



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 04:03 PM UTC

PDB ID : 7Y85 / pdb_00007y85
EMDB ID : EMD-33681
Title : CryoEM structure of type III-E CRISPR Craspase gRAMP-crRNA in complex with TPR-CHAT protease bound to self RNA target
Authors : Zhang, J.T.; Cui, N.; Huang, H.D.; Jia, N.
Deposited on : 2022-06-22
Resolution : 2.73 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
MolProbity : 4-5-2 with Phenix2.0
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

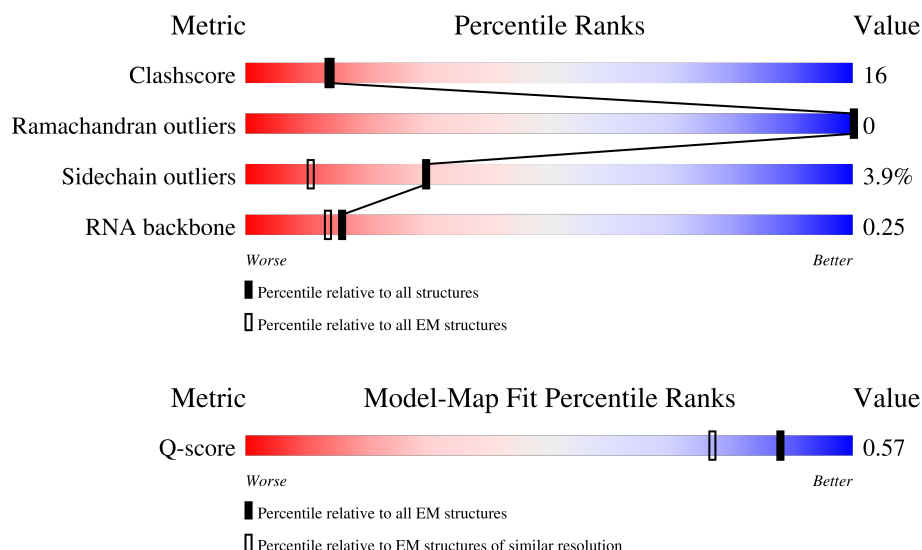
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.73 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	10432 (2.23 - 3.23)

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	A	1728	7% 45% 25% 29%
2	B	110	7% 14% 10% 67%
3	C	56	7% 12% 18% 11% 59%

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Mol	Chain	Length	Quality of chain
4	D	746	25% 53% 33% • 13%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 16502 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 RAMP superfamily protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1232	9941	6333	1734	1843	31	0	0

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	initiating methionine	UNP A0A0B0EGF3
A	-4	HIS	-	expression tag	UNP A0A0B0EGF3
A	-3	HIS	-	expression tag	UNP A0A0B0EGF3
A	-2	HIS	-	expression tag	UNP A0A0B0EGF3
A	-1	HIS	-	expression tag	UNP A0A0B0EGF3
A	0	HIS	-	expression tag	UNP A0A0B0EGF3
A	1	HIS	-	expression tag	UNP A0A0B0EGF3
A	2	LYS	-	expression tag	UNP A0A0B0EGF3
A	3	SER	-	expression tag	UNP A0A0B0EGF3
A	4	ASN	-	expression tag	UNP A0A0B0EGF3
A	5	ASP	-	expression tag	UNP A0A0B0EGF3

- Molecule 2 is a RNA chain called crRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
2	B	36	757	339	128	254	36	0	0

- Molecule 3 is a RNA chain called Self RNA target.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
3	C	23	496	221	93	159	23	0	0

- Molecule 4 is a protein called CHAT domain protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	652	Total	C	N	O	S	0	0
			5303	3391	907	980	25		

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	717	GLY	-	expression tag	UNP A0A0B0EKL4
D	718	GLY	-	expression tag	UNP A0A0B0EKL4
D	719	SER	-	expression tag	UNP A0A0B0EKL4
D	720	GLY	-	expression tag	UNP A0A0B0EKL4
D	721	TRP	-	expression tag	UNP A0A0B0EKL4
D	722	SER	-	expression tag	UNP A0A0B0EKL4
D	723	HIS	-	expression tag	UNP A0A0B0EKL4
D	724	PRO	-	expression tag	UNP A0A0B0EKL4
D	725	GLN	-	expression tag	UNP A0A0B0EKL4
D	726	PHE	-	expression tag	UNP A0A0B0EKL4
D	727	GLU	-	expression tag	UNP A0A0B0EKL4
D	728	LYS	-	expression tag	UNP A0A0B0EKL4
D	729	GLY	-	expression tag	UNP A0A0B0EKL4
D	730	GLY	-	expression tag	UNP A0A0B0EKL4
D	731	GLY	-	expression tag	UNP A0A0B0EKL4
D	732	SER	-	expression tag	UNP A0A0B0EKL4
D	733	GLY	-	expression tag	UNP A0A0B0EKL4
D	734	GLY	-	expression tag	UNP A0A0B0EKL4
D	735	GLY	-	expression tag	UNP A0A0B0EKL4
D	736	SER	-	expression tag	UNP A0A0B0EKL4
D	737	GLY	-	expression tag	UNP A0A0B0EKL4
D	738	GLY	-	expression tag	UNP A0A0B0EKL4
D	739	TRP	-	expression tag	UNP A0A0B0EKL4
D	740	SER	-	expression tag	UNP A0A0B0EKL4
D	741	HIS	-	expression tag	UNP A0A0B0EKL4
D	742	PRO	-	expression tag	UNP A0A0B0EKL4
D	743	GLN	-	expression tag	UNP A0A0B0EKL4
D	744	PHE	-	expression tag	UNP A0A0B0EKL4
D	745	GLU	-	expression tag	UNP A0A0B0EKL4
D	746	LYS	-	expression tag	UNP A0A0B0EKL4

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
5	A	4	Total	Zn	0
			4	4	

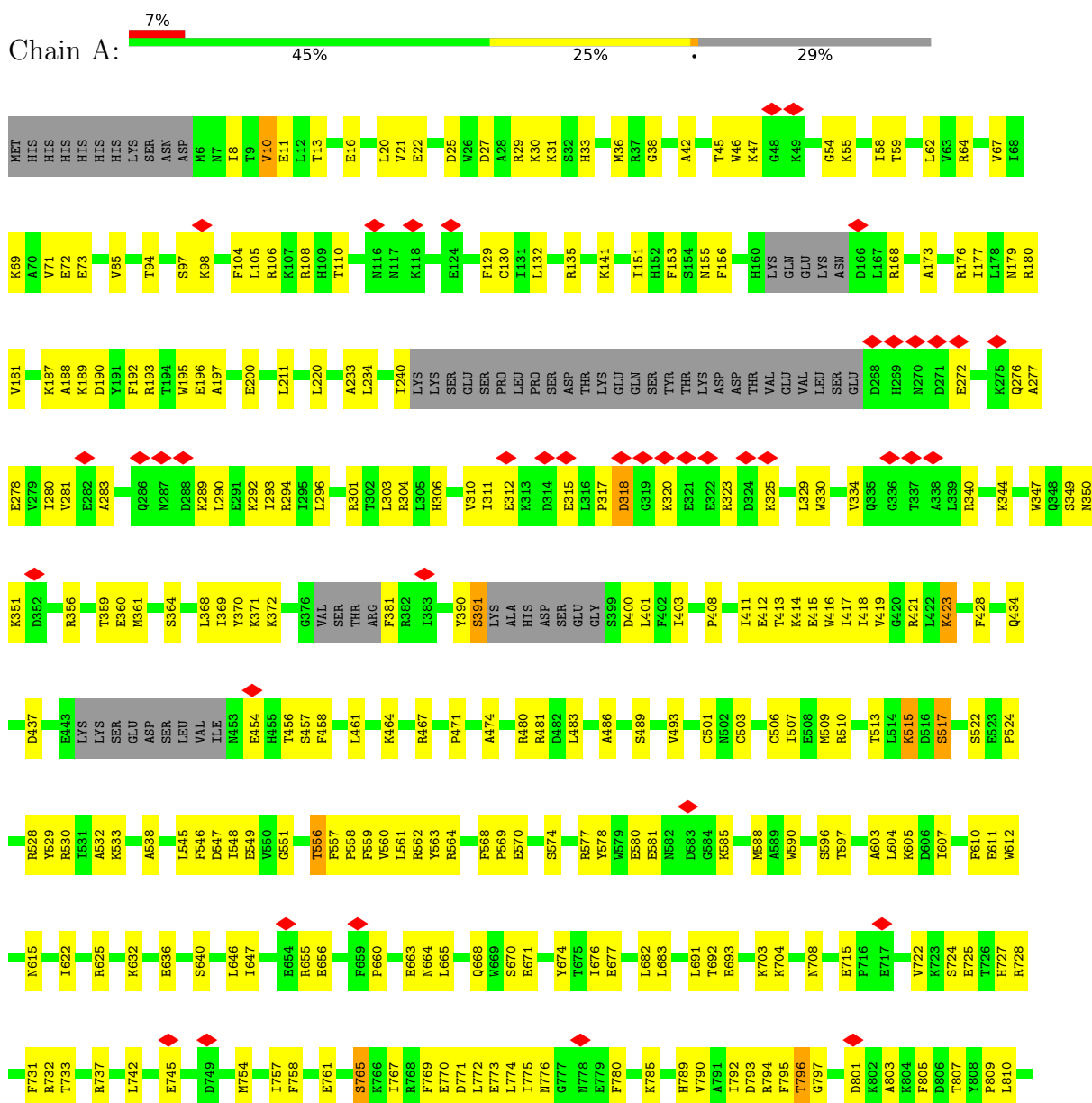
- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
6	A	1	Total 1	Mg 1	0

3 Residue-property plots

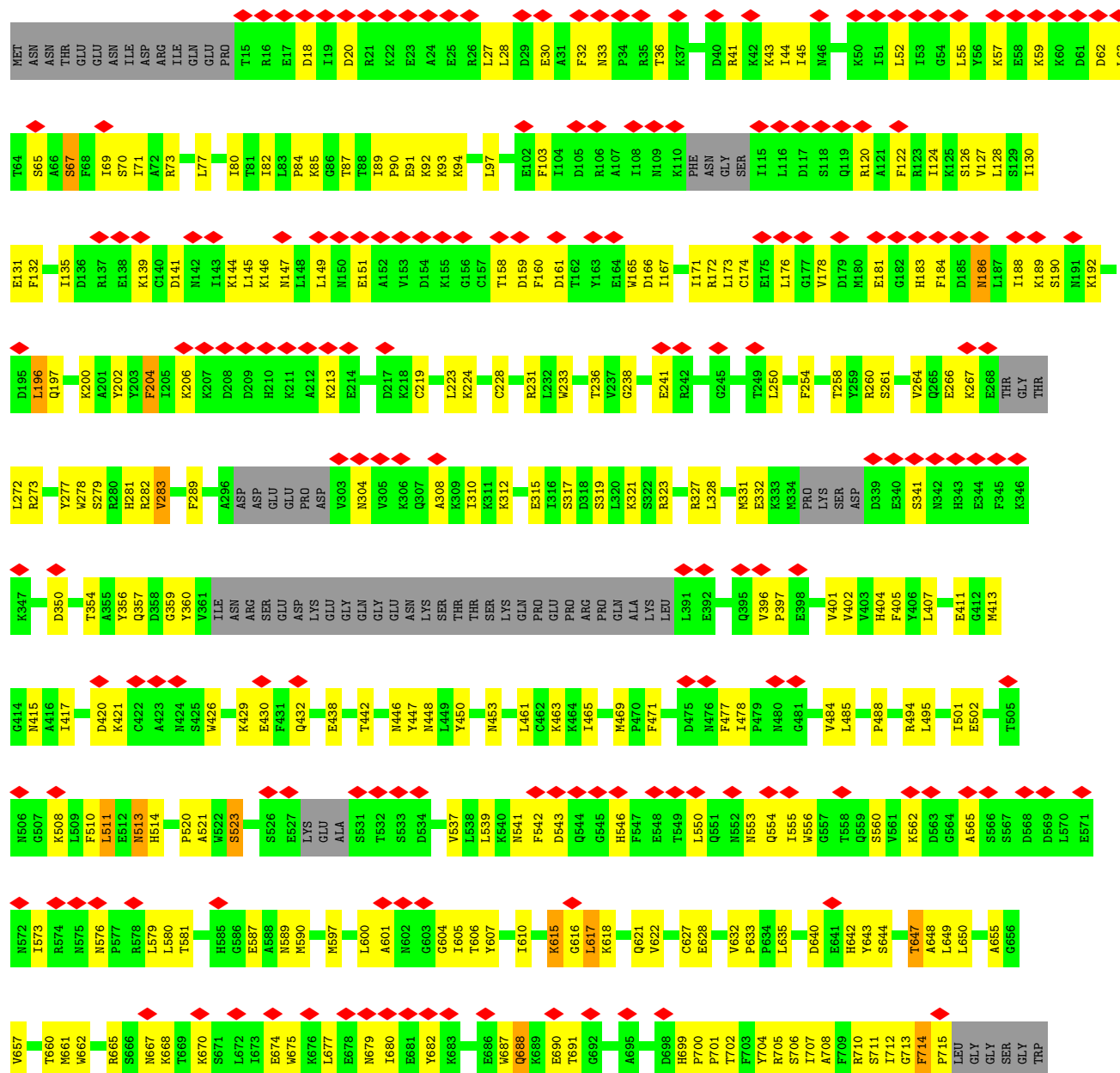
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: RAMP superfamily protein





G U U U U U G A A A C C A A A G G A A G G
G19 A20 C21 U22 A25 U26 U27 U28 C29 A30 A31 C32 C33 C34 U35 C36 C37 C38 C39 A40 A41 U42 U43 U44 U45 U46 C47 C48 C49 C50 C51 C52 G53 A54 C



SER HIS PRO GLN PHE GLU LYS GLY GLY GLY SER GLY GLY GLY SER GLY GLY TRP SER HIS PRO GLN PHE GLU LYS

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	123862	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	9.560	Depositor
Minimum map value	-6.045	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.218	Depositor
Recommended contour level	1	Depositor
Map size ($\text{\AA}$)	281.6, 281.6, 281.6	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.1, 1.1, 1.1	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, MG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.18	0/10174	0.30	0/13710
2	B	0.33	0/843	0.50	3/1309 (0.2%)
3	C	0.14	0/555	0.25	0/864
4	D	0.14	0/5412	0.31	1/7287 (0.0%)
All	All	0.18	0/16984	0.31	4/23170 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	45	C	P-O3'-C3'	-5.71	111.64	120.20
2	B	46	U	C2'-C3'-O3'	-5.28	101.58	109.50
4	D	601	ALA	CB-CA-C	-5.19	110.17	117.23
2	B	47	G	O4'-C1'-N9	5.04	115.75	108.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	9941	0	9818	316	0
2	B	757	0	387	43	0
3	C	496	0	250	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	5303	0	5299	180	0
5	A	4	0	0	0	0
6	A	1	0	0	0	0
All	All	16502	0	15754	507	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 16.

All (507) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1595:TYR:CD2	1:A:1614:PRO:HB3	1.44	1.52
1:A:1595:TYR:CD2	1:A:1614:PRO:CB	2.20	1.23
1:A:1595:TYR:HD2	1:A:1614:PRO:CB	1.57	1.13
1:A:1595:TYR:CE2	1:A:1614:PRO:HB3	2.02	0.95
1:A:1595:TYR:CE2	1:A:1614:PRO:CB	2.62	0.81
1:A:580:GLU:HG3	1:A:607:ILE:HD12	1.62	0.79
1:A:1595:TYR:HD2	1:A:1614:PRO:HB2	1.46	0.79
4:D:579:LEU:HD23	4:D:680:ILE:HD11	1.65	0.77
4:D:59:LYS:HB3	4:D:62:ASP:HB3	1.70	0.73
1:A:1026:SER:HB2	1:A:1395:LEU:HD22	1.71	0.72
4:D:494:ARG:NH1	4:D:633:PRO:O	2.17	0.71
4:D:69:ILE:HG21	4:D:120:ARG:HE	1.55	0.71
1:A:501:CYS:O	1:A:510:ARG:NH2	2.24	0.70
1:A:456:THR:HG21	1:A:549:GLU:HG3	1.75	0.69
1:A:772:LEU:HD23	1:A:822:GLY:HA3	1.74	0.69
1:A:401:LEU:HB3	4:D:438:GLU:HG3	1.73	0.69
1:A:785:LYS:NZ	1:A:991:ASN:O	2.25	0.69
1:A:1004:ARG:NH1	1:A:1429:VAL:O	2.27	0.67
1:A:1644:ARG:HH12	1:A:1656:ILE:HD13	1.59	0.67
1:A:1689:ASN:HB3	1:A:1690:PRO:HD3	1.77	0.67
1:A:581:GLU:HB3	1:A:603:ALA:HB1	1.76	0.67
1:A:683:LEU:HB2	1:A:809:PRO:HB3	1.76	0.66
4:D:266:GLU:O	4:D:323:ARG:NH1	2.27	0.66
1:A:1595:TYR:CD2	1:A:1614:PRO:HB2	2.21	0.66
1:A:1426:LYS:HD3	2:B:43:U:H5"	1.77	0.66
1:A:728:ARG:HH22	1:A:765:SER:HB2	1.59	0.66
1:A:489:SER:HB3	4:D:587:GLU:HG3	1.78	0.66
1:A:794:ARG:NH1	2:B:46:U:OP1	2.25	0.66
4:D:321:LYS:O	4:D:643:TYR:OH	2.14	0.66
1:A:757:ILE:HG21	1:A:838:ILE:HD11	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1529:GLU:HB3	1:A:1593:LYS:HD3	1.77	0.65
1:A:417:ILE:HD13	1:A:562:ARG:HG2	1.77	0.65
1:A:304:ARG:NH1	1:A:360:GLU:OE2	2.30	0.65
1:A:361:MET:HE2	1:A:361:MET:HA	1.78	0.65
1:A:757:ILE:HG23	1:A:830:PHE:HZ	1.60	0.65
1:A:1631:VAL:HB	1:A:1634:LEU:HB2	1.79	0.65
1:A:306:HIS:NE2	1:A:692:THR:OG1	2.31	0.64
1:A:16:GLU:OE2	2:B:31:C:N4	2.31	0.64
1:A:434:GLN:NE2	1:A:454:GLU:OE1	2.31	0.64
1:A:310:VAL:HA	1:A:315:GLU:HB2	1.79	0.64
1:A:674:TYR:OH	1:A:727:HIS:NE2	2.26	0.64
4:D:665:ARG:HG2	4:D:705:ARG:HD3	1.79	0.64
1:A:1004:ARG:NH2	1:A:1420:PHE:O	2.32	0.63
1:A:296:LEU:HB2	1:A:329:LEU:HD11	1.80	0.63
4:D:272:LEU:HD11	4:D:341:SER:HB2	1.80	0.62
4:D:537:VAL:HG11	4:D:573:ILE:HG12	1.81	0.62
1:A:180:ARG:NH2	1:A:190:ASP:OD2	2.24	0.62
1:A:320:LYS:HB2	1:A:323:ARG:HB2	1.80	0.62
1:A:190:ASP:OD1	2:B:33:G:N2	2.33	0.62
1:A:400:ASP:HB2	2:B:27:G:H22	1.64	0.62
1:A:417:ILE:HD11	1:A:560:VAL:HG13	1.81	0.62
1:A:761:GLU:HG3	3:C:17:G:H21	1.63	0.62
1:A:975:LEU:HD21	1:A:1510:GLU:HG3	1.82	0.61
1:A:1647:VAL:HG21	1:A:1655:TYR:HB2	1.82	0.61
4:D:130:ILE:HG12	4:D:172:ARG:HG2	1.81	0.61
1:A:277:ALA:O	1:A:281:VAL:HG23	1.99	0.61
1:A:1444:ILE:HG23	1:A:1451:GLU:HB3	1.82	0.61
4:D:655:ALA:HB1	4:D:715:PRO:HG3	1.82	0.61
1:A:421:ARG:HE	1:A:556:THR:HB	1.66	0.61
3:C:5:G:H2'	3:C:6:G:H8	1.64	0.61
4:D:687:TRP:O	4:D:691:THR:OG1	2.18	0.61
1:A:47:LYS:H	1:A:54:GLY:HA2	1.66	0.60
1:A:408:PRO:HG2	1:A:411:ILE:HB	1.83	0.60
1:A:141:LYS:HD2	2:B:25:A:H5'	1.82	0.60
1:A:925:TYR:OH	1:A:1686:SER:O	2.19	0.60
1:A:925:TYR:HD2	1:A:1482:LEU:HD21	1.65	0.60
4:D:171:ILE:HD12	4:D:197:GLN:HE21	1.66	0.60
4:D:465:ILE:HG12	4:D:495:LEU:HD13	1.83	0.60
1:A:177:ILE:HG12	1:A:193:ARG:HG2	1.84	0.60
3:C:22:C:OP1	4:D:87:THR:N	2.32	0.59
1:A:312:GLU:OE2	1:A:351:LYS:NZ	2.35	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:461:LEU:HD23	1:A:471:PRO:HG3	1.85	0.59
1:A:515:LYS:HB2	1:A:560:VAL:HB	1.84	0.59
4:D:91:GLU:OE2	4:D:94:LYS:NZ	2.26	0.59
4:D:465:ILE:O	4:D:469:MET:HB2	2.02	0.59
4:D:622:VAL:HG11	4:D:649:LEU:HD13	1.84	0.59
1:A:391:SER:OG	2:B:30:A:N1	2.30	0.59
4:D:661:MET:HG2	4:D:710:ARG:HD2	1.83	0.59
1:A:483:LEU:HD12	1:A:509:MET:HE3	1.85	0.59
1:A:1598:LEU:HD13	1:A:1622:ILE:HD11	1.84	0.59
1:A:1660:ASP:HB3	1:A:1665:LYS:HD2	1.85	0.59
1:A:797:GLY:H	1:A:1479:LYS:HD2	1.67	0.59
1:A:8:ILE:HG22	1:A:10:VAL:HG12	1.83	0.58
1:A:765:SER:OG	1:A:767:ILE:O	2.21	0.58
4:D:173:LEU:HD22	4:D:178:VAL:HG21	1.85	0.58
4:D:87:THR:HA	4:D:277:TYR:CE2	2.39	0.58
1:A:1615:GLN:HB3	1:A:1617:HIS:NE2	2.19	0.58
4:D:627:CYS:HA	4:D:660:THR:HG23	1.85	0.58
1:A:104:PHE:O	2:B:27:G:O2'	2.19	0.57
1:A:329:LEU:HD22	1:A:691:LEU:HD13	1.86	0.57
1:A:998:ILE:HB	1:A:1434:ALA:HB3	1.86	0.57
1:A:356:ARG:HD2	1:A:524:PRO:HG3	1.87	0.57
1:A:1426:LYS:HB2	2:B:43:U:H4'	1.86	0.57
4:D:635:LEU:HD21	4:D:640:ASP:HB2	1.86	0.57
4:D:139:LYS:O	4:D:144:LYS:NZ	2.35	0.57
1:A:173:ALA:HB1	1:A:195:TRP:HB3	1.86	0.57
1:A:731:PHE:HZ	1:A:842:LEU:HD13	1.68	0.57
4:D:27:LEU:HD22	4:D:44:ILE:HG23	1.86	0.57
4:D:223:LEU:HD11	4:D:236:THR:HG21	1.86	0.57
4:D:312:LYS:HA	4:D:315:GLU:HG3	1.86	0.57
1:A:708:ASN:HB2	1:A:715:GLU:HG2	1.87	0.56
1:A:676:ILE:HG21	1:A:852:LEU:HD11	1.87	0.56
2:B:42:U:H2'	2:B:43:U:H6	1.69	0.56
4:D:520:PRO:HG2	4:D:647:THR:HB	1.86	0.56
1:A:801:ASP:O	3:C:7:C:O2'	2.15	0.56
2:B:47:G:O2'	2:B:48:C:OP2	2.22	0.56
3:C:9:G:H2'	3:C:10:A:H8	1.71	0.56
4:D:675:TRP:CE2	4:D:687:TRP:HB2	2.40	0.56
4:D:159:ASP:N	4:D:166:ASP:OD2	2.38	0.56
4:D:643:TYR:HB3	4:D:647:THR:HG21	1.88	0.56
1:A:423:LYS:HB2	1:A:605:LYS:HD3	1.88	0.56
1:A:538:ALA:HB3	1:A:1426:LYS:HB3	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:310:ILE:HG22	4:D:426:TRP:HB3	1.87	0.56
4:D:124:ILE:HA	4:D:127:VAL:HG22	1.89	0.55
1:A:72:GLU:OE2	1:A:106:ARG:NH2	2.39	0.55
1:A:728:ARG:NH2	1:A:758:PHE:O	2.40	0.55
1:A:1536:LEU:HD13	1:A:1587:ILE:HG23	1.87	0.55
4:D:442:THR:O	4:D:446:ASN:ND2	2.34	0.55
1:A:481:ARG:NH2	1:A:590:TRP:O	2.28	0.55
1:A:954:ILE:HG12	1:A:1567:LEU:HD12	1.88	0.55
4:D:158:THR:HG21	4:D:183:HIS:CE1	2.42	0.55
1:A:434:GLN:NE2	1:A:457:SER:O	2.40	0.54
2:B:35:A:H1'	2:B:36:C:C5	2.42	0.54
1:A:547:ASP:OD1	1:A:547:ASP:N	2.37	0.54
1:A:106:ARG:HB2	2:B:27:G:H1'	1.89	0.54
1:A:774:LEU:HD22	1:A:810:LEU:HD13	1.88	0.54
1:A:94:THR:HG23	2:B:27:G:N7	2.23	0.54
4:D:539:LEU:N	4:D:581:THR:O	2.34	0.54
1:A:745:GLU:CD	1:A:745:GLU:H	2.15	0.54
4:D:264:VAL:HA	4:D:267:LYS:HE2	1.89	0.54
1:A:1500:THR:HG21	1:A:1653:LYS:HE2	1.89	0.54
4:D:650:LEU:HD21	4:D:713:GLY:HA3	1.89	0.54
1:A:1642:ASP:OD2	1:A:1644:ARG:NH2	2.42	0.53
4:D:73:ARG:HG3	4:D:124:ILE:HD12	1.90	0.53
4:D:580:LEU:HD22	4:D:617:LEU:HD22	1.89	0.53
1:A:192:PHE:HB2	1:A:390:TYR:CE1	2.44	0.53
4:D:282:ARG:NH2	4:D:407:LEU:O	2.35	0.53
4:D:597:MET:HA	4:D:606:THR:HA	1.91	0.53
1:A:16:GLU:HG3	1:A:233:ALA:HB1	1.91	0.53
1:A:180:ARG:NH1	3:C:22:C:O2	2.36	0.53
1:A:317:PRO:O	1:A:340:ARG:NH2	2.26	0.53
1:A:1452:ARG:NE	1:A:1510:GLU:OE2	2.42	0.53
1:A:793:ASP:OD2	1:A:1023:GLY:N	2.41	0.53
1:A:984:HIS:ND1	1:A:1503:PRO:O	2.40	0.53
1:A:1598:LEU:HD12	1:A:1616:TRP:HZ3	1.73	0.53
1:A:604:LEU:HB3	1:A:607:ILE:HD11	1.90	0.53
1:A:1623:ASP:O	1:A:1627:LYS:HG2	2.08	0.53
1:A:188:ALA:O	3:C:19:U:O2'	2.27	0.53
1:A:1449:ASN:HB3	1:A:1450:PRO:HD3	1.90	0.52
1:A:940:ILE:HG13	1:A:1418:HIS:CE1	2.45	0.52
4:D:45:ILE:HD13	4:D:77:LEU:HB3	1.90	0.52
4:D:401:VAL:HG23	4:D:420:ASP:HA	1.90	0.52
1:A:1455:THR:O	1:A:1494:GLN:NE2	2.37	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:790:VAL:HG21	2:B:47:G:H21	1.74	0.52
4:D:120:ARG:HG3	4:D:120:ARG:HH11	1.75	0.52
4:D:213:LYS:HG3	4:D:250:LEU:HD11	1.92	0.52
1:A:954:ILE:HB	1:A:1522:VAL:HB	1.92	0.52
1:A:1544:ASN:HA	1:A:1634:LEU:HD23	1.92	0.52
4:D:200:LYS:HD2	4:D:204:PHE:HZ	1.74	0.52
2:B:42:U:H2'	2:B:43:U:C6	2.45	0.52
4:D:668:LYS:HG2	4:D:705:ARG:HB3	1.92	0.52
1:A:597:THR:O	1:A:724:SER:OG	2.19	0.51
4:D:82:ILE:HD13	4:D:92:LYS:HG2	1.92	0.51
4:D:321:LYS:HB2	4:D:521:ALA:HB2	1.92	0.51
1:A:30:LYS:HD3	1:A:36:MET:HE2	1.91	0.51
1:A:1002:GLU:HB3	1:A:1549:LEU:HD11	1.92	0.51
3:C:5:G:H2'	3:C:6:G:C8	2.43	0.51
4:D:679:ASN:HB3	4:D:682:TYR:CD2	2.45	0.51
1:A:596:SER:OG	2:B:38:C:OP1	2.23	0.51
1:A:1460:GLU:O	2:B:49:C:O2'	2.26	0.51
1:A:181:VAL:HA	1:A:188:ALA:HA	1.91	0.51
1:A:1450:PRO:HB2	1:A:1453:SER:HB3	1.93	0.51
1:A:278:GLU:OE2	1:A:369:ILE:HG12	2.11	0.51
2:B:19:G:H3'	2:B:20:A:H8	1.75	0.51
1:A:418:ILE:HD12	1:A:561:LEU:HD22	1.93	0.51
1:A:1428:ARG:NH2	1:A:1531:GLU:OE1	2.43	0.51
1:A:1513:ASP:O	1:A:1516:ASN:ND2	2.35	0.51
1:A:1616:TRP:HB3	1:A:1622:ILE:HG21	1.93	0.50
1:A:731:PHE:CZ	1:A:842:LEU:HD13	2.46	0.50
4:D:289:PHE:HE1	4:D:415:ASN:HD21	1.59	0.50
1:A:193:ARG:NH2	1:A:437:ASP:OD1	2.44	0.50
1:A:722:VAL:HB	1:A:772:LEU:HB2	1.93	0.50
4:D:513:ASN:OD1	4:D:513:ASN:N	2.44	0.50
4:D:661:MET:HG3	4:D:708:ALA:O	2.12	0.50
1:A:510:ARG:HA	2:B:32:G:OP2	2.11	0.50
1:A:1017:SER:HB3	1:A:1414:CYS:SG	2.52	0.50
4:D:127:VAL:HG11	4:D:165:TRP:HB3	1.92	0.50
4:D:607:TYR:CD2	4:D:648:ALA:HB2	2.47	0.50
1:A:959:GLU:HG3	1:A:1517:VAL:HG22	1.93	0.50
1:A:33:HIS:HB3	1:A:390:TYR:HB2	1.93	0.49
1:A:155:ASN:HD22	2:B:22:U:H1'	1.76	0.49
1:A:674:TYR:HB3	1:A:868:ILE:HD13	1.94	0.49
1:A:1499:GLU:HG3	1:A:1500:THR:HG23	1.94	0.49
1:A:1596:GLN:O	1:A:1600:GLU:HG3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:LYS:HD3	1:A:292:LYS:HD2	1.94	0.49
1:A:294:ARG:NE	3:C:14:U:O2'	2.44	0.49
1:A:403:ILE:HG23	4:D:442:THR:HA	1.93	0.49
4:D:450:TYR:HB3	4:D:453:ASN:HB2	1.95	0.49
4:D:513:ASN:HB2	4:D:514:HIS:CE1	2.46	0.49
4:D:553:ASN:OD1	4:D:554:GLN:N	2.46	0.49
1:A:312:GLU:HA	1:A:344:LYS:HE2	1.94	0.49
4:D:184:PHE:CE2	4:D:206:LYS:HD2	2.48	0.49
1:A:604:LEU:HD13	1:A:607:ILE:HD11	1.95	0.49
4:D:173:LEU:HA	4:D:176:LEU:HD12	1.93	0.49
1:A:176:ARG:NH2	1:A:196:GLU:OE2	2.45	0.49
4:D:550:LEU:HD23	4:D:674:GLU:HG2	1.94	0.49
4:D:657:VAL:N	4:D:712:ILE:O	2.42	0.49
1:A:998:ILE:HG21	1:A:1549:LEU:HD21	1.95	0.48
4:D:502:GLU:OE1	4:D:508:LYS:NZ	2.37	0.48
1:A:318:ASP:OD1	1:A:318:ASP:N	2.39	0.48
1:A:682:LEU:HD22	1:A:820:LEU:HD11	1.96	0.48
1:A:570:GLU:O	1:A:574:SER:OG	2.26	0.48
1:A:922:TYR:OH	1:A:1663:TYR:OH	2.19	0.48
1:A:108:ARG:HB3	2:B:29:C:H5'	1.96	0.48
1:A:942:HIS:NE2	1:A:1426:LYS:O	2.46	0.48
1:A:1014:LEU:HD12	1:A:1622:ILE:HG12	1.95	0.48
4:D:231:ARG:HG2	4:D:359:GLY:HA3	1.94	0.48
1:A:1470:ASP:OD1	1:A:1473:SER:OG	2.26	0.48
1:A:1480:PHE:HB2	1:A:1628:LEU:HD13	1.95	0.48
2:B:48:C:H2'	2:B:49:C:C6	2.48	0.48
4:D:167:ILE:HD12	4:D:167:ILE:H	1.79	0.48
4:D:508:LYS:HE2	4:D:513:ASN:HD21	1.78	0.48
1:A:458:PHE:HB2	1:A:548:ILE:HG23	1.96	0.48
1:A:775:ILE:HD13	1:A:821:LYS:HB2	1.95	0.48
1:A:910:SER:HB2	1:A:911:PRO:HD3	1.95	0.48
4:D:141:ASP:H	4:D:144:LYS:HE3	1.79	0.48
4:D:310:ILE:CG2	4:D:426:TRP:HB3	2.43	0.48
1:A:151:ILE:HD13	1:A:211:LEU:HD13	1.95	0.48
1:A:369:ILE:HA	1:A:372:LYS:HE3	1.94	0.48
1:A:1459:LEU:HD22	2:B:49:C:H2'	1.96	0.48
1:A:179:ASN:HB3	2:B:34:U:H3'	1.95	0.48
1:A:418:ILE:HD11	1:A:568:PHE:HZ	1.78	0.48
4:D:85:LYS:HA	4:D:328:LEU:HB3	1.96	0.48
1:A:323:ARG:HH22	1:A:325:LYS:HZ3	1.61	0.48
4:D:33:ASN:O	4:D:41:ARG:NH2	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:238:GLY:HA2	4:D:241:GLU:HB2	1.95	0.48
1:A:132:LEU:HD21	1:A:220:LEU:HB2	1.96	0.47
1:A:533:LYS:HE2	2:B:42:U:O2'	2.14	0.47
4:D:63:LEU:HG	4:D:67:SER:HB3	1.96	0.47
1:A:803:ALA:HA	3:C:9:G:H21	1.79	0.47
1:A:155:ASN:ND2	2:B:22:U:H1'	2.30	0.47
2:B:30:A:H4'	2:B:31:C:H5''	1.96	0.47
4:D:675:TRP:CZ2	4:D:682:TYR:HB3	2.49	0.47
4:D:397:PRO:HD2	4:D:714:PHE:HZ	1.78	0.47
4:D:579:LEU:HD23	4:D:680:ILE:CD1	2.41	0.47
4:D:605:ILE:HG12	4:D:610:ILE:HG13	1.95	0.47
1:A:20:LEU:HB3	1:A:195:TRP:HB2	1.95	0.47
1:A:276:GLN:O	1:A:280:ILE:HG13	2.15	0.47
1:A:950:LEU:HD13	1:A:1532:LEU:HD23	1.96	0.47
4:D:126:SER:HB2	4:D:149:LEU:HD22	1.96	0.47
4:D:171:ILE:HA	4:D:174:CYS:SG	2.54	0.47
4:D:317:SER:O	4:D:321:LYS:HG3	2.15	0.47
4:D:405:PHE:HB2	4:D:488:PRO:HA	1.96	0.47
4:D:589:ASN:O	4:D:642:HIS:NE2	2.45	0.47
4:D:701:PRO:HA	4:D:704:TYR:HB3	1.96	0.47
1:A:21:VAL:HG11	1:A:38:GLY:O	2.14	0.47
1:A:58:ILE:HG23	1:A:62:LEU:HD23	1.97	0.47
1:A:805:PHE:HB3	3:C:9:G:C4	2.50	0.47
1:A:1010:HIS:CE1	1:A:1538:SER:HA	2.49	0.47
4:D:429:LYS:HG3	4:D:477:PHE:CD2	2.50	0.47
1:A:272:GLU:OE1	1:A:350:ASN:ND2	2.48	0.47
1:A:733:THR:O	1:A:737:ARG:HG3	2.15	0.47
1:A:776:ASN:HB3	1:A:780:PHE:HE1	1.80	0.47
4:D:304:ASN:O	4:D:308:ALA:N	2.47	0.47
1:A:517:SER:HB2	1:A:557:PHE:HB3	1.97	0.47
4:D:447:TYR:CE1	4:D:662:TRP:HB2	2.50	0.47
4:D:546:HIS:O	4:D:670:LYS:NZ	2.31	0.47
1:A:732:ARG:CZ	2:B:39:A:H5''	2.45	0.46
4:D:402:VAL:HG13	4:D:485:LEU:HD23	1.97	0.46
1:A:168:ARG:HG3	1:A:168:ARG:HH11	1.80	0.46
4:D:94:LYS:HB2	4:D:94:LYS:HE3	1.73	0.46
4:D:188:ILE:HG13	4:D:189:LYS:N	2.29	0.46
4:D:224:LYS:HB2	4:D:224:LYS:HE2	1.67	0.46
4:D:628:GLU:O	4:D:632:VAL:HG22	2.15	0.46
1:A:817:PRO:HG2	1:A:819:LYS:HE3	1.96	0.46
1:A:1006:MET:HE1	1:A:1562:ILE:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:30:GLU:O	4:D:36:THR:OG1	2.33	0.46
4:D:122:PHE:HE2	4:D:151:GLU:HB2	1.79	0.46
4:D:700:PRO:N	4:D:701:PRO:HD2	2.30	0.46
1:A:11:GLU:HB2	1:A:240:ILE:HD13	1.98	0.46
1:A:1548:LYS:HB3	1:A:1553:LYS:HG2	1.96	0.46
1:A:347:TRP:CZ2	1:A:351:LYS:HD3	2.51	0.46
1:A:471:PRO:HD2	1:A:474:ALA:HB3	1.97	0.46
1:A:578:TYR:CZ	1:A:588:MET:HG2	2.50	0.46
1:A:1553:LYS:HE2	2:B:49:C:OP1	2.16	0.46
4:D:65:SER:O	4:D:69:ILE:N	2.45	0.46
4:D:84:PRO:HA	4:D:90:PRO:HG3	1.97	0.46
4:D:196:LEU:HB2	4:D:219:CYS:SG	2.56	0.46
4:D:404:HIS:N	4:D:417:ILE:O	2.44	0.46
1:A:180:ARG:O	1:A:189:LYS:N	2.45	0.46
1:A:1009:THR:HG21	2:B:46:U:C5	2.51	0.46
4:D:485:LEU:HD13	4:D:714:PHE:HB3	1.97	0.46
1:A:976:LYS:HA	1:A:976:LYS:HD3	1.63	0.46
4:D:350:ASP:O	4:D:354:THR:HG23	2.16	0.46
1:A:972:GLU:HG3	1:A:985:LYS:HG2	1.98	0.46
1:A:371:LYS:HD2	1:A:371:LYS:HA	1.70	0.46
1:A:795:PHE:CE1	1:A:1476:PRO:HD2	2.51	0.46
1:A:1582:VAL:HG12	1:A:1587:ILE:HG13	1.98	0.46
4:D:20:ASP:CG	4:D:55:LEU:HD12	2.41	0.46
4:D:146:LYS:HA	4:D:176:LEU:HD13	1.97	0.46
4:D:228:CYS:HA	4:D:233:TRP:CD1	2.51	0.46
4:D:224:LYS:HG2	4:D:261:SER:HA	1.97	0.45
1:A:58:ILE:HB	1:A:156:PHE:HB2	1.98	0.45
1:A:200:GLU:OE1	1:A:467:ARG:NH2	2.49	0.45
1:A:187:LYS:HE2	1:A:187:LYS:HB3	1.60	0.45
1:A:1005:GLY:O	1:A:1009:THR:HG23	2.16	0.45
4:D:171:ILE:HD12	4:D:197:GLN:HB3	1.97	0.45
1:A:771:ASP:OD1	1:A:771:ASP:N	2.50	0.45
1:A:966:ILE:O	1:A:990:PHE:N	2.47	0.45
1:A:1027:THR:HG22	1:A:1394:THR:HG22	1.99	0.45
4:D:429:LYS:HG3	4:D:477:PHE:HD2	1.81	0.45
1:A:188:ALA:HB2	2:B:36:C:H1'	1.98	0.45
1:A:1469:PRO:HB2	1:A:1674:TYR:CD2	2.52	0.45
1:A:957:LYS:NZ	1:A:1565:ASP:OD2	2.40	0.45
1:A:1022:PHE:CZ	1:A:1395:LEU:HB3	2.51	0.45
1:A:1022:PHE:HD2	1:A:1024:GLU:HG3	1.82	0.45
1:A:1395:LEU:HD21	1:A:1464:PRO:HG2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1530:TRP:HB2	1:A:1594:GLY:HA2	1.98	0.45
1:A:1661:TYR:HE1	1:A:1688:TRP:HB3	1.81	0.45
4:D:94:LYS:HG2	4:D:132:PHE:HE1	1.82	0.45
4:D:228:CYS:HB3	4:D:278:TRP:NE1	2.31	0.45
1:A:415:GLU:OE1	1:A:564:ARG:NH1	2.50	0.45
1:A:612:TRP:CE2	1:A:646:LEU:HD13	2.52	0.45
1:A:1483:HIS:HE1	1:A:1546:ALA:C	2.24	0.45
4:D:600:LEU:H	4:D:604:GLY:HA2	1.82	0.45
1:A:632:LYS:O	1:A:636:GLU:HG2	2.17	0.45
1:A:939:ILE:HG13	1:A:1409:LEU:HB3	1.99	0.45
4:D:43:LYS:HE3	4:D:43:LYS:HB2	1.70	0.45
4:D:273:ARG:NH2	4:D:523:SER:HB2	2.32	0.45
4:D:665:ARG:HD2	4:D:667:ASN:HB2	1.98	0.45
1:A:20:LEU:N	1:A:195:TRP:O	2.44	0.45
1:A:20:LEU:HD12	1:A:42:ALA:HB3	1.98	0.44
1:A:130:CYS:SG	1:A:135:ARG:HD3	2.57	0.44
1:A:507:ILE:HD11	4:D:448:ASN:HB3	1.98	0.44
1:A:517:SER:HB3	1:A:558:PRO:O	2.17	0.44
1:A:577:ARG:NH1	1:A:580:GLU:OE1	2.38	0.44
1:A:585:LYS:HD2	1:A:655:ARG:CD	2.46	0.44
1:A:737:ARG:HD2	1:A:942:HIS:O	2.17	0.44
4:D:71:ILE:HG21	4:D:103:PHE:CD2	2.51	0.44
1:A:413:THR:N	1:A:615:ASN:OD1	2.49	0.44
1:A:793:ASP:HB3	1:A:796:THR:O	2.17	0.44
4:D:52:LEU:HD11	4:D:70:SER:OG	2.16	0.44
4:D:278:TRP:CE2	4:D:281:HIS:HB2	2.52	0.44
4:D:632:VAL:HG13	4:D:644:SER:HB3	1.99	0.44
1:A:64:ARG:HG3	1:A:153:PHE:CD1	2.51	0.44
1:A:428:PHE:CZ	1:A:551:GLY:HA3	2.53	0.44
1:A:952:GLY:HA2	1:A:1570:PHE:CE2	2.53	0.44
4:D:80:ILE:HG22	4:D:357:GLN:HG2	1.99	0.44
4:D:556:TRP:CH2	4:D:677:LEU:HD21	2.51	0.44
1:A:303:LEU:HD22	1:A:310:VAL:HG11	1.98	0.44
1:A:412:GLU:HG2	1:A:414:LYS:HE2	2.00	0.44
1:A:67:VAL:O	1:A:71:VAL:HG23	2.17	0.44
1:A:1448:ASN:ND2	1:A:1453:SER:O	2.46	0.44
4:D:565:ALA:HB3	4:D:600:LEU:HD22	2.00	0.44
4:D:579:LEU:HD22	4:D:621:GLN:O	2.18	0.44
1:A:47:LYS:HG3	1:A:55:LYS:HG3	2.00	0.44
1:A:989:PHE:HD2	1:A:996:LEU:HD22	1.81	0.44
4:D:411:GLU:O	4:D:413:MET:N	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:677:GLU:O	1:A:864:ALA:N	2.49	0.44
1:A:1582:VAL:HG13	1:A:1586:ASP:HB3	2.00	0.44
4:D:71:ILE:HD13	4:D:71:ILE:HA	1.87	0.44
4:D:413:MET:HE1	4:D:432:GLN:NE2	2.32	0.44
1:A:200:GLU:OE1	1:A:464:LYS:NZ	2.50	0.44
1:A:281:VAL:HG13	1:A:293:ILE:HG13	2.00	0.44
4:D:405:PHE:CE1	4:D:469:MET:HE3	2.53	0.44
4:D:643:TYR:HB3	4:D:647:THR:CG2	2.48	0.44
1:A:25:ASP:O	1:A:29:ARG:N	2.50	0.44
1:A:922:TYR:OH	2:B:51:C:OP2	2.29	0.44
1:A:923:PRO:HG2	1:A:1645:TYR:CD2	2.53	0.44
4:D:463:LYS:HG2	4:D:501:ILE:HB	1.99	0.44
4:D:485:LEU:HD22	4:D:714:PHE:CD2	2.52	0.44
4:D:213:LYS:HA	4:D:254:PHE:CZ	2.53	0.43
4:D:241:GLU:OE2	4:D:283:VAL:HG11	2.18	0.43
4:D:555:ILE:HG21	4:D:677:LEU:HD22	2.00	0.43
1:A:179:ASN:O	2:B:33:G:O2'	2.22	0.43
1:A:364:SER:O	1:A:368:LEU:HG	2.18	0.43
1:A:693:GLU:OE2	1:A:703:LYS:NZ	2.47	0.43
1:A:528:ARG:NH2	1:A:771:ASP:OD2	2.50	0.43
1:A:908:HIS:ND1	1:A:911:PRO:HD3	2.34	0.43
1:A:31:LYS:HA	1:A:31:LYS:HD3	1.94	0.43
1:A:98:LYS:HB2	1:A:98:LYS:HE2	1.63	0.43
4:D:120:ARG:HG3	4:D:120:ARG:NH1	2.33	0.43
4:D:213:LYS:HA	4:D:254:PHE:HZ	1.83	0.43
4:D:260:ARG:O	4:D:264:VAL:HG23	2.18	0.43
4:D:541:ASN:HB2	4:D:600:LEU:HD23	2.01	0.43
1:A:69:LYS:O	1:A:73:GLU:HG3	2.19	0.43
2:B:47:G:H1'	2:B:48:C:C6	2.53	0.43
4:D:543:ASP:HB3	4:D:546:HIS:ND1	2.33	0.43
1:A:303:LEU:HD23	1:A:303:LEU:HA	1.84	0.43
1:A:532:ALA:HB2	2:B:39:A:H1'	2.00	0.43
1:A:728:ARG:HB2	1:A:769:PHE:HB2	2.01	0.43
4:D:160:PHE:O	4:D:190:SER:OG	2.31	0.43
4:D:471:PHE:HB2	4:D:477:PHE:CD2	2.54	0.43
1:A:1545:ILE:HG12	1:A:1630:TRP:HA	2.00	0.43
1:A:545:LEU:HD22	3:C:13:A:C2	2.53	0.43
4:D:80:ILE:HG23	4:D:360:TYR:CD2	2.53	0.43
3:C:10:A:H2'	3:C:11:A:H8	1.83	0.43
4:D:28:LEU:O	4:D:32:PHE:N	2.47	0.43
4:D:699:HIS:CD2	4:D:701:PRO:HG2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1018:CYS:SG	1:A:1405:PRO:HA	2.59	0.42
2:B:43:U:H2'	2:B:44:U:C6	2.54	0.42
4:D:328:LEU:O	4:D:332:GLU:HG2	2.19	0.42
1:A:683:LEU:HD22	1:A:807:THR:HB	2.01	0.42
1:A:176:ARG:NE	1:A:196:GLU:OE1	2.44	0.42
1:A:304:ARG:HB3	1:A:359:THR:OG1	2.19	0.42
1:A:480:ARG:HB2	1:A:509:MET:SD	2.59	0.42
1:A:493:VAL:HG11	2:B:33:G:H5'	2.01	0.42
4:D:541:ASN:ND2	4:D:541:ASN:O	2.52	0.42
4:D:556:TRP:HB2	4:D:562:LYS:HE2	2.01	0.42
1:A:1003:LEU:HD11	1:A:1434:ALA:HB2	2.01	0.42
1:A:1461:SER:HB3	3:C:5:G:H21	1.84	0.42
4:D:501:ILE:HD12	4:D:501:ILE:N	2.34	0.42
1:A:503:CYS:SG	1:A:506:CYS:HB2	2.59	0.42
1:A:660:PRO:O	1:A:664:ASN:HB2	2.19	0.42
1:A:1019:PHE:O	1:A:1405:PRO:HG3	2.19	0.42
4:D:97:LEU:HD23	4:D:132:PHE:HA	2.01	0.42
4:D:690:GLU:HG3	4:D:691:THR:HG23	2.00	0.42
4:D:71:ILE:HG21	4:D:103:PHE:CG	2.55	0.42
4:D:590:MET:HA	4:D:633:PRO:HG3	2.01	0.42
1:A:458:PHE:HD2	1:A:548:ILE:HG12	1.85	0.42
1:A:935:ARG:NH2	1:A:1404:HIS:O	2.53	0.42
1:A:1483:HIS:ND1	1:A:1631:VAL:HG22	2.35	0.42
1:A:64:ARG:HG3	1:A:153:PHE:CE1	2.55	0.42
1:A:301:ARG:NH2	1:A:529:TYR:OH	2.52	0.42
1:A:950:LEU:HD22	1:A:1569:THR:HG21	2.01	0.42
1:A:842:LEU:HD12	1:A:842:LEU:HA	1.90	0.42
1:A:992:ILE:O	1:A:995:GLU:HG2	2.20	0.42
1:A:1022:PHE:CD2	1:A:1024:GLU:HG3	2.55	0.41
1:A:1512:MET:HE2	1:A:1518:PHE:CZ	2.55	0.41
4:D:331:MET:H	4:D:331:MET:HG2	1.62	0.41
1:A:330:TRP:NE1	1:A:691:LEU:O	2.42	0.41
1:A:578:TYR:OH	1:A:588:MET:HG2	2.20	0.41
1:A:656:GLU:H	1:A:656:GLU:CD	2.28	0.41
1:A:1015:THR:HG22	1:A:1534:LEU:HD13	2.01	0.41
4:D:120:ARG:O	4:D:124:ILE:HG22	2.20	0.41
4:D:161:ASP:O	4:D:192:LYS:NZ	2.43	0.41
1:A:173:ALA:HA	1:A:197:ALA:HA	2.01	0.41
4:D:89:ILE:HG13	4:D:90:PRO:HD2	2.02	0.41
1:A:283:ALA:HB1	1:A:334:VAL:HA	2.01	0.41
1:A:919:ASN:HB2	1:A:1687:PRO:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:327:ARG:O	4:D:331:MET:HG2	2.21	0.41
4:D:415:ASN:HB3	4:D:430:GLU:HG2	2.02	0.41
1:A:106:ARG:NH1	1:A:110:THR:OG1	2.46	0.41
1:A:588:MET:HE2	1:A:588:MET:HB2	1.64	0.41
3:C:3:G:H2'	3:C:4:G:C8	2.55	0.41
1:A:530:ARG:HB2	1:A:546:PHE:CZ	2.56	0.41
1:A:914:ASP:OD2	1:A:918:LYS:N	2.54	0.41
3:C:2:C:H4'	3:C:3:G:OP1	2.19	0.41
4:D:537:VAL:HG13	4:D:560:SER:HB3	2.03	0.41
1:A:234:LEU:HD12	1:A:515:LYS:HG3	2.03	0.41
1:A:611:GLU:N	1:A:647:ILE:O	2.52	0.41
1:A:1462:PRO:HG3	2:B:49:C:O3'	2.21	0.41
4:D:484:VAL:HG21	4:D:510:PHE:HZ	1.85	0.41
1:A:737:ARG:HG2	1:A:742:LEU:HD22	2.01	0.41
1:A:72:GLU:HG3	1:A:129:PHE:HZ	1.86	0.41
1:A:293:ILE:HG22	1:A:370:TYR:HB2	2.03	0.41
1:A:417:ILE:CG2	1:A:610:PHE:HB2	2.51	0.41
1:A:530:ARG:CD	1:A:725:GLU:HG2	2.50	0.41
1:A:622:ILE:O	1:A:625:ARG:NH1	2.54	0.41
2:B:35:A:H1'	2:B:36:C:C6	2.56	0.41
4:D:147:ASN:O	4:D:151:GLU:HG3	2.20	0.41
4:D:461:LEU:HD11	4:D:495:LEU:HD22	2.02	0.41
4:D:573:ILE:O	4:D:616:GLY:HA3	2.21	0.41
4:D:590:MET:HA	4:D:633:PRO:CG	2.51	0.41
4:D:668:LYS:HA	4:D:668:LYS:HD2	1.81	0.41
1:A:22:GLU:CD	1:A:46:TRP:HE1	2.28	0.41
1:A:272:GLU:OE2	1:A:349:SER:OG	2.39	0.41
1:A:671:GLU:HB3	1:A:825:TRP:CZ3	2.56	0.41
1:A:754:MET:HE2	1:A:754:MET:HB3	1.90	0.41
1:A:789:HIS:HA	2:B:47:G:O5'	2.20	0.41
4:D:80:ILE:HG23	4:D:360:TYR:HD2	1.86	0.41
4:D:254:PHE:O	4:D:258:THR:HG22	2.21	0.41
1:A:381:PHE:HB3	4:D:57:LYS:HD3	2.02	0.40
1:A:559:PHE:CD2	1:A:604:LEU:HD21	2.55	0.40
1:A:767:ILE:HD11	1:A:824:PHE:CD1	2.56	0.40
1:A:924:HIS:HB2	1:A:1479:LYS:HG3	2.03	0.40
1:A:1646:PRO:HD2	1:A:1663:TYR:CZ	2.56	0.40
1:A:1673:PRO:HG2	1:A:1676:THR:HG22	2.02	0.40
4:D:97:LEU:HD21	4:D:131:GLU:HB2	2.03	0.40
4:D:186:ASN:OD1	4:D:186:ASN:N	2.53	0.40
4:D:615:LYS:HG3	4:D:618:LYS:NZ	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:TRP:CE3	1:A:563:TYR:HD2	2.40	0.40
1:A:704:LYS:NZ	1:A:773:GLU:OE2	2.54	0.40
1:A:727:HIS:ND1	1:A:770:GLU:O	2.54	0.40
1:A:1646:PRO:HG3	1:A:1661:TYR:CE1	2.56	0.40
4:D:396:VAL:O	4:D:421:LYS:HE3	2.21	0.40
4:D:511:LEU:HD21	4:D:688:GLN:HB2	2.03	0.40
1:A:1502:GLN:HB2	1:A:1505:ARG:HB2	2.04	0.40
1:A:1674:TYR:O	1:A:1678:VAL:HG23	2.21	0.40
4:D:93:LYS:NZ	4:D:356:TYR:OH	2.32	0.40
4:D:223:LEU:HD22	4:D:233:TRP:CE2	2.56	0.40
4:D:494:ARG:NH2	4:D:640:ASP:OD1	2.48	0.40
1:A:59:THR:HA	1:A:155:ASN:OD1	2.22	0.40
1:A:437:ASP:OD1	1:A:437:ASP:C	2.64	0.40
1:A:486:ALA:HB2	1:A:665:LEU:HD11	2.03	0.40
1:A:663:GLU:O	1:A:828:LYS:NZ	2.53	0.40
1:A:1007:LEU:HA	1:A:1007:LEU:HD22	1.88	0.40
1:A:1409:LEU:HD21	1:A:1425:TYR:HB2	2.02	0.40
1:A:1467:SER:HB3	1:A:1476:PRO:HG2	2.03	0.40
1:A:1648:LEU:HD13	2:B:51:C:C2	2.57	0.40
4:D:714:PHE:HA	4:D:715:PRO:HD3	1.83	0.40
1:A:45:THR:OG1	2:B:22:U:OP1	2.35	0.40
1:A:461:LEU:CD2	1:A:471:PRO:HG3	2.49	0.40
1:A:563:TYR:CZ	1:A:569:PRO:HD3	2.57	0.40
1:A:668:GLN:O	1:A:828:LYS:N	2.54	0.40
4:D:89:ILE:HG12	4:D:135:ILE:HB	2.03	0.40
4:D:579:LEU:HD13	4:D:580:LEU:N	2.36	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	
1	A	1212/1728 (70%)	1189 (98%)	23 (2%)	0	100	100
4	D	638/746 (86%)	619 (97%)	19 (3%)	0	100	100
All	All	1850/2474 (75%)	1808 (98%)	42 (2%)	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	
1	A	1073/1531 (70%)	1036 (97%)	37 (3%)	32	55
4	D	580/657 (88%)	553 (95%)	27 (5%)	23	43
All	All	1653/2188 (76%)	1589 (96%)	64 (4%)	30	50

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	VAL
1	A	13	THR
1	A	27	ASP
1	A	85	VAL
1	A	97	SER
1	A	105	LEU
1	A	290	LEU
1	A	311	ILE
1	A	318	ASP
1	A	391	SER
1	A	419	VAL
1	A	423	LYS
1	A	513	THR
1	A	515	LYS
1	A	517	SER
1	A	522	SER
1	A	556	THR
1	A	640	SER

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Mol	Chain	Res	Type
1	A	670	SER
1	A	765	SER
1	A	792	ILE
1	A	796	THR
1	A	830	PHE
1	A	836	LEU
1	A	842	LEU
1	A	951	SER
1	A	1001	SER
1	A	1007	LEU
1	A	1422	THR
1	A	1426	LYS
1	A	1454	LEU
1	A	1458	VAL
1	A	1467	SER
1	A	1511	VAL
1	A	1531	GLU
1	A	1595	TYR
1	A	1613	LEU
4	D	18	ASP
4	D	67	SER
4	D	128	LEU
4	D	145	LEU
4	D	181	GLU
4	D	186	ASN
4	D	196	LEU
4	D	202	TYR
4	D	204	PHE
4	D	279	SER
4	D	283	VAL
4	D	319	SER
4	D	478	ILE
4	D	511	LEU
4	D	513	ASN
4	D	523	SER
4	D	542	PHE
4	D	576	ASN
4	D	615	LYS
4	D	617	LEU
4	D	647	THR
4	D	688	GLN
4	D	702	THR

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Mol	Chain	Res	Type
4	D	706	SER
4	D	707	ILE
4	D	711	SER
4	D	714	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	109	HIS
1	A	270	ASN
1	A	582	ASN
1	A	673	GLN
1	A	1010	HIS
1	A	1483	HIS
1	A	1635	ASN
1	A	1671	ASN
4	D	197	GLN
4	D	215	HIS
4	D	295	GLN
4	D	551	GLN
4	D	572	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	35/110 (31%)	17 (48%)	0
3	C	22/56 (39%)	8 (36%)	1 (4%)
All	All	57/166 (34%)	25 (43%)	1 (1%)

All (25) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	20	A
2	B	22	U
2	B	25	A
2	B	27	G
2	B	28	U
2	B	32	G
2	B	33	G
2	B	35	A

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Mol	Chain	Res	Type
2	B	38	C
2	B	39	A
2	B	40	A
2	B	46	U
2	B	47	G
2	B	49	C
2	B	50	C
2	B	51	C
2	B	52	G
3	C	2	C
3	C	3	G
3	C	5	G
3	C	9	G
3	C	15	U
3	C	17	G
3	C	21	C
3	C	22	C

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	C	2	C

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](#)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

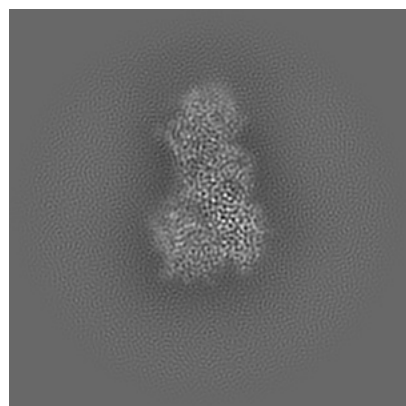
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-33681. These allow visual inspection of the internal detail of the map and identification of artifacts.

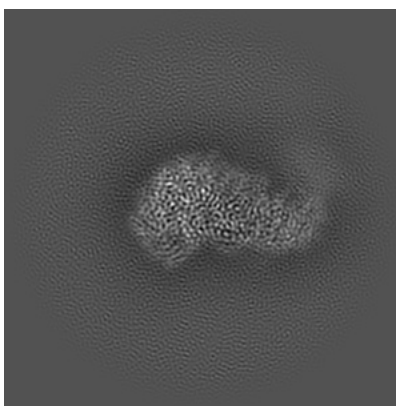
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

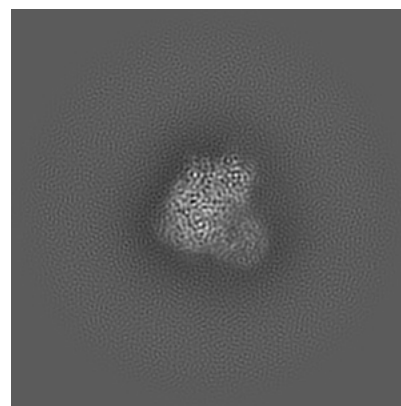
6.1.1 Primary map



X

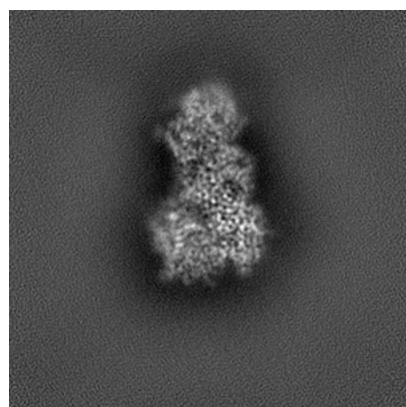


Y

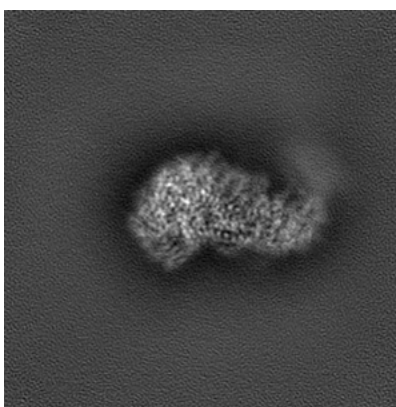


Z

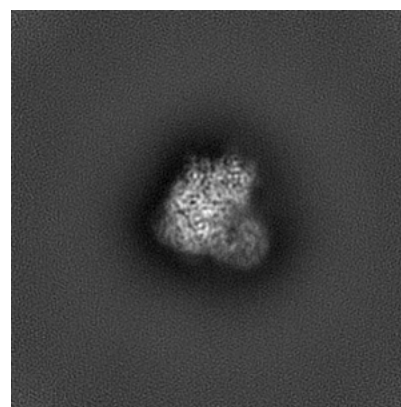
6.1.2 Raw map



X



Y

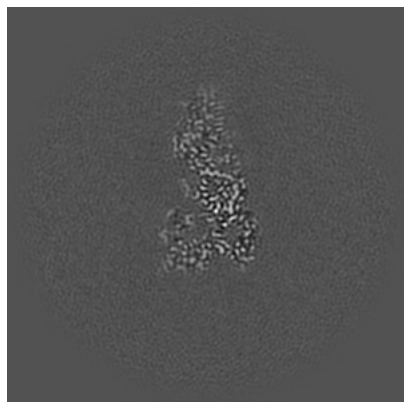


Z

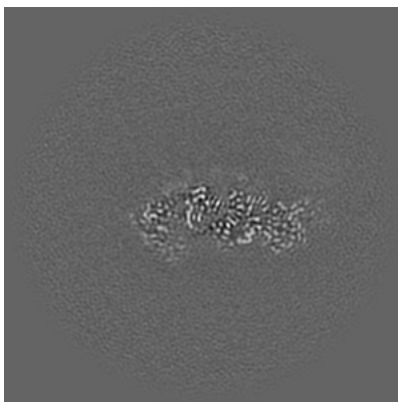
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

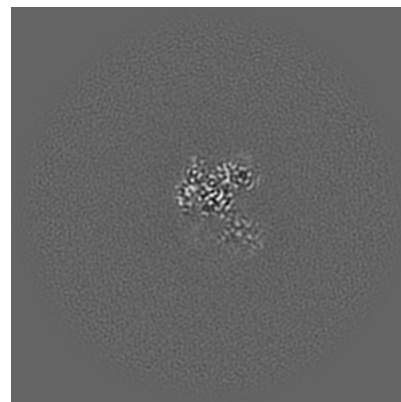
6.2.1 Primary map



X Index: 128

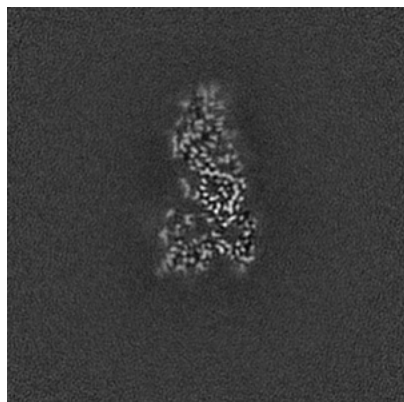


Y Index: 128

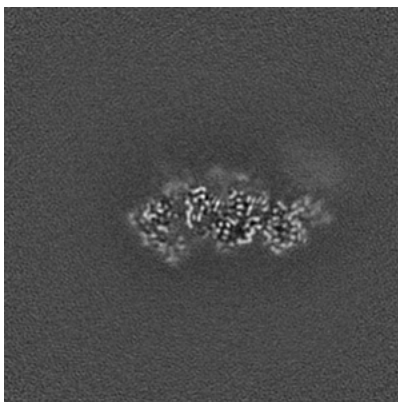


Z Index: 128

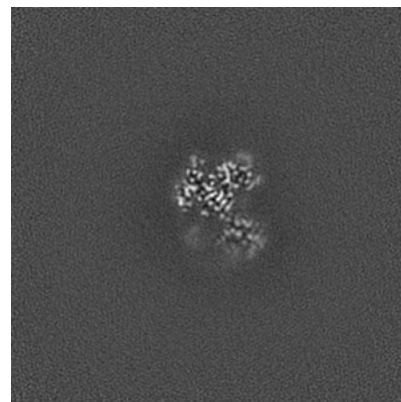
6.2.2 Raw map



X Index: 128



Y Index: 128

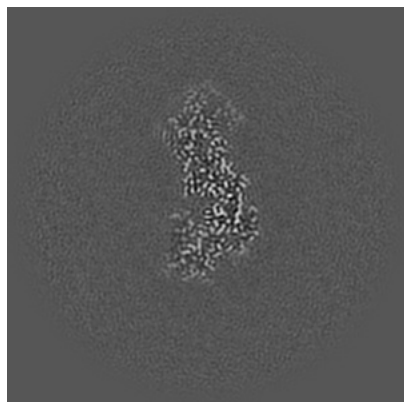


Z Index: 128

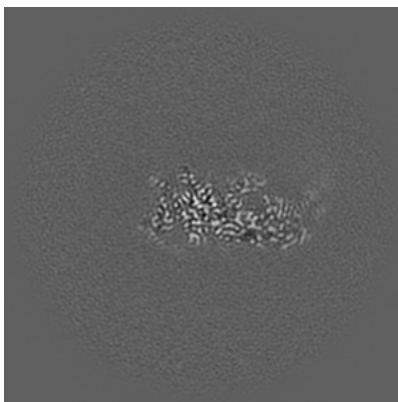
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

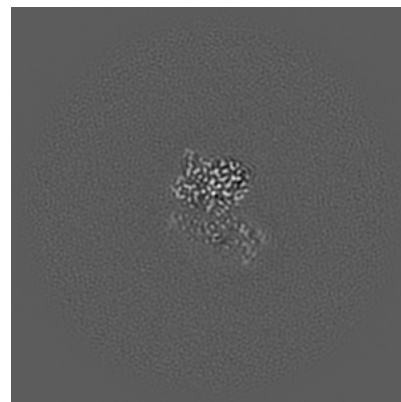
6.3.1 Primary map



X Index: 121

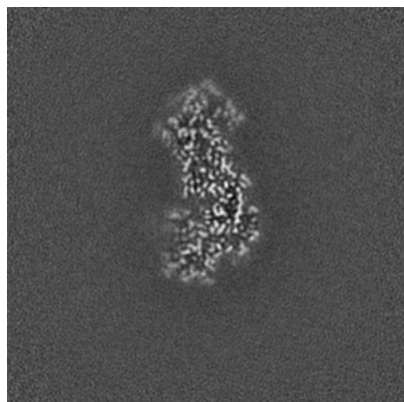


Y Index: 134

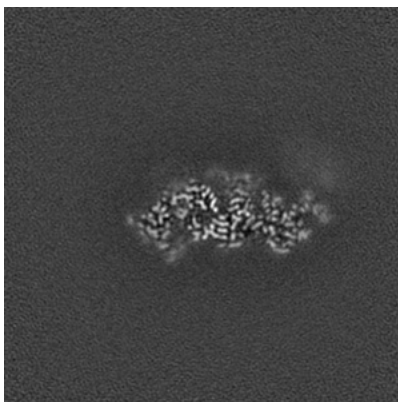


Z Index: 120

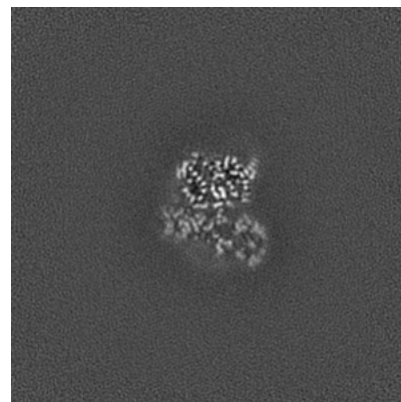
6.3.2 Raw map



X Index: 121



Y Index: 130

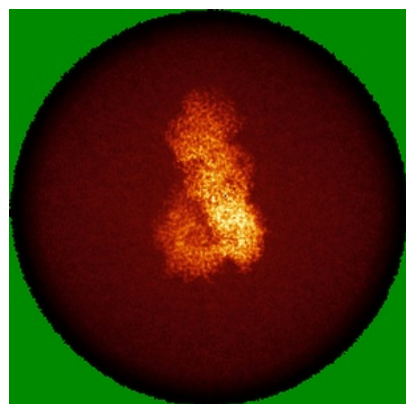


Z Index: 117

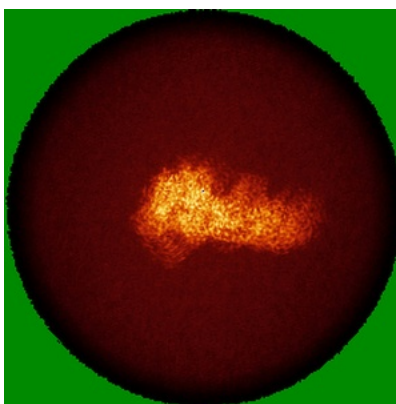
The images above show the largest variance slices of the map in three orthogonal directions.

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

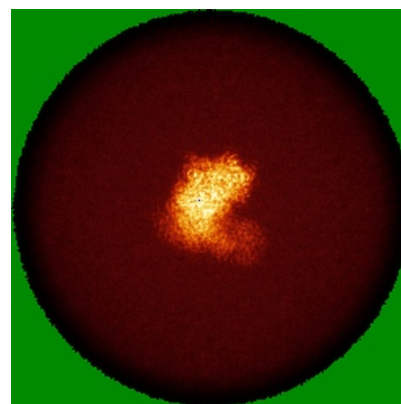
6.4.1 Primary map



X

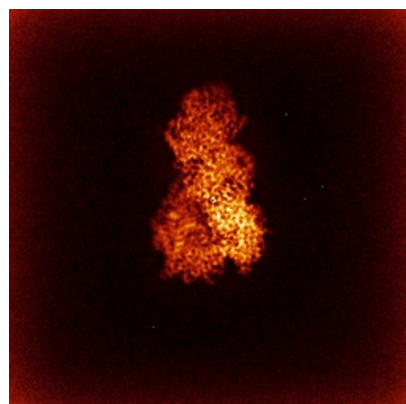


Y

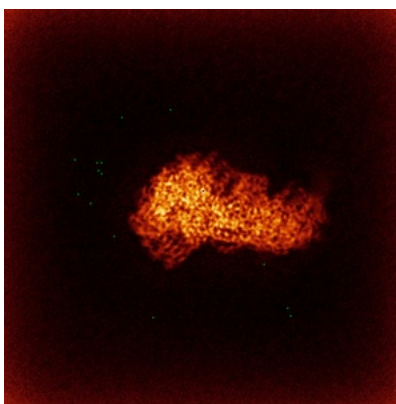


Z

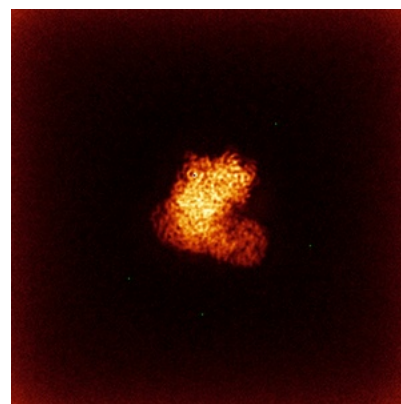
6.4.2 Raw map



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](#)

This section was not generated.

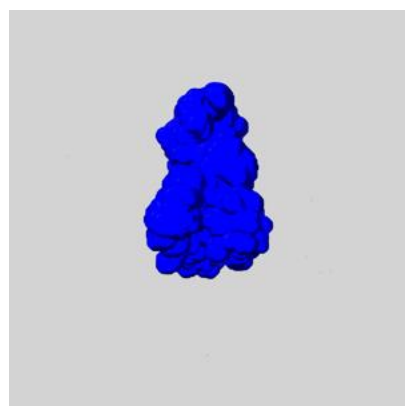
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

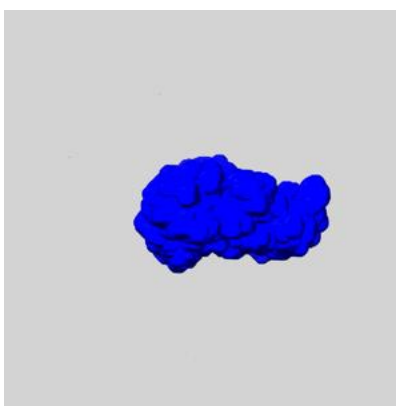
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

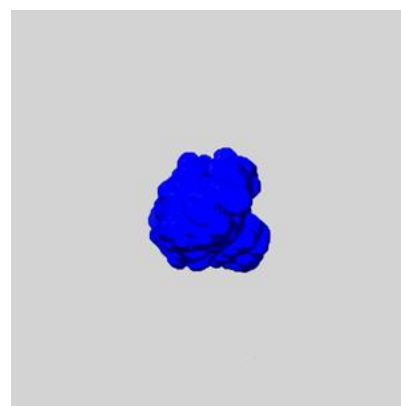
6.6.1 emd_33681_msk_1.map [i](#)



X



Y

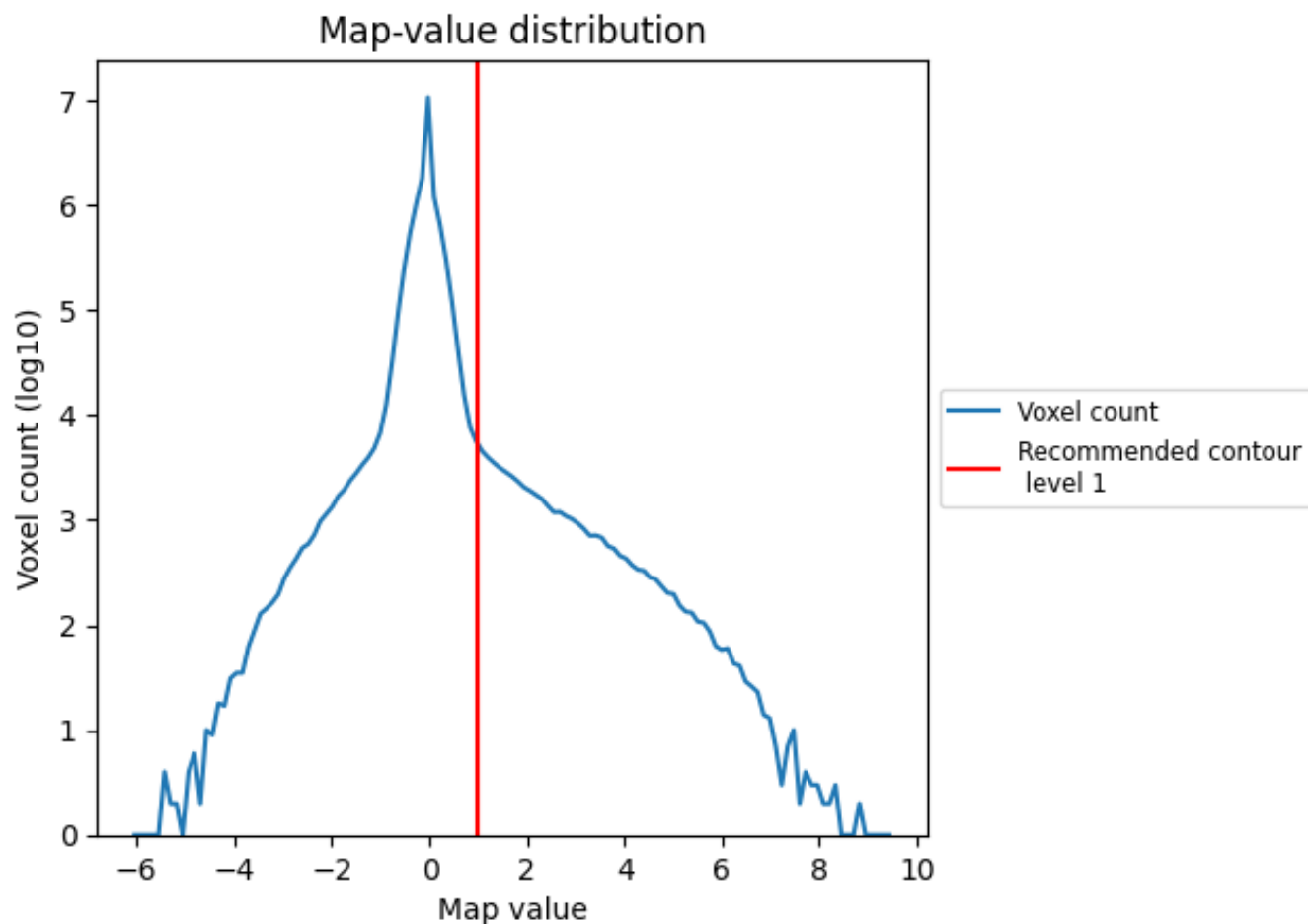


Z

7 Map analysis [i](#)

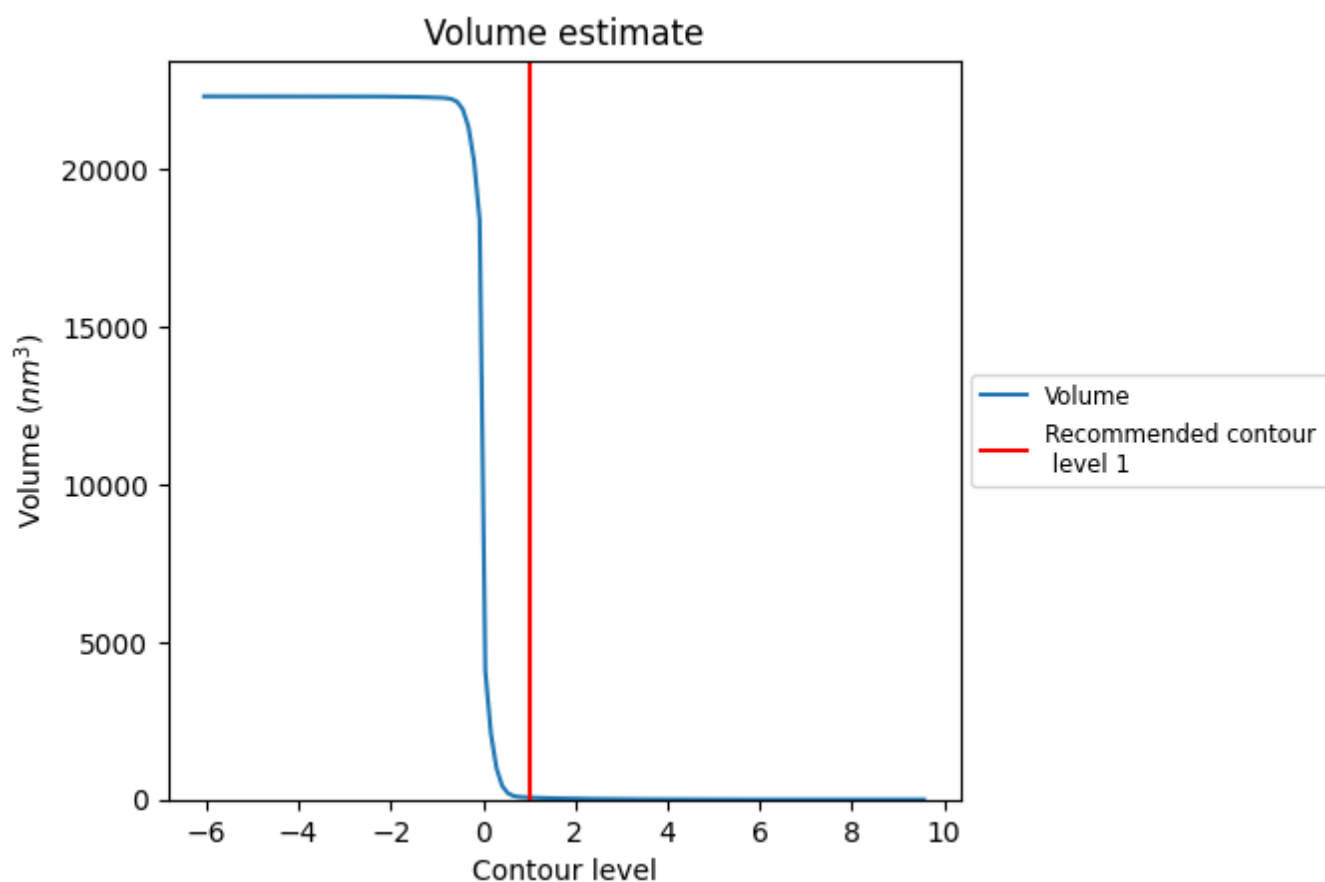
This section contains the results of statistical analysis of the map.

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.

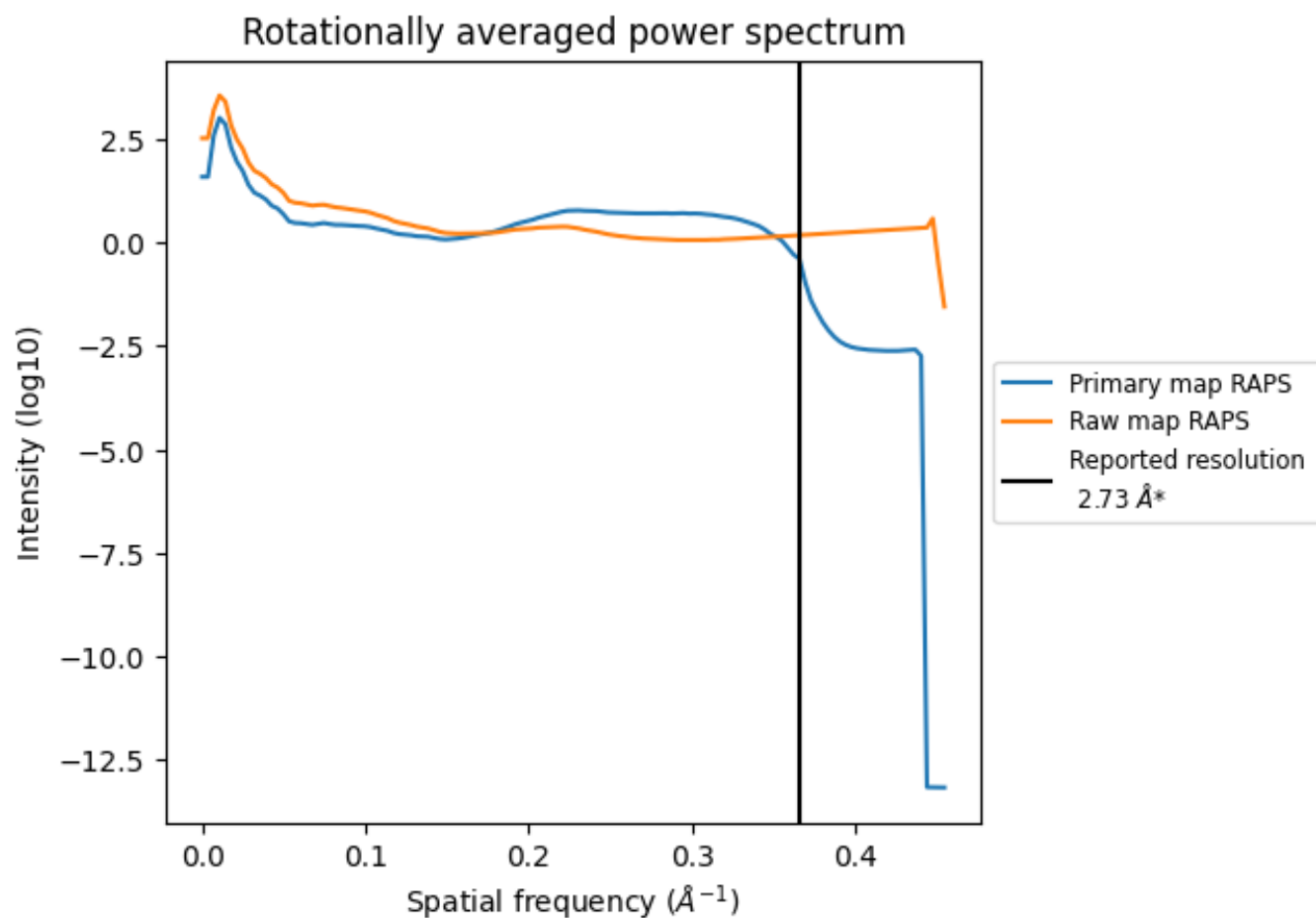
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 65 nm^3 ; this corresponds to an approximate mass of 59 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 ⓘ

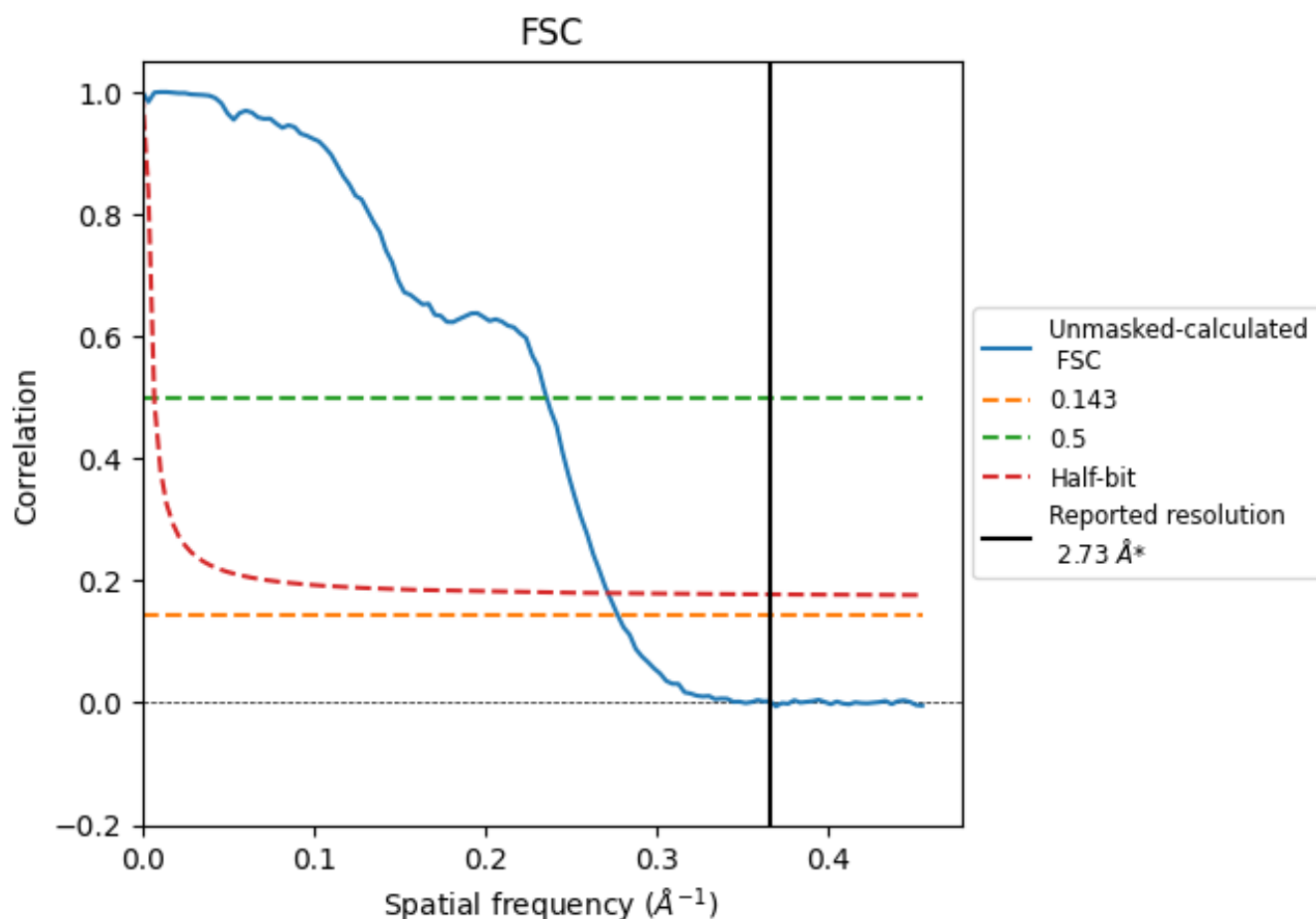


*Reported resolution corresponds to spatial frequency of 0.366 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.366 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.73	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.61	4.24	3.68

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.61 differs from the reported value 2.73 by more than 10 %

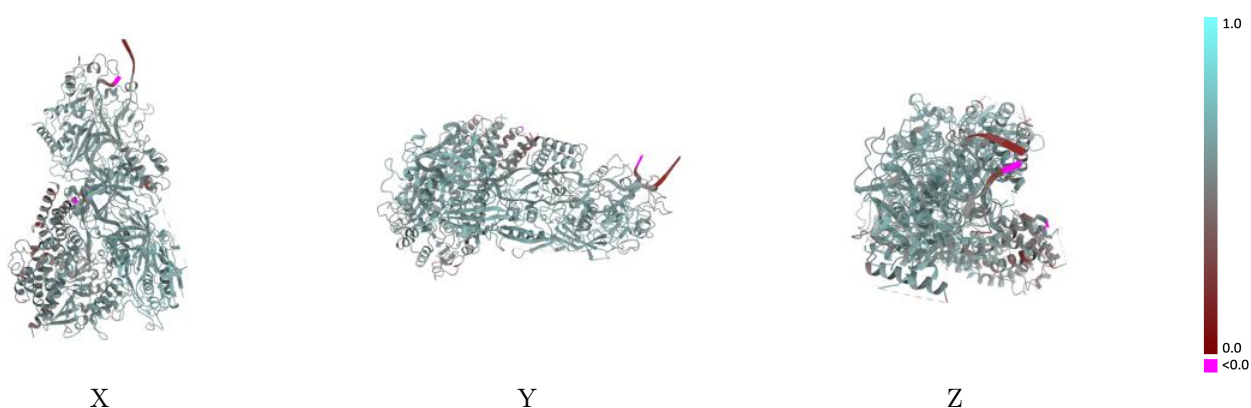
9 Map-model fit [i](#)

This section contains information regarding the fit between EMD map EMD-33681 and PDB model 7Y85. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)

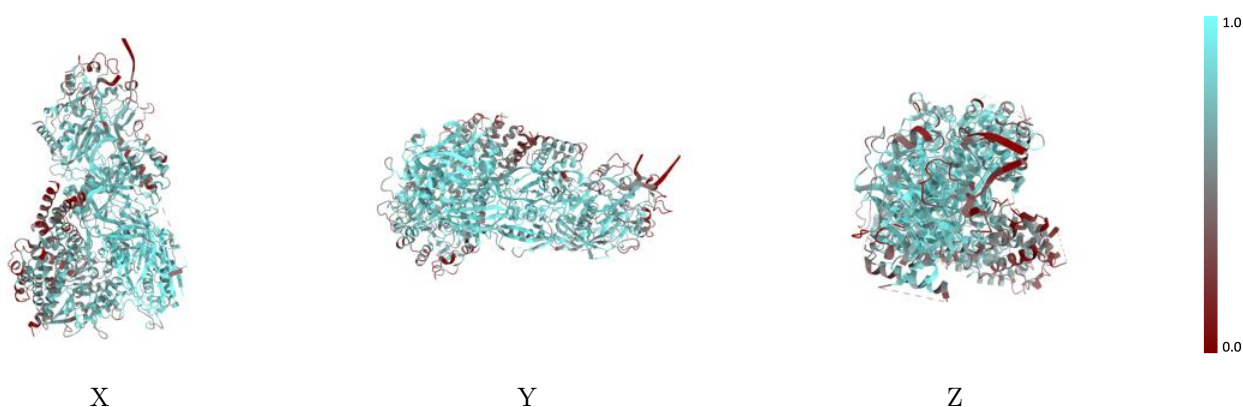
This section was not generated.

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



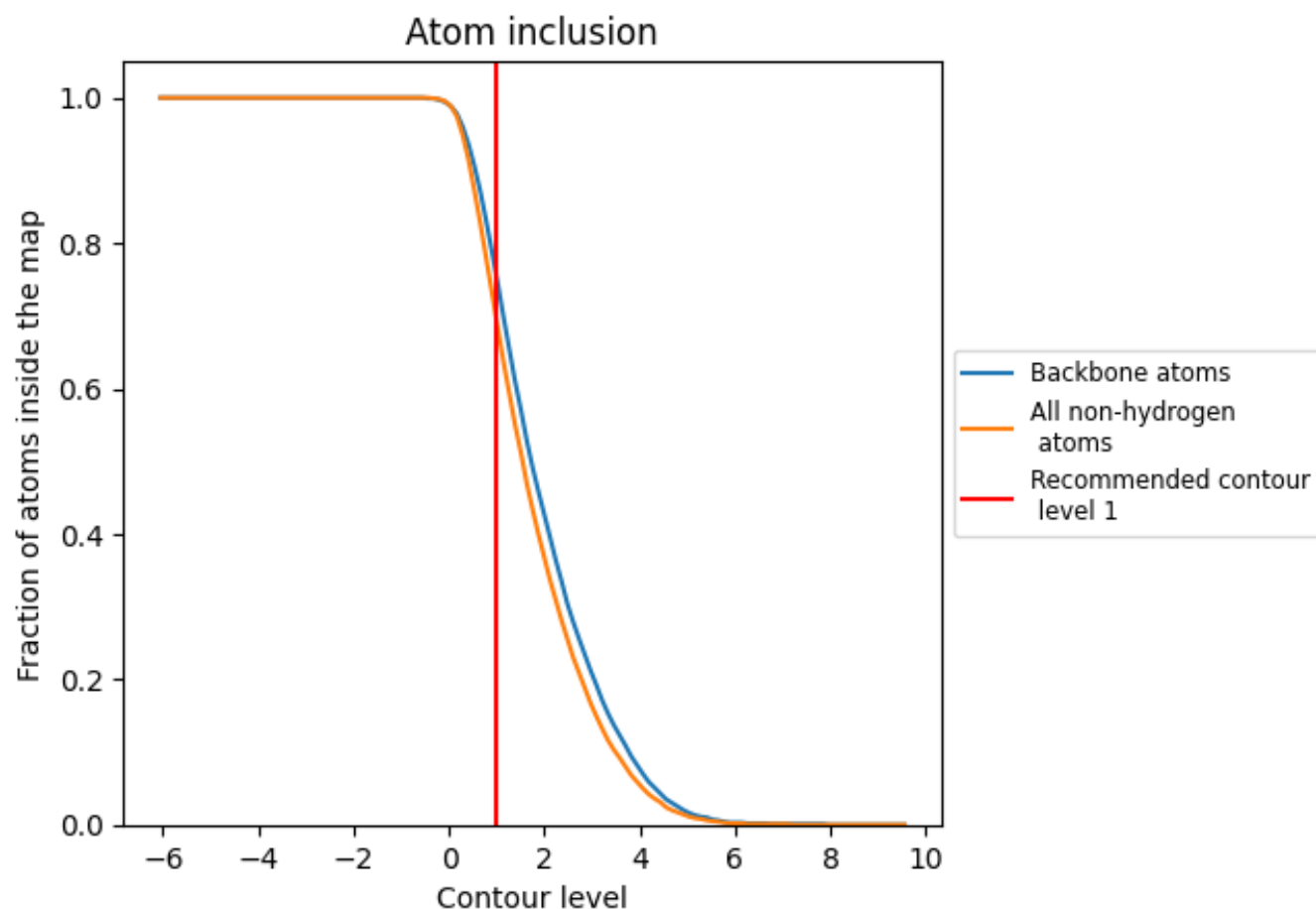
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 70% 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 (1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6950	0.5700
A	0.7520	0.5910
B	0.8300	0.5700
C	0.6530	0.5100
D	0.5730	0.5360

