



# Full wwPDB/EMDatabank EM Map/Model Validation Report ⓘ

Jan 22, 2018 – 12:03 PM EST

PDB ID : 6F41  
EMDB ID: : EMD-4181  
Title : RNA Polymerase III initially transcribing complex  
Authors : Vorlaender, M.K.; Khatter, H.; Wetzels, R.; Hagen, W.J.H.; Mueller, C.W.  
Deposited on : 2017-11-29  
Resolution : 4.30 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/EMValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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MolProbity : 4.02b-467  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030736

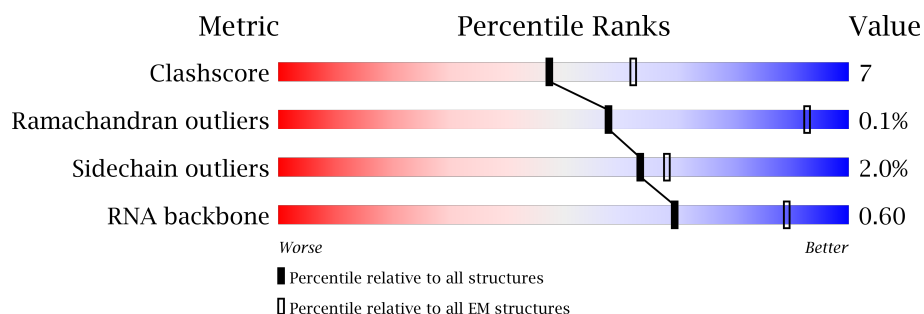
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 4.30 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	125131	1336
Ramachandran outliers	121729	1120
Sidechain outliers	121581	1026
RNA backbone	3398	335

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	1460	78% 17% . .
2	B	1149	75% 21% . .
3	C	335	85% 14% .
4	D	161	61% 12% 26%
5	E	215	80% 20%
6	F	155	45% 8% 46%
7	G	212	69% 16% 15%
8	H	146	78% 18% .

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Mol	Chain	Length	Quality of chain
9	I	110	
10	J	70	
11	K	142	
12	L	70	
13	M	282	
14	N	422	
15	O	654	
16	P	317	
17	Q	251	
18	U	240	
19	V	596	
20	W	594	
21	R	6	
22	X	81	
23	Y	81	

## 2 Entry composition

There are 24 unique types of molecules in this entry. The entry contains 47788 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 DNA-directed RNA polymerase III subunit RPC1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1398	Total	C	N	O	S	0	0
			10972	6919	1936	2059	58		

- Molecule 2 is a protein called DNA-directed RNA polymerase III subunit RPC2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	1114	Total	C	N	O	S	0	0
			8788	5558	1516	1654	60		

- Molecule 3 is a protein called DNA-directed RNA polymerases I and III subunit RPAC1.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	335	Total	C	N	O	S	0	0
			2655	1681	454	511	9		

- Molecule 4 is a protein called DNA-directed RNA polymerase III subunit RPC9.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	119	Total	C	N	O	S	0	0
			977	628	156	187	6		

- Molecule 5 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC1.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	214	Total	C	N	O	S	0	0
			1751	1111	309	320	11		

- Molecule 6 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC2.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	83	Total	C	N	O	S	0	0
			671	429	114	125	3		

- Molecule 7 is a protein called DNA-directed RNA polymerase III subunit RPC8.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	180	Total	C	N	O	S	0	0
			1448	950	231	261	6		

- Molecule 8 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC3.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	140	Total	C	N	O	S	0	0
			1120	703	188	224	5		

- Molecule 9 is a protein called DNA-directed RNA polymerase III subunit RPC10.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	34	Total	C	N	O	S	0	0
			255	161	39	49	6		

- Molecule 10 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC5.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	67	Total	C	N	O	S	0	0
			549	350	95	98	6		

- Molecule 11 is a protein called DNA-directed RNA polymerases I and III subunit RPAC2.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	101	Total	C	N	O	S	0	0
			792	496	130	161	5		

- Molecule 12 is a protein called DNA-directed RNA polymerases I, II, and III subunit RPABC4.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	45	Total	C	N	O	S	0	0
			358	221	71	62	4		

- Molecule 13 is a protein called DNA-directed RNA polymerase III subunit RPC5.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	183	Total	C	N	O	S	0	0
			1484	942	257	283	2		

- Molecule 14 is a protein called DNA-directed RNA polymerase III subunit RPC4.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	104	Total	C	N	O	S	0	0
			797	505	143	146	3		

- Molecule 15 is a protein called DNA-directed RNA polymerase III subunit RPC3.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	O	534	Total	C	N	O	S	0	0
			4293	2733	736	806	18		

- Molecule 16 is a protein called DNA-directed RNA polymerase III subunit RPC6.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	P	246	Total	C	N	O	S	0	0
			1990	1276	326	377	11		

- Molecule 17 is a protein called DNA-directed RNA polymerase III subunit RPC7.

Mol	Chain	Residues	Atoms				AltConf	Trace
17	Q	35	Total	C	N	O	0	0
			273	181	45	47		

- Molecule 18 is a protein called TATA-box-binding protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	U	180	Total	C	N	O	S	0	0
			1416	921	242	247	6		

- Molecule 19 is a protein called Transcription factor IIIB 70 kDa subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	V	337	Total	C	N	O	S	0	0
			2686	1682	486	504	14		

- Molecule 20 is a protein called Transcription factor TFIIB component B”.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	W	218	Total	C	N	O	S	0	0
			1798	1131	319	339	9		

- Molecule 21 is a RNA chain called RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	R	6	Total	C	N	O	P	0	0
			126	56	20	44	6		

- Molecule 22 is a DNA chain called Non-Template DNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	X	60	Total	C	N	O	P	0	0
			1221	588	204	369	60		

- Molecule 23 is a DNA chain called Template-DNA.

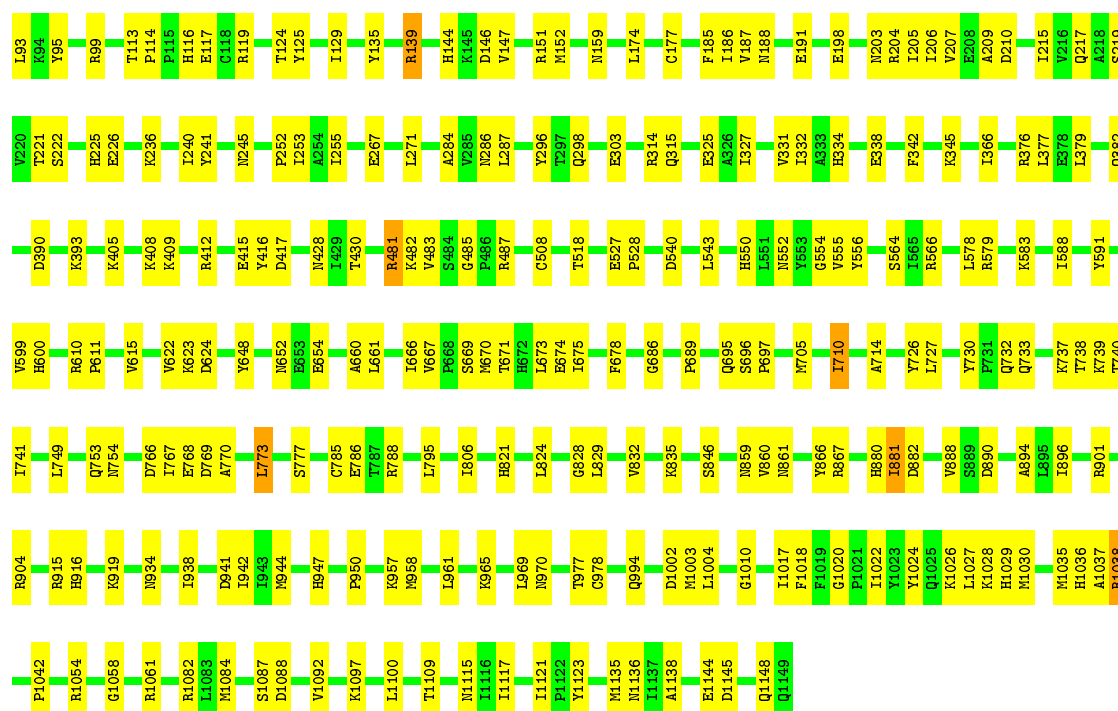
Mol	Chain	Residues	Atoms					AltConf	Trace
23	Y	66	Total	C	N	O	P	0	0
			1361	649	260	386	66		

- Molecule 24 is ZINC ION (three-letter code: ZN) (formula: Zn).

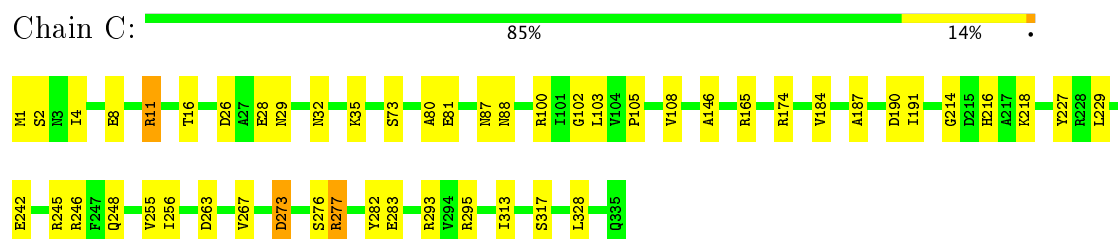
Mol	Chain	Residues	Atoms		AltConf
24	J	1	Total	Zn	0
			1	1	
24	B	1	Total	Zn	0
			1	1	
24	I	1	Total	Zn	0
			1	1	
24	V	1	Total	Zn	0
			1	1	
24	A	2	Total	Zn	0
			2	2	
24	L	1	Total	Zn	0
			1	1	







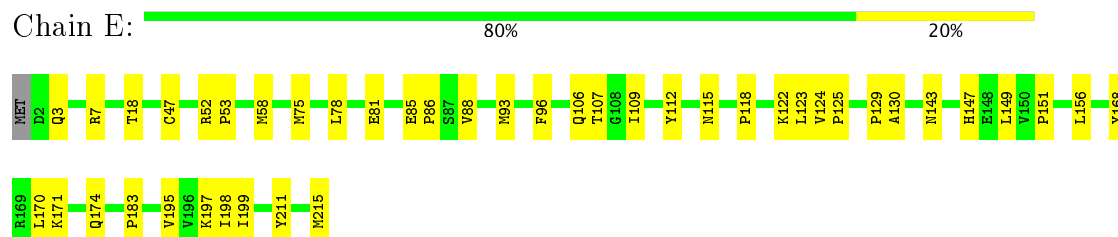
- Molecule 3: DNA-directed RNA polymerases I and III subunit RPAC1



- Molecule 4: DNA-directed RNA polymerase III subunit RPC9



- Molecule 5: DNA-directed RNA polymerases I, II, and III subunit RPABC1



- |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |     |
|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| MET | THR | GLU | ASP | ILE | GLU | GLN | LYS | LYS | LYS | THR | ALA | THR | VAL | THR | PRO | GLN | GLY | PHE | PRO | LYS | HIS | ILE | GLN | GLY | GLU | GLY | GLU | GLN | ASP | VAL | ASP | MET |
| A67 | S89 | F83 | G64 | V66 | E67 | NH4 | V79 | D86 | K61 | R73 | T62 | C63 | A61 | S62 | S65 | S60 | A61 | S62 | F63 | Q64 | I65 | V66 | E67 | NH4 | V79 | D86 | K61 | R73 | T62 | C63 | A61 |     |



Response	Percentage
Yes	67%
No	14%
Don't know	18%



Category	Percentage
Very good	59%
Good	17%
Not good	•
Very bad	22%

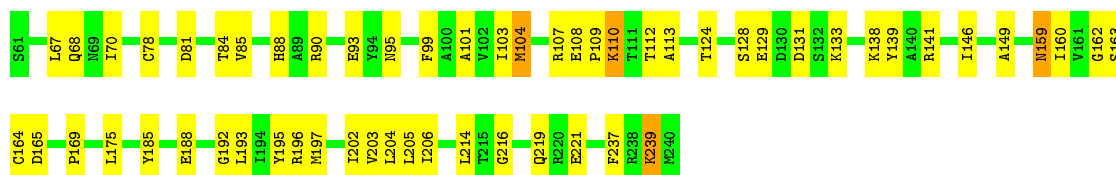


Response	Percentage
Yes, the U.S. is a democracy	11%
No, the U.S. is not a democracy	86%

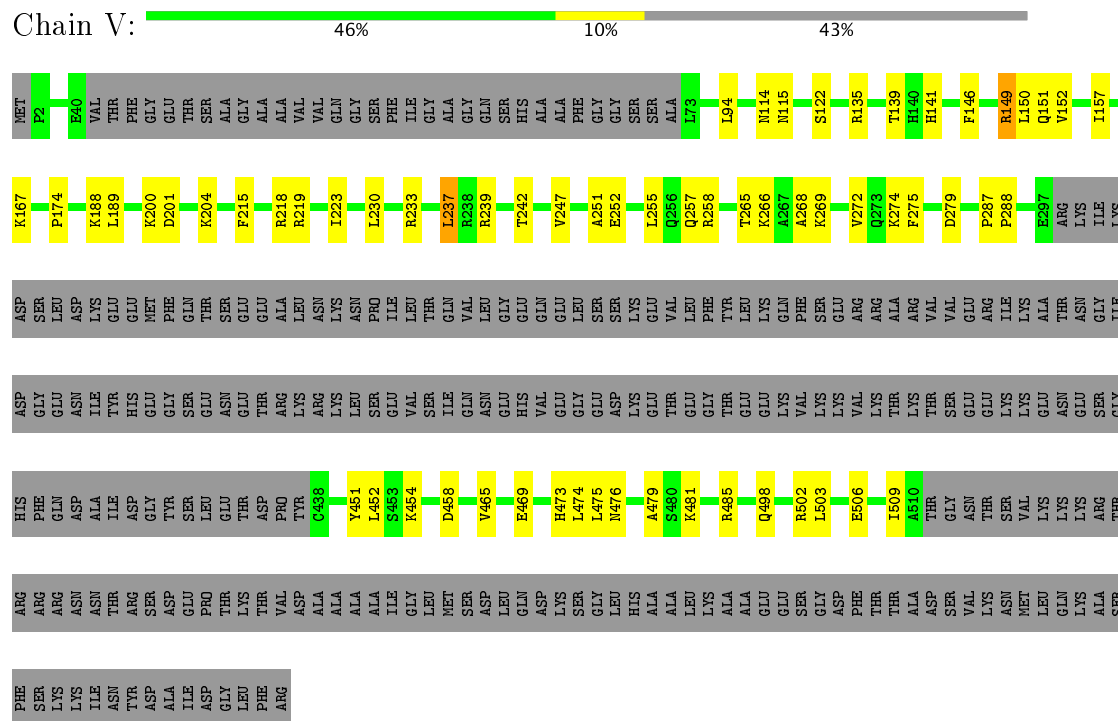


Response	Percentage
Yes	51%
No	22%
Don't know	25%

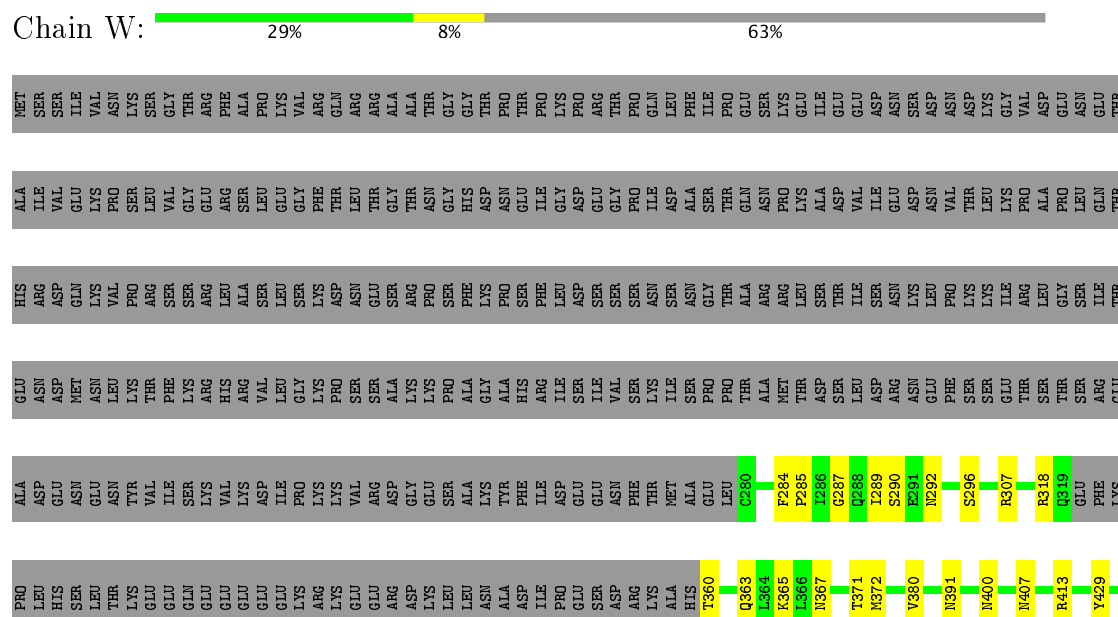


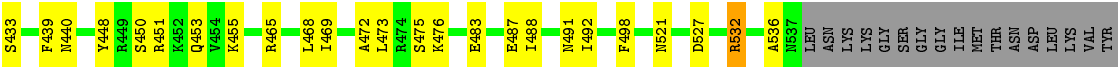


- Molecule 19: Transcription factor IIIB 70 kDa subunit



- Molecule 20: Transcription factor TFIIIB component B”





• Molecule 21: RNA



• Molecule 22: Non-Template DNA



• Molecule 23: Template-DNA



## 4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	29951	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 ( $e^-/\text{\AA}^2$ )	61.3	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

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	0.23	0/11168	0.47	1/15086 (0.0%)
10	J	0.23	0/558	0.43	0/750
11	K	0.23	0/803	0.43	0/1083
12	L	0.22	0/360	0.47	0/478
13	M	0.23	0/1518	0.46	0/2054
14	N	0.23	0/805	0.48	0/1081
15	O	0.23	0/4358	0.44	0/5879
16	P	0.24	0/2020	0.47	1/2718 (0.0%)
17	Q	0.26	0/281	0.46	0/381
18	U	0.25	0/1443	0.47	0/1942
19	V	0.23	0/2728	0.43	0/3676
2	B	0.24	0/8943	0.48	0/12068
20	W	0.23	0/1831	0.43	0/2454
21	R	0.10	0/139	0.66	0/214
22	X	0.55	0/1363	1.05	0/2098
23	Y	0.54	0/1530	0.95	0/2357
3	C	0.24	0/2711	0.46	0/3676
4	D	0.23	0/991	0.43	0/1328
5	E	0.23	0/1787	0.44	0/2406
6	F	0.23	0/683	0.45	0/923
7	G	0.24	0/1486	0.44	0/2017
8	H	0.24	0/1138	0.51	0/1540
9	I	0.25	0/261	0.56	0/354
All	All	0.26	0/48905	0.52	2/66563 (0.0%)

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	1
18	U	0	1
3	C	0	2
All	All	0	4

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	1039	LEU	CA-CB-CG	7.11	131.66	115.30
16	P	265	LEU	CA-CB-CG	5.07	126.96	115.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	70	CYS	Peptide
3	C	146	ALA	Peptide
3	C	273	ASP	Peptide
18	U	70	ILE	Peptide

## 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	10972	0	11096	173	0
2	B	8788	0	8904	160	0
3	C	2655	0	2628	32	0
4	D	977	0	983	9	0
5	E	1751	0	1776	26	0
6	F	671	0	692	8	0
7	G	1448	0	1446	18	0
8	H	1120	0	1089	19	0
9	I	255	0	242	5	0
10	J	549	0	559	11	0
11	K	792	0	790	16	0
12	L	358	0	383	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	M	1484	0	1448	39	0
14	N	797	0	846	20	0
15	O	4293	0	4456	54	0
16	P	1990	0	2030	39	0
17	Q	273	0	285	6	0
18	U	1416	0	1493	42	0
19	V	2686	0	2703	49	0
20	W	1798	0	1789	35	0
21	R	126	0	65	1	0
22	X	1221	0	686	8	0
23	Y	1361	0	744	16	0
24	A	2	0	0	0	0
24	B	1	0	0	0	0
24	I	1	0	0	0	0
24	J	1	0	0	0	0
24	L	1	0	0	0	0
24	V	1	0	0	0	0
All	All	47788	0	47133	681	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 7.

All (681) 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:C:88:ASN:H	12:L:60:ARG:HH22	1.19	0.88
15:O:201:ILE:HD13	15:O:283:ASN:HD22	1.48	0.79
16:P:256:VAL:HG22	16:P:257:THR:HG23	1.67	0.76
1:A:666:LYS:HD2	1:A:670:GLY:HA3	1.66	0.76
13:M:148:LEU:HA	13:M:182:PHE:H	1.52	0.75
11:K:66:VAL:HG12	11:K:67:GLU:HG3	1.69	0.75
2:B:904:ARG:NH1	2:B:1030:MET:SD	2.61	0.74
13:M:140:TRP:HD1	13:M:142:GLY:H	1.37	0.73
13:M:164:LYS:HB3	13:M:167:GLN:HB2	1.69	0.73
13:M:111:ARG:HB3	13:M:120:GLU:HB3	1.70	0.73
5:E:3:GLN:HE21	5:E:7:ARG:HH22	1.35	0.73
19:V:146:PHE:O	19:V:149:ARG:NH1	2.21	0.73
20:W:407:ASN:ND2	23:Y:65:DG:OP1	2.22	0.73
15:O:549:GLN:HE21	15:O:567:ARG:HH22	1.38	0.72
5:E:151:PRO:HG2	5:E:198:ILE:HD11	1.72	0.72
14:N:366:HIS:HB2	14:N:370:LYS:HB2	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1162:GLU:HA	1:A:1278:ILE:HD11	1.72	0.71
1:A:520:HIS:HB3	2:B:1082:ARG:HH12	1.56	0.70
1:A:577:THR:HG21	11:K:89:CYS:H	1.56	0.70
2:B:695:GLN:HG2	2:B:697:PRO:HD2	1.73	0.70
6:F:75:PRO:HD2	6:F:78:GLN:HE21	1.57	0.68
15:O:40:ARG:NH2	16:P:314:GLU:OE1	2.25	0.68
1:A:217:ARG:HH22	15:O:555:ALA:HA	1.59	0.68
1:A:1369:LEU:HD11	1:A:1379:MET:HG3	1.77	0.67
13:M:92:ASN:HD21	13:M:181:PRO:HG3	1.60	0.67
20:W:450:SER:H	20:W:453:GLN:HE22	1.42	0.66
1:A:373:VAL:HG23	2:B:1087:SER:HB3	1.77	0.66
2:B:667:VAL:HG12	2:B:669:SER:H	1.59	0.66
1:A:1023:ARG:NH2	1:A:1028:MET:SD	2.69	0.66
15:O:233:SER:HB2	15:O:236:LYS:HD3	1.78	0.66
19:V:237:LEU:HB3	20:W:290:SER:HB3	1.78	0.66
1:A:630:ASN:ND2	1:A:650:GLY:O	2.28	0.65
9:I:7:SER:O	13:M:146:GLN:NE2	2.28	0.65
1:A:124:LEU:HA	1:A:127:LEU:HD23	1.80	0.64
19:V:215:PHE:HB3	19:V:223:ILE:HG21	1.78	0.64
15:O:163:VAL:HG13	15:O:164:ILE:HG23	1.79	0.64
2:B:1061:ARG:NH1	23:Y:26:DG:OP2	2.30	0.64
18:U:84:THR:O	18:U:88:HIS:NE2	2.29	0.64
15:O:512:ARG:NH1	15:O:570:GLU:OE1	2.30	0.64
1:A:1177:TYR:OH	1:A:1260:MET:SD	2.56	0.64
15:O:580:ASN:O	15:O:584:ASN:ND2	2.31	0.64
7:G:130:TRP:HE3	7:G:137:LYS:HB3	1.61	0.63
3:C:8:GLU:OE2	3:C:11:ARG:NE	2.30	0.63
16:P:135:LYS:NZ	16:P:150:LEU:O	2.28	0.63
18:U:68:GLN:NE2	18:U:163:SER:OG	2.31	0.63
2:B:221:THR:HG21	2:B:334:HIS:H	1.62	0.63
18:U:175:LEU:HD11	18:U:193:LEU:HD21	1.81	0.63
16:P:170:LEU:HD13	16:P:172:ILE:H	1.64	0.62
1:A:366:GLY:HA2	2:B:1061:ARG:HH21	1.65	0.62
2:B:83:ILE:HD11	2:B:93:LEU:HB2	1.81	0.62
1:A:264:PRO:O	1:A:269:ARG:NH1	2.32	0.62
1:A:473:LEU:HB2	1:A:520:HIS:HB2	1.81	0.62
2:B:610:ARG:NH2	2:B:654:GLU:OE2	2.32	0.62
15:O:289:LYS:HD2	15:O:323:SER:HB2	1.82	0.62
18:U:104:MET:HB2	18:U:113:ALA:HB3	1.82	0.61
5:E:168:TYR:HB2	5:E:170:LEU:HD22	1.82	0.61
1:A:389:ILE:HG22	1:A:537:VAL:HG11	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1145:ASP:OD2	2:B:1148:GLN:NE2	2.34	0.61
2:B:888:VAL:HG11	12:L:54:ARG:HH21	1.64	0.61
2:B:766:ASP:HA	2:B:770:ALA:HB3	1.82	0.61
18:U:107:ARG:NH2	20:W:469:ILE:O	2.29	0.61
3:C:81:GLU:OE2	12:L:70:ARG:NH2	2.30	0.61
16:P:142:PHE:HB3	16:P:145:ARG:HD2	1.81	0.61
1:A:1045:ASP:O	1:A:1053:LYS:NZ	2.34	0.60
2:B:66:ASN:HD21	2:B:159:ASN:HD21	1.49	0.60
1:A:1261:GLN:OE1	1:A:1265:ARG:NH1	2.34	0.60
3:C:248:GLN:HA	3:C:256:ILE:HD11	1.83	0.60
13:M:218:SER:HB2	20:W:380:VAL:HG12	1.81	0.60
2:B:55:LYS:HG2	2:B:59:LYS:HD3	1.83	0.60
2:B:1028:LYS:HG2	2:B:1029:HIS:H	1.66	0.60
11:K:62:SER:OG	11:K:104:ARG:NH1	2.34	0.60
19:V:258:ARG:NH2	23:Y:63:DT:OP2	2.35	0.60
1:A:117:GLU:HG2	15:O:212:GLU:HG2	1.82	0.59
2:B:944:MET:SD	2:B:1024:TYR:OH	2.60	0.59
1:A:1187:ARG:NH1	1:A:1189:ASP:OD1	2.35	0.59
11:K:64:GLN:HE22	11:K:100:LEU:HD13	1.67	0.59
1:A:1166:LEU:HD21	1:A:1268:PRO:HA	1.85	0.59
1:A:348:LYS:HG3	1:A:349:PRO:HD3	1.84	0.59
15:O:288:MET:HG3	15:O:326:ILE:HD13	1.85	0.58
1:A:559:THR:HG21	2:B:947:HIS:HE1	1.69	0.58
16:P:313:ASP:O	16:P:314:GLU:HG3	2.04	0.58
1:A:70:CYS:SG	1:A:71:HIS:N	2.76	0.58
14:N:394:VAL:HG23	14:N:412:VAL:HB	1.84	0.58
18:U:129:GLU:HG2	18:U:133:LYS:HE2	1.86	0.58
1:A:668:VAL:HG22	1:A:677:VAL:HG23	1.84	0.58
22:X:9:DA:H61	23:Y:73:DT:H3	1.50	0.58
1:A:895:ASP:OD1	1:A:1380:ARG:NH2	2.36	0.58
2:B:47:LEU:HD11	2:B:675:ILE:HD11	1.86	0.58
1:A:546:SER:OG	1:A:1349:SER:O	2.20	0.58
6:F:85:MET:HB2	6:F:151:LEU:HB3	1.86	0.57
17:Q:49:ASN:HA	17:Q:52:ARG:HG2	1.86	0.57
18:U:193:LEU:HD23	18:U:206:ILE:HD11	1.85	0.57
20:W:465:ARG:HD3	20:W:468:LEU:HD22	1.84	0.57
2:B:829:LEU:HD11	2:B:896:ILE:HD12	1.87	0.57
1:A:568:ASP:O	8:H:22:LYS:NZ	2.36	0.57
18:U:107:ARG:HA	18:U:110:LYS:HA	1.84	0.57
2:B:686:GLY:HA3	2:B:740:THR:HG21	1.87	0.57
1:A:599:LYS:HB2	8:H:96:VAL:HG22	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:785:CYS:SG	2:B:1026:LYS:NZ	2.74	0.57
2:B:832:VAL:HA	2:B:881:ILE:HD11	1.86	0.57
2:B:1097:LYS:NZ	2:B:1115:ASN:OD1	2.35	0.57
20:W:450:SER:H	20:W:453:GLN:NE2	2.03	0.57
3:C:73:SER:O	3:C:214:GLY:N	2.38	0.57
16:P:173:GLU:CD	16:P:174:PHE:H	2.08	0.57
18:U:93:GLU:OE2	20:W:439:PHE:N	2.32	0.57
18:U:195:TYR:HD2	18:U:204:LEU:HD21	1.70	0.57
1:A:1374:PHE:O	1:A:1378:LYS:NZ	2.37	0.56
8:H:63:LEU:HG	8:H:89:LEU:HD23	1.86	0.56
1:A:475:ASN:HA	1:A:508:TYR:HE1	1.70	0.56
1:A:714:ILE:HA	1:A:717:VAL:HG12	1.87	0.56
5:E:112:TYR:OH	5:E:115:ASN:O	2.23	0.56
15:O:199:TYR:HA	15:O:283:ASN:HB3	1.87	0.56
2:B:1038:ARG:NH2	2:B:1042:PRO:O	2.30	0.56
2:B:58:VAL:HG12	2:B:61:HIS:HB2	1.85	0.56
9:I:3:SER:HB2	9:I:12:LEU:HB2	1.87	0.56
1:A:252:ARG:HB2	15:O:42:LEU:HD12	1.87	0.56
1:A:1350:VAL:HG23	1:A:1351:ASP:H	1.71	0.56
7:G:39:ILE:HD11	7:G:45:CYS:HB3	1.87	0.56
16:P:154:GLN:NE2	20:W:527:ASP:OD1	2.39	0.56
1:A:632:VAL:HG23	1:A:633:PHE:H	1.71	0.56
2:B:124:THR:HG22	2:B:188:ASN:H	1.71	0.56
13:M:217:LYS:NZ	20:W:360:THR:O	2.37	0.55
2:B:961:LEU:HD23	2:B:1018:PHE:CE2	2.41	0.55
19:V:275:PHE:CE2	19:V:279:ASP:HB2	2.41	0.55
15:O:468:LEU:HD22	15:O:478:VAL:HB	1.87	0.55
3:C:80:ALA:HB3	3:C:102:GLY:HA2	1.87	0.55
5:E:47:CYS:HB2	5:E:53:PRO:HA	1.89	0.55
2:B:207:VAL:HG23	2:B:366:ILE:HD12	1.88	0.55
14:N:291:LEU:HD12	14:N:294:LEU:HD11	1.88	0.55
1:A:166:LYS:HA	1:A:178:ILE:HG22	1.89	0.55
1:A:1092:ILE:HA	1:A:1095:GLN:HE21	1.72	0.55
2:B:332:ILE:O	2:B:345:LYS:NZ	2.36	0.55
3:C:165:ARG:NH2	3:C:190:ASP:OD1	2.40	0.55
3:C:229:LEU:HB2	3:C:293:ARG:HD3	1.89	0.55
16:P:172:ILE:HG22	16:P:173:GLU:H	1.72	0.55
2:B:821:HIS:HA	2:B:824:LEU:HD13	1.89	0.55
3:C:282:TYR:O	3:C:283:GLU:HG3	2.07	0.54
13:M:250:GLY:O	13:M:251:THR:HG22	2.08	0.54
1:A:1050:ASP:HA	1:A:1053:LYS:HE2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1100:LEU:HD21	2:B:1109:THR:HG21	1.88	0.54
14:N:305:MET:HA	14:N:412:VAL:HG13	1.88	0.54
1:A:999:ASP:OD1	1:A:1002:ARG:NH1	2.41	0.54
8:H:11:GLN:O	8:H:29:ALA:N	2.40	0.54
1:A:1373:ARG:HD3	1:A:1391:LYS:HD3	1.90	0.54
2:B:773:LEU:HB3	2:B:942:ILE:HG12	1.89	0.54
1:A:1355:ILE:O	1:A:1359:GLY:N	2.40	0.54
20:W:472:ALA:O	20:W:475:SER:OG	2.26	0.54
22:X:60:DG:H1	23:Y:22:DC:H2'	1.72	0.54
2:B:1036:HIS:CE1	2:B:1058:GLY:HA3	2.43	0.54
11:K:88:PHE:HB3	11:K:106:GLN:HB3	1.89	0.54
2:B:1027:LEU:HD23	2:B:1028:LYS:H	1.72	0.53
13:M:110:ALA:HB3	13:M:245:LEU:HB2	1.90	0.53
15:O:199:TYR:HB3	15:O:286:ARG:HB3	1.91	0.53
19:V:188:LYS:HD2	19:V:247:VAL:HG13	1.89	0.53
1:A:1038:GLU:OE1	1:A:1039:LEU:HD13	2.08	0.53
1:A:834:HIS:NE2	2:B:674:GLU:OE2	2.41	0.53
2:B:129:ILE:HD11	2:B:152:MET:HB2	1.90	0.53
2:B:252:PRO:HG2	2:B:255:ILE:HD13	1.89	0.53
8:H:106:GLU:HA	8:H:112:ILE:HA	1.90	0.53
2:B:591:TYR:HB2	2:B:652:ASN:HB3	1.89	0.53
13:M:137:GLU:O	13:M:141:ASN:ND2	2.42	0.53
13:M:95:ARG:HH11	14:N:390:PHE:HE2	1.57	0.53
16:P:253:LEU:HD12	16:P:261:TYR:HB3	1.90	0.53
1:A:1148:ASP:O	1:A:1291:ARG:NH1	2.42	0.53
1:A:789:ASN:ND2	1:A:791:PRO:HD2	2.24	0.53
15:O:640:ARG:NH2	16:P:308:GLU:O	2.42	0.53
2:B:678:PHE:HB2	2:B:978:CYS:HB2	1.90	0.53
12:L:31:CYS:SG	12:L:32:ALA:N	2.82	0.53
1:A:1121:LEU:O	1:A:1342:THR:OG1	2.27	0.53
2:B:666:ILE:HD11	2:B:673:LEU:HD12	1.89	0.53
3:C:11:ARG:HH22	8:H:21:ASN:HB3	1.74	0.53
1:A:1141:ILE:HD11	1:A:1295:VAL:HB	1.92	0.52
1:A:948:ASN:OD1	1:A:950:GLN:NE2	2.42	0.52
1:A:974:LEU:HD21	1:A:1002:ARG:HH21	1.74	0.52
1:A:554:THR:OG1	1:A:555:GLN:OE1	2.25	0.52
4:D:141:CYS:SG	4:D:144:ARG:NH2	2.83	0.52
13:M:74:PHE:N	14:N:363:ILE:O	2.33	0.52
18:U:109:PRO:O	18:U:110:LYS:HG3	2.10	0.52
19:V:265:THR:HG23	19:V:266:LYS:H	1.74	0.52
4:D:15:GLU:HA	4:D:18:LYS:HB3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:884:TYR:OH	1:A:888:ARG:NH2	2.42	0.52
2:B:151:ARG:H	2:B:430:THR:HG22	1.74	0.52
2:B:210:ASP:H	2:B:215:ILE:HG22	1.74	0.52
1:A:859:PRO:HD3	2:B:661:LEU:HD21	1.90	0.52
8:H:38:LEU:HB3	8:H:125:LEU:HB3	1.92	0.52
18:U:159:ASN:HD21	18:U:216:GLY:H	1.58	0.52
2:B:730:TYR:OH	3:C:100:ARG:NH1	2.43	0.52
5:E:156:LEU:HD21	5:E:195:VAL:HG23	1.91	0.52
2:B:754:ASN:HD21	10:J:52:THR:HG21	1.75	0.52
1:A:1186:VAL:HG23	1:A:1230:ILE:HG23	1.91	0.52
19:V:251:ALA:HB2	20:W:407:ASN:HD22	1.74	0.52
15:O:192:VAL:HG22	15:O:274:VAL:HG13	1.92	0.51
1:A:1278:ILE:HG22	1:A:1297:GLY:HA3	1.91	0.51
1:A:1301:ARG:HE	1:A:1325:VAL:HG13	1.75	0.51
1:A:19:SER:O	2:B:1138:ALA:N	2.37	0.51
1:A:386:ASN:O	1:A:699:LYS:NZ	2.43	0.51
1:A:464:ARG:NH1	1:A:467:GLU:OE2	2.43	0.51
1:A:550:ILE:HD12	1:A:551:ILE:HD12	1.92	0.51
1:A:716:ASP:O	1:A:789:ASN:ND2	2.38	0.51
1:A:502:GLU:HG2	2:B:767:ILE:HG13	1.92	0.51
13:M:105:PRO:HB2	13:M:123:ILE:HD11	1.93	0.51
14:N:374:LYS:NZ	14:N:377:ASN:O	2.29	0.51
18:U:164:CYS:SG	18:U:165:ASP:N	2.83	0.51
5:E:107:THR:HB	5:E:130:ALA:HB3	1.92	0.51
15:O:354:GLU:HG2	15:O:355:LYS:HG2	1.93	0.51
5:E:78:LEU:HD11	5:E:109:ILE:HD12	1.93	0.51
1:A:607:LYS:NZ	8:H:119:GLY:O	2.43	0.51
3:C:105:PRO:HB2	3:C:187:ALA:HB3	1.92	0.51
12:L:48:CYS:HB3	12:L:56:LEU:HD21	1.91	0.51
17:Q:47:ILE:O	17:Q:52:ARG:NH2	2.44	0.51
1:A:475:ASN:HB3	1:A:485:ILE:HG13	1.93	0.51
1:A:12:ARG:N	2:B:1144:GLU:O	2.42	0.51
10:J:31:ASP:HB3	10:J:34:THR:HG22	1.93	0.51
2:B:113:THR:OG1	2:B:117:GLU:OE1	2.29	0.51
2:B:806:ILE:HG22	2:B:828:GLY:HA3	1.92	0.51
1:A:1315:THR:OG1	1:A:1337:ARG:NH2	2.44	0.50
18:U:133:LYS:HE3	19:V:452:LEU:HA	1.93	0.50
18:U:88:HIS:HB2	19:V:475:LEU:HD22	1.91	0.50
2:B:969:LEU:HD11	2:B:994:GLN:HG2	1.94	0.50
17:Q:38:SER:C	17:Q:39:ILE:HG13	2.31	0.50
19:V:239:ARG:NH2	20:W:285:PRO:HD2	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:W:483:GLU:O	20:W:487:GLU:N	2.43	0.50
1:A:1335:ILE:HD11	1:A:1359:GLY:HA2	1.94	0.50
2:B:198:GLU:HB3	2:B:377:LEU:HD23	1.93	0.50
7:G:207:LEU:HG	7:G:209:SER:H	1.76	0.50
2:B:81:GLN:NE2	2:B:95:TYR:O	2.45	0.50
2:B:483:VAL:HG12	2:B:485:GLY:H	1.76	0.50
2:B:846:SER:HB3	2:B:866:TYR:HB3	1.92	0.50
3:C:16:THR:O	3:C:295:ARG:NH1	2.45	0.50
19:V:200:LYS:HG3	19:V:204:LYS:NZ	2.26	0.50
15:O:130:LEU:H	15:O:130:LEU:HD23	1.75	0.50
18:U:188:GLU:O	19:V:219:ARG:NH2	2.40	0.50
2:B:961:LEU:HB2	2:B:1022:ILE:HD11	1.94	0.50
2:B:409:LYS:HB2	2:B:412:ARG:HB3	1.94	0.50
13:M:195:GLN:HG3	13:M:199:ARG:HH11	1.77	0.50
8:H:7:ASP:OD1	8:H:58:THR:OG1	2.23	0.50
14:N:299:ASN:C	14:N:300:LYS:HD2	2.33	0.50
1:A:379:THR:HG22	2:B:1035:MET:HA	1.94	0.49
19:V:201:ASP:HA	19:V:204:LYS:HE2	1.93	0.49
1:A:117:GLU:HA	1:A:120:LYS:HE3	1.94	0.49
1:A:1275:LEU:HD11	1:A:1302:ASP:HB3	1.93	0.49
3:C:28:GLU:CD	3:C:29:ASN:H	2.15	0.49
2:B:271:LEU:O	2:B:550:HIS:NE2	2.44	0.49
13:M:95:ARG:NH1	14:N:412:VAL:O	2.42	0.49
1:A:153:ARG:HH12	15:O:339:LEU:HB3	1.77	0.49
15:O:464:ASN:ND2	16:P:254:GLU:OE2	2.45	0.49
2:B:705:MET:HB3	2:B:1027:LEU:HD11	1.93	0.49
1:A:703:ARG:NH2	11:K:93:ILE:O	2.41	0.49
2:B:738:THR:HG23	2:B:977:THR:HA	1.95	0.49
3:C:191:ILE:HG23	10:J:15:GLY:HA3	1.95	0.49
16:P:102:ARG:HH21	16:P:154:GLN:HE21	1.61	0.49
16:P:311:TYR:CD2	17:Q:37:PRO:HG2	2.47	0.49
3:C:245:ARG:NH1	3:C:263:ASP:OD2	2.44	0.49
15:O:222:HIS:HA	15:O:225:ASN:HD22	1.78	0.49
15:O:37:LEU:HA	15:O:40:ARG:HE	1.77	0.49
18:U:169:PRO:O	18:U:239:LYS:NZ	2.30	0.49
2:B:615:VAL:HB	2:B:671:THR:HA	1.94	0.49
4:D:140:GLU:HG3	4:D:144:ARG:HH12	1.77	0.49
8:H:56:THR:HB	8:H:145:ARG:HG2	1.93	0.49
16:P:236:THR:HG23	16:P:239:ASN:H	1.76	0.49
18:U:196:ARG:HG3	18:U:203:VAL:HG12	1.94	0.49
19:V:485:ARG:HG3	20:W:498:PHE:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1432:LYS:HB2	7:G:59:LEU:HD11	1.95	0.49
1:A:90:VAL:HG21	1:A:323:VAL:HG11	1.95	0.49
1:A:583:MET:HB2	1:A:696:ARG:HG2	1.94	0.49
2:B:527:GLU:HG2	2:B:528:PRO:HD3	1.94	0.49
1:A:787:ASN:ND2	8:H:21:ASN:OD1	2.45	0.48
2:B:186:ILE:HA	2:B:191:GLU:HA	1.94	0.48
4:D:99:SER:OG	4:D:101:GLU:OE1	2.30	0.48
16:P:225:ILE:HD12	16:P:235:LEU:HG	1.94	0.48
16:P:297:PHE:O	16:P:298:LYS:HD3	2.13	0.48
1:A:1384:LEU:HD23	1:A:1416:ILE:HD11	1.94	0.48
15:O:43:ASN:HB3	15:O:47:PHE:HB2	1.95	0.48
15:O:83:ILE:HD13	15:O:86:MET:HB2	1.95	0.48
18:U:128:SER:HB2	18:U:131:ASP:HB2	1.95	0.48
19:V:451:TYR:HA	19:V:454:LYS:HG2	1.95	0.48
1:A:468:ASP:OD1	1:A:468:ASP:N	2.47	0.48
2:B:240:ILE:HG12	2:B:286:ASN:HD21	1.77	0.48
5:E:170:LEU:HB2	5:E:174:GLN:HE21	1.77	0.48
13:M:123:ILE:HG22	13:M:146:GLN:HB3	1.93	0.48
2:B:1004:LEU:HD23	2:B:1017:ILE:HB	1.95	0.48
2:B:379:LEU:H	2:B:379:LEU:HD23	1.77	0.48
1:A:91:PHE:CE2	1:A:224:PRO:HG3	2.49	0.48
2:B:965:LYS:HA	10:J:44:TYR:OH	2.13	0.48
15:O:80:VAL:HG21	15:O:88:VAL:HG22	1.96	0.48
1:A:1373:ARG:NH2	23:Y:22:DC:O5'	2.45	0.48
7:G:4:LEU:HD13	7:G:73:ARG:HG3	1.95	0.48
1:A:1286:ARG:HH12	1:A:1291:ARG:HA	1.77	0.48
1:A:530:GLU:HG2	6:F:91:ALA:HB1	1.95	0.48
1:A:997:GLN:HG3	1:A:998:TYR:HD1	1.78	0.48
2:B:1088:ASP:OD2	2:B:1123:TYR:N	2.47	0.48
2:B:412:ARG:HE	2:B:416:TYR:HB2	1.79	0.48
13:M:174:GLU:OE1	13:M:175:ARG:NH1	2.46	0.48
15:O:254:ASN:O	15:O:258:GLU:N	2.46	0.48
2:B:296:TYR:OH	9:I:27:ARG:NH2	2.46	0.48
2:B:390:ASP:HA	2:B:393:LYS:HE3	1.94	0.48
7:G:46:ILE:HB	7:G:75:VAL:HG23	1.94	0.48
16:P:105:ILE:HD11	16:P:113:ARG:HH22	1.79	0.48
19:V:275:PHE:HE2	19:V:279:ASP:HB2	1.79	0.48
1:A:406:GLU:OE2	1:A:412:ASN:ND2	2.47	0.48
19:V:257:GLN:OE1	19:V:258:ARG:NH1	2.47	0.48
2:B:314:ARG:O	2:B:315:GLN:HG2	2.14	0.48
2:B:730:TYR:O	2:B:753:GLN:NE2	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:149:ARG:HH21	7:G:201:GLN:HG2	1.79	0.48
13:M:163:VAL:HG23	13:M:168:VAL:HB	1.96	0.48
1:A:1320:LEU:HD23	1:A:1320:LEU:H	1.78	0.47
2:B:240:ILE:HG22	2:B:253:ILE:HG13	1.95	0.47
2:B:338:GLU:HG3	2:B:342:PHE:CD1	2.49	0.47
1:A:794:MET:SD	2:B:950:PRO:HG2	2.54	0.47
5:E:81:GLU:HB2	5:E:96:PHE:HE1	1.78	0.47
15:O:266:ASP:HB3	15:O:270:SER:HB3	1.96	0.47
1:A:903:THR:HG23	1:A:914:PHE:HA	1.95	0.47
3:C:313:ILE:O	3:C:317:SER:N	2.46	0.47
4:D:126:GLN:HE22	7:G:86:THR:HG23	1.79	0.47
1:A:111:SER:HB2	1:A:234:ILE:HG22	1.95	0.47
14:N:366:HIS:NE2	14:N:372:SER:OG	2.34	0.47
16:P:53:GLN:NE2	20:W:363:GLN:OE1	2.42	0.47
1:A:11:LYS:HD2	2:B:1117:ILE:HG21	1.96	0.47
2:B:236:LYS:HB2	2:B:241:TYR:CE2	2.49	0.47
19:V:122:SER:HB3	22:X:34:DT:H5"	1.97	0.47
20:W:532:ARG:HH12	20:W:536:ALA:HB2	1.79	0.47
20:W:307:ARG:NH2	23:Y:66:DC:OP2	2.47	0.47
1:A:433:LEU:HD23	1:A:444:LEU:HD21	1.95	0.47
2:B:934:ASN:HD21	2:B:938:ILE:HD13	1.78	0.47
4:D:15:GLU:HG3	4:D:18:LYS:HE3	1.97	0.47
16:P:247:LEU:HD12	16:P:252:LYS:HG3	1.97	0.47
1:A:393:ALA:HB3	1:A:499:ARG:HB2	1.96	0.47
2:B:59:LYS:NZ	2:B:518:THR:O	2.42	0.47
15:O:595:LEU:HD11	15:O:633:ARG:HH21	1.79	0.47
18:U:85:VAL:HG12	18:U:146:ILE:HD13	1.97	0.47
19:V:458:ASP:OD1	19:V:458:ASP:N	2.46	0.47
1:A:1066:SER:O	1:A:1070:PHE:N	2.44	0.47
2:B:417:ASP:H	19:V:151:GLN:HE22	1.62	0.47
2:B:795:LEU:HD12	2:B:894:ALA:HB3	1.96	0.47
18:U:78:CYS:HB2	18:U:149:ALA:HB3	1.96	0.47
1:A:815:GLN:NE2	1:A:816:GLN:O	2.48	0.47
1:A:895:ASP:OD1	1:A:895:ASP:N	2.45	0.47
1:A:409:THR:HG22	1:A:410:ARG:HD2	1.96	0.47
2:B:116:HIS:HA	2:B:119:ARG:HG2	1.97	0.47
2:B:667:VAL:HB	2:B:670:MET:HB2	1.96	0.47
11:K:85:ASP:O	11:K:107:THR:OG1	2.30	0.47
18:U:192:GLY:HA3	18:U:205:LEU:HD21	1.97	0.47
1:A:470:ASP:HA	1:A:528:ARG:HH11	1.80	0.47
2:B:552:ASN:OD1	2:B:566:ARG:HD2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:578:LEU:HD22	2:B:583:LYS:HD3	1.97	0.47
2:B:622:VAL:HG23	2:B:623:LYS:H	1.79	0.47
7:G:27:THR:HG22	7:G:31:ASN:HD21	1.80	0.47
1:A:683:ARG:HH12	1:A:925:GLU:HG2	1.80	0.47
1:A:374:ASP:OD1	2:B:1038:ARG:NE	2.48	0.47
16:P:111:LYS:HG2	16:P:116:LEU:O	2.15	0.47
23:Y:62:DG:H2"	23:Y:63:DT:H5"	1.96	0.47
10:J:10:CYS:SG	10:J:11:GLY:N	2.87	0.46
1:A:41:ASP:HB3	1:A:48:PRO:HB2	1.97	0.46
20:W:487:GLU:O	20:W:491:ASN:ND2	2.48	0.46
3:C:256:ILE:HG22	3:C:267:VAL:HA	1.96	0.46
6:F:115:THR:HG22	6:F:116:ASP:H	1.80	0.46
3:C:242:GLU:OE1	3:C:246:ARG:NH2	2.48	0.46
1:A:1064:GLU:OE2	1:A:1068:ARG:NE	2.42	0.46
1:A:1254:ASN:HD21	9:I:13:LEU:HD21	1.80	0.46
1:A:757:ASN:HD22	1:A:760:GLN:NE2	2.13	0.46
2:B:835:LYS:HB3	2:B:880:HIS:CE1	2.51	0.46
2:B:99:ARG:NH2	2:B:146:ASP:OD1	2.49	0.46
19:V:498:GLN:OE1	19:V:502:ARG:NH1	2.48	0.46
2:B:705:MET:HE1	2:B:919:LYS:HD3	1.96	0.46
2:B:81:GLN:HG3	2:B:95:TYR:H	1.81	0.46
8:H:58:THR:HB	8:H:143:LEU:HB2	1.98	0.46
15:O:212:GLU:OE2	15:O:216:GLN:NE2	2.42	0.46
15:O:346:GLN:HA	15:O:349:ALA:HB3	1.97	0.46
20:W:365:LYS:NZ	20:W:367:ASN:OD1	2.48	0.46
1:A:1316:THR:OG1	1:A:1317:ASN:N	2.49	0.46
2:B:961:LEU:HD11	2:B:1020:GLY:HA3	1.96	0.46
3:C:276:SER:O	3:C:277:ARG:HD2	2.15	0.46
3:C:229:LEU:HD23	3:C:295:ARG:HA	1.97	0.46
16:P:266:GLU:O	16:P:270:GLN:N	2.45	0.46
2:B:1010:GLY:O	3:C:227:TYR:OH	2.33	0.46
2:B:284:ALA:HA	2:B:287:LEU:HD13	1.98	0.46
1:A:555:GLN:NE2	2:B:768:GLU:HB3	2.31	0.46
19:V:114:ASN:O	19:V:167:LYS:NZ	2.40	0.46
2:B:81:GLN:HG2	2:B:82:LEU:HD12	1.98	0.46
10:J:26:GLN:HG2	10:J:27:GLU:N	2.31	0.46
15:O:259:LEU:H	15:O:259:LEU:HD23	1.81	0.46
18:U:185:TYR:CD1	18:U:193:LEU:HD13	2.51	0.46
19:V:149:ARG:NH1	19:V:150:LEU:HD13	2.31	0.46
1:A:1092:ILE:O	1:A:1095:GLN:NE2	2.48	0.45
2:B:139:ARG:HD3	2:B:139:ARG:H	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:125:TYR:HB2	2:B:186:ILE:HD11	1.98	0.45
19:V:503:LEU:HD22	19:V:506:GLU:HB3	1.98	0.45
2:B:40:THR:HB	2:B:624:ASP:HB3	1.98	0.45
18:U:162:GLY:O	18:U:214:LEU:N	2.40	0.45
19:V:135:ARG:NH2	19:V:174:PRO:O	2.49	0.45
1:A:398:VAL:HG23	2:B:1037:ALA:HB2	1.97	0.45
2:B:689:PRO:HG3	2:B:915:ARG:NH2	2.31	0.45
13:M:100:LYS:HD3	13:M:101:PRO:HD2	1.98	0.45
15:O:316:LEU:O	15:O:320:GLU:HG3	2.16	0.45
15:O:527:LEU:HD23	16:P:175:ILE:HG21	1.98	0.45
20:W:371:THR:HG22	20:W:372:MET:HG3	1.98	0.45
1:A:95:TYR:O	1:A:99:THR:N	2.50	0.45
5:E:124:VAL:HB	5:E:125:PRO:HD3	1.97	0.45
16:P:263:VAL:HG12	16:P:265:LEU:H	1.81	0.45
18:U:204:LEU:HB3	18:U:214:LEU:HD13	1.99	0.45
16:P:132:ARG:HD2	19:V:503:LEU:HD21	1.98	0.45
1:A:12:ARG:HG2	1:A:13:ILE:H	1.81	0.45
2:B:204:ARG:CZ	2:B:376:ARG:HH21	2.30	0.45
2:B:832:VAL:HG13	12:L:60:ARG:HA	1.99	0.45
15:O:191:PHE:O	15:O:195:CYS:N	2.47	0.45
13:M:216:VAL:HG11	16:P:95:SER:OG	2.17	0.45
2:B:144:HIS:HD2	2:B:147:VAL:HG22	1.81	0.45
11:K:47:ILE:HD12	11:K:47:ILE:H	1.82	0.45
18:U:219:GLN:HG3	18:U:221:GLU:H	1.82	0.45
23:Y:69:DA:H2"	23:Y:70:DA:C8	2.52	0.45
1:A:434:LEU:HD21	1:A:441:ARG:HG2	1.98	0.45
1:A:440:ALA:O	1:A:441:ARG:HG3	2.17	0.45
3:C:255:VAL:HG13	3:C:256:ILE:HG23	1.98	0.45
7:G:115:LEU:HD11	7:G:200:CYS:HB2	1.97	0.45
19:V:274:LYS:HG2	19:V:275:PHE:H	1.82	0.45
22:X:18:DT:O2	23:Y:65:DG:N2	2.50	0.45
2:B:696:SER:OG	2:B:697:PRO:HD3	2.16	0.45
8:H:12:VAL:HA	8:H:28:ALA:HA	1.99	0.45
15:O:484:MET:HG3	15:O:485:PRO:HD3	1.98	0.45
1:A:1333:TYR:HD2	5:E:149:LEU:HD13	1.82	0.45
1:A:1319:VAL:HG21	1:A:1335:ILE:HG22	1.99	0.45
2:B:1027:LEU:HD23	2:B:1028:LYS:N	2.31	0.45
2:B:331:VAL:HA	2:B:334:HIS:CE1	2.52	0.45
2:B:890:ASP:OD1	2:B:890:ASP:N	2.49	0.45
4:D:134:LEU:HD13	4:D:137:ILE:HD12	1.98	0.45
13:M:137:GLU:OE2	13:M:143:VAL:HG11	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:M:161:ALA:HB2	13:M:170:LEU:HG	1.98	0.45
1:A:352:GLY:H	1:A:355:GLN:HB2	1.82	0.45
2:B:916:HIS:HD1	2:B:957:LYS:HG3	1.82	0.45
15:O:222:HIS:HA	15:O:225:ASN:ND2	2.31	0.45
15:O:471:THR:OG1	15:O:475:VAL:O	2.29	0.45
19:V:268:ALA:O	19:V:272:VAL:HG23	2.17	0.45
1:A:1318:HIS:O	1:A:1321:GLU:HG3	2.17	0.44
1:A:579:LEU:HD23	1:A:609:VAL:HG21	1.98	0.44
1:A:757:ASN:HD22	1:A:760:GLN:HE21	1.65	0.44
14:N:386:ALA:HB2	14:N:416:ILE:HG22	1.99	0.44
18:U:197:MET:HB3	18:U:202:ILE:HG22	1.98	0.44
19:V:476:ASN:HB3	19:V:479:ALA:HB3	2.00	0.44
1:A:105:GLY:HA2	1:A:148:CYS:SG	2.57	0.44
1:A:949:ASN:HA	1:A:1061:ARG:NH2	2.32	0.44
3:C:26:ASP:OD1	3:C:26:ASP:N	2.50	0.44
10:J:26:GLN:HG2	10:J:27:GLU:H	1.82	0.44
11:K:55:SER:HB3	11:K:60:SER:HB2	1.99	0.44
18:U:81:ASP:HB3	18:U:84:THR:HG22	2.00	0.44
19:V:242:THR:OG1	20:W:296:SER:OG	2.26	0.44
18:U:138:LYS:NZ	19:V:465:VAL:O	2.42	0.44
19:V:474:LEU:HD22	19:V:475:LEU:H	1.82	0.44
2:B:405:LYS:HA	2:B:408:LYS:HE3	1.99	0.44
2:B:412:ARG:HH22	2:B:415:GLU:HG2	1.82	0.44
10:J:17:LYS:HB3	10:J:39:LEU:HD21	2.00	0.44
13:M:96:LEU:HA	13:M:101:PRO:HA	2.00	0.44
13:M:89:GLN:HG2	14:N:394:VAL:HG12	1.99	0.44
18:U:103:ILE:HD11	18:U:112:THR:HB	2.00	0.44
18:U:88:HIS:HB2	19:V:475:LEU:HD13	1.98	0.44
1:A:929:ASN:O	1:A:931:GLN:NE2	2.50	0.44
18:U:108:GLU:HA	18:U:109:PRO:HA	1.89	0.44
19:V:237:LEU:HG	20:W:287:GLY:H	1.82	0.44
1:A:4:VAL:HA	7:G:38:ILE:HG22	1.99	0.44
2:B:206:ILE:O	2:B:217:GLN:HB2	2.17	0.44
2:B:555:VAL:HA	2:B:599:VAL:HG23	2.00	0.44
3:C:216:HIS:CE1	3:C:218:LYS:HG3	2.53	0.44
13:M:79:SER:HB2	13:M:258:THR:HG23	1.99	0.44
13:M:245:LEU:HD12	14:N:406:ALA:HB2	2.00	0.44
1:A:763:GLU:HG2	1:A:822:ARG:HE	1.83	0.43
2:B:1002:ASP:OD1	2:B:1003:MET:N	2.50	0.43
3:C:103:LEU:HB3	10:J:6:ARG:HE	1.83	0.43
11:K:126:ASP:HA	11:K:129:ASP:HB2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:O:596:LYS:O	15:O:600:SER:OG	2.36	0.43
16:P:102:ARG:NH1	16:P:152:SER:OG	2.50	0.43
16:P:176:ASN:HA	16:P:179:LEU:HB3	1.99	0.43
1:A:1125:ARG:HH22	1:A:1129:ILE:HD11	1.83	0.43
1:A:1365:LYS:NZ	1:A:1378:LYS:O	2.35	0.43
1:A:15:GLY:HA3	1:A:1407:ALA:HA	2.00	0.43
1:A:597:ILE:HG22	8:H:97:MET:HG2	2.00	0.43
2:B:579:ARG:NH1	2:B:588:ILE:O	2.51	0.43
2:B:556:TYR:HD2	2:B:600:HIS:CE1	2.37	0.43
2:B:710:ILE:HD13	2:B:726:TYR:HB3	2.00	0.43
13:M:71:ILE:N	14:N:365:VAL:O	2.51	0.43
1:A:894:GLU:OE2	1:A:1385:GLN:NE2	2.52	0.43
1:A:822:ARG:HH12	1:A:845:LYS:HD2	1.81	0.43
2:B:1004:LEU:HB2	2:B:1017:ILE:HD12	1.99	0.43
11:K:95:HIS:CD2	11:K:97:SER:HB3	2.54	0.43
15:O:212:GLU:OE1	15:O:333:THR:OG1	2.32	0.43
15:O:356:THR:HG22	15:O:358:GLY:H	1.83	0.43
2:B:327:ILE:O	2:B:331:VAL:HG22	2.19	0.43
7:G:93:THR:HG22	7:G:94:ALA:H	1.83	0.43
12:L:26:THR:O	12:L:62:LYS:NZ	2.51	0.43
13:M:160:ALA:HB2	13:M:173:ILE:HD11	2.00	0.43
14:N:303:ARG:HD2	14:N:413:ASP:HB2	1.99	0.43
19:V:252:GLU:O	19:V:255:LEU:HG	2.18	0.43
1:A:880:ALA:HA	23:Y:24:DT:C2	2.54	0.43
2:B:90:GLU:HB2	20:W:391:ASN:HD21	1.84	0.43
15:O:308:THR:HG21	15:O:459:ILE:HD11	1.99	0.43
3:C:103:LEU:HD23	3:C:218:LYS:HG2	2.00	0.43
16:P:66:GLN:HB3	16:P:69:GLU:OE2	2.18	0.43
1:A:1122:GLY:HA2	1:A:1346:HIS:HE1	1.83	0.43
5:E:147:HIS:HE1	5:E:149:LEU:HD23	1.82	0.43
13:M:251:THR:HA	14:N:409:LEU:HD22	2.00	0.43
18:U:99:PHE:HE2	18:U:101:ALA:HB3	1.84	0.43
15:O:192:VAL:HG21	15:O:273:ILE:HA	2.01	0.43
15:O:472:LYS:HB3	15:O:475:VAL:HG12	2.00	0.43
23:Y:58:DT:H6	23:Y:58:DT:H2'	1.68	0.43
1:A:949:ASN:HA	1:A:1061:ARG:HH21	1.84	0.43
1:A:1227:LYS:H	1:A:1227:LYS:HD2	1.83	0.43
1:A:134:ASN:HB3	1:A:1381:ASP:OD2	2.19	0.43
2:B:860:VAL:HG23	2:B:861:ASN:H	1.83	0.43
19:V:218:ARG:HH22	19:V:288:PRO:HG2	1.83	0.43
20:W:289:ILE:H	20:W:289:ILE:HD12	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:ASP:HA	1:A:149:LYS:NZ	2.34	0.43
7:G:120:TYR:O	7:G:128:TRP:HB3	2.18	0.43
11:K:47:ILE:HG13	11:K:65:ILE:HG12	2.01	0.43
13:M:125:LEU:HD21	13:M:146:GLN:HB2	2.01	0.43
1:A:832:LEU:HD13	1:A:833:PRO:HD2	2.00	0.42
2:B:727:LEU:O	2:B:786:GLU:N	2.50	0.42
10:J:41:LEU:HD22	10:J:46:CYS:HB3	2.00	0.42
11:K:126:ASP:N	11:K:126:ASP:OD1	2.52	0.42
19:V:237:LEU:O	19:V:237:LEU:HD13	2.19	0.42
19:V:287:PRO:HA	19:V:288:PRO:HD3	1.86	0.42
18:U:107:ARG:NH1	20:W:469:ILE:HG22	2.34	0.42
2:B:408:LYS:HB2	16:P:119:HIS:CD2	2.55	0.42
12:L:29:TYR:HE1	12:L:58:LYS:HZ2	1.67	0.42
14:N:285:ALA:O	14:N:289:HIS:ND1	2.34	0.42
18:U:95:ASN:ND2	20:W:440:ASN:OD1	2.52	0.42
22:X:28:DA:H4'	22:X:29:DA:OP1	2.19	0.42
1:A:591:ASP:N	1:A:591:ASP:OD1	2.52	0.42
2:B:788:ARG:NH1	2:B:882:ASP:OD2	2.51	0.42
15:O:199:TYR:CD1	15:O:286:ARG:HD2	2.55	0.42
15:O:106:TYR:CE1	15:O:208:TYR:HE2	2.37	0.42
19:V:139:THR:HG23	19:V:141:HIS:H	1.85	0.42
2:B:114:PRO:HB3	2:B:174:LEU:HD22	2.00	0.42
5:E:3:GLN:HB3	5:E:7:ARG:NH2	2.34	0.42
5:E:85:GLU:HA	5:E:86:PRO:HD3	1.91	0.42
18:U:67:LEU:HD12	18:U:160:ILE:HG13	2.00	0.42
1:A:1261:GLN:HG3	1:A:1265:ARG:HH12	1.84	0.42
2:B:487:ARG:HD2	2:B:508:CYS:HB2	2.01	0.42
13:M:256:LYS:HD2	13:M:259:ILE:HB	2.01	0.42
19:V:94:LEU:HG	19:V:146:PHE:HE1	1.85	0.42
20:W:318:ARG:NH1	20:W:448:TYR:HB2	2.35	0.42
5:E:118:PRO:O	5:E:122:LYS:NZ	2.44	0.42
15:O:300:ALA:O	15:O:304:VAL:HG12	2.20	0.42
15:O:313:LYS:HE2	15:O:313:LYS:HB3	1.94	0.42
20:W:429:TYR:O	20:W:433:SER:N	2.53	0.42
1:A:403:THR:OG1	1:A:464:ARG:O	2.28	0.42
1:A:91:PHE:CD2	1:A:224:PRO:HG3	2.55	0.42
1:A:914:PHE:HE2	5:E:211:TYR:H	1.68	0.42
13:M:78:ILE:HG22	13:M:170:LEU:HB2	2.00	0.42
1:A:1076:PHE:CZ	1:A:1080:LYS:HD2	2.55	0.42
1:A:1325:VAL:HG12	1:A:1326:LEU:HD22	2.00	0.42
2:B:737:LYS:HB3	2:B:741:ILE:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:410:ARG:NE	6:F:104:ASN:OD1	2.53	0.42
9:I:18:ASP:N	9:I:18:ASP:OD1	2.53	0.42
16:P:173:GLU:O	16:P:174:PHE:HB2	2.20	0.42
18:U:107:ARG:HH21	20:W:473:LEU:HB2	1.85	0.42
18:U:175:LEU:HB3	18:U:237:PHE:CE2	2.54	0.42
22:X:70:DG:H1	23:Y:12:DC:H42	1.67	0.42
20:W:455:LYS:NZ	23:Y:57:DT:OP2	2.36	0.42
23:Y:61:DA:H4'	23:Y:62:DG:OP1	2.19	0.42
1:A:135:LEU:HD21	1:A:1378:LYS:HG2	2.00	0.42
1:A:558:ILE:HG12	1:A:666:LYS:HB2	2.01	0.42
1:A:610:PHE:CD1	1:A:613:LEU:HD21	2.55	0.42
1:A:856:LEU:HD23	1:A:861:PHE:HA	2.01	0.42
2:B:267:GLU:HG3	2:B:271:LEU:HD23	2.02	0.42
2:B:303:GLU:HA	2:B:325:GLU:HG2	2.01	0.42
8:H:107:VAL:N	8:H:111:LEU:O	2.52	0.42
19:V:509:ILE:H	19:V:509:ILE:HG13	1.71	0.42
2:B:540:ASP:HB3	2:B:543:LEU:HD13	2.01	0.42
2:B:777:SER:OG	2:B:941:ASP:OD1	2.37	0.42
15:O:159:ILE:HA	15:O:159:ILE:HD13	1.94	0.42
22:X:16:DG:H2''	22:X:17:DC:OP2	2.20	0.42
1:A:585:ASP:OD1	1:A:585:ASP:N	2.53	0.41
1:A:993:GLU:O	5:E:197:LYS:NZ	2.50	0.41
2:B:205:ILE:HG22	2:B:219:SER:HB2	2.00	0.41
2:B:225:HIS:O	2:B:226:GLU:HG2	2.20	0.41
2:B:732:GLN:HB2	2:B:753:GLN:HA	2.01	0.41
3:C:255:VAL:HG21	3:C:273:ASP:OD2	2.20	0.41
7:G:88:TRP:HE1	7:G:145:LYS:HG2	1.85	0.41
15:O:201:ILE:HG21	15:O:256:PRO:HB2	2.02	0.41
18:U:203:VAL:HG21	22:X:24:DA:H5'	2.01	0.41
20:W:488:ILE:HG22	20:W:492:ILE:HG23	2.01	0.41
1:A:1222:VAL:HG22	1:A:1232:ILE:HD13	2.03	0.41
1:A:498:PHE:HZ	1:A:519:LEU:HD22	1.85	0.41
2:B:203:ASN:HD21	2:B:222:SER:H	1.68	0.41
2:B:611:PRO:HA	2:B:648:TYR:HA	2.02	0.41
5:E:147:HIS:CE1	5:E:149:LEU:HD23	2.55	0.41
6:F:81:THR:OG1	6:F:144:GLU:OE2	2.25	0.41
13:M:170:LEU:HD21	14:N:307:PHE:CZ	2.54	0.41
16:P:90:GLU:HA	16:P:124:CYS:SG	2.60	0.41
1:A:429:GLY:O	1:A:465:HIS:ND1	2.53	0.41
1:A:808:GLN:HA	1:A:812:VAL:O	2.21	0.41
2:B:554:GLY:HA2	2:B:564:SER:HA	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:766:ASP:N	2:B:766:ASP:OD1	2.48	0.41
5:E:106:GLN:HA	5:E:129:PRO:HG2	2.02	0.41
15:O:108:GLN:HG2	15:O:117:THR:O	2.19	0.41
16:P:59:ASN:HB2	16:P:83:LYS:HZ1	1.84	0.41
18:U:90:ARG:HB3	19:V:473:HIS:CE1	2.55	0.41
19:V:233:ARG:HH21	20:W:285:PRO:HG3	1.85	0.41
1:A:607:LYS:HB3	1:A:659:ILE:HB	2.01	0.41
2:B:733:GLN:HG2	10:J:52:THR:O	2.20	0.41
3:C:328:LEU:HG	11:K:121:LEU:HD21	2.02	0.41
15:O:240:GLN:O	15:O:244:ASN:ND2	2.53	0.41
15:O:620:LEU:HD23	15:O:622:SER:H	1.85	0.41
2:B:325:GLU:N	2:B:325:GLU:OE1	2.54	0.41
2:B:135:TYR:OH	2:B:417:ASP:OD1	2.25	0.41
3:C:2:SER:O	3:C:4:ILE:HG13	2.20	0.41
1:A:15:GLY:HA2	1:A:1408:VAL:HG22	2.03	0.41
1:A:823:VAL:HG11	1:A:863:PHE:CD1	2.55	0.41
4:D:111:ASN:HB2	4:D:117:LYS:HE3	2.01	0.41
6:F:74:ILE:HG23	6:F:78:GLN:HG3	2.01	0.41
13:M:121:ILE:O	13:M:148:LEU:HD23	2.21	0.41
1:A:1301:ARG:HH21	1:A:1325:VAL:HA	1.86	0.41
1:A:1336:ILE:HD13	1:A:1355:ILE:HG22	2.02	0.41
1:A:892:SER:HB2	1:A:1371:ILE:HG23	2.03	0.41
2:B:379:LEU:HD23	2:B:382:GLN:NE2	2.36	0.41
2:B:882:ASP:HB2	2:B:901:ARG:HB2	2.02	0.41
7:G:31:ASN:O	7:G:35:ALA:N	2.53	0.41
16:P:106:TRP:CZ3	16:P:108:LYS:HB3	2.55	0.41
21:R:3:U:H2'	21:R:4:G:C8	2.56	0.41
19:V:230:LEU:HD22	19:V:233:ARG:HD3	2.02	0.41
1:A:498:PHE:CZ	1:A:519:LEU:HD22	2.56	0.41
1:A:511:ASP:OD1	1:A:511:ASP:N	2.54	0.41
1:A:11:LYS:HA	2:B:1145:ASP:HA	2.02	0.41
2:B:45:TRP:CE2	2:B:739:LYS:HD3	2.55	0.41
7:G:124:GLU:O	7:G:125:GLU:HG2	2.20	0.41
8:H:108:SER:OG	8:H:111:LEU:HD12	2.21	0.41
13:M:164:LYS:N	13:M:167:GLN:O	2.41	0.41
19:V:189:LEU:HD23	19:V:189:LEU:O	2.21	0.41
1:A:113:ILE:H	1:A:113:ILE:HG13	1.68	0.41
1:A:607:LYS:NZ	8:H:120:GLY:HA3	2.36	0.41
3:C:108:VAL:HB	3:C:184:VAL:HG12	2.03	0.41
1:A:1327:GLY:HA2	5:E:183:PRO:HD2	2.03	0.41
7:G:100:SER:HB2	7:G:108:ILE:HB	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:N:364:ARG:HH22	14:N:374:LYS:HB2	1.85	0.41
18:U:113:ALA:HB2	18:U:139:TYR:CE2	2.56	0.41
1:A:974:LEU:HG	1:A:1002:ARG:HE	1.87	0.41
2:B:185:PHE:HB3	2:B:187:VAL:HG23	2.03	0.41
13:M:242:ASN:OD1	13:M:243:ILE:N	2.54	0.41
15:O:137:ILE:HG21	17:Q:57:LYS:HG2	2.03	0.41
1:A:1079:ARG:HD3	6:F:84:TYR:CE2	2.56	0.40
2:B:64:SER:HB2	2:B:379:LEU:CD1	2.50	0.40
2:B:660:ALA:O	2:B:674:GLU:N	2.54	0.40
2:B:916:HIS:ND1	2:B:957:LYS:HG3	2.36	0.40
5:E:199:ILE:HG13	5:E:199:ILE:H	1.79	0.40
11:K:79:VAL:HG21	11:K:124:LEU:HB3	2.03	0.40
11:K:93:ILE:HA	11:K:94:PRO:HD3	1.95	0.40
13:M:164:LYS:HG3	13:M:165:ASP:N	2.36	0.40
1:A:39:LEU:HD12	1:A:53:ALA:HB1	2.03	0.40
1:A:580:LEU:HD11	1:A:590:PHE:CG	2.56	0.40
1:A:974:LEU:HD22	1:A:975:VAL:H	1.86	0.40
3:C:32:ASN:HB3	3:C:35:LYS:HB2	2.03	0.40
8:H:22:LYS:HD2	8:H:45:GLU:OE1	2.21	0.40
1:A:818:ILE:HD12	1:A:823:VAL:HA	2.02	0.40
2:B:1084:MET:HE2	2:B:1121:ILE:HA	2.02	0.40
2:B:209:ALA:HA	2:B:215:ILE:HG22	2.04	0.40
16:P:310:VAL:HG12	17:Q:41:LEU:HD23	2.02	0.40
16:P:64:VAL:HG22	16:P:65:LYS:H	1.86	0.40
18:U:112:THR:OG1	18:U:124:THR:O	2.26	0.40
19:V:152:VAL:HG21	19:V:157:ILE:HD11	2.04	0.40
19:V:239:ARG:HH22	20:W:284:PHE:H	1.70	0.40
19:V:469:GLU:OE2	20:W:476:LYS:HG3	2.22	0.40
23:Y:10:DC:H2'	23:Y:11:DA:C8	2.57	0.40
1:A:1386:LEU:HB3	1:A:1395:HIS:ND1	2.37	0.40
1:A:789:ASN:HD22	1:A:790:ALA:N	2.20	0.40
1:A:520:HIS:CB	2:B:1082:ARG:HH12	2.31	0.40
2:B:1092:VAL:HG21	2:B:1121:ILE:HD11	2.03	0.40
2:B:177:CYS:HB2	2:B:714:ALA:HB1	2.04	0.40
1:A:714:ILE:HD13	2:B:958:MET:HE2	2.03	0.40
4:D:68:ILE:HG23	4:D:72:PHE:HD2	1.86	0.40
5:E:171:LYS:H	5:E:174:GLN:NE2	2.20	0.40
7:G:10:LEU:HA	7:G:69:ASN:HA	2.04	0.40
13:M:159:TYR:CE2	14:N:309:LEU:HB2	2.56	0.40
16:P:116:LEU:HB3	16:P:120:VAL:HG13	2.03	0.40
16:P:179:LEU:HG	16:P:247:LEU:HD21	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:840:LYS:HG2	1:A:845:LYS:NZ	2.37	0.40
2:B:203:ASN:ND2	2:B:221:THR:OG1	2.54	0.40
2:B:481:ARG:NH2	2:B:482:LYS:HB2	2.37	0.40
5:E:18:THR:HG23	5:E:143:ASN:HB3	2.04	0.40
5:E:88:VAL:HG13	5:E:112:TYR:CE1	2.56	0.40
5:E:93:MET:HB3	5:E:123:LEU:HD11	2.03	0.40
8:H:58:THR:O	8:H:143:LEU:N	2.55	0.40
12:L:26:THR:OG1	12:L:27:LEU:N	2.54	0.40
16:P:102:ARG:NH2	16:P:154:GLN:HB2	2.37	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	1388/1460 (95%)	1233 (89%)	155 (11%)	0	100	100
2	B	1112/1149 (97%)	983 (88%)	128 (12%)	1 (0%)	55	89
3	C	333/335 (99%)	303 (91%)	30 (9%)	0	100	100
4	D	113/161 (70%)	101 (89%)	12 (11%)	0	100	100
5	E	212/215 (99%)	192 (91%)	20 (9%)	0	100	100
6	F	81/155 (52%)	72 (89%)	9 (11%)	0	100	100
7	G	174/212 (82%)	151 (87%)	22 (13%)	1 (1%)	28	71
8	H	136/146 (93%)	121 (89%)	14 (10%)	1 (1%)	25	68
9	I	32/110 (29%)	27 (84%)	5 (16%)	0	100	100
10	J	65/70 (93%)	59 (91%)	6 (9%)	0	100	100
11	K	99/142 (70%)	93 (94%)	6 (6%)	0	100	100
12	L	43/70 (61%)	39 (91%)	4 (9%)	0	100	100
13	M	179/282 (64%)	160 (89%)	19 (11%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
14	N	100/422 (24%)	93 (93%)	7 (7%)	0	100	100
15	O	528/654 (81%)	496 (94%)	32 (6%)	0	100	100
16	P	238/317 (75%)	219 (92%)	19 (8%)	0	100	100
17	Q	33/251 (13%)	29 (88%)	4 (12%)	0	100	100
18	U	178/240 (74%)	163 (92%)	15 (8%)	0	100	100
19	V	331/596 (56%)	305 (92%)	26 (8%)	0	100	100
20	W	214/594 (36%)	195 (91%)	19 (9%)	0	100	100
All	All	5589/7581 (74%)	5034 (90%)	552 (10%)	3 (0%)	58	89

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
7	G	79	PRO
2	B	769	ASP
8	H	110	ASP

### 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	1214/1257 (97%)	1189 (98%)	25 (2%)	59	81
2	B	975/1006 (97%)	958 (98%)	17 (2%)	66	85
3	C	296/296 (100%)	291 (98%)	5 (2%)	66	85
4	D	110/145 (76%)	104 (94%)	6 (6%)	25	60
5	E	196/197 (100%)	192 (98%)	4 (2%)	60	82
6	F	73/137 (53%)	72 (99%)	1 (1%)	71	86
7	G	160/190 (84%)	158 (99%)	2 (1%)	73	87
8	H	123/128 (96%)	123 (100%)	0	100	100
9	I	31/98 (32%)	30 (97%)	1 (3%)	44	72
10	J	62/65 (95%)	62 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
11	K	91/130 (70%)	90 (99%)	1 (1%)	78	89
12	L	40/57 (70%)	40 (100%)	0	100	100
13	M	160/249 (64%)	156 (98%)	4 (2%)	53	77
14	N	88/360 (24%)	87 (99%)	1 (1%)	78	89
15	O	490/593 (83%)	478 (98%)	12 (2%)	54	78
16	P	227/285 (80%)	220 (97%)	7 (3%)	45	73
17	Q	31/212 (15%)	31 (100%)	0	100	100
18	U	152/205 (74%)	147 (97%)	5 (3%)	43	72
19	V	295/513 (58%)	290 (98%)	5 (2%)	66	85
20	W	194/534 (36%)	188 (97%)	6 (3%)	45	73
All	All	5008/6657 (75%)	4906 (98%)	102 (2%)	63	82

All (102) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	129	ARG
1	A	136	ARG
1	A	137	ARG
1	A	213	ARG
1	A	274	MET
1	A	303	LEU
1	A	310	ASN
1	A	348	LYS
1	A	410	ARG
1	A	483	LEU
1	A	499	ARG
1	A	783	ASN
1	A	789	ASN
1	A	848	VAL
1	A	944	ASN
1	A	1039	LEU
1	A	1155	ARG
1	A	1187	ARG
1	A	1227	LYS
1	A	1265	ARG
1	A	1286	ARG
1	A	1294	LEU
1	A	1332	ARG

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Mol	Chain	Res	Type
1	A	1350	VAL
1	A	1439	LYS
2	B	66	ASN
2	B	139	ARG
2	B	245	ASN
2	B	298	GLN
2	B	428	ASN
2	B	481	ARG
2	B	710	ILE
2	B	749	LEU
2	B	773	LEU
2	B	859	ASN
2	B	867	ARG
2	B	881	ILE
2	B	970	ASN
2	B	1038	ARG
2	B	1054	ARG
2	B	1135	MET
2	B	1136	ASN
3	C	1	MET
3	C	11	ARG
3	C	87	ASN
3	C	174	ARG
3	C	277	ARG
4	D	60	ARG
4	D	71	ASN
4	D	98	MET
4	D	120	LYS
4	D	130	ASN
4	D	153	MET
5	E	52	ARG
5	E	58	MET
5	E	75	MET
5	E	215	MET
6	F	90	ARG
7	G	73	ARG
7	G	113	ASN
9	I	2	LEU
11	K	74	ASN
13	M	144	ASN
13	M	163	VAL
13	M	166	MET

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Mol	Chain	Res	Type
13	M	220	ASN
14	N	300	LYS
15	O	73	ARG
15	O	107	LEU
15	O	184	LYS
15	O	205	LYS
15	O	238	ARG
15	O	243	MET
15	O	252	ILE
15	O	268	LYS
15	O	286	ARG
15	O	580	ASN
15	O	631	ASN
15	O	636	ASN
16	P	23	MET
16	P	55	LEU
16	P	83	LYS
16	P	189	ASN
16	P	239	ASN
16	P	271	MET
16	P	298	LYS
18	U	104	MET
18	U	110	LYS
18	U	141	ARG
18	U	159	ASN
18	U	239	LYS
19	V	115	ASN
19	V	149	ARG
19	V	237	LEU
19	V	269	LYS
19	V	481	LYS
20	W	292	ASN
20	W	400	ASN
20	W	413	ARG
20	W	451	ARG
20	W	521	ASN
20	W	532	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (81) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	310	ASN

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Mol	Chain	Res	Type
1	A	361	GLN
1	A	367	ASN
1	A	481	HIS
1	A	520	HIS
1	A	578	GLN
1	A	760	GLN
1	A	783	ASN
1	A	815	GLN
1	A	828	GLN
1	A	950	GLN
1	A	1095	GLN
1	A	1131	ASN
1	A	1197	GLN
1	A	1346	HIS
1	A	1354	HIS
1	A	1419	GLN
2	B	66	ASN
2	B	81	GLN
2	B	144	HIS
2	B	159	ASN
2	B	203	ASN
2	B	225	HIS
2	B	245	ASN
2	B	334	HIS
2	B	428	ASN
2	B	434	ASN
2	B	595	HIS
2	B	596	GLN
2	B	600	HIS
2	B	693	HIS
2	B	821	HIS
2	B	859	ASN
2	B	947	HIS
2	B	970	ASN
2	B	994	GLN
2	B	1029	HIS
2	B	1136	ASN
2	B	1149	GLN
3	C	87	ASN
3	C	99	HIS
3	C	296	ASN
4	D	71	ASN

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Mol	Chain	Res	Type
4	D	126	GLN
4	D	130	ASN
5	E	3	GLN
5	E	174	GLN
5	E	179	GLN
6	F	78	GLN
7	G	31	ASN
7	G	113	ASN
9	I	9	ASN
10	J	53	HIS
11	K	64	GLN
13	M	92	ASN
13	M	141	ASN
13	M	144	ASN
13	M	190	ASN
13	M	220	ASN
13	M	224	GLN
15	O	152	HIS
15	O	161	GLN
15	O	225	ASN
15	O	240	GLN
15	O	244	ASN
15	O	283	ASN
15	O	549	GLN
15	O	572	HIS
15	O	580	ASN
15	O	631	ASN
16	P	154	GLN
16	P	189	ASN
16	P	239	ASN
18	U	91	ASN
18	U	159	ASN
19	V	115	ASN
19	V	473	HIS
20	W	292	ASN
20	W	391	ASN
20	W	400	ASN
20	W	521	ASN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
21	R	5/6 (83%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 7 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.