



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 05:16 AM UTC

PDB ID : 6OGZ / pdb_00006ogz
EMDB ID : EMD-20060
Title : In situ structure of Rotavirus RNA-dependent RNA polymerase at transcript-elongated state
Authors : Ding, K.; Chang, T.; Shen, W.; Roy, P.; Zhou, Z.H.
Deposited on : 2019-04-03
Resolution : 3.60 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

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>.

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

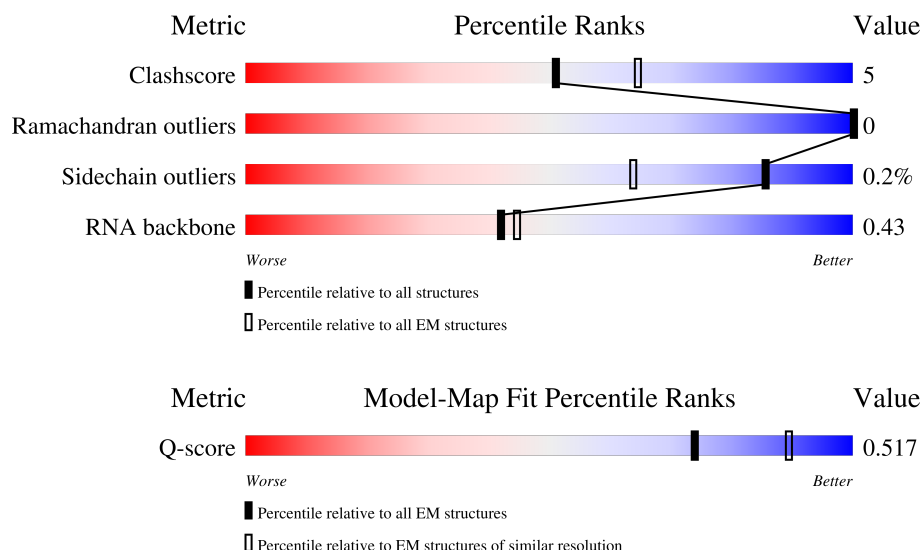
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
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 i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.60 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.



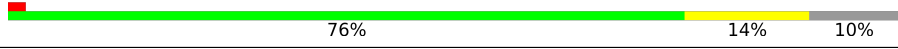
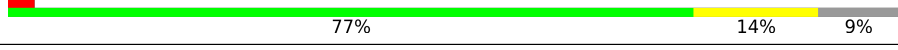
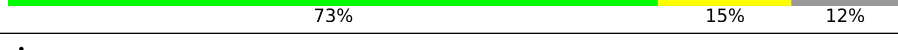
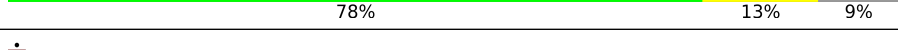
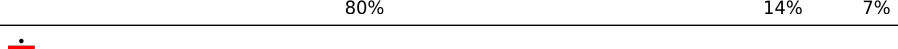
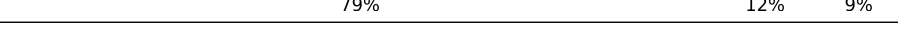
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	12797 (3.10 - 4.10)

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 ≥ 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	B	18	
2	C	17	
3	A	1088	

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Mol	Chain	Length	Quality of chain
4	E	887	 76%13%10%
4	F	887	 79%12%9%
4	G	887	 76%14%10%
4	H	887	 77%14%9%
4	I	887	 71%17%12%
4	J	887	 79%12%9%
4	K	887	 73%15%12%
4	L	887	 78%13%9%
4	M	887	 80%14%7%
4	N	887	 79%12%9%

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 RNA chain called RNA (5'-R(P*UP*AP*UP*AP*UP*AP*UP*AP*UP*AP*UP*AP*UP*AP*UP*AP*UP*A)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	18	Total	C	N	O	P	0	0
			378	171	63	126	18		

- Molecule 2 is a RNA chain called RNA (5'-R(P*AP*UP*AP*UP*AP*UP*AP*UP*AP*UP*AP*UP*AP*UP*AP*UP*A)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
2	C	17	Total	C	N	O	P	0	0
			358	162	61	118	17		

- Molecule 3 is a protein called RNA-dependent RNA polymerase of rotavirus A.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	A	1082	Total	C	N	O	S	0	0
			8754	5609	1450	1656	39		

- Molecule 4 is a protein called Inner capsid protein VP2.

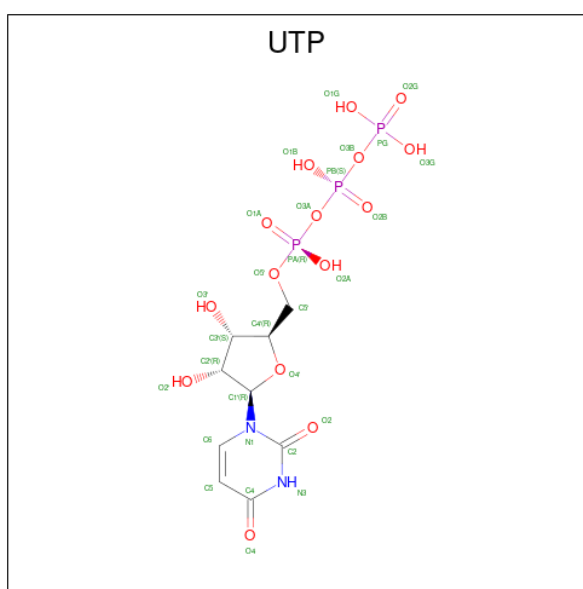
Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	794	Total 6492	C 4127	N 1117	O 1212	S 36	0	0
4	F	803	Total 6570	C 4174	N 1131	O 1229	S 36	0	0
4	G	794	Total 6487	C 4120	N 1119	O 1212	S 36	0	0
4	I	781	Total 6379	C 4052	N 1101	O 1190	S 36	0	0
4	H	805	Total 6587	C 4186	N 1134	O 1231	S 36	0	0
4	J	805	Total 6587	C 4186	N 1134	O 1231	S 36	0	0

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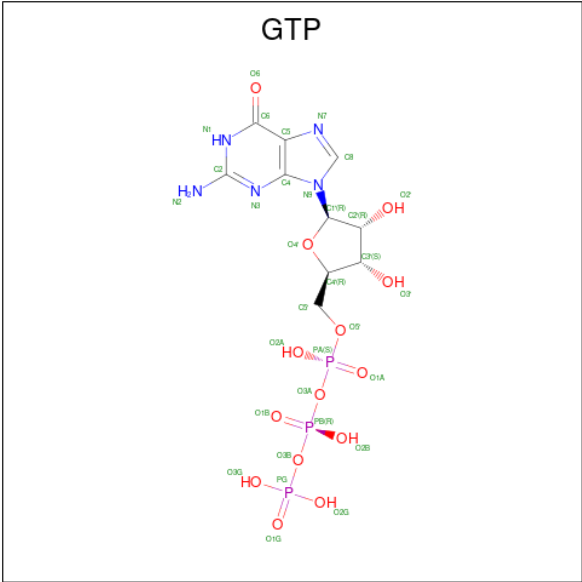
Mol	Chain	Residues	Atoms					AltConf	Trace
4	K	781	Total	C	N	O	S	0	0
			6379	4052	1101	1190	36		
4	L	805	Total	C	N	O	S	0	0
			6587	4186	1134	1231	36		
4	M	826	Total	C	N	O	S	0	0
			6756	4292	1162	1266	36		
4	N	805	Total	C	N	O	S	0	0
			6587	4186	1134	1231	36		

- Molecule 5 is URIDINE 5'-TRIPHOSPHATE (CCD ID: UTP) (formula: $C_9H_{15}N_2O_{15}P_3$).

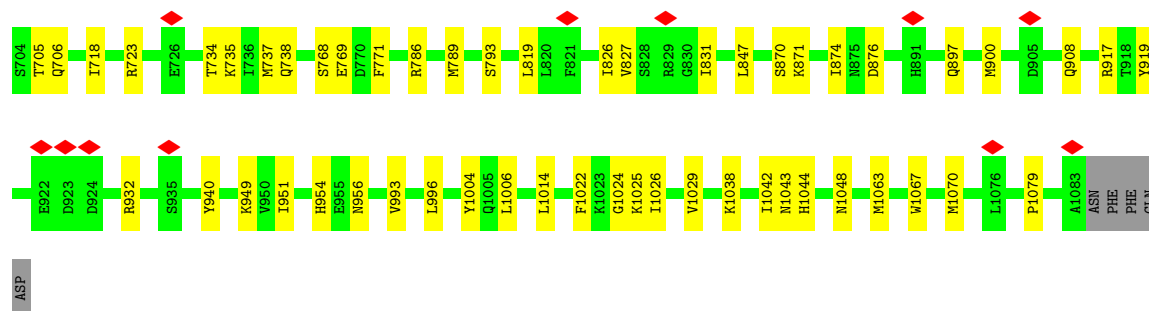


Mol	Chain	Residues	Atoms					AltConf
5	A	1	Total	C	N	O	P	0
			29	9	2	15	3	

- Molecule 6 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).

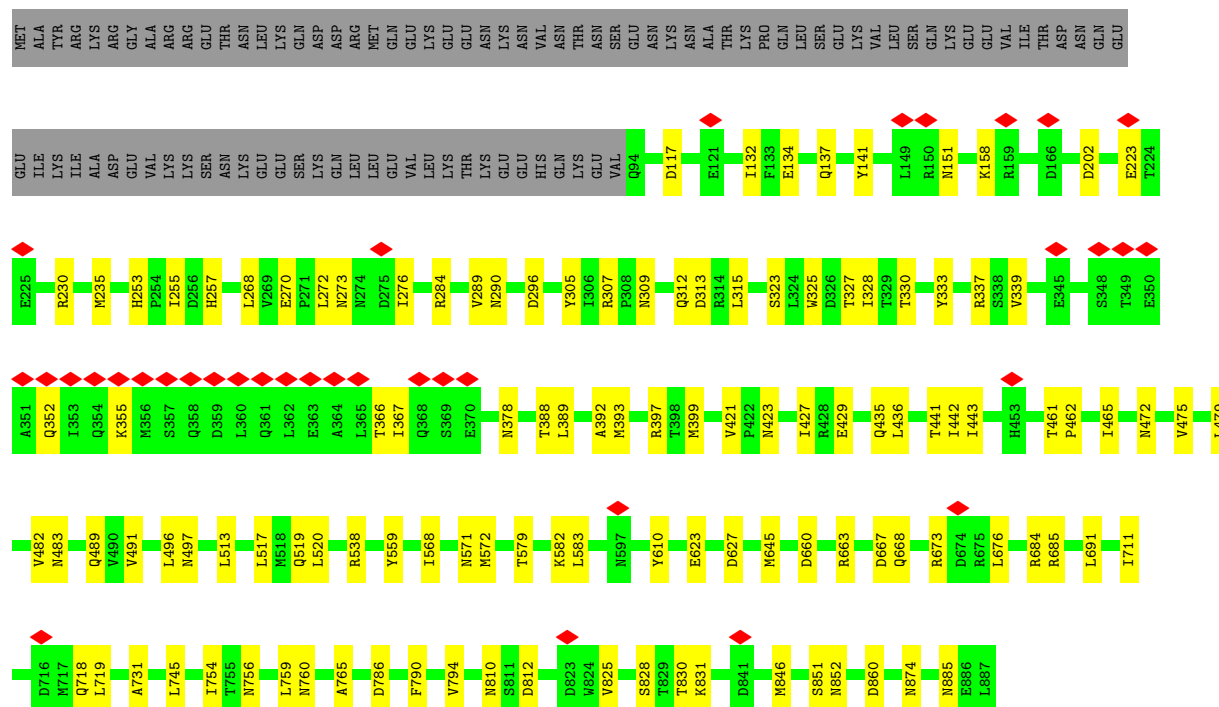


Mol	Chain	Residues	Atoms					AltConf
6	A	1	Total	C	N	O	P	0
			32	10	5	14	3	



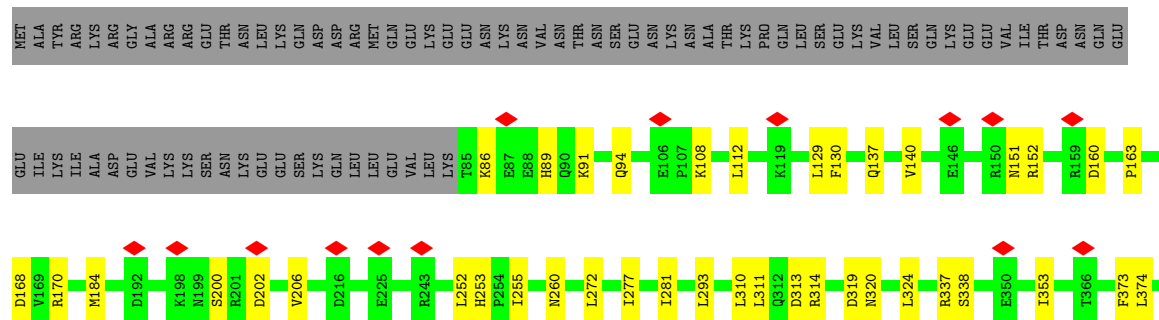
• Molecule 4: Inner capsid protein VP2

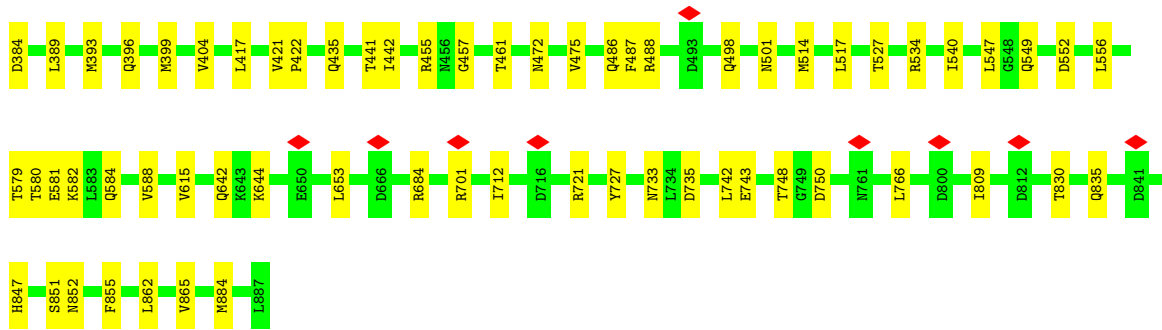
Chain E: 76% 13% 10%



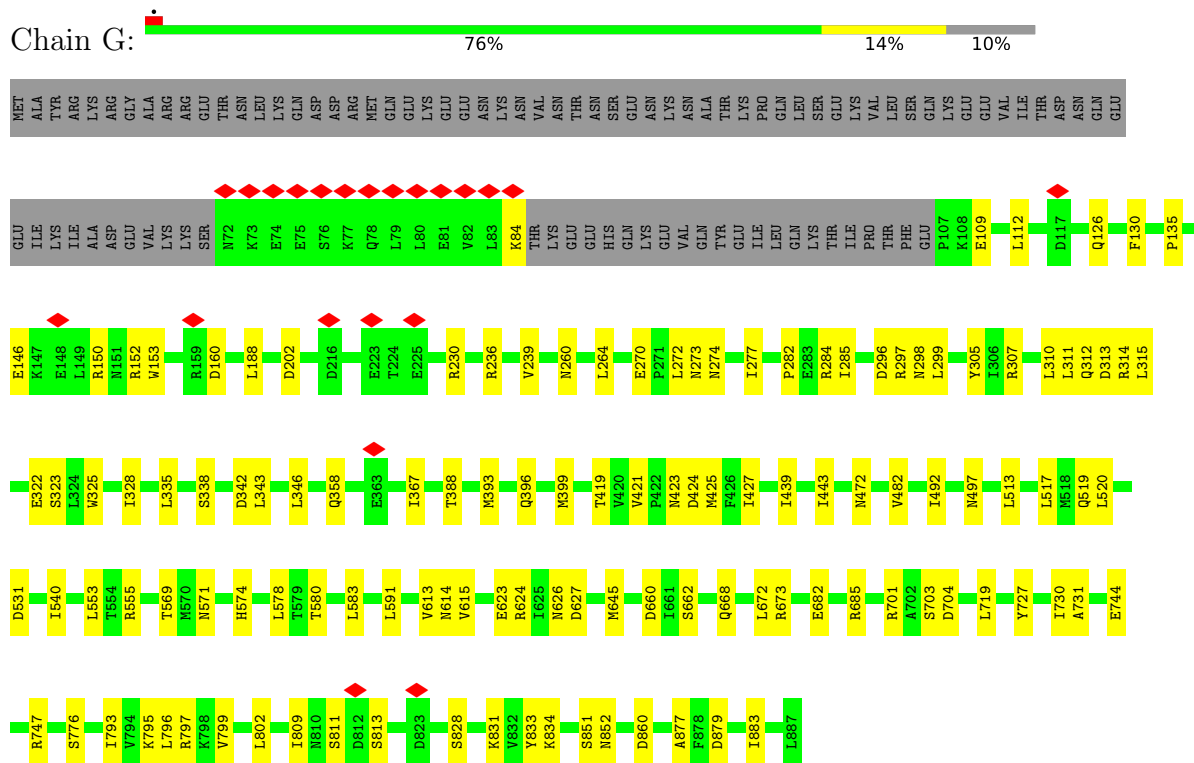
• Molecule 4: Inner capsid protein VP2

Chain F: 79% 12% 9%

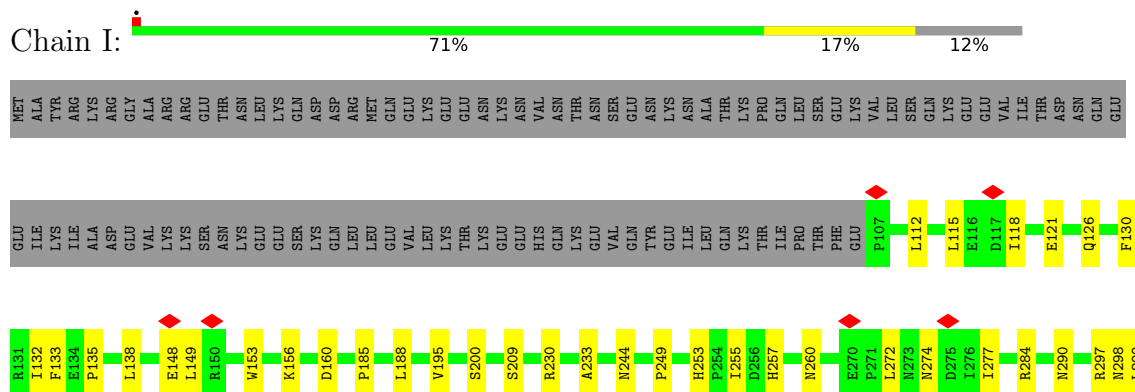




• Molecule 4: Inner capsid protein VP2



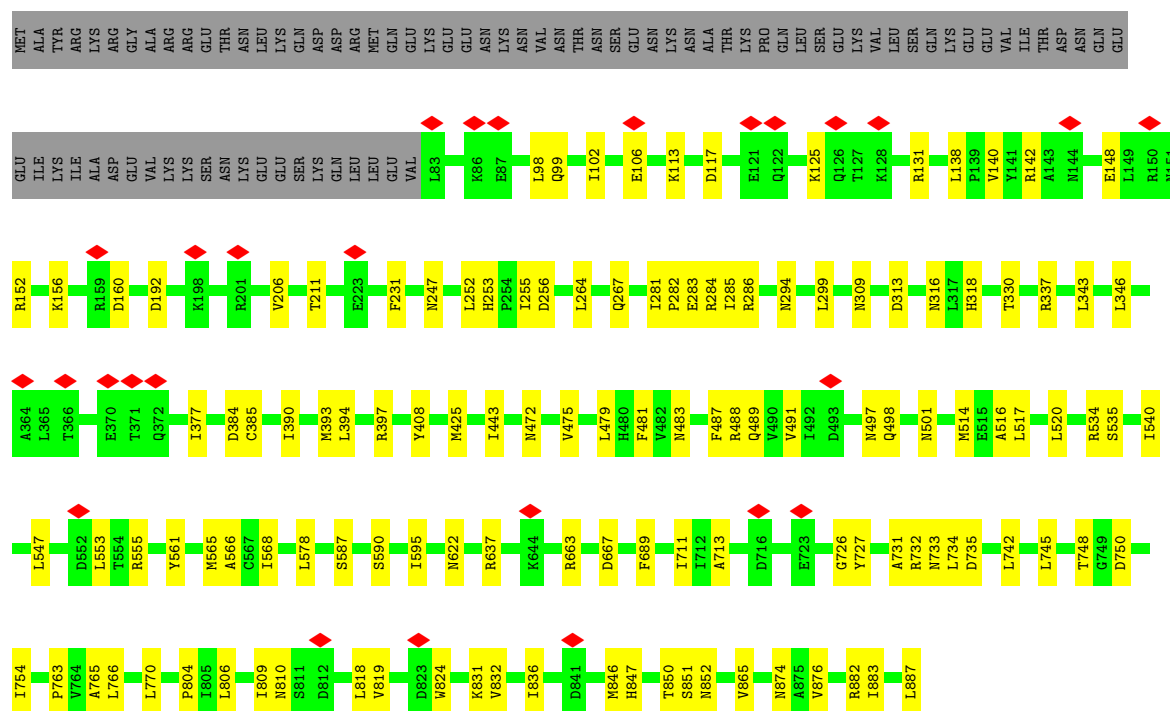
• Molecule 4: Inner capsid protein VP2





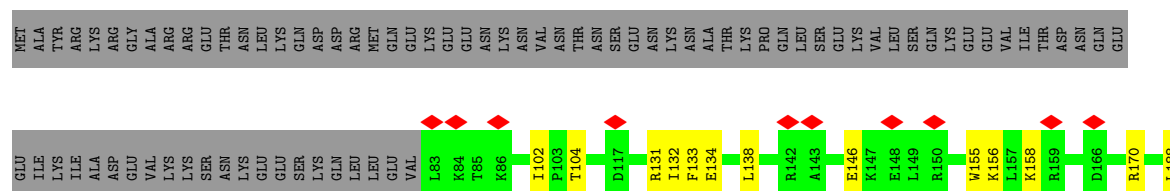
• Molecule 4: Inner capsid protein VP2

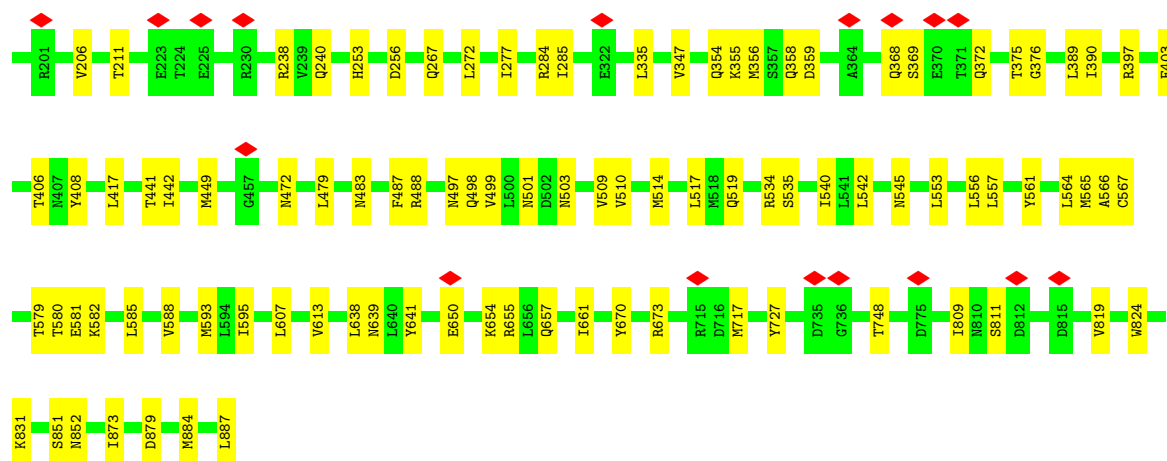
Chain H: 77% 14% 9%



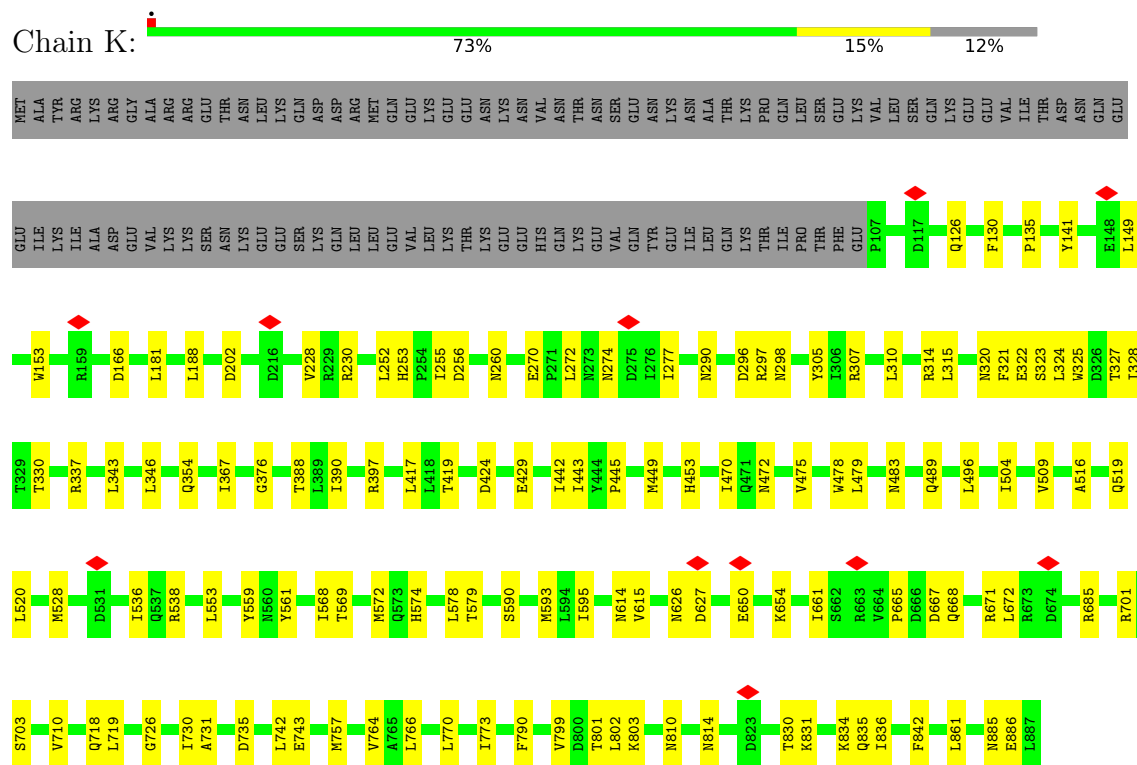
• Molecule 4: Inner capsid protein VP2

Chain J: 79% 12% 9%

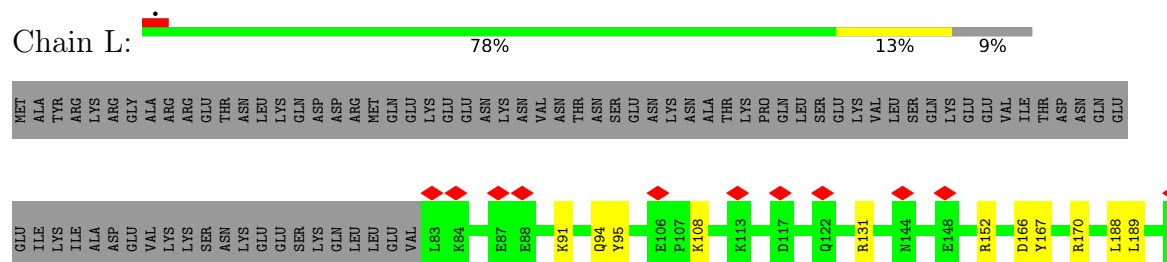


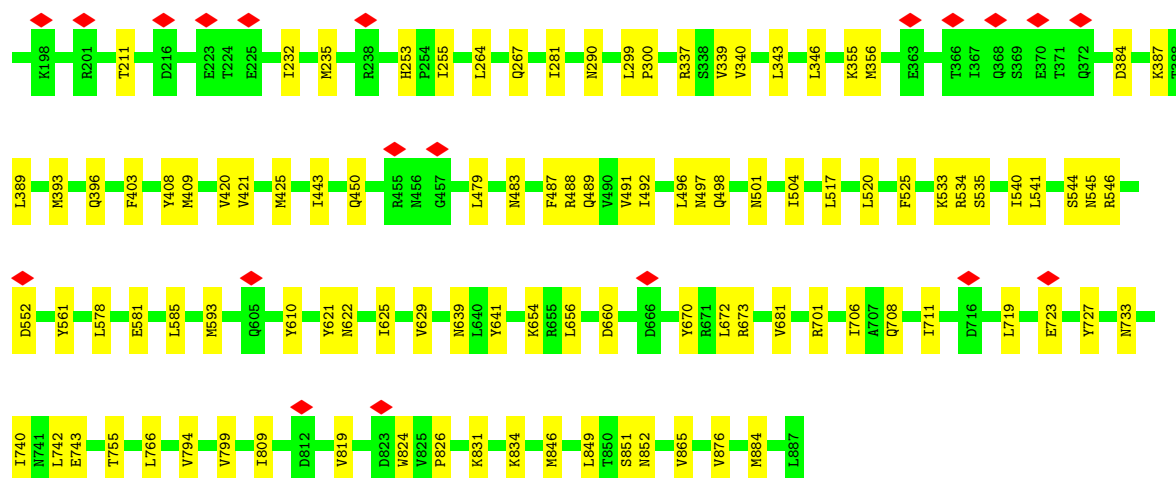


• Molecule 4: Inner capsid protein VP2



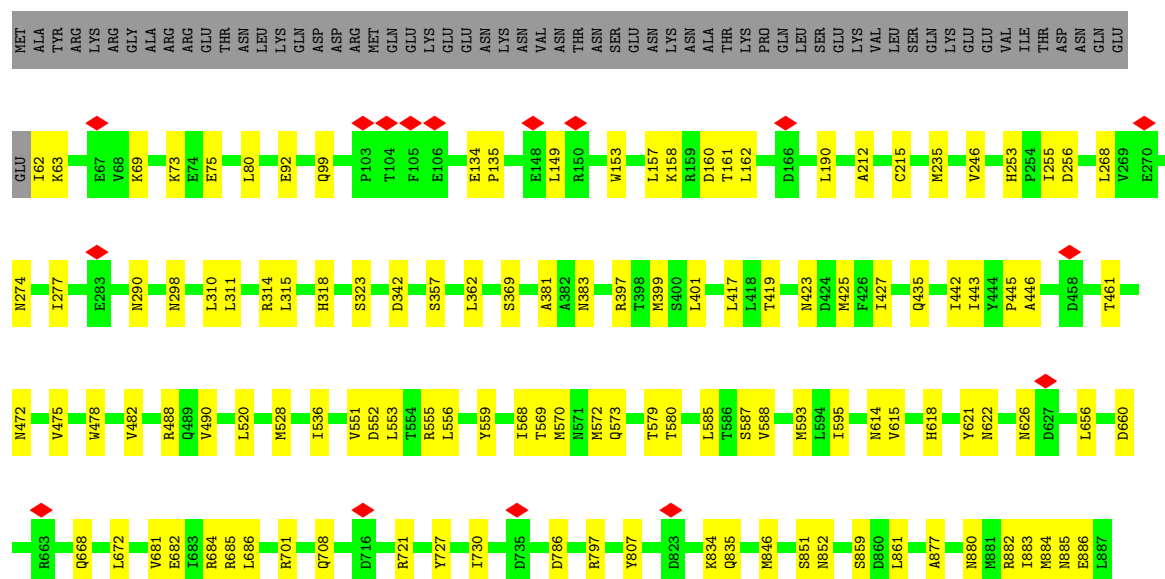
• Molecule 4: Inner capsid protein VP2





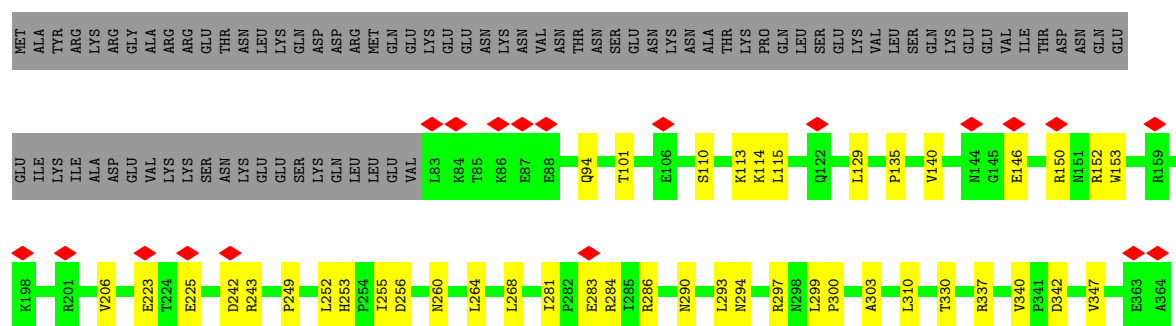
• Molecule 4: Inner capsid protein VP2

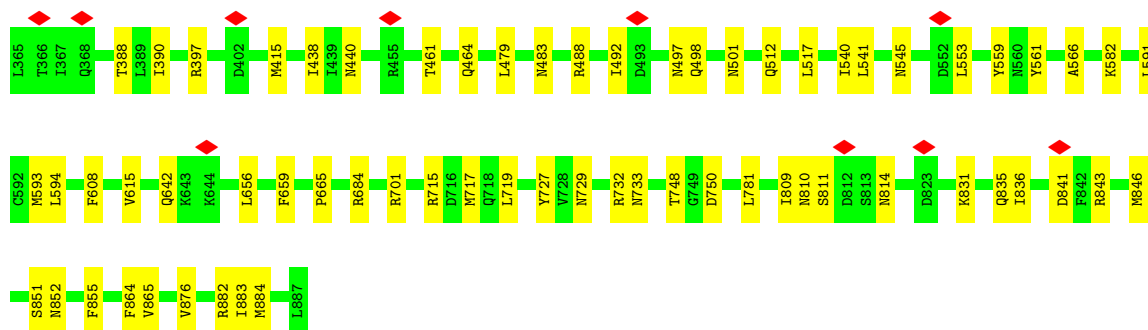
Chain M: 80% 14% 7%



• Molecule 4: Inner capsid protein VP2

Chain N: 79% 12% 9%





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	411438	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 ($e^-/\text{\AA}^2$)	18	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.135	Depositor
Minimum map value	-0.096	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.0175	Depositor
Map size (Å)	353.76, 353.76, 353.76	wwPDB
Map dimensions	264, 264, 264	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.34, 1.34, 1.34	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: GTP, UTP

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	B	0.20	0/422	0.29	0/653
2	C	0.21	0/400	0.31	0/619
3	A	0.32	0/8926	0.48	0/12069
4	E	0.35	0/6611	0.53	0/8969
4	F	0.33	0/6690	0.49	0/9074
4	G	0.36	0/6602	0.51	0/8950
4	H	0.33	0/6707	0.50	0/9096
4	I	0.36	0/6495	0.51	0/8810
4	J	0.32	0/6707	0.49	0/9096
4	K	0.36	0/6495	0.53	0/8810
4	L	0.33	0/6707	0.50	0/9096
4	M	0.35	0/6876	0.51	0/9320
4	N	0.33	0/6707	0.51	0/9096
All	All	0.34	0/76345	0.50	0/103658

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	B	378	0	190	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	358	0	180	0	0
3	A	8754	0	8840	105	0
4	E	6492	0	6520	75	0
4	F	6570	0	6595	65	0
4	G	6487	0	6525	80	0
4	H	6587	0	6619	80	0
4	I	6379	0	6408	95	0
4	J	6587	0	6619	76	0
4	K	6379	0	6408	91	0
4	L	6587	0	6619	73	0
4	M	6756	0	6803	72	0
4	N	6587	0	6619	65	0
5	A	29	0	11	1	0
6	A	32	0	12	1	0
All	All	74962	0	74968	817	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 5.

All (817) 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)
4:I:148:GLU:O	4:I:149:LEU:HG	1.45	1.13
3:A:1026:ILE:HG22	3:A:1029:VAL:HG12	1.48	0.94
4:H:637:ARG:HD2	4:H:689:PHE:HE2	1.33	0.92
4:H:637:ARG:HD2	4:H:689:PHE:CE2	2.14	0.82
4:F:488:ARG:HE	4:F:501:ASN:HB3	1.52	0.75
3:A:1026:ILE:CG2	3:A:1029:VAL:HG12	2.19	0.72
4:K:322:GLU:HB3	4:K:578:LEU:HD13	1.70	0.72
4:L:488:ARG:HE	4:L:501:ASN:HB3	1.55	0.71
4:K:320:ASN:CG	4:K:671:ARG:NH1	2.49	0.71
4:J:188:LEU:H	4:J:267:GLN:HE22	1.38	0.71
4:H:488:ARG:HE	4:H:501:ASN:HB3	1.56	0.70
4:L:420:VAL:HA	4:L:504:ILE:HD11	1.74	0.69
4:J:488:ARG:HE	4:J:501:ASN:HB3	1.58	0.69
4:H:637:ARG:CD	4:H:689:PHE:HE2	2.05	0.68
4:J:650:GLU:HG3	4:J:661:ILE:HG21	1.74	0.68
4:K:323:SER:HB3	4:K:668:GLN:HE22	1.60	0.67
4:N:140:VAL:HG11	4:N:152:ARG:HE	1.60	0.67
4:L:188:LEU:H	4:L:267:GLN:HE22	1.44	0.65
3:A:1026:ILE:HG22	3:A:1029:VAL:CG1	2.24	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:343:LEU:HD23	4:G:346:LEU:HD21	1.79	0.65
3:A:517:LEU:HD11	3:A:655:VAL:HG21	1.79	0.65
3:A:819:LEU:HD11	3:A:940:TYR:HB2	1.79	0.64
3:A:184:ARG:H	3:A:195:VAL:HG21	1.63	0.64
3:A:900:MET:HE1	3:A:1006:LEU:HD22	1.79	0.64
4:L:253:HIS:HD2	4:L:255:ILE:H	1.46	0.64
4:E:323:SER:HB3	4:E:668:GLN:HE22	1.63	0.64
4:G:126:GLN:HE21	4:G:188:LEU:HD11	1.63	0.63
4:J:284:ARG:HE	4:J:566:ALA:HB2	1.63	0.63
4:K:324:LEU:HG	4:K:668:GLN:HE21	1.64	0.62
4:K:322:GLU:HA	4:K:322:GLU:OE1	2.00	0.62
4:J:389:LEU:HD23	4:J:557:LEU:HD11	1.82	0.62
4:K:320:ASN:CG	4:K:671:ARG:HH12	2.08	0.62
4:K:320:ASN:ND2	4:L:545:ASN:OD1	2.28	0.62
3:A:244:VAL:HA	3:A:295:MET:HE3	1.82	0.62
4:F:517:LEU:HG	4:F:540:ILE:HG23	1.82	0.62
4:I:663:ARG:HG3	4:J:354:GLN:HE21	1.63	0.62
4:I:272:LEU:HD22	4:I:305:TYR:HB3	1.82	0.61
4:M:274:ASN:HD21	4:M:298:ASN:HA	1.64	0.61
4:N:488:ARG:HE	4:N:501:ASN:HB3	1.65	0.61
4:N:310:LEU:HD21	4:N:615:VAL:HG13	1.81	0.61
4:N:253:HIS:HD2	4:N:255:ILE:H	1.49	0.61
4:E:389:LEU:HD23	4:E:399:MET:HE1	1.83	0.61
4:E:230:ARG:NH2	4:N:750:ASP:OD2	2.34	0.60
4:M:290:ASN:ND2	4:M:559:TYR:OH	2.34	0.60
4:H:727:TYR:HB2	4:H:809:ILE:HB	1.83	0.60
4:F:94:GLN:HE22	4:G:860:ASP:H	1.50	0.60
4:J:390:ILE:HD11	4:J:557:LEU:HG	1.83	0.60
4:M:323:SER:HB3	4:M:668:GLN:HE22	1.67	0.60
4:H:487:PHE:HB3	4:H:498:GLN:HB3	1.84	0.60
4:H:102:ILE:HD11	4:H:578:LEU:HD22	1.84	0.59
4:H:281:ILE:HD11	4:H:865:VAL:HG11	1.83	0.59
4:K:126:GLN:HE21	4:K:188:LEU:HD11	1.67	0.59
4:E:284:ARG:HH12	4:E:497:ASN:HD21	1.48	0.59
4:K:290:ASN:ND2	4:K:559:TYR:OH	2.36	0.59
4:I:130:PHE:H	4:I:260:ASN:HD21	1.50	0.59
4:I:552:ASP:OD2	4:I:884:MET:SD	2.61	0.59
4:K:320:ASN:ND2	4:K:671:ARG:NH1	2.50	0.59
4:E:429:GLU:OE1	4:F:534:ARG:NH2	2.36	0.59
4:K:253:HIS:HD2	4:K:255:ILE:H	1.51	0.59
4:N:249:PRO:HB3	4:N:846:MET:HE1	1.83	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:J:487:PHE:HB3	4:J:498:GLN:HB3	1.85	0.58
4:I:505:ARG:NE	4:I:562:GLU:OE2	2.37	0.58
4:J:368:GLN:HE22	4:J:406:THR:HG21	1.68	0.58
4:M:423:ASN:ND2	4:M:482:VAL:O	2.37	0.58
4:G:272:LEU:HD22	4:G:305:TYR:HB3	1.85	0.58
4:G:311:LEU:HD11	4:G:314:ARG:HH21	1.68	0.58
4:L:94:GLN:HE22	4:M:859:SER:H	1.50	0.58
3:A:531:THR:OG1	3:A:532:GLN:N	2.36	0.58
4:N:492:ILE:HD11	4:N:497:ASN:HD21	1.68	0.58
3:A:1025:LYS:HB3	4:I:365:LEU:HD12	1.86	0.58
4:G:682:GLU:OE2	4:G:685:ARG:NH1	2.37	0.58
4:J:285:ILE:HD13	4:J:607:LEU:HD21	1.85	0.58
4:M:397:ARG:NH2	4:M:579:THR:OG1	2.37	0.58
4:J:654:LYS:HB2	4:J:661:ILE:HD11	1.86	0.58
4:K:429:GLU:OE2	4:L:534:ARG:NH2	2.36	0.58
4:N:281:ILE:HD11	4:N:865:VAL:HG11	1.84	0.58
4:I:290:ASN:ND2	4:I:559:TYR:OH	2.37	0.58
4:J:517:LEU:HG	4:J:540:ILE:HG23	1.84	0.58
4:J:819:VAL:HA	4:J:824:TRP:HE1	1.68	0.58
4:F:163:PRO:HB2	4:F:168:ASP:HB3	1.86	0.57
4:I:442:ILE:HG12	4:I:536:ILE:HD11	1.84	0.57
1:B:585:A:O2'	4:E:538:ARG:NH1	2.36	0.57
4:F:527:THR:O	4:F:527:THR:HG22	2.04	0.57
4:I:320:ASN:O	4:I:671:ARG:NH1	2.36	0.57
3:A:352:THR:HG22	3:A:354:ASP:H	1.69	0.57
4:G:793:ILE:HD12	4:G:799:VAL:HG12	1.87	0.57
4:J:564:LEU:HD23	4:J:607:LEU:HD13	1.87	0.57
4:L:517:LEU:HG	4:L:540:ILE:HG23	1.85	0.57
3:A:1026:ILE:CG2	3:A:1029:VAL:CG1	2.82	0.57
4:K:141:TYR:CE2	4:K:810:ASN:ND2	2.73	0.57
4:G:270:GLU:OE1	4:G:307:ARG:NH2	2.38	0.57
3:A:951:ILE:O	4:M:99:GLN:NE2	2.38	0.57
4:M:730:ILE:HD12	4:M:834:LYS:HB3	1.87	0.57
4:H:479:LEU:O	4:H:483:ASN:ND2	2.38	0.57
4:L:492:ILE:HD11	4:L:497:ASN:HD21	1.70	0.57
4:M:701:ARG:HD3	4:M:835:GLN:HG3	1.86	0.57
4:E:270:GLU:OE1	4:E:307:ARG:NH2	2.38	0.56
4:E:333:TYR:OH	4:E:337:ARG:NH2	2.35	0.56
4:G:472:ASN:ND2	4:G:519:GLN:OE1	2.38	0.56
4:G:310:LEU:HD11	4:G:615:VAL:HG13	1.87	0.56
3:A:1026:ILE:HD12	3:A:1063:MET:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:290:ASN:ND2	4:E:559:TYR:OH	2.38	0.56
4:F:337:ARG:NH2	4:F:384:ASP:OD1	2.36	0.56
4:K:479:LEU:O	4:K:483:ASN:ND2	2.38	0.56
4:K:553:LEU:HD12	4:K:595:ILE:HG21	1.86	0.56
4:F:579:THR:HG22	4:G:297:ARG:HE	1.69	0.56
4:I:623:GLU:O	4:I:624:ARG:NH2	2.39	0.56
4:F:170:ARG:NH2	4:F:743:GLU:OE2	2.38	0.56
4:K:343:LEU:HD23	4:K:346:LEU:HD21	1.88	0.56
4:L:131:ARG:NH2	4:L:211:THR:OG1	2.39	0.56
4:M:314:ARG:NH1	4:M:626:ASN:OD1	2.39	0.56
4:L:443:ILE:HG12	4:L:520:LEU:HD11	1.88	0.56
4:G:427:ILE:HD11	4:H:534:ARG:HH12	1.71	0.56
3:A:516:VAL:HG23	3:A:636:VAL:HG22	1.87	0.55
4:E:462:PRO:HA	4:E:465:ILE:HG22	1.88	0.55
4:E:568:ILE:HD13	4:E:572:MET:HE3	1.86	0.55
3:A:1026:ILE:HG23	3:A:1067:TRP:HB2	1.88	0.55
4:F:472:ASN:HB3	4:F:475:VAL:HG12	1.88	0.55
4:L:337:ARG:NH2	4:L:384:ASP:OD1	2.39	0.55
3:A:1026:ILE:O	3:A:1029:VAL:HG12	2.06	0.55
4:E:230:ARG:HG3	4:N:748:THR:HB	1.89	0.55
4:G:274:ASN:HD21	4:G:298:ASN:HA	1.71	0.55
4:K:320:ASN:O	4:K:671:ARG:NH1	2.38	0.55
4:H:253:HIS:HD2	4:H:255:ILE:H	1.54	0.55
4:J:654:LYS:CA	4:J:661:ILE:HD11	2.37	0.55
4:M:381:ALA:HB1	4:M:587:SER:HB2	1.89	0.55
4:M:721:ARG:NH1	4:M:727:TYR:OH	2.40	0.55
4:E:312:GLN:HG2	4:E:571:ASN:HD22	1.72	0.54
4:I:249:PRO:HB3	4:I:846:MET:HE3	1.90	0.54
4:H:555:ARG:HB3	4:H:883:ILE:HG22	1.88	0.54
4:K:489:GLN:HB3	4:K:496:LEU:HD11	1.89	0.54
4:F:281:ILE:HD11	4:F:865:VAL:HG11	1.90	0.54
4:L:727:TYR:HB2	4:L:809:ILE:HB	1.89	0.54
3:A:460:ARG:HH22	5:A:1101:UTP:H3'	1.72	0.54
4:J:579:THR:HG22	4:K:297:ARG:HE	1.72	0.54
4:E:397:ARG:NH2	4:E:579:THR:O	2.40	0.54
4:F:487:PHE:HB3	4:F:498:GLN:HB3	1.89	0.54
4:N:852:ASN:HD22	4:N:855:PHE:HE1	1.53	0.54
4:F:642:GLN:HG2	4:G:202:ASP:HA	1.89	0.54
4:M:69:LYS:HA	4:M:80:LEU:HD11	1.88	0.54
4:N:150:ARG:NH1	4:N:223:GLU:OE1	2.40	0.54
4:J:369:SER:HB2	4:J:372:GLN:HB2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:479:LEU:O	4:L:483:ASN:ND2	2.41	0.54
4:N:701:ARG:HD2	4:N:835:GLN:HG3	1.90	0.54
4:F:580:THR:OG1	4:G:296:ASP:OD1	2.26	0.54
4:H:106:GLU:OE2	4:H:663:ARG:NH2	2.40	0.54
4:J:138:LEU:HD13	4:J:156:LYS:HZ1	1.72	0.54
4:J:654:LYS:CB	4:J:661:ILE:HD11	2.38	0.54
4:I:553:LEU:HD12	4:I:595:ILE:HG21	1.88	0.54
4:K:504:ILE:HG12	4:K:509:VAL:HG21	1.89	0.54
4:N:719:LEU:HD11	4:N:781:LEU:HD11	1.90	0.54
4:M:445:PRO:HB3	4:M:528:MET:HG3	1.88	0.54
3:A:54:SER:OG	3:A:55:LEU:N	2.41	0.54
4:F:313:ASP:OD1	4:F:313:ASP:N	2.40	0.54
4:I:313:ASP:N	4:I:313:ASP:OD1	2.41	0.54
4:K:568:ILE:HD13	4:K:572:MET:HE3	1.90	0.54
4:M:472:ASN:HB3	4:M:475:VAL:HG12	1.90	0.54
4:N:101:THR:HB	4:N:665:PRO:HG3	1.90	0.54
4:N:290:ASN:ND2	4:N:559:TYR:OH	2.41	0.54
4:I:427:ILE:HG23	4:J:534:ARG:HH12	1.72	0.53
4:E:272:LEU:HD22	4:E:305:TYR:HB3	1.90	0.53
4:I:274:ASN:HD21	4:I:298:ASN:HA	1.74	0.53
4:I:488:ARG:HH11	4:I:490:VAL:HG11	1.74	0.53
4:M:553:LEU:HD12	4:M:595:ILE:HG21	1.89	0.53
3:A:789:MET:SD	3:A:871:LYS:NZ	2.81	0.53
4:E:731:ALA:HB2	4:E:831:LYS:HD3	1.89	0.53
4:K:654:LYS:HB2	4:K:661:ILE:HD11	1.90	0.53
3:A:1038:LYS:NZ	4:G:367:ILE:O	2.40	0.53
4:F:281:ILE:HD13	4:F:293:LEU:HD13	1.89	0.53
4:G:150:ARG:HD2	4:G:152:ARG:HH12	1.73	0.53
4:G:314:ARG:NH1	4:G:626:ASN:OD1	2.42	0.53
4:I:449:MET:HE1	4:I:470:ILE:HA	1.91	0.53
4:E:309:ASN:ND2	4:E:623:GLU:OE2	2.42	0.53
4:E:479:LEU:O	4:E:483:ASN:ND2	2.41	0.53
4:G:627:ASP:OD1	4:G:685:ARG:NH2	2.42	0.53
4:I:874:ASN:OD1	4:I:882:ARG:NH1	2.41	0.53
4:K:354:GLN:HG3	4:M:75:GLU:HG3	1.90	0.53
4:I:149:LEU:C	4:I:149:LEU:HD12	2.34	0.53
4:L:264:LEU:HD13	4:L:849:LEU:HD12	1.91	0.53
4:L:819:VAL:HA	4:L:824:TRP:HE1	1.74	0.53
3:A:265:GLU:OE2	3:A:488:ARG:NH2	2.42	0.53
3:A:340:LEU:HA	3:A:575:ILE:HD11	1.90	0.53
4:H:732:ARG:HH11	4:H:836:ILE:HD11	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:701:ARG:NH1	4:G:833:TYR:OH	2.42	0.53
4:G:703:SER:OG	4:G:704:ASP:N	2.42	0.53
4:I:244:ASN:HB2	4:I:853:LEU:HB2	1.90	0.53
4:I:489:GLN:HB3	4:I:496:LEU:HD11	1.91	0.53
4:J:851:SER:OG	4:J:852:ASN:N	2.42	0.53
4:I:399:MET:HE2	4:I:425:MET:HE2	1.91	0.52
4:J:748:THR:HB	4:K:230:ARG:HG3	1.91	0.52
3:A:263:THR:O	3:A:508:ARG:NH1	2.40	0.52
4:E:392:ALA:HB3	4:E:399:MET:HE2	1.90	0.52
4:F:108:LYS:NZ	4:F:581:GLU:OE2	2.42	0.52
4:J:355:LYS:NZ	4:J:359:ASP:OD2	2.43	0.52
4:K:327:THR:HA	4:K:330:THR:HG22	1.90	0.52
4:K:442:ILE:HD13	4:K:536:ILE:HD11	1.91	0.52
4:L:706:ILE:O	4:L:834:LYS:NZ	2.42	0.52
3:A:786:ARG:NH2	3:A:876:ASP:OD2	2.41	0.52
4:G:277:ILE:HD11	4:G:299:LEU:HD11	1.91	0.52
4:G:555:ARG:NH1	4:G:879:ASP:OD1	2.42	0.52
4:I:132:ILE:HG23	4:I:133:PHE:HD1	1.74	0.52
4:H:330:THR:OG1	4:H:397:ARG:NH2	2.42	0.52
4:E:645:MET:HE3	4:E:673:ARG:HD2	1.91	0.52
4:G:731:ALA:HB2	4:G:831:LYS:HD3	1.91	0.52
4:I:627:ASP:OD1	4:I:685:ARG:NH1	2.40	0.52
4:H:472:ASN:HB2	4:H:475:VAL:HG22	1.90	0.52
4:H:851:SER:OG	4:H:852:ASN:N	2.43	0.52
4:E:327:THR:HA	4:E:330:THR:HG22	1.90	0.52
4:H:140:VAL:HG21	4:H:152:ARG:HH12	1.75	0.52
4:K:397:ARG:NH2	4:K:579:THR:OG1	2.42	0.52
4:L:170:ARG:NH1	4:L:639:ASN:O	2.43	0.52
4:E:756:ASN:O	4:E:760:ASN:ND2	2.42	0.52
4:I:415:MET:HG2	4:I:438:ILE:HD13	1.91	0.52
3:A:1004:TYR:OH	3:A:1043:ASN:OD1	2.24	0.52
4:E:339:VAL:HG11	4:E:610:TYR:HB2	1.91	0.52
4:H:393:MET:HA	4:H:425:MET:HE3	1.91	0.52
4:H:443:ILE:HG12	4:H:520:LEU:HD11	1.92	0.52
4:M:660:ASP:OD1	4:M:660:ASP:N	2.43	0.52
4:F:253:HIS:HD2	4:F:255:ILE:H	1.57	0.52
4:H:313:ASP:OD1	4:H:313:ASP:N	2.42	0.52
4:N:110:SER:OG	4:N:114:LYS:NZ	2.42	0.52
4:N:883:ILE:HG22	4:N:884:MET:HG3	1.92	0.52
4:F:89:HIS:HD2	4:F:91:LYS:H	1.58	0.51
4:G:315:LEU:HD13	4:G:672:LEU:HD11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:403:PHE:HB3	4:L:585:LEU:HB2	1.91	0.51
4:L:851:SER:OG	4:L:852:ASN:N	2.39	0.51
4:M:401:LEU:HD22	4:M:585:LEU:HD13	1.92	0.51
4:G:730:ILE:HD12	4:G:834:LYS:HB3	1.91	0.51
4:J:408:TYR:OH	4:J:535:SER:OG	2.28	0.51
4:K:314:ARG:NH1	4:K:626:ASN:OD1	2.42	0.51
4:K:742:LEU:HD21	4:K:766:LEU:HD23	1.92	0.51
4:H:284:ARG:HE	4:H:566:ALA:HB2	1.76	0.51
4:F:552:ASP:HB3	4:F:884:MET:HA	1.91	0.51
4:H:309:ASN:HD22	4:H:622:ASN:HB3	1.75	0.51
4:J:132:ILE:HG13	4:J:133:PHE:HD1	1.74	0.51
4:E:268:LEU:HD21	4:E:684:ARG:HD2	1.92	0.51
4:F:750:ASP:OD2	4:G:230:ARG:NH2	2.39	0.51
4:I:253:HIS:HD2	4:I:255:ILE:H	1.58	0.51
4:I:453:HIS:NE2	4:K:886:GLU:OE2	2.43	0.51
4:K:445:PRO:HB3	4:K:528:MET:HG3	1.93	0.51
3:A:268:GLU:HG2	3:A:919:TYR:HB3	1.92	0.51
3:A:507:THR:HG23	3:A:510:LEU:HD12	1.92	0.51
4:I:531:ASP:OD2	4:K:538:ARG:NH2	2.43	0.51
4:H:517:LEU:HG	4:H:540:ILE:HG23	1.92	0.51
4:H:553:LEU:HD12	4:H:595:ILE:HG21	1.91	0.51
4:L:450:GLN:OE1	4:M:880:ASN:ND2	2.42	0.51
4:L:487:PHE:HB3	4:L:498:GLN:HB3	1.93	0.51
3:A:73:SER:OG	3:A:74:ILE:N	2.44	0.51
4:F:748:THR:HB	4:G:230:ARG:HG3	1.93	0.51
4:G:660:ASP:N	4:G:660:ASP:OD1	2.44	0.51
4:J:131:ARG:HH22	4:J:211:THR:HB	1.76	0.51
4:K:322:GLU:HB3	4:K:578:LEU:CD1	2.40	0.51
4:N:225:GLU:OE2	4:N:843:ARG:NH1	2.43	0.51
4:I:126:GLN:HE21	4:I:188:LEU:HD11	1.76	0.51
4:L:541:LEU:O	4:L:545:ASN:ND2	2.44	0.51
1:B:587:A:H62	4:G:358:GLN:HE21	1.58	0.51
4:E:427:ILE:HD11	4:F:534:ARG:HH12	1.75	0.51
3:A:169:LEU:HD22	3:A:230:ILE:HD12	1.93	0.51
3:A:262:ILE:HD11	3:A:508:ARG:HG3	1.93	0.51
4:E:235:MET:HE2	4:E:846:MET:HG3	1.93	0.51
4:E:366:THR:HG23	4:E:367:ILE:HG13	1.92	0.51
3:A:22:VAL:HG12	3:A:77:TYR:HB2	1.93	0.50
4:G:146:GLU:HA	4:G:811:SER:HB2	1.93	0.50
4:K:130:PHE:H	4:K:260:ASN:HD21	1.58	0.50
4:I:300:PRO:HG3	4:H:578:LEU:HD11	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:385:CYS:SG	4:H:587:SER:OG	2.62	0.50
4:J:479:LEU:O	4:J:483:ASN:ND2	2.44	0.50
4:F:253:HIS:CD2	4:F:255:ILE:H	2.29	0.50
4:K:321:PHE:CD1	4:K:324:LEU:HD12	2.46	0.50
4:K:472:ASN:HB3	4:K:475:VAL:HG12	1.93	0.50
3:A:614:ARG:O	3:A:617:ASN:ND2	2.45	0.50
4:H:282:PRO:HG2	4:H:285:ILE:HG12	1.94	0.50
4:H:742:LEU:HD21	4:H:766:LEU:HD23	1.93	0.50
4:J:727:TYR:HB2	4:J:809:ILE:HB	1.92	0.50
4:K:277:ILE:HG12	4:K:861:LEU:HB3	1.92	0.50
3:A:1025:LYS:HE3	4:I:363:GLU:OE1	2.12	0.50
4:H:131:ARG:NH2	4:H:211:THR:OG1	2.45	0.50
4:F:701:ARG:HD2	4:F:835:GLN:HG3	1.93	0.50
4:G:531:ASP:OD2	4:I:361:GLN:NE2	2.45	0.50
4:I:112:LEU:HD12	4:I:334:ILE:HG22	1.94	0.50
4:I:327:THR:HA	4:I:330:THR:HG22	1.93	0.50
4:J:582:LYS:NZ	4:K:296:ASP:OD2	2.44	0.50
4:M:851:SER:OG	4:M:852:ASN:N	2.44	0.50
4:N:851:SER:OG	4:N:852:ASN:N	2.44	0.50
3:A:605:LEU:HB2	3:A:628:VAL:HG11	1.93	0.50
3:A:687:ILE:HG23	3:A:900:MET:HG2	1.94	0.50
4:H:283:GLU:OE2	4:H:286:ARG:NH2	2.45	0.50
4:K:141:TYR:CZ	4:K:810:ASN:ND2	2.80	0.50
4:M:552:ASP:HA	4:M:555:ARG:HG3	1.92	0.50
4:H:337:ARG:NH2	4:H:384:ASP:OD1	2.41	0.50
4:H:734:LEU:HD12	4:H:765:ALA:HB1	1.94	0.50
4:M:682:GLU:OE1	4:M:685:ARG:NH1	2.44	0.50
4:G:424:ASP:HB3	4:G:574:HIS:HE1	1.77	0.50
4:I:314:ARG:NH1	4:I:626:ASN:OD1	2.45	0.50
4:J:498:GLN:O	4:J:561:TYR:OH	2.30	0.50
4:K:270:GLU:OE2	4:K:307:ARG:NH2	2.44	0.50
4:K:320:ASN:ND2	4:K:671:ARG:HH11	2.10	0.50
4:M:399:MET:HE2	4:M:425:MET:HE2	1.94	0.50
4:F:137:GLN:NE2	4:F:151:ASN:O	2.45	0.49
4:H:343:LEU:HD23	4:H:346:LEU:HD21	1.94	0.49
4:K:770:LEU:HD21	4:K:802:LEU:HD22	1.93	0.49
4:E:851:SER:OG	4:E:852:ASN:N	2.45	0.49
4:G:264:LEU:HD22	4:G:851:SER:HB2	1.94	0.49
4:M:568:ILE:HD13	4:M:572:MET:HE3	1.94	0.49
3:A:209:SER:OG	3:A:697:ASN:ND2	2.46	0.49
4:F:455:ARG:NH2	4:F:457:GLY:O	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:K:726:GLY:HA3	4:K:810:ASN:OD1	2.11	0.49
4:M:160:ASP:OD1	4:M:160:ASP:N	2.44	0.49
3:A:238:SER:HB2	3:A:335:ARG:HH11	1.78	0.49
4:F:311:LEU:HD11	4:F:314:ARG:HE	1.76	0.49
4:G:744:GLU:HG2	4:G:747:ARG:HH21	1.76	0.49
4:I:626:ASN:HA	4:I:629:VAL:HG12	1.95	0.49
4:M:315:LEU:HD13	4:M:672:LEU:HD11	1.95	0.49
3:A:85:VAL:HG22	3:A:143:ASN:HD21	1.78	0.49
4:E:790:PHE:HE2	4:E:830:THR:HG21	1.78	0.49
4:F:140:VAL:HG21	4:F:152:ARG:HE	1.77	0.49
4:G:877:ALA:HB2	4:G:883:ILE:HG23	1.94	0.49
4:K:472:ASN:ND2	4:K:519:GLN:OE1	2.35	0.49
4:L:340:VAL:HG13	4:L:387:LYS:HD3	1.95	0.49
4:M:235:MET:HE2	4:M:846:MET:HG3	1.93	0.49
4:E:860:ASP:O	4:N:94:GLN:NE2	2.45	0.49
4:L:290:ASN:HD21	4:L:876:VAL:H	1.59	0.49
4:L:489:GLN:HB3	4:L:496:LEU:HD11	1.95	0.49
4:N:479:LEU:O	4:N:483:ASN:ND2	2.45	0.49
3:A:827:VAL:O	3:A:831:ILE:N	2.42	0.49
4:G:797:ARG:HE	4:H:294:ASN:HD22	1.61	0.49
4:I:567:CYS:SG	4:I:611:TYR:HB2	2.52	0.49
4:F:310:LEU:HD11	4:F:615:VAL:HG13	1.95	0.49
4:F:556:LEU:HD22	4:F:884:MET:HE3	1.95	0.49
4:I:472:ASN:HB3	4:I:475:VAL:HG12	1.94	0.49
4:L:525:PHE:H	4:L:533:LYS:HD2	1.77	0.49
4:N:876:VAL:HA	4:N:882:ARG:HA	1.93	0.49
4:E:137:GLN:HE21	4:E:151:ASN:HD22	1.60	0.48
4:F:441:THR:HG23	4:F:442:ILE:HG13	1.94	0.48
4:G:313:ASP:N	4:G:313:ASP:OD1	2.45	0.48
4:L:408:TYR:OH	4:L:535:SER:OG	2.31	0.48
3:A:608:ILE:HA	3:A:611:VAL:HG12	1.94	0.48
4:I:472:ASN:HD21	4:I:519:GLN:HG2	1.78	0.48
4:H:99:GLN:O	4:H:663:ARG:NH1	2.46	0.48
4:J:253:HIS:HB3	4:J:256:ASP:HB2	1.94	0.48
4:L:166:ASP:HB3	4:L:743:GLU:HG2	1.95	0.48
3:A:368:ARG:HG3	3:A:540:MET:HE2	1.95	0.48
4:I:148:GLU:C	4:I:149:LEU:HG	2.29	0.48
4:M:357:SER:O	4:M:357:SER:OG	2.30	0.48
3:A:233:SER:OG	3:A:237:ARG:NH1	2.46	0.48
4:I:514:MET:HG3	4:I:547:LEU:HD13	1.94	0.48
4:H:514:MET:HE3	4:H:547:LEU:HA	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:J:441:THR:HG23	4:J:442:ILE:HG13	1.95	0.48
3:A:529:HIS:HD1	3:A:664:THR:HG1	1.59	0.48
3:A:622:ALA:HB2	4:F:404:VAL:HG21	1.94	0.48
4:E:435:GLN:OE1	4:E:461:THR:OG1	2.29	0.48
4:J:397:ARG:NH2	4:J:581:GLU:OE1	2.47	0.48
4:N:146:GLU:HA	4:N:811:SER:HB2	1.94	0.48
3:A:396:SER:O	3:A:396:SER:OG	2.32	0.48
4:F:851:SER:OG	4:F:852:ASN:N	2.47	0.48
4:I:673:ARG:NH2	4:J:887:LEU:OXT	2.46	0.48
4:J:238:ARG:HE	4:J:240:GLN:HE21	1.62	0.48
4:L:356:MET:HE3	4:L:409:MET:HG2	1.95	0.48
4:N:727:TYR:HB2	4:N:809:ILE:HB	1.95	0.48
4:I:427:ILE:HG13	4:I:580:THR:HG21	1.96	0.48
4:L:641:TYR:OH	4:L:743:GLU:OE1	2.32	0.48
4:I:504:ILE:HG13	4:I:509:VAL:HG21	1.95	0.48
4:I:851:SER:OG	4:I:852:ASN:N	2.45	0.48
4:K:718:GLN:NE2	4:K:719:LEU:O	2.47	0.48
4:E:202:ASP:HA	4:N:642:GLN:HG3	1.96	0.48
4:H:377:ILE:HA	4:H:590:SER:HB3	1.96	0.48
4:M:253:HIS:HD2	4:M:255:ILE:H	1.61	0.48
4:N:129:LEU:HD23	4:N:252:LEU:HD21	1.96	0.48
4:I:320:ASN:HD21	4:J:545:ASN:HD21	1.61	0.47
4:H:408:TYR:OH	4:H:535:SER:OG	2.31	0.47
4:J:655:ARG:O	4:J:657:GLN:NE2	2.47	0.47
4:K:449:MET:HE1	4:K:470:ILE:HA	1.95	0.47
4:K:569:THR:HG22	4:K:614:ASN:HB3	1.95	0.47
4:I:277:ILE:HD11	4:I:299:LEU:HD11	1.96	0.47
3:A:140:SER:OG	3:A:738:GLN:OE1	2.32	0.47
4:K:443:ILE:HG12	4:K:520:LEU:HD11	1.96	0.47
4:M:342:ASP:HB2	4:M:383:ASN:HD21	1.79	0.47
4:M:882:ARG:HG3	4:M:885:ASN:HB3	1.96	0.47
3:A:628:VAL:HG23	3:A:633:ASN:HB3	1.96	0.47
3:A:874:ILE:HD13	3:A:996:LEU:HG	1.96	0.47
4:K:627:ASP:OD1	4:K:685:ARG:NH2	2.42	0.47
4:N:541:LEU:O	4:N:545:ASN:ND2	2.47	0.47
4:K:202:ASP:OD1	4:K:202:ASP:N	2.38	0.47
4:F:417:LEU:HD22	4:F:588:VAL:HG21	1.97	0.47
4:L:152:ARG:NH1	4:L:723:GLU:OE1	2.47	0.47
3:A:335:ARG:O	3:A:459:THR:OG1	2.33	0.47
3:A:1079:PRO:HB3	4:M:92:GLU:HG3	1.96	0.47
4:E:472:ASN:HD21	4:E:519:GLN:HG2	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:721:ARG:NH1	4:F:727:TYR:OH	2.48	0.47
4:L:393:MET:O	4:L:396:GLN:NE2	2.48	0.47
4:L:719:LEU:HD13	4:L:826:PRO:HB2	1.97	0.47
4:M:435:GLN:OE1	4:M:461:THR:OG1	2.28	0.47
4:N:330:THR:OG1	4:N:397:ARG:NH2	2.48	0.47
3:A:826:ILE:HD11	3:A:949:LYS:HD2	1.96	0.47
3:A:1026:ILE:CD1	3:A:1063:MET:HB3	2.45	0.47
4:E:289:VAL:HG13	4:N:440:ASN:HD21	1.79	0.47
4:E:627:ASP:OD1	4:E:685:ARG:NH2	2.48	0.47
4:F:130:PHE:H	4:F:260:ASN:HD21	1.63	0.47
4:G:427:ILE:HD12	4:G:580:THR:HG21	1.96	0.47
4:I:786:ASP:OD1	4:I:786:ASP:N	2.45	0.47
4:H:818:LEU:O	4:H:824:TRP:NE1	2.47	0.47
4:J:879:ASP:OD1	4:J:879:ASP:N	2.47	0.47
4:K:272:LEU:HD22	4:K:305:TYR:HB3	1.97	0.47
4:M:593:MET:HE2	4:M:593:MET:HB3	1.70	0.47
4:N:593:MET:HE2	4:N:593:MET:HB3	1.77	0.47
3:A:141:ASN:ND2	3:A:208:ASN:OD1	2.47	0.47
4:G:719:LEU:HD23	4:G:828:SER:HB3	1.96	0.47
4:I:725:TYR:O	4:I:810:ASN:ND2	2.47	0.47
4:K:149:LEU:O	4:K:149:LEU:HG	2.15	0.47
4:K:735:ASP:OD1	4:K:735:ASP:N	2.48	0.47
4:M:311:LEU:O	4:M:622:ASN:ND2	2.44	0.47
4:N:337:ARG:NH1	4:N:388:THR:OG1	2.46	0.47
4:I:806:LEU:HD12	4:I:832:VAL:HG21	1.97	0.47
4:J:503:ASN:HB3	4:J:509:VAL:HG23	1.97	0.47
4:L:740:ILE:HB	4:L:766:LEU:HD23	1.96	0.47
4:K:274:ASN:HD21	4:K:298:ASN:HA	1.80	0.46
4:L:621:TYR:HD1	4:L:656:LEU:HD21	1.80	0.46
4:L:755:THR:HG22	4:L:794:VAL:HG12	1.97	0.46
4:N:461:THR:HB	4:N:464:GLN:HG3	1.97	0.46
4:E:325:TRP:HA	4:E:328:ILE:HG22	1.97	0.46
4:K:253:HIS:CD2	4:K:255:ILE:H	2.32	0.46
4:M:427:ILE:HD12	4:M:580:THR:HG21	1.97	0.46
3:A:1044:HIS:CE1	3:A:1048:ASN:HD22	2.33	0.46
4:E:388:THR:HG23	4:E:583:LEU:HD12	1.97	0.46
4:H:113:LYS:HE3	4:H:117:ASP:HB2	1.97	0.46
4:L:552:ASP:HB3	4:L:884:MET:HA	1.97	0.46
3:A:545:LEU:HB3	3:A:558:LEU:HD21	1.97	0.46
4:E:874:ASN:OD1	4:E:874:ASN:N	2.49	0.46
4:G:388:THR:HG23	4:G:583:LEU:HD12	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:I:253:HIS:CD2	4:I:255:ILE:H	2.33	0.46
4:I:310:LEU:HD21	4:I:615:VAL:HG13	1.97	0.46
4:J:553:LEU:HD12	4:J:595:ILE:HG21	1.96	0.46
3:A:113:LYS:HB2	3:A:118:LEU:HD12	1.97	0.46
4:E:352:GLN:HA	4:E:355:LYS:HB2	1.96	0.46
4:I:660:ASP:N	4:I:660:ASP:OD1	2.48	0.46
4:H:497:ASN:OD1	4:H:497:ASN:N	2.47	0.46
4:J:102:ILE:HG22	4:J:104:THR:H	1.79	0.46
4:J:567:CYS:SG	4:J:607:LEU:HD11	2.55	0.46
4:L:742:LEU:HD22	4:L:799:VAL:HG21	1.97	0.46
4:E:423:ASN:HD21	4:E:482:VAL:HG12	1.81	0.46
4:H:882:ARG:HH11	4:H:887:LEU:HD11	1.80	0.46
4:J:556:LEU:HD22	4:J:884:MET:HE3	1.97	0.46
4:N:283:GLU:OE2	4:N:286:ARG:NH2	2.48	0.46
3:A:870:SER:O	3:A:870:SER:OG	2.31	0.46
4:I:115:LEU:HD21	4:I:331:SER:HB3	1.97	0.46
4:I:230:ARG:HG3	4:H:748:THR:HB	1.98	0.46
4:I:665:PRO:HG2	4:J:355:LYS:HD2	1.97	0.46
4:H:489:GLN:NE2	4:H:498:GLN:OE1	2.42	0.46
4:H:264:LEU:O	4:H:267:GLN:NE2	2.49	0.46
3:A:539:ILE:HG23	3:A:562:LYS:HG3	1.98	0.46
4:F:393:MET:HE2	4:F:421:VAL:HG13	1.97	0.46
4:I:323:SER:HB3	4:I:668:GLN:HE22	1.81	0.46
4:J:641:TYR:HB2	4:K:202:ASP:HB3	1.98	0.46
4:K:757:MET:HE2	4:K:764:VAL:HG22	1.98	0.46
4:L:91:LYS:HD3	4:L:95:TYR:HE2	1.81	0.46
4:M:446:ALA:CB	4:M:520:LEU:CD1	2.94	0.46
3:A:91:LYS:NZ	3:A:150:GLU:OE2	2.49	0.46
4:K:665:PRO:HG3	4:L:355:LYS:HD2	1.97	0.46
4:M:556:LEU:HD22	4:M:884:MET:HE3	1.98	0.46
3:A:1014:LEU:HD11	3:A:1042:ILE:HD11	1.98	0.45
4:F:184:MET:HE3	4:F:684:ARG:HH21	1.81	0.45
4:F:742:LEU:HD21	4:F:766:LEU:HD22	1.97	0.45
4:G:399:MET:HE3	4:G:425:MET:HE2	1.97	0.45
4:I:121:GLU:OE1	4:I:624:ARG:NH1	2.50	0.45
4:N:135:PRO:HB2	4:N:153:TRP:HE3	1.81	0.45
3:A:387:GLN:NE2	3:A:551:ASP:OD2	2.47	0.45
4:I:230:ARG:NH2	4:H:750:ASP:OD2	2.46	0.45
4:H:475:VAL:HG12	4:H:516:ALA:HB1	1.98	0.45
4:L:498:GLN:O	4:L:561:TYR:OH	2.32	0.45
4:G:273:ASN:OD1	4:G:273:ASN:N	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:423:ASN:ND2	4:G:482:VAL:O	2.49	0.45
4:I:118:ILE:HG12	4:I:657:GLN:HG3	1.97	0.45
4:J:510:VAL:HG22	4:J:514:MET:HE2	1.98	0.45
4:L:253:HIS:CD2	4:L:255:ILE:H	2.31	0.45
4:M:149:LEU:O	4:M:149:LEU:HG	2.16	0.45
4:N:264:LEU:HD11	4:N:851:SER:HB2	1.98	0.45
4:G:130:PHE:N	4:G:260:ASN:OD1	2.47	0.45
4:I:325:TRP:HA	4:I:328:ILE:HG22	1.99	0.45
4:H:667:ASP:OD1	4:H:667:ASP:N	2.46	0.45
4:K:710:VAL:HG11	4:K:790:PHE:HZ	1.81	0.45
4:N:340:VAL:HG11	4:N:390:ILE:HB	1.98	0.45
4:H:394:LEU:HG	4:H:568:ILE:HD11	1.98	0.45
4:J:403:PHE:HB3	4:J:585:LEU:HB2	1.97	0.45
4:J:472:ASN:HD21	4:J:519:GLN:HG2	1.81	0.45
4:K:376:GLY:O	4:K:590:SER:OG	2.35	0.45
4:N:110:SER:HG	4:N:114:LYS:NZ	2.15	0.45
4:E:273:ASN:H	4:E:276:ILE:HG22	1.81	0.45
4:I:253:HIS:O	4:I:257:HIS:N	2.49	0.45
4:L:167:TYR:OH	4:L:641:TYR:O	2.27	0.45
4:F:393:MET:O	4:F:396:GLN:NE2	2.49	0.45
4:F:514:MET:HB2	4:F:547:LEU:HD13	1.99	0.45
4:G:325:TRP:HA	4:G:328:ILE:HG22	1.98	0.45
4:K:166:ASP:HB3	4:K:743:GLU:HB2	1.97	0.45
4:L:281:ILE:HD11	4:L:865:VAL:HG11	1.99	0.45
4:L:622:ASN:HD22	4:L:625:ILE:HD11	1.80	0.45
3:A:618:LYS:O	4:F:584:GLN:NE2	2.50	0.45
4:I:342:ASP:OD1	4:I:342:ASP:N	2.45	0.45
4:J:580:THR:OG1	4:K:296:ASP:OD1	2.33	0.45
4:K:730:ILE:HG13	4:K:834:LYS:HB3	1.98	0.45
4:M:488:ARG:HH11	4:M:490:VAL:HG11	1.82	0.45
4:E:134:GLU:HB2	4:E:158:LYS:HG3	1.98	0.45
4:E:810:ASN:OD1	4:E:810:ASN:N	2.48	0.45
4:I:877:ALA:HB2	4:I:883:ILE:HG23	1.98	0.45
4:K:701:ARG:HD3	4:K:835:GLN:HG3	1.98	0.45
4:M:419:THR:HB	4:M:478:TRP:HE1	1.82	0.45
4:N:284:ARG:HD2	4:N:566:ALA:HB2	1.97	0.45
4:F:129:LEU:HD13	4:F:252:LEU:HD11	1.99	0.44
4:I:393:MET:HE2	4:I:421:VAL:HG13	1.98	0.44
4:H:731:ALA:HB2	4:H:831:LYS:HD2	1.99	0.44
4:K:325:TRP:HA	4:K:328:ILE:HG22	1.99	0.44
4:N:517:LEU:HG	4:N:540:ILE:HG23	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:443:ILE:HG12	4:E:520:LEU:HD11	1.99	0.44
4:J:375:THR:OG1	4:J:376:GLY:N	2.50	0.44
4:K:650:GLU:HG3	4:K:661:ILE:HG21	1.99	0.44
4:K:814:ASN:OD1	4:K:814:ASN:N	2.51	0.44
4:N:810:ASN:OD1	4:N:810:ASN:N	2.47	0.44
4:F:712:ILE:HA	4:F:830:THR:HG22	1.98	0.44
4:G:112:LEU:HD11	4:G:338:SER:HB3	2.00	0.44
4:G:202:ASP:OD2	4:G:236:ARG:NE	2.50	0.44
4:G:393:MET:HE2	4:G:421:VAL:HG13	1.99	0.44
4:G:851:SER:OG	4:G:852:ASN:N	2.51	0.44
4:H:253:HIS:HB3	4:H:256:ASP:HB2	1.98	0.44
4:J:564:LEU:HD23	4:J:607:LEU:CD1	2.47	0.44
4:N:115:LEU:HD23	4:N:115:LEU:HA	1.84	0.44
4:E:489:GLN:HB3	4:E:496:LEU:HD11	1.99	0.44
4:E:667:ASP:HA	4:F:549:GLN:HE22	1.82	0.44
4:F:353:ILE:HG12	4:F:374:LEU:HD23	1.99	0.44
4:L:541:LEU:HA	4:L:544:SER:HB3	2.00	0.44
3:A:623:THR:HG21	4:F:373:PHE:HE1	1.83	0.44
3:A:900:MET:O	3:A:908:GLN:NE2	2.49	0.44
4:E:472:ASN:HB3	4:E:475:VAL:HG12	1.98	0.44
4:F:435:GLN:OE1	4:F:461:THR:OG1	2.36	0.44
4:G:284:ARG:NH1	4:G:497:ASN:HD21	2.15	0.44
4:I:284:ARG:HE	4:I:566:ALA:HB2	1.82	0.44
4:N:498:GLN:O	4:N:561:TYR:OH	2.32	0.44
4:I:135:PRO:HB2	4:I:153:TRP:HE3	1.83	0.44
4:I:475:VAL:HG21	4:I:520:LEU:HD12	1.99	0.44
4:H:138:LEU:HD11	4:H:156:LYS:HG3	2.00	0.44
4:K:731:ALA:HB2	4:K:831:LYS:HE2	2.00	0.44
4:N:293:LEU:HD11	4:N:608:PHE:HZ	1.82	0.44
4:E:393:MET:HE2	4:E:421:VAL:HG13	2.00	0.44
4:G:662:SER:O	4:G:662:SER:OG	2.36	0.44
4:I:707:ALA:N	4:I:770:LEU:O	2.50	0.44
4:H:819:VAL:HA	4:H:824:TRP:HE1	1.83	0.44
4:J:417:LEU:HD22	4:J:588:VAL:HG21	2.00	0.44
4:K:390:ILE:HD12	4:K:561:TYR:HB2	2.00	0.44
3:A:705:THR:OG1	3:A:706:GLN:N	2.50	0.44
4:H:142:ARG:NH2	4:H:148:GLU:OE1	2.50	0.44
4:H:231:PHE:HZ	4:H:846:MET:HE2	1.82	0.44
4:J:717:MET:HE3	4:J:831:LYS:HG3	1.99	0.44
4:L:660:ASP:OD1	4:L:660:ASP:N	2.51	0.44
4:M:446:ALA:CB	4:M:520:LEU:HD11	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:450:GLY:O	3:A:462:ILE:N	2.50	0.44
3:A:498:ASN:HD22	3:A:917:ARG:HB3	1.83	0.44
4:E:253:HIS:CD2	4:E:255:ILE:H	2.36	0.44
4:E:441:THR:HG22	4:E:442:ILE:HG13	2.00	0.44
4:K:419:THR:HB	4:K:478:TRP:HE1	1.83	0.44
4:K:667:ASP:OD2	4:L:546:ARG:NH1	2.50	0.44
4:M:701:ARG:HA	4:M:708:GLN:HE22	1.83	0.44
4:F:399:MET:HE3	4:F:399:MET:HB2	1.76	0.43
4:G:517:LEU:HD13	4:G:540:ILE:HG12	2.00	0.43
4:J:335:LEU:HD11	4:J:613:VAL:HG11	2.00	0.43
4:K:337:ARG:NH2	4:K:388:THR:OG1	2.51	0.43
4:N:297:ARG:HD3	4:N:864:PHE:HD1	1.83	0.43
3:A:847:LEU:HD23	3:A:847:LEU:HA	1.89	0.43
3:A:1026:ILE:CG2	3:A:1067:TRP:HB2	2.47	0.43
4:I:160:ASP:N	4:I:160:ASP:OD1	2.46	0.43
4:I:297:ARG:HD3	4:H:98:LEU:HD22	2.00	0.43
4:I:719:LEU:HD22	4:I:729:ASN:HB3	2.01	0.43
4:J:170:ARG:NH1	4:J:638:LEU:O	2.50	0.43
4:L:629:VAL:HG12	4:L:681:VAL:HG12	1.99	0.43
3:A:185:HIS:H	3:A:185:HIS:CD2	2.36	0.43
3:A:232:LEU:HD23	3:A:232:LEU:HA	1.84	0.43
4:F:582:LYS:NZ	4:G:296:ASP:OD2	2.39	0.43
4:I:426:PHE:HB2	4:I:431:LEU:HB2	2.01	0.43
4:L:670:TYR:HA	4:L:673:ARG:HG2	2.00	0.43
4:E:141:TYR:OH	4:E:812:ASP:OD2	2.34	0.43
4:E:313:ASP:OD1	4:E:313:ASP:N	2.50	0.43
4:F:200:SER:OG	4:F:202:ASP:O	2.28	0.43
4:H:160:ASP:OD1	4:H:160:ASP:N	2.47	0.43
4:H:806:LEU:HD12	4:H:832:VAL:HG11	1.99	0.43
4:N:342:ASP:N	4:N:342:ASP:OD1	2.41	0.43
4:E:137:GLN:NE2	4:E:223:GLU:OE2	2.43	0.43
4:E:491:VAL:HA	4:E:496:LEU:HA	2.01	0.43
4:F:112:LEU:HD11	4:F:338:SER:HB3	2.00	0.43
4:M:253:HIS:HB3	4:M:256:ASP:HB2	2.00	0.43
3:A:184:ARG:NH1	6:A:1102:GTP:O1G	2.51	0.43
4:G:160:ASP:OD1	4:G:160:ASP:N	2.43	0.43
4:I:185:PRO:HG3	4:I:188:LEU:HD12	2.00	0.43
4:I:637:ARG:HD2	4:I:689:PHE:HE2	1.84	0.43
4:J:873:ILE:HD12	4:J:873:ILE:HA	1.92	0.43
4:L:672:LEU:HD12	4:L:672:LEU:HA	1.84	0.43
3:A:6:LEU:HD11	4:G:84:LYS:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:735:LYS:HD2	3:A:735:LYS:HA	1.86	0.43
4:E:786:ASP:N	4:E:786:ASP:OD1	2.49	0.43
4:I:118:ILE:HG23	4:I:620:ASN:HD22	1.84	0.43
4:I:773:ILE:HG23	4:I:803:LYS:HD2	2.00	0.43
4:J:654:LYS:HB2	4:J:661:ILE:CD1	2.49	0.43
4:J:670:TYR:HA	4:J:673:ARG:HG2	2.01	0.43
4:L:708:GLN:H	4:L:708:GLN:HG2	1.51	0.43
4:M:318:HIS:CE1	4:M:573:GLN:HG3	2.54	0.43
3:A:359:MET:HE3	3:A:359:MET:HB2	1.96	0.43
4:G:323:SER:HB3	4:G:668:GLN:HE22	1.84	0.43
4:H:113:LYS:HA	4:H:113:LYS:HD2	1.91	0.43
4:H:498:GLN:O	4:H:561:TYR:OH	2.35	0.43
4:H:745:LEU:HD21	4:H:754:ILE:HG21	1.99	0.43
4:L:339:VAL:HG11	4:L:610:TYR:HB2	2.00	0.43
4:L:497:ASN:OD1	4:L:497:ASN:N	2.52	0.43
4:L:593:MET:HE2	4:L:593:MET:HB3	1.76	0.43
1:B:588:U:H4'	3:A:932:ARG:HH22	1.83	0.43
4:I:195:VAL:HG21	4:I:209:SER:HA	2.01	0.43
4:I:759:LEU:HD21	4:I:794:VAL:HG11	2.01	0.43
4:H:253:HIS:CD2	4:H:255:ILE:H	2.35	0.43
4:M:135:PRO:HB2	4:M:153:TRP:HE3	1.84	0.43
4:M:681:VAL:HG13	4:M:686:LEU:HB2	2.01	0.43
3:A:648:VAL:HG13	3:A:670:VAL:HG23	2.01	0.43
3:A:793:SER:O	3:A:793:SER:OG	2.36	0.43
3:A:954:HIS:CD2	3:A:956:ASN:H	2.36	0.43
4:G:553:LEU:HD21	4:G:591:LEU:HG	2.00	0.43
4:I:593:MET:HE2	4:I:593:MET:HB3	1.80	0.43
4:J:565:MET:HB3	4:J:565:MET:HE2	1.78	0.43
4:J:873:ILE:HD11	4:J:887:LEU:HD11	2.01	0.43
4:E:719:LEU:HD23	4:E:828:SER:HB3	2.01	0.42
4:H:481:PHE:HZ	4:H:501:ASN:HD21	1.67	0.42
4:K:228:VAL:HB	4:K:842:PHE:HD2	1.84	0.42
4:K:453:HIS:NE2	4:M:886:GLU:OE2	2.52	0.42
3:A:248:MET:HE3	3:A:681:TYR:HD2	1.83	0.42
3:A:519:THR:HG22	3:A:668:ALA:HA	2.01	0.42
4:E:296:ASP:OD1	4:N:582:LYS:NZ	2.45	0.42
4:F:862:LEU:HD23	4:F:862:LEU:HA	1.83	0.42
4:H:876:VAL:HA	4:H:882:ARG:HA	2.00	0.42
4:M:369:SER:O	4:M:369:SER:OG	2.31	0.42
4:F:644:LYS:HB3	4:F:644:LYS:HE2	1.84	0.42
4:F:733:ASN:ND2	4:F:735:ASP:OD2	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:645:MET:HB2	4:G:673:ARG:HD2	2.01	0.42
4:I:427:ILE:HG22	4:I:429:GLU:HG2	2.00	0.42
4:I:790:PHE:HE2	4:I:830:THR:HG21	1.83	0.42
4:K:453:HIS:HE1	4:M:551:VAL:HG11	1.85	0.42
4:K:703:SER:O	4:K:834:LYS:NZ	2.35	0.42
4:K:830:THR:OG1	4:K:831:LYS:N	2.53	0.42
4:L:343:LEU:HB3	4:L:346:LEU:HD21	2.01	0.42
4:M:570:MET:HG3	4:M:618:HIS:CD2	2.55	0.42
4:E:378:ASN:N	4:E:378:ASN:OD1	2.52	0.42
4:E:711:ILE:HA	4:E:765:ALA:HA	2.01	0.42
4:I:244:ASN:OD1	4:I:244:ASN:N	2.51	0.42
4:J:356:MET:HE3	4:J:356:MET:HB2	1.86	0.42
4:L:701:ARG:HD3	4:L:708:GLN:HE22	1.85	0.42
4:M:62:ILE:HG13	4:M:63:LYS:HD3	2.00	0.42
4:F:727:TYR:HB2	4:F:809:ILE:HB	2.01	0.42
4:N:715:ARG:HE	4:N:717:MET:HE3	1.85	0.42
4:N:733:ASN:OD1	4:N:733:ASN:N	2.53	0.42
3:A:254:LEU:O	4:F:86:LYS:NZ	2.50	0.42
3:A:680:ARG:HB3	3:A:687:ILE:HD11	2.01	0.42
4:G:312:GLN:HB3	4:G:571:ASN:HB2	2.01	0.42
4:I:730:ILE:HG23	4:I:834:LYS:HB3	2.02	0.42
4:L:733:ASN:OD1	4:L:733:ASN:N	2.52	0.42
4:N:732:ARG:HH22	4:N:836:ILE:HD11	1.85	0.42
3:A:644:THR:OG1	3:A:645:LYS:N	2.53	0.42
4:E:885:ASN:OD1	4:E:885:ASN:N	2.52	0.42
4:G:135:PRO:HB2	4:G:153:TRP:HE3	1.84	0.42
4:G:443:ILE:HG12	4:G:520:LEU:HD11	2.01	0.42
4:H:316:ASN:HD22	4:H:318:HIS:CE1	2.38	0.42
4:H:390:ILE:O	4:H:394:LEU:N	2.51	0.42
4:M:253:HIS:CD2	4:M:255:ILE:H	2.37	0.42
4:M:277:ILE:HG12	4:M:861:LEU:HD12	2.02	0.42
4:M:310:LEU:HD21	4:M:615:VAL:HG22	2.02	0.42
4:N:300:PRO:HD2	4:N:303:ALA:HB2	2.01	0.42
4:N:512:GLN:O	4:N:512:GLN:NE2	2.53	0.42
3:A:521:VAL:HA	3:A:666:VAL:HG23	2.01	0.42
4:J:449:MET:H	4:J:449:MET:HG2	1.64	0.42
4:K:593:MET:HE2	4:K:593:MET:HB3	1.72	0.42
4:K:773:ILE:HG23	4:K:803:LYS:HE2	2.02	0.42
4:K:801:THR:O	4:K:801:THR:OG1	2.37	0.42
4:M:621:TYR:HA	4:M:656:LEU:HD21	2.02	0.42
4:N:110:SER:OG	4:N:110:SER:O	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:549:THR:OG1	3:A:550:ASN:N	2.48	0.42
3:A:768:SER:OG	3:A:769:GLU:N	2.53	0.42
3:A:993:VAL:HG22	3:A:1070:MET:HE3	2.02	0.42
4:F:272:LEU:HD21	4:F:277:ILE:HD11	2.01	0.42
4:I:745:LEU:HD21	4:I:754:ILE:HG21	2.02	0.42
4:J:593:MET:HB3	4:J:593:MET:HE2	1.87	0.42
4:M:786:ASP:HB3	4:M:807:TYR:HE2	1.84	0.42
1:B:589:A:H61	3:A:482:ILE:HG23	1.85	0.42
3:A:401:SER:OG	3:A:402:ASN:N	2.52	0.42
4:E:132:ILE:HD12	4:E:132:ILE:HA	1.93	0.42
4:E:315:LEU:HD21	4:E:676:LEU:HD12	2.01	0.42
4:H:252:LEU:HB2	4:H:847:HIS:CE1	2.55	0.42
4:J:497:ASN:OD1	4:J:497:ASN:N	2.53	0.42
3:A:262:ILE:HG13	3:A:508:ARG:HA	2.02	0.41
4:E:691:LEU:HD23	4:E:691:LEU:HA	1.87	0.41
4:E:759:LEU:HD21	4:E:794:VAL:HG11	2.02	0.41
4:G:322:GLU:HB3	4:G:578:LEU:HD11	2.02	0.41
4:G:393:MET:O	4:G:396:GLN:NE2	2.53	0.41
4:G:439:ILE:HD13	4:G:439:ILE:HA	1.91	0.41
4:J:272:LEU:HD21	4:J:277:ILE:HD11	2.02	0.41
4:L:299:LEU:HA	4:L:300:PRO:HD3	1.94	0.41
4:M:157:LEU:HB3	4:M:160:ASP:HB3	2.02	0.41
3:A:102:LEU:O	3:A:111:ASN:ND2	2.42	0.41
3:A:633:ASN:OD1	3:A:633:ASN:N	2.53	0.41
4:F:324:LEU:HD11	4:F:653:LEU:HD21	2.02	0.41
4:G:727:TYR:HB2	4:G:809:ILE:HB	2.01	0.41
4:G:795:LYS:HG2	4:G:796:LEU:HD12	2.02	0.41
4:I:565:MET:HE2	4:I:565:MET:HB3	1.98	0.41
4:I:666:ASP:HB2	4:J:347:VAL:HG22	2.02	0.41
4:H:711:ILE:HA	4:H:765:ALA:HA	2.03	0.41
4:H:733:ASN:ND2	4:H:735:ASP:OD2	2.44	0.41
4:H:770:LEU:HD23	4:H:804:PRO:HD3	2.01	0.41
4:L:188:LEU:HD23	4:L:188:LEU:HA	1.93	0.41
4:L:232:ILE:HA	4:L:235:MET:HG3	2.01	0.41
4:M:569:THR:HG22	4:M:614:ASN:HB3	2.02	0.41
3:A:121:THR:OG1	3:A:122:ALA:N	2.53	0.41
4:E:517:LEU:HD12	4:E:517:LEU:HA	1.89	0.41
4:J:356:MET:HG3	4:J:542:LEU:HD13	2.02	0.41
4:L:393:MET:HA	4:L:425:MET:HE3	2.01	0.41
4:L:504:ILE:HD13	4:L:504:ILE:HG21	1.85	0.41
4:N:268:LEU:HD22	4:N:684:ARG:HH12	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:1022:PHE:CZ	3:A:1024:GLY:HA2	2.54	0.41
4:E:745:LEU:HD21	4:E:754:ILE:HG21	2.02	0.41
4:H:713:ALA:HA	4:H:763:PRO:HB3	2.00	0.41
4:J:134:GLU:HG2	4:J:158:LYS:HB3	2.01	0.41
4:J:170:ARG:NH1	4:J:639:ASN:O	2.41	0.41
4:M:268:LEU:HD21	4:M:684:ARG:HD2	2.02	0.41
4:N:729:ASN:HD22	4:N:831:LYS:HA	1.85	0.41
4:N:814:ASN:OD1	4:N:814:ASN:N	2.50	0.41
3:A:718:ILE:HD11	3:A:771:PHE:HA	2.01	0.41
4:F:389:LEU:HD11	4:F:421:VAL:HG21	2.02	0.41
4:K:315:LEU:HD13	4:K:672:LEU:HD11	2.02	0.41
4:L:389:LEU:HD11	4:L:421:VAL:HG21	2.03	0.41
4:L:578:LEU:HD23	4:L:578:LEU:HA	1.90	0.41
4:N:113:LYS:HA	4:N:113:LYS:HD2	1.85	0.41
3:A:407:GLN:NE2	3:A:414:THR:OG1	2.54	0.41
4:I:138:LEU:HD11	4:I:156:LYS:HE2	2.02	0.41
4:I:452:MET:HE1	4:I:462:PRO:HA	2.01	0.41
4:H:565:MET:HB3	4:H:565:MET:HE2	1.77	0.41
4:K:252:LEU:HD23	4:K:256:ASP:HB3	2.02	0.41
4:L:711:ILE:HG23	4:L:831:LYS:HG3	2.02	0.41
4:L:846:MET:HE3	4:L:846:MET:HB3	1.90	0.41
4:M:417:LEU:HG	4:M:588:VAL:HG11	2.03	0.41
3:A:368:ARG:HH22	4:E:663:ARG:HG3	1.86	0.41
4:F:422:PRO:HB3	4:F:486:GLN:HA	2.03	0.41
4:G:282:PRO:HD2	4:G:285:ILE:HD12	2.01	0.41
4:G:419:THR:HG21	4:G:513:LEU:HD22	2.02	0.41
4:H:299:LEU:HD13	4:H:299:LEU:HA	1.94	0.41
4:H:726:GLY:HA3	4:H:810:ASN:ND2	2.35	0.41
4:L:108:LYS:NZ	4:L:581:GLU:OE1	2.53	0.41
4:L:742:LEU:HD21	4:L:766:LEU:HD22	2.03	0.41
4:N:347:VAL:HG22	4:N:594:LEU:HA	2.03	0.41
3:A:560:LEU:HD23	3:A:560:LEU:HA	1.83	0.41
4:E:253:HIS:O	4:E:257:HIS:N	2.48	0.41
4:I:777:SER:HB3	4:I:780:SER:HB3	2.02	0.41
4:H:711:ILE:HG23	4:H:831:LYS:HG3	2.02	0.41
4:J:146:GLU:HA	4:J:811:SER:HB2	2.02	0.41
4:K:135:PRO:HB2	4:K:153:TRP:HE3	1.86	0.41
4:M:157:LEU:HD12	4:M:157:LEU:HA	1.94	0.41
3:A:457:ARG:NH1	3:A:458:ARG:O	2.54	0.41
4:E:513:LEU:O	4:E:517:LEU:N	2.51	0.41
4:E:718:GLN:HE21	4:E:825:VAL:HG21	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:342:ASP:OD1	4:G:342:ASP:N	2.50	0.41
4:I:318:HIS:CE1	4:I:573:GLN:HG3	2.55	0.41
4:H:247:ASN:HB3	4:H:850:THR:HB	2.03	0.41
4:H:726:GLY:HA3	4:H:810:ASN:HD21	1.86	0.41
4:K:475:VAL:HG23	4:K:516:ALA:HB1	2.03	0.41
4:L:189:LEU:HD23	4:L:189:LEU:HA	1.90	0.41
4:M:134:GLU:HG2	4:M:158:LYS:HB2	2.02	0.41
4:M:212:ALA:HA	4:M:215:CYS:HB2	2.02	0.41
4:M:797:ARG:HH21	4:N:294:ASN:HB3	1.86	0.41
4:N:299:LEU:HA	4:N:300:PRO:HD3	1.91	0.41
4:N:415:MET:HG2	4:N:438:ILE:HD13	2.03	0.41
3:A:897:GLN:NE2	4:G:109:GLU:OE2	2.52	0.41
4:E:582:LYS:HE2	4:E:582:LYS:HB2	1.97	0.41
4:I:663:ARG:HD3	4:J:358:GLN:HE22	1.86	0.41
4:K:367:ILE:HD12	4:M:362:LEU:HB2	2.03	0.41
4:K:424:ASP:HB3	4:K:574:HIS:HE1	1.84	0.41
4:M:442:ILE:HA	4:M:536:ILE:HD11	2.01	0.41
4:M:877:ALA:HB2	4:M:883:ILE:HG23	2.03	0.41
4:N:553:LEU:HD11	4:N:591:LEU:HG	2.03	0.41
4:G:569:THR:OG1	4:G:614:ASN:O	2.39	0.40
4:J:499:VAL:HG12	4:J:565:MET:HE1	2.02	0.40
4:K:181:LEU:HD11	4:K:685:ARG:HG3	2.03	0.40
4:K:836:ILE:HA	4:K:836:ILE:HD13	1.76	0.40
4:N:841:ASP:OD1	4:N:841:ASP:N	2.48	0.40
3:A:142:LEU:HD23	3:A:142:LEU:HA	1.94	0.40
3:A:691:ALA:HB2	3:A:723:ARG:HB3	2.03	0.40
4:E:117:ASP:OD1	4:E:117:ASP:N	2.54	0.40
4:E:436:LEU:HB2	4:E:462:PRO:HG3	2.03	0.40
4:F:319:ASP:OD1	4:F:320:ASN:N	2.54	0.40
4:G:335:LEU:HD11	4:G:613:VAL:HG11	2.03	0.40
4:G:802:LEU:HD12	4:G:802:LEU:HA	1.92	0.40
4:J:285:ILE:CD1	4:J:607:LEU:HD21	2.48	0.40
4:K:310:LEU:HD21	4:K:615:VAL:HG13	2.03	0.40
4:L:654:LYS:HB3	4:L:654:LYS:HE3	1.82	0.40
4:N:242:ASP:OD1	4:N:243:ARG:N	2.54	0.40
3:A:734:THR:HA	3:A:737:MET:HE2	2.04	0.40
4:G:555:ARG:HB3	4:G:883:ILE:HG22	2.02	0.40
4:I:315:LEU:HD13	4:I:672:LEU:HD11	2.04	0.40
4:J:155:TRP:CD2	4:J:253:HIS:HD2	2.39	0.40
4:J:335:LEU:HD21	4:J:613:VAL:HG13	2.03	0.40
4:K:417:LEU:HD23	4:K:417:LEU:HA	1.91	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:343:LEU:HD23	4:L:346:LEU:HD21	2.03	0.40
4:M:162:LEU:HA	4:M:162:LEU:HD23	1.82	0.40
4:N:656:LEU:HD23	4:N:659:PHE:HE2	1.85	0.40
4:F:160:ASP:OD1	4:F:160:ASP:N	2.45	0.40
4:F:852:ASN:HD22	4:F:855:PHE:HE1	1.68	0.40
4:G:776:SER:HB3	4:G:813:SER:HB2	2.03	0.40
4:I:200:SER:HA	4:I:233:ALA:HA	2.02	0.40
4:H:156:LYS:HE2	4:H:156:LYS:HB3	1.88	0.40
4:J:661:ILE:O	4:J:661:ILE:HG22	2.20	0.40
4:N:256:ASP:O	4:N:260:ASN:ND2	2.55	0.40
4:E:660:ASP:OD1	4:E:660:ASP:N	2.47	0.40
4:G:284:ARG:HH11	4:G:492:ILE:HD12	1.87	0.40
4:G:623:GLU:O	4:G:624:ARG:NH2	2.51	0.40
4:G:624:ARG:HD3	4:G:624:ARG:HA	1.76	0.40
4:G:793:ILE:HD11	4:G:802:LEU:HD13	2.04	0.40
4:H:125:LYS:NZ	4:H:192:ASP:OD2	2.54	0.40
4:K:885:ASN:OD1	4:K:885:ASN:N	2.49	0.40
4:M:73:LYS:H	4:M:73:LYS:HG2	1.63	0.40
4:M:190:LEU:HG	4:M:851:SER:HB3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	
3	A	1080/1088 (99%)	1025 (95%)	55 (5%)	0	100	100
4	E	792/887 (89%)	762 (96%)	30 (4%)	0	100	100
4	F	801/887 (90%)	774 (97%)	27 (3%)	0	100	100
4	G	790/887 (89%)	755 (96%)	35 (4%)	0	100	100
4	H	803/887 (90%)	769 (96%)	34 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	I	779/887 (88%)	745 (96%)	34 (4%)	0	100	100
4	J	803/887 (90%)	766 (95%)	37 (5%)	0	100	100
4	K	779/887 (88%)	744 (96%)	35 (4%)	0	100	100
4	L	803/887 (90%)	771 (96%)	32 (4%)	0	100	100
4	M	824/887 (93%)	790 (96%)	34 (4%)	0	100	100
4	N	803/887 (90%)	773 (96%)	30 (4%)	0	100	100
All	All	9057/9958 (91%)	8674 (96%)	383 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	
3	A	983/989 (99%)	983 (100%)	0	100	100
4	E	730/818 (89%)	730 (100%)	0	100	100
4	F	739/818 (90%)	737 (100%)	2 (0%)	86	83
4	G	730/818 (89%)	729 (100%)	1 (0%)	88	87
4	H	741/818 (91%)	738 (100%)	3 (0%)	84	81
4	I	717/818 (88%)	716 (100%)	1 (0%)	88	87
4	J	741/818 (91%)	740 (100%)	1 (0%)	88	87
4	K	717/818 (88%)	716 (100%)	1 (0%)	88	87
4	L	741/818 (91%)	740 (100%)	1 (0%)	88	87
4	M	761/818 (93%)	758 (100%)	3 (0%)	84	81
4	N	741/818 (91%)	740 (100%)	1 (0%)	88	87
All	All	8341/9169 (91%)	8327 (100%)	14 (0%)	85	85

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	F	206	VAL
4	F	847	HIS
4	G	239	VAL
4	I	557	LEU
4	H	206	VAL
4	H	491	VAL
4	H	874	ASN
4	J	206	VAL
4	K	799	VAL
4	L	491	VAL
4	M	161	THR
4	M	246	VAL
4	M	443	ILE
4	N	206	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (178) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	143	ASN
3	A	283	GLN
3	A	407	GLN
3	A	530	ASN
3	A	547	ASN
3	A	653	ASN
3	A	697	ASN
3	A	862	GLN
3	A	908	GLN
3	A	920	GLN
3	A	954	HIS
3	A	983	GLN
3	A	1048	ASN
4	E	122	GLN
4	E	151	ASN
4	E	175	ASN
4	E	247	ASN
4	E	332	ASN
4	E	396	GLN
4	E	423	ASN
4	E	450	GLN
4	E	480	HIS
4	E	497	ASN
4	E	573	GLN
4	E	614	ASN

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Mol	Chain	Res	Type
4	E	668	GLN
4	E	718	GLN
4	E	760	ASN
4	E	835	GLN
4	F	126	GLN
4	F	260	ASN
4	F	267	GLN
4	F	320	ASN
4	F	352	GLN
4	F	380	GLN
4	F	453	HIS
4	F	512	GLN
4	F	573	GLN
4	F	695	ASN
4	F	760	ASN
4	G	126	GLN
4	G	199	ASN
4	G	240	GLN
4	G	247	ASN
4	G	265	GLN
4	G	274	ASN
4	G	358	GLN
4	G	361	GLN
4	G	480	HIS
4	G	545	ASN
4	G	573	GLN
4	G	574	HIS
4	G	614	ASN
4	G	636	ASN
4	G	668	GLN
4	G	708	GLN
4	G	718	GLN
4	G	762	GLN
4	G	835	GLN
4	G	880	ASN
4	I	175	ASN
4	I	257	HIS
4	I	260	ASN
4	I	274	ASN
4	I	287	ASN
4	I	290	ASN
4	I	361	GLN

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Mol	Chain	Res	Type
4	I	380	GLN
4	I	396	GLN
4	I	573	GLN
4	I	574	HIS
4	I	614	ASN
4	I	620	ASN
4	I	668	GLN
4	I	708	GLN
4	I	753	GLN
4	I	762	GLN
4	I	821	ASN
4	I	835	GLN
4	H	151	ASN
4	H	274	ASN
4	H	287	ASN
4	H	294	ASN
4	H	309	ASN
4	H	316	ASN
4	H	320	ASN
4	H	352	GLN
4	H	450	GLN
4	H	453	HIS
4	H	512	GLN
4	H	636	ASN
4	H	695	ASN
4	H	738	GLN
4	H	753	GLN
4	H	760	ASN
4	H	838	GLN
4	H	847	HIS
4	J	89	HIS
4	J	94	GLN
4	J	240	GLN
4	J	267	GLN
4	J	312	GLN
4	J	352	GLN
4	J	354	GLN
4	J	358	GLN
4	J	368	GLN
4	J	407	ASN
4	J	484	ASN
4	J	512	GLN

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Mol	Chain	Res	Type
4	J	545	ASN
4	J	574	HIS
4	J	612	ASN
4	J	657	GLN
4	J	708	GLN
4	J	761	ASN
4	J	762	GLN
4	J	810	ASN
4	K	126	GLN
4	K	260	ASN
4	K	274	ASN
4	K	372	GLN
4	K	396	GLN
4	K	450	GLN
4	K	483	ASN
4	K	545	ASN
4	K	573	GLN
4	K	614	ASN
4	K	668	GLN
4	K	760	ASN
4	K	835	GLN
4	K	880	ASN
4	L	94	GLN
4	L	260	ASN
4	L	265	GLN
4	L	267	GLN
4	L	312	GLN
4	L	368	GLN
4	L	380	GLN
4	L	453	HIS
4	L	512	GLN
4	L	622	ASN
4	L	626	ASN
4	L	636	ASN
4	L	690	ASN
4	L	698	GLN
4	L	756	ASN
4	L	760	ASN
4	M	90	GLN
4	M	199	ASN
4	M	260	ASN
4	M	274	ASN

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Mol	Chain	Res	Type
4	M	320	ASN
4	M	396	GLN
4	M	423	ASN
4	M	506	ASN
4	M	573	GLN
4	M	574	HIS
4	M	614	ASN
4	M	620	ASN
4	M	668	GLN
4	M	708	GLN
4	M	760	ASN
4	M	835	GLN
4	N	90	GLN
4	N	220	GLN
4	N	287	ASN
4	N	309	ASN
4	N	407	ASN
4	N	453	HIS
4	N	497	ASN
4	N	512	GLN
4	N	545	ASN
4	N	639	ASN
4	N	695	ASN
4	N	741	ASN
4	N	760	ASN
4	N	835	GLN
4	N	852	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	B	17/18 (94%)	5 (29%)	0
2	C	16/17 (94%)	5 (31%)	0
All	All	33/35 (94%)	10 (30%)	0

All (10) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	B	585	A
1	B	586	U
1	B	588	U

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Mol	Chain	Res	Type
1	B	589	A
1	B	590	U
2	C	591	A
2	C	592	U
2	C	597	A
2	C	604	U
2	C	605	A

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	UTP	A	1101	-	29,30,30	0.55	0	43,47,47	0.66	0
6	GTP	A	1102	-	33,34,34	1.01	4 (12%)	50,54,54	1.61	8 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	UTP	A	1101	-	-	0/22/38/38	0/2/2/2
6	GTP	A	1102	-	-	6/22/38/38	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1102	GTP	PB-O3A	2.25	1.61	1.59
6	A	1102	GTP	PB-O3B	2.12	1.61	1.59
6	A	1102	GTP	C2-N3	2.08	1.38	1.33
6	A	1102	GTP	PA-O3A	2.07	1.61	1.59

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1102	GTP	C5-C4-N3	-4.89	120.61	128.39
6	A	1102	GTP	C2-N3-C4	4.55	120.14	112.30
6	A	1102	GTP	C2-N1-C6	-3.06	119.57	125.11
6	A	1102	GTP	N9-C4-N3	2.98	131.92	125.95
6	A	1102	GTP	N9-C8-N7	-2.79	108.23	113.40
6	A	1102	GTP	C5-C6-N1	2.69	120.10	113.25
6	A	1102	GTP	C8-N7-C5	2.60	108.89	104.26
6	A	1102	GTP	O6-C6-C5	-2.21	120.69	126.53

There are no chirality outliers.

All (6) torsion outliers are listed below:

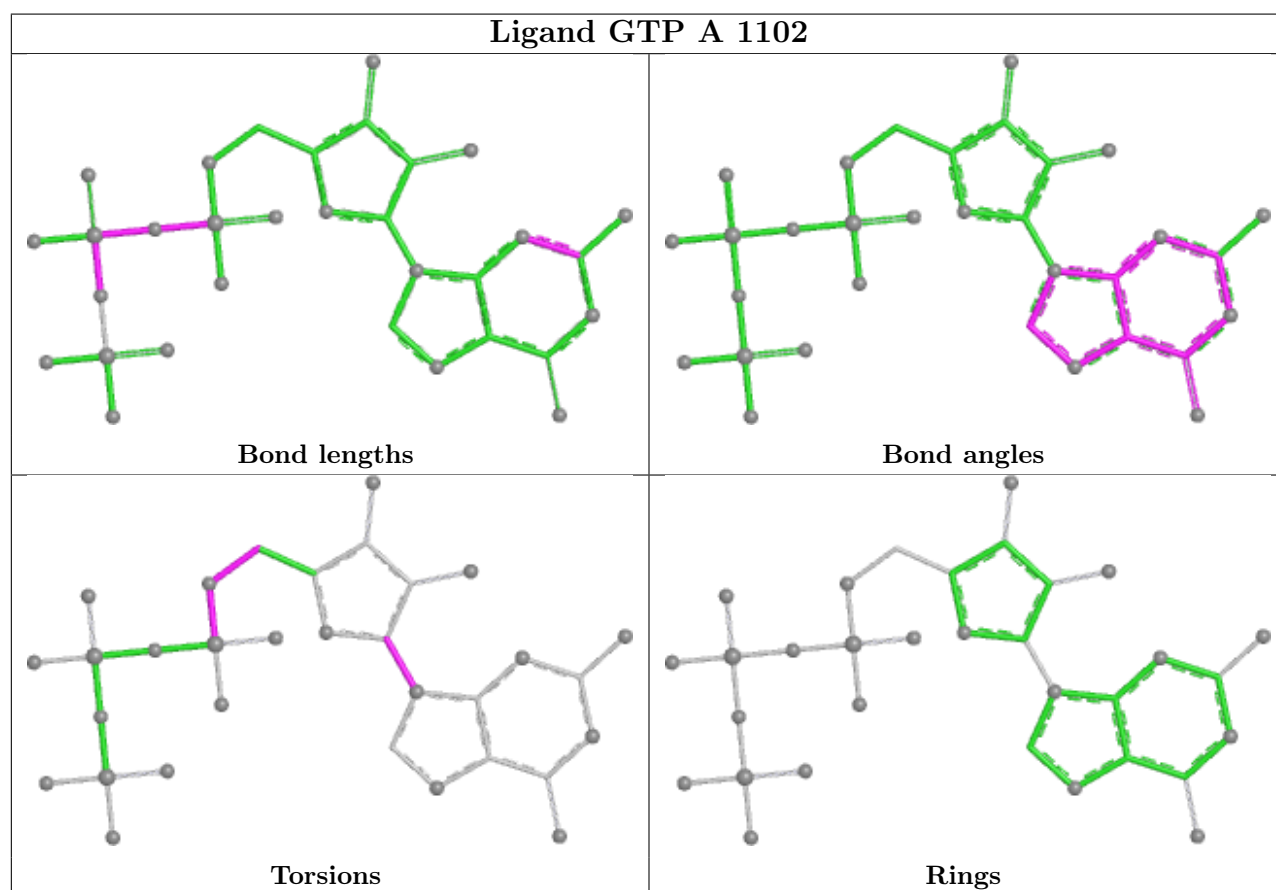
Mol	Chain	Res	Type	Atoms
6	A	1102	GTP	C5'-O5'-PA-O3A
6	A	1102	GTP	C5'-O5'-PA-O1A
6	A	1102	GTP	C5'-O5'-PA-O2A
6	A	1102	GTP	C4'-C5'-O5'-PA
6	A	1102	GTP	O4'-C1'-N9-C4
6	A	1102	GTP	O4'-C1'-N9-C8

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1101	UTP	1	0
6	A	1102	GTP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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.

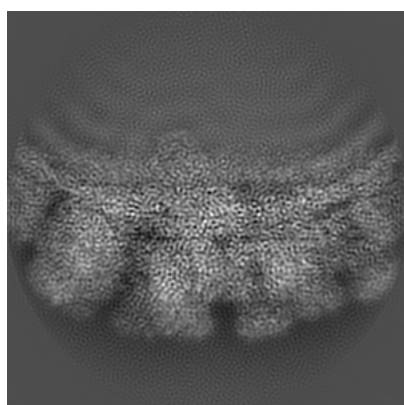
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-20060. These allow visual inspection of the internal detail of the map and identification of artifacts.

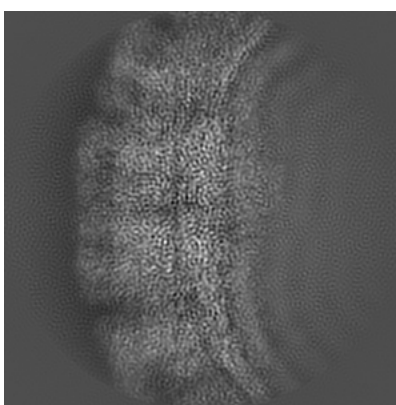
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

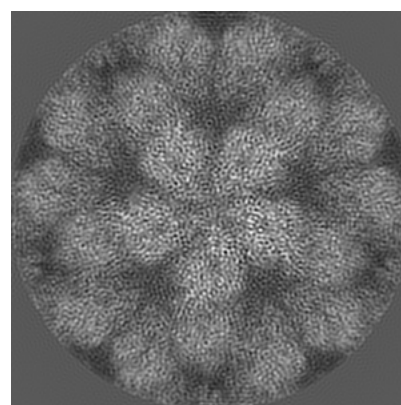
6.1.1 Primary map



X



Y

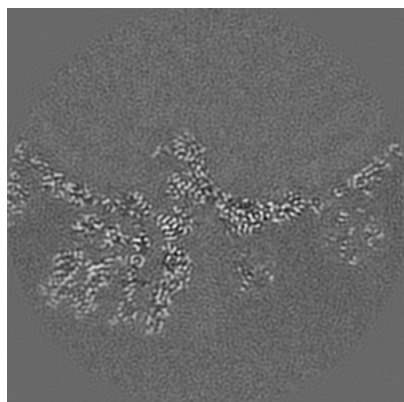


Z

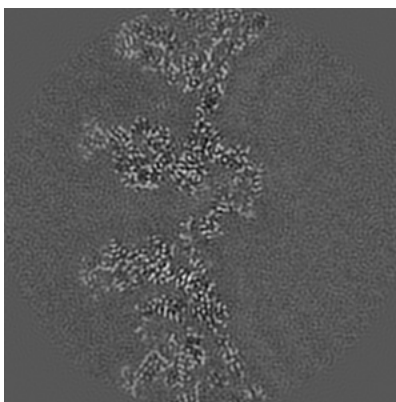
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

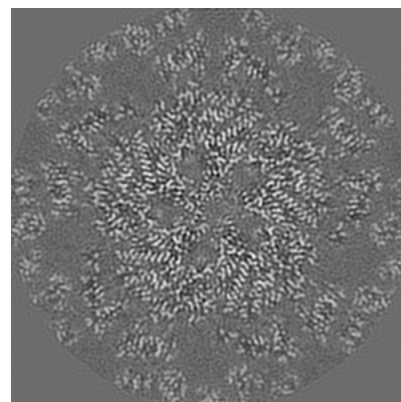
6.2.1 Primary map



X Index: 132



Y Index: 132

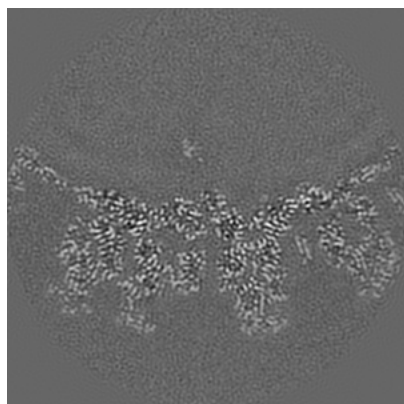


Z Index: 132

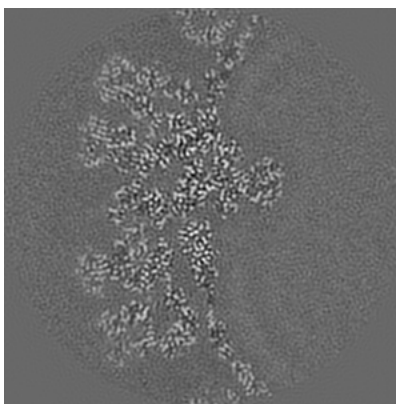
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

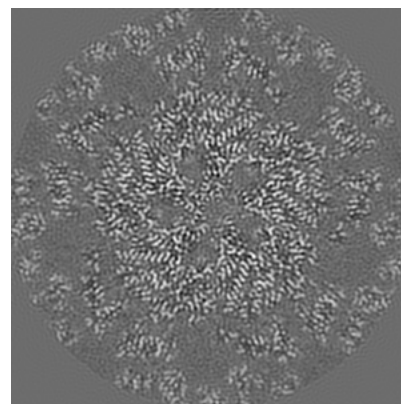
6.3.1 Primary map



X Index: 117



Y Index: 109

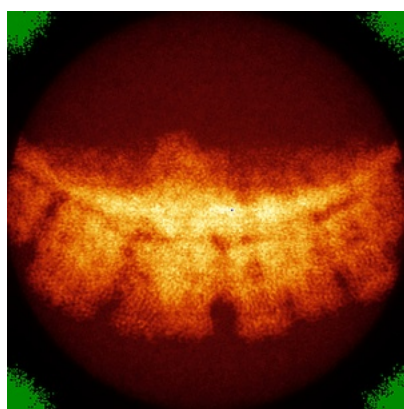


Z Index: 132

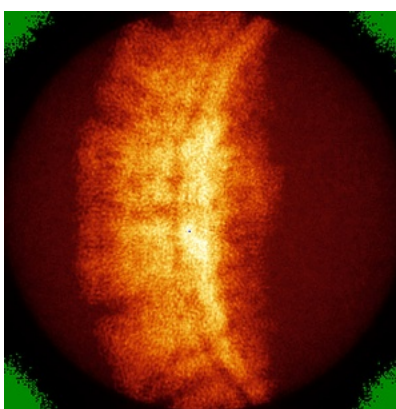
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

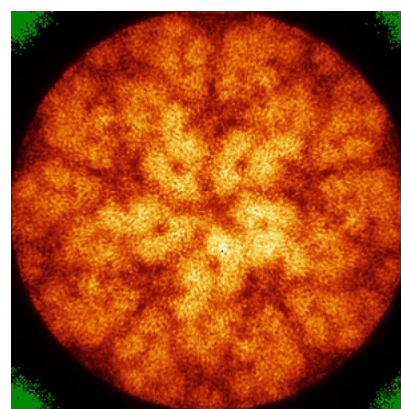
6.4.1 Primary map



X



Y

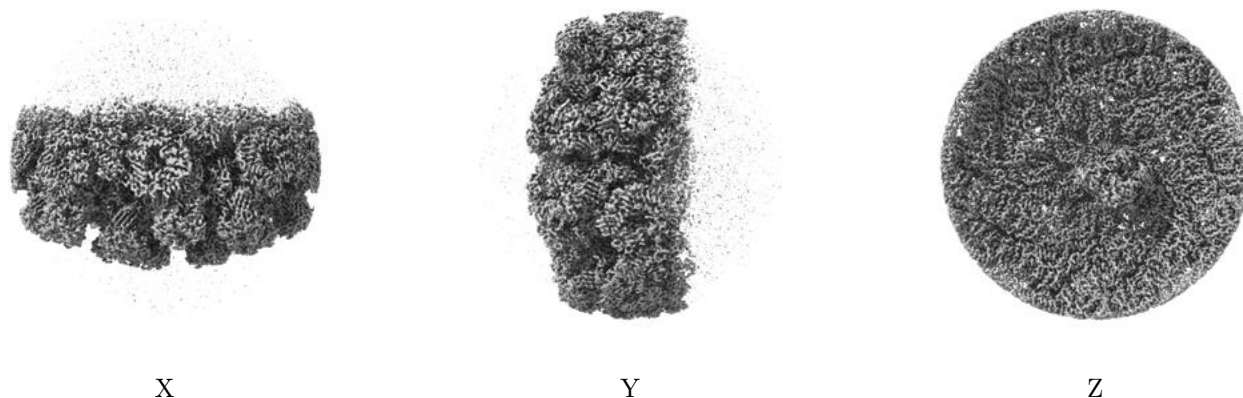


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0175. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

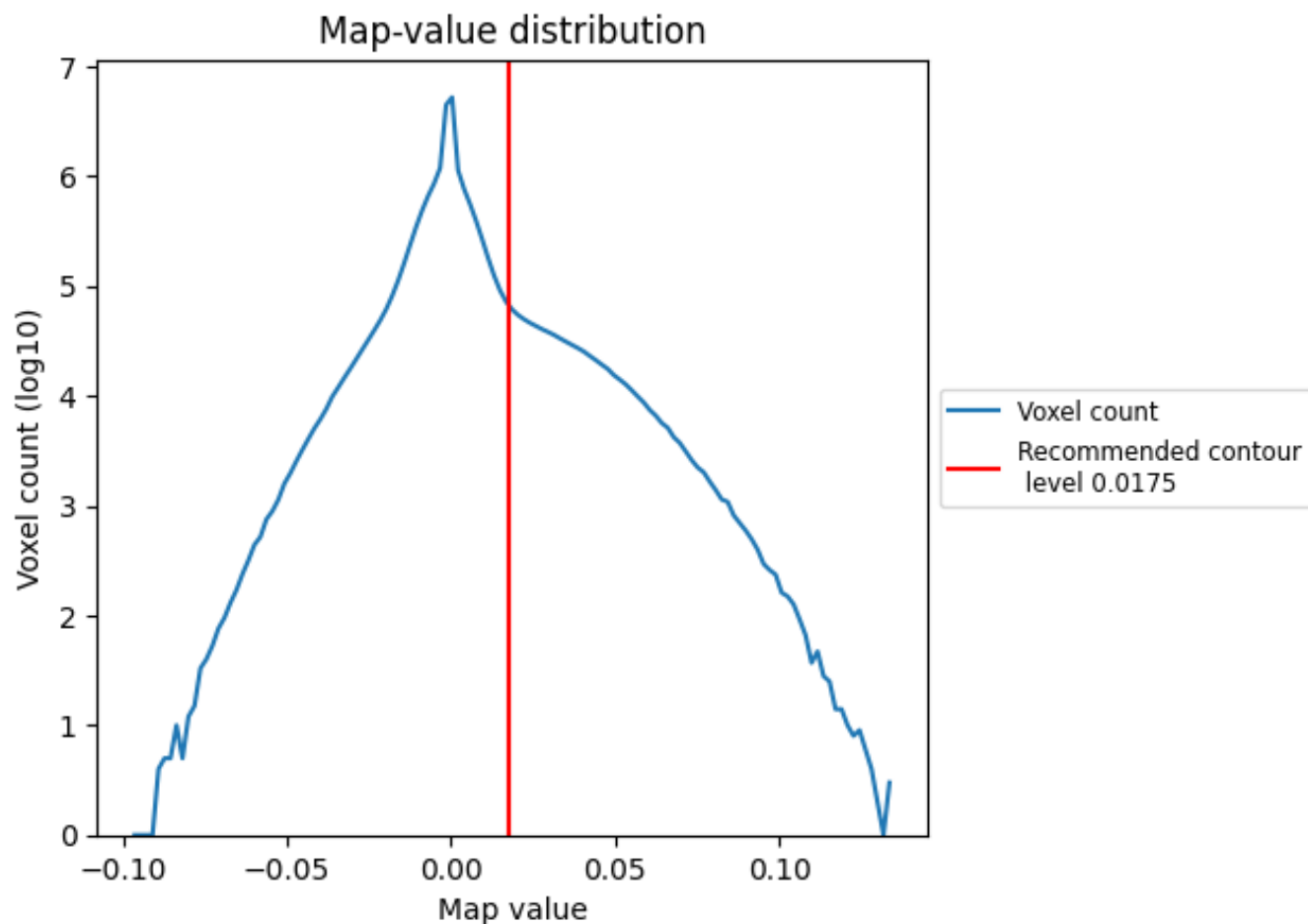
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

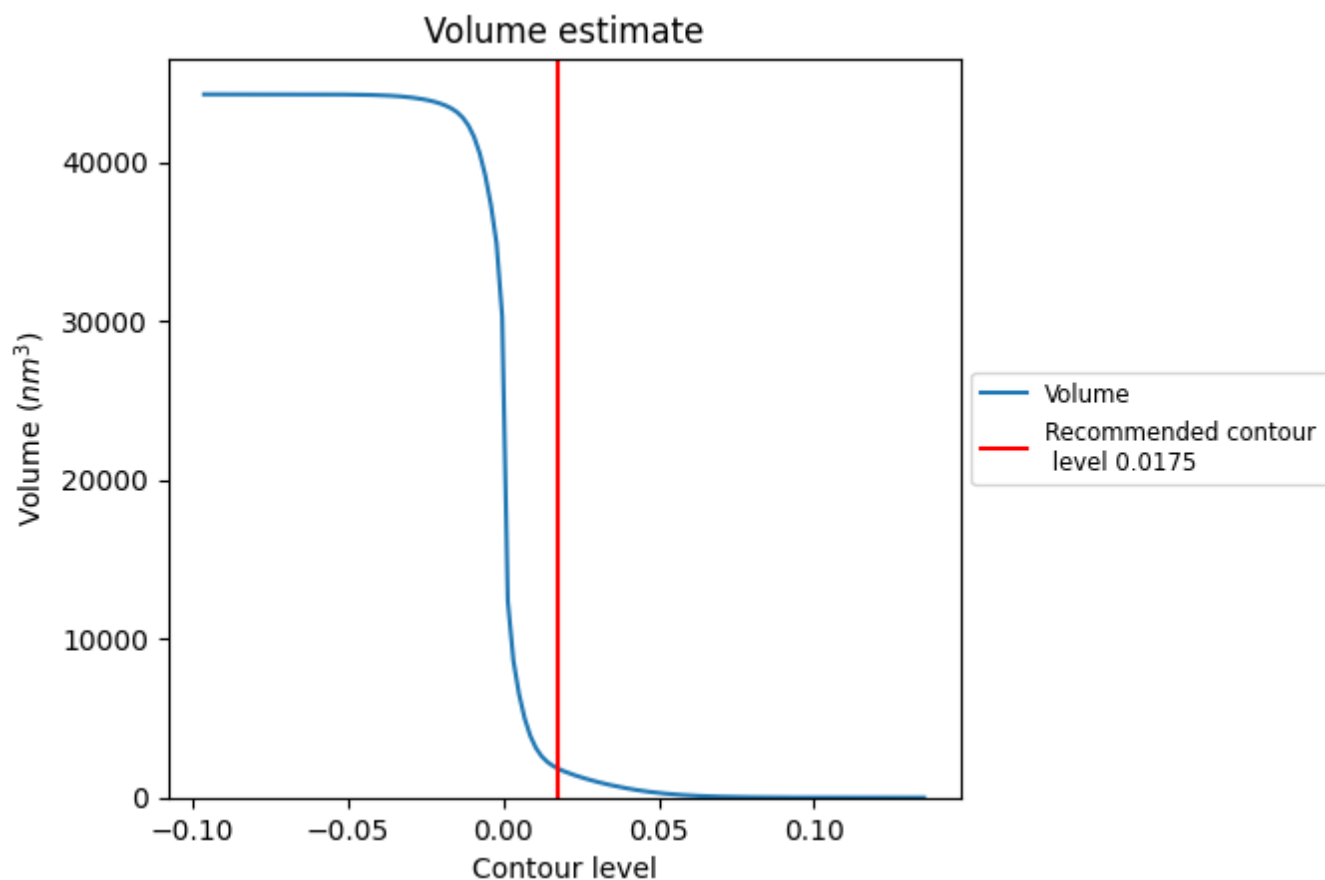
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

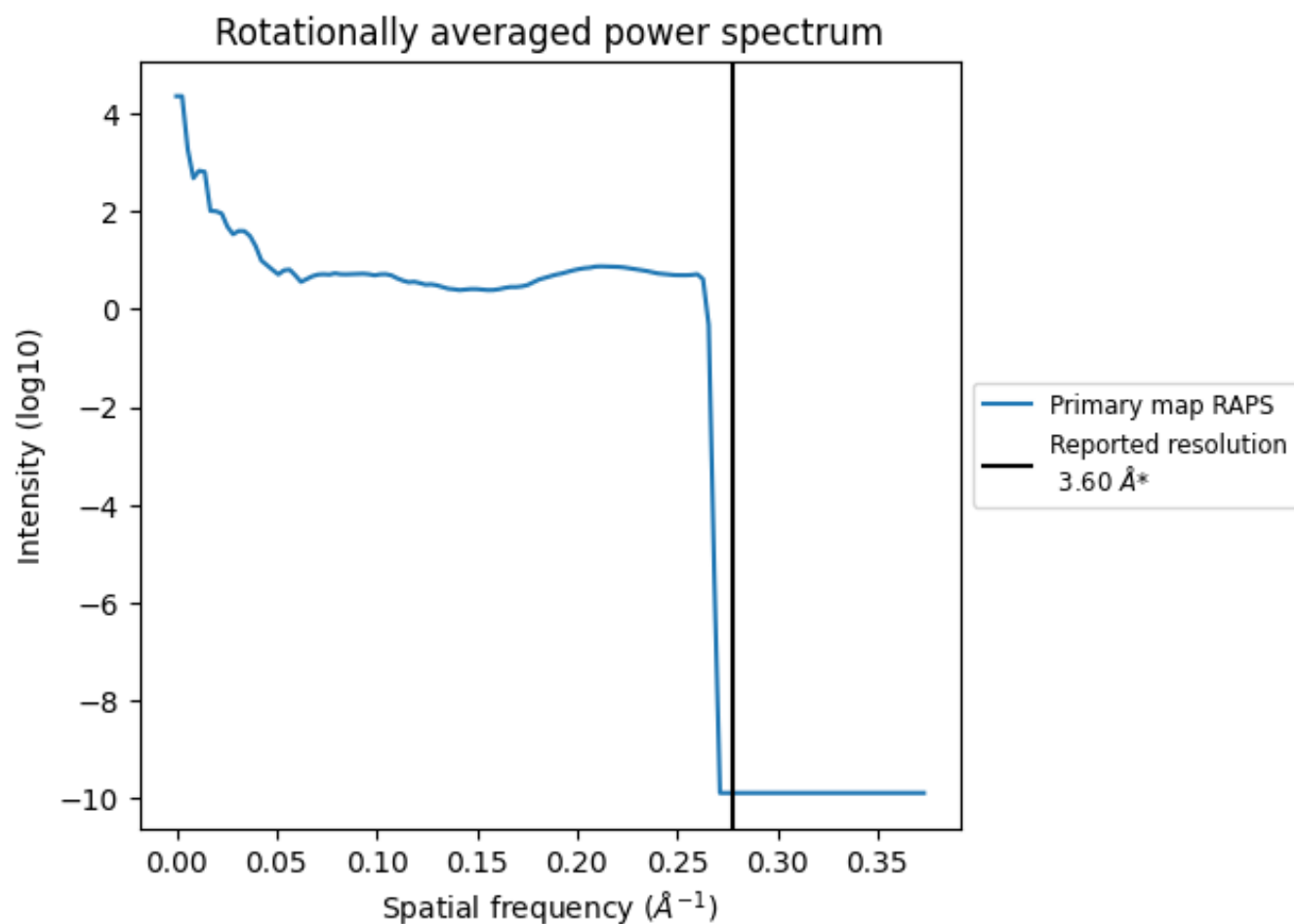
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1833 nm³; this corresponds to an approximate mass of 1656 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.278 Å⁻¹

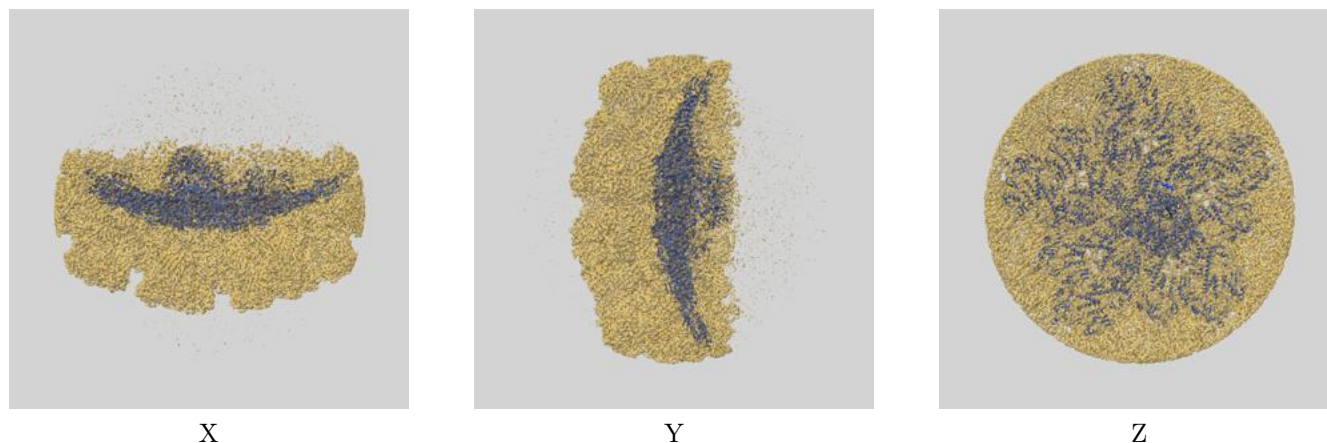
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

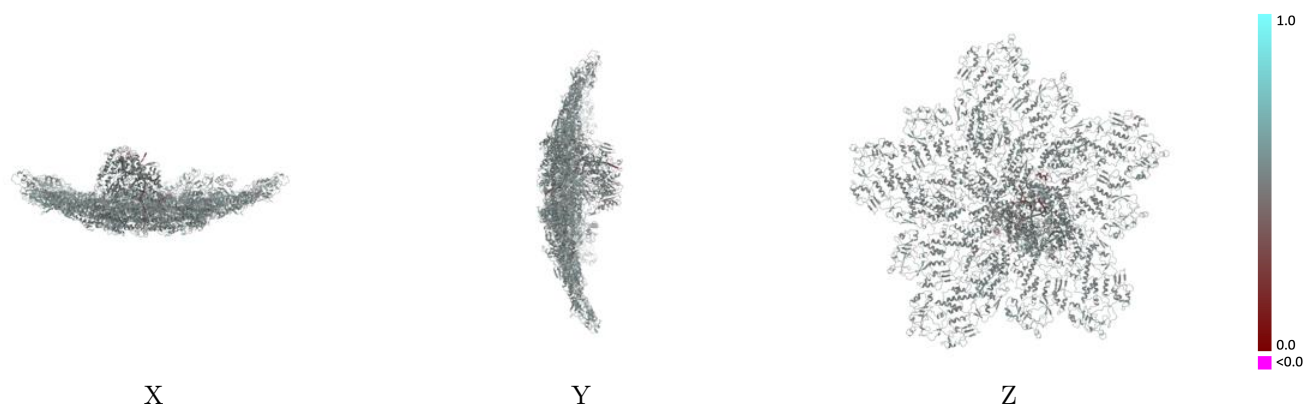
This section contains information regarding the fit between EMDB map EMD-20060 and PDB model 6OGZ. Per-residue inclusion information can be found in [section 3](#) on [page 7](#).

9.1 Map-model overlay [i](#)



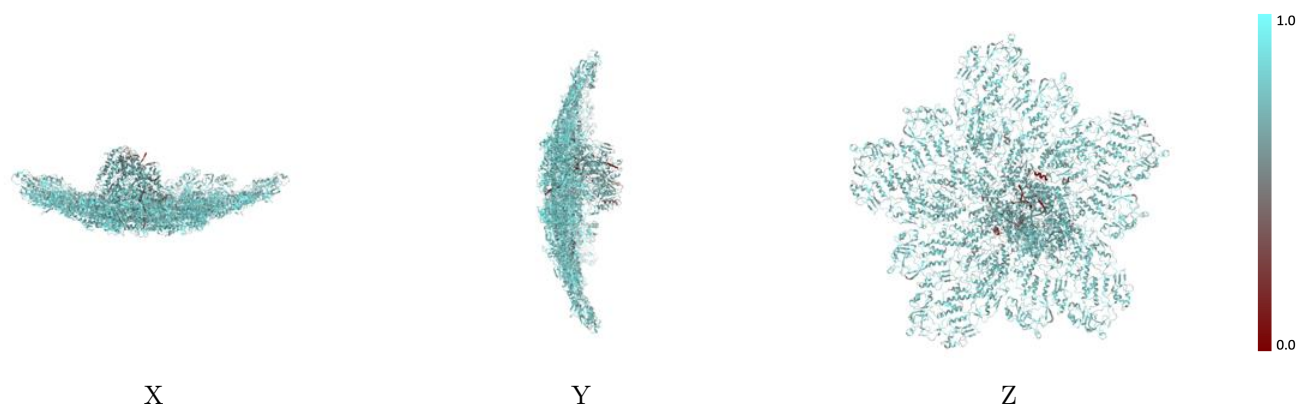
The images above show the 3D surface view of the map at the recommended contour level 0.0175 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



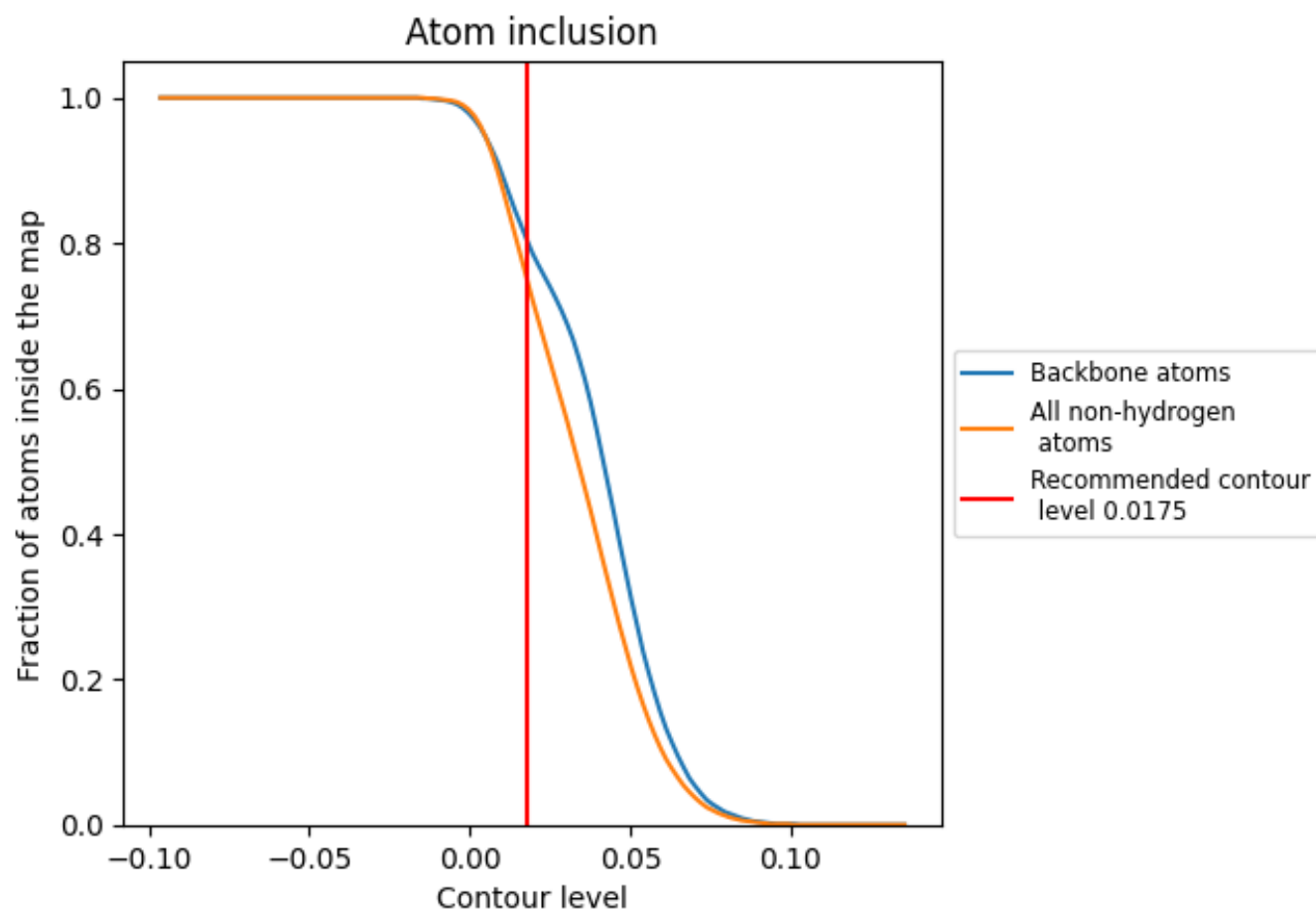
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.0175).

9.4 Atom inclusion [i](#)



At the recommended contour level, 81% of all backbone atoms, 76% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.0175) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7550	 0.5170
A	 0.7250	 0.5130
B	 0.4000	 0.3480
C	 0.4860	 0.3940
E	 0.7530	 0.5160
F	 0.7580	 0.5190
G	 0.7650	 0.5210
H	 0.7510	 0.5160
I	 0.7850	 0.5240
J	 0.7490	 0.5180
K	 0.7830	 0.5220
L	 0.7520	 0.5170
M	 0.7750	 0.5210
N	 0.7530	 0.5160

