



## Full wwPDB EM Validation Report ⓘ

Apr 6, 2026 – 12:30 AM UTC

PDB ID : 9JHD / pdb\_00009jhd  
EMDB ID : EMD-61479  
Title : Cryo-EM structure of RacrIC1-Cascade IC1-complex III  
Authors : Wang, H.; Zhang, S.; Li, S.; Zhang, K.; Feng, Y.  
Deposited on : 2024-09-09  
Resolution : 2.98 Å(reported)

This is a Full wwPDB EM 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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

EMDB validation analysis : 0.0.1.dev132  
MolProbity : 4-5-2 with Phenix2.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
EM percentile statistics : **NOT EXECUTED**  
MapQ : **FAILED**  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

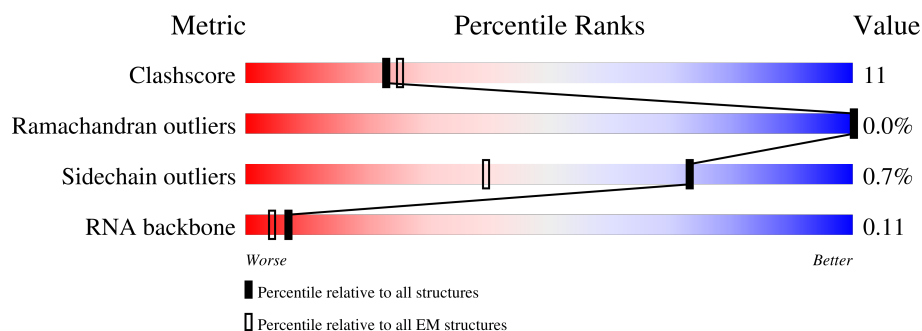
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 2.98 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.













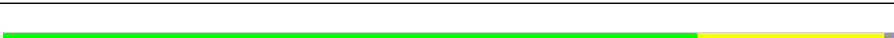

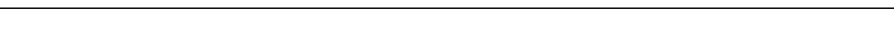
Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102
RNA backbone	8273	3508

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	289	72% 26% ..
1	B	289	74% 25% .
1	C	289	78% 20% .
1	D	289	72% 27% .
1	E	289	78% 21% .
1	F	289	78% 21% .
1	G	289	78% 20% .

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Mol	Chain	Length	Quality of chain
1	H	289	 76% 22% ..
1	I	289	 76% 22% .
1	J	289	 75% 24% .
1	K	289	 75% 24% .
2	L	116	 76% 22% .
2	M	116	 68% 29% ..
2	N	116	 73% 25% .
2	O	116	 75% 23% .
2	P	116	 73% 25% .
2	Q	116	 76% 22% .
2	R	116	 78% 21% .
2	S	116	 91% 8% .
3	T	124	 23% 25% 52%

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 64916 atoms, of which 31760 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 CRISPR-associated protein, TM1801 family.

Mol	Chain	Residues	Atoms						AltConf	Trace
1	A	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0
1	B	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0
1	C	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0
1	D	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0
1	E	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0
1	F	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0
1	G	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0
1	H	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0
1	I	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0
1	J	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0
1	K	286	Total 4472	C 1408	H 2232	N 401	O 418	S 13	0	0

- Molecule 2 is a protein called CRISPR-associated protein, CT1133 family.

Mol	Chain	Residues	Atoms						AltConf	Trace
2	L	114	Total 1806	C 572	H 901	N 159	O 169	S 5	0	0
2	M	114	Total 1806	C 572	H 901	N 159	O 169	S 5	0	0
2	N	114	Total 1806	C 572	H 901	N 159	O 169	S 5	0	0
2	O	114	Total 1806	C 572	H 901	N 159	O 169	S 5	0	0

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Mol	Chain	Residues	Atoms						AltConf	Trace
2	P	114	Total	C	H	N	O	S	0	0
			1806	572	901	159	169	5		
2	Q	114	Total	C	H	N	O	S	0	0
			1806	572	901	159	169	5		
2	R	114	Total	C	H	N	O	S	0	0
			1806	572	901	159	169	5		
2	S	114	Total	C	H	N	O	S	0	0
			1806	572	901	159	169	5		

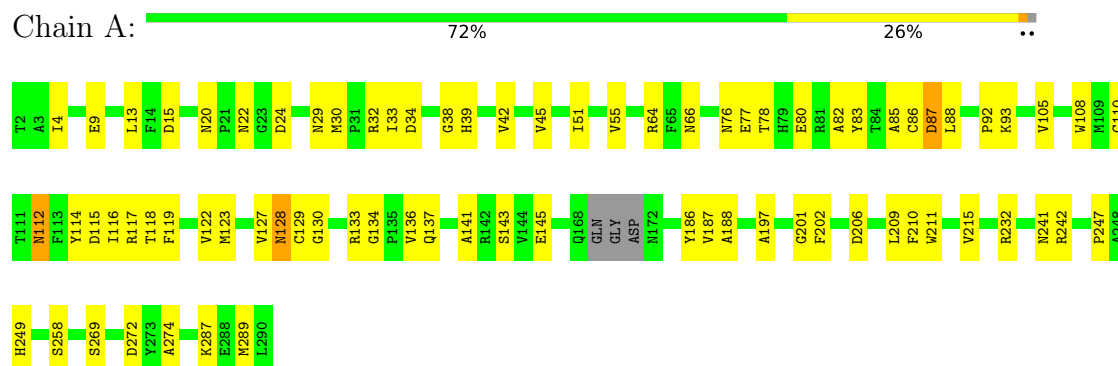
- Molecule 3 is a RNA chain called RacrIC1 (124-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
3	T	60	Total	C	N	O	P	0	0
			1276	571	230	415	60		

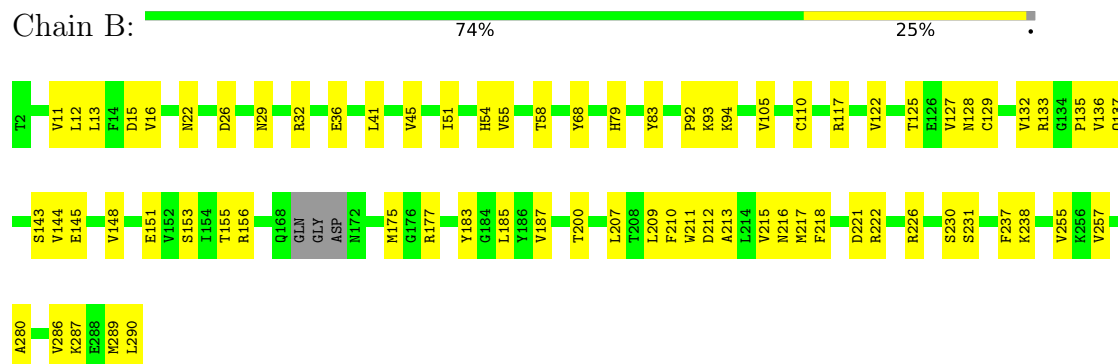
### 3 Residue-property plots [i](#)

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. 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.

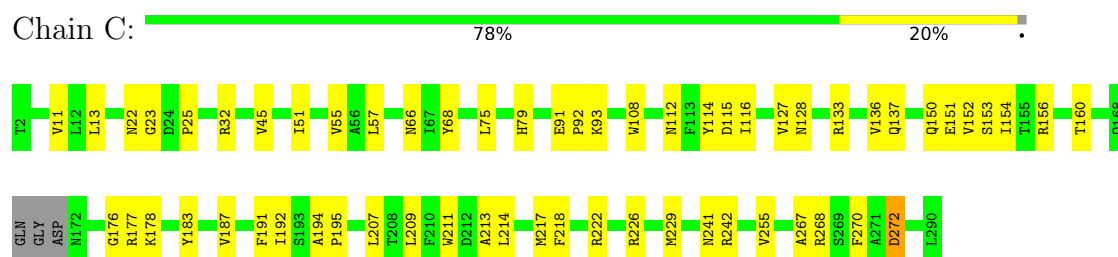
- Molecule 1: CRISPR-associated protein, TM1801 family



- Molecule 1: CRISPR-associated protein, TM1801 family

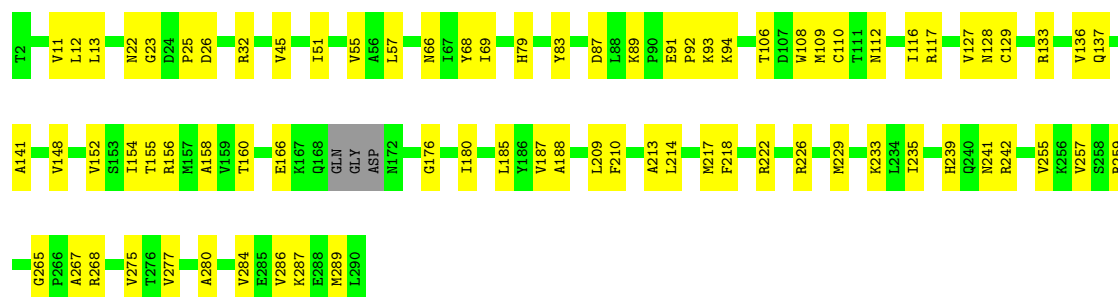


- Molecule 1: CRISPR-associated protein, TM1801 family



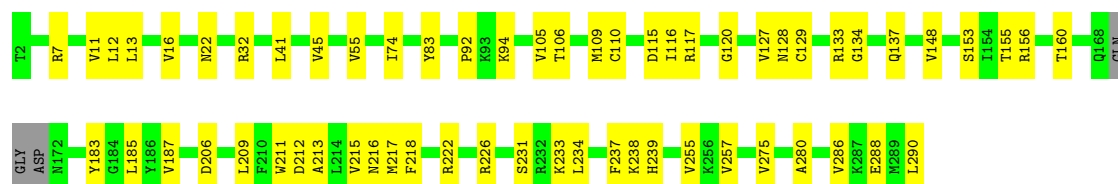
- Molecule 1: CRISPR-associated protein, TM1801 family

Chain D:  72% 27%




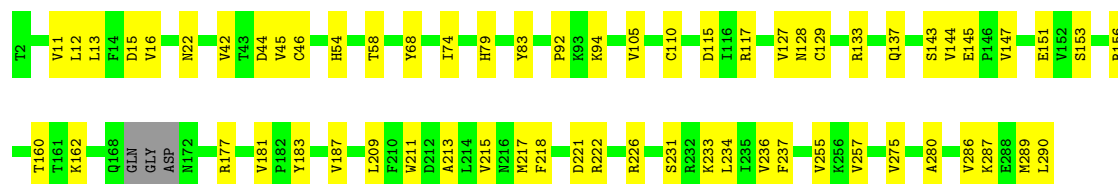
- Molecule 1: CRISPR-associated protein, TM1801 family

Chain E:  78% 21%




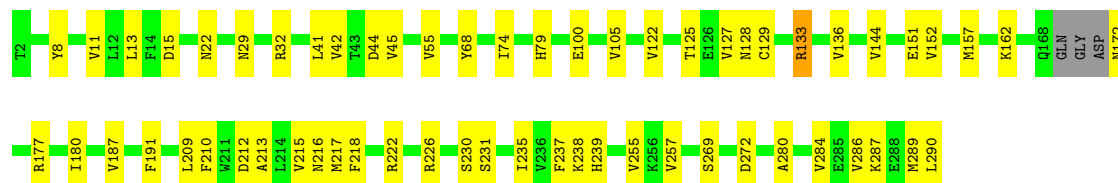
- Molecule 1: CRISPR-associated protein, TM1801 family

Chain F:  78% 21%



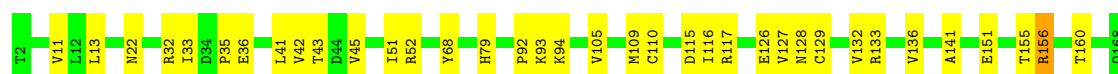
- Molecule 1: CRISPR-associated protein, TM1801 family

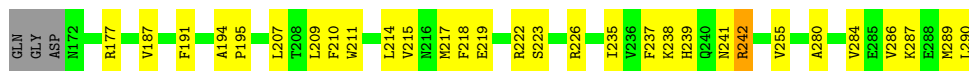
Chain G:  78% 20%



- Molecule 1: CRISPR-associated protein, TM1801 family

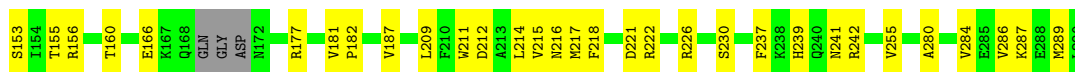
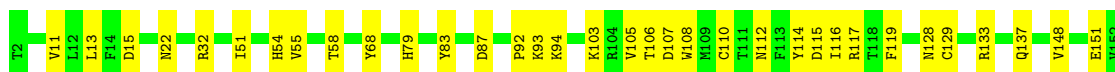
Chain H:  76% 22%





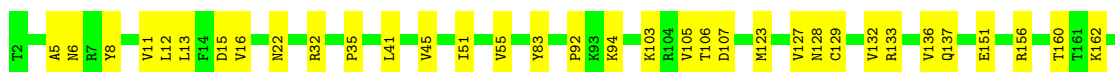
- Molecule 1: CRISPR-associated protein, TM1801 family

Chain I: 76% 22%



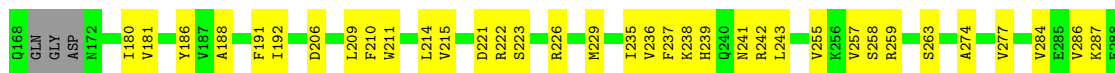
- Molecule 1: CRISPR-associated protein, TM1801 family

Chain J: 75% 24%



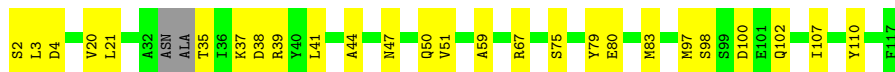
- Molecule 1: CRISPR-associated protein, TM1801 family

Chain K: 75% 24%



- Molecule 2: CRISPR-associated protein, CT1133 family

Chain L: 76% 22%



- Molecule 2: CRISPR-associated protein, CT1133 family

Chain M: 68% 29%







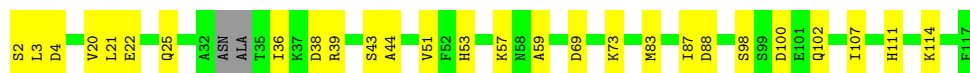
- Molecule 2: CRISPR-associated protein, CT1133 family

Chain N: 73% 25% .



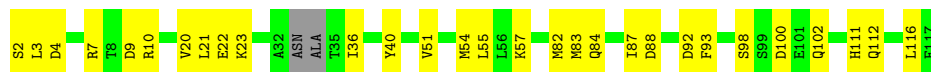
- Molecule 2: CRISPR-associated protein, CT1133 family

Chain O: 75% 23% .



- Molecule 2: CRISPR-associated protein, CT1133 family

Chain P: 73% 25% .



- Molecule 2: CRISPR-associated protein, CT1133 family

Chain Q: 76% 22% .



- Molecule 2: CRISPR-associated protein, CT1133 family

Chain R: 78% 21% .



- Molecule 2: CRISPR-associated protein, CT1133 family

Chain S: 91% 8% .



- Molecule 3: RacIC1 (124-MER)

Chain T: 23% 25% 52% .

A U A A	G32	C33	A34	C35	A36	U37	G38	C39	A40	C41	U42	C43	G44	U45	C46	U47	A48	C49	U50	A51	G52	C53	G54	C55	A56	G57	C58	U59	U60	A	A1	A2	G3	A4	A5	A6	C7	C8	G9	C10	U11	G12	C13	U14	G15	C16	G17	A18	C19	A20	U21	U22	U23	G24	A25	C26	U27	G28	C29	C30	A31

## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	82089	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	58	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor

## 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	0.15	0/2285	0.27	0/3087
1	B	0.35	0/2285	0.30	0/3087
1	C	0.24	0/2285	0.31	0/3087
1	D	0.30	0/2285	0.31	0/3087
1	E	0.37	0/2285	0.30	0/3087
1	F	0.38	0/2285	0.30	0/3087
1	G	0.37	0/2285	0.31	0/3087
1	H	0.36	0/2285	0.31	0/3087
1	I	0.30	0/2285	0.31	0/3087
1	J	0.38	0/2285	0.31	0/3087
1	K	0.20	0/2285	0.30	0/3087
2	L	0.15	0/923	0.25	0/1241
2	M	0.17	0/923	0.25	0/1241
2	N	0.17	0/923	0.24	0/1241
2	O	0.15	0/923	0.23	0/1241
2	P	0.13	0/923	0.22	0/1241
2	Q	0.14	0/923	0.23	0/1241
2	R	0.12	0/923	0.22	0/1241
2	S	0.10	0/923	0.19	0/1241
3	T	0.42	0/1426	0.39	0/2219
All	All	0.30	0/33945	0.29	0/46104

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	2240	2232	2228	53	0
1	B	2240	2232	2228	55	0
1	C	2240	2232	2228	46	0
1	D	2240	2232	2228	62	0
1	E	2240	2232	2228	49	0
1	F	2240	2232	2228	46	0
1	G	2240	2232	2228	51	0
1	H	2240	2232	2228	51	0
1	I	2240	2232	2228	48	0
1	J	2240	2232	2228	55	0
1	K	2240	2232	2228	50	0
2	L	905	901	897	21	0
2	M	905	901	897	31	0
2	N	905	901	897	24	0
2	O	905	901	897	24	0
2	P	905	901	897	24	0
2	Q	905	901	897	20	0
2	R	905	901	897	20	0
2	S	905	901	897	7	0
3	T	1276	0	649	138	0
All	All	33156	31760	32333	725	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 11.

All (725) 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)
3:T:31:A:O2'	3:T:32:G:H4'	1.62	0.97
2:Q:2:SER:N	2:R:88:ASP:O	2.01	0.93
3:T:2:A:O2'	3:T:3:G:OP2	1.88	0.91
1:H:45:VAL:CG1	3:T:13:C:H5'	2.02	0.90
1:A:143:SER:OG	1:A:145:GLU:O	1.90	0.89
3:T:1:A:H2'	3:T:2:A:H4'	1.54	0.88
1:H:222:ARG:O	1:J:133:ARG:NH1	2.07	0.86
1:I:133:ARG:NH1	1:K:222:ARG:O	2.10	0.85
3:T:59:U:H2'	3:T:60:U:H5'	1.59	0.84
2:R:98:SER:OG	2:R:100:ASP:OD1	1.94	0.83
1:H:133:ARG:NH1	1:I:222:ARG:O	2.11	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:2:SER:N	2:O:88:ASP:O	2.12	0.82
2:R:2:SER:N	2:S:88:ASP:O	2.13	0.82
1:B:222:ARG:O	1:D:133:ARG:NH1	2.13	0.82
2:O:2:SER:N	2:P:88:ASP:O	2.13	0.81
1:A:110:CYS:O	1:A:117:ARG:NH1	2.13	0.81
1:C:68:TYR:OH	1:C:79:HIS:NE2	2.13	0.81
3:T:56:A:O2'	3:T:57:G:O5'	1.99	0.81
1:D:26:ASP:OD1	2:M:39:ARG:NH1	2.14	0.80
1:D:45:VAL:CG1	3:T:49:C:H5'	2.11	0.80
2:N:98:SER:OG	2:N:100:ASP:OD1	2.00	0.79
1:E:133:ARG:NH1	1:F:222:ARG:O	2.16	0.79
1:G:151:GLU:OE2	1:G:177:ARG:NE	2.16	0.79
2:P:2:SER:N	2:Q:88:ASP:O	2.15	0.79
1:K:45:VAL:HG12	3:T:1:A:H5''	1.63	0.78
2:L:2:SER:N	2:M:88:ASP:O	2.17	0.78
1:D:137:GLN:NE2	3:T:47:U:OP1	2.17	0.78
2:O:25:GLN:OE1	2:O:36:ILE:N	2.17	0.78
2:S:98:SER:OG	2:S:100:ASP:OD1	2.02	0.77
1:H:68:TYR:OH	1:H:79:HIS:NE2	2.17	0.77
1:I:137:GLN:NE2	3:T:5:A:OP1	2.17	0.77
1:K:143:SER:OG	1:K:145:GLU:O	2.02	0.77
1:A:108:TRP:O	1:A:112:ASN:ND2	2.18	0.77
1:I:151:GLU:OE2	1:I:177:ARG:NE	2.18	0.77
1:I:166:GLU:OE1	1:I:166:GLU:N	2.16	0.77
3:T:31:A:H2'	3:T:31:A:N3	1.99	0.77
1:H:156:ARG:NH2	1:H:160:THR:OG1	2.17	0.77
1:F:156:ARG:NH2	1:F:160:THR:OG1	2.17	0.76
2:P:98:SER:OG	2:P:100:ASP:OD1	2.02	0.76
1:A:258:SER:O	1:A:274:ALA:N	2.19	0.76
1:J:177:ARG:NH1	2:Q:40:TYR:OH	2.18	0.75
2:Q:98:SER:OG	2:Q:100:ASP:OD1	2.02	0.75
1:A:133:ARG:NH1	1:C:222:ARG:O	2.20	0.75
2:M:2:SER:N	2:N:88:ASP:O	2.18	0.75
1:C:213:ALA:O	1:C:217:MET:N	2.20	0.75
3:T:1:A:C2'	3:T:2:A:H4'	2.16	0.75
2:O:98:SER:OG	2:O:100:ASP:OD1	2.04	0.74
1:I:177:ARG:NH2	2:S:54:MET:SD	2.61	0.73
1:D:154:ILE:O	1:D:176:GLY:N	2.21	0.73
1:J:212:ASP:OD1	1:J:216:ASN:ND2	2.22	0.73
2:R:7:ARG:NE	2:R:9:ASP:OD2	2.21	0.73
1:F:217:MET:HE3	1:F:218:PHE:CE2	2.24	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:98:SER:OG	2:M:100:ASP:OD1	2.02	0.72
1:K:259:ARG:NH2	1:K:263:SER:O	2.22	0.72
1:J:137:GLN:NE2	3:T:17:G:OP1	2.23	0.71
1:F:45:VAL:HG12	3:T:31:A:H4'	1.72	0.71
3:T:56:A:O2'	3:T:57:G:O4'	2.04	0.71
2:P:102:GLN:OE1	2:Q:57:LYS:NZ	2.23	0.71
1:D:83:TYR:OH	1:D:128:ASN:O	2.08	0.71
1:E:83:TYR:OH	1:E:128:ASN:O	2.07	0.71
1:E:133:ARG:HD3	3:T:34:A:H5''	1.71	0.71
2:S:10:ARG:NH1	2:S:92:ASP:OD1	2.23	0.71
1:J:151:GLU:OE2	1:J:177:ARG:NE	2.24	0.70
2:P:7:ARG:NE	2:P:9:ASP:OD2	2.23	0.70
1:F:68:TYR:OH	1:F:79:HIS:NE2	2.23	0.70
1:E:11:VAL:HG11	1:E:255:VAL:HG21	1.73	0.69
1:G:11:VAL:HG11	1:G:255:VAL:HG21	1.73	0.69
1:D:136:VAL:HG11	1:D:210:PHE:CE2	2.27	0.69
1:A:24:ASP:N	1:A:30:MET:O	2.26	0.69
2:L:98:SER:OG	2:L:100:ASP:OD1	2.07	0.69
2:L:102:GLN:OE1	2:M:57:LYS:NZ	2.23	0.69
1:A:66:ASN:ND2	1:A:112:ASN:O	2.26	0.68
1:A:141:ALA:HB2	1:A:188:ALA:HB2	1.75	0.68
1:C:137:GLN:NE2	3:T:53:C:OP1	2.26	0.68
3:T:59:U:C2'	3:T:60:U:H5'	2.24	0.68
1:B:226:ARG:NH2	3:T:45:U:H5''	2.09	0.68
1:F:133:ARG:HD3	3:T:28:G:H5''	1.76	0.68
1:I:217:MET:HE3	1:I:218:PHE:CE2	2.28	0.68
1:B:212:ASP:OD1	1:B:216:ASN:ND2	2.27	0.67
1:H:13:LEU:CD2	1:H:187:VAL:HG13	2.23	0.67
1:G:133:ARG:NH1	1:J:222:ARG:O	2.28	0.67
1:G:287:LYS:NZ	1:G:289:MET:SD	2.67	0.67
3:T:5:A:O2'	3:T:6:A:H5'	1.95	0.67
1:F:133:ARG:NH1	1:G:222:ARG:O	2.27	0.67
3:T:41:C:O2'	3:T:42:U:O2	2.13	0.67
1:F:11:VAL:HG11	1:F:255:VAL:HG21	1.76	0.67
1:A:211:TRP:O	1:A:215:VAL:HG23	1.94	0.67
2:N:102:GLN:OE1	2:O:53:HIS:ND1	2.28	0.66
2:O:25:GLN:CD	2:O:36:ILE:HG23	2.20	0.66
1:I:239:HIS:HD2	1:I:284:VAL:HG22	1.60	0.66
1:B:151:GLU:OE2	1:B:177:ARG:NE	2.29	0.66
1:B:280:ALA:HB2	1:B:286:VAL:HG23	1.76	0.66
1:C:133:ARG:NH1	1:D:222:ARG:O	2.28	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:11:VAL:HG11	1:I:255:VAL:HG21	1.77	0.66
1:K:155:THR:O	3:T:7:C:H5''	1.96	0.66
2:R:10:ARG:NH1	2:R:92:ASP:OD1	2.28	0.66
1:D:287:LYS:NZ	1:D:289:MET:SD	2.69	0.66
1:G:217:MET:HE3	1:G:218:PHE:CE2	2.31	0.65
1:K:24:ASP:N	1:K:30:MET:O	2.29	0.65
1:C:151:GLU:OE2	1:C:177:ARG:NE	2.29	0.65
1:D:13:LEU:CD2	1:D:187:VAL:HG13	2.26	0.65
1:E:133:ARG:CD	3:T:34:A:H5''	2.27	0.65
1:C:267:ALA:O	1:C:268:ARG:NH1	2.28	0.65
2:Q:7:ARG:NE	2:Q:9:ASP:OD2	2.28	0.65
1:D:136:VAL:HG11	1:D:210:PHE:CZ	2.32	0.64
2:M:69:ASP:O	2:M:73:LYS:N	2.30	0.64
1:K:62:ALA:O	1:K:66:ASN:ND2	2.29	0.64
2:P:111:HIS:NE2	2:Q:87:ILE:O	2.29	0.64
1:A:45:VAL:HG13	3:T:60:U:H3'	1.80	0.64
1:B:110:CYS:O	1:B:117:ARG:NH1	2.31	0.64
1:B:137:GLN:NE2	3:T:41:C:OP1	2.30	0.64
1:G:55:VAL:HG22	1:G:209:LEU:HD22	1.80	0.64
1:F:226:ARG:NH2	3:T:34:A:OP1	2.31	0.64
3:T:44:G:O2'	3:T:45:U:OP2	2.15	0.64
1:D:110:CYS:O	1:D:117:ARG:NH1	2.31	0.63
1:E:155:THR:O	3:T:43:C:H5''	1.98	0.63
1:D:23:GLY:O	3:T:50:U:H3'	1.99	0.63
1:H:45:VAL:HG13	3:T:13:C:H5'	1.80	0.63
1:E:217:MET:HE3	1:E:218:PHE:CE2	2.32	0.63
1:K:223:SER:OG	3:T:4:A:OP1	2.09	0.63
1:J:15:ASP:OD1	1:J:230:SER:OG	2.14	0.63
1:D:11:VAL:HG11	1:D:255:VAL:HG21	1.81	0.63
1:G:212:ASP:OD1	1:G:216:ASN:ND2	2.32	0.62
2:M:23:LYS:HB2	2:M:116:LEU:HD13	1.82	0.62
2:N:10:ARG:NH1	2:N:92:ASP:OD1	2.32	0.62
3:T:1:A:H2'	3:T:1:A:N3	2.14	0.62
1:F:110:CYS:O	1:F:117:ARG:NH1	2.32	0.62
1:A:241:ASN:OD1	1:A:242:ARG:N	2.33	0.62
1:E:156:ARG:NH2	1:E:160:THR:OG1	2.33	0.62
1:B:83:TYR:OH	1:B:128:ASN:O	2.12	0.61
1:B:213:ALA:O	1:B:217:MET:N	2.32	0.61
1:I:221:ASP:OD2	1:I:226:ARG:NH1	2.33	0.61
1:J:103:LYS:NZ	1:J:107:ASP:OD2	2.31	0.61
1:E:137:GLN:NE2	3:T:35:C:OP1	2.31	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:THR:O	3:T:49:C:H5''	2.00	0.61
1:H:287:LYS:NZ	1:H:289:MET:SD	2.73	0.61
2:N:44:ALA:HB2	2:N:51:VAL:HG11	1.81	0.61
1:F:151:GLU:OE2	1:F:177:ARG:NE	2.34	0.61
2:O:111:HIS:NE2	2:P:87:ILE:O	2.28	0.61
1:C:150:GLN:O	1:C:152:VAL:HG23	2.01	0.61
1:E:22:ASN:OD1	1:E:32:ARG:NE	2.31	0.61
1:E:109:MET:HE3	1:E:116:ILE:CD1	2.31	0.61
1:D:213:ALA:O	1:D:217:MET:N	2.34	0.61
2:L:35:THR:OG1	2:L:38:ASP:OD1	2.15	0.61
1:A:34:ASP:O	1:A:38:GLY:N	2.32	0.60
1:K:241:ASN:OD1	1:K:243:LEU:N	2.29	0.60
1:A:122:VAL:H	3:T:59:U:H1'	1.65	0.60
1:B:45:VAL:HG13	3:T:42:U:H3'	1.83	0.60
1:F:22:ASN:ND2	3:T:32:G:OP1	2.34	0.60
1:G:13:LEU:CD2	1:G:187:VAL:HG13	2.31	0.60
3:T:55:C:O3'	3:T:56:A:H4'	2.01	0.60
1:C:22:ASN:OD1	1:C:32:ARG:NE	2.30	0.60
1:G:22:ASN:HB2	1:G:42:VAL:HG13	1.83	0.60
1:J:215:VAL:HG13	1:J:290:LEU:HD22	1.82	0.60
1:B:11:VAL:HG11	1:B:255:VAL:HG21	1.84	0.60
1:B:144:VAL:HG11	1:B:257:VAL:HG11	1.82	0.60
1:F:12:LEU:HD21	1:F:217:MET:HE2	1.82	0.60
1:B:175:MET:SD	2:M:39:ARG:NH2	2.75	0.60
1:E:105:VAL:HG12	1:E:129:CYS:SG	2.42	0.60
2:O:25:GLN:OE1	2:O:36:ILE:HG23	2.02	0.60
3:T:43:C:H5'	3:T:44:G:OP2	2.02	0.60
1:C:183:TYR:OH	1:C:270:PHE:N	2.35	0.59
1:G:133:ARG:HD2	3:T:22:U:H5'	1.84	0.59
1:I:13:LEU:CD2	1:I:187:VAL:HG13	2.31	0.59
1:J:287:LYS:NZ	1:J:289:MET:SD	2.75	0.59
1:D:235:ILE:HG12	1:D:277:VAL:HG21	1.83	0.59
1:H:151:GLU:OE2	1:H:177:ARG:NE	2.36	0.59
1:J:11:VAL:HG11	1:J:255:VAL:HG21	1.84	0.59
2:O:102:GLN:OE1	2:P:57:LYS:NZ	2.32	0.59
1:H:223:SER:OG	3:T:16:C:OP1	2.21	0.59
1:K:287:LYS:NZ	1:K:289:MET:SD	2.76	0.59
1:J:231:SER:O	1:J:290:LEU:HD21	2.03	0.59
1:G:15:ASP:OD1	1:G:230:SER:OG	2.20	0.58
1:I:110:CYS:O	1:I:117:ARG:NH1	2.34	0.58
2:Q:20:VAL:HG11	2:Q:83:MET:HG3	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:ILE:O	1:C:55:VAL:HG23	2.03	0.58
2:N:43:SER:C	2:N:51:VAL:HG21	2.28	0.58
1:K:255:VAL:HG22	1:K:277:VAL:HG22	1.85	0.58
1:H:110:CYS:O	1:H:117:ARG:NH1	2.36	0.58
1:K:50:LYS:NZ	1:K:221:ASP:OD2	2.24	0.58
1:K:49:ARG:HB2	3:T:1:A:H5'	1.85	0.58
1:A:32:ARG:NH2	1:C:153:SER:O	2.37	0.58
1:K:114:TYR:O	1:K:118:THR:N	2.37	0.57
2:Q:10:ARG:NH1	2:Q:92:ASP:OD1	2.37	0.57
1:I:13:LEU:HD21	1:I:187:VAL:HG13	1.86	0.57
1:G:45:VAL:CG1	3:T:25:A:H5'	2.35	0.57
1:J:235:ILE:HG12	1:J:277:VAL:HG21	1.86	0.57
1:C:11:VAL:HG11	1:C:255:VAL:HG21	1.86	0.57
1:E:280:ALA:HB2	1:E:286:VAL:HG23	1.86	0.57
1:G:213:ALA:O	1:G:217:MET:N	2.38	0.57
1:I:156:ARG:NH2	1:I:160:THR:OG1	2.38	0.57
1:A:13:LEU:CD2	1:A:187:VAL:HG13	2.35	0.57
2:P:10:ARG:NH1	2:P:92:ASP:O	2.36	0.57
1:H:215:VAL:HG13	1:H:290:LEU:HD22	1.86	0.57
2:M:51:VAL:O	2:M:54:MET:N	2.38	0.57
1:G:13:LEU:HD21	1:G:187:VAL:HG13	1.86	0.57
1:H:92:PRO:O	1:H:94:LYS:N	2.38	0.57
2:O:114:LYS:N	2:P:84:GLN:OE1	2.38	0.57
1:E:117:ARG:NE	1:E:206:ASP:OD2	2.30	0.56
1:E:213:ALA:O	1:E:217:MET:N	2.37	0.56
1:G:55:VAL:HG22	1:G:209:LEU:CD2	2.35	0.56
1:H:226:ARG:NH2	3:T:15:G:O3'	2.38	0.56
1:A:13:LEU:HD21	1:A:187:VAL:HG13	1.86	0.56
1:I:280:ALA:HB2	1:I:286:VAL:HG23	1.86	0.56
2:M:35:THR:OG1	2:M:38:ASP:OD1	2.22	0.56
1:F:137:GLN:NE2	3:T:29:C:OP1	2.39	0.56
2:Q:3:LEU:HD22	2:Q:115:ALA:HB2	1.87	0.56
1:A:115:ASP:O	1:A:119:PHE:N	2.36	0.56
2:O:20:VAL:HG11	2:O:83:MET:HG3	1.87	0.56
1:A:87:ASP:OD1	1:A:87:ASP:N	2.39	0.56
1:K:241:ASN:OD1	1:K:242:ARG:N	2.38	0.56
2:M:67:ARG:NH2	2:M:80:GLU:OE1	2.39	0.56
2:R:20:VAL:HG21	2:R:86:ILE:HD12	1.88	0.56
1:J:12:LEU:HD21	1:J:217:MET:HE2	1.88	0.56
3:T:3:G:O2'	3:T:4:A:H5'	2.05	0.56
1:D:22:ASN:ND2	3:T:50:U:OP1	2.35	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:45:VAL:HG13	3:T:49:C:H5'	1.88	0.55
1:J:280:ALA:HB2	1:J:286:VAL:HG23	1.87	0.55
1:A:269:SER:OG	1:A:272:ASP:OD1	2.18	0.55
1:B:12:LEU:HD21	1:B:217:MET:HE2	1.89	0.55
1:I:155:THR:O	3:T:13:C:H5''	2.06	0.55
1:K:258:SER:O	1:K:274:ALA:N	2.39	0.55
1:I:287:LYS:NZ	1:I:289:MET:SD	2.80	0.55
2:R:7:ARG:NH1	2:S:91:SER:O	2.39	0.55
2:M:38:ASP:OD1	2:M:38:ASP:N	2.39	0.55
2:P:10:ARG:NH1	2:P:92:ASP:OD1	2.40	0.55
1:F:54:HIS:CE1	1:F:58:THR:HG21	2.42	0.55
1:K:121:ALA:N	1:K:132:VAL:O	2.40	0.55
1:B:215:VAL:HG13	1:B:290:LEU:HD22	1.88	0.55
1:F:151:GLU:OE2	2:O:39:ARG:NH1	2.40	0.55
1:C:152:VAL:N	1:C:178:LYS:O	2.35	0.54
1:D:241:ASN:OD1	1:D:242:ARG:N	2.40	0.54
1:H:280:ALA:HB2	1:H:286:VAL:HG23	1.89	0.54
1:K:46:CYS:HA	3:T:1:A:H4'	1.88	0.54
1:C:217:MET:HE3	1:C:218:PHE:CZ	2.42	0.54
2:M:14:LEU:HD21	2:M:93:PHE:HD1	1.71	0.54
1:D:83:TYR:O	1:D:87:ASP:N	2.41	0.54
1:D:166:GLU:OE1	1:D:166:GLU:N	2.36	0.54
1:E:106:THR:HG22	1:E:129:CYS:O	2.06	0.54
2:Q:102:GLN:OE1	2:R:53:HIS:ND1	2.38	0.54
1:C:241:ASN:OD1	1:C:242:ARG:N	2.40	0.54
2:N:25:GLN:CD	2:N:36:ILE:HG23	2.33	0.54
1:J:83:TYR:OH	1:J:128:ASN:O	2.20	0.54
2:M:20:VAL:HG11	2:M:83:MET:HG3	1.90	0.54
1:B:231:SER:O	1:B:290:LEU:HD21	2.08	0.54
1:G:125:THR:HG21	1:J:160:THR:HG23	1.90	0.54
1:I:15:ASP:OD1	1:I:230:SER:OG	2.20	0.54
1:J:92:PRO:O	1:J:94:LYS:N	2.41	0.54
1:C:45:VAL:HG12	3:T:55:C:H5''	1.89	0.53
1:H:127:VAL:HG12	1:H:127:VAL:O	2.09	0.53
1:A:34:ASP:N	1:A:39:HIS:O	2.36	0.53
1:C:136:VAL:HG22	1:C:192:ILE:HG23	1.91	0.53
1:H:32:ARG:HB2	1:H:41:LEU:O	2.08	0.53
3:T:28:G:H2'	3:T:29:C:H5'	1.89	0.53
1:G:68:TYR:OH	1:G:79:HIS:NE2	2.37	0.53
1:G:157:MET:O	1:G:172:ASN:ND2	2.41	0.53
1:K:34:ASP:OD2	1:K:37:THR:OG1	2.17	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:37:LYS:O	2:M:60:SER:OG	2.26	0.53
2:M:22:GLU:N	2:M:36:ILE:HD11	2.24	0.53
1:A:114:TYR:O	1:A:118:THR:N	2.42	0.53
1:I:13:LEU:HD22	1:I:187:VAL:HG22	1.91	0.53
1:K:16:VAL:O	1:K:16:VAL:HG13	2.09	0.53
1:I:68:TYR:HH	1:I:79:HIS:CE1	2.26	0.53
2:O:22:GLU:N	2:O:36:ILE:HD11	2.24	0.53
1:A:287:LYS:NZ	1:A:289:MET:SD	2.82	0.52
1:C:55:VAL:HG22	1:C:209:LEU:CD2	2.39	0.52
1:E:32:ARG:NH2	1:F:153:SER:O	2.38	0.52
1:G:269:SER:OG	1:G:272:ASP:OD1	2.24	0.52
1:H:22:ASN:HB2	1:H:42:VAL:HG13	1.91	0.52
1:J:45:VAL:CG1	3:T:19:A:H5'	2.38	0.52
1:K:148:VAL:HG11	2:S:96:THR:HG21	1.91	0.52
2:Q:17:LEU:HB2	2:Q:86:ILE:HG21	1.91	0.52
1:B:13:LEU:CD2	1:B:187:VAL:HG13	2.38	0.52
1:C:108:TRP:O	1:C:112:ASN:ND2	2.38	0.52
1:F:105:VAL:HG12	1:F:129:CYS:SG	2.49	0.52
1:H:11:VAL:HG11	1:H:255:VAL:HG21	1.91	0.52
1:G:45:VAL:HG13	3:T:24:G:H3'	1.91	0.52
1:I:212:ASP:OD1	1:I:216:ASN:ND2	2.42	0.52
1:A:20:ASN:ND2	1:A:29:ASN:O	2.40	0.52
1:B:133:ARG:NH1	1:E:222:ARG:O	2.42	0.52
2:N:107:ILE:HD13	2:O:87:ILE:HG23	1.91	0.52
1:D:127:VAL:HG12	1:D:127:VAL:O	2.09	0.52
1:H:237:PHE:CE2	1:H:286:VAL:HG22	2.44	0.52
1:C:51:ILE:HD11	1:C:214:LEU:HD21	1.91	0.52
1:D:155:THR:OG1	3:T:55:C:O2'	2.23	0.52
1:C:55:VAL:HG22	1:C:209:LEU:HD22	1.91	0.52
1:F:115:ASP:OD1	1:F:115:ASP:N	2.41	0.52
1:H:51:ILE:HD11	1:H:214:LEU:CD2	2.39	0.52
2:M:8:THR:HG22	2:M:8:THR:O	2.10	0.52
1:G:11:VAL:HG12	1:G:235:ILE:HB	1.92	0.52
1:I:133:ARG:HD3	3:T:4:A:H4'	1.91	0.52
1:G:215:VAL:HG13	1:G:290:LEU:HD22	1.92	0.52
1:K:42:VAL:HG23	1:K:186:TYR:CD2	2.46	0.52
3:T:58:C:OP2	3:T:58:C:H2'	2.10	0.52
1:I:106:THR:HG22	1:I:129:CYS:O	2.10	0.51
1:F:231:SER:O	1:F:290:LEU:HD21	2.10	0.51
1:I:105:VAL:HG12	1:I:129:CYS:SG	2.50	0.51
1:K:237:PHE:CD1	1:K:286:VAL:HG22	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:148:VAL:HG11	2:L:50:GLN:NE2	2.26	0.51
1:C:272:ASP:N	1:C:272:ASP:OD1	2.44	0.51
1:D:22:ASN:ND2	3:T:51:A:OP1	2.44	0.51
1:A:134:GLY:O	1:A:137:GLN:NE2	2.43	0.51
1:D:152:VAL:HG21	1:D:180:ILE:HD13	1.91	0.51
3:T:50:U:H4'	3:T:51:A:OP1	2.09	0.51
1:F:92:PRO:O	1:F:94:LYS:N	2.43	0.51
1:F:215:VAL:HG13	1:F:290:LEU:HD22	1.93	0.51
1:G:74:ILE:HD13	1:J:162:LYS:HD2	1.93	0.51
1:K:18:ASN:OD1	1:K:180:ILE:HD12	2.11	0.51
1:C:13:LEU:CD2	1:C:187:VAL:HG13	2.41	0.51
1:D:13:LEU:HD21	1:D:187:VAL:HG13	1.91	0.51
2:M:21:LEU:HD21	2:M:56:LEU:HD23	1.93	0.51
1:C:25:PRO:O	2:L:39:ARG:NH1	2.38	0.50
1:A:82:ALA:O	1:A:85:ALA:N	2.43	0.50
1:C:115:ASP:OD1	1:C:116:ILE:N	2.44	0.50
2:N:110:TYR:HB2	2:O:87:ILE:HG21	1.93	0.50
3:T:59:U:H2'	3:T:60:U:C5'	2.38	0.50
1:E:127:VAL:HG12	1:E:127:VAL:O	2.11	0.50
1:F:74:ILE:HD13	1:G:162:LYS:HD2	1.93	0.50
1:H:219:GLU:OE1	1:H:222:ARG:NH2	2.44	0.50
3:T:37:U:H2'	3:T:39:G:H5''	1.92	0.50
1:I:156:ARG:NH1	3:T:12:G:H21	2.10	0.50
1:K:121:ALA:O	1:K:132:VAL:N	2.43	0.50
1:A:82:ALA:O	1:A:86:CYS:N	2.35	0.50
1:B:127:VAL:O	1:B:127:VAL:HG12	2.11	0.50
1:I:115:ASP:O	1:I:119:PHE:N	2.43	0.50
1:D:233:LYS:HD3	1:D:275:VAL:HG11	1.94	0.50
1:B:13:LEU:HD21	1:B:187:VAL:HG13	1.93	0.49
1:F:213:ALA:O	1:F:217:MET:N	2.45	0.49
1:G:237:PHE:CD1	1:G:286:VAL:HG22	2.47	0.49
1:H:51:ILE:HD11	1:H:214:LEU:HD21	1.94	0.49
1:J:217:MET:HE3	1:J:218:PHE:CE2	2.46	0.49
2:Q:36:ILE:HG22	2:Q:40:TYR:CE2	2.47	0.49
3:T:47:U:HO2'	3:T:48:A:H8	1.59	0.49
1:H:13:LEU:HD22	1:H:187:VAL:HG22	1.94	0.49
1:H:194:ALA:HB3	1:H:195:PRO:HD3	1.94	0.49
2:Q:10:ARG:NH1	2:Q:92:ASP:O	2.45	0.49
1:A:114:TYR:OH	1:A:206:ASP:O	2.30	0.49
1:D:89:LYS:O	1:D:91:GLU:N	2.45	0.49
2:R:23:LYS:HB2	2:R:116:LEU:HD13	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:259:ARG:NH2	1:D:265:GLY:O	2.41	0.49
3:T:8:C:O2	3:T:8:C:C2'	2.61	0.49
3:T:43:C:H2'	3:T:45:U:H5'	1.95	0.49
1:E:55:VAL:HG22	1:E:209:LEU:CD2	2.42	0.49
1:I:22:ASN:ND2	3:T:9:G:OP1	2.46	0.49
1:J:11:VAL:HG12	1:J:235:ILE:HB	1.93	0.49
2:O:69:ASP:O	2:O:73:LYS:N	2.44	0.49
2:Q:114:LYS:N	2:R:84:GLN:OE1	2.46	0.49
1:J:106:THR:HG22	1:J:129:CYS:O	2.12	0.49
2:M:14:LEU:HD21	2:M:93:PHE:CD1	2.47	0.49
1:B:92:PRO:O	1:B:94:LYS:N	2.45	0.49
1:G:100:GLU:OE1	1:G:100:GLU:N	2.41	0.49
1:J:156:ARG:NH2	1:J:160:THR:OG1	2.46	0.49
2:O:3:LEU:HD12	2:O:4:ASP:N	2.28	0.49
1:D:108:TRP:NE1	1:D:112:ASN:OD1	2.46	0.49
1:F:237:PHE:CE1	1:F:286:VAL:HG22	2.47	0.49
2:L:67:ARG:NH2	2:L:80:GLU:OE1	2.46	0.49
2:N:21:LEU:HD22	2:N:59:ALA:CB	2.43	0.49
1:B:36:GLU:OE2	1:E:148:VAL:HG21	2.13	0.49
1:J:32:ARG:HB2	1:J:41:LEU:O	2.13	0.49
2:L:97:MET:O	2:L:102:GLN:NE2	2.46	0.49
1:I:115:ASP:OD1	1:I:116:ILE:N	2.45	0.48
1:J:51:ILE:HD11	1:J:214:LEU:HD21	1.96	0.48
3:T:1:A:H2	3:T:2:A:H8	1.59	0.48
1:D:12:LEU:HD21	1:D:217:MET:HE2	1.94	0.48
2:M:21:LEU:HD22	2:M:59:ALA:HB2	1.94	0.48
3:T:37:U:H2'	3:T:39:G:C5'	2.42	0.48
1:B:29:ASN:ND2	3:T:45:U:H5	2.11	0.48
1:A:4:ILE:N	1:A:202:PHE:O	2.44	0.48
1:E:237:PHE:CD1	1:E:286:VAL:HG22	2.49	0.48
1:F:127:VAL:HG12	1:F:127:VAL:O	2.13	0.48
1:G:144:VAL:HG11	1:G:257:VAL:HG11	1.96	0.48
1:G:209:LEU:HD23	1:G:209:LEU:O	2.14	0.48
2:O:38:ASP:OD1	2:O:38:ASP:N	2.46	0.48
3:T:37:U:C2	3:T:39:G:H4'	2.49	0.48
3:T:55:C:O2	3:T:55:C:C2'	2.61	0.48
1:F:133:ARG:CD	3:T:28:G:H5''	2.44	0.48
1:H:217:MET:HE3	1:H:218:PHE:CE2	2.49	0.48
1:A:15:ASP:OD2	1:A:232:ARG:NE	2.39	0.48
1:J:237:PHE:CD2	1:J:286:VAL:HG22	2.49	0.48
1:C:55:VAL:HG21	1:C:114:TYR:CE2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:233:LYS:HD3	1:F:275:VAL:HG11	1.95	0.48
1:G:127:VAL:O	1:G:127:VAL:HG12	2.13	0.48
1:H:223:SER:CB	3:T:16:C:OP1	2.62	0.48
1:A:197:ALA:O	1:A:201:GLY:N	2.36	0.47
1:C:66:ASN:O	1:C:114:TYR:N	2.45	0.47
1:J:105:VAL:HG12	1:J:129:CYS:SG	2.54	0.47
2:L:107:ILE:HD11	2:M:93:PHE:CZ	2.49	0.47
2:P:22:GLU:N	2:P:36:ILE:HD11	2.29	0.47
1:B:175:MET:HE1	1:D:25:PRO:HB3	1.95	0.47
1:I:211:TRP:O	1:I:215:VAL:HG23	2.14	0.47
2:N:44:ALA:N	2:N:51:VAL:HG21	2.29	0.47
1:C:45:VAL:CG1	3:T:55:C:H5'	2.45	0.47
1:E:115:ASP:OD1	1:E:116:ILE:N	2.47	0.47
1:G:29:ASN:OD1	3:T:26:A:O2'	2.32	0.47
2:O:21:LEU:HD22	2:O:59:ALA:HB2	1.96	0.47
1:D:267:ALA:O	1:D:268:ARG:NH1	2.33	0.47
1:E:209:LEU:HD23	1:E:209:LEU:O	2.14	0.47
2:N:23:LYS:HB2	2:N:116:LEU:HD13	1.95	0.47
1:C:154:ILE:N	1:C:176:GLY:O	2.47	0.47
1:H:239:HIS:CD2	1:H:284:VAL:HG22	2.49	0.47
1:B:105:VAL:HG12	1:B:129:CYS:SG	2.55	0.47
1:H:11:VAL:HG12	1:H:235:ILE:HB	1.96	0.47
3:T:23:U:HO2'	3:T:24:G:H8	1.62	0.47
3:T:31:A:N3	3:T:31:A:C2'	2.75	0.47
3:T:53:C:O2'	3:T:54:G:H5'	2.12	0.47
1:D:155:THR:O	3:T:55:C:H3'	2.14	0.47
1:I:103:LYS:NZ	1:I:107:ASP:OD2	2.46	0.47
1:J:13:LEU:CD2	1:J:187:VAL:HG22	2.44	0.47
2:P:84:GLN:O	2:P:88:ASP:N	2.46	0.47
1:B:54:HIS:NE2	1:B:58:THR:HG21	2.30	0.47
1:B:148:VAL:HG13	2:M:50:GLN:NE2	2.29	0.47
1:D:69:ILE:HG21	3:T:49:C:OP1	2.15	0.47
1:E:233:LYS:HD3	1:E:275:VAL:HG11	1.97	0.47
1:F:16:VAL:HG13	1:F:16:VAL:O	2.15	0.47
1:I:22:ASN:ND2	3:T:8:C:OP1	2.43	0.47
1:I:217:MET:HE3	1:I:218:PHE:CZ	2.49	0.47
2:M:107:ILE:HD13	2:N:87:ILE:HG23	1.97	0.47
1:C:153:SER:HA	1:C:176:GLY:O	2.14	0.47
1:F:15:ASP:OD1	1:F:15:ASP:N	2.47	0.47
1:J:55:VAL:HG22	1:J:209:LEU:HD22	1.97	0.47
1:J:127:VAL:HG12	1:J:127:VAL:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:133:ARG:HD2	3:T:46:C:H5'	1.97	0.46
1:F:143:SER:OG	1:F:145:GLU:O	2.31	0.46
1:E:156:ARG:HB2	3:T:42:U:H1'	1.97	0.46
1:H:93:LYS:NZ	1:H:126:GLU:OE1	2.44	0.46
1:I:55:VAL:HG22	1:I:209:LEU:HD21	1.97	0.46
1:I:92:PRO:O	1:I:94:LYS:N	2.49	0.46
1:E:212:ASP:OD1	1:E:216:ASN:ND2	2.48	0.46
2:M:107:ILE:CD1	2:N:87:ILE:HG23	2.45	0.46
2:O:44:ALA:HB2	2:O:51:VAL:HG11	1.96	0.46
1:B:26:ASP:OD2	2:N:39:ARG:NH1	2.48	0.46
1:E:120:GLY:O	3:T:35:C:H4'	2.15	0.46
1:G:136:VAL:HG11	1:G:210:PHE:CZ	2.51	0.46
1:K:55:VAL:HG21	1:K:114:TYR:CD2	2.51	0.46
3:T:50:U:H1'	3:T:51:A:C2	2.51	0.46
1:D:239:HIS:CD2	1:D:284:VAL:HG22	2.50	0.46
1:H:207:LEU:HD21	1:H:211:TRP:CZ2	2.51	0.46
1:J:35:PRO:O	2:Q:50:GLN:NE2	2.49	0.46
1:B:221:ASP:OD2	1:B:226:ARG:NH1	2.48	0.46
1:C:226:ARG:NH2	3:T:57:G:H5'	2.30	0.46
1:D:209:LEU:HD23	1:D:209:LEU:O	2.16	0.46
1:E:92:PRO:O	1:E:94:LYS:N	2.48	0.46
2:O:43:SER:C	2:O:51:VAL:HG21	2.40	0.46
2:Q:21:LEU:HD22	2:Q:59:ALA:HB2	1.96	0.46
2:Q:23:LYS:HB2	2:Q:116:LEU:HD13	1.98	0.46
2:R:20:VAL:HG21	2:R:86:ILE:CD1	2.46	0.46
1:D:13:LEU:HD22	1:D:187:VAL:HG22	1.97	0.46
1:F:221:ASP:OD2	1:F:226:ARG:NH1	2.49	0.46
1:J:217:MET:HE3	1:J:218:PHE:CZ	2.51	0.46
1:A:128:ASN:OD1	1:A:130:GLY:N	2.43	0.46
1:E:7:ARG:O	1:E:239:HIS:ND1	2.46	0.46
1:G:237:PHE:CE1	1:G:286:VAL:HG22	2.51	0.46
1:I:108:TRP:NE1	1:I:112:ASN:OD1	2.49	0.46
1:B:68:TYR:HH	1:B:79:HIS:HE2	1.64	0.46
1:H:110:CYS:SG	1:H:132:VAL:HG21	2.55	0.46
1:K:222:ARG:NH1	1:K:229:MET:O	2.41	0.46
2:M:25:GLN:CD	2:M:36:ILE:HG23	2.41	0.46
3:T:2:A:H1'	3:T:3:G:C8	2.50	0.46
3:T:38:G:H1'	3:T:39:G:C8	2.51	0.46
2:M:22:GLU:CA	2:M:36:ILE:HD11	2.46	0.45
3:T:38:G:H4'	3:T:39:G:H5'	1.98	0.45
1:E:55:VAL:HG22	1:E:209:LEU:HD21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:R:21:LEU:HD22	2:R:59:ALA:HB2	1.99	0.45
1:A:92:PRO:O	1:A:93:LYS:C	2.60	0.45
1:C:91:GLU:HB3	1:C:92:PRO:HD2	1.98	0.45
1:F:280:ALA:HB2	1:F:286:VAL:HG23	1.98	0.45
1:H:109:MET:HE3	1:H:116:ILE:CD1	2.47	0.45
2:L:21:LEU:HD22	2:L:59:ALA:CB	2.47	0.45
2:O:107:ILE:HD11	2:P:93:PHE:CZ	2.51	0.45
1:A:136:VAL:HG11	1:A:210:PHE:CE1	2.52	0.45
1:B:122:VAL:CG2	3:T:40:A:H2'	2.46	0.45
1:B:287:LYS:NZ	1:B:289:MET:SD	2.90	0.45
1:G:133:ARG:CD	3:T:22:U:H5'	2.45	0.45
3:T:8:C:OP1	3:T:9:G:OP1	2.34	0.45
1:H:105:VAL:HG12	1:H:129:CYS:SG	2.57	0.45
1:H:136:VAL:HG11	1:H:210:PHE:CZ	2.51	0.45
1:K:221:ASP:OD1	1:K:226:ARG:NH1	2.49	0.45
2:Q:20:VAL:O	2:Q:23:LYS:N	2.50	0.45
1:D:45:VAL:HG13	3:T:48:A:H3'	1.99	0.45
1:D:222:ARG:NH1	1:D:229:MET:O	2.47	0.45
1:F:234:LEU:HD23	1:F:289:MET:HB2	1.97	0.45
1:A:128:ASN:OD1	1:A:128:ASN:C	2.58	0.45
1:C:156:ARG:NH2	1:C:160:THR:OG1	2.49	0.45
1:E:185:LEU:HD23	1:E:257:VAL:HG11	1.99	0.45
1:E:12:LEU:HD21	1:E:217:MET:HE2	1.99	0.45
2:N:55:LEU:O	2:N:58:ASN:N	2.49	0.45
2:N:79:TYR:O	2:N:83:MET:N	2.36	0.45
1:E:226:ARG:NH2	3:T:39:G:O3'	2.49	0.44
1:I:93:LYS:O	1:I:128:ASN:ND2	2.43	0.44
1:I:181:VAL:O	1:I:182:PRO:C	2.60	0.44
1:J:11:VAL:O	1:J:11:VAL:HG13	2.18	0.44
1:C:222:ARG:NH1	1:C:229:MET:O	2.47	0.44
1:E:233:LYS:NZ	1:E:288:GLU:OE1	2.45	0.44
1:I:83:TYR:O	1:I:87:ASP:N	2.50	0.44
2:R:10:ARG:NH1	2:R:92:ASP:O	2.45	0.44
1:A:33:ILE:HD12	1:A:33:ILE:C	2.42	0.44
1:B:32:ARG:NH2	1:E:153:SER:O	2.42	0.44
1:G:13:LEU:HD22	1:G:187:VAL:HG22	1.99	0.44
1:G:127:VAL:O	1:G:128:ASN:C	2.60	0.44
2:L:44:ALA:HA	2:L:51:VAL:HG21	2.00	0.44
2:P:21:LEU:HD21	2:P:83:MET:HE3	1.98	0.44
2:R:21:LEU:HD22	2:R:59:ALA:CB	2.47	0.44
1:G:136:VAL:HG11	1:G:210:PHE:CE1	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:104:LEU:O	2:M:108:GLY:N	2.43	0.44
1:A:127:VAL:O	1:A:128:ASN:C	2.60	0.44
1:D:51:ILE:HD11	1:D:214:LEU:HD21	1.99	0.44
1:D:127:VAL:O	1:D:128:ASN:C	2.60	0.44
1:H:136:VAL:HG11	1:H:210:PHE:CE2	2.53	0.44
1:I:209:LEU:HD23	1:I:209:LEU:O	2.18	0.44
1:B:226:ARG:NE	3:T:46:C:OP1	2.50	0.44
1:D:68:TYR:HH	1:D:79:HIS:CE1	2.32	0.44
1:D:141:ALA:HB2	1:D:188:ALA:HB2	1.99	0.44
1:D:280:ALA:HB2	1:D:286:VAL:HG23	2.00	0.44
1:F:287:LYS:NZ	1:F:289:MET:SD	2.91	0.44
2:P:20:VAL:HG11	2:P:83:MET:HG3	1.99	0.44
2:R:67:ARG:NH2	2:R:80:GLU:OE1	2.47	0.44
1:C:128:ASN:C	1:C:128:ASN:OD1	2.61	0.44
1:D:92:PRO:O	1:D:94:LYS:N	2.51	0.44
1:F:22:ASN:HB2	1:F:42:VAL:HG13	1.99	0.44
1:K:257:VAL:HA	1:K:274:ALA:O	2.18	0.44
1:B:15:ASP:OD1	1:B:230:SER:OG	2.29	0.44
1:I:237:PHE:CE1	1:I:286:VAL:HG22	2.53	0.44
1:K:25:PRO:HG3	3:T:2:A:H2'	2.00	0.44
1:B:125:THR:HG21	1:E:160:THR:HG23	1.99	0.43
1:D:106:THR:HG22	1:D:129:CYS:O	2.17	0.43
1:F:22:ASN:ND2	1:F:46:CYS:SG	2.91	0.43
1:F:83:TYR:OH	1:F:128:ASN:O	2.25	0.43
1:H:41:LEU:HA	1:H:141:ALA:O	2.18	0.43
1:J:123:MET:HE3	1:J:132:VAL:HG23	2.00	0.43
1:J:127:VAL:O	1:J:128:ASN:C	2.61	0.43
2:Q:22:GLU:N	2:Q:36:ILE:HD11	2.33	0.43
1:B:122:VAL:HG22	3:T:40:A:H2'	1.98	0.43
1:B:209:LEU:HD23	1:B:209:LEU:O	2.18	0.43
1:E:12:LEU:HD13	1:E:234:LEU:HD13	2.00	0.43
1:H:133:ARG:CD	3:T:10:C:H5''	2.49	0.43
1:H:155:THR:N	3:T:20:A:OP2	2.44	0.43
1:J:13:LEU:CD2	1:J:187:VAL:HG13	2.47	0.43
2:R:22:GLU:CA	2:R:36:ILE:HD11	2.48	0.43
3:T:37:U:C2'	3:T:39:G:H5''	2.48	0.43
1:F:144:VAL:HG11	1:F:257:VAL:HG11	2.00	0.43
1:G:105:VAL:HG12	1:G:129:CYS:SG	2.58	0.43
1:G:231:SER:O	1:G:290:LEU:HD21	2.19	0.43
1:H:13:LEU:HD21	1:H:187:VAL:HG13	1.98	0.43
1:K:157:MET:N	3:T:7:C:OP1	2.39	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:226:ARG:NH2	3:T:3:G:H4'	2.33	0.43
3:T:47:U:O2'	3:T:48:A:H8	2.00	0.43
1:C:127:VAL:O	1:C:128:ASN:C	2.61	0.43
1:K:211:TRP:O	1:K:215:VAL:HG23	2.18	0.43
2:P:51:VAL:O	2:P:54:MET:N	2.51	0.43
1:A:22:ASN:OD1	1:A:32:ARG:NE	2.50	0.43
1:A:64:ARG:O	1:A:66:ASN:N	2.48	0.43
1:G:32:ARG:HB2	1:G:41:LEU:O	2.18	0.43
1:J:251:LEU:HD21	1:J:284:VAL:HG21	1.99	0.43
2:N:22:GLU:HA	2:N:36:ILE:HD11	2.00	0.43
2:N:38:ASP:OD1	2:N:38:ASP:N	2.51	0.43
3:T:11:U:O2'	3:T:12:G:O5'	2.36	0.43
1:E:45:VAL:CG1	3:T:37:U:H5'	2.48	0.43
1:G:122:VAL:HG22	3:T:22:U:H2'	2.01	0.43
1:G:226:ARG:NH2	3:T:28:G:OP1	2.51	0.43
1:K:141:ALA:HB2	1:K:188:ALA:HB2	2.01	0.43
2:L:21:LEU:HD22	2:L:59:ALA:HB2	2.00	0.43
3:T:43:C:O2	3:T:43:C:H3'	2.17	0.43
1:B:51:ILE:O	1:B:55:VAL:HG23	2.18	0.43
1:D:185:LEU:HD23	1:D:257:VAL:HG11	2.00	0.43
1:H:52:ARG:HE	1:H:115:ASP:HB2	1.83	0.43
2:P:21:LEU:HD13	2:P:55:LEU:HB3	2.00	0.43
1:A:45:VAL:CG1	3:T:60:U:H3'	2.47	0.43
1:A:247:PRO:HB2	1:A:249:HIS:CE1	2.53	0.43
1:E:237:PHE:CE1	1:E:286:VAL:HG22	2.54	0.43
1:G:8:TYR:CE2	1:G:238:LYS:HD2	2.54	0.43
1:H:35:PRO:O	2:R:50:GLN:NE2	2.52	0.43
3:T:2:A:C2'	3:T:3:G:OP2	2.65	0.43
1:A:77:GLU:O	1:A:80:GLU:N	2.51	0.43
1:B:153:SER:O	1:D:32:ARG:NH2	2.36	0.43
1:G:11:VAL:CG1	1:G:235:ILE:HD13	2.49	0.43
1:H:209:LEU:O	1:H:209:LEU:HD23	2.19	0.43
1:I:51:ILE:HD11	1:I:214:LEU:CD2	2.48	0.43
1:B:237:PHE:CE2	1:B:286:VAL:HG22	2.54	0.43
1:I:241:ASN:OD1	1:I:242:ARG:N	2.52	0.43
1:J:16:VAL:O	1:J:16:VAL:HG13	2.18	0.43
1:J:207:LEU:HD21	1:J:211:TRP:CH2	2.54	0.43
2:L:20:VAL:HG11	2:L:83:MET:HG3	2.00	0.43
1:B:211:TRP:O	1:B:215:VAL:HG23	2.19	0.42
1:K:237:PHE:HD1	1:K:286:VAL:HG22	1.83	0.42
1:B:185:LEU:HD23	1:B:257:VAL:HG11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:45:VAL:HG21	3:T:12:G:H2'	2.02	0.42
2:L:107:ILE:HD13	2:M:87:ILE:HG23	2.01	0.42
2:L:110:TYR:HB2	2:M:87:ILE:HG21	2.00	0.42
1:E:74:ILE:HD13	1:F:162:LYS:HD2	2.01	0.42
1:J:22:ASN:OD1	1:J:32:ARG:NE	2.48	0.42
3:T:5:A:N6	3:T:6:A:N6	2.67	0.42
1:B:45:VAL:CG1	3:T:42:U:H3'	2.48	0.42
1:E:16:VAL:O	1:E:183:TYR:HA	2.18	0.42
1:E:211:TRP:O	1:E:215:VAL:HG23	2.20	0.42
1:I:209:LEU:HD23	1:I:209:LEU:C	2.44	0.42
1:J:241:ASN:OD1	1:J:242:ARG:N	2.52	0.42
1:A:42:VAL:HG23	1:A:186:TYR:CE2	2.53	0.42
1:A:115:ASP:OD1	1:A:116:ILE:N	2.52	0.42
1:B:217:MET:HE3	1:B:218:PHE:CZ	2.55	0.42
1:D:12:LEU:HD12	1:D:233:LYS:O	2.19	0.42
1:K:235:ILE:HG12	1:K:277:VAL:HG21	2.02	0.42
2:O:21:LEU:C	2:O:36:ILE:HD11	2.45	0.42
2:R:107:ILE:HD11	2:S:93:PHE:CZ	2.55	0.42
1:F:211:TRP:CZ2	1:F:236:VAL:HG11	2.54	0.42
1:G:280:ALA:HB2	1:G:286:VAL:HG23	2.01	0.42
1:B:32:ARG:HB2	1:B:41:LEU:O	2.20	0.42
1:C:207:LEU:HD21	1:C:211:TRP:CZ2	2.55	0.42
1:G:45:VAL:HG11	3:T:25:A:H5'	2.01	0.42
1:B:143:SER:OG	1:B:145:GLU:O	2.36	0.42
1:E:110:CYS:O	1:E:117:ARG:NH1	2.48	0.42
1:J:239:HIS:CD2	1:J:284:VAL:HG22	2.55	0.42
1:K:45:VAL:CG1	3:T:1:A:H5''	2.43	0.42
1:D:11:VAL:CG1	1:D:255:VAL:HG21	2.50	0.42
1:D:55:VAL:HG22	1:D:209:LEU:CD2	2.50	0.42
1:D:217:MET:HE3	1:D:218:PHE:CZ	2.54	0.42
2:L:47:ASN:O	2:L:51:VAL:HG23	2.20	0.42
1:A:123:MET:O	1:A:128:ASN:HA	2.20	0.42
1:C:207:LEU:HD21	1:C:211:TRP:CH2	2.54	0.42
1:E:127:VAL:O	1:E:128:ASN:C	2.63	0.42
2:L:75:SER:O	2:L:79:TYR:N	2.38	0.42
1:D:93:LYS:O	1:D:128:ASN:CB	2.68	0.41
1:E:231:SER:O	1:E:290:LEU:HD21	2.20	0.41
1:K:8:TYR:O	1:K:192:ILE:N	2.53	0.41
1:K:238:LYS:O	1:K:284:VAL:HA	2.20	0.41
2:P:36:ILE:HG22	2:P:40:TYR:CD2	2.55	0.41
3:T:59:U:C4	3:T:60:U:C4	3.08	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:VAL:O	1:B:128:ASN:C	2.62	0.41
1:B:132:VAL:CG1	1:B:200:THR:HG22	2.50	0.41
1:C:194:ALA:N	1:C:195:PRO:CD	2.83	0.41
1:H:32:ARG:NE	1:H:43:THR:HG23	2.35	0.41
1:H:36:GLU:OE2	1:I:148:VAL:HG11	2.20	0.41
1:I:54:HIS:CE1	1:I:58:THR:HG21	2.55	0.41
1:K:209:LEU:C	1:K:209:LEU:HD23	2.45	0.41
2:L:100:ASP:OD1	2:L:100:ASP:N	2.53	0.41
2:N:22:GLU:CA	2:N:36:ILE:HD11	2.49	0.41
2:N:45:SER:HB2	2:O:57:LYS:HZ3	1.85	0.41
3:T:33:C:C4	3:T:34:A:N6	2.88	0.41
3:T:37:U:N1	3:T:39:G:H5'	2.35	0.41
1:A:76:ASN:OD1	1:C:160:THR:HG22	2.20	0.41
1:C:152:VAL:O	1:C:178:LYS:N	2.37	0.41
1:F:44:ASP:OD1	1:F:44:ASP:N	2.53	0.41
1:G:44:ASP:OD1	1:G:44:ASP:N	2.53	0.41
1:G:217:MET:HE3	1:G:218:PHE:CZ	2.55	0.41
1:H:127:VAL:O	1:H:128:ASN:C	2.63	0.41
1:J:55:VAL:HG22	1:J:209:LEU:CD2	2.51	0.41
1:K:44:ASP:OD1	1:K:45:VAL:N	2.53	0.41
2:P:22:GLU:CA	2:P:36:ILE:HD11	2.50	0.41
2:R:19:ALA:HB1	2:R:113:ARG:HD2	2.01	0.41
1:B:135:PRO:HD3	1:B:200:THR:HG21	2.02	0.41
1:B:207:LEU:HD21	1:B:211:TRP:CH2	2.56	0.41
1:C:92:PRO:O	1:C:93:LYS:C	2.63	0.41
1:D:55:VAL:HG22	1:D:209:LEU:HD21	2.02	0.41
1:I:32:ARG:NH2	1:K:153:SER:O	2.54	0.41
1:J:136:VAL:HG11	1:J:210:PHE:CZ	2.54	0.41
2:L:41:LEU:HD23	2:M:60:SER:HB2	2.02	0.41
1:A:86:CYS:SG	1:A:105:VAL:HG22	2.60	0.41
1:A:209:LEU:HD23	1:A:209:LEU:O	2.20	0.41
1:F:209:LEU:O	1:F:209:LEU:HD23	2.20	0.41
1:H:45:VAL:HG12	3:T:13:C:H5'	1.97	0.41
1:H:241:ASN:OD1	1:H:242:ARG:N	2.53	0.41
3:T:50:U:O2	3:T:50:U:C2'	2.69	0.41
1:A:13:LEU:HD22	1:A:187:VAL:HG22	2.03	0.41
1:J:8:TYR:CZ	1:J:238:LYS:HD2	2.56	0.41
1:K:42:VAL:HG23	1:K:186:TYR:CE2	2.55	0.41
2:N:69:ASP:O	2:N:73:LYS:N	2.51	0.41
3:T:1:A:H2'	3:T:2:A:C4'	2.37	0.41
3:T:50:U:H1'	3:T:51:A:N3	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:MET:O	1:A:128:ASN:OD1	2.39	0.41
1:G:239:HIS:CD2	1:G:284:VAL:HG22	2.55	0.41
1:J:5:ALA:O	1:J:6:ASN:ND2	2.54	0.41
1:J:45:VAL:HG13	3:T:18:A:H3'	2.01	0.41
2:P:3:LEU:HD12	2:P:4:ASP:N	2.36	0.41
3:T:49:C:N4	3:T:51:A:H1'	2.35	0.41
1:A:128:ASN:OD1	1:A:129:CYS:N	2.54	0.41
1:H:32:ARG:NH2	1:I:153:SER:O	2.48	0.41
1:I:55:VAL:HG21	1:I:114:TYR:CE1	2.55	0.41
1:J:226:ARG:NE	3:T:22:U:OP1	2.54	0.41
3:T:2:A:H1'	3:T:3:G:P	2.61	0.41
1:B:136:VAL:HG11	1:B:210:PHE:CZ	2.56	0.41
1:D:66:ASN:HB2	1:D:112:ASN:O	2.20	0.41
1:F:13:LEU:CD2	1:F:187:VAL:HG22	2.51	0.41
1:F:147:VAL:HB	1:F:181:VAL:HG13	2.01	0.41
1:G:152:VAL:HG21	1:G:180:ILE:HD13	2.03	0.41
1:K:210:PHE:CE2	1:K:214:LEU:HD11	2.56	0.41
2:P:40:TYR:CG	2:P:55:LEU:HD21	2.56	0.41
2:R:23:LYS:HG2	2:R:116:LEU:HD22	2.02	0.41
3:T:29:C:O2'	3:T:30:C:P	2.79	0.41
1:D:109:MET:HE3	1:D:116:ILE:HD12	2.03	0.41
1:D:226:ARG:NE	3:T:52:G:OP1	2.54	0.41
1:K:19:GLY:O	1:K:181:VAL:N	2.47	0.41
1:K:239:HIS:CD2	1:K:284:VAL:HG22	2.56	0.41
2:P:82:MET:SD	2:P:112:GLN:NE2	2.94	0.41
3:T:5:A:C2'	3:T:6:A:H5'	2.51	0.41
1:C:127:VAL:O	1:C:127:VAL:HG12	2.20	0.40
1:E:13:LEU:CD2	1:E:187:VAL:HG22	2.50	0.40
1:H:51:ILE:HD13	1:H:210:PHE:CE2	2.56	0.40
1:K:16:VAL:HG23	1:K:229:MET:HG2	2.03	0.40
1:A:51:ILE:O	1:A:55:VAL:HG23	2.21	0.40
1:A:122:VAL:HB	3:T:59:U:C6	2.56	0.40
1:C:75:LEU:HD21	1:D:158:ALA:HB3	2.02	0.40
1:E:32:ARG:HB2	1:E:41:LEU:O	2.21	0.40
1:E:134:GLY:O	1:E:137:GLN:NE2	2.53	0.40
1:J:251:LEU:HD22	1:J:284:VAL:HG11	2.03	0.40
2:L:3:LEU:HD12	2:L:4:ASP:N	2.36	0.40
3:T:19:A:H2'	3:T:21:U:OP1	2.21	0.40
3:T:55:C:O2'	3:T:55:C:O2	2.39	0.40
1:C:23:GLY:H	3:T:57:G:P	2.43	0.40
1:D:156:ARG:NH2	1:D:160:THR:OG1	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:16:VAL:O	1:F:183:TYR:HA	2.20	0.40
1:J:51:ILE:HD13	1:J:210:PHE:CE1	2.57	0.40
1:J:213:ALA:O	1:J:217:MET:N	2.55	0.40
1:K:10:PHE:HB2	1:K:236:VAL:HG22	2.04	0.40
2:N:25:GLN:OE1	2:N:36:ILE:HG23	2.22	0.40
2:P:23:LYS:HB2	2:P:116:LEU:HD13	2.02	0.40
1:A:83:TYR:O	1:A:88:LEU:N	2.54	0.40
1:B:16:VAL:O	1:B:183:TYR:HA	2.20	0.40
1:B:22:ASN:OD1	1:B:32:ARG:NE	2.49	0.40
1:G:133:ARG:HD3	3:T:22:U:H4'	2.03	0.40
1:J:45:VAL:CG1	3:T:19:A:C5'	2.99	0.40
1:J:194:ALA:N	1:J:195:PRO:CD	2.84	0.40
1:J:226:ARG:NH2	3:T:22:U:OP1	2.55	0.40
1:K:8:TYR:CE2	1:K:238:LYS:HD3	2.56	0.40
1:K:127:VAL:O	1:K:128:ASN:C	2.64	0.40
2:M:102:GLN:O	2:M:105:PHE:HB3	2.21	0.40
3:T:59:U:O2	3:T:59:U:O4'	2.40	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	282/289 (98%)	263 (93%)	19 (7%)	0	100	100
1	B	282/289 (98%)	263 (93%)	18 (6%)	1 (0%)	30	62
1	C	282/289 (98%)	247 (88%)	35 (12%)	0	100	100
1	D	282/289 (98%)	258 (92%)	24 (8%)	0	100	100
1	E	282/289 (98%)	262 (93%)	20 (7%)	0	100	100
1	F	282/289 (98%)	256 (91%)	26 (9%)	0	100	100
1	G	282/289 (98%)	266 (94%)	16 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	282/289 (98%)	256 (91%)	26 (9%)	0	100	100
1	I	282/289 (98%)	252 (89%)	30 (11%)	0	100	100
1	J	282/289 (98%)	257 (91%)	25 (9%)	0	100	100
1	K	282/289 (98%)	259 (92%)	23 (8%)	0	100	100
2	L	110/116 (95%)	109 (99%)	1 (1%)	0	100	100
2	M	110/116 (95%)	108 (98%)	2 (2%)	0	100	100
2	N	110/116 (95%)	109 (99%)	1 (1%)	0	100	100
2	O	110/116 (95%)	110 (100%)	0	0	100	100
2	P	110/116 (95%)	110 (100%)	0	0	100	100
2	Q	110/116 (95%)	107 (97%)	3 (3%)	0	100	100
2	R	110/116 (95%)	109 (99%)	1 (1%)	0	100	100
2	S	110/116 (95%)	110 (100%)	0	0	100	100
All	All	3982/4107 (97%)	3711 (93%)	270 (7%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	93	LYS

### 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	239/241 (99%)	234 (98%)	5 (2%)	47	74
1	B	239/241 (99%)	237 (99%)	2 (1%)	73	86
1	C	239/241 (99%)	236 (99%)	3 (1%)	61	81
1	D	239/241 (99%)	238 (100%)	1 (0%)	84	90
1	E	239/241 (99%)	238 (100%)	1 (0%)	84	90
1	F	239/241 (99%)	239 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	239/241 (99%)	237 (99%)	2 (1%)	73	86
1	H	239/241 (99%)	234 (98%)	5 (2%)	47	74
1	I	239/241 (99%)	239 (100%)	0	100	100
1	J	239/241 (99%)	237 (99%)	2 (1%)	73	86
1	K	239/241 (99%)	237 (99%)	2 (1%)	73	86
2	L	96/97 (99%)	96 (100%)	0	100	100
2	M	96/97 (99%)	95 (99%)	1 (1%)	68	83
2	N	96/97 (99%)	96 (100%)	0	100	100
2	O	96/97 (99%)	96 (100%)	0	100	100
2	P	96/97 (99%)	96 (100%)	0	100	100
2	Q	96/97 (99%)	96 (100%)	0	100	100
2	R	96/97 (99%)	96 (100%)	0	100	100
2	S	96/97 (99%)	96 (100%)	0	100	100
All	All	3397/3427 (99%)	3373 (99%)	24 (1%)	73	87

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLU
1	A	78	THR
1	A	87	ASP
1	A	112	ASN
1	A	128	ASN
1	B	156	ARG
1	B	238	LYS
1	C	57	LEU
1	C	191	PHE
1	C	272	ASP
1	D	57	LEU
1	E	238	LYS
1	G	133	ARG
1	G	191	PHE
1	H	33	ILE
1	H	156	ARG
1	H	191	PHE
1	H	238	LYS
1	H	242	ARG

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Mol	Chain	Res	Type
1	J	191	PHE
1	J	238	LYS
1	K	191	PHE
1	K	206	ASP
2	M	38	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	ASN
1	A	131	GLN
1	B	66	ASN
1	B	150	GLN
1	B	189	HIS
1	D	189	HIS
1	D	239	HIS
1	E	20	ASN
1	E	150	GLN
1	E	179	HIS
1	F	54	HIS
1	F	150	GLN
1	F	179	HIS
1	F	245	ASN
1	H	20	ASN
1	H	150	GLN
1	I	150	GLN
1	I	179	HIS
1	I	220	HIS
1	J	150	GLN
1	J	189	HIS
1	J	245	ASN
1	K	189	HIS
2	L	50	GLN
2	M	50	GLN
2	M	62	HIS
2	N	62	HIS

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	T	59/124 (47%)	34 (57%)	3 (5%)

All (34) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	T	2	A
3	T	3	G
3	T	8	C
3	T	12	G
3	T	13	C
3	T	14	U
3	T	18	A
3	T	20	A
3	T	21	U
3	T	22	U
3	T	23	U
3	T	24	G
3	T	26	A
3	T	27	C
3	T	29	C
3	T	30	C
3	T	32	G
3	T	36	A
3	T	38	G
3	T	39	G
3	T	40	A
3	T	42	U
3	T	44	G
3	T	45	U
3	T	48	A
3	T	49	C
3	T	50	U
3	T	51	A
3	T	54	G
3	T	55	C
3	T	56	A
3	T	57	G
3	T	58	C
3	T	59	U

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	T	2	A
3	T	50	U
3	T	56	A

## 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 oligosaccharides 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 Map visualisation

This section contains visualisations of the EMDB entry EMD-61479. These allow visual inspection of the internal detail of the map and identification of artifacts.

Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections

This section was not generated.

### 6.2 Central slices

This section was not generated.

### 6.3 Largest variance slices

This section was not generated.

### 6.4 Orthogonal standard-deviation projections (False-color)

This section was not generated.

### 6.5 Orthogonal surface views

This section was not generated.

### 6.6 Mask visualisation

This section was not generated. No masks/segmentation were deposited.

## 7 Map analysis ⓘ

This section contains the results of statistical analysis of the map.

### 7.1 Map-value distribution ⓘ

This section was not generated.

### 7.2 Volume estimate versus contour level ⓘ

This section was not generated.

### 7.3 Rotationally averaged power spectrum ⓘ

This section was not generated. The rotationally averaged power spectrum had issues being displayed.

## 8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

## 9 Map-model fit

This section was not generated.