	D	344	HIS	N-CA-C	-7.99	89.44	111.00
2	D	347	GLN	CA-CB-CG	-7.68	96.50	113.40
2	B	172	TRP	C-N-CA	-7.48	103.00	121.70
2	D	344	HIS	ND1-CE1-NE2	-7.19	94.09	109.90
5	G	51	DC	O5'-P-OP2	-7.01	99.39	105.70
1	A	910	GLN	CB-CG-CD	-6.83	93.83	111.60
1	A	762	ASN	OD1-CG-ND2	-6.79	106.29	121.90
2	B	173	ASN	N-CA-CB	6.64	122.55	110.60
1	A	524	ASN	CA-CB-CG	-6.19	99.79	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	909	LEU	C-N-CA	-6.13	106.39	121.70
1	A	910	GLN	CG-CD-OE1	6.00	133.61	121.60
1	A	524	ASN	N-CA-C	5.98	127.15	111.00
2	B	173	ASN	CA-CB-CG	-5.97	100.26	113.40
2	D	344	HIS	CB-CA-C	-5.91	98.58	110.40
1	C	956	ASN	N-CA-C	-5.89	95.09	111.00
2	D	343	TYR	C-N-CA	-5.88	107.01	121.70
1	C	956	ASN	CA-CB-CG	-5.82	100.61	113.40
2	D	347	GLN	N-CA-CB	5.66	120.78	110.60
1	C	955	THR	C-N-CA	-5.59	107.73	121.70
2	D	346	TYR	C-N-CA	5.48	135.40	121.70
1	C	954	ILE	CA-CB-CG1	5.35	121.16	111.00
1	A	414	HIS	N-CA-C	-5.25	96.83	111.00
1	C	956	ASN	N-CA-CB	5.17	119.90	110.60
1	C	604	ASP	CB-CG-OD1	-5.09	113.72	118.30
3	E	2	DA	O4'-C4'-C3'	-5.07	102.47	104.50
2	D	236	LEU	CA-CB-CG	5.01	126.83	115.30

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	478	GLY	Peptide
1	A	523	ARG	Peptide
1	A	762	ASN	Sidechain,Mainchain
1	A	910	GLN	Sidechain
2	D	344	HIS	Sidechain
2	D	347	GLN	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	A	5008	0	4967	189	0
1	C	5008	0	4967	153	0
2	B	2714	0	2665	97	0
2	D	2714	0	2664	70	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	697	0	379	32	0
4	F	1027	0	566	42	0
5	G	1256	0	686	34	0
6	H	919	0	506	23	0
7	I	326	0	182	7	0
7	J	326	0	183	10	0
8	A	1	0	0	0	0
8	C	1	0	0	0	0
9	A	2	0	0	0	0
9	C	2	0	0	0	0
All	All	20001	0	17765	565	0

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 15.

All (565) 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:762:ASN:CB	1:A:762:ASN:CG	1.78	1.50
1:A:910:GLN:NE2	1:A:910:GLN:CD	1.76	1.37
1:C:448:SER:CB	1:A:415:LEU:CD1	2.02	1.35
1:C:448:SER:CB	1:A:415:LEU:HD11	1.58	1.32
1:A:458:LEU:HB3	1:A:462:ASN:OD1	1.41	1.19
1:A:462:ASN:HB3	1:A:466:GLN:CG	1.76	1.16
1:C:831:GLN:OE1	1:C:852:ARG:CB	2.00	1.08
1:C:448:SER:OG	1:A:415:LEU:HD11	1.53	1.07
1:C:448:SER:HB2	1:A:415:LEU:CD1	1.76	1.06
1:C:831:GLN:OE1	1:C:852:ARG:HB2	1.57	1.05
1:A:462:ASN:CB	1:A:466:GLN:HG2	1.86	1.05
1:A:910:GLN:OE1	1:A:910:GLN:NE2	1.89	1.04
1:C:448:SER:HB2	1:A:415:LEU:HD12	1.37	1.01
1:C:424:LYS:NZ	4:F:14:DG:N7	2.08	1.01
1:C:448:SER:HB3	1:A:415:LEU:HD11	1.43	0.99
2:B:108:MET:SD	2:B:127:LYS:NZ	2.36	0.97
1:A:762:ASN:CB	1:A:762:ASN:ND2	2.28	0.97
1:A:998:ALA:HA	1:A:1008:ASP:OD1	1.64	0.96
1:A:910:GLN:OE1	1:A:910:GLN:CD	0.77	0.96
2:B:86:LYS:NZ	2:B:110:SER:OG	1.99	0.95
1:A:462:ASN:HB3	1:A:466:GLN:HG2	0.96	0.94
1:A:459:ARG:NH1	3:E:16:DT:OP1	2.01	0.94
2:B:79:HIS:NE2	2:B:88:GLU:OE2	2.03	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:448:SER:CB	1:A:415:LEU:HD12	1.93	0.92
1:C:831:GLN:OE1	1:C:852:ARG:HG3	1.70	0.91
1:C:831:GLN:OE1	1:C:852:ARG:CG	2.18	0.91
3:E:15:DC:O2	4:F:20:DG:N2	2.04	0.90
1:A:462:ASN:CB	1:A:466:GLN:CG	2.47	0.87
1:C:513:ARG:HH11	1:A:513:ARG:HE	1.20	0.87
1:C:448:SER:HB3	1:A:415:LEU:CD1	2.00	0.87
1:C:448:SER:OG	1:A:415:LEU:CD1	2.14	0.86
1:A:860:ARG:HH21	1:A:866:LYS:NZ	1.75	0.85
1:A:462:ASN:CG	1:A:466:GLN:HB3	1.98	0.84
3:E:18:DG:H1	4:F:17:DC:H42	1.24	0.83
2:B:76:ALA:HB3	2:B:93:HIS:O	1.77	0.83
1:A:860:ARG:HH21	1:A:866:LYS:HZ1	1.24	0.83
2:D:76:ALA:HB3	2:D:93:HIS:O	1.78	0.83
1:C:598:ARG:NH1	1:C:604:ASP:OD1	2.10	0.83
1:A:458:LEU:CB	1:A:462:ASN:OD1	2.27	0.82
3:E:15:DC:N3	4:F:20:DG:N1	2.26	0.82
3:E:21:DC:O2	4:F:14:DG:N2	2.11	0.82
1:A:996:MET:HB2	1:A:997:ASN:OD1	1.80	0.81
2:B:284:ARG:HB2	2:B:284:ARG:HH11	1.45	0.81
1:C:513:ARG:NH1	1:A:513:ARG:HE	1.78	0.80
1:A:846:GLU:OE2	1:A:850:ARG:NH2	2.12	0.79
1:A:420:ARG:NH1	6:H:28:DT:OP2	2.16	0.79
1:C:1000:GLN:HG2	4:F:31:DT:O4'	1.81	0.79
2:B:284:ARG:NH1	2:B:284:ARG:HB2	1.98	0.78
1:C:479:PHE:N	1:C:480:GLY:O	2.16	0.78
2:B:58:ARG:HH12	7:I:7:DG:H5''	1.46	0.78
1:A:462:ASN:HB2	1:A:466:GLN:NE2	1.99	0.78
2:D:142:LEU:HD11	2:D:153:CYS:HB3	1.66	0.78
1:A:997:ASN:HD22	1:A:1012:HIS:CE1	2.01	0.77
1:A:910:GLN:OE1	1:A:910:GLN:CG	2.31	0.77
2:B:258:ILE:HD13	2:B:284:ARG:NH1	2.00	0.76
1:A:997:ASN:OD1	1:A:997:ASN:N	2.19	0.75
1:C:459:ARG:HG2	6:H:27:DC:H5'	1.69	0.75
1:A:420:ARG:NH1	6:H:27:DC:H5''	2.03	0.74
1:C:608:THR:HB	1:C:718:ARG:HG2	1.70	0.73
1:C:426:ARG:NH1	4:F:12:DT:OP1	2.21	0.73
1:A:641:ARG:NH1	1:A:983:ASN:OD1	2.22	0.73
1:C:415:LEU:HD11	1:A:452:THR:HG21	1.69	0.73
1:A:788:SER:HB2	2:B:65:ASN:HA	1.69	0.73
1:C:473:MET:HG3	1:C:474:MET:HG2	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:ARG:HA	1:A:431:LYS:HZ3	1.55	0.72
1:A:426:ARG:NH2	5:G:12:DT:OP1	2.20	0.72
1:A:860:ARG:HE	1:A:866:LYS:HE3	1.54	0.71
2:B:283:LYS:NZ	2:B:314:SER:O	2.21	0.71
1:A:776:LEU:HD23	1:A:805:PRO:HD2	1.73	0.70
1:C:688:GLU:OE1	2:D:73:ARG:NH1	2.24	0.70
1:A:1011:LYS:NZ	6:H:8:DG:OP1	2.24	0.70
1:A:997:ASN:ND2	1:A:1012:HIS:CE1	2.60	0.70
1:A:478:GLY:HA2	1:A:479:PHE:O	1.90	0.70
1:A:762:ASN:CA	1:A:762:ASN:ND2	2.54	0.70
2:B:58:ARG:NH1	7:I:7:DG:H5''	2.06	0.70
5:G:36:DT:H2'	5:G:37:DA:C8	2.28	0.69
2:D:58:ARG:HH12	7:J:7:DG:H5''	1.56	0.69
5:G:53:DG:N2	7:J:9:DC:O2	2.18	0.69
2:D:49:ARG:HH21	2:D:58:ARG:HD2	1.57	0.69
2:B:232:ARG:HE	2:B:234:ILE:HD11	1.56	0.68
1:A:424:LYS:NZ	1:A:431:LYS:HE3	2.08	0.68
2:B:270:GLU:HB3	2:B:289:TYR:HE1	1.59	0.67
1:C:630:GLY:O	4:F:30:DC:OP1	2.12	0.67
5:G:2:DT:H2'	5:G:3:DG:C8	2.30	0.67
1:C:643:SER:OG	1:C:983:ASN:ND2	2.26	0.67
2:B:165:THR:OG1	2:B:166:GLU:OE2	2.10	0.67
2:D:1:MET:HB3	2:D:305:PRO:HG3	1.76	0.67
1:A:610:GLY:HA3	1:A:653:LEU:HG	1.76	0.67
5:G:2:DT:H2'	5:G:3:DG:H8	1.59	0.66
1:A:452:THR:HA	1:A:455:LEU:HD13	1.77	0.66
1:A:462:ASN:CB	1:A:466:GLN:CD	2.63	0.66
2:B:142:LEU:HD11	2:B:153:CYS:HB3	1.76	0.66
1:A:981:GLU:OE2	4:F:34:DG:N2	2.28	0.66
3:E:17:DG:H2''	3:E:18:DG:N7	2.10	0.66
1:C:631:SER:HA	1:A:864:LYS:HD2	1.77	0.66
1:C:897:ARG:HH11	1:C:945:MET:HE2	1.59	0.66
1:A:418:LEU:HD12	1:A:426:ARG:HH22	1.60	0.66
1:A:735:ARG:NH2	1:A:748:ILE:O	2.28	0.65
2:D:338:THR:HG22	2:D:340:PRO:HA	1.78	0.65
1:C:748:ILE:HD12	1:C:749:CYS:N	2.11	0.65
2:B:258:ILE:HD13	2:B:284:ARG:HH11	1.60	0.65
2:B:197:LEU:HD13	2:B:200:LEU:HD13	1.78	0.65
2:D:222:HIS:HB2	2:D:259:THR:HG21	1.79	0.65
1:A:774:GLU:OE2	1:A:778:ARG:NH2	2.30	0.64
1:C:463:GLU:HA	5:G:22:DA:H4'	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:H:40:DC:H2'	6:H:41:DT:C6	2.31	0.64
1:C:598:ARG:NH1	1:C:604:ASP:CG	2.50	0.64
1:C:621:GLY:HA3	1:C:684:GLU:OE2	1.97	0.64
1:A:462:ASN:CB	1:A:466:GLN:NE2	2.61	0.64
1:A:415:LEU:O	1:A:415:LEU:HG	1.97	0.64
2:B:217:TYR:OH	2:B:235:ARG:NH2	2.30	0.64
1:C:832:ASP:HB3	1:C:837:VAL:HG11	1.80	0.64
2:D:167:ARG:NH2	2:D:172:TRP:O	2.31	0.64
2:D:230:PRO:HB2	2:D:232:ARG:HG2	1.80	0.64
3:E:12:DA:H2'	3:E:13:DG:C8	2.32	0.63
5:G:33:DT:O4	5:G:34:DA:N6	2.31	0.63
1:A:840:LYS:HE3	1:A:843:PRO:HA	1.80	0.63
1:A:910:GLN:NE2	1:A:910:GLN:CG	2.61	0.63
2:B:226:SER:HG	2:B:228:CYS:HG	1.46	0.63
1:C:598:ARG:HH12	1:C:604:ASP:CG	2.00	0.63
1:C:422:ALA:HB1	4:F:12:DT:H5'	1.81	0.63
1:A:529:LEU:HD13	1:A:1021:LEU:HD11	1.81	0.62
2:B:15:VAL:HG21	2:B:345:PHE:HZ	1.64	0.62
1:A:462:ASN:HB2	1:A:466:GLN:CD	2.20	0.62
1:C:512:GLY:HA2	1:C:513:ARG:HB3	1.81	0.62
1:C:864:LYS:HD2	1:A:631:SER:HA	1.79	0.62
1:C:738:GLU:OE2	1:C:809:THR:OG1	2.17	0.62
1:C:951:ASP:N	1:C:951:ASP:OD1	2.33	0.62
1:A:462:ASN:CG	1:A:466:GLN:CB	2.68	0.61
2:D:266:ILE:HG22	2:D:325:GLY:HA2	1.82	0.61
4:F:6:DG:H2''	4:F:7:DG:H8	1.65	0.61
1:C:656:GLU:HG2	1:C:657:ASP:H	1.64	0.61
1:A:762:ASN:HD22	1:A:763:MET:N	1.99	0.60
5:G:8:DG:H2'	5:G:9:DT:C6	2.36	0.60
1:A:672:LEU:HB3	1:A:1022:GLN:HE22	1.66	0.60
1:C:622:MET:SD	1:C:991:ARG:NH2	2.74	0.60
1:C:749:CYS:HB3	1:C:752:CYS:SG	2.41	0.60
2:D:184:LEU:HD13	2:D:245:PRO:HB2	1.83	0.60
5:G:53:DG:N1	7:J:9:DC:N3	2.39	0.60
1:A:762:ASN:CA	1:A:762:ASN:CG	2.64	0.60
2:B:159:ARG:NH1	2:B:205:SER:OG	2.34	0.60
1:A:911:MET:HB2	1:A:931:TYR:CE2	2.36	0.60
2:B:123:ARG:NH1	2:B:125:GLU:OE2	2.35	0.59
1:C:649:ILE:HG23	1:C:663:PHE:HB3	1.84	0.59
4:F:18:DC:H2'	4:F:19:DA:C8	2.38	0.59
2:B:150:LYS:NZ	2:B:240:LEU:HB3	2.18	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:290:VAL:HG13	2:B:297:VAL:HG23	1.84	0.59
4:F:31:DT:H2''	4:F:32:DG:C8	2.38	0.59
1:A:1000:GLN:HG2	5:G:42:DT:O4'	2.03	0.59
1:C:658:ASP:OD1	1:C:658:ASP:N	2.34	0.59
6:H:32:DC:H2''	6:H:33:DA:H5''	1.83	0.59
1:A:455:LEU:O	1:A:459:ARG:HB2	2.03	0.59
1:A:473:MET:O	1:A:477:ARG:NH1	2.36	0.59
4:F:9:DT:H2'	4:F:10:DT:C6	2.38	0.59
1:A:419:THR:HG22	1:A:420:ARG:H	1.68	0.59
1:A:410:ARG:O	1:A:412:ARG:NH1	2.37	0.58
1:C:513:ARG:N	1:A:471:GLU:OE2	2.33	0.58
2:D:208:VAL:HG21	2:D:261:ALA:HB3	1.85	0.58
5:G:17:DC:H2'	5:G:18:DA:C8	2.39	0.57
1:C:590:GLU:HA	1:C:593:ILE:HG22	1.85	0.57
1:A:462:ASN:OD1	1:A:466:GLN:HB3	2.04	0.57
2:B:274:PHE:CZ	2:B:348:VAL:HG21	2.39	0.57
1:A:420:ARG:HH12	6:H:28:DT:P	2.27	0.57
1:C:451:LEU:HA	1:A:454:PHE:HZ	1.70	0.57
2:D:270:GLU:HB3	2:D:289:TYR:HE1	1.70	0.57
1:A:410:ARG:HG2	1:A:411:PRO:HD2	1.87	0.57
2:B:338:THR:HG22	2:B:340:PRO:HA	1.85	0.57
1:A:551:VAL:HG21	1:A:577:ARG:HD3	1.85	0.57
2:B:148:ARG:NH2	2:B:240:LEU:O	2.29	0.57
1:C:546:SER:OG	1:C:577:ARG:NH1	2.37	0.57
1:A:729:TYR:HB2	1:A:734:VAL:HG23	1.86	0.57
2:D:232:ARG:HD2	2:D:234:ILE:HD11	1.85	0.57
1:A:472:ALA:C	1:A:477:ARG:HH12	2.09	0.56
2:D:152:ALA:HB2	2:D:240:LEU:HD21	1.88	0.56
2:B:307:TRP:HB3	2:B:311:ILE:HG23	1.88	0.56
1:C:831:GLN:OE1	1:C:852:ARG:CA	2.54	0.56
2:B:232:ARG:NE	2:B:234:ILE:HD11	2.20	0.56
1:C:667:LYS:NZ	6:H:4:DA:OP2	2.29	0.56
1:A:423:GLN:NE2	1:A:426:ARG:HH11	2.04	0.56
1:A:984:GLU:HB2	4:F:34:DG:H21	1.69	0.55
2:B:1:MET:HA	2:B:349:SER:O	2.06	0.55
2:D:102:LEU:HB3	2:D:136:ALA:HB1	1.88	0.55
2:D:249:CYS:SG	2:D:250:THR:N	2.79	0.55
5:G:41:DC:H2''	5:G:42:DT:H71	1.88	0.55
1:A:812:THR:OG1	1:A:813:LEU:N	2.40	0.55
1:C:445:ASP:OD2	1:A:415:LEU:HD21	2.07	0.55
2:B:233:LEU:HB3	2:B:252:LEU:HB3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:554:ILE:HD12	1:A:577:ARG:HG3	1.88	0.55
1:A:649:ILE:HG23	1:A:663:PHE:HB3	1.89	0.55
1:A:561:THR:O	1:A:563:SER:N	2.39	0.54
1:A:658:ASP:N	1:A:658:ASP:OD1	2.40	0.54
1:A:889:CYS:HB2	1:A:898:ARG:HG3	1.88	0.54
1:A:996:MET:CB	1:A:997:ASN:OD1	2.53	0.54
3:E:15:DC:N4	4:F:20:DG:O6	2.37	0.54
4:F:7:DG:H2'	4:F:8:DG:C8	2.42	0.54
2:B:289:TYR:HD2	2:B:300:GLU:HG2	1.71	0.54
2:D:256:LEU:CD2	2:D:284:ARG:HH12	2.20	0.54
1:A:423:GLN:HA	1:A:426:ARG:HD2	1.89	0.54
4:F:2:DT:H2'	4:F:3:DG:C8	2.42	0.54
1:A:708:ARG:NH2	1:A:719:SER:OG	2.41	0.54
2:D:43:THR:HG23	2:D:62:PHE:CE2	2.42	0.54
1:A:667:LYS:NZ	3:E:4:DA:OP2	2.40	0.54
3:E:18:DG:H1	4:F:17:DC:N4	1.98	0.54
1:A:462:ASN:OD1	1:A:466:GLN:CB	2.56	0.54
3:E:28:DC:H4'	3:E:29:DC:OP1	2.08	0.54
5:G:42:DT:H2''	5:G:43:DG:C8	2.43	0.54
1:A:512:GLY:HA2	1:A:513:ARG:HB3	1.90	0.53
1:A:795:ARG:NE	2:B:39:ARG:NH1	2.56	0.53
1:A:626:SER:O	1:A:994:ARG:NH2	2.42	0.53
2:D:80:PHE:CE1	2:D:151:THR:HG21	2.44	0.53
1:C:1000:GLN:CG	4:F:31:DT:C4'	2.86	0.53
2:D:58:ARG:NH1	7:J:7:DG:H5''	2.22	0.53
1:A:536:PHE:CD1	1:A:583:VAL:HG21	2.44	0.53
2:B:148:ARG:N	2:B:149:GLY:HA2	2.23	0.53
2:B:162:MET:HG3	2:B:163:PRO:HD2	1.89	0.53
4:F:13:DT:H4'	4:F:14:DG:OP1	2.08	0.53
1:A:424:LYS:HZ1	1:A:431:LYS:HE3	1.73	0.53
1:C:424:LYS:NZ	4:F:14:DG:C5	2.77	0.53
2:B:11:CYS:HB3	2:B:14:LEU:HD12	1.90	0.53
2:B:232:ARG:HH11	2:B:234:ILE:HD11	1.74	0.53
1:C:613:VAL:HG22	1:C:649:ILE:HD13	1.90	0.53
1:C:652:ARG:NH1	1:C:655:GLY:O	2.42	0.53
1:C:776:LEU:HD23	1:C:805:PRO:HD2	1.91	0.53
1:A:428:ARG:HA	1:A:431:LYS:NZ	2.21	0.52
1:A:410:ARG:HD3	6:H:36:DA:N3	2.24	0.52
2:D:104:SER:HB2	2:D:136:ALA:HB2	1.90	0.52
7:I:1:DG:H2''	7:I:2:DA:C8	2.45	0.52
2:D:262:ILE:HB	2:D:274:PHE:HB2	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:326:THR:HG21	2:B:347:GLN:HE21	1.74	0.52
1:C:732:LYS:HD2	7:J:16:DA:H4'	1.92	0.52
1:A:882:ARG:O	1:A:885:VAL:HG12	2.09	0.52
5:G:7:DG:H2'	5:G:8:DG:H8	1.73	0.52
1:C:486:CYS:SG	1:C:500:TYR:OH	2.47	0.52
1:C:730:ASP:OD1	1:C:730:ASP:N	2.38	0.52
6:H:33:DA:H2''	6:H:34:DA:C8	2.45	0.52
2:B:266:ILE:HG12	2:B:325:GLY:HA2	1.92	0.52
2:D:202:ASP:OD1	2:D:203:GLY:N	2.42	0.52
1:A:748:ILE:HG21	1:A:756:ARG:HE	1.75	0.52
1:A:583:VAL:HG12	1:A:709:LEU:HD21	1.91	0.51
2:D:43:THR:HG22	2:D:45:ILE:H	1.76	0.51
6:H:28:DT:H2'	6:H:29:DG:C8	2.45	0.51
2:B:94:GLY:O	2:B:103:SER:OG	2.27	0.51
2:D:80:PHE:HD2	2:D:89:CYS:HB2	1.74	0.51
2:B:278:GLN:HB3	2:B:284:ARG:HH12	1.74	0.51
1:C:561:THR:O	1:C:563:SER:N	2.43	0.51
1:C:833:GLU:O	1:C:834:ILE:C	2.49	0.51
2:D:307:TRP:HB3	2:D:311:ILE:HG23	1.91	0.51
3:E:2:DA:C8	3:E:2:DA:H5'	2.46	0.51
1:A:910:GLN:NE2	1:A:910:GLN:CB	2.73	0.51
2:B:96:ARG:HH11	2:B:100:ASN:HD21	1.57	0.51
1:C:883:GLU:OE1	1:C:883:GLU:N	2.43	0.51
2:D:289:TYR:HD2	2:D:300:GLU:HB3	1.75	0.51
1:A:811:PRO:O	1:A:971:ARG:NH1	2.43	0.51
2:D:82:ALA:HA	2:D:87:PRO:HB3	1.92	0.51
1:A:603:ASP:OD2	1:A:606:MET:HG2	2.11	0.51
1:A:773:ASP:N	1:A:773:ASP:OD1	2.42	0.51
2:B:154:VAL:HG11	2:B:216:VAL:HG21	1.93	0.51
1:A:795:ARG:CZ	2:B:39:ARG:NH1	2.74	0.51
2:B:237:HIS:CD2	2:B:239:GLU:OE2	2.63	0.50
1:C:610:GLY:HA3	1:C:653:LEU:HG	1.91	0.50
1:C:897:ARG:NH1	1:C:945:MET:HE2	2.26	0.50
2:D:256:LEU:HD21	2:D:284:ARG:HH12	1.76	0.50
5:G:13:DT:H3	6:H:33:DA:H2	1.57	0.50
6:H:28:DT:N3	6:H:29:DG:O6	2.44	0.50
3:E:18:DG:H2''	3:E:19:DA:C8	2.46	0.50
5:G:30:DG:H1'	5:G:31:DA:H5'	1.92	0.50
1:C:1000:GLN:HG2	4:F:31:DT:C4'	2.40	0.50
1:A:866:LYS:HD3	1:A:867:PRO:HD2	1.93	0.50
1:C:746:THR:HG22	1:C:755:THR:HB	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:22:DA:H2''	3:E:23:DA:C8	2.47	0.50
5:G:5:DA:H2''	5:G:6:DG:N7	2.27	0.50
2:B:150:LYS:HZ3	2:B:240:LEU:HB3	1.76	0.50
1:C:999:ARG:HD2	3:E:7:DG:N3	2.27	0.50
1:A:424:LYS:HZ2	1:A:431:LYS:HE3	1.76	0.50
1:C:489:ILE:HD11	1:A:503:MET:HA	1.94	0.50
2:D:239:GLU:OE2	2:D:246:VAL:HG13	2.12	0.50
1:C:735:ARG:NH1	1:C:743:SER:HA	2.27	0.49
1:A:751:LEU:HD23	1:A:811:PRO:HB2	1.93	0.49
1:C:673:SER:O	1:C:673:SER:OG	2.31	0.49
1:C:926:ASP:N	1:C:926:ASP:OD1	2.46	0.49
1:A:955:THR:HG23	1:A:958:LEU:HB2	1.94	0.49
2:B:198:PRO:HB2	2:B:199:GLU:OE2	2.12	0.49
1:C:656:GLU:N	1:C:656:GLU:OE1	2.33	0.49
1:A:472:ALA:C	1:A:477:ARG:NH1	2.66	0.49
2:B:15:VAL:HG12	2:B:30:LEU:HD11	1.95	0.49
2:B:208:VAL:HG21	2:B:261:ALA:HB3	1.94	0.49
1:C:587:LYS:HZ1	1:C:711:ILE:HG23	1.78	0.49
2:D:290:VAL:HG13	2:D:297:VAL:HG23	1.95	0.49
2:D:58:ARG:HH22	7:J:7:DG:H5''	1.76	0.49
1:A:748:ILE:CG2	1:A:756:ARG:HE	2.26	0.49
2:B:96:ARG:NH1	2:B:100:ASN:HD21	2.11	0.49
2:B:279:SER:HB3	2:B:282:GLN:HG2	1.94	0.49
1:C:709:LEU:O	1:C:719:SER:HA	2.13	0.49
1:C:520:HIS:NE2	4:F:23:DT:OP2	2.44	0.49
1:C:858:GLN:NE2	1:C:862:LYS:NZ	2.61	0.49
2:D:49:ARG:NH2	2:D:58:ARG:HD2	2.27	0.49
1:C:1000:GLN:HG3	4:F:31:DT:H4'	1.94	0.49
1:C:911:MET:HB2	1:C:931:TYR:CE2	2.48	0.48
1:A:414:HIS:ND1	1:A:414:HIS:N	2.60	0.48
1:C:448:SER:OG	1:A:415:LEU:HD13	2.07	0.48
1:C:789:GLU:HB3	1:C:793:GLU:OE1	2.13	0.48
1:A:611:PHE:HB2	1:A:720:PHE:HD1	1.78	0.48
1:C:553:ILE:HD12	1:C:1007:GLU:HG3	1.93	0.48
1:C:586:LEU:HD21	1:C:720:PHE:CD2	2.48	0.48
1:C:955:THR:OG1	1:C:956:ASN:N	2.46	0.48
1:C:735:ARG:HH12	1:C:743:SER:HA	1.79	0.48
2:D:65:ASN:HB2	2:D:123:ARG:NH1	2.28	0.48
3:E:21:DC:H2''	3:E:22:DA:C8	2.49	0.48
3:E:25:DA:H2''	3:E:26:DA:C8	2.48	0.48
1:C:630:GLY:C	4:F:30:DC:OP1	2.51	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:473:MET:HA	1:C:479:PHE:CG	2.49	0.48
2:B:187:LEU:HD12	2:B:188:GLU:H	1.79	0.48
4:F:43:DG:H1'	4:F:44:DC:H5'	1.96	0.48
4:F:6:DG:H2''	4:F:7:DG:C8	2.48	0.48
1:C:415:LEU:HB2	1:C:426:ARG:NH2	2.28	0.48
1:C:531:PRO:HB2	1:C:580:VAL:HG12	1.96	0.48
2:B:210:LEU:HD13	2:B:271:TYR:HD2	1.79	0.48
3:E:13:DG:H2''	3:E:14:DA:H8	1.79	0.48
5:G:3:DG:H2'	5:G:4:DC:C6	2.49	0.48
1:A:551:VAL:HB	1:A:554:ILE:HD11	1.96	0.48
1:C:481:LEU:HD11	1:C:516:PHE:HA	1.96	0.48
7:J:4:DC:H2''	7:J:5:DT:C6	2.49	0.48
1:C:414:HIS:HD2	1:C:416:LEU:HG	1.79	0.47
2:D:148:ARG:N	2:D:149:GLY:HA2	2.29	0.47
1:A:870:ARG:HD2	7:I:16:DA:C6	2.50	0.47
1:A:420:ARG:CZ	6:H:27:DC:H5''	2.44	0.47
1:A:620:ASP:OD1	1:A:621:GLY:N	2.47	0.47
2:D:218:PHE:HB2	2:D:234:ILE:HB	1.95	0.47
2:B:31:PHE:HD1	2:B:45:ILE:HG12	1.78	0.47
2:B:93:HIS:HA	2:B:94:GLY:HA2	1.64	0.47
1:C:469:GLU:OE2	1:C:479:PHE:HZ	1.96	0.47
1:C:955:THR:HG23	1:C:958:LEU:HB2	1.96	0.47
2:B:82:ALA:HA	2:B:87:PRO:HB3	1.96	0.47
1:C:626:SER:O	1:C:994:ARG:NH2	2.48	0.47
2:D:179:PRO:HB2	2:D:181:GLN:OE1	2.14	0.47
2:D:154:VAL:HG11	2:D:216:VAL:HG21	1.96	0.47
4:F:7:DG:C6	4:F:8:DG:C6	3.03	0.47
1:C:458:LEU:HD21	1:C:466:GLN:HB2	1.97	0.47
5:G:15:DT:H2''	5:G:16:DA:H5''	1.97	0.47
1:A:762:ASN:HD22	1:A:762:ASN:C	2.18	0.47
1:C:713:VAL:CG1	1:C:718:ARG:HD2	2.45	0.47
2:B:10:ASN:HD21	2:B:56:LYS:NZ	2.13	0.46
2:B:10:ASN:ND2	2:B:56:LYS:NZ	2.63	0.46
2:B:207:HIS:ND1	2:B:207:HIS:O	2.48	0.46
2:B:249:CYS:SG	2:B:250:THR:N	2.88	0.46
2:B:128:GLU:OE2	2:B:130:VAL:HG22	2.14	0.46
2:B:222:HIS:HB2	2:B:259:THR:HG21	1.97	0.46
1:C:420:ARG:HH11	1:A:459:ARG:NH2	2.14	0.46
1:C:912:LYS:HA	1:C:915:TRP:CZ3	2.51	0.46
1:A:445:ASP:OD2	1:A:448:SER:HB3	2.15	0.46
3:E:22:DA:H2''	3:E:23:DA:H8	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:350:PHE:CD2	2:B:351:GLN:HG3	2.50	0.46
2:B:232:ARG:HH11	2:B:234:ILE:CD1	2.29	0.46
4:F:26:DA:H2''	4:F:27:DG:C8	2.51	0.46
1:A:449:VAL:O	1:A:453:LEU:HG	2.16	0.46
2:D:56:LYS:HD2	2:D:56:LYS:HA	1.79	0.46
1:C:1000:GLN:HG3	4:F:31:DT:C4'	2.46	0.46
1:A:679:LEU:HD13	1:A:1010:LEU:HD13	1.97	0.46
1:C:579:ASP:OD1	1:C:579:ASP:N	2.48	0.46
2:B:1:MET:SD	2:B:302:ARG:NH1	2.89	0.46
1:A:424:LYS:O	1:A:428:ARG:HG3	2.16	0.46
2:B:1:MET:HB3	2:B:305:PRO:HG3	1.98	0.46
2:B:104:SER:HB2	2:B:136:ALA:HB2	1.97	0.45
1:A:625:VAL:HB	1:A:637:GLU:O	2.15	0.45
2:D:114:ARG:HB2	2:D:115:GLY:HA2	1.99	0.45
2:D:70:PRO:HA	2:D:71:PRO:HD3	1.87	0.45
5:G:12:DT:H2''	5:G:13:DT:H5'	1.97	0.45
1:C:448:SER:O	1:C:452:THR:HG22	2.17	0.45
1:C:481:LEU:HD23	1:A:507:VAL:HG23	1.99	0.45
4:F:36:DA:H2''	4:F:37:DA:O4'	2.16	0.45
1:A:438:ALA:O	1:A:443:GLY:N	2.50	0.45
1:C:481:LEU:HG	1:C:515:ILE:HG22	1.98	0.45
1:C:649:ILE:CG2	1:C:663:PHE:HB3	2.46	0.45
1:C:680:MET:HB3	1:C:682:VAL:HG23	1.97	0.45
1:A:462:ASN:HB3	1:A:466:GLN:CD	2.29	0.45
2:B:262:ILE:HB	2:B:274:PHE:HB3	1.97	0.45
1:C:707:SER:OG	1:C:708:ARG:N	2.50	0.45
1:C:742:ALA:HA	1:C:743:SER:HA	1.65	0.45
3:E:13:DG:H2''	3:E:14:DA:C8	2.52	0.45
5:G:21:DC:H2''	5:G:22:DA:C8	2.52	0.45
5:G:7:DG:H2'	5:G:8:DG:C8	2.51	0.45
1:A:507:VAL:HG11	1:A:516:PHE:HD2	1.80	0.45
2:B:284:ARG:HH11	2:B:284:ARG:CB	2.23	0.45
1:C:416:LEU:HD21	1:A:451:LEU:HD11	1.99	0.45
1:A:762:ASN:CB	1:A:762:ASN:HD22	2.24	0.45
6:H:16:DC:H2'	6:H:17:DC:C5	2.52	0.45
2:B:239:GLU:OE1	2:B:239:GLU:N	2.50	0.45
1:C:462:ASN:N	1:C:462:ASN:OD1	2.49	0.45
2:D:145:ILE:HD11	2:D:152:ALA:HB3	1.98	0.45
3:E:16:DT:H2''	3:E:17:DG:H5'	1.98	0.45
6:H:33:DA:H2''	6:H:34:DA:H8	1.81	0.45
2:B:242:LEU:HA	2:B:243:GLY:HA2	1.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:265:PRO:HA	2:B:271:TYR:HD1	1.82	0.44
1:C:611:PHE:HB2	1:C:720:PHE:HD1	1.82	0.44
2:D:87:PRO:HA	2:D:88:GLU:HA	1.57	0.44
1:C:410:ARG:HD2	3:E:25:DA:H2	1.81	0.44
1:A:821:GLY:O	1:A:824:THR:HG22	2.17	0.44
1:A:872:ASN:OD1	1:A:872:ASN:N	2.49	0.44
2:B:232:ARG:NH1	2:B:234:ILE:HD11	2.31	0.44
2:B:289:TYR:CD2	2:B:300:GLU:HG2	2.51	0.44
1:C:748:ILE:HD12	1:C:749:CYS:HB2	1.97	0.44
1:A:421:ARG:HE	1:A:424:LYS:HB2	1.82	0.44
2:B:107:TYR:CE1	2:B:126:GLU:HG2	2.51	0.44
2:D:212:ARG:HG3	2:D:269:HIS:NE2	2.33	0.44
2:D:313:HIS:CE1	1:A:891:LEU:HD23	2.52	0.44
2:B:202:ASP:OD1	2:B:203:GLY:N	2.50	0.44
2:B:81:GLU:OE1	2:B:82:ALA:N	2.50	0.44
1:C:656:GLU:HG2	1:C:657:ASP:N	2.30	0.44
2:D:14:LEU:HD13	2:D:46:PHE:CZ	2.53	0.44
3:E:17:DG:H2''	3:E:18:DG:C8	2.53	0.44
1:A:649:ILE:CG2	1:A:663:PHE:HB3	2.47	0.44
6:H:34:DA:H2''	6:H:35:DA:H8	1.83	0.44
6:H:5:DG:H2''	6:H:6:DT:OP1	2.16	0.44
2:D:235:ARG:HB3	2:D:250:THR:HG23	2.00	0.44
5:G:29:DG:H1'	5:G:30:DG:O4'	2.17	0.44
6:H:39:DC:H4'	6:H:40:DC:OP1	2.18	0.44
1:A:590:GLU:HG3	1:A:713:VAL:HG23	2.00	0.44
1:A:746:THR:O	1:A:756:ARG:N	2.50	0.44
2:B:256:LEU:HD21	2:B:284:ARG:NE	2.33	0.44
1:C:536:PHE:CD2	1:C:583:VAL:HG21	2.52	0.44
1:C:918:THR:O	1:C:974:SER:OG	2.20	0.44
2:D:311:ILE:HG13	2:D:331:ILE:HD11	2.00	0.44
2:D:28:VAL:O	2:D:48:VAL:HG12	2.18	0.44
1:A:628:LYS:HE2	1:A:1005:GLU:OE2	2.18	0.44
1:A:738:GLU:OE2	1:A:775:ASN:ND2	2.51	0.44
1:C:813:LEU:HD11	1:C:818:CYS:HB2	1.98	0.44
1:C:822:ASN:HD22	1:C:908:TYR:HE1	1.66	0.43
1:C:849:ARG:NH1	1:C:849:ARG:HB3	2.33	0.43
1:C:895:GLU:N	1:C:895:GLU:OE1	2.51	0.43
1:A:412:ARG:HD3	1:A:412:ARG:HA	1.76	0.43
2:B:339:PRO:HB2	2:B:341:GLU:N	2.34	0.43
1:C:489:ILE:HA	1:C:489:ILE:HD12	1.84	0.43
1:C:582:LEU:HD23	1:C:582:LEU:HA	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:844:SER:N	1:C:847:GLU:OE1	2.33	0.43
2:D:148:ARG:NH1	2:D:240:LEU:O	2.50	0.43
1:A:790:SER:OG	1:A:793:GLU:OE1	2.36	0.43
1:A:463:GLU:HA	4:F:22:DC:H4'	2.00	0.43
2:D:93:HIS:HA	2:D:94:GLY:HA2	1.61	0.43
1:A:423:GLN:HA	1:A:426:ARG:HB2	1.99	0.43
1:A:951:ASP:N	1:A:951:ASP:OD1	2.52	0.43
1:A:465:LYS:H	1:A:465:LYS:HD2	1.82	0.43
1:A:530:LEU:HD11	1:A:1011:LYS:HG3	2.01	0.43
1:C:465:LYS:HG3	1:C:466:GLN:N	2.34	0.43
1:C:590:GLU:HG3	1:C:713:VAL:HG23	2.00	0.43
1:C:587:LYS:NZ	1:C:711:ILE:HG23	2.33	0.43
1:C:834:ILE:HD11	1:C:891:LEU:HD13	2.00	0.43
1:A:628:LYS:NZ	5:G:42:DT:H4'	2.34	0.43
2:B:41:CYS:HG	2:B:46:PHE:HE1	1.66	0.43
2:D:99:ASN:O	2:D:101:GLU:N	2.51	0.43
1:A:423:GLN:NE2	1:A:426:ARG:NH1	2.67	0.43
1:A:476:GLY:HA3	1:A:477:ARG:HA	1.46	0.43
1:A:472:ALA:HA	1:A:477:ARG:HH12	1.83	0.43
1:A:958:LEU:HA	1:A:961:THR:HG22	2.00	0.43
2:B:108:MET:HB2	2:B:127:LYS:NZ	2.33	0.43
2:B:137:ARG:HG3	2:B:140:HIS:ND1	2.34	0.43
1:C:456:LEU:HG	1:A:420:ARG:HE	1.83	0.43
1:A:490:ARG:NH2	1:A:497:CYS:SG	2.91	0.43
2:B:141:THR:HG21	2:B:209:ALA:HB3	2.01	0.43
1:C:475:GLN:HG2	1:C:476:GLY:H	1.84	0.43
1:C:725:ARG:HD2	1:C:810:GLN:OE1	2.19	0.43
1:C:748:ILE:HG21	1:C:756:ARG:HD3	2.01	0.43
2:D:58:ARG:NH2	2:D:118:ARG:O	2.52	0.43
1:A:926:ASP:OD1	1:A:926:ASP:N	2.49	0.42
2:B:140:HIS:CD2	2:B:155:LEU:HD11	2.54	0.42
3:E:33:DA:H2	4:F:2:DT:H3	1.67	0.42
5:G:4:DC:H2''	5:G:5:DA:H8	1.84	0.42
1:A:830:PHE:O	1:A:834:ILE:HG12	2.19	0.42
2:B:226:SER:OG	2:B:228:CYS:SG	2.61	0.42
1:C:420:ARG:NH1	1:A:459:ARG:HE	2.17	0.42
1:A:421:ARG:NH2	1:A:424:LYS:HD3	2.33	0.42
5:G:6:DG:H2'	5:G:7:DG:C8	2.55	0.42
6:H:40:DC:H2'	6:H:41:DT:H6	1.78	0.42
7:I:12:DT:C4	7:I:13:DC:N4	2.87	0.42
2:B:317:TRP:HB3	2:B:331:ILE:HD13	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:871:MET:SD	1:C:875:TYR:HD2	2.42	0.42
3:E:4:DA:H2''	3:E:5:DG:H5''	2.01	0.42
7:J:7:DG:H2''	7:J:8:DC:OP2	2.20	0.42
2:B:33:GLN:OE1	2:B:44:GLY:HA2	2.19	0.42
2:D:242:LEU:HA	2:D:243:GLY:HA2	1.52	0.42
2:B:75:PRO:HG2	2:B:77:ILE:CD1	2.50	0.42
1:C:625:VAL:HB	1:C:637:GLU:O	2.19	0.42
3:E:24:DA:H2''	3:E:25:DA:OP2	2.20	0.42
3:E:22:DA:C6	4:F:14:DG:N2	2.88	0.42
1:C:1006:LEU:HD23	1:C:1006:LEU:HA	1.80	0.42
1:C:598:ARG:CZ	1:C:604:ASP:OD2	2.67	0.42
2:D:148:ARG:CZ	2:D:148:ARG:HB3	2.49	0.42
3:E:20:DA:H2	4:F:15:DT:H3	1.68	0.42
7:I:14:DT:H2'	7:I:14:DT:H6	1.66	0.42
1:A:628:LYS:HZ2	5:G:42:DT:H4'	1.85	0.42
1:C:475:GLN:HG2	1:C:476:GLY:N	2.35	0.42
2:D:274:PHE:CE2	2:D:348:VAL:HG21	2.54	0.42
1:A:460:ALA:HA	3:E:16:DT:H4'	2.01	0.42
3:E:20:DA:H2''	3:E:21:DC:C6	2.54	0.42
5:G:39:DC:H2''	5:G:40:DA:C8	2.55	0.42
5:G:7:DG:H22	6:H:39:DC:H42	1.65	0.42
1:A:628:LYS:NZ	5:G:42:DT:H5''	2.34	0.41
1:A:742:ALA:HA	1:A:743:SER:HA	1.61	0.41
2:B:14:LEU:HD13	2:B:46:PHE:CE2	2.55	0.41
2:D:25:GLU:HB3	2:D:114:ARG:HH12	1.85	0.41
4:F:11:DT:C6	4:F:12:DT:H72	2.56	0.41
1:A:478:GLY:HA2	1:A:479:PHE:C	2.40	0.41
1:C:653:LEU:HA	1:C:654:GLU:HA	1.82	0.41
2:D:110:SER:OG	2:D:111:VAL:N	2.53	0.41
2:D:141:THR:HG21	2:D:207:HIS:O	2.21	0.41
2:D:14:LEU:HD13	2:D:46:PHE:CE2	2.55	0.41
3:E:31:DG:C2	4:F:5:DA:C2	3.08	0.41
5:G:56:DC:H2''	5:G:57:DA:C8	2.55	0.41
2:B:311:ILE:HG13	2:B:331:ILE:HD11	2.00	0.41
2:B:84:ASP:HA	2:B:85:GLY:HA2	1.49	0.41
1:C:507:VAL:HG23	1:A:481:LEU:HD11	2.03	0.41
2:D:156:PHE:CE1	2:D:200:LEU:HD23	2.55	0.41
3:E:31:DG:H2''	3:E:32:DC:C6	2.55	0.41
1:A:675:ARG:NH2	1:A:1018:SER:HA	2.35	0.41
2:D:85:GLY:O	2:D:87:PRO:HD3	2.19	0.41
1:A:410:ARG:HH22	5:G:12:DT:C4'	2.33	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:730:ASP:OD1	1:A:730:ASP:N	2.51	0.41
1:C:420:ARG:NH1	1:A:459:ARG:NE	2.69	0.41
1:C:834:ILE:HD12	1:C:855:LEU:HB2	2.02	0.41
4:F:16:DT:C4	4:F:17:DC:N4	2.89	0.41
1:A:473:MET:HB2	1:A:475:GLN:NE2	2.36	0.41
1:C:833:GLU:OE1	1:C:833:GLU:HA	2.19	0.41
1:C:860:ARG:HH21	1:C:867:PRO:HD2	1.86	0.41
2:D:207:HIS:O	2:D:207:HIS:ND1	2.53	0.41
1:A:429:ASP:N	1:A:429:ASP:OD1	2.51	0.41
1:A:984:GLU:O	1:A:987:ASN:HB2	2.21	0.41
1:C:682:VAL:HG12	1:C:683:ASP:O	2.21	0.41
5:G:28:DT:H1'	5:G:29:DG:C8	2.56	0.41
6:H:14:DC:H2''	6:H:15:DT:C6	2.56	0.41
5:G:8:DG:N2	6:H:39:DC:C2	2.89	0.41
1:A:587:LYS:HE3	1:A:712:SER:O	2.21	0.41
2:B:87:PRO:HA	2:B:88:GLU:HA	1.69	0.41
1:C:495:LEU:HD23	1:C:495:LEU:HA	1.89	0.41
1:C:748:ILE:CG2	1:C:756:ARG:HD3	2.51	0.41
2:D:131:GLY:HA3	2:D:132:ASP:C	2.41	0.41
1:A:489:ILE:O	1:A:493:THR:HG22	2.20	0.41
1:A:594:MET:O	1:A:598:ARG:HG3	2.20	0.41
1:A:815:ALA:HB1	1:A:916:ARG:HH11	1.85	0.41
2:B:145:ILE:HD11	2:B:152:ALA:HB3	2.02	0.41
1:C:672:LEU:HB3	1:C:1022:GLN:HE22	1.85	0.41
1:C:435:LYS:HA	1:C:435:LYS:HD3	1.93	0.41
2:D:150:LYS:HE3	2:D:150:LYS:HB2	1.83	0.41
2:D:22:LEU:HD22	2:D:90:TYR:CD2	2.56	0.41
2:D:72:LEU:HG	2:D:72:LEU:H	1.64	0.41
1:C:519:LEU:HA	1:C:519:LEU:HD23	1.91	0.41
1:C:795:ARG:NE	2:D:39:ARG:HE	2.19	0.41
1:C:955:THR:OG1	7:J:15:DT:OP1	2.38	0.41
1:A:472:ALA:CA	1:A:477:ARG:HH12	2.33	0.40
1:A:859:LEU:HD23	1:A:859:LEU:HA	1.88	0.40
2:B:152:ALA:CB	2:B:240:LEU:HD21	2.51	0.40
2:B:51:LYS:HG2	2:B:52:LYS:HG3	2.02	0.40
1:A:738:GLU:HA	1:A:805:PRO:HB3	2.02	0.40
1:A:866:LYS:CD	1:A:867:PRO:HD2	2.51	0.40
1:A:981:GLU:OE2	4:F:34:DG:C2	2.75	0.40
2:B:350:PHE:HD2	2:B:351:GLN:HG3	1.86	0.40
1:C:944:SER:OG	1:C:945:MET:N	2.55	0.40
4:F:7:DG:N1	4:F:8:DG:C6	2.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:531:PRO:HB3	1:A:580:VAL:HG12	2.04	0.40
1:A:942:LEU:HD23	1:A:942:LEU:HA	1.95	0.40
2:B:184:LEU:HD12	2:B:245:PRO:HB2	2.02	0.40
1:C:437:PHE:HD1	1:C:440:LYS:HZ2	1.67	0.40
7:I:5:DT:H2''	7:I:6:DG:N7	2.36	0.40
1:A:579:ASP:OD1	1:A:580:VAL:N	2.54	0.40
1:C:672:LEU:HB3	1:C:1022:GLN:NE2	2.36	0.40
1:A:981:GLU:OE2	4:F:34:DG:N1	2.54	0.40
7:J:5:DT:H2''	7:J:6:DG:C8	2.57	0.40

There are no symmetry-related clashes.

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 PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	620/1159 (54%)	592 (96%)	26 (4%)	2 (0%)	43	76
1	C	620/1159 (54%)	593 (96%)	27 (4%)	0	100	100
2	B	349/533 (66%)	330 (95%)	19 (5%)	0	100	100
2	D	349/533 (66%)	322 (92%)	27 (8%)	0	100	100
All	All	1938/3384 (57%)	1837 (95%)	99 (5%)	2 (0%)	56	85

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	524	ASN
1	A	562	VAL

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 PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	548/1000 (55%)	528 (96%)	20 (4%)	38	70
1	C	548/1000 (55%)	534 (97%)	14 (3%)	49	77
2	B	303/465 (65%)	297 (98%)	6 (2%)	58	81
2	D	303/465 (65%)	297 (98%)	6 (2%)	58	81
All	All	1702/2930 (58%)	1656 (97%)	46 (3%)	51	77

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	428	ARG
1	C	514	GLN
1	C	560	TRP
1	C	582	LEU
1	C	628	LYS
1	C	654	GLU
1	C	756	ARG
1	C	833	GLU
1	C	870	ARG
1	C	915	TRP
1	C	917	SER
1	C	992	ARG
1	C	999	ARG
1	C	1000	GLN
2	D	79	HIS
2	D	116	CYS
2	D	172	TRP
2	D	300	GLU
2	D	302	ARG
2	D	347	GLN
1	A	414	HIS
1	A	416	LEU
1	A	417	SER
1	A	464	HIS

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Mol	Chain	Res	Type
1	A	466	GLN
1	A	479	PHE
1	A	514	GLN
1	A	560	TRP
1	A	588	ASP
1	A	702	LYS
1	A	729	TYR
1	A	752	CYS
1	A	860	ARG
1	A	869	MET
1	A	947	LYS
1	A	997	ASN
1	A	999	ARG
1	A	1000	GLN
1	A	1008	ASP
1	A	1029	LYS
2	B	106	LEU
2	B	116	CYS
2	B	127	LYS
2	B	162	MET
2	B	232	ARG
2	B	299	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	858	GLN
1	C	983	ASN
1	C	1012	HIS
2	D	207	HIS
1	A	762	ASN
1	A	1000	GLN
1	A	1012	HIS
1	A	1022	GLN
2	B	10	ASN
2	B	173	ASN
2	B	237	HIS

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 6 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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.