



Full wwPDB EM Validation Report ⓘ

Mar 28, 2026 – 04:33 AM UTC

PDB ID : 7TID / pdb_00007tid
EMDB ID : EMD-25617
Title : Structure of the yeast clamp loader (Replication Factor C RFC) bound to the sliding clamp (Proliferating Cell Nuclear Antigen PCNA) and primer-template DNA
Authors : Gaubitz, C.; Liu, X.; Pajak, J.; Stone, N.; Hayes, J.; Demo, G.; Kelch, B.A.
Deposited on : 2022-01-13
Resolution : 3.30 Å (reported)
Based on initial model : 1SXJ

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

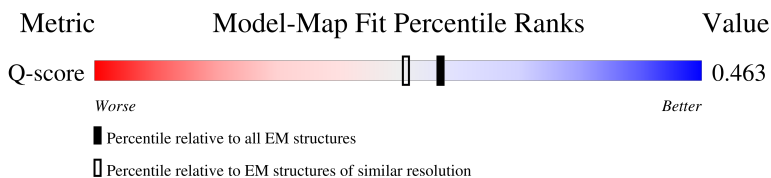
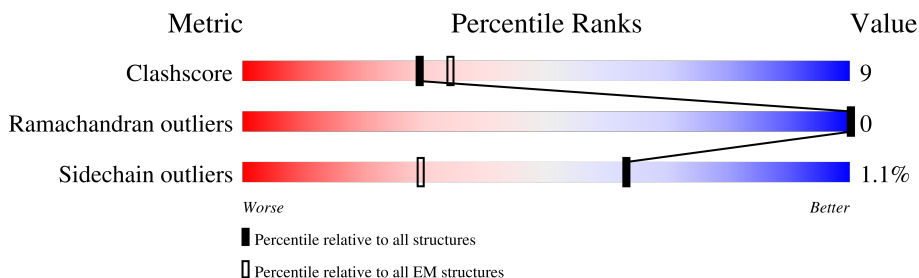
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	15087 (2.80 - 3.80)

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	861	
2	B	323	
3	C	340	

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Mol	Chain	Length	Quality of chain
4	D	353	73%23%.
5	E	354	71%25%5%
6	F	264	77%19%..
6	G	264	6% 66%30%..
6	H	264	23% 65%31%.
7	I	30	17% 47%33%20%
8	J	20	50%40%10%

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 21289 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 Replication factor C subunit 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	486	Total	C	N	O	S	0	0
			3848	2441	675	715	17		

- Molecule 2 is a protein called Replication factor C subunit 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	319	Total	C	N	O	S	0	0
			2505	1580	446	466	13		

- Molecule 3 is a protein called Replication factor C subunit 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	327	Total	C	N	O	S	0	0
			2586	1628	451	499	8		

- Molecule 4 is a protein called Replication factor C subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	340	Total	C	N	O	S	0	0
			2687	1696	464	517	10		

- Molecule 5 is a protein called Replication factor C subunit 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	338	Total	C	N	O	S	0	0
			2672	1693	466	495	18		

- Molecule 6 is a protein called Proliferating cell nuclear antigen.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	255	Total	C	N	O	S	0	0
			2003	1281	316	396	10		

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Mol	Chain	Residues	Atoms					AltConf	Trace
6	G	254	Total	C	N	O	S	0	0
			1994	1275	315	395	9		
6	H	253	Total	C	N	O	S	0	0
			1984	1268	313	393	10		

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	-5	GLY	-	expression tag	UNP P15873
F	-4	PRO	-	expression tag	UNP P15873
F	-3	HIS	-	expression tag	UNP P15873
F	-2	MET	-	expression tag	UNP P15873
F	-1	ALA	-	expression tag	UNP P15873
F	0	SER	-	expression tag	UNP P15873
G	-5	GLY	-	expression tag	UNP P15873
G	-4	PRO	-	expression tag	UNP P15873
G	-3	HIS	-	expression tag	UNP P15873
G	-2	MET	-	expression tag	UNP P15873
G	-1	ALA	-	expression tag	UNP P15873
G	0	SER	-	expression tag	UNP P15873
H	-5	GLY	-	expression tag	UNP P15873
H	-4	PRO	-	expression tag	UNP P15873
H	-3	HIS	-	expression tag	UNP P15873
H	-2	MET	-	expression tag	UNP P15873
H	-1	ALA	-	expression tag	UNP P15873
H	0	SER	-	expression tag	UNP P15873

- Molecule 7 is a DNA chain called DNA (5'-D(P*TP*TP*TP*TP*TP*TP*TP*AP*TP*GP*TP*AP*CP*TP*CP*GP*TP*AP*GP*TP*GP*TP*CP*T)-3').

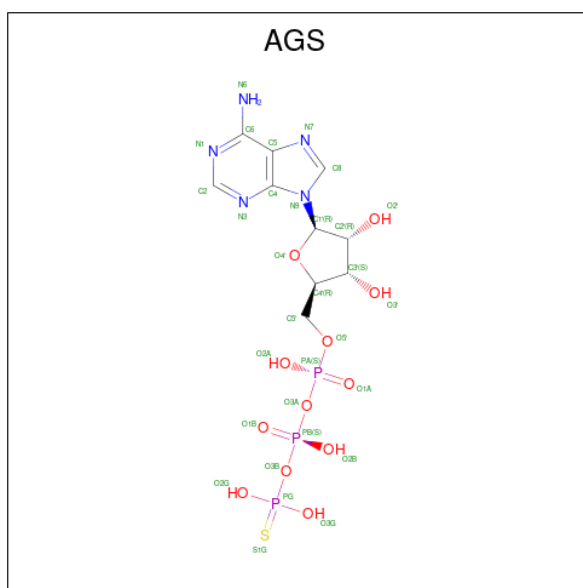
Mol	Chain	Residues	Atoms					AltConf	Trace
7	I	24	Total	C	N	O	P	0	0
			488	237	72	155	24		

- Molecule 8 is a DNA chain called DNA (5'-D(*AP*GP*AP*CP*AP*CP*TP*AP*CP*GP*AP*GP*TP*AP*CP*AP*TP*A)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
8	J	18	Total	C	N	O	P	0	0
			367	176	73	101	17		

- Molecule 9 is PHOSPHOTHIOPHOSPHORIC ACID-ADENYLATE ESTER (CCD ID:

AGS) (formula: C₁₀H₁₆N₅O₁₂P₃S) (labeled as "Ligand of Interest" by depositor).

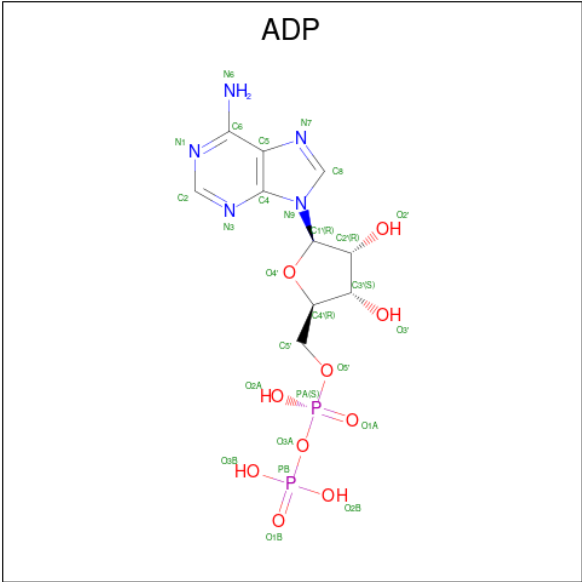


Mol	Chain	Residues	Atoms						AltConf
9	A	1	Total	C	N	O	P	S	0
			31	10	5	12	3	1	
9	B	1	Total	C	N	O	P	S	0
			31	10	5	12	3	1	
9	C	1	Total	C	N	O	P	S	0
			31	10	5	12	3	1	
9	D	1	Total	C	N	O	P	S	0
			31	10	5	12	3	1	

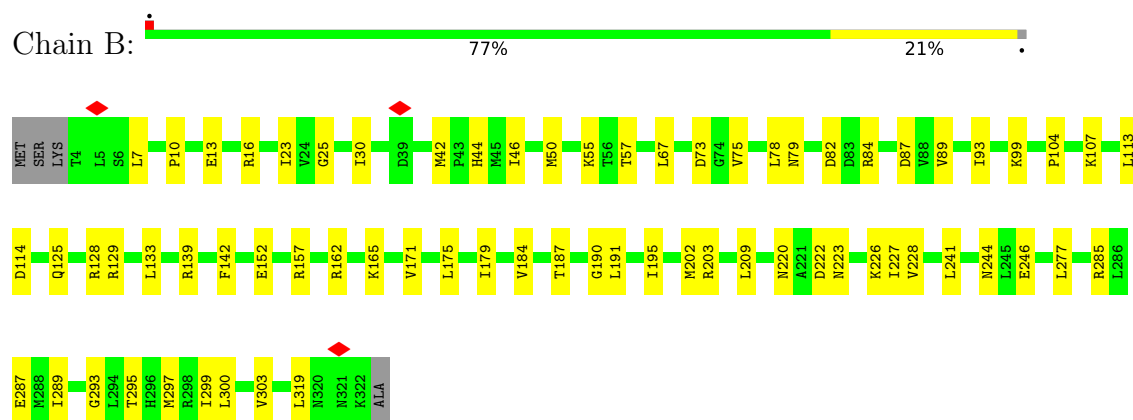
- Molecule 10 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
10	A	1	Total	Mg	0
			1	1	
10	B	1	Total	Mg	0
			1	1	
10	C	1	Total	Mg	0
			1	1	
10	D	1	Total	Mg	0
			1	1	

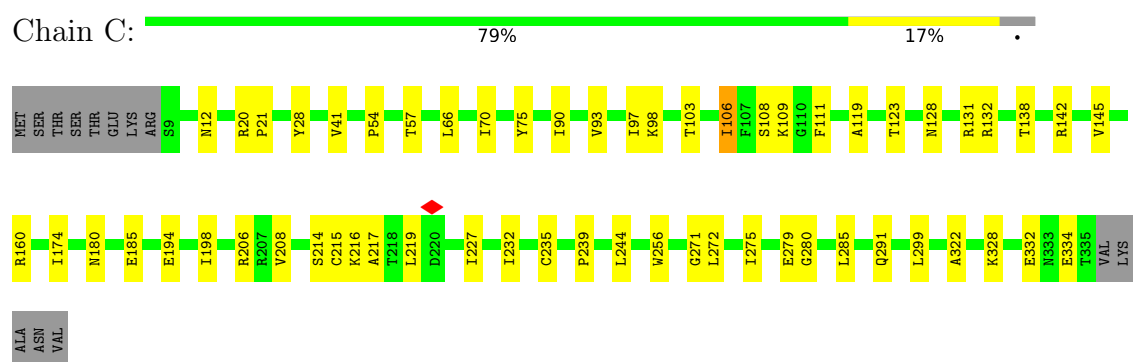
- Molecule 11 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂) (labeled as "Ligand of Interest" by depositor).



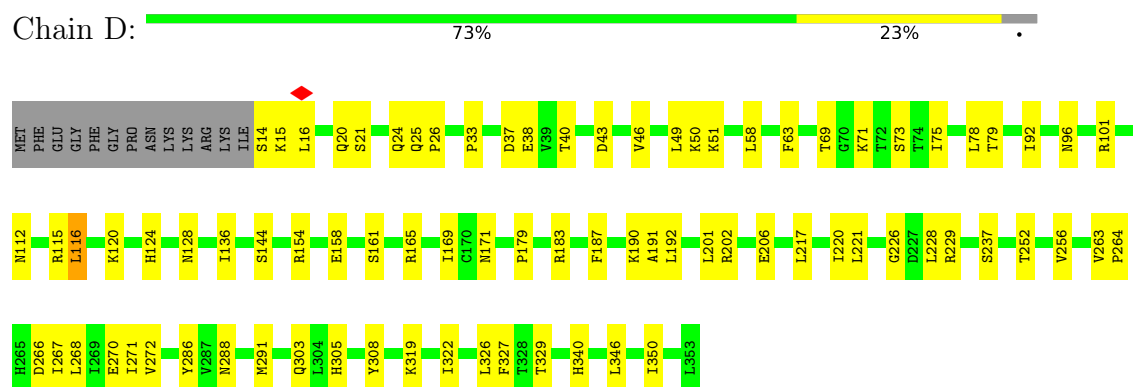
Mol	Chain	Residues	Atoms					AltConf
11	E	1	Total	C	N	O	P	0
			27	10	5	10	2	



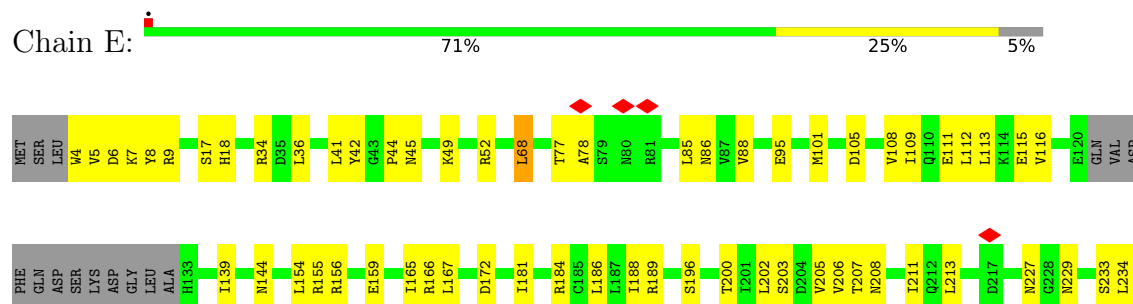
• Molecule 3: Replication factor C subunit 3



• Molecule 4: Replication factor C subunit 2

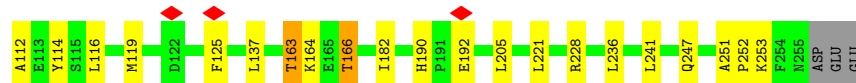
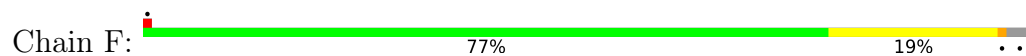


• Molecule 5: Replication factor C subunit 5

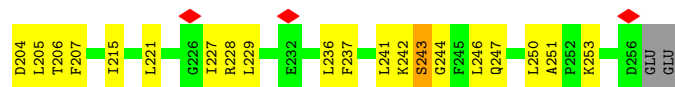
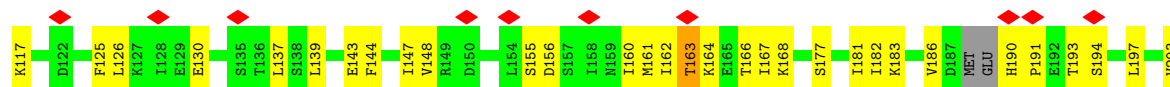
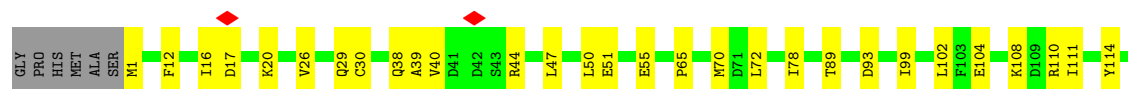




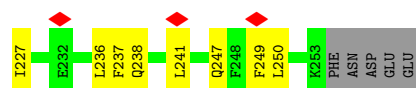
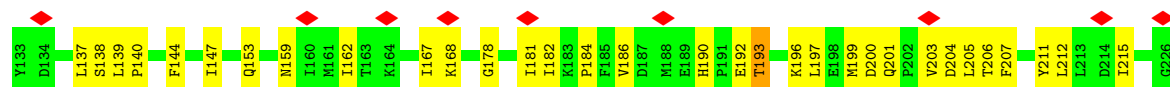
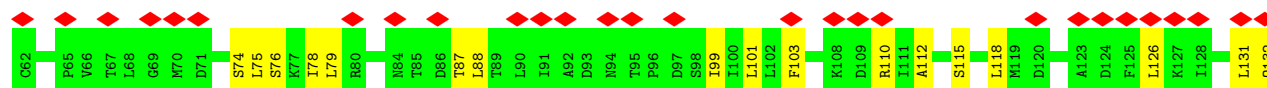
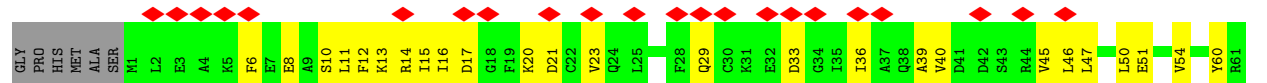
- Molecule 6: Proliferating cell nuclear antigen



- Molecule 6: Proliferating cell nuclear antigen



- Molecule 6: Proliferating cell nuclear antigen



- Molecule 7: DNA (5'-D(P*TP*TP*TP*TP*TP*TP*TP*AP*TP*GP*TP*AP*CP*TP*CP*GP*TP*AP*GP*TP*GP*TP*CP*T)-3')



- Molecule 8: DNA (5'-D(*AP*GP*AP*CP*AP*CP*TP*AP*CP*GP*AP*GP*TP*AP*CP*AP*TP*A)-3')



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	76270	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2300	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	19.399	Depositor
Minimum map value	-9.166	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	0.7	Depositor
Map size (Å)	139.92, 148.4, 143.09999	wwPDB
Map dimensions	135, 140, 132	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.06, 1.06, 1.06	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: ADP, MG, AGS

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	0.12	0/3924	0.33	0/5301
2	B	0.10	0/2540	0.29	0/3425
3	C	0.11	0/2626	0.32	0/3552
4	D	0.09	0/2732	0.28	0/3694
5	E	0.11	0/2710	0.31	0/3661
6	F	0.12	0/2033	0.33	0/2742
6	G	0.13	0/2023	0.35	0/2728
6	H	0.14	0/2013	0.38	0/2715
7	I	0.16	0/542	0.40	0/835
8	J	0.15	0/413	0.32	0/635
All	All	0.12	0/21556	0.33	0/29288

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3848	0	3889	79	0
2	B	2505	0	2600	53	0
3	C	2586	0	2614	37	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	2687	0	2721	56	0
5	E	2672	0	2797	61	0
6	F	2003	0	2025	34	0
6	G	1994	0	2013	54	0
6	H	1984	0	2010	55	0
7	I	488	0	279	7	0
8	J	367	0	203	8	0
9	A	31	0	12	3	0
9	B	31	0	12	3	0
9	C	31	0	12	2	0
9	D	31	0	12	2	0
10	A	1	0	0	0	0
10	B	1	0	0	0	0
10	C	1	0	0	0	0
10	D	1	0	0	0	0
11	E	27	0	12	2	0
All	All	21289	0	21211	394	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 9.

All (394) 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:D:288:ASN:HD21	5:E:227:ASN:HD21	1.27	0.81
3:C:41:VAL:HG21	3:C:66:LEU:HD11	1.64	0.77
4:D:263:VAL:H	4:D:303:GLN:HE22	1.32	0.76
6:G:51:GLU:O	6:G:244:GLY:HA3	1.85	0.76
5:E:4:TRP:HD1	5:E:6:ASP:H	1.32	0.76
3:C:174:ILE:HD11	3:C:208:VAL:HG21	1.69	0.74
6:G:177:SER:H	6:H:115:SER:HB3	1.54	0.72
6:H:103:PHE:HB2	6:H:112:ALA:HB3	1.72	0.72
6:H:168:LYS:HG2	6:H:181:ILE:HD12	1.71	0.71
1:A:328:GLU:HA	1:A:331:LYS:HG2	1.72	0.71
6:H:45:VAL:HG23	6:H:46:LEU:HD12	1.73	0.70
6:G:161:MET:HB3	6:G:168:LYS:HB3	1.72	0.70
1:A:550:LYS:HE3	1:A:551:PRO:HD2	1.75	0.69
6:G:30:CYS:O	6:G:65:PRO:HA	1.92	0.69
1:A:594:GLU:HG3	1:A:667:THR:HA	1.75	0.68
6:G:221:LEU:HD23	6:G:241:LEU:HD21	1.77	0.67
2:B:82:ASP:OD1	3:C:98:LYS:NZ	2.29	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:CYS:H	9:A:901:AGS:HN62	1.42	0.66
2:B:203:ARG:HE	9:B:401:AGS:H5'1	1.60	0.65
3:C:271:GLY:HA3	4:D:190:LYS:HD3	1.79	0.65
2:B:13:GLU:HG3	3:C:138:THR:HG21	1.79	0.65
1:A:755:THR:HG22	1:A:759:ILE:HG23	1.77	0.65
6:F:13:LYS:HG2	6:F:79:LEU:HB3	1.79	0.64
6:F:49:SER:HB3	6:F:247:GLN:HG3	1.80	0.64
6:H:23:VAL:HG11	6:H:39:ALA:HB1	1.79	0.64
1:A:295:ASP:O	2:B:139:ARG:NH2	2.32	0.63
1:A:402:VAL:HG11	6:G:44:ARG:HE	1.63	0.63
2:B:46:ILE:HG12	2:B:142:PHE:HB2	1.80	0.62
5:E:45:ASN:O	5:E:229:ASN:ND2	2.31	0.62
5:E:101:MET:HE3	5:E:105:ASP:HA	1.81	0.62
1:A:399:MET:HG3	6:G:251:ALA:HB1	1.80	0.62
6:H:6:PHE:O	6:H:87:THR:OG1	2.18	0.62
6:H:137:LEU:HD13	6:H:139:LEU:HD23	1.82	0.62
5:E:205:VAL:HG21	5:E:234:LEU:HD21	1.82	0.62
6:F:221:LEU:HB3	6:F:241:LEU:HD21	1.81	0.61
1:A:375:LEU:HD22	1:A:421:ILE:HD12	1.82	0.61
6:H:88:LEU:HD11	6:H:101:LEU:HD12	1.81	0.61
6:H:36:ILE:HD12	6:H:51:GLU:HG3	1.83	0.61
2:B:84:ARG:NH1	2:B:114:ASP:O	2.34	0.61
1:A:461:PRO:HA	1:A:464:ARG:HG2	1.82	0.61
2:B:128:ARG:HE	2:B:157:ARG:HH21	1.49	0.60
5:E:4:TRP:HB3	5:E:7:LYS:HG2	1.83	0.60
4:D:120:LYS:HD2	6:F:95:THR:HB	1.83	0.60
4:D:192:LEU:HD12	4:D:228:LEU:HB2	1.82	0.59
2:B:175:LEU:HD22	2:B:209:LEU:HD22	1.84	0.59
3:C:12:ASN:O	4:D:165:ARG:NH2	2.36	0.59
5:E:36:LEU:O	5:E:166:ARG:NH1	2.36	0.59
6:G:228:ARG:HD2	6:G:236:LEU:HD23	1.84	0.58
2:B:57:THR:HG21	9:B:401:AGS:H3'	1.86	0.58
2:B:67:LEU:HD12	2:B:75:VAL:HG23	1.84	0.58
1:A:477:ARG:HE	1:A:512:GLY:HA3	1.68	0.57
6:H:88:LEU:HD13	6:H:103:PHE:HE2	1.68	0.57
1:A:314:LYS:HE3	1:A:318:MET:HE2	1.85	0.57
1:A:351:LEU:HD22	1:A:475:PHE:HE1	1.69	0.57
1:A:464:ARG:HH12	1:A:467:ASP:HB2	1.68	0.57
3:C:285:LEU:HD13	3:C:299:LEU:HD21	1.84	0.57
4:D:158:GLU:OE2	4:D:183:ARG:NH1	2.38	0.57
4:D:25:GLN:HG3	4:D:26:PRO:HD2	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:510:THR:HG21	1:A:517:VAL:HG23	1.87	0.57
6:H:8:GLU:HB3	6:H:11:LEU:HD23	1.87	0.57
2:B:73:ASP:OD1	2:B:73:ASP:N	2.38	0.57
6:F:228:ARG:HB2	6:F:236:LEU:HB3	1.87	0.57
6:F:89:THR:HB	6:F:102:LEU:HB2	1.86	0.56
6:F:228:ARG:HD2	6:F:236:LEU:HD23	1.86	0.56
1:A:320:LEU:HD13	1:A:351:LEU:HD21	1.86	0.56
1:A:487:LEU:HA	1:A:490:ILE:HG12	1.86	0.56
1:A:637:LEU:HD22	2:B:300:LEU:HD13	1.86	0.56
5:E:244:ASN:HD22	5:E:248:LEU:HA	1.69	0.56
6:G:38:GLN:HE22	6:G:126:LEU:H	1.51	0.56
6:G:161:MET:HE1	6:G:204:ASP:HB2	1.87	0.56
1:A:420:VAL:HG12	1:A:450:PRO:HB2	1.87	0.56
1:A:515:ARG:HD2	2:B:157:ARG:HA	1.88	0.56
1:A:712:ARG:HG3	5:E:4:TRP:HZ3	1.71	0.56
2:B:171:VAL:HA	2:B:202:MET:HE1	1.88	0.56
3:C:185:GLU:OE1	3:C:216:LYS:NZ	2.38	0.56
5:E:17:SER:H	11:E:401:ADP:HN62	1.53	0.56
5:E:263:ILE:O	5:E:267:THR:HG23	2.06	0.56
5:E:41:LEU:HD13	5:E:188:ILE:HB	1.88	0.56
6:H:197:LEU:HG	6:H:199:MET:HE3	1.88	0.56
1:A:714:LEU:HD13	1:A:746:MET:HE2	1.88	0.55
6:F:80:ARG:NH2	6:H:153:GLN:OE1	2.39	0.55
1:A:322:ASN:O	1:A:326:ASN:ND2	2.40	0.55
1:A:524:ILE:HD13	1:A:539:ILE:HD12	1.89	0.55
5:E:267:THR:HG22	5:E:303:LEU:HD12	1.88	0.55
6:F:236:LEU:HD11	6:F:247:GLN:HB2	1.88	0.55
1:A:394:ASN:O	1:A:398:ASN:ND2	2.39	0.55
3:C:219:LEU:HD21	3:C:227:ILE:HG12	1.88	0.55
3:C:20:ARG:NH2	4:D:158:GLU:OE1	2.40	0.55
3:C:54:PRO:HD2	3:C:57:THR:HG21	1.88	0.55
6:G:70:MET:HE2	6:G:99:ILE:HG21	1.89	0.55
1:A:642:PRO:O	1:A:645:ALA:HB3	2.06	0.55
1:A:593:GLN:O	5:E:283:ARG:NH2	2.40	0.54
6:G:160:ILE:HD13	6:G:229:LEU:HD11	1.88	0.54
8:J:18:DA:H8	8:J:18:DA:H5"	1.72	0.54
1:A:611:LEU:HD21	5:E:350:MET:HB3	1.89	0.54
2:B:179:ILE:HD12	2:B:184:VAL:HB	1.90	0.54
6:G:139:LEU:HD11	6:G:144:PHE:HB2	1.89	0.54
5:E:4:TRP:CD1	5:E:6:ASP:H	2.21	0.54
4:D:319:LYS:HA	4:D:322:ILE:HG22	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:H:227:ILE:HD12	6:H:237:PHE:HD2	1.73	0.54
1:A:681:ARG:O	1:A:685:GLU:HG3	2.08	0.53
5:E:206:VAL:HG21	5:E:213:LEU:HD11	1.90	0.53
7:I:26:DT:H2''	7:I:27:DC:C4	2.44	0.53
3:C:90:ILE:N	7:I:15:DT:OP1	2.42	0.53
4:D:79:THR:HG22	4:D:136:ILE:HD12	1.89	0.53
4:D:268:LEU:O	4:D:272:VAL:HG23	2.09	0.53
6:H:17:ASP:HA	6:H:20:LYS:HD3	1.91	0.53
5:E:154:LEU:HD11	5:E:167:LEU:HD11	1.90	0.52
6:F:38:GLN:HB2	6:F:125:PHE:HB3	1.90	0.52
2:B:50:MET:O	2:B:55:LYS:NZ	2.42	0.52
6:F:103:PHE:HB2	6:F:112:ALA:HB3	1.91	0.52
6:G:190:HIS:HB2	6:G:193:THR:HG22	1.91	0.52
1:A:428:GLY:HA3	2:B:125:GLN:HG2	1.92	0.52
3:C:119:ALA:HB3	3:C:145:VAL:HG13	1.90	0.52
1:A:292:ARG:HH22	1:A:294:GLU:HG2	1.74	0.52
6:F:205:LEU:HD13	6:F:252:PRO:HB3	1.92	0.52
2:B:125:GLN:OE1	2:B:128:ARG:NH1	2.40	0.52
6:G:163:THR:O	6:G:164:LYS:HG2	2.10	0.52
5:E:44:PRO:O	5:E:49:LYS:NZ	2.42	0.52
6:G:205:LEU:HD12	6:G:229:LEU:HD12	1.91	0.52
5:E:271:VAL:HG11	5:E:310:VAL:HG11	1.91	0.52
4:D:69:THR:OG1	4:D:71:LYS:NZ	2.42	0.51
1:A:320:LEU:HD21	1:A:366:VAL:HG21	1.92	0.51
2:B:89:VAL:HG22	2:B:93:ILE:HD12	1.92	0.51
6:H:12:PHE:HA	6:H:15:ILE:HG22	1.91	0.51
6:F:66:VAL:HG13	6:F:96:PRO:HG3	1.92	0.51
6:H:12:PHE:O	6:H:16:ILE:HG12	2.09	0.51
6:F:114:TYR:HD1	6:H:178:GLY:HA2	1.76	0.51
6:F:54:VAL:HA	6:F:57:PHE:HD2	1.76	0.51
1:A:413:ASN:OD1	1:A:416:GLY:N	2.44	0.51
1:A:722:ALA:HB2	1:A:766:VAL:HG22	1.92	0.51
4:D:171:ASN:ND2	9:D:401:AGS:S1G	2.77	0.51
6:H:131:LEU:HD12	6:H:132:GLN:H	1.76	0.51
1:A:696:THR:HG22	1:A:697:ASP:H	1.75	0.51
3:C:93:VAL:HG22	3:C:97:ILE:HD12	1.92	0.51
6:H:20:LYS:HE3	6:H:76:SER:HB3	1.93	0.51
5:E:68:LEU:HD13	5:E:88:VAL:HG13	1.93	0.51
2:B:23:ILE:HG22	2:B:25:GLY:H	1.75	0.50
4:D:266:ASP:O	4:D:270:GLU:HG3	2.12	0.50
1:A:304:PRO:HD3	9:A:901:AGS:H2'	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:503:ILE:H	1:A:503:ILE:HD12	1.74	0.50
6:H:227:ILE:HD12	6:H:237:PHE:CD2	2.47	0.50
6:G:167:ILE:HB	6:G:182:ILE:HB	1.93	0.50
2:B:295:THR:O	2:B:299:ILE:HG12	2.10	0.50
6:G:215:ILE:HG21	6:G:237:PHE:HD2	1.77	0.50
4:D:271:ILE:HG12	4:D:286:TYR:HE2	1.77	0.50
6:H:8:GLU:HG3	6:H:10:SER:H	1.76	0.50
7:I:24:DT:H2"	7:I:25:DG:C5	2.46	0.50
3:C:108:SER:OG	3:C:109:LYS:N	2.45	0.50
6:F:41:ASP:OD2	6:F:43:SER:OG	2.30	0.50
4:D:308:TYR:O	4:D:319:LYS:NZ	2.45	0.49
5:E:203:SER:O	5:E:207:THR:HG23	2.12	0.49
3:C:217:ALA:HA	4:D:51:LYS:HD2	1.93	0.49
6:F:2:LEU:HB3	6:F:92:ALA:HB3	1.93	0.49
6:H:74:SER:O	6:H:78:ILE:HG12	2.13	0.49
6:G:143:GLU:OE2	6:H:110:ARG:NH1	2.45	0.49
6:G:144:PHE:O	6:G:148:VAL:HG23	2.13	0.49
1:A:547:ILE:HD13	2:B:152:GLU:HB2	1.95	0.49
5:E:112:LEU:O	5:E:116:VAL:HG23	2.12	0.49
6:F:88:LEU:HD11	6:F:101:LEU:HD23	1.94	0.49
6:H:212:LEU:HA	6:H:215:ILE:HD12	1.94	0.49
5:E:144:ASN:HD21	5:E:172:ASP:H	1.59	0.49
3:C:194:GLU:O	3:C:198:ILE:HG12	2.12	0.49
3:C:328:LYS:O	3:C:332:GLU:HG2	2.11	0.49
4:D:264:PRO:HD2	4:D:267:ILE:HD12	1.95	0.49
6:G:78:ILE:HG23	6:G:114:TYR:HD2	1.78	0.49
5:E:95:GLU:HG2	5:E:139:ILE:HB	1.94	0.49
1:A:402:VAL:HG21	6:G:44:ARG:HH21	1.78	0.49
1:A:646:VAL:HG21	2:B:289:ILE:HG21	1.94	0.49
1:A:646:VAL:HA	1:A:650:VAL:HB	1.94	0.49
6:F:163:THR:HG22	6:F:164:LYS:H	1.77	0.49
1:A:544:GLU:O	2:B:162:ARG:NH2	2.46	0.49
4:D:16:LEU:O	4:D:20:GLN:HG2	2.13	0.49
5:E:259:TRP:HZ3	5:E:302:GLU:HG3	1.78	0.49
6:H:138:SER:HB3	6:H:196:LYS:HB2	1.95	0.49
6:G:166:THR:HB	6:G:181:ILE:HD11	1.95	0.48
2:B:87:ASP:OD2	2:B:87:ASP:N	2.46	0.48
2:B:227:ILE:HG23	2:B:228:VAL:HG23	1.95	0.48
1:A:396:LEU:HD13	1:A:447:THR:HB	1.95	0.48
6:G:186:VAL:HA	6:G:194:SER:HB2	1.95	0.48
1:A:483:ILE:HD12	1:A:507:ILE:HG21	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:H:200:ASP:HB2	6:H:201:GLN:OE1	2.13	0.48
2:B:220:ASN:HD22	2:B:222:ASP:H	1.61	0.48
6:H:137:LEU:HD22	6:H:167:ILE:HD13	1.96	0.48
5:E:109:ILE:HG23	5:E:113:LEU:HD12	1.96	0.48
6:G:155:SER:OG	6:G:156:ASP:N	2.47	0.48
3:C:131:ARG:HB2	3:C:160:ARG:HE	1.77	0.48
5:E:88:VAL:HB	5:E:95:GLU:HB2	1.94	0.48
8:J:18:DA:H2''	8:J:19:DT:H5'	1.95	0.48
1:A:351:LEU:HD23	1:A:473:ILE:HB	1.95	0.47
2:B:220:ASN:ND2	2:B:222:ASP:H	2.12	0.47
4:D:58:LEU:O	4:D:165:ARG:NE	2.46	0.47
2:B:104:PRO:HB2	2:B:107:LYS:HG2	1.96	0.47
3:C:21:PRO:HD3	9:C:401:AGS:H2'	1.96	0.47
1:A:517:VAL:O	1:A:521:LEU:HG	2.14	0.47
1:A:643:LEU:HD13	2:B:293:GLY:HA3	1.95	0.47
4:D:329:THR:HG21	4:D:346:LEU:HD22	1.96	0.47
5:E:267:THR:HG21	5:E:306:SER:OG	2.14	0.47
6:F:7:GLU:HG2	6:F:8:GLU:N	2.29	0.47
6:G:55:GLU:CD	6:G:55:GLU:H	2.22	0.47
8:J:3:DA:H2''	8:J:4:DG:C8	2.49	0.47
6:G:55:GLU:HG2	6:G:243:SER:HB2	1.97	0.47
6:H:13:LYS:HA	6:H:79:LEU:HD23	1.95	0.47
8:J:17:DC:H2''	8:J:18:DA:C8	2.50	0.47
1:A:691:ARG:NH2	5:E:95:GLU:OE1	2.48	0.46
6:F:83:ASN:ND2	6:F:85:THR:H	2.13	0.46
2:B:277:LEU:HB3	2:B:285:ARG:HD3	1.97	0.46
6:H:54:VAL:HG13	6:H:60:TYR:HB3	1.97	0.46
8:J:9:DT:H2''	8:J:10:DA:N7	2.30	0.46
5:E:257:PRO:HD2	5:E:260:ILE:HD12	1.97	0.46
7:I:20:DG:H2''	7:I:21:DT:C5	2.51	0.46
2:B:79:ASN:HB3	3:C:132:ARG:HG2	1.96	0.46
1:A:702:ARG:O	1:A:702:ARG:NH1	2.44	0.46
4:D:21:SER:HA	4:D:24:GLN:HG2	1.98	0.46
4:D:171:ASN:OD1	5:E:155:ARG:NH1	2.49	0.46
6:G:143:GLU:O	6:G:147:ILE:HG13	2.16	0.46
4:D:33:PRO:HB3	4:D:38:GLU:HB2	1.97	0.46
6:G:12:PHE:O	6:G:16:ILE:HG12	2.16	0.46
5:E:202:LEU:HD22	5:E:237:LEU:HD22	1.98	0.46
6:H:207:PHE:HD2	6:H:250:LEU:HD22	1.80	0.46
9:A:901:AGS:S1G	2:B:128:ARG:NH2	2.89	0.46
2:B:129:ARG:HG3	2:B:133:LEU:HG	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:G:206:THR:HB	6:G:253:LYS:HE3	1.96	0.46
2:B:42:MET:O	2:B:139:ARG:NH1	2.42	0.45
4:D:201:LEU:HD23	4:D:217:LEU:HB3	1.97	0.45
6:G:104:GLU:HG3	6:G:111:ILE:HG22	1.97	0.45
1:A:378:ASN:HD21	2:B:129:ARG:HG2	1.81	0.45
7:I:25:DG:H22	8:J:7:DA:H61	1.63	0.45
1:A:352:TYR:CZ	1:A:474:GLN:HB2	2.50	0.45
5:E:258:ASP:OD1	5:E:258:ASP:N	2.49	0.45
1:A:352:TYR:HE1	1:A:472:ASP:HB2	1.82	0.45
1:A:713:LEU:HB3	1:A:742:TRP:HE1	1.81	0.45
1:A:738:THR:H	1:A:741:ASP:HB2	1.82	0.45
4:D:37:ASP:OD2	4:D:50:LYS:NZ	2.49	0.45
5:E:8:TYR:HE2	5:E:208:ASN:HD22	1.63	0.45
6:G:207:PHE:HB3	6:G:250:LEU:HD21	1.99	0.45
2:B:187:THR:HG23	2:B:190:GLY:H	1.81	0.45
4:D:112:ASN:OD1	4:D:115:ARG:NH2	2.45	0.45
1:A:384:SER:O	1:A:388:LEU:HB3	2.17	0.45
1:A:581:TYR:HE1	1:A:589:PRO:HD3	1.82	0.45
6:G:16:ILE:O	6:G:20:LYS:HB3	2.17	0.45
1:A:619:ASN:OD1	5:E:331:ARG:NH1	2.42	0.45
4:D:43:ASP:HA	4:D:46:VAL:HG22	1.99	0.45
6:G:29:GLN:HG3	6:G:125:PHE:HE2	1.81	0.45
1:A:458:ARG:HH21	1:A:472:ASP:HB3	1.82	0.44
2:B:241:LEU:HD21	2:B:277:LEU:HD21	1.99	0.44
1:A:587:PHE:HZ	7:I:8:DT:H3	1.65	0.44
4:D:220:ILE:HG12	4:D:256:VAL:HG11	1.99	0.44
6:H:40:VAL:HG22	6:H:47:LEU:HD22	1.99	0.44
1:A:412:GLN:OE1	1:A:413:ASN:ND2	2.51	0.44
2:B:30:ILE:HD13	2:B:30:ILE:HA	1.84	0.44
5:E:116:VAL:HG12	5:E:165:ILE:HD11	1.98	0.44
6:G:130:GLU:CD	6:G:130:GLU:H	2.25	0.44
6:H:162:ILE:HD12	6:H:167:ILE:HG12	1.99	0.44
2:B:223:ASN:O	2:B:227:ILE:HG22	2.16	0.44
6:G:55:GLU:O	6:G:242:LYS:NZ	2.50	0.44
1:A:383:ARG:NH2	1:A:425:GLU:O	2.51	0.44
2:B:223:ASN:HA	2:B:226:LYS:HE3	1.99	0.44
2:B:13:GLU:OE2	2:B:16:ARG:NH2	2.43	0.44
3:C:128:ASN:HA	3:C:131:ARG:HE	1.82	0.44
4:D:229:ARG:HD3	5:E:184:ARG:HA	2.00	0.44
6:G:162:ILE:HG23	6:G:203:VAL:H	1.83	0.44
6:H:46:LEU:O	6:H:47:LEU:HD23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:695:SER:OG	5:E:9:ARG:NH1	2.51	0.44
6:H:238:GLN:HG2	6:H:247:GLN:HG3	1.99	0.44
1:A:612:GLU:O	1:A:616:GLU:HG3	2.17	0.43
2:B:202:MET:HG2	9:B:401:AGS:C4	2.48	0.43
5:E:267:THR:O	5:E:271:VAL:HG22	2.17	0.43
6:G:1:MET:HG2	6:G:93:ASP:HA	1.99	0.43
6:G:236:LEU:HD11	6:G:247:GLN:HB3	2.00	0.43
8:J:9:DT:H2''	8:J:10:DA:C5	2.52	0.43
3:C:256:TRP:CD1	3:C:328:LYS:HZ2	2.36	0.43
6:F:182:ILE:HG22	6:G:110:ARG:HG3	1.99	0.43
6:H:33:ASP:N	6:H:33:ASP:OD2	2.50	0.43
2:B:244:ASN:ND2	2:B:246:GLU:H	2.15	0.43
3:C:322:ALA:HB2	4:D:327:PHE:HB2	2.00	0.43
4:D:49:LEU:HD13	4:D:78:LEU:HD22	2.01	0.43
1:A:626:ILE:HD11	5:E:331:ARG:HA	2.00	0.43
3:C:106:ILE:HG12	6:F:251:ALA:HB2	2.01	0.43
4:D:14:SER:OG	4:D:15:LYS:N	2.47	0.43
4:D:346:LEU:O	4:D:350:ILE:HG23	2.19	0.43
6:G:163:THR:O	6:G:166:THR:OG1	2.27	0.43
1:A:722:ALA:HA	1:A:766:VAL:HG13	2.01	0.43
4:D:154:ARG:HD3	4:D:179:PRO:HB2	1.99	0.43
6:F:163:THR:HB	6:F:166:THR:OG1	2.19	0.43
6:H:184:PRO:HG3	6:H:197:LEU:HD23	2.00	0.43
1:A:313:ASN:OD1	1:A:476:ARG:N	2.32	0.43
2:B:223:ASN:HA	2:B:226:LYS:HG2	1.99	0.43
6:H:15:ILE:HG13	6:H:50:LEU:HD11	2.01	0.43
6:H:99:ILE:HG12	6:H:118:LEU:HD11	2.01	0.43
6:H:190:HIS:HB3	6:H:192:GLU:OE2	2.18	0.43
3:C:28:TYR:OH	3:C:180:ASN:ND2	2.51	0.43
6:H:236:LEU:HD13	6:H:249:PHE:CE2	2.54	0.43
1:A:729:ILE:HD12	1:A:774:TYR:HD1	1.84	0.43
4:D:191:ALA:HB1	4:D:226:GLY:HA3	1.99	0.43
5:E:85:LEU:HD13	5:E:108:VAL:HG11	2.00	0.43
5:E:283:ARG:NE	5:E:343:GLU:OE2	2.40	0.43
6:H:29:GLN:HE22	6:H:126:LEU:HD11	1.84	0.43
6:H:236:LEU:HD12	6:H:237:PHE:N	2.33	0.43
5:E:313:LEU:HD22	5:E:353:LEU:HD22	2.00	0.42
1:A:688:TYR:CE1	5:E:86:ASN:HB2	2.54	0.42
5:E:77:THR:OG1	5:E:78:ALA:N	2.52	0.42
5:E:159:GLU:OE1	5:E:184:ARG:NH2	2.52	0.42
6:G:16:ILE:HG23	6:G:72:LEU:HD22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:H:45:VAL:HG21	6:H:211:TYR:CZ	2.54	0.42
1:A:323:TRP:HD1	1:A:343:SER:HB2	1.83	0.42
6:F:107:LYS:HD2	6:F:107:LYS:HA	1.84	0.42
6:G:137:LEU:HB3	6:G:197:LEU:HG	2.02	0.42
3:C:28:TYR:O	9:C:401:AGS:N6	2.50	0.42
3:C:206:ARG:HH12	4:D:158:GLU:CD	2.28	0.42
4:D:201:LEU:HD22	4:D:221:LEU:HG	2.01	0.42
4:D:101:ARG:NH2	4:D:144:SER:OG	2.52	0.42
5:E:18:HIS:CE1	5:E:52:ARG:HH21	2.37	0.42
6:F:83:ASN:HD21	6:F:85:THR:HG22	1.85	0.42
6:G:183:LYS:HD2	6:G:183:LYS:HA	1.87	0.42
1:A:327:TRP:CD2	1:A:450:PRO:HG3	2.54	0.42
2:B:99:LYS:HD2	6:G:117:LYS:HD3	2.01	0.42
3:C:214:SER:HB3	3:C:235:CYS:SG	2.59	0.42
4:D:116:LEU:HA	6:F:119:MET:HA	2.00	0.42
6:H:140:PRO:HG3	6:H:193:THR:HA	2.01	0.42
6:F:13:LYS:HA	6:F:79:LEU:HD13	2.01	0.42
6:F:98:SER:HA	6:F:116:LEU:O	2.20	0.42
6:G:40:VAL:HG22	6:G:47:LEU:HD13	2.01	0.42
7:I:18:DT:H2''	7:I:19:DC:C6	2.54	0.42
8:J:18:DA:H2'	8:J:19:DT:C6	2.55	0.42
2:B:191:LEU:O	2:B:195:ILE:HG12	2.19	0.42
3:C:70:ILE:HD13	3:C:111:PHE:HZ	1.84	0.42
6:H:10:SER:O	6:H:14:ARG:HG3	2.20	0.42
6:H:144:PHE:HA	6:H:147:ILE:HG22	2.01	0.42
6:H:159:ASN:OD1	6:H:206:THR:HG22	2.20	0.42
2:B:10:PRO:HD3	3:C:142:ARG:HH22	1.84	0.42
2:B:165:LYS:HD3	2:B:165:LYS:HA	1.72	0.42
5:E:42:TYR:CZ	5:E:189:ARG:HB3	2.55	0.42
6:H:15:ILE:HD12	6:H:15:ILE:HA	1.94	0.42
3:C:244:LEU:HD21	3:C:280:GLY:HA3	2.02	0.41
3:C:275:ILE:O	3:C:279:GLU:HG2	2.20	0.41
6:F:16:ILE:HG23	6:F:72:LEU:HD22	2.02	0.41
6:G:26:VAL:HG13	6:G:39:ALA:HB2	2.01	0.41
1:A:561:GLY:HA2	1:A:657:ALA:HB1	2.02	0.41
4:D:96:ASN:CG	5:E:156:ARG:HG2	2.45	0.41
6:F:68:LEU:HD13	6:F:99:ILE:HD12	2.01	0.41
1:A:506:LEU:HD21	1:A:517:VAL:HG13	2.03	0.41
2:B:287:GLU:HB3	2:B:319:LEU:HD13	2.01	0.41
4:D:24:GLN:HB2	5:E:34:ARG:HD2	2.02	0.41
5:E:9:ARG:HD3	11:E:401:ADP:H3'	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:190:HIS:HB3	6:F:192:GLU:OE2	2.20	0.41
6:G:89:THR:HB	6:G:102:LEU:HB2	2.02	0.41
6:H:167:ILE:HB	6:H:182:ILE:CG2	2.50	0.41
4:D:237:SER:HB3	5:E:186:LEU:HD11	2.03	0.41
5:E:111:GLU:O	5:E:115:GLU:HG3	2.20	0.41
5:E:196:SER:O	5:E:200:THR:HG23	2.21	0.41
6:G:190:HIS:HA	6:G:191:PRO:HD3	1.95	0.41
1:A:505:ARG:O	1:A:509:THR:HG23	2.20	0.41
2:B:44:HIS:CE1	2:B:139:ARG:HA	2.55	0.41
5:E:229:ASN:O	5:E:233:SER:HB3	2.21	0.41
5:E:270:ILE:HD13	5:E:350:MET:HE3	2.03	0.41
6:H:21:ASP:OD1	6:H:21:ASP:N	2.52	0.41
4:D:161:SER:O	4:D:165:ARG:NH1	2.53	0.41
5:E:206:VAL:HB	5:E:211:ILE:HB	2.02	0.41
6:H:236:LEU:HD13	6:H:249:PHE:HE2	1.86	0.41
1:A:640:LEU:HD21	2:B:297:MET:HG2	2.02	0.41
3:C:291:GLN:HG2	3:C:334:GLU:HG3	2.03	0.41
4:D:16:LEU:HD12	4:D:16:LEU:H	1.86	0.41
4:D:79:THR:HB	4:D:92:ILE:HD11	2.03	0.41
4:D:124:HIS:CE1	4:D:128:ASN:HD21	2.39	0.41
5:E:240:MET:HE1	5:E:254:ILE:HD13	2.03	0.41
5:E:264:HIS:O	5:E:268:ARG:HG2	2.20	0.41
6:G:17:ASP:OD1	6:G:20:LYS:NZ	2.53	0.41
6:G:168:LYS:NZ	6:G:181:ILE:HB	2.36	0.41
1:A:479:ASP:OD1	1:A:482:SER:OG	2.25	0.41
4:D:202:ARG:O	4:D:206:GLU:HG2	2.20	0.41
4:D:228:LEU:HD23	9:D:401:AGS:C4	2.51	0.41
6:G:50:LEU:HD12	6:G:246:LEU:HD13	2.02	0.41
6:H:45:VAL:O	6:H:250:LEU:HA	2.21	0.41
1:A:347:ARG:NH2	1:A:447:THR:H	2.18	0.40
3:C:103:THR:HA	6:F:253:LYS:HA	2.02	0.40
4:D:63:PHE:HD1	4:D:187:PHE:HB2	1.86	0.40
4:D:291:MET:HG3	4:D:340:HIS:CG	2.56	0.40
5:E:275:SER:OG	5:E:276:VAL:N	2.54	0.40
6:H:162:ILE:HB	6:H:203:VAL:HB	2.03	0.40
6:H:204:ASP:C	6:H:205:LEU:HD12	2.46	0.40
4:D:63:PHE:HB3	4:D:71:LYS:HG2	2.03	0.40
4:D:75:ILE:HG21	4:D:169:ILE:HD11	2.03	0.40
4:D:305:HIS:HB2	4:D:326:LEU:HD13	2.03	0.40
6:F:163:THR:HB	6:F:166:THR:HG1	1.87	0.40
6:G:227:ILE:HG23	6:G:237:PHE:CE1	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:640:LEU:HG	1:A:643:LEU:HD23	2.03	0.40
1:A:690:THR:HG22	1:A:735:TYR:HB3	2.03	0.40
2:B:78:LEU:HB2	2:B:113:LEU:HD23	2.03	0.40
3:C:215:CYS:SG	3:C:232:ILE:HD13	2.61	0.40
3:C:239:PRO:HB3	3:C:272:LEU:HD13	2.03	0.40
4:D:33:PRO:HD2	4:D:73:SER:HB3	2.03	0.40
6:G:108:LYS:HA	6:G:108:LYS:HD3	1.81	0.40
1:A:389:ASN:OD1	1:A:439:GLN:NE2	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	484/861 (56%)	474 (98%)	10 (2%)	0	100	100
2	B	317/323 (98%)	308 (97%)	9 (3%)	0	100	100
3	C	325/340 (96%)	316 (97%)	9 (3%)	0	100	100
4	D	338/353 (96%)	333 (98%)	5 (2%)	0	100	100
5	E	334/354 (94%)	330 (99%)	4 (1%)	0	100	100
6	F	253/264 (96%)	247 (98%)	6 (2%)	0	100	100
6	G	250/264 (95%)	247 (99%)	3 (1%)	0	100	100
6	H	251/264 (95%)	242 (96%)	9 (4%)	0	100	100
All	All	2552/3023 (84%)	2497 (98%)	55 (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	422/742 (57%)	418 (99%)	4 (1%)	70	78
2	B	280/283 (99%)	278 (99%)	2 (1%)	76	80
3	C	284/296 (96%)	281 (99%)	3 (1%)	65	76
4	D	301/312 (96%)	298 (99%)	3 (1%)	68	76
5	E	310/324 (96%)	307 (99%)	3 (1%)	68	76
6	F	230/237 (97%)	226 (98%)	4 (2%)	53	71
6	G	229/237 (97%)	227 (99%)	2 (1%)	70	78
6	H	228/237 (96%)	224 (98%)	4 (2%)	51	70
All	All	2284/2668 (86%)	2259 (99%)	25 (1%)	63	76

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	320	LEU
1	A	386	THR
1	A	389	ASN
1	A	714	LEU
2	B	7	LEU
2	B	303	VAL
3	C	75	TYR
3	C	106	ILE
3	C	123	THR
4	D	40	THR
4	D	116	LEU
4	D	252	THR
5	E	5	VAL
5	E	68	LEU
5	E	181	ILE
6	F	19	PHE
6	F	137	LEU
6	F	163	THR
6	F	166	THR

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Mol	Chain	Res	Type
6	G	163	THR
6	G	243	SER
6	H	75	LEU
6	H	186	VAL
6	H	193	THR
6	H	241	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	364	HIS
1	A	368	GLN
1	A	398	ASN
1	A	415	ASN
1	A	546	ASN
1	A	571	ASN
1	A	575	ASN
1	A	608	GLN
2	B	18	GLN
2	B	60	HIS
2	B	92	GLN
2	B	148	ASN
2	B	204	GLN
2	B	216	HIS
2	B	276	ASN
2	B	320	ASN
3	C	180	ASN
3	C	291	GLN
3	C	317	GLN
4	D	96	ASN
4	D	124	HIS
4	D	128	ASN
4	D	236	GLN
4	D	288	ASN
4	D	303	GLN
4	D	345	ASN
4	D	352	GLN
5	E	80	ASN
5	E	86	ASN
5	E	118	GLN
5	E	144	ASN
5	E	208	ASN

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Mol	Chain	Res	Type
5	E	244	ASN
5	E	292	HIS
6	F	27	ASN
6	F	83	ASN
6	G	24	GLN
6	G	38	GLN
6	G	201	GLN
6	G	238	GLN
6	H	27	ASN
6	H	29	GLN
6	H	64	HIS
6	H	84	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 for Mogul analysis.

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
9	AGS	C	401	10	32,33,33	0.45	1 (3%)	45,52,52	0.66	1 (2%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	AGS	D	401	10	32,33,33	0.47	1 (3%)	45,52,52	0.64	1 (2%)
11	ADP	E	401	-	28,29,29	1.42	4 (14%)	43,45,45	1.84	9 (20%)
9	AGS	A	901	10	32,33,33	0.46	1 (3%)	45,52,52	0.66	1 (2%)
9	AGS	B	401	10	32,33,33	0.46	1 (3%)	45,52,52	0.66	1 (2%)

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
9	AGS	C	401	10	-	2/21/38/38	0/3/3/3
9	AGS	D	401	10	-	2/21/38/38	0/3/3/3
11	ADP	E	401	-	-	0/16/32/32	0/3/3/3
9	AGS	A	901	10	-	0/21/38/38	0/3/3/3
9	AGS	B	401	10	-	5/21/38/38	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	E	401	ADP	C5-C4	4.71	1.47	1.39
11	E	401	ADP	C5-C6	2.75	1.48	1.41
11	E	401	ADP	C8-N7	2.39	1.36	1.31
11	E	401	ADP	C5-N7	-2.25	1.35	1.39
9	A	901	AGS	PG-S1G	2.16	1.95	1.90
9	D	401	AGS	PG-S1G	2.16	1.95	1.90
9	B	401	AGS	PG-S1G	2.14	1.95	1.90
9	C	401	AGS	PG-S1G	2.12	1.95	1.90

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	E	401	ADP	C5-C4-N3	-5.86	118.64	126.72
11	E	401	ADP	N3-C4-N9	4.60	134.98	127.17
11	E	401	ADP	C2-N3-C4	3.71	120.90	111.83
9	A	901	AGS	PB-O3B-PG	-3.56	120.12	133.17
9	C	401	AGS	PB-O3B-PG	-3.56	120.16	133.17
9	D	401	AGS	PB-O3B-PG	-3.53	120.27	133.17
9	B	401	AGS	PB-O3B-PG	-3.52	120.27	133.17
11	E	401	ADP	C4-C5-N7	-3.51	106.57	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	E	401	ADP	N3-C2-N1	-3.21	123.72	128.58
11	E	401	ADP	C4-N9-C8	2.60	108.47	105.74
11	E	401	ADP	C5-N7-C8	2.56	107.47	103.45
11	E	401	ADP	C3'-C2'-C1'	2.30	105.81	101.46
11	E	401	ADP	C6-C5-N7	2.09	136.11	132.09

There are no chirality outliers.

All (9) torsion outliers are listed below:

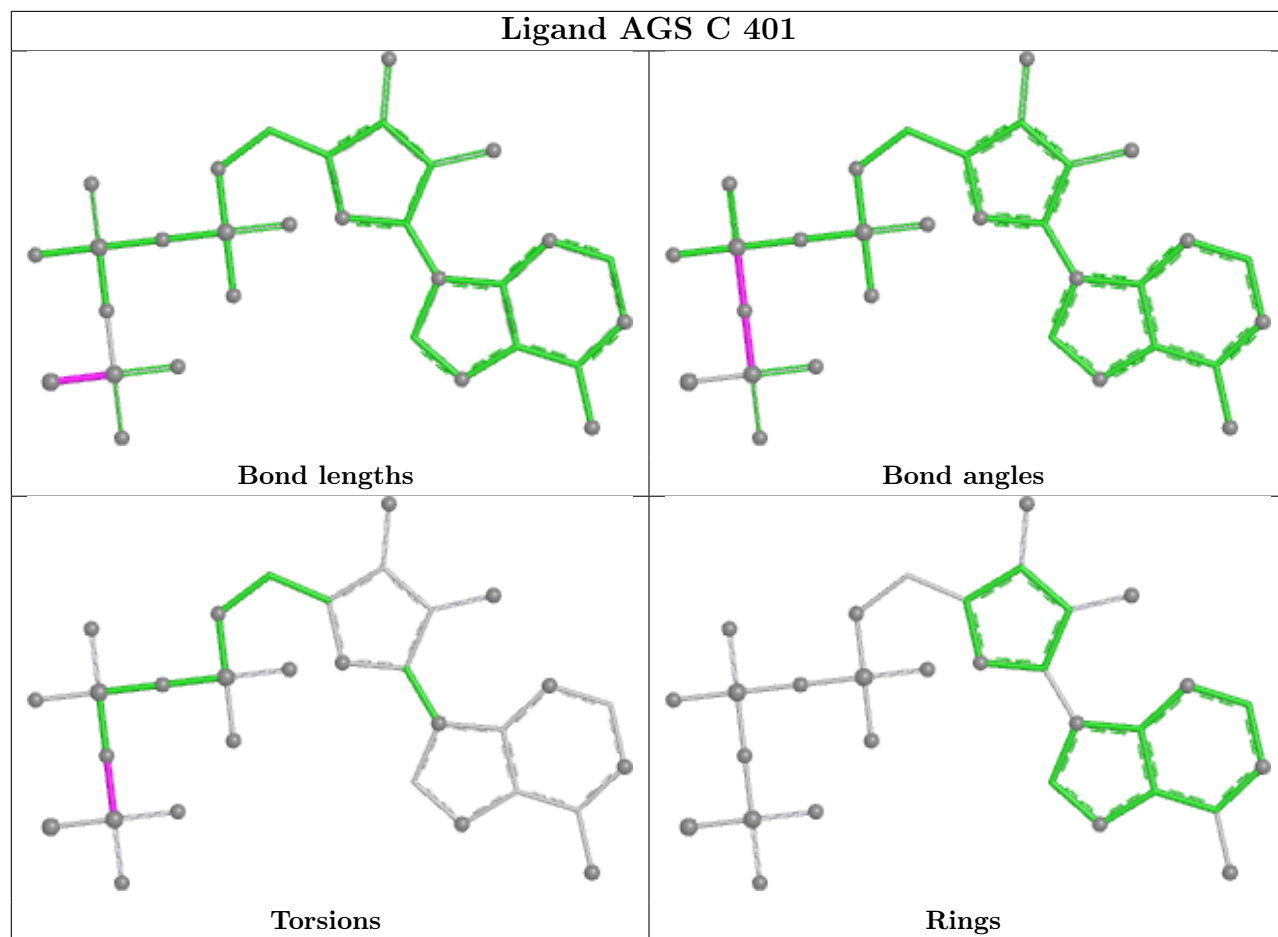
Mol	Chain	Res	Type	Atoms
9	B	401	AGS	PB-O3B-PG-O2G
9	B	401	AGS	PB-O3B-PG-O3G
9	B	401	AGS	C5'-O5'-PA-O1A
9	C	401	AGS	PB-O3B-PG-O2G
9	C	401	AGS	PB-O3B-PG-O3G
9	B	401	AGS	C5'-O5'-PA-O2A
9	B	401	AGS	C5'-O5'-PA-O3A
9	D	401	AGS	PA-O3A-PB-O1B
9	D	401	AGS	PA-O3A-PB-O2B

There are no ring outliers.

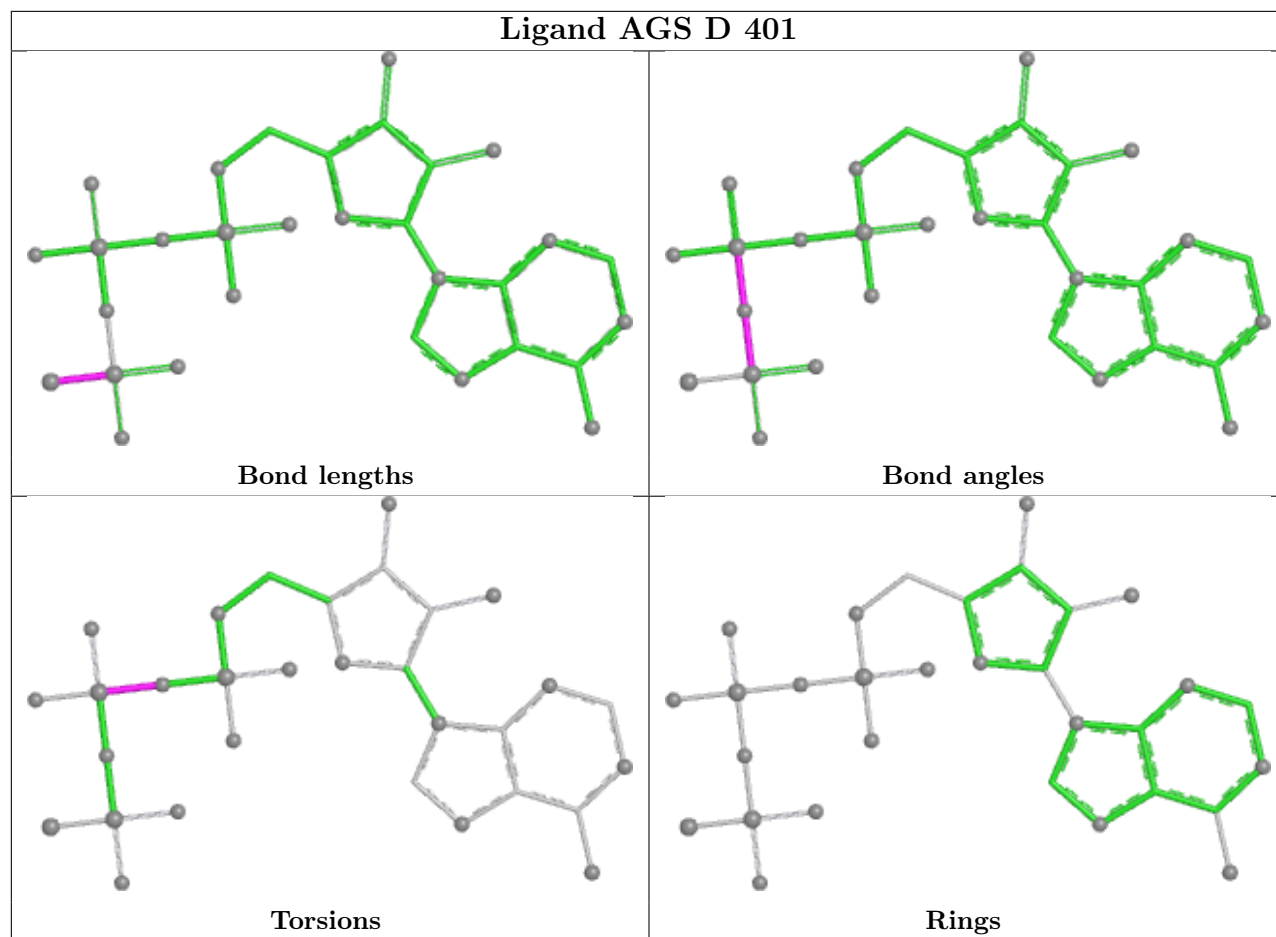
5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	C	401	AGS	2	0
9	D	401	AGS	2	0
11	E	401	ADP	2	0
9	A	901	AGS	3	0
9	B	401	AGS	3	0

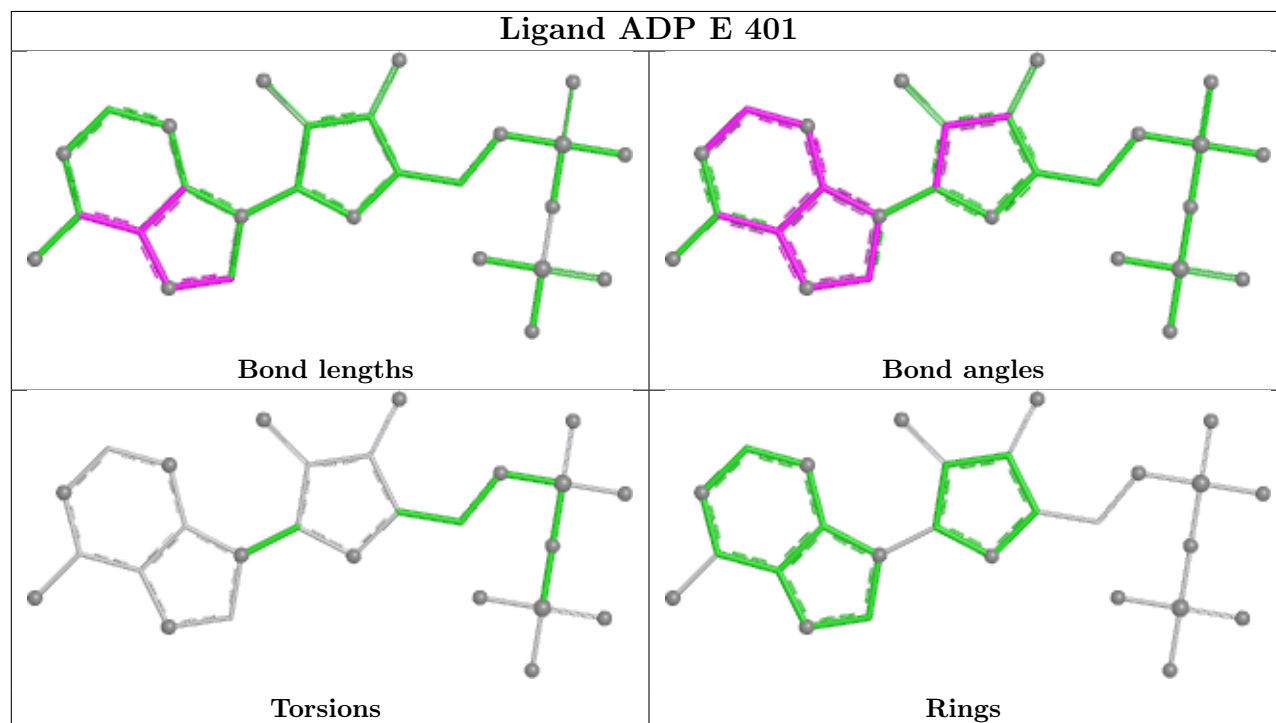
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

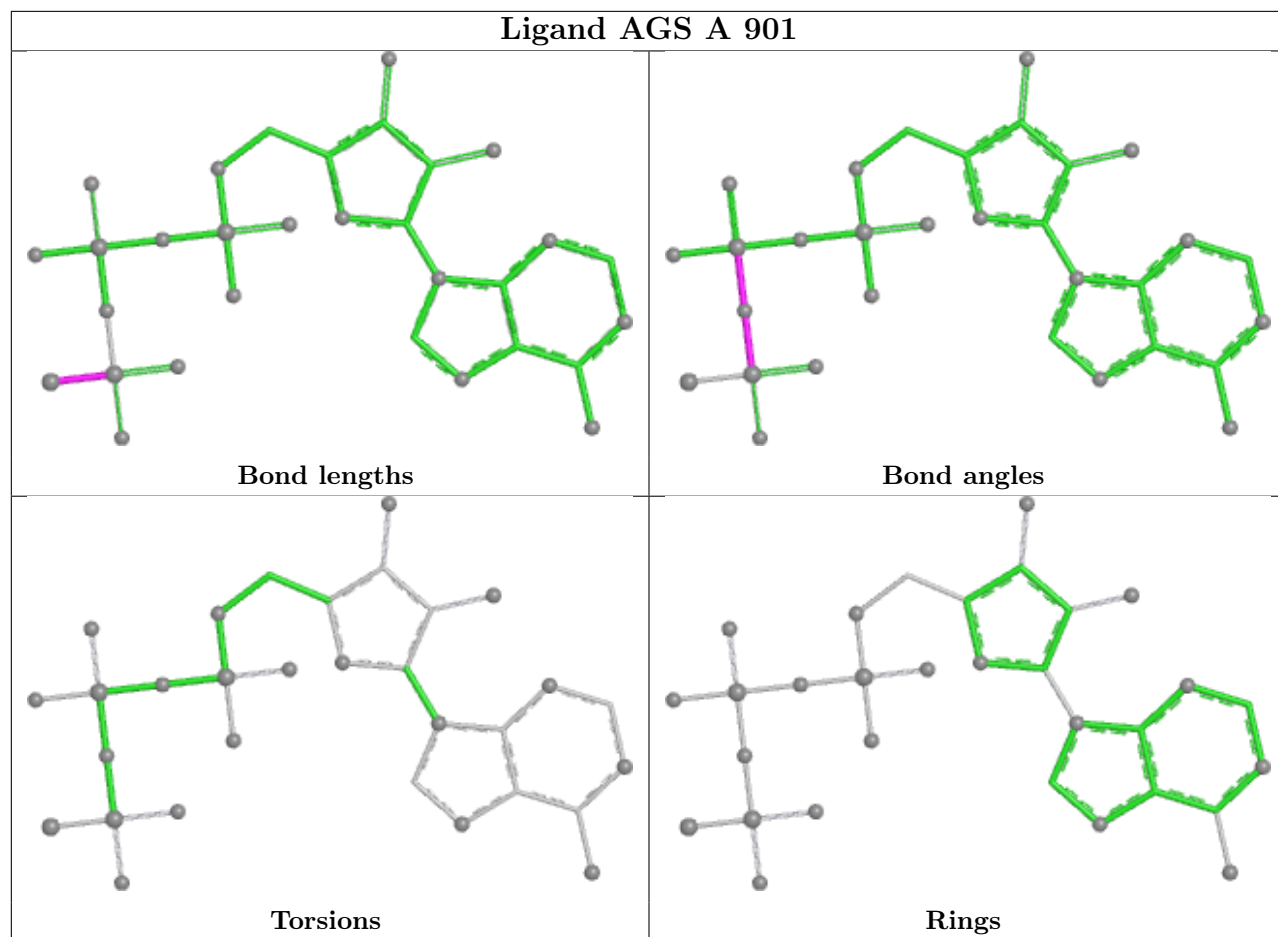


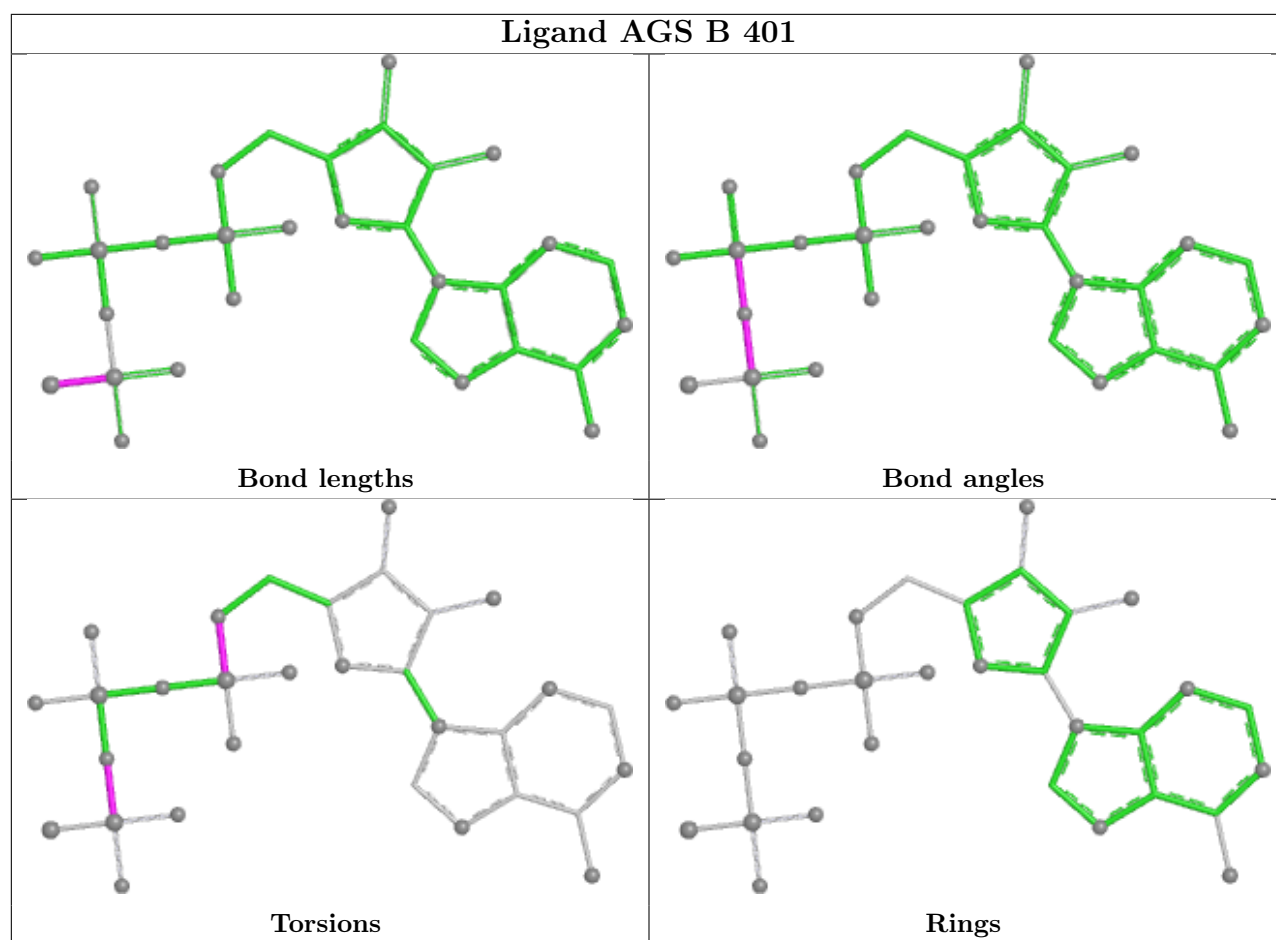
Ligand AGS D 401



Ligand ADP E 401







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.

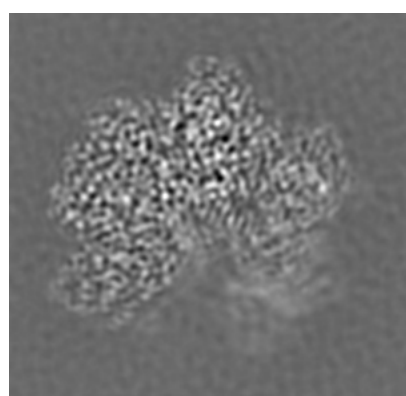
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-25617. These allow visual inspection of the internal detail of the map and identification of artifacts.

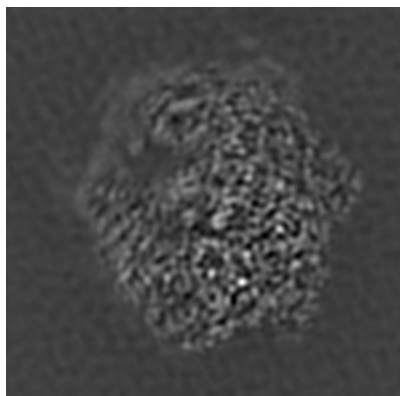
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

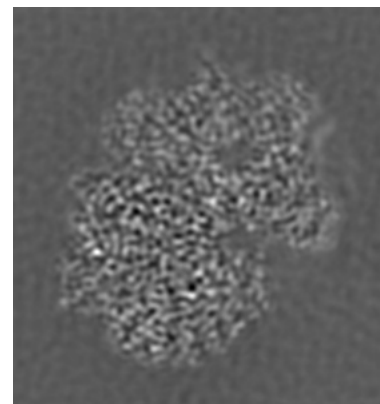
6.1.1 Primary map



X

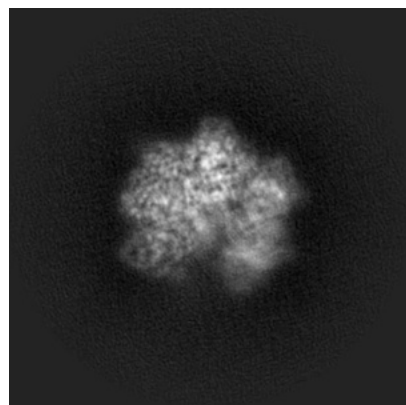


Y

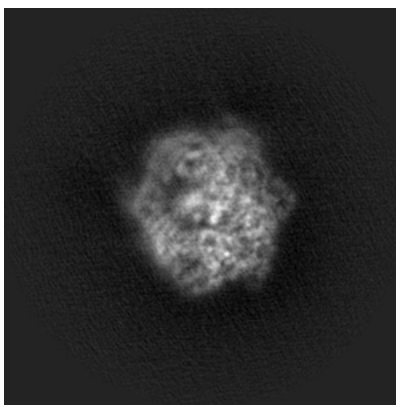


Z

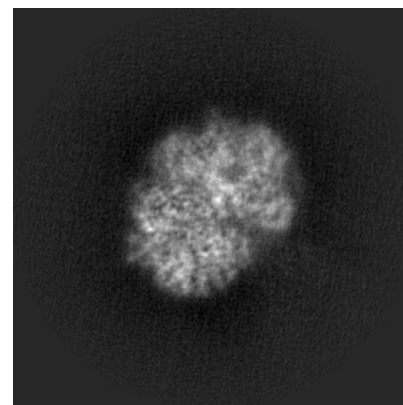
6.1.2 Raw map



X



Y

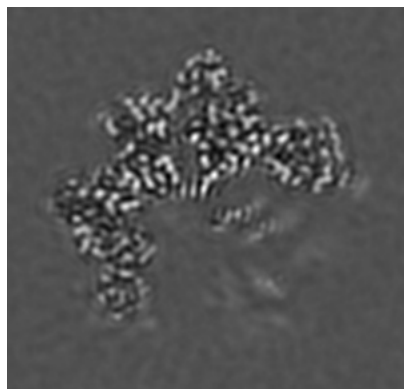


Z

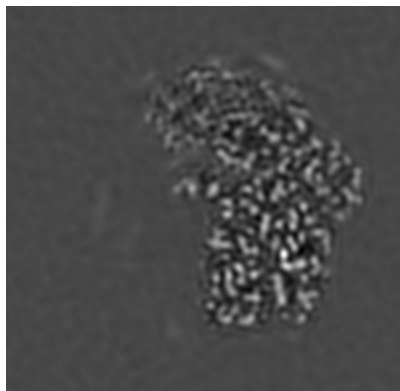
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

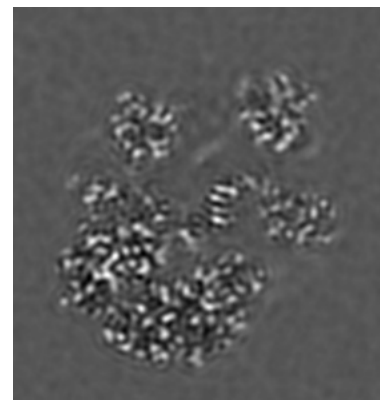
6.2.1 Primary map



X Index: 66

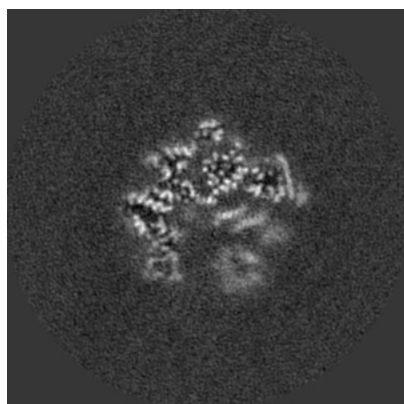


Y Index: 70

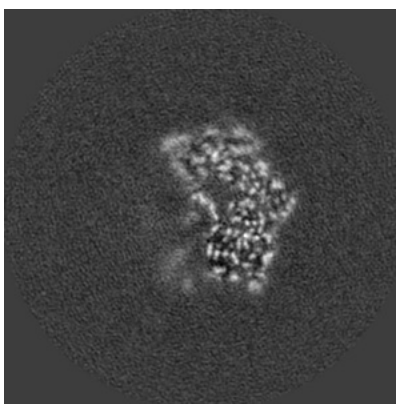


Z Index: 67

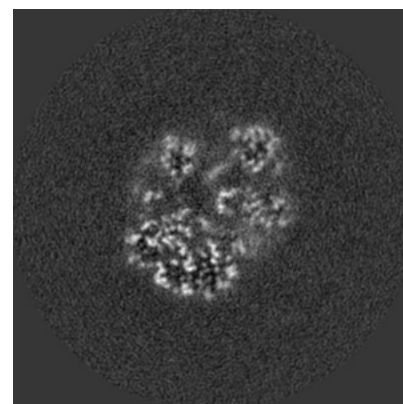
6.2.2 Raw map



X Index: 120



Y Index: 120

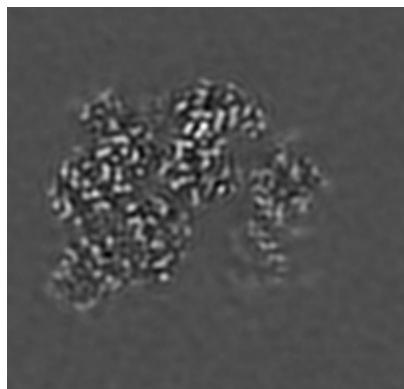


Z Index: 120

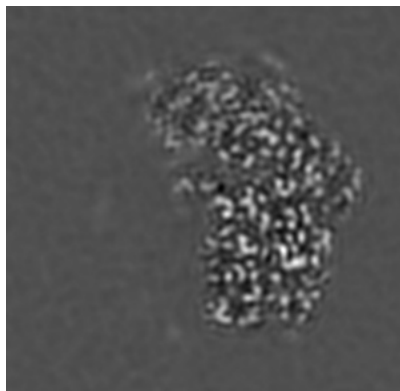
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

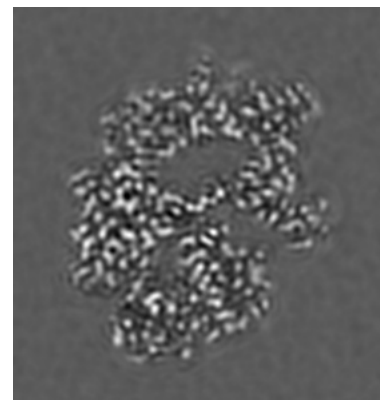
6.3.1 Primary map



X Index: 50

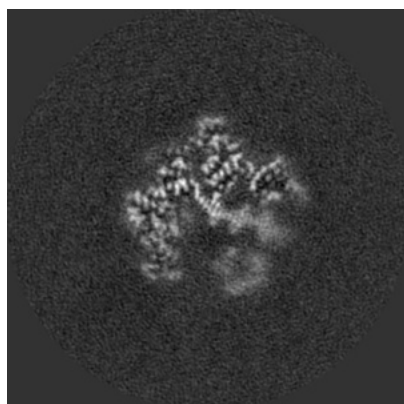


Y Index: 69

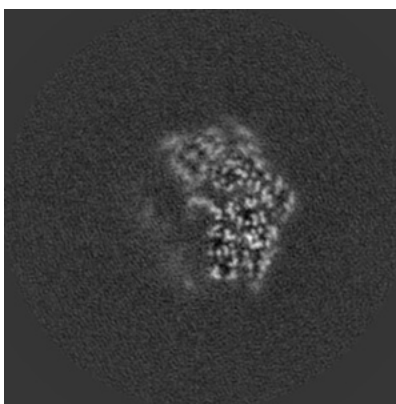


Z Index: 75

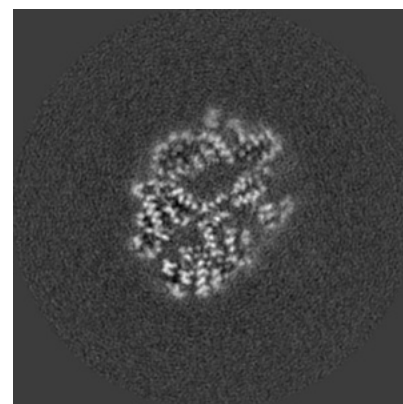
6.3.2 Raw map



X Index: 123



Y Index: 122

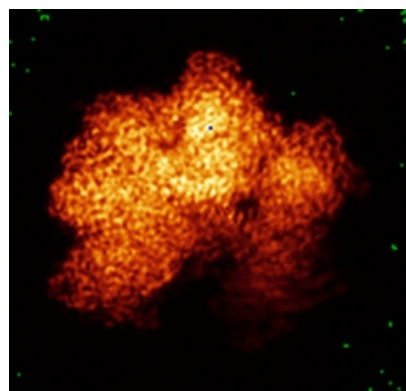


Z Index: 127

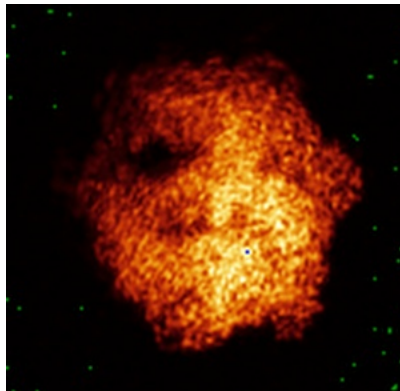
The images above show the largest variance slices of the map in three orthogonal directions.

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

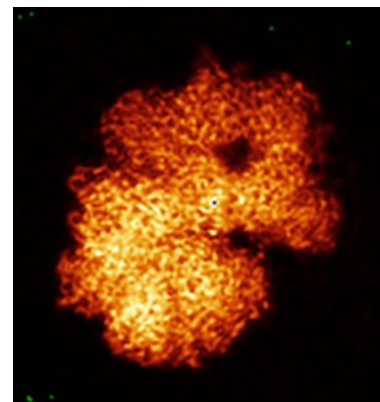
6.4.1 Primary map



X

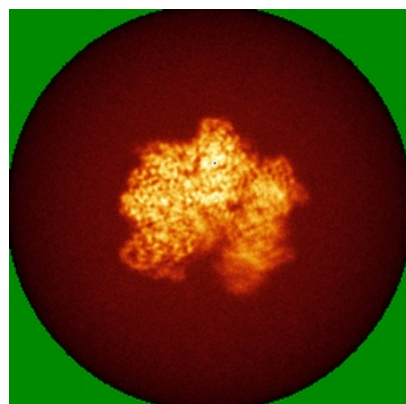


Y

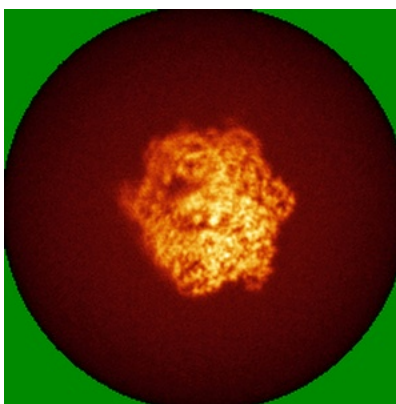


Z

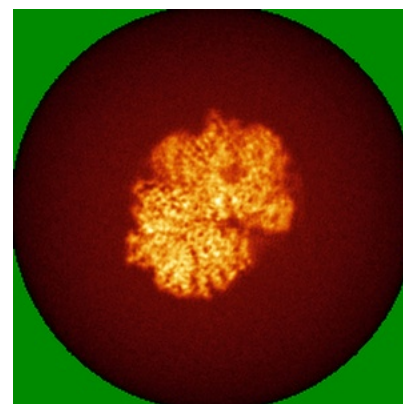
6.4.2 Raw map



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.

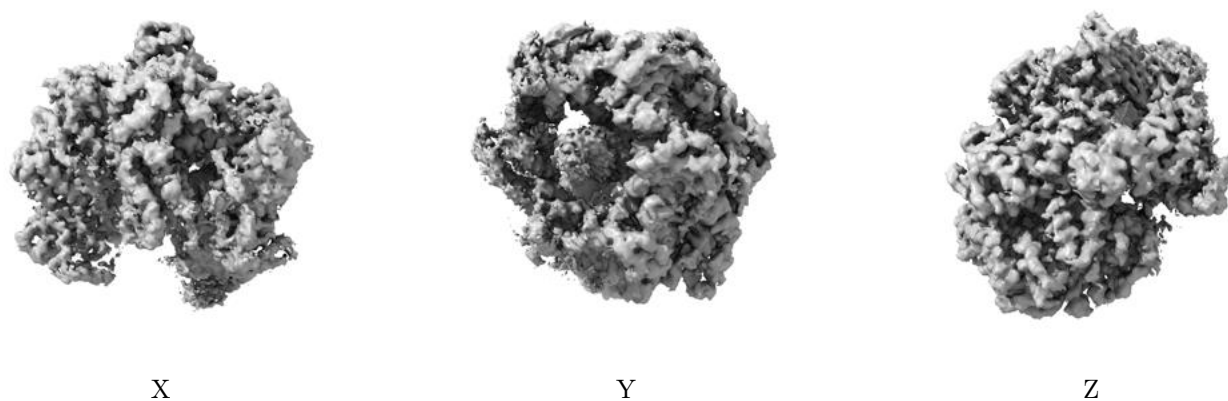
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.7. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

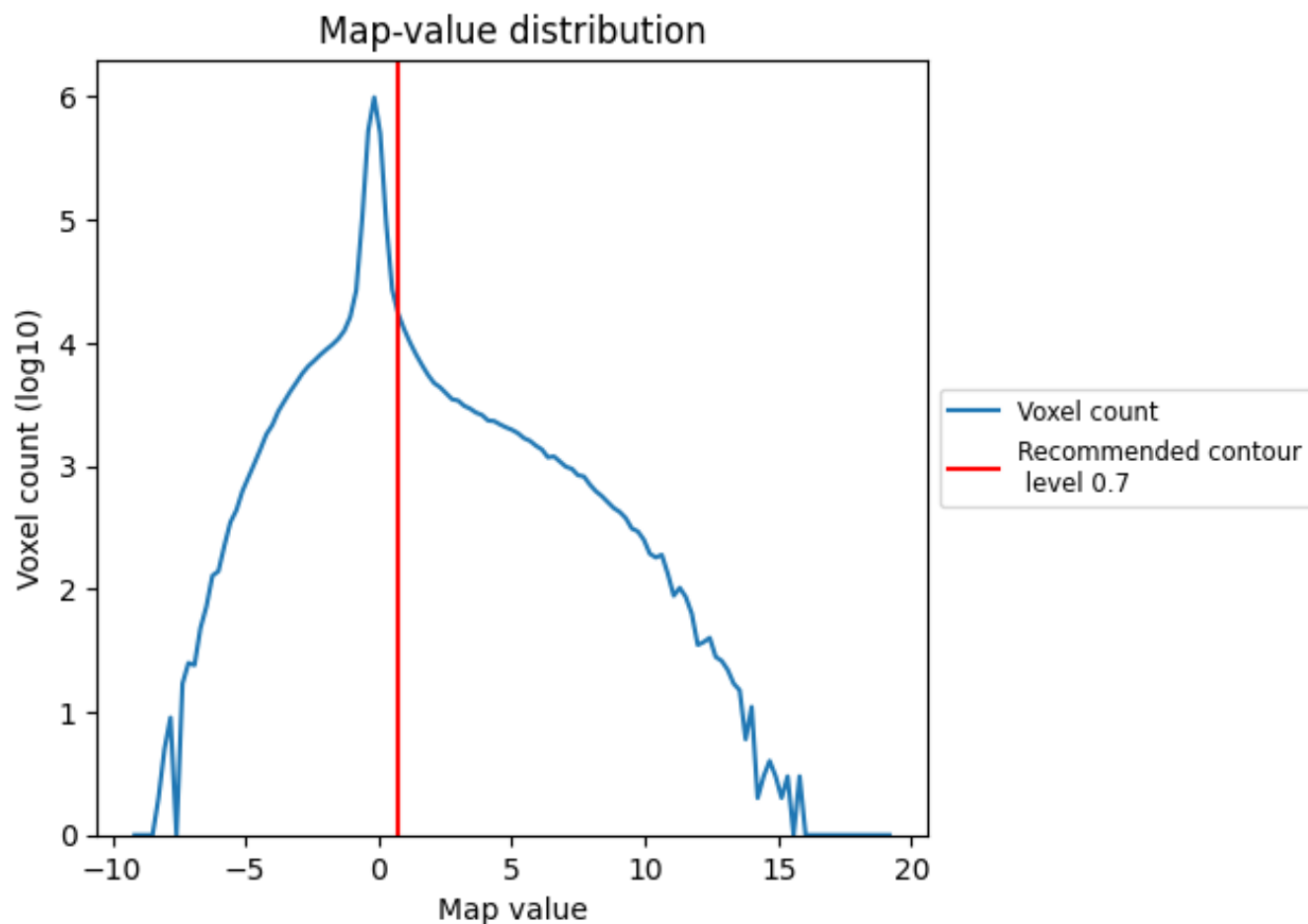
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

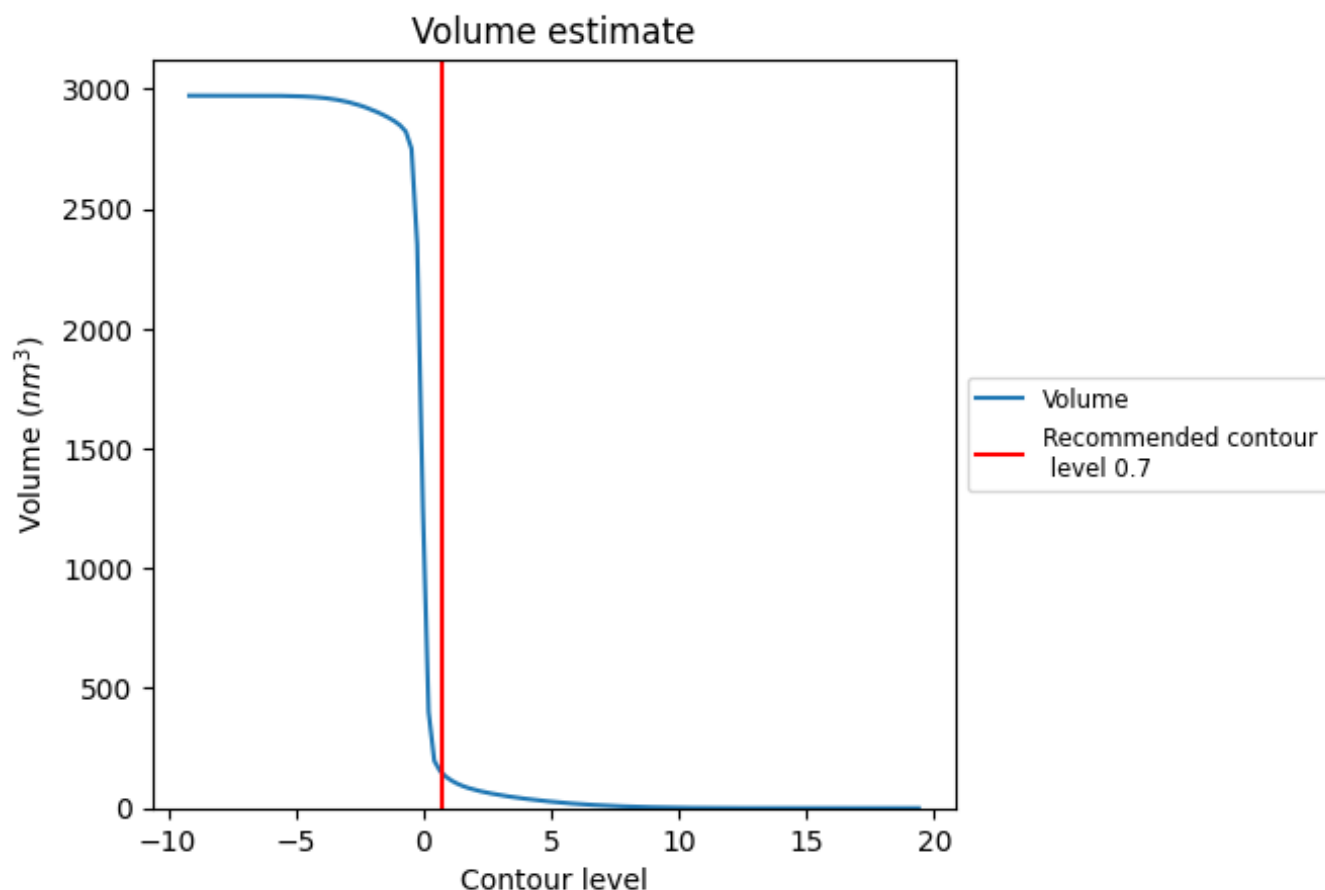
This section contains the results of statistical analysis of the map.

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 150 nm^3 ; this corresponds to an approximate mass of 136 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.

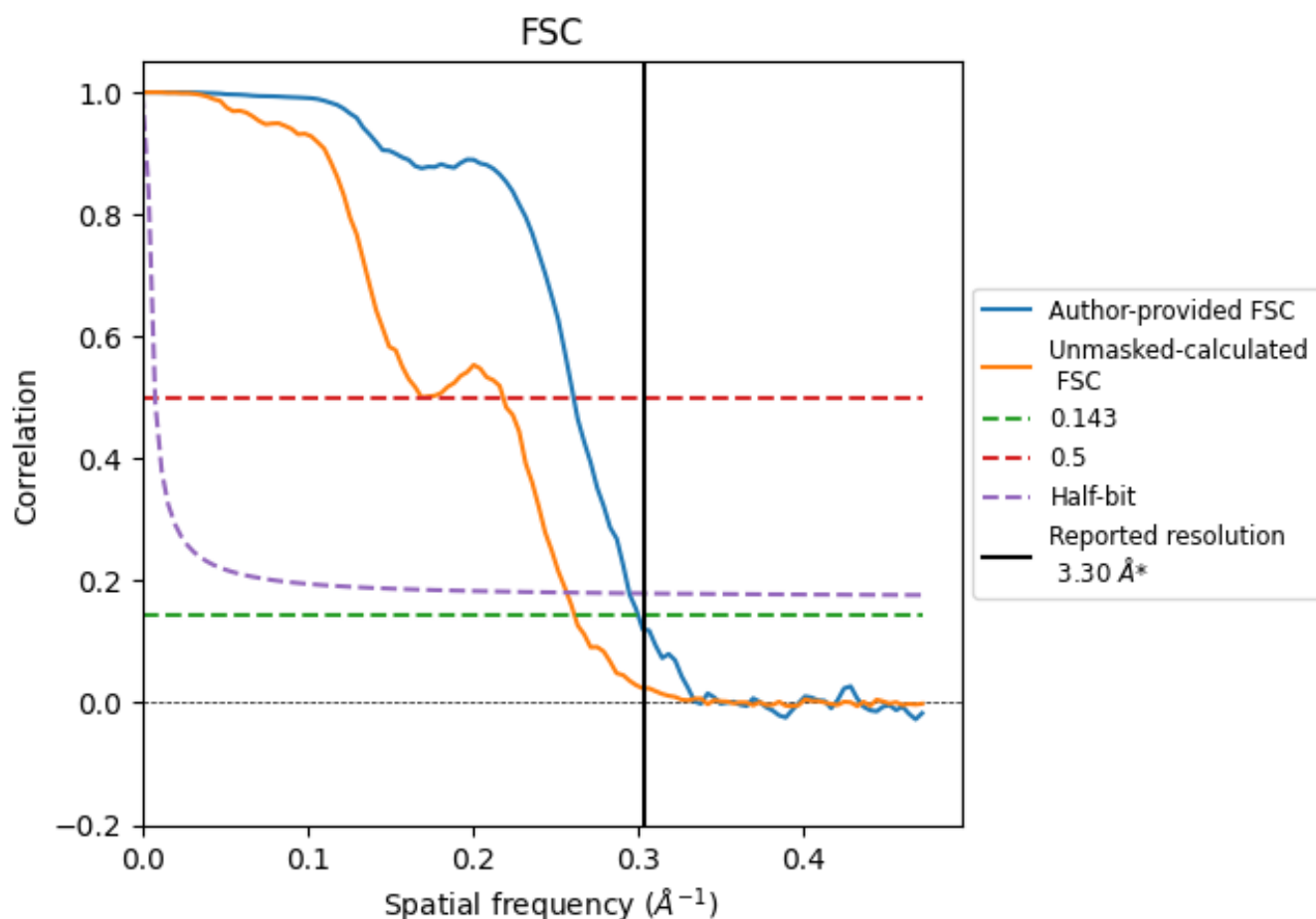
7.3 Rotationally averaged power spectrum [i](#)

This section was not generated. The rotationally averaged power spectrum is only generated for cubic maps.

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 $0.303 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

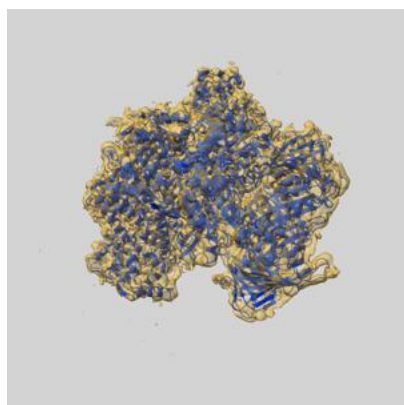
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.30	-	-
Author-provided FSC curve	3.33	3.83	3.39
Unmasked-calculated*	3.82	4.58	3.90

*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 3.82 differs from the reported value 3.3 by more than 10 %

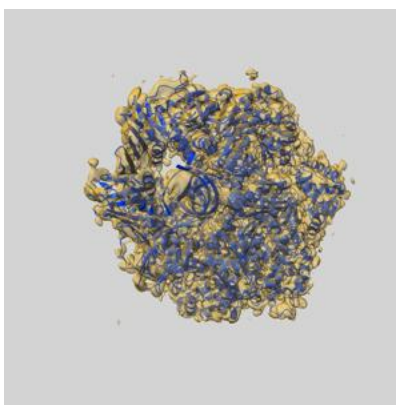
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-25617 and PDB model 7TID. Per-residue inclusion information can be found in [section 3](#) on [page 8](#).

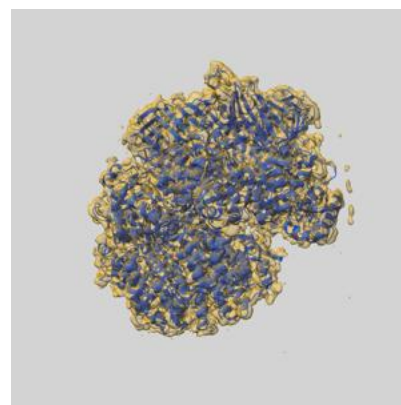
9.1 Map-model overlay [i](#)



X



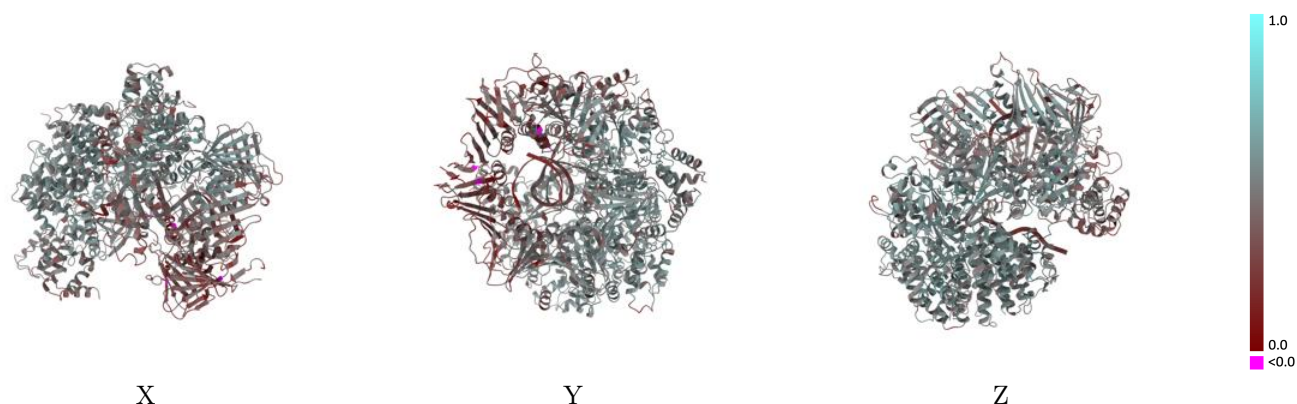
Y



Z

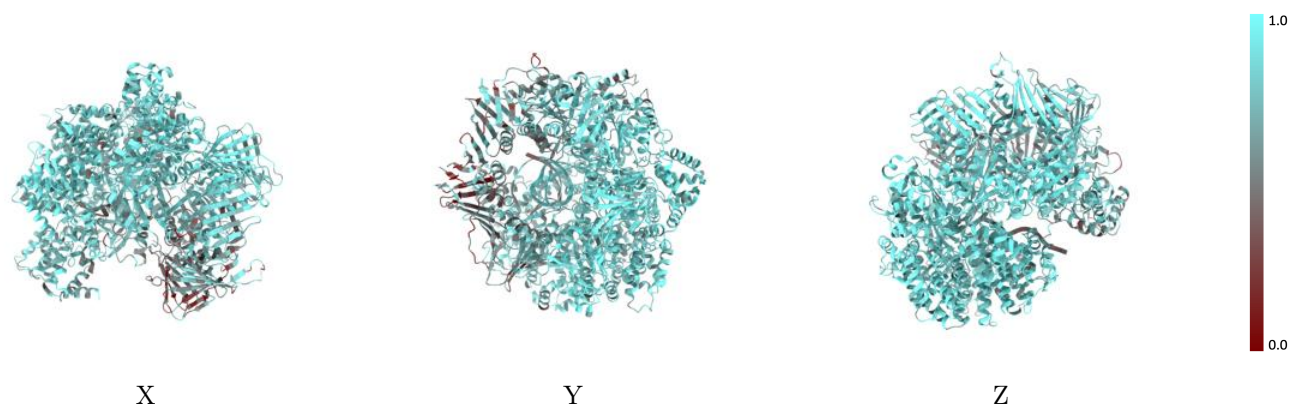
The images above show the 3D surface view of the map at the recommended contour level 0.7 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



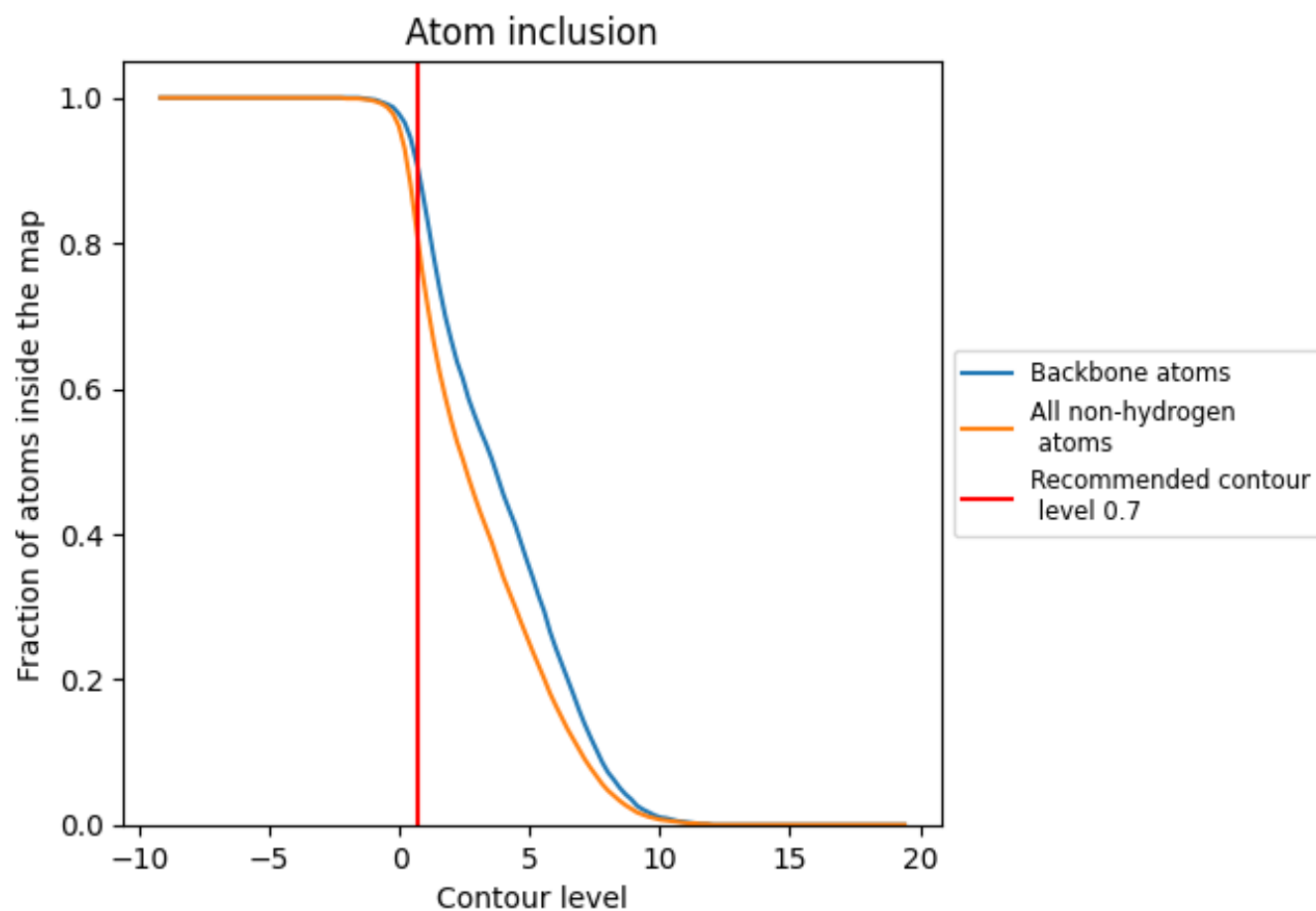
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.7).

9.4 Atom inclusion [i](#)



At the recommended contour level, 91% of all backbone atoms, 81% 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 (0.7) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8080	 0.4630
A	 0.7620	 0.4290
B	 0.8890	 0.5160
C	 0.8990	 0.5320
D	 0.9110	 0.5410
E	 0.8620	 0.4990
F	 0.8310	 0.4710
G	 0.7280	 0.4000
H	 0.5410	 0.3240
I	 0.6910	 0.3170
J	 0.8470	 0.3920

